Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-XLB

Level ID	File ID
A	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R022.D
B	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R023.D
C	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R024.D
D	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R025.D
E	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R026.D
F	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R027.D
G	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R028.D
H	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R029.D
I	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R030.D
J	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R031.D
K	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R032.D
L	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R033.D
M	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R034.D
N	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R035.D
O	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R036.D
P	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R037.D
Q	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R038.D
R	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R039.D
S	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R040.D
T	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R041.D
U	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R042.D
V	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R043.D
W	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R044.D
X	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R045.D
Y	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R046.D
Z	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R047.D
AA	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R048.D
AB	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R049.D
AC	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R050.D
AD	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R051.D

Analyte Name	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF
Decachlorobiphenyl	A	2.5	5660	B	5.0	5610	C	50	5290	D	100	5190	E	200	4790
	F	500	4260												
Aroclor 1016 {1}	A	25	198	B	50	195	C	500	190	D	1000	188	E	2000	173
	F	5000	155												
Aroclor 1016 {2}	A	25	375	B	50	377	C	500	322	D	1000	320	E	2000	291
	F	5000	255												
Aroclor 1016 {3}	A	25	148	B	50	139	C	500	173	D	1000	176	E	2000	163
	F	5000	148												
Aroclor 1016 {4}	A	25	142	B	50	136	C	500	140	D	1000	137	E	2000	124
	F	5000	105												
Aroclor 1016 {5}	A	25	152	B	50	152	C	500	159	D	1000	156	E	2000	139
	F	5000	125												
Aroclor 1260 {1}	A	25	294	B	50	311	C	500	290	D	1000	281	E	2000	257
	F	5000	232												
Aroclor 1260 {2}	A	25	320	B	50	344	C	500	308	D	1000	296	E	2000	270
	F	5000	243												
Aroclor 1260 {3}	A	25	392	B	50	383	C	500	383	D	1000	382	E	2000	358
	F	5000	334												
Aroclor 1260 {4}	A	25	188	B	50	237	C	500	253	D	1000	256	E	2000	245
	F	5000	234												

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-XLB

Analyte Name	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF
Aroclor 1260 {5}	A	25	456	B	50	453	C	500	449	D	1000	450	E	2000	428
	F	5000	401												

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-XLB

Analyte Name	Compound Type	Calibration Evaluation				
		Fit Type	Eval.	Eval. Result	Q	Control Criteria
Decachlorobiphenyl	SURR	AverageRF	% RSD	10.4		≤ 20
Aroclor 1016 {1}	MULTI	AverageRF	% RSD	9.0		≤ 20
Aroclor 1016 {2}	MULTI	AverageRF	% RSD	14.7		≤ 20
Aroclor 1016 {3}	MULTI	AverageRF	% RSD	9.6		≤ 20
Aroclor 1016 {4}	MULTI	AverageRF	% RSD	10.7		≤ 20
Aroclor 1016 {5}	MULTI	AverageRF	% RSD	8.7		≤ 20
Aroclor 1260 {1}	MULTI	AverageRF	% RSD	10.3		≤ 20
Aroclor 1260 {2}	MULTI	AverageRF	% RSD	12.1		≤ 20
Aroclor 1260 {3}	MULTI	AverageRF	% RSD	5.9		≤ 20
Aroclor 1260 {4}	MULTI	AverageRF	% RSD	10.5		≤ 20
Aroclor 1260 {5}	MULTI	AverageRF	% RSD	4.9		≤ 20

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010
Date Analyzed: 05/28/2010

Second Source Calibration Verification
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082

Calibration ID: CAL9510
Units: ng/mL

File ID:
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R052.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R053.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R054.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R055.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R056.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R057.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R058.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R059.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R060.D

Column ID: DB-XLB

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Aroclor 1016 {1}	1000	1100	183	195	6	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1000	323	326	1	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	158	181	15	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1000	131	137	5	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1100	147	163	10	NA	± 100 %	AverageRF
Aroclor 1016	1000	1100	NA	NA	NA	-7	± 15 %	NA
Aroclor 1260 {1}	1000	1000	277	280	1	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	980	297	292	-2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	910	372	339	-9	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	1200	235	282	20	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1200	440	506	15	NA	± 100 %	AverageRF
Aroclor 1260	1000	1100	NA	NA	NA	-5	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/08/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F003.D

Column ID: DB-35MS

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	97	5900	5740	-3	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	1000	142	146	3	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1000	286	300	5	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1000	243	245	1	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	970	207	201	-3	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	154	158	2	NA	± 100 %	AverageRF
Aroclor 1016	1000	1000	NA	NA	NA	2	± 15 %	NA
Aroclor 1260 {1}	1000	950	302	288	-5	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	930	370	346	-7	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	940	453	425	-6	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	990	433	429	-1	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	990	571	567	-1	NA	± 100 %	AverageRF
Aroclor 1260	1000	960	NA	NA	NA	-4	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/08/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R003.D

Column ID: DB-XLB

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	100	5130	5210	1	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	990	183	181	-1	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	960	323	311	-4	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	158	172	9	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1000	131	135	3	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	147	154	5	NA	± 100 %	AverageRF
Aroclor 1016	1000	1000	NA	NA	NA	2	± 15 %	NA
Aroclor 1260 {1}	1000	990	277	275	-1	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	980	297	290	-2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	1000	372	374	1	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	1100	235	252	7	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1000	440	442	1	NA	± 100 %	AverageRF
Aroclor 1260	1000	1000	NA	NA	NA	1	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/09/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F019.D

Column ID: DB-35MS

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	98	5900	5760	-2	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	1000	142	143	1	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1000	286	295	3	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	960	243	232	-4	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	970	207	201	-3	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	154	156	1	NA	± 100 %	AverageRF
Aroclor 1016	1000	990	NA	NA	NA	-1	± 15 %	NA
Aroclor 1260 {1}	1000	940	302	284	-6	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	930	370	344	-7	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	930	453	421	-7	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	990	433	427	-1	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1000	571	571	0	NA	± 100 %	AverageRF
Aroclor 1260	1000	960	NA	NA	NA	-4	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/09/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R019.D

Column ID: DB-XLB

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	100	5130	5190	1	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	970	183	178	-3	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	950	323	306	-5	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	158	170	7	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1000	131	133	2	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	147	152	3	NA	± 100 %	AverageRF
Aroclor 1016	1000	1000	NA	NA	NA	1	± 15 %	NA
Aroclor 1260 {1}	1000	990	277	274	-1	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	980	297	292	-2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	1000	372	374	0	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	1100	235	251	7	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1000	440	443	1	NA	± 100 %	AverageRF
Aroclor 1260	1000	1000	NA	NA	NA	1	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244

Analysis Run Log
Polychlorinated Biphenyls (PCBs)

Analysis Method: 8082

Analysis Lot: KWG1005547
Instrument ID: GC09.i
Column: DB-35MS

File ID	Sample Name	Lab Code	Date Analysis Started	Start Time	Q	Date Analysis Finished	Finish Time
0608F002.D	Instrument Blank	KWG1005547-1	6/8/2010	18:01		6/8/2010	18:01
0608F003.D	Continuing Calibration Verification	KWG1005547-2	6/8/2010	18:27		6/8/2010	18:27
0608F004.D	ZZZZZZ	ZZZZZZ	6/8/2010	18:53		6/8/2010	18:53
0608F005.D	Method Blank	KWG1005541-2	6/8/2010	19:19		6/8/2010	19:19
0608F006.D	Lab Control Sample	KWG1005541-1	6/8/2010	19:45		6/8/2010	19:45
0608F007.D	D-4-16	K1005244-003	6/8/2010	20:12		6/8/2010	20:12
0608F008.D	D-4-16MS	KWG1005541-3	6/8/2010	20:38		6/8/2010	20:38
0608F009.D	D-4-16DMS	KWG1005541-4	6/8/2010	21:04		6/8/2010	21:04
0608F010.D	ZZZZZZ	ZZZZZZ	6/8/2010	21:30		6/8/2010	21:30
0608F011.D	ZZZZZZ	ZZZZZZ	6/8/2010	21:56		6/8/2010	21:56
0608F012.D	ZZZZZZ	ZZZZZZ	6/8/2010	22:22		6/8/2010	22:22
0608F013.D	ZZZZZZ	ZZZZZZ	6/8/2010	22:48		6/8/2010	22:48
0608F014.D	ZZZZZZ	ZZZZZZ	6/8/2010	23:14		6/8/2010	23:14
0608F015.D	ZZZZZZ	ZZZZZZ	6/8/2010	23:40		6/8/2010	23:40
0608F016.D	ZZZZZZ	ZZZZZZ	6/9/2010	00:06		6/9/2010	00:06
0608F018.D	Instrument Blank	KWG1005547-3	6/9/2010	00:59		6/9/2010	00:59
0608F019.D	Continuing Calibration Verification	KWG1005547-4	6/9/2010	01:25		6/9/2010	01:25
0608F020.D	ZZZZZZ	ZZZZZZ	6/9/2010	01:51		6/9/2010	01:51
0608F021.D	ZZZZZZ	ZZZZZZ	6/9/2010	02:17		6/9/2010	02:17
0608F022.D	ZZZZZZ	ZZZZZZ	6/9/2010	02:43		6/9/2010	02:43
0608F024.D	ZZZZZZ	ZZZZZZ	6/9/2010	03:35		6/9/2010	03:35
0608F025.D	ZZZZZZ	ZZZZZZ	6/9/2010	04:02		6/9/2010	04:02
0608F026.D	ZZZZZZ	ZZZZZZ	6/9/2010	04:28		6/9/2010	04:28
0608F027.D	ZZZZZZ	ZZZZZZ	6/9/2010	04:54		6/9/2010	04:54
0608F028.D	ZZZZZZ	ZZZZZZ	6/9/2010	05:20		6/9/2010	05:20
0608F029.D	ZZZZZZ	ZZZZZZ	6/9/2010	05:46		6/9/2010	05:46
0608F031.D	Instrument Blank	KWG1005547-5	6/9/2010	06:38		6/9/2010	06:38
0608F032.D	Continuing Calibration Verification	KWG1005547-6	6/9/2010	07:04		6/9/2010	07:04

Results flagged with an asterisk (*) indicate the holding time was exceeded for the analysis

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Extracted: 06/02/2010

Extraction Prep Log
Polychlorinated Biphenyls (PCBs)

Extraction Method: EPA 3541
Analysis Method: 8082

Extraction Lot: KWG1005541
Level: Low

Sample Name	Lab Code	Date Collected	Date Received	Sample Amount	Final Volume	% Solids	Note
D-4-16	K1005244-003	05/19/10	05/21/10	22.55g	4mL	90.7	
Method Blank	KWG1005541-2	NA	NA	22.78g	4mL	NA	
D-4-16MS	KWG1005541-3	05/19/10	05/21/10	22.78g	4mL	90.7	
D-4-16DMS	KWG1005541-4	05/19/10	05/21/10	22.33g	4mL	90.7	
Lab Control Sample	KWG1005541-1	NA	NA	20.00g	4mL	NA	

Results flagged with an asterisk (*) indicate the holding time was exceeded for the analysis

COLUMBIA ANALYTICAL SERVICES, INC.

Confirmation Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010
Date Extracted: 06/02/2010

Polychlorinated Biphenyls (PCBs)

Sample Name: D-4-16
Lab Code: K1005244-003

Extraction Method: EPA 3541
Analysis Method: 8082

Units: ug/Kg
Basis: Dry
Level: Low

Analyte Name	MRL	MDL	Primary Result	Confirmation Result	RPD	Q	Dilution Factor	Date Analyzed
Aroclor 1248	9.8	2.1	14	14	0.0		1	06/08/10

Organic Analysis:
Polychlorinated Biphenyls (PCBs)

Validation Package

Organic Analysis:
Polychlorinated Biphenyls (PCBs)

Validation Package

QC Reports

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244

Surrogate Recovery Summary
Polychlorinated Biphenyls (PCBs)

Extraction Method: EPA 3541
Analysis Method: 8082

Units: PERCENT
Level: Low

Sample Name	Lab Code	Sur1
D-4-16	K1005244-003	106
Method Blank	KWG1005541-2	78
D-4-16MS	KWG1005541-3	103
D-4-16DMS	KWG1005541-4	87
Lab Control Sample	KWG1005541-1	93

Surrogate Recovery Control Limits (%)

Sur1 = Decachlorobiphenyl	35-133

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	06/02/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	06/08/2010

Matrix Spike/Duplicate Matrix Spike Summary
Polychlorinated Biphenyls (PCBs)

Sample Name:	D-4-16	**Units:**	ug/Kg
Lab Code:	K1005244-003	**Basis:**	Dry
Extraction Method:	EPA 3541	**Level:**	Low
Analysis Method:	8082	**Extraction Lot:**	KWG1005541

		D-4-16MS KWG1005541-3 Matrix Spike			D-4-16DMS KWG1005541-4 Duplicate Matrix Spike					
Analyte Name	**Sample Result**	**Result**	**Expected**	**%Rec**	**Result**	**Expected**	**%Rec**	**%Rec Limits**	**RPD**	**RPD Limit**
Aroclor 1016	ND	219	194	113	188	197	95	27-174	15	40
Aroclor 1260	ND	215	194	111	183	197	93	20-185	16	40

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	06/02/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	06/08/2010

Lab Control Spike Summary
Polychlorinated Biphenyls (PCBs)

Extraction Method:	EPA 3541	**Units:**	ug/Kg
Analysis Method:	8082	**Basis:**	Dry
		Level:	Low
		Extraction Lot:	KWG1005541

	Lab Control Sample KWG1005541-1 Lab Control Spike			
Analyte Name	**Result**	**Expected**	**%Rec**	**%Rec Limits**
Aroclor 1016	203	200	101	48-121
Aroclor 1260	204	200	102	53-129

Results flagged with an asterisk (*) indicate values outside control criteria.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	06/02/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	06/08/2010
		Time Analyzed:	19:19

Method Blank Summary
Polychlorinated Biphenyls (PCBs)

Sample Name:	Method Blank	**File ID:**	J:\GC09\DATA\060810.B\0608F005.D
Lab Code:	KWG1005541-2	**Instrument ID:**	GC09.i
Extraction Method:	EPA 3541	**Level:**	Low
Analysis Method:	8082	**Extraction Lot:**	KWG1005541

This Method Blank applies to the following analyses:

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed
Lab Control Sample	KWG1005541-1	J:\GC09\DATA\060810.B\0608F006.D	06/08/10	19:45
D-4-16	K1005244-003	J:\GC09\DATA\060810.B\0608F007.D	06/08/10	20:12
D-4-16MS	KWG1005541-3	J:\GC09\DATA\060810.B\0608F008.D	06/08/10	20:38
D-4-16DMS	KWG1005541-4	J:\GC09\DATA\060810.B\0608F009.D	06/08/10	21:04

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	06/02/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	06/08/2010
		Time Analyzed:	19:45

Lab Control Sample Summary
Polychlorinated Biphenyls (PCBs)

Sample Name:	Lab Control Sample	**File ID:**	J:\GC09\DATA\060810.B\0608F006.D
Lab Code:	KWG1005541-1	**Instrument ID:**	GC09.i
Extraction Method:	EPA 3541	**Level:**	Low
Analysis Method:	8082	**Extraction Lot:**	KWG1005541

This Lab Control Sample applies to the following analyses:

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed
Method Blank	KWG1005541-2	J:\GC09\DATA\060810.B\0608F005.D	06/08/10	19:19
D-4-16	K1005244-003	J:\GC09\DATA\060810.B\0608F007.D	06/08/10	20:12
D-4-16MS	KWG1005541-3	J:\GC09\DATA\060810.B\0608F008.D	06/08/10	20:38
D-4-16DMS	KWG1005541-4	J:\GC09\DATA\060810.B\0608F009.D	06/08/10	21:04

Organic Analysis:
Polychlorinated Biphenyls (PCBs)

Validation Package

Raw Data

COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Polychlorinated Biphenyls (PCBs)

Sample Name: D-4-16
Lab Code: K1005244-003

Extraction Method: EPA 3541
Analysis Method: 8082

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Aroclor 1016	ND	U	9.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1221	ND	U	20	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1232	ND	U	9.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1242	ND	U	9.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1248	**14**		9.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1254	ND	U	9.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1260	ND	U	9.8	2.1	1	06/02/10	06/08/10	KWG1005541	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Decachlorobiphenyl	106	35-133	06/08/10	Acceptable

Comments:

Exception Report

Data File: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F007.D
Lab ID: K1005244-003
RunType: SMPL
Matrix: MISC. SOLID

Date Acquired: 06/08/2010 20:12
Date Quantitated: 06/09/2010 10:29
Batch ID: KWG1005547
Analysis Method: 8082
MethodJoinID: MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
Preparation Holding Time	NA	NA	NA	x	
Pre-Preparation Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Method Blank	NA	NA	NA	x	
MB Surrogate Recovery	NA	NA	NA	x	
Lab Control Spike	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature]

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R007.D	**Date Acquired:**	06/08/2010 20:12
Lab ID:	K1005244-003	**Date Quantitated:**	06/09/2010 10:30
RunType:	SMPL	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
Preparation Holding Time	NA	NA	NA	x	
Pre-Preparation Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Method Blank	NA	NA	NA	x	
MB Surrogate Recovery	NA	NA	NA	x	
Lab Control Spike	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: MWS

Secondary Review: 6/10/10

Quantitation Report

Bottle ID:		Tier:	V	Matrix:	MISC. SOLID
Prod Code:	8082 PCB_LL	Collect Date:	05/19/2010	Receive Date:	05/21/2010

Analysis Lot:	KWG1005547	Prep Lot:	KWG1005541	Report Group:	K1005244
Analysis Method:	8082	Prep Method:	EPA 3541		
Prep Ref:	917270	Prep Date:	06/02/2010		

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:	Polychlorinated Biphenyls (PCBs)	Report List ID:	LJ10990
MB Ref:	J:\GC09\DATA\060810.B\0608F005.D	Method ID:	MJ696
		Quant based on Report List	

Data File #1:	J:\GC09\DATA\060810.B\0608F007.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R007.D			Vial:	6
Acqu Date:	06/08/2010 20:12	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	SMPL			Soln Conc. Units:	ng/mL
Lab ID:	K1005244-003				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2			Rpt
Decachlorobiphenyl	17.39 0.00	19.02 -0.01	581513	541680	98.50	105.53			106OK
			%Recovery =		99OK	106OK	Limits =	35-133	

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1016			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1016 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1221 {1}			0d	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {2}			0d	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {3}			0d	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {4}			0d	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1232 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1242 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F007.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R007.D			Vial:	6
Acqu Date:	06/08/2010 20:12	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	SMPL			Soln Conc. Units:	ng/mL
Lab ID:	K1005244-003				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1242 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248			0	0	73.85	73.09	14	14	14
Aroclor 1248 {1}	6.52	7.60	30061	14096m	72.94	70.51	14	14	
Aroclor 1248 {2}	7.16	7.71	12031	14043m	57.33	70.66	11	14	
Aroclor 1248 {3}	7.82	7.99	18674	13494m	74.38	67.44	15	13	
Aroclor 1248 {4}	8.19	8.84	29364	31438m	84.52	82.85	17	16	
Aroclor 1248 {5}	8.29	8.94	33970	20337m	80.09	73.99	16	14	
Aroclor 1254			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1254 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1260 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	

The -/+ after Retention Time symbolize the direction of the RT shift

Prep Amount:	22.55 g	Dilution:	1.0
Prep Final Vol:	4 mL	Unit Factor:	1
Solids:	90.7 %		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F007.D
Report Date: 09-Jun-2010 10:29

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F007.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R007.D
Inj Date : 08-JUN-2010 20:12
Sample Info: K1005244-003 | D-4-16
Misc Info : KELSO\SEMIVOA GC\8082-PCB_LL\05244003.H | F=4
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

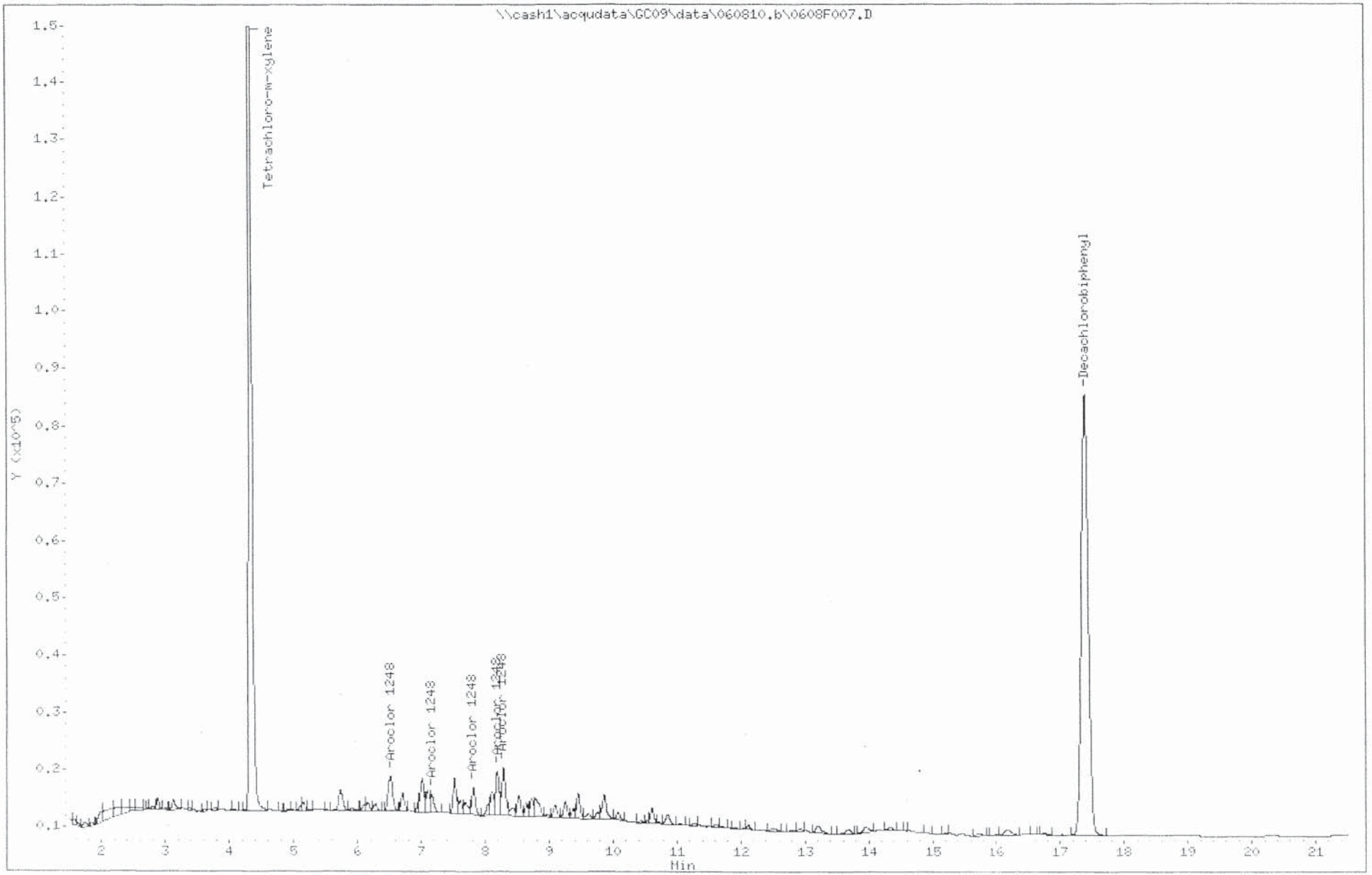
Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.333	5.017	651130	587572	91.6	99.8		100.00
Aroclor 1248	6.523	7.603	30061	14096	72.9	70.5	80.00- 120.00	100.00
	7.163	7.707	12031	14043	57.3	70.7	39.24- 58.85	40.02
	7.817	7.993	18674	13494	74.4	67.4	43.38- 65.07	62.12
	8.190	8.843	29364	31438	84.5	82.8	66.78- 100.17	97.68
	8.287	8.943	33970	20337	80.1	74.0	84.10- 126.14	113.00
			Average of Peak Amounts =		73.8	73.1		
[illegible]	17.587	19.017	581513	541680	98.5	106		100.00

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F007.D
Date : 08-JUN-2010 20:12
Client ID:
Sample Info: K1005244-003 | D-4-16

Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

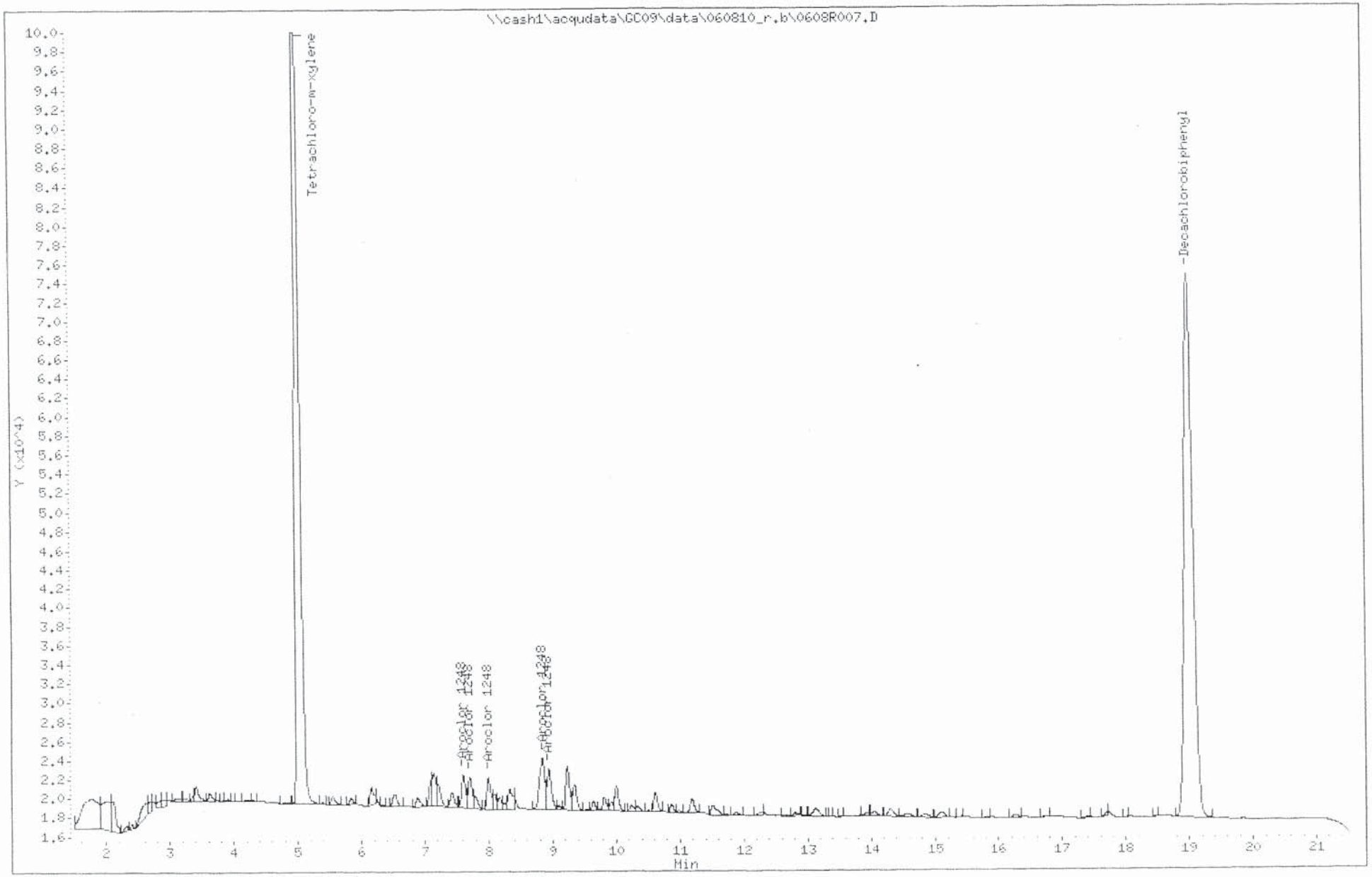


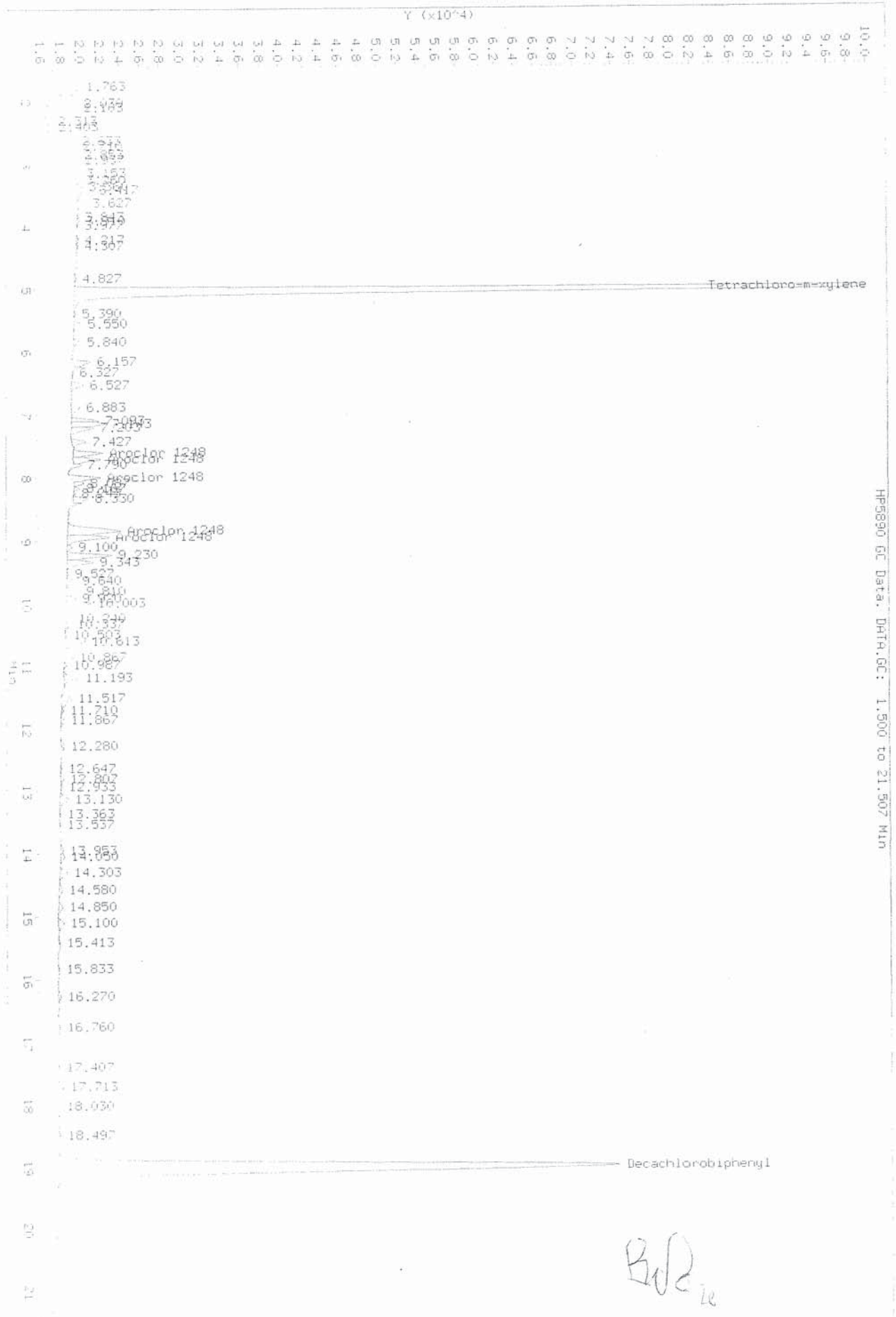
Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R007.D
Date : 08-JUN-2010 20:12
Client ID:
Sample Info: K1005244-003 | D-4-16

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB




Data File: \\cash1\acqdata\GC09\data\060810_r.b\0508R007.D
Injection Date: 08-JUN-2010 20:12
Instrument: GC09.1
Client Sample ID:
Y (x10^4)
1.763
4.827
Tetrachloro-m-xylene
5.390
5.550
5.840
6.157
6.327
6.527
6.883
7.427
Aroclor 1248
7.790
Aroclor 1248
Aroclor 1248
9.100
9.230
9.343
10.003
10.503
10.613
10.867
10.987
11.193
11.517
11.710
11.867
12.280
12.647
12.802
12.933
13.130
13.363
13.537
14.303
14.580
14.850
15.100
15.413
15.833
16.270
16.760
17.407
17.713
18.030
18.497
Decachlorobiphenyl
HP5890 GC Data. DATA.GC: 1.500 to 21.507 Min

Data File: \\crash\acqudata\GC09\data\060810_r.b\06080007.D
Injection Date: 08-JUN-2010 20:12
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.507 Min
Tetrachloro-m-xylene
Decachlorobiphenyl
Aroclor 1248
Min

COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Polychlorinated Biphenyls (PCBs)

Sample Name: Method Blank
Lab Code: KWG1005541-2

Extraction Method: EPA 3541
Analysis Method: 8082

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Aroclor 1016	ND	U	8.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1221	ND	U	18	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1232	ND	U	8.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1242	ND	U	8.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1248	ND	U	8.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1254	ND	U	8.8	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1260	ND	U	8.8	2.1	1	06/02/10	06/08/10	KWG1005541	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Decachlorobiphenyl	78	35-133	06/08/10	Acceptable

Comments:

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F005.D	**Date Acquired:**	06/08/2010 19:19
Lab ID:	KWG1005541-2	**Date Quantitated:**	06/09/2010 10:28
RunType:	MB	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: WWW

Secondary Review: 6/10/10

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R005.D	**Date Acquired:**	06/08/2010 19:19
Lab ID:	KWG1005541-2	**Date Quantitated:**	06/09/2010 10:29
RunType:	MB	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature] 6/10/10

Quantitation Report

Bottle ID:		Tier:		Matrix:	MISC_SOLID
Prod Code:	8082 PCB_LL	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:	KWG1005541	Report Group:	
Analysis Method:	8082	Prep Method:	EPA 3541		
Prep Ref:	917272	Prep Date:	06/02/2010		

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:		Method ID:	MJ696
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F005.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R005.D			Vial:	4
Acqu Date:	06/08/2010 19:19	Quant Date:	06/09/2010 10:28	Dilution:	1.0
Run Type:	MB			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-2				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	4.33	5.02	447608	384158	62.99	65.24			65OK
			%Recovery =		63OK	65OK	Limits =	10-135	
Decachlorobiphenyl	17.39	19.02 0.00	424178	398568	71.85	77.65			78OK
			%Recovery =		72OK	78OK	Limits =	35-133	

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1016			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1016 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1016 {5}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1221 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1232 {1}			0	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {2}			0	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {3}			0	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {4}			0	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1242 {1}			0	0	0.0000	0.0000	2.1U	2.1U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F005.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R005.D			Vial:	4
Acqu Date:	06/08/2010 19:19	Quant Date:	06/09/2010 10:28	Dilution:	1.0
Run Type:	MB			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-2				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds — Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1242 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {5}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1248 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {5}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1254 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {5}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1260 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260 {5}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1262 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {5}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1268 {1}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {2}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {3}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {4}			0	0	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {5}			0	0	0.0000	0.0000	2.1U	2.1U	

The -/+ after Retention Time symbolize the direction of the RT shift

Prep Amount:	22.78 g	Dilution:	1.0
Prep Final Vol:	4 mL	Unit Factor:	1
Solids:	%		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F005.D
Report Date: 09-Jun-2010 10:28

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F005.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R005.D
Inj Date : 08-JUN-2010 19:19
Sample Info: KWG1005541-2 | MB | 8082 PCB_LL | MISC. SOLID
Misc Info : SEMIVOA GC\W1005541\2-MB.H | F=4 D=1 A=22.78
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

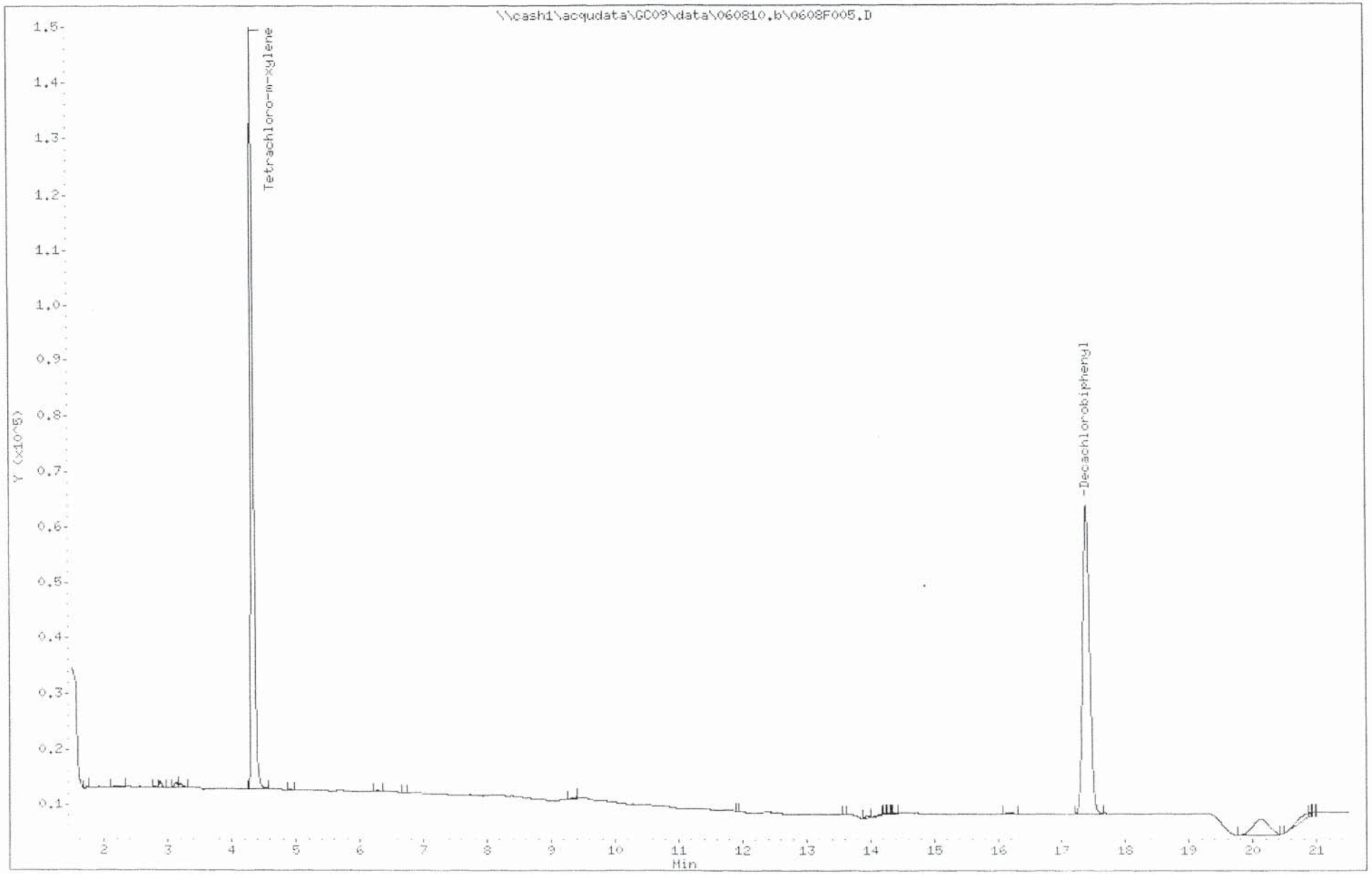
Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.333	5.017	447608	384158	63.0	65.2		100.00
Decachlorobiphenyl	17.390	19.020	424178	398568	71.8	77.6		100.00

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F005.D
Date : 08-JUN-2010 19:19
Client ID:
Sample Info: KWG1005541-2 | MB | 8082 PCB_LL | MISC. SOLID

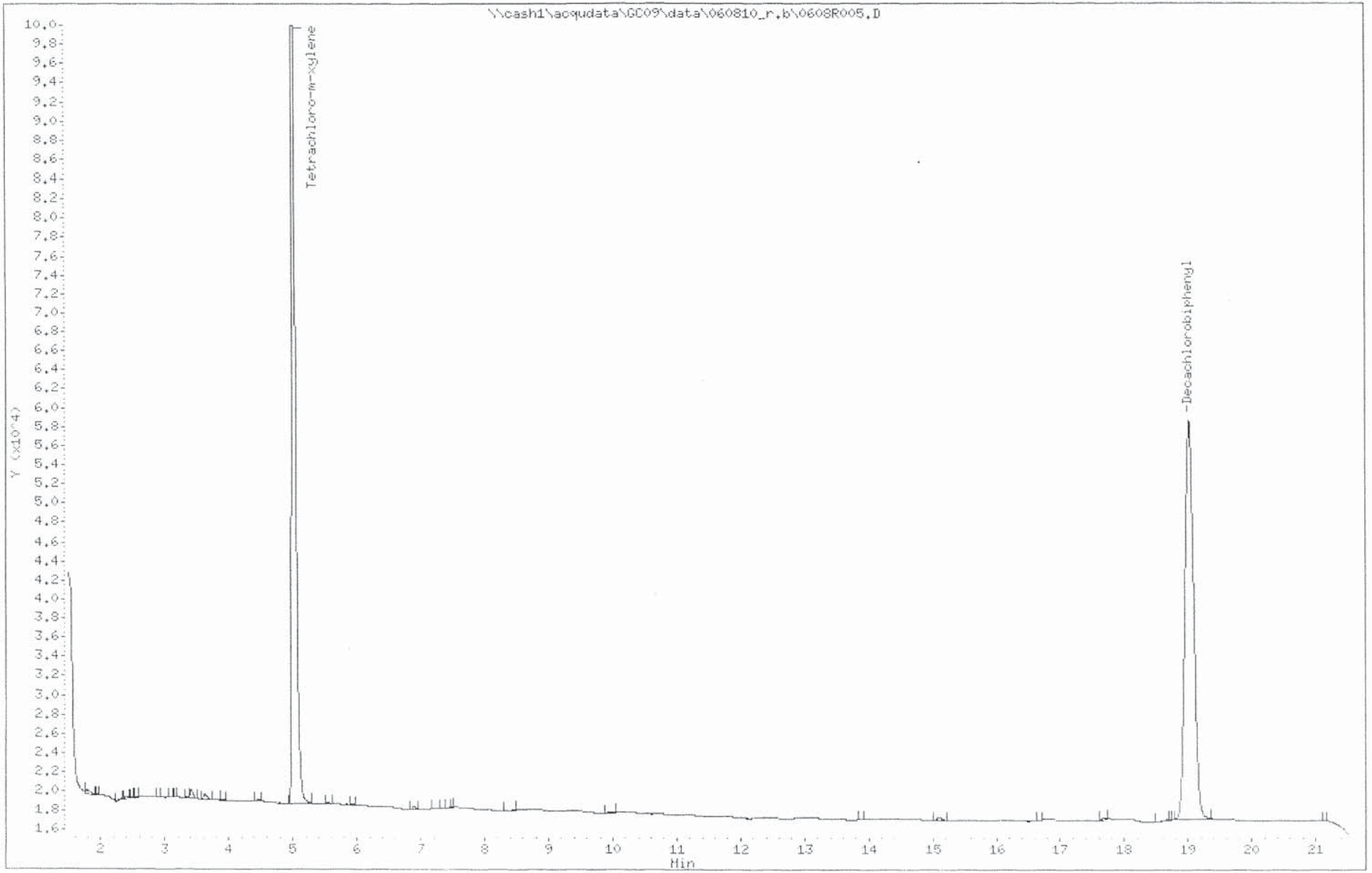
Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R005.D
Date : 08-JUN-2010 19:19
Client ID:
Sample Info: KWG1005541-2 | MB | 8082 PCB_LL | MISC. SOLID

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Polychlorinated Biphenyls (PCBs)

Sample Name: D-4-16MS
Lab Code: KWG1005541-3

Extraction Method: EPA 3541
Analysis Method: 8082

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Aroclor 1016	**219**		9.7	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1221	ND	U	20	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1232	ND	U	9.7	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1242	ND	U	9.7	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1248	ND	U	9.7	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1254	ND	U	9.7	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1260	**215**		9.7	2.1	1	06/02/10	06/08/10	KWG1005541	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Decachlorobiphenyl	103	35-133	06/08/10	Acceptable

Comments:

Exception Report

Data File: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F008.D
Lab ID: KWG1005541-3 -- K1005244-003MS
RunType: MS
Matrix: MISC. SOLID

Date Acquired: 06/08/2010 20:38
Date Quantitated: 06/09/2010 10:29
Batch ID: KWG1005547
Analysis Method: 8082
MethodJoinID: MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature]

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R008.D	**Date Acquired:**	06/08/2010 20:38
Lab ID:	KWG1005541-3 -- K1005244-003MS	**Date Quantitated:**	06/09/2010 10:30
RunType:	MS	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature]

Quantitation Report

Bottle ID:		Tier:		Matrix:	MISC. SOLID
Prod Code:	8082 PCB_LL	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:	KWG1005541	Report Group:	
Analysis Method:	8082	Prep Method:	EPA 3541		
Prep Ref:	917273	Prep Date:	06/02/2010		

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:	J:\GC09\DATA\060810.B\0608F005.D	Method ID:	MJ696
			Quant based on Method

Data File #1:	J:\GC09\DATA\060810.B\0608F008.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R008.D			Vial:	7
Acqu Date:	06/08/2010 20:38	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	MS			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-3 -- K1005244-003MS				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	4.33	5.02	642771	577056	90.46	97.99			98OK
			%Recovery =		90OK	98OK	Limits =	10-135	
Decachlorobiphenyl	17.39	19.02 -0.01	566222	527067	95.91	102.68			103OK
			%Recovery =		96OK	103OK	Limits =	35-133	

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1016			0	0	1,110	1,134	215	219	219
Aroclor 1016 {1}	6.16 0.00	6.53	149695	204802	1,057	1,118	205	216	
Aroclor 1016 {2}	6.71	7.14 0.00	311457	352426	1,088	1,090	211	211	
Aroclor 1016 {3}	7.01 0.00	7.43 0.00	246779	185890	1,014	1,177	196	228	
Aroclor 1016 {4}	7.52	7.60	227969	146599	1,100	1,122	213	217	
Aroclor 1016 {5}	7.82 0.00	7.71	199113	171398	1,291	1,162	250	225	
Aroclor 1221			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1221 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1232 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1242 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
P: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F008.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R008.D			Vial:	7
Acqu Date:	06/08/2010 20:38	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	MS			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-3 -- K1005244-003MS				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1242 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1248 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1254 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260			0	0	965.95	1,112	187	215	215
Aroclor 1260 {1}	9.47	10.34	281136	284961	929.88	1,027	180	199	
Aroclor 1260 {2}	9.75 0.00	10.87 0.00	346804	310728	937.88	1,047	182	203	
Aroclor 1260 {3}	10.08	11.52 0.00	419823	366027	927.10	983.65	179	190	
Aroclor 1260 {4}	10.85 0.00	12.28 0.00	388841	298377	897.59	1,267	174	245	
Aroclor 1260 {5}	12.12 0.00	13.11 0.00	649374	542898	1,137	1,235	220	239	
Aroclor 1262			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1262 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1268 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	

The -/+ after Retention Time symbolize the direction of the RT shift

Prep Amount:	22.78 g	Dilution:	1.0
Prep Final Vol:	4 mL	Unit Factor:	1
Solids:	90.7 %		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F008.D
Report Date: 09-Jun-2010 10:29

Laboratory Name

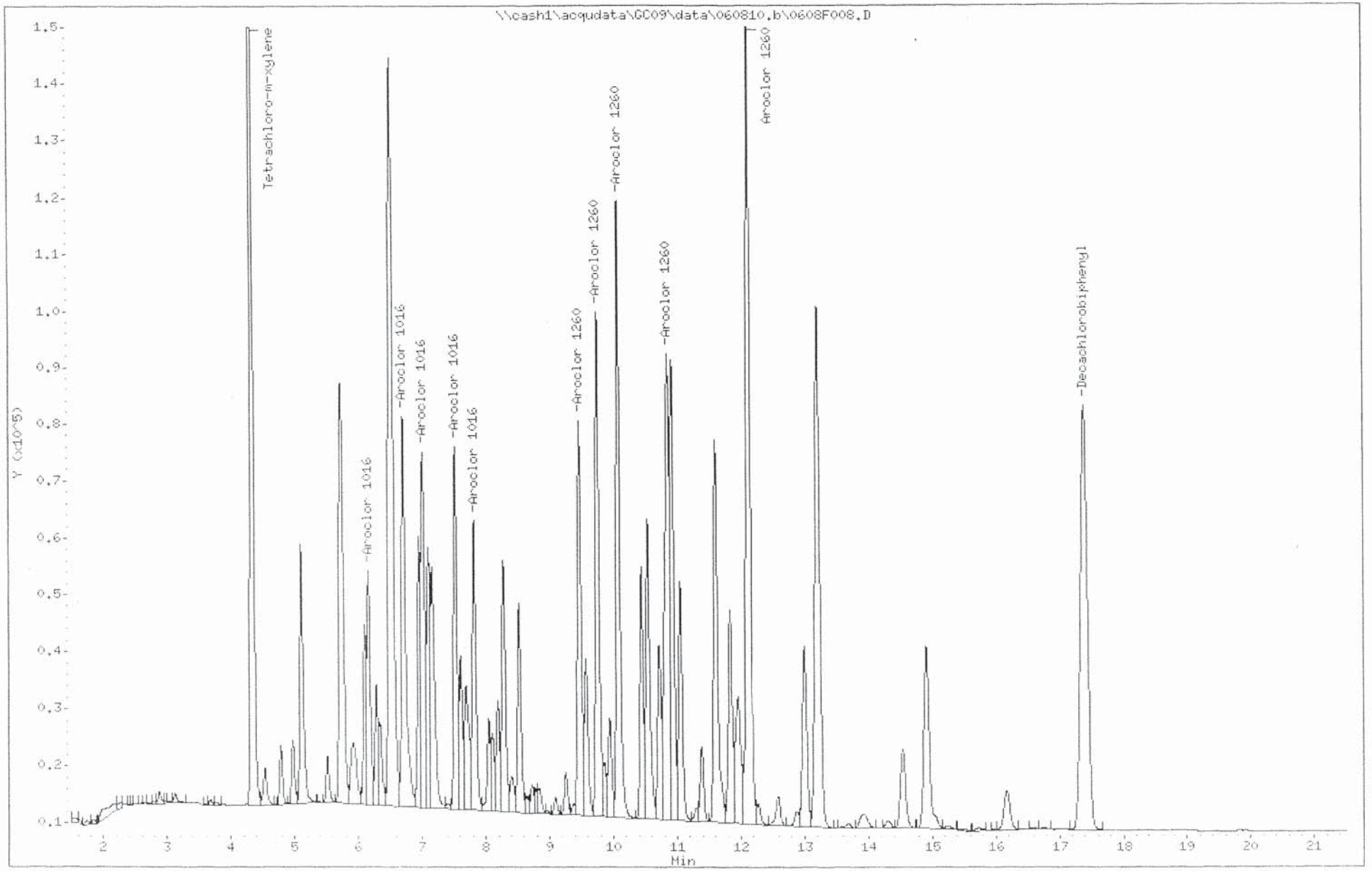
Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F008.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R008.D
Inj Date : 08-JUN-2010 20:38
Sample Info: KWG1005541-3 | K1005244-003MS | 8082 PCB_LL
Misc Info : SEMIVOA GC\W1005541\3-MS.H | F=4 D=1 A=22.78
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro m xylene	4.333	5.017	642771	577056	90.4	98.0		100.00
Aroclor 1016	6.163	6.527	149695	204802	1060	1120	80.00- 120.00	100.00
	6.707	7.143	311457	352426	1090	1090	164.25- 246.38	208.06
	7.013	7.427	246779	185890	1010	1180	130.18- 195.27	164.85
	7.523	7.603	227969	146599	1100	1120	112.90- 169.35	152.29
	7.817	7.707	199113	171398	1290	1160	87.15- 130.73	133.01
		Average of Peak Amounts =			1110	1130		
Aroclor 1260	9.470	10.343	281136	284961	930	1030	80.00- 120.00	100.00
	9.750	10.870	346804	310728	938	1050	99.43- 149.15	123.36
	10.077	11.520	419823	366027	927	984	120.99- 181.49	149.33
	10.853	12.283	388841	298377	898	1270	123.90- 185.85	138.31
	12.120	13.113	649374	542898	1140	1240	165.84- 248.76	230.98
		Average of Peak Amounts =			967	1110		
Decachlorobiphenyl	17.390	19.017	566222	527067	95.9	103		100.00

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F008.D
Date : 08-JUN-2010 20:38
Client ID:
Sample Info: KWG1005541-3 | K1005244-003MS | 8082 PCB_LL

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

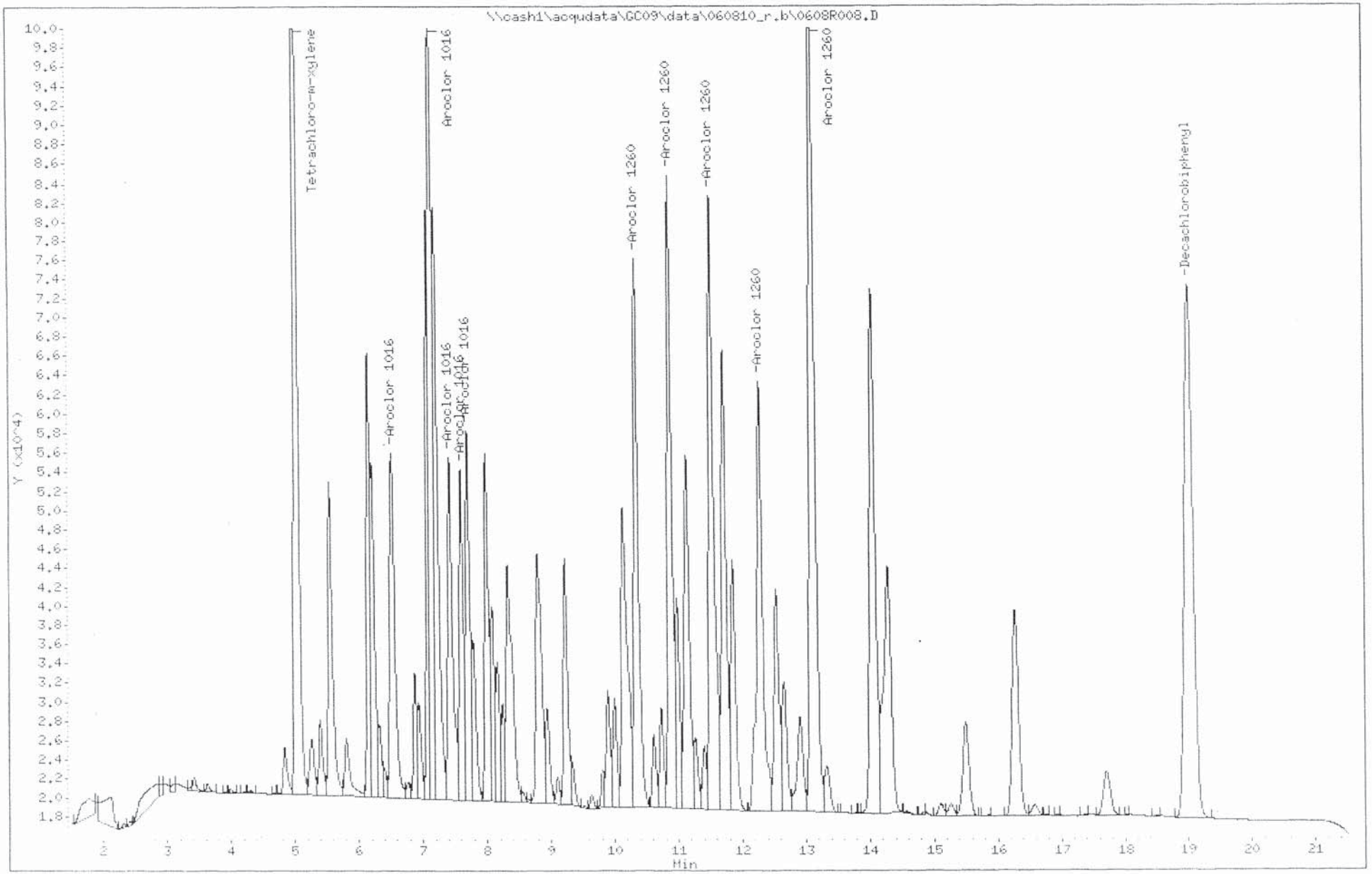


Data File: \\oash1\acqudata\GC09\data\060810_r.b\0608R008.D
Date : 08-JUN-2010 20:38
Client ID:
Sample Info: KWG1005541-3 | K1005244-003MS | 8082 PCB_LL

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Collected:**	05/19/2010
Sample Matrix:	Misc. solid	**Date Received:**	05/21/2010

Polychlorinated Biphenyls (PCBs)

Sample Name:	D-4-16DMS	**Units:**	ug/Kg
Lab Code:	KWG1005541-4	**Basis:**	Dry
Extraction Method:	EPA 3541	**Level:**	Low
Analysis Method:	8082		

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Aroclor 1016	**188**		9.9	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1221	ND	U	20	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1232	ND	U	9.9	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1242	ND	U	9.9	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1248	ND	U	9.9	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1254	ND	U	9.9	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1260	**183**		9.9	2.1	1	06/02/10	06/08/10	KWG1005541	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Decachlorobiphenyl	87	35-133	06/08/10	Acceptable

Comments:

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F009.D	**Date Acquired:**	06/08/2010 21:04
Lab ID:	KWG1005541-4 -- K1005244-003DMS	**Date Quantitated:**	06/09/2010 10:29
RunType:	DMS	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature]

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R009.D	**Date Acquired:**	06/08/2010 21:04
Lab ID:	KWG1005541-4 -- K1005244-003DMS	**Date Quantitated:**	06/09/2010 10:30
RunType:	DMS	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: 6/10/10

Quantitation Report

Bottle ID:		Tier:		Matrix:	MISC. SOLID
Prod Code:	8082 PCB_LL	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:	KWG1005541	Report Group:	
Analysis Method:	8082	Prep Method:	EPA 3541		
Prep Ref:	917274	Prep Date:	06/02/2010		

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:	J:\GC09\DATA\060810.B\0608F005.D	Method ID:	MJ696
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F009.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R009.D			Vial:	8
Acqu Date:	06/08/2010 21:04	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	DMS			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-4 -- K1005244-003DMS				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	4.33	5.02	528135	462740	74.33	78.58			79OK
			%Recovery =		74OK	79OK	Limits =	10-135	
Decachlorobiphenyl	17.38 -0.01	19.01 -0.01	491950	447392	83.33	87.16			87OK
			%Recovery =		83OK	87OK	Limits =	35-133	

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1016			0	0	910.65	952.90	180	188	188
Aroclor 1016 {1}	6.16 0.00	6.53	122841	174689	867.53	953.69	171	188	
Aroclor 1016 {2}	6.71	7.14 0.00	257995	294583	900.89	910.92	178	180	
Aroclor 1016 {3}	7.01 0.00	7.43 0.00	205400	155562	844.24	985.04	167	195	
Aroclor 1016 {4}	7.52	7.60	187829	123512	906.72	945.17	179	187	
Aroclor 1016 {5}	7.82 0.00	7.70 0.00	159421	143011	1,034	969.69	204	192	
Aroclor 1221			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1221 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1232 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1242 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F009.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R009.D			Vial:	8
Acqu Date:	06/08/2010 21:04	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	DMS			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-4 -- K1005244-003DMS				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1242 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1248 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1254 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260			0	0	814.19	927.14	161	183	183
Aroclor 1260 {1}	9.47	10.34 0.00	235953	240771	780.43	867.84	154	171	
Aroclor 1260 {2}	9.75 0.00	10.87 0.00	293217	259856	792.96	875.57	157	173	
Aroclor 1260 {3}	10.07 0.00	11.52 0.00	355446	304734	784.93	818.93	155	162	
Aroclor 1260 {4}	10.85 -0.01	12.28 0.00	328849	246561	759.10	1,047	150	207	
Aroclor 1260 {5}	12.12 -0.01	13.11 -0.01	544445	451119	953.53	1,026	188	203	
Aroclor 1262			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1262 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1268 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	

The -/+ after Retention Time symbolize the direction of the RT shift

Prep Amount:	22.33 g	Dilution:	1.0
Prep Final Vol:	4 mL	Unit Factor:	1
Solids:	90.7 %		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F009.D
Report Date: 09-Jun-2010 10:29

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F009.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R009.D
Inj Date : 08-JUN-2010 21:04
Sample Info: KWG1005541-4 | K1005244-003DMS | 8082 PCB_LL
Misc Info : SEMIVOA GC\W1005541\4-DMS.H | F=4 D=1 A=22.33
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

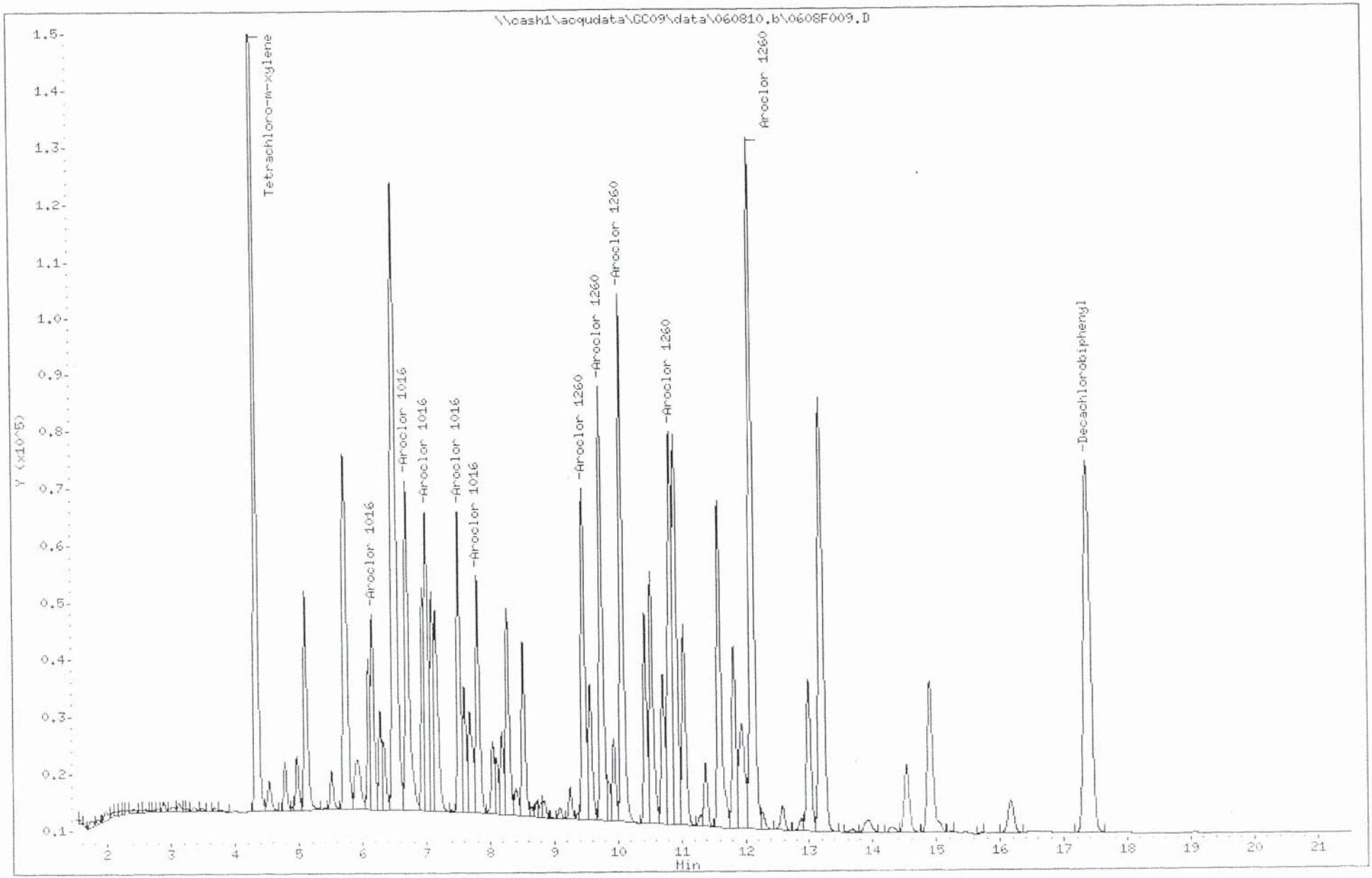
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro m-xylene	4.333	5.017	528135	462740	74.3	78.6		100.00
Aroclor 1016	6.163	6.527	122841	174689	868	954	80.00- 120.00	100.00
	6.707	7.143	257995	294583	901	911	164.25- 246.38	210.02
	7.013	7.427	206400	155562	844	985	130.18- 195.27	167.21
	7.523	7.603	187829	123512	907	945	112.90- 169.35	152.90
	7.817	7.703	159421	143011	1030	970	87.15- 130.73	129.78
			Average of Peak Amounts =		910	953		
Aroclor 1260	9.470	10.340	235953	240771	780	868	80.00- 120.00	100.00
	9.750	10.870	293217	259856	793	876	99.43- 149.15	124.27
	10.073	11.520	355446	304734	785	819	120.99- 181.49	150.64
	10.850	12.283	328849	246561	759	1050	123.90- 185.85	139.37
	12.117	13.110	544445	451119	954	1030	165.84- 248.76	230.74
			Average of Peak Amounts =		814	929		
Decachlorobiphenyl	17.383	19.013	491950	447392	83.3	87.2		100.00

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F009.D
Date : 08-JUN-2010 21:04
Client ID:
Sample Info: KWG1005541-4 | K1005244-003DMS | 8082 PCB_LL

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

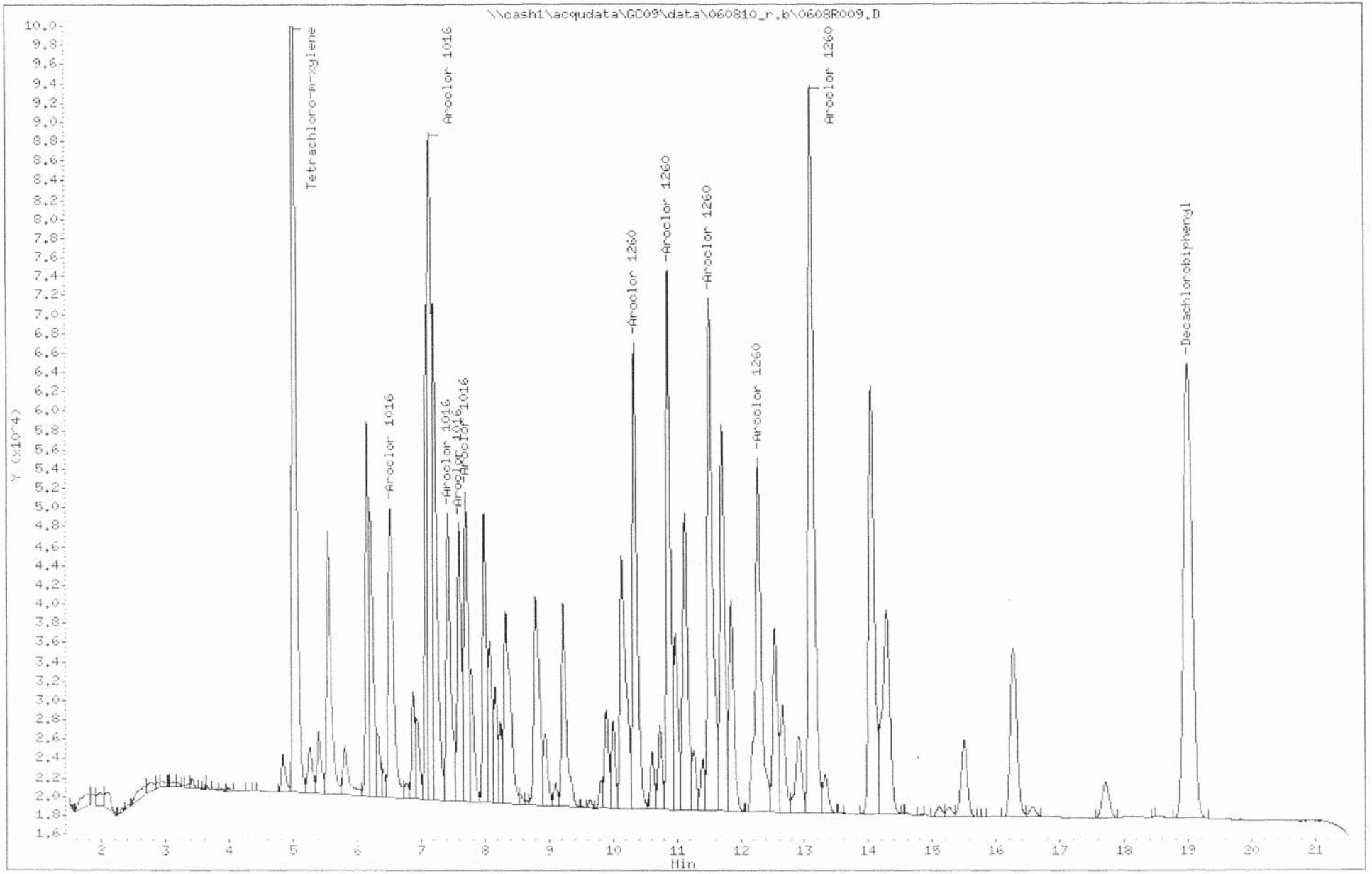
Column phase: DB-35MS



Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R009.D
Date : 08-JUN-2010 21:04
Client ID:
Sample Info: KWG1005541-4 | K1005244-003DMS | 8082 PCB_LL

Instrument: GC09.i
Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Polychlorinated Biphenyls (PCBs)

Sample Name: Lab Control Sample
Lab Code: KWG1005541-1

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 3541
Analysis Method: 8082

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Aroclor 1016	**203**		10	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1221	ND	U	20	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1232	ND	U	10	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1242	ND	U	10	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1248	ND	U	10	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1254	ND	U	10	2.1	1	06/02/10	06/08/10	KWG1005541	
Aroclor 1260	**204**		10	2.1	1	06/02/10	06/08/10	KWG1005541	

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Decachlorobiphenyl	93	35-133	06/08/10	Acceptable

Comments:

Exception Report

Data File: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F006.D
Lab ID: KWG1005541-1
RunType: LCS
Matrix: MISC. SOLID

Date Acquired: 06/08/2010 19:45
Date Quantitated: 06/09/2010 10:28
Batch ID: KWG1005547
Analysis Method: 8082
MethodJoinID: MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature]

Exception Report

Data File:	\\CASHI\ACQUDATA\GC09\DATA\060810_R.B\0608R006.D	**Date Acquired:**	06/08/2010 19:45
Lab ID:	KWG1005541-1	**Date Quantitated:**	06/09/2010 10:29
RunType:	LCS	**Batch ID:**	KWG1005547
Matrix:	MISC. SOLID	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Analytical Holding Time	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Calibration Verification Ave %D	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Calibration Verification Ave %D (Closing)	NA	NA	NA	x	
Continuing Calibration Recovery (Closing)	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature] 6/10/10

Quantitation Report

Bottle ID:		Tier:		Matrix:	MISC_SOLID
Prod Code:	8082 PCB_LL	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:	KWG1005541	Report Group:	
Analysis Method:	8082	Prep Method:	EPA 3541		
Prep Ref:	917271	Prep Date:	06/02/2010		

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:	J:\GC09\DATA\060810.B\0608F005.D	Method ID:	MJ696
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F006.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R006.D			Vial:	5
Acqu Date:	06/08/2010 19:45	Quant Date:	06/09/2010 10:28	Dilution:	1.0
Run Type:	LCS			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-1				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2		Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	4.33	5.02		591845	518809	83.29	88.10			88OK
				%Recovery =		83OK	88OK	Limits =	10-135	
Decachlorobiphenyl	17.39	19.02	0.00	514364	476905	87.13	92.91			93OK
				%Recovery =		87OK	93OK	Limits =	35-133	

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT #1	RT #2		Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1016				0	0	1,005	1,014	201	203	203
Aroclor 1016 {1}	6.16	6.53	0.00	138279	183793	976.56	1,003	195	201	
Aroclor 1016 {2}	6.71	7.14	0.00	291435	310282	1,018	959.46	204	192	
Aroclor 1016 {3}	7.02	7.43	0.00	219819	170835	903.50	1,082	181	216	
Aroclor 1016 {4}	7.52	7.60		200038	130192	965.66	996.29	193	199	
Aroclor 1016 {5}	7.82	7.71		178814	151789	1,160	1,029	232	206	
Aroclor 1221				0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1221 {1}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {2}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {3}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1221 {4}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232				0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1232 {1}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {2}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {3}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1232 {4}				0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242				0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1242 {1}				0d	0d	0.0000	0.0000	2.1U	2.1U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL, also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F006.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R006.D			Vial:	5
Acqu Date:	06/08/2010 19:45	Quant Date:	06/09/2010 10:28	Dilution:	1.0
Run Type:	LCS			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005541-1				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/Kg #1	ug/Kg #2	Rpt
Aroclor 1242 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1242 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1248 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1248 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1254 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1254 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1260			0	0	896.32	1,021	179	204	204
Aroclor 1260 {1}	9.47	10.34	251339	265246	831.32	956.06	166	191	
Aroclor 1260 {2}	9.75 0.00	10.87 0.00	334314	285279	904.10	961.24	181	192	
Aroclor 1260 {3}	10.08	11.52	390006	333213	861.25	895.46	172	179	
Aroclor 1260 {4}	10.85 0.00	12.28 0.00	353725	271578	816.53	1,153	163	231	
Aroclor 1260 {5}	12.12 0.00	13.12	610026	499913	1,068	1,137	214	227	
Aroclor 1262			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1262 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1262 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268			0	0	0.0000	0.0000	2.1U	2.1U	2.1U
Aroclor 1268 {1}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {2}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {3}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {4}			0d	0d	0.0000	0.0000	2.1U	2.1U	
Aroclor 1268 {5}			0d	0d	0.0000	0.0000	2.1U	2.1U	

The -/+ after Retention Time symbolize the direction of the RT shift

Prep Amount:	20.00 g	Dilution:	1.0
Prep Final Vol:	4 mL	Unit Factor:	1
Solids:	%		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F006.D
Report Date: 09-Jun-2010 10:28

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F006.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R006.D
Inj Date : 08-JUN-2010 19:45
Sample Info: KWG1005541-1 | LCS | 8082 PCB_LL | MISC. SOLI
Misc Info : SEMIVOA GC\W1005541\1-LCS.H | F=4 D=1 A=20.00
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

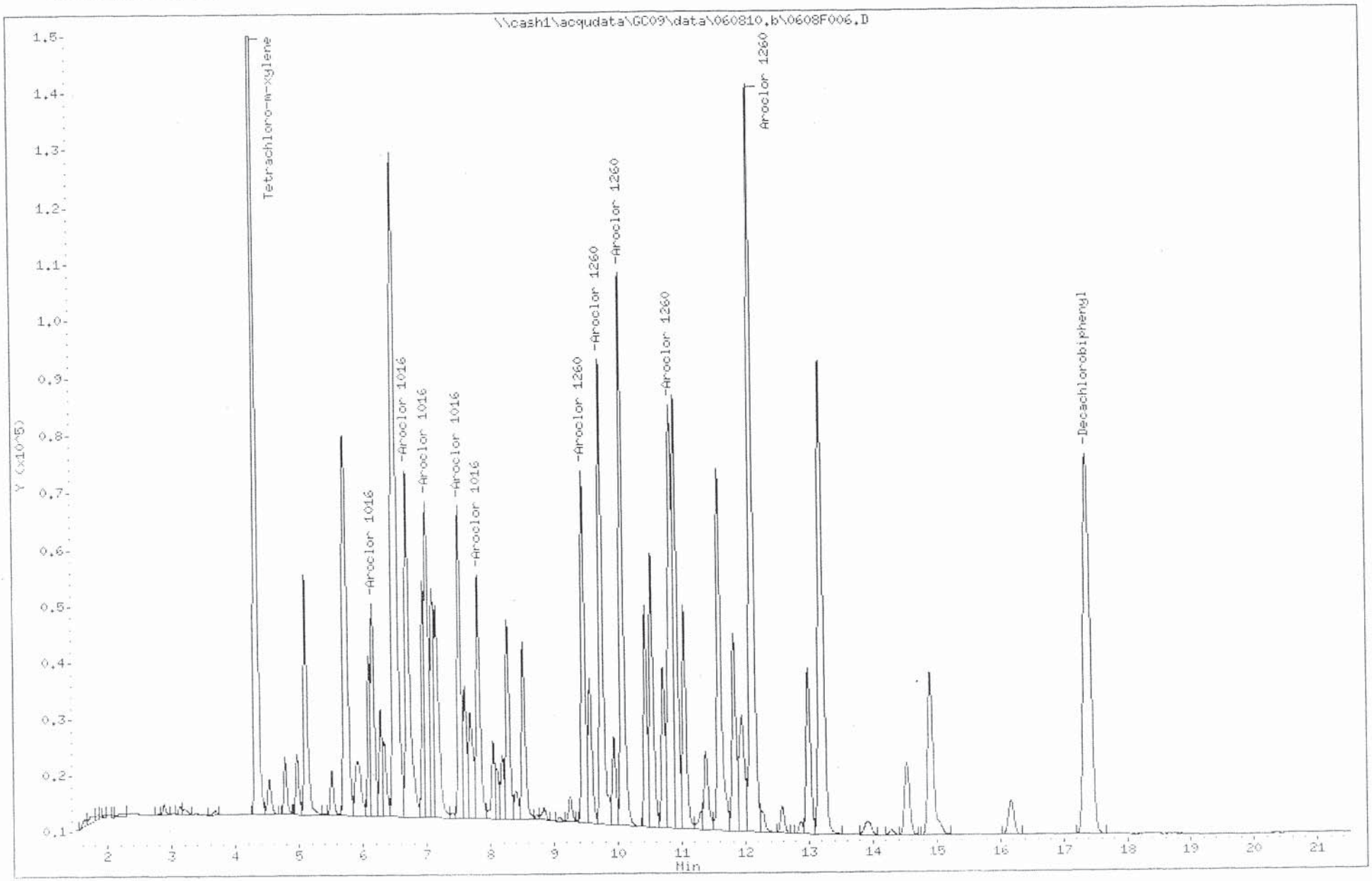
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.333	5.017	591845	518809	83.3	88.1		100.00
Aroclor 1016	6.163	6.527	138279	183793	976	1000	80.00- 120.00	100.00
	6.707	7.143	291435	310282	1020	959	164.25- 246.38	210.76
	7.017	7.427	219819	170835	904	1080	130.18- 195.27	158.97
	7.523	7.603	200038	130192	966	996	112.90- 169.35	144.66
	7.820	7.707	178814	151789	1160	1030	87.15- 130.73	129.31
			Average of Peak Amounts =		1000	1010		
Aroclor 1260	9.470	10.343	251339	265246	831	956	80.00- 120.00	100.00
	9.750	10.870	334314	285279	904	961	99.43- 149.15	133.01
	10.077	11.523	390006	333213	661	895	120.99- 181.49	155.17
	10.853	12.283	353725	271578	816	1150	123.90- 185.85	140.74
	12.120	13.117	610026	499913	1070	1140	165.84- 248.76	242.71
			Average of Peak Amounts =		896	1020		
Decachlorobiphenyl	17.390	19.020	514364	476905	87.1	92.9		100.00

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F006.D
Date : 08-JUN-2010 19:45
Client ID:
Sample Info: KWG1005541-1 | LCS | 8082 PCB_LL | MISC. SOLI

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

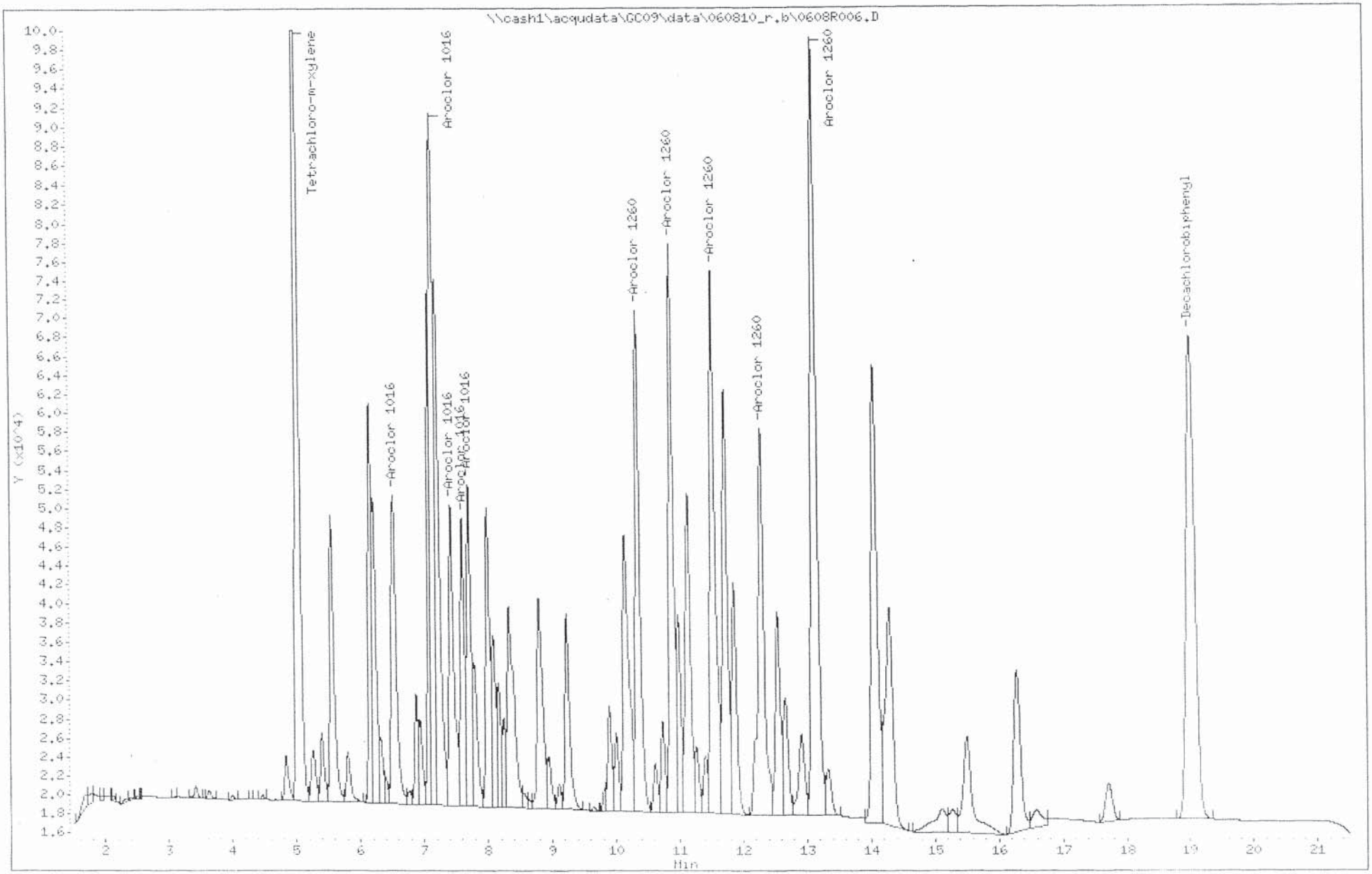
Column phase: DB-35MS



Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R006.D
Date : 08-JUN-2010 19:45
Client ID:
Sample Info: KWG1005541-1 | LCS | 8082 PCB_LL | MISC, SOLI

Instrument: GC09.i
Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



Organic Analysis:
Polychlorinated Biphenyls (PCBs)

Validation Package

Standards Data

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-35MS

Level ID	File ID
A	\\Cash1\Acqudata\GC09\data\052710A.B\0527F022.D
B	\\Cash1\Acqudata\GC09\data\052710A.B\0527F023.D
C	\\Cash1\Acqudata\GC09\data\052710A.B\0527F024.D
D	\\Cash1\Acqudata\GC09\data\052710A.B\0527F025.D
E	\\Cash1\Acqudata\GC09\data\052710A.B\0527F026.D
F	\\Cash1\Acqudata\GC09\data\052710A.B\0527F027.D
G	\\Cash1\Acqudata\GC09\data\052710A.B\0527F028.D
H	\\Cash1\Acqudata\GC09\data\052710A.B\0527F029.D
I	\\Cash1\Acqudata\GC09\data\052710A.B\0527F030.D
J	\\Cash1\Acqudata\GC09\data\052710A.B\0527F031.D
K	\\Cash1\Acqudata\GC09\data\052710A.B\0527F032.D
L	\\Cash1\Acqudata\GC09\data\052710A.B\0527F033.D
M	\\Cash1\Acqudata\GC09\data\052710A.B\0527F034.D
N	\\Cash1\Acqudata\GC09\data\052710A.B\0527F035.D
O	\\Cash1\Acqudata\GC09\data\052710A.B\0527F036.D
P	\\Cash1\Acqudata\GC09\data\052710A.B\0527F037.D
Q	\\Cash1\Acqudata\GC09\data\052710A.B\0527F038.D
R	\\Cash1\Acqudata\GC09\data\052710A.B\0527F039.D
S	\\Cash1\Acqudata\GC09\data\052710A.B\0527F040.D
T	\\Cash1\Acqudata\GC09\data\052710A.B\0527F041.D
U	\\Cash1\Acqudata\GC09\data\052710A.B\0527F042.D
V	\\Cash1\Acqudata\GC09\data\052710A.B\0527F043.D
W	\\Cash1\Acqudata\GC09\data\052710A.B\0527F044.D
X	\\Cash1\Acqudata\GC09\data\052710A.B\0527F045.D
Y	\\Cash1\Acqudata\GC09\data\052710A.B\0527F046.D
Z	\\Cash1\Acqudata\GC09\data\052710A.B\0527F047.D
AA	\\Cash1\Acqudata\GC09\data\052710A.B\0527F048.D
AB	\\Cash1\Acqudata\GC09\data\052710A.B\0527F049.D
AC	\\Cash1\Acqudata\GC09\data\052710A.B\0527F050.D
AD	\\Cash1\Acqudata\GC09\data\052710A.B\0527F051.D

Analyte Name	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF
Decachlorobiphenyl	A	2.5	6920	B	5.0	6750	C	50	6030	D	100	5800	E	200	5260
	F	500	4670												
Aroclor 1016 {1}	A	25	153	B	50	165	C	500	152	D	1000	143	E	2000	128
	F	5000	109												
Aroclor 1016 {2}	A	25	282	B	50	335	C	500	312	D	1000	299	E	2000	265
	F	5000	224												
Aroclor 1016 {3}	A	25	284	B	50	298	C	500	261	D	1000	234	E	2000	207
	F	5000	176												
Aroclor 1016 {4}	A	25	245	B	50	236	C	500	212	D	1000	205	E	2000	185
	F	5000	158												
Aroclor 1016 {5}	A	25	161	B	50	158	C	500	162	D	1000	163	E	2000	149
	F	5000	132												
Aroclor 1260 {1}	A	25	353	B	50	358	C	500	308	D	1000	292	E	2000	267
	F	5000	235												
Aroclor 1260 {2}	A	25	459	B	50	446	C	500	373	D	1000	351	E	2000	317
	F	5000	274												
Aroclor 1260 {3}	A	25	549	B	50	546	C	500	454	D	1000	432	E	2000	393
	F	5000	342												
Aroclor 1260 {4}	A	25	431	B	50	489	C	500	453	D	1000	440	E	2000	413
	F	5000	373												

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-35MS

Analyte Name	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF
Aroclor 1260 {5}	A	25	583	B	50	607	C	500	601	D	1000	593	E	2000	546
	F	5000	496												

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-35MS

Analyte Name	Compound Type	Calibration Evaluation				
		Fit Type	Eval.	Eval. Result	Q	Control Criteria
Decachlorobiphenyl	SURR	AverageRF	% RSD	14.6		≤ 20
Aroclor 1016 {1}	MULTI	AverageRF	% RSD	14.3		≤ 20
Aroclor 1016 {2}	MULTI	AverageRF	% RSD	13.5		≤ 20
Aroclor 1016 {3}	MULTI	AverageRF	% RSD	19.2		≤ 20
Aroclor 1016 {4}	MULTI	AverageRF	% RSD	15.6		≤ 20
Aroclor 1016 {5}	MULTI	AverageRF	% RSD	7.8		≤ 20
Aroclor 1260 {1}	MULTI	AverageRF	% RSD	15.8		≤ 20
Aroclor 1260 {2}	MULTI	AverageRF	% RSD	19.5		≤ 20
Aroclor 1260 {3}	MULTI	AverageRF	% RSD	18.3		≤ 20
Aroclor 1260 {4}	MULTI	AverageRF	% RSD	9.0		≤ 20
Aroclor 1260 {5}	MULTI	AverageRF	% RSD	7.5		≤ 20

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010
Date Analyzed: 05/28/2010

Second Source Calibration Verification
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082

Calibration ID: CAL9510
Units: ng/mL

File ID:
\\Cash1\Acqudata\GC09\data\052710A.B\0527F052.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F053.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F054.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F055.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F056.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F057.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F058.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F059.D
\\Cash1\Acqudata\GC09\data\052710A.B\0527F060.D

Column ID: DB-35MS

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Aroclor 1016 {1}	1000	1100	142	160	13	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1200	286	334	17	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	243	257	6	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1100	207	224	8	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1300	154	198	29	NA	± 100 %	AverageRF
Aroclor 1016	1000	1100	NA	NA	NA	-14	± 15 %	NA
Aroclor 1260 {1}	1000	950	302	287	-5	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	1000	370	377	2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	990	453	450	-1	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	860	433	374	-14	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1200	571	671	18	NA	± 100 %	AverageRF
Aroclor 1260	1000	1000	NA	NA	NA	0	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-XLB

Level ID	File ID
A	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R022.D
B	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R023.D
C	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R024.D
D	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R025.D
E	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R026.D
F	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R027.D
G	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R028.D
H	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R029.D
I	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R030.D
J	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R031.D
K	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R032.D
L	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R033.D
M	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R034.D
N	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R035.D
O	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R036.D
P	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R037.D
Q	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R038.D
R	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R039.D
S	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R040.D
T	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R041.D
U	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R042.D
V	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R043.D
W	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R044.D
X	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R045.D
Y	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R046.D
Z	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R047.D
AA	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R048.D
AB	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R049.D
AC	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R050.D
AD	\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R051.D

Analyte Name	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF
Decachlorobiphenyl	A	2.5	5660	B	5.0	5610	C	50	5290	D	100	5190	E	200	4790
	F	500	4260												
Aroclor 1016 {1}	A	25	198	B	50	195	C	500	190	D	1000	188	E	2000	173
	F	5000	155												
Aroclor 1016 {2}	A	25	375	B	50	377	C	500	322	D	1000	320	E	2000	291
	F	5000	255												
Aroclor 1016 {3}	A	25	148	B	50	139	C	500	173	D	1000	176	E	2000	163
	F	5000	148												
Aroclor 1016 {4}	A	25	142	B	50	136	C	500	140	D	1000	137	E	2000	124
	F	5000	105												
Aroclor 1016 {5}	A	25	152	B	50	152	C	500	159	D	1000	156	E	2000	139
	F	5000	125												
Aroclor 1260 {1}	A	25	294	B	50	311	C	500	290	D	1000	281	E	2000	257
	F	5000	232												
Aroclor 1260 {2}	A	25	320	B	50	344	C	500	308	D	1000	296	E	2000	270
	F	5000	243												
Aroclor 1260 {3}	A	25	392	B	50	383	C	500	383	D	1000	382	E	2000	358
	F	5000	334												
Aroclor 1260 {4}	A	25	188	B	50	237	C	500	253	D	1000	256	E	2000	245
	F	5000	234												

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-XLB

Analyte Name	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF	Level ID	Amt	RF
Aroclor 1260 {5}	A	25	456	B	50	453	C	500	449	D	1000	450	E	2000	428
	F	5000	401												

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010

Initial Calibration Summary
Polychlorinated Biphenyls (PCBs)

Calibration ID: CAL9510
Instrument ID: GC09.i

Column: DB-XLB

Analyte Name	Compound Type	Calibration Evaluation				
		Fit Type	Eval.	Eval. Result	Q	Control Criteria
Decachlorobiphenyl	SURR	AverageRF	% RSD	10.4		≤ 20
Aroclor 1016 {1}	MULTI	AverageRF	% RSD	9.0		≤ 20
Aroclor 1016 {2}	MULTI	AverageRF	% RSD	14.7		≤ 20
Aroclor 1016 {3}	MULTI	AverageRF	% RSD	9.6		≤ 20
Aroclor 1016 {4}	MULTI	AverageRF	% RSD	10.7		≤ 20
Aroclor 1016 {5}	MULTI	AverageRF	% RSD	8.7		≤ 20
Aroclor 1260 {1}	MULTI	AverageRF	% RSD	10.3		≤ 20
Aroclor 1260 {2}	MULTI	AverageRF	% RSD	12.1		≤ 20
Aroclor 1260 {3}	MULTI	AverageRF	% RSD	5.9		≤ 20
Aroclor 1260 {4}	MULTI	AverageRF	% RSD	10.5		≤ 20
Aroclor 1260 {5}	MULTI	AverageRF	% RSD	4.9		≤ 20

Results flagged with an asterisk (*) indicate values outside control criteria.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 05/27/2010
Date Analyzed: 05/28/2010

Second Source Calibration Verification
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082

Calibration ID: CAL9510
Units: ng/mL

File ID: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R052.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R053.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R054.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R055.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R056.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R057.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R058.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R059.D
\\Cash1\Acqudata\GC09\data\052710A_r.b\0527R060.D

Column ID: DB-XLB

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Aroclor 1016 {1}	1000	1100	183	195	6	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1000	323	326	1	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	158	181	15	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1000	131	137	5	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1100	147	163	10	NA	± 100 %	AverageRF
Aroclor 1016	1000	1100	NA	NA	NA	-7	± 15 %	NA
Aroclor 1260 {1}	1000	1000	277	280	1	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	980	297	292	-2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	910	372	339	-9	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	1200	235	282	20	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1200	440	506	15	NA	± 100 %	AverageRF
Aroclor 1260	1000	1100	NA	NA	NA	-5	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound** ‡ **CCC Compound**

Sequence Name: D:\GC09\SEQUENCE\052710A.S
Comment: PCB Aroclors by EPA 8082
Operator: LHarris
Data Path: D:\GC09\DATA\052710A.B\
Pre-Seq Cmd:
Post-Seq Cmd:

Method Sections To Run
(X) Full Method
( ) Reprocessing Only

On A Barcode Mismatch
(X) Inject Anyway
( ) Don't Inject

ICAL 9510

Do not use to report AR1221 for any DOD 4.1 hits
CH 5/28/10

Line	Type	Vial	DataFile	Method	Sample Name
1	DeleteGC				
2	MaskName -		----r---		
3	Clocksrt -		19:29		
4	IB	1	0527F020	PCB_REG	IB
5	IB	2	0527F021	PCB_REG	IB
6	ICAL	3	0527F022	PCB_REG	1016/1260 @ 25ppb | PCB5-55F
7	ICAL	4	0527F023	PCB_REG	1016/1260 @ 50ppb | PCB5-55G
8	ICAL	5	0527F024	PCB_REG	1016/1260 @ 500ppb | PCB5-55H
9	ICAL	6	0527F025	PCB_REG	1016/1260 @ 1000ppb | PCB5-55I
10	ICAL	7	0527F026	PCB_REG	1016/1260 @ 2000ppb | PCB5-55J
11	ICAL	8	0527F027	PCB_REG	1016/1260 @ 5000ppb | PCB5-55K
12	ICAL	9	0527F028	PCB_REG	1221/1254 @ 50/25ppb | PCB5-52D
13	ICAL	10	0527F029	PCB_REG	1221/1254 @ 100/50ppb | PCB5-52E
14	ICAL	11	0527F030	PCB_REG	1221/1254 @ 1000/500ppb | PCB5-52F
15	ICAL	12	0527F031	PCB_REG	1221/1254 @ 2000/1000ppb | PCB5-52G
16	ICAL	13	0527F032	PCB_REG	1221/1254 @ 4000/2000ppb | PCB5-52H
17	ICAL	14	0527F033	PCB_REG	1221/1254 @ 10000/5000ppb | PCB5-52I
18	ICAL	15	0527F034	PCB_REG	1232/1262 @ 25ppb | PCB5-52J
19	ICAL	16	0527F035	PCB_REG	1232/1262 @ 50ppb | PCB5-52K
20	ICAL	17	0527F036	PCB_REG	1232/1262 @ 500ppb | PCB5-52L
21	ICAL	18	0527F037	PCB_REG	1232/1262 @ 1000ppb | PCB5-52M
22	ICAL	19	0527F038	PCB_REG	1232/1262 @ 2000ppb | PCB5-52N
23	ICAL	20	0527F039	PCB_REG	1232/1262 @ 5000ppb | PCB5-52O
24	ICAL	21	0527F040	PCB_REG	1242/1268 @ 25ppb | PCB5-52P
25	ICAL	22	0527F041	PCB_REG	1242/1268 @ 50ppb | PCB5-52Q
26	ICAL	23	0527F042	PCB_REG	1242/1268 @ 500ppb | PCB5-52R
27	ICAL	24	0527F043	PCB_REG	1242/1268 @ 1000ppb | PCB5-52S
28	ICAL	25	0527F044	PCB_REG	1242/1268 @ 2000ppb | PCB5-52T
29	ICAL	26	0527F045	PCB_REG	1242/1268 @ 5000ppb | PCB5-53A
30	ICAL	27	0527F046	PCB_REG	1248 @ 25ppb | PCB5-53B
31	ICAL	28	0527F047	PCB_REG	1248 @ 50ppb | PCB5-53C
32	ICAL	29	0527F048	PCB_REG	1248 @ 500ppb | PCB5-53D
33	ICAL	30	0527F049	PCB_REG	1248 @ 1000ppb | PCB5-53E
34	ICAL	31	0527F050	PCB_REG	1248 @ 2000ppb | PCB5-53F
35	ICAL	32	0527F051	PCB_REG	1248 @ 5000ppb | PCB5-53G
36	ICV	33	0527F052	PCB_REG	1016 @ 1000ppb | PCB5-49E
37	ICV	34	0527F053	PCB_REG	1221 @ 1000ppb | PCB5-53J
38	ICV	35	0527F054	PCB_REG	1232 @ 1000ppb | PCB5-49F
39	ICV	36	0527F055	PCB_REG	1242 @ 1000ppb | PCB5-49G
40	ICV	37	0527F056	PCB_REG	1248 @ 1000ppb | PCB5-49H
41	ICV	38	0527F057	PCB_REG	1254 @ 1000ppb | PCB5-53K
42	ICV	39	0527F058	PCB_REG	1260 @ 1000ppb | PCB5-49I
43	ICV	40	0527F059	PCB_REG	1262 @ 1000ppb | PCB5-49J

Last Modified: Thu May 27 18:26:17 2010 Page: 1

Sequence Name: D:\GC09\SEQUENCE\052710A.S

Line	Type	Vial	DataFile	Method	Sample Name
44	ICV	41	0527F060	PCB_REG	1268 @ 1000ppb | PCB5-49K
45					

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F021.D
Report Date: 28-May-2010 17:28

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F021.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R021.D
Inj Date : 27-MAY-2010 23:27
Sample Info: IB
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F021.D
Date : 27-MAY-2010 23:27
Client ID:
Sample Info: IB

Instrument: GC09.i

Operator: LHarris

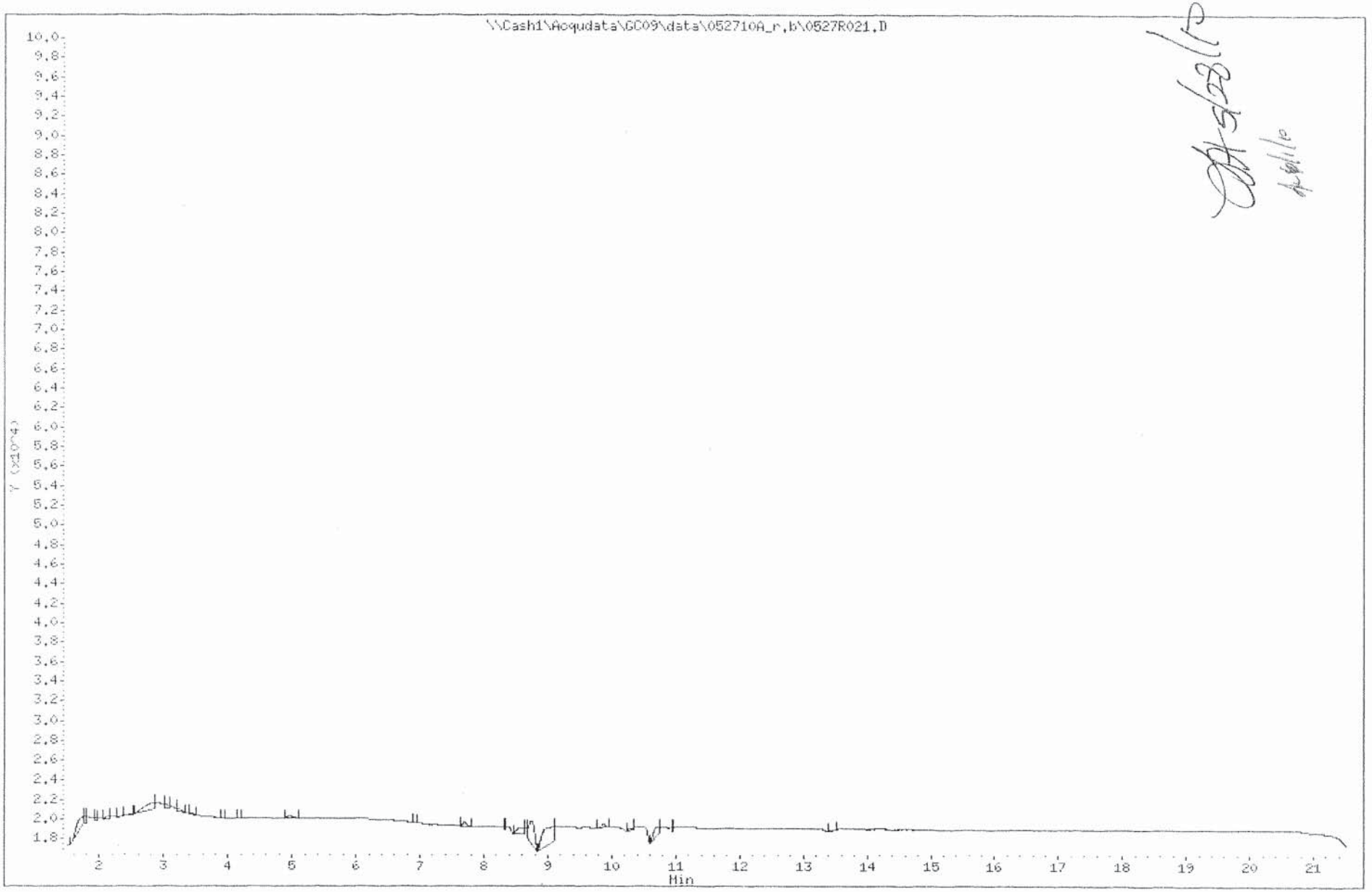
Column phase: DB-35MS

Column diameter: 0.53

Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R021.D
Date : 27-MAY-2010 23:27
Client ID:
Sample Info: IB

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F022.D
Report Date: 28-May-2010 17:28

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F022.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R022.D
Inj Date : 27-MAY-2010 23:53
Sample Info: 1016/1260 @ 25ppb | PCB5-55F
Misc Info :
Cal Date : 28-MAY-2010 13:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.343	5.030	22176	16014	3.12	2.72		100.00
Aroclor 1016	6.180	6.540	3830	4938	27.0	29.8	80.00- 120.00	100.00 (M)
	6.720	7.163	7061	9382	24.6	29.0	162.91- 244.36	184.36 (M)
	7.033	7.447	7093	3695	28.9	23.4	145.03- 217.55	185.21 (M)
	7.540	7.623	6135	3546	29.4	27.1	115.01- 172.51	160.18 (M)
	7.837	7.727	4022	3811	25.7	25.8	76.85- 115.27	105.01 (M)
			Average of Peak Amounts =		27.1	27.0		
Aroclor 1260	9.490	10.367	8826	7343	29.2	26.5	80.00- 120.00	100.00
	9.770	10.893	11469	7990	31.0	26.9	99.64- 149.46	129.94
	10.097	11.547	13727	9798	30.3	26.3	122.12- 183.18	155.51
	10.873	12.313	10765	4708	24.8	20.0	109.38- 164.06	121.97
	12.147	13.147	14576	11411	25.5	26.0	135.68- 203.52	165.14
			Average of Peak Amounts =		28.2	25.1		
Decachlorobiphenyl	17.433	19.073	17308	14157	2.93	2.76		100.00

QC Flag Legend

M - Compound response manually integrated.

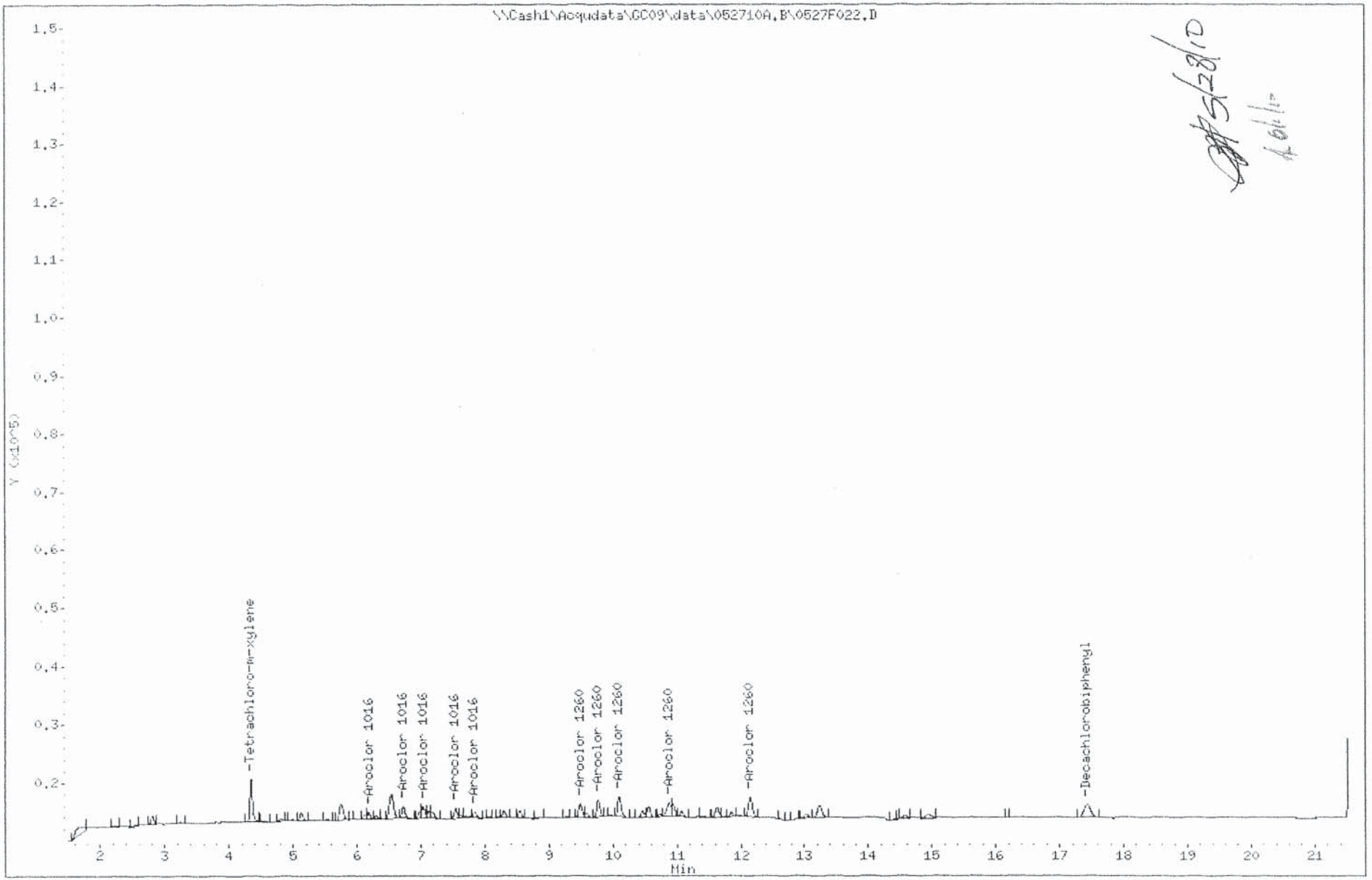
5/28/10
6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F022.D
Date : 27-MAY-2010 23:53
Client ID:
Sample Info: 1016/1260 @ 25ppb | PCB5-55F

Instrument: GC09.i

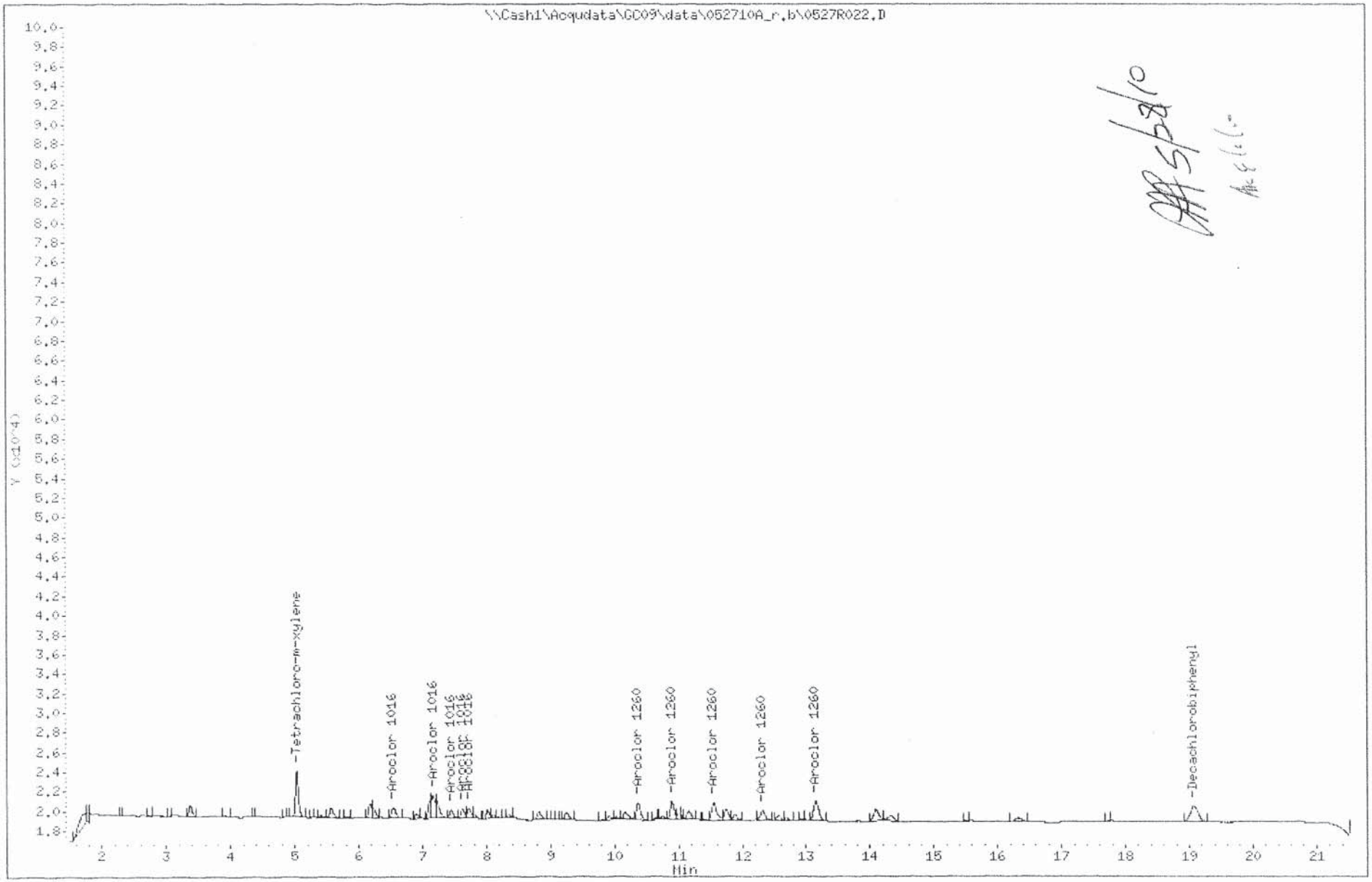
Operator: LHarris
Column diameter: 0.53

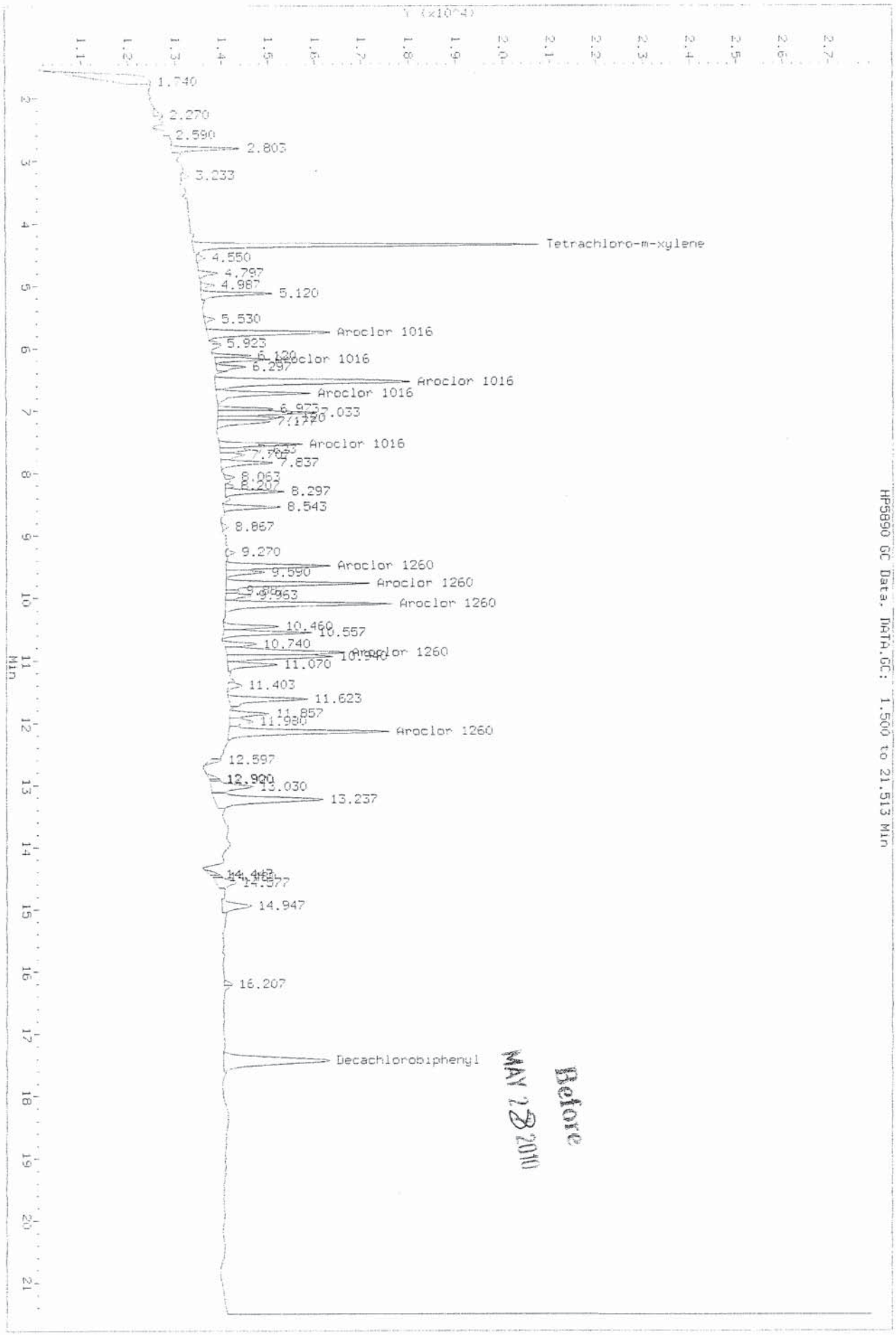
Column phase: DB-35MS



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R022.D
Date : 27-MAY-2010 23:53
Client ID:
Sample Info: 1016/1260 @ 25ppb | PCB5-55F

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

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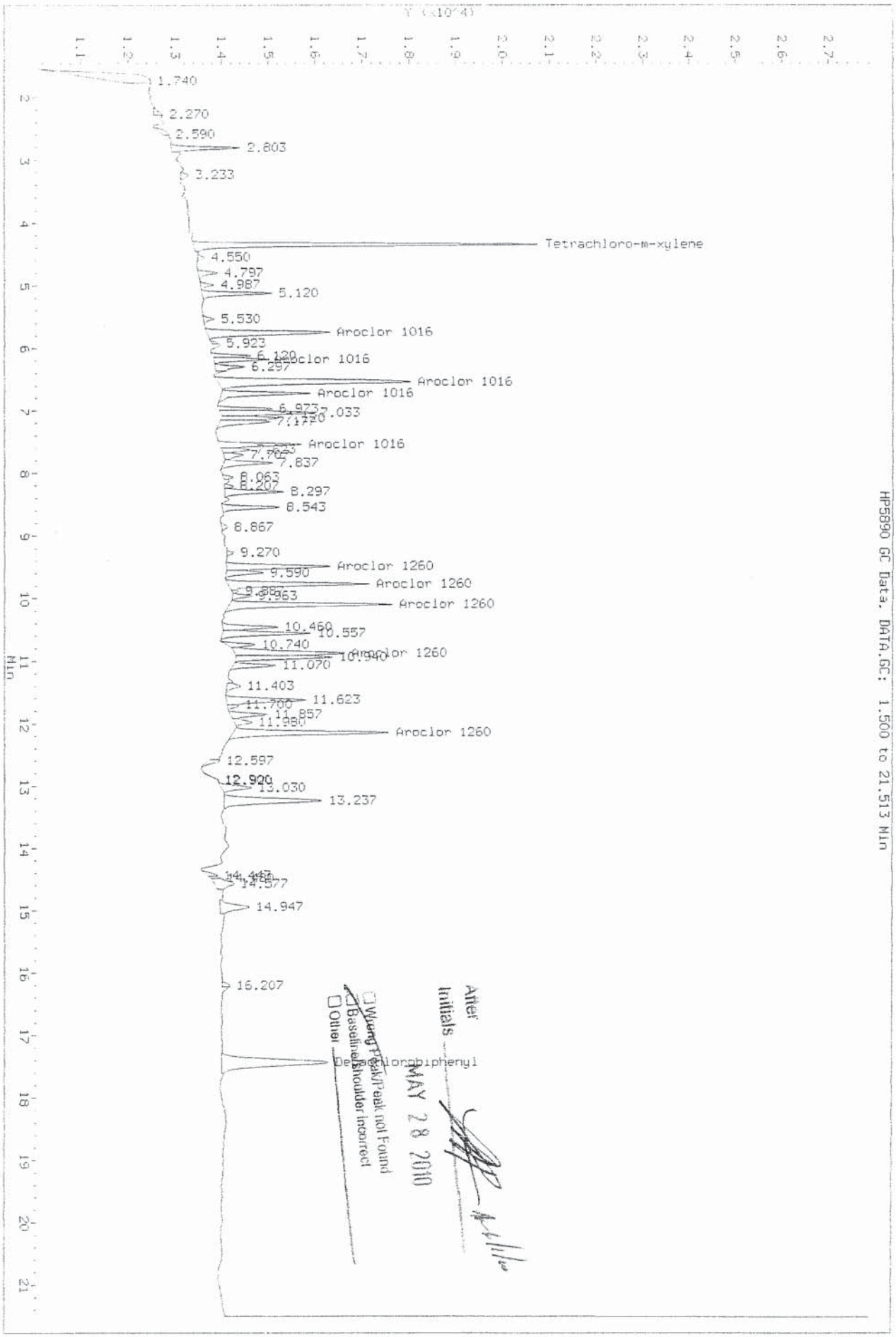
Injection Date: 27-MAY-2010 23:53
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.513 Min
Y (x10^4)
Min
1.740
2.270
2.590
2.803
3.233
Tetrachloro-m-xylene
4.550
4.797
4.987
5.120
5.530
Aroclor 1016
5.923
6.130
Aroclor 1016
6.297
Aroclor 1016
Aroclor 1016
6.923
7.033
Aroclor 1016
7.837
8.063
8.297
8.543
8.867
9.270
Aroclor 1260
9.590
Aroclor 1260
9.963
Aroclor 1260
10.460
10.557
10.740
Aroclor 1260
11.070
11.403
11.623
11.857
Aroclor 1260
12.597
12.900
13.030
13.237
14.947
16.207
Decachlorobiphenyl
Before
MAY 2 8 2010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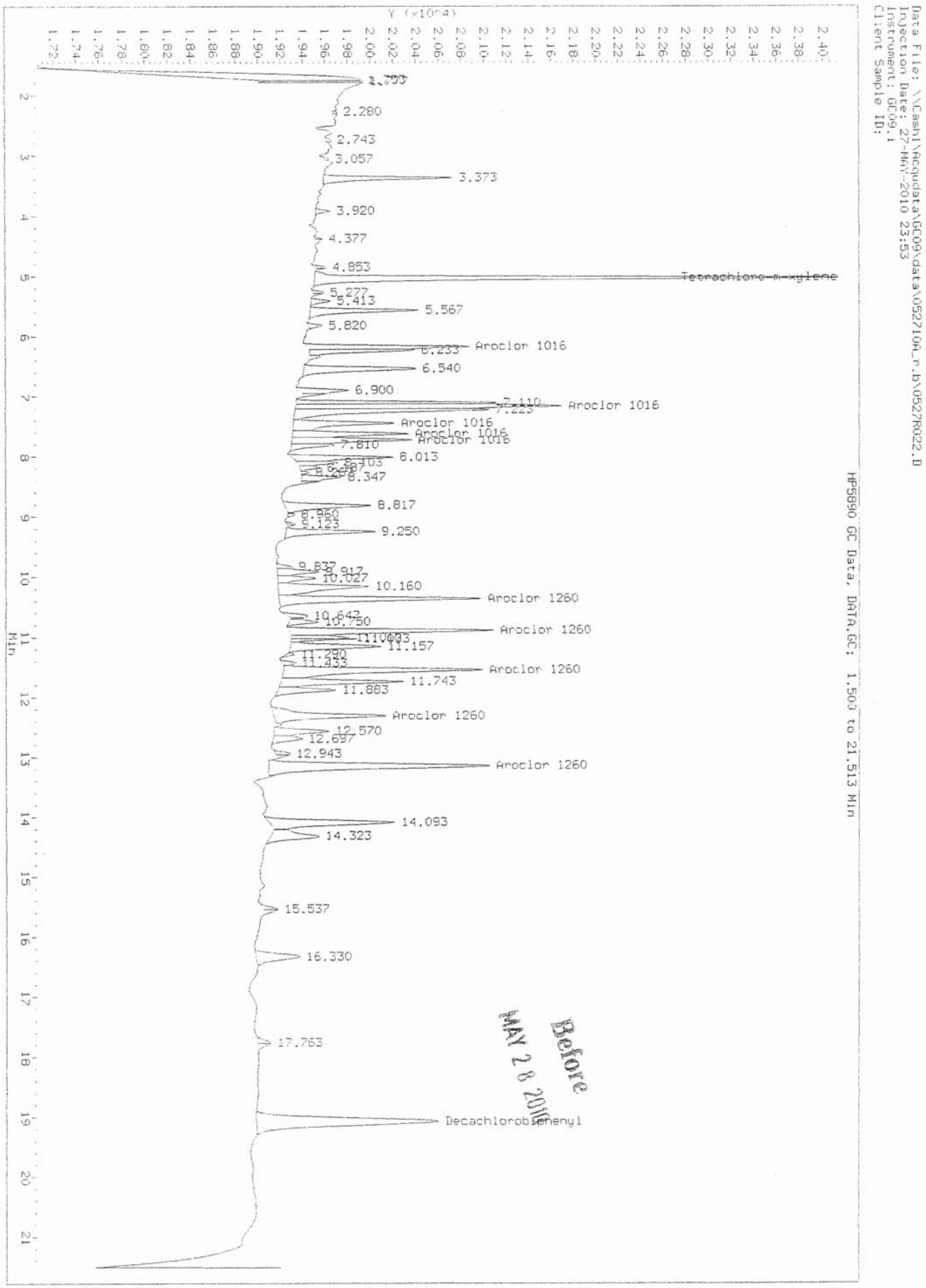
Injection Date: 27-MAY-2010 23:53
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.513 Min
Y (x10^4)
Min
1.740
2.270
2.590
2.803
3.233
Tetrachloro-m-xylene
4.550
4.797
4.987
5.120
5.530
Aroclor 1016
5.923
6.120
Aroclor 1016
6.297
Aroclor 1016
Aroclor 1016
6.973
7.033
Aroclor 1016
7.837
8.063
8.207
8.297
8.543
8.867
9.270
Aroclor 1260
9.590
Aroclor 1260
Aroclor 1260
10.460
10.557
10.740
Aroclor 1260
11.070
11.403
11.623
11.857
11.980
Aroclor 1260
12.597
12.900
13.030
13.237
14.947
16.207
Decachlorobiphenyl
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/Shoulder incorrect
Other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Injection Date: 27-MAY-2010 23:53
Instrument: GC09.i
Client Sample ID:
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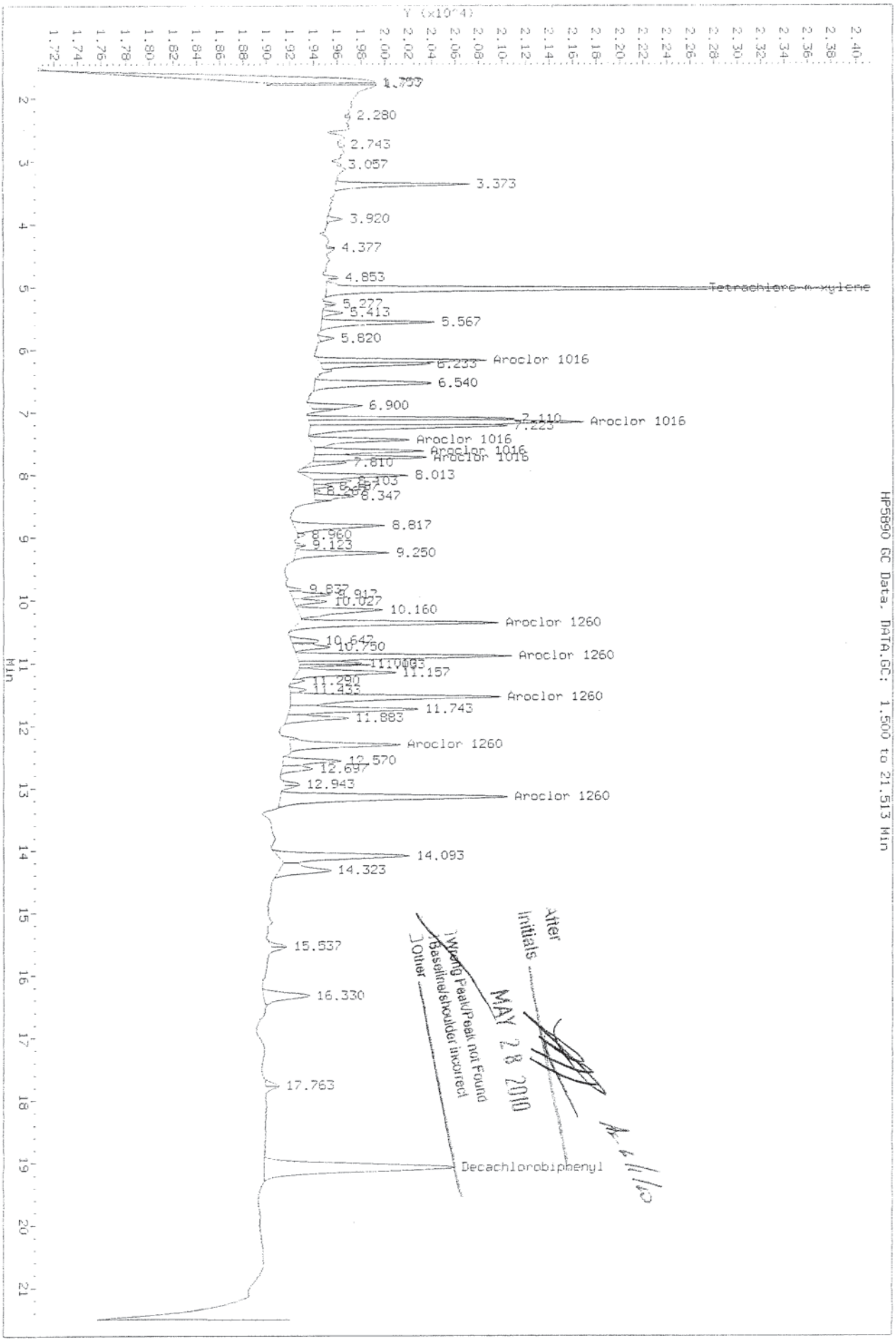
Y (x10^4)
Min
1.700
2.280
2.743
3.057
3.373
3.920
4.377
4.853
Tetrachloro-m-xylene
5.277
5.413
5.567
5.820
6.233
Aroclor 1016
6.540
6.900
7.110
7.223
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.810
8.013
8.347
8.817
8.960
9.123
9.250
9.837
9.917
10.027
10.160
Aroclor 1260
10.642
10.750
Aroclor 1260
11.083
11.157
11.290
11.453
Aroclor 1260
11.743
11.883
Aroclor 1260
12.570
12.697
12.943
Aroclor 1260
14.093
14.323
15.537
16.330
17.763
Decachlorobiphenyl
Before
MAY 28 2010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Instrument: 6609.i
Client Sample ID:
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Y (x10^4)
Min
1.709
2.280
2.743
3.057
3.373
3.920
4.377
4.853
Tetrachloro-m-xylene
5.227
5.413
5.567
5.820
6.233
Aroclor 1016
6.540
6.900
7.110
Aroclor 1016
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.810
8.013
8.103
8.347
8.817
8.960
9.123
9.250
9.837
9.912
10.027
10.160
Aroclor 1260
10.642
10.750
Aroclor 1260
11.157
11.290
11.453
Aroclor 1260
11.743
11.883
Aroclor 1260
12.570
12.697
12.943
Aroclor 1260
14.093
14.323
15.537
16.330
17.763
Decachlorobiphenyl
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/Shoulder Incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F023.D
Report Date: 28-May-2010 17:28

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F023.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R023.D
Inj Date : 28-MAY-2010 00:19
Sample Info: 1016/1260 @ 50ppb | PCB5-55G
Misc Info :
Cal Date : 28-MAY-2010 13:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.343	5.030	42292	30894	5.95	5.25		100.00
Aroclor 1016	6.180	6.543	8225	9771	58.1	59.1	80.00- 120.00	100.00(M)
	6.723	7.163	16749	18833	58.5	58.2	162.91- 244.36	203.64(M)
	7.033	7.447	14911	6955	61.3	44.0	145.03- 217.55	181.29(M)
	7.540	7.620	11824	6794	57.1	52.0	115.01- 172.51	143.76(M)
	7.837	7.727	7901	7616	51.2	51.6	76.85- 115.27	96.06(M)
			Average of Peak Amounts =		57.2	53.0		
Aroclor 1260	9.490	10.367	17890	15555	59.2	56.1	80.00- 120.00	100.00
	9.770	10.893	22282	17181	60.2	57.9	99.64- 149.46	124.55
	10.097	11.547	27309	19173	60.3	51.5	122.12- 183.18	152.65
	10.877	12.313	24459	11860	56.5	50.4	109.38- 164.06	136.72
	12.147	13.147	30342	22668	53.1	51.6	135.68- 203.52	169.61
			Average of Peak Amounts =		57.9	53.5		
Decachlorobiphenyl	17.433	19.073	33726	28071	5.71	5.47		100.00

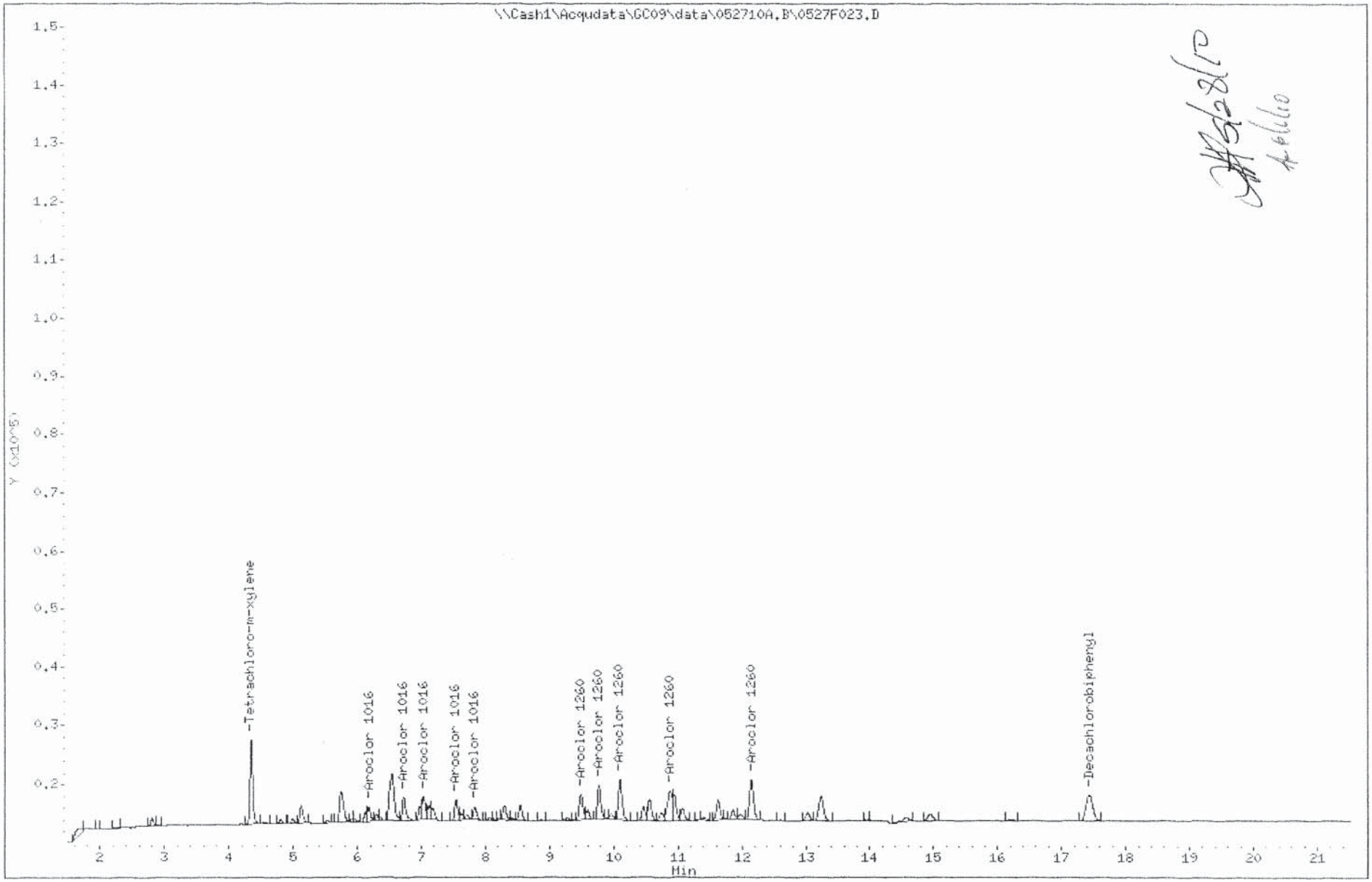
QC Flag Legend

M - Compound response manually integrated.

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F023.D
Date : 28-MAY-2010 00:19
Client ID:
Sample Info: 1016/1260 @ 50ppb | PCB5-55G

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



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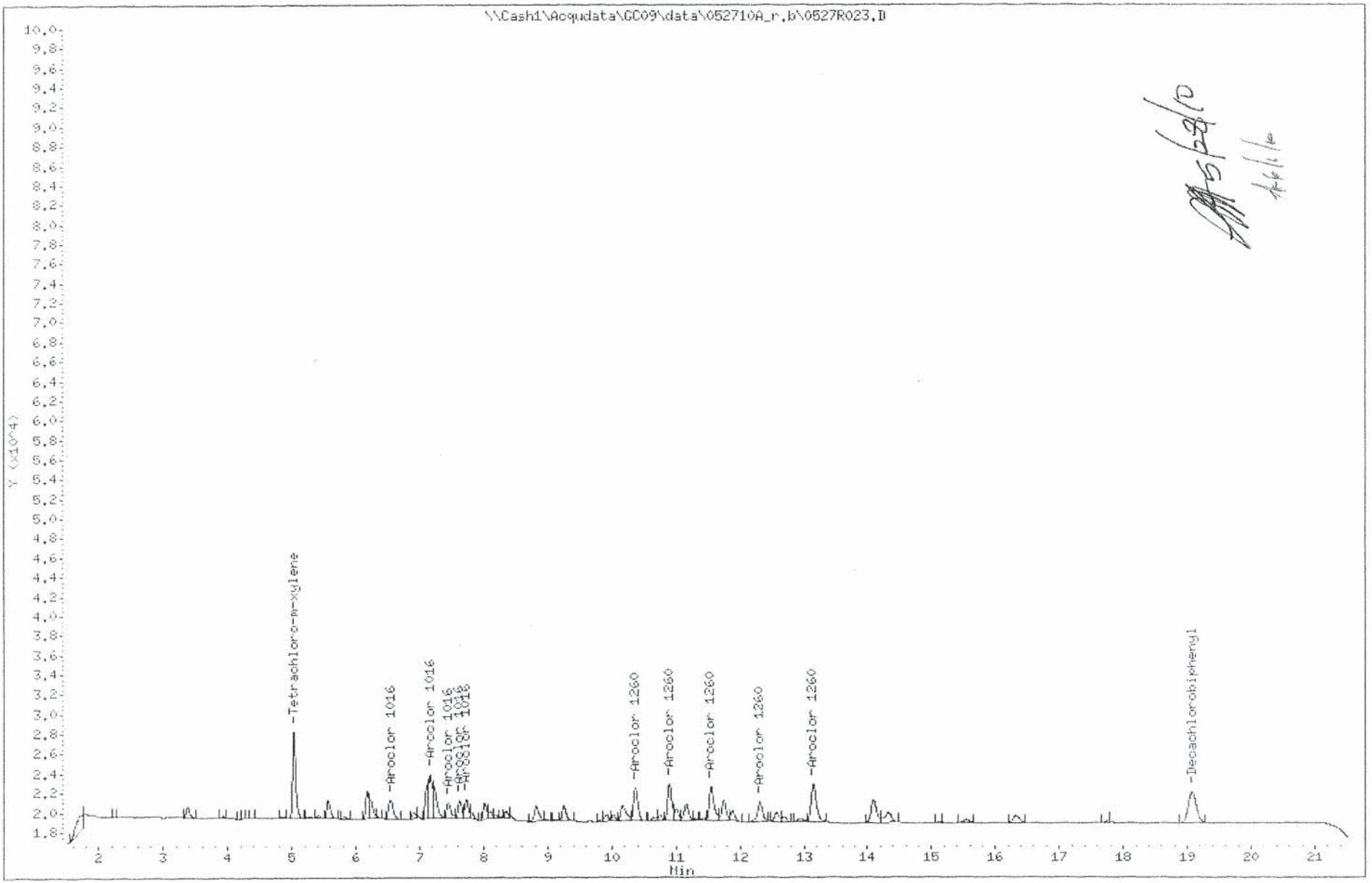
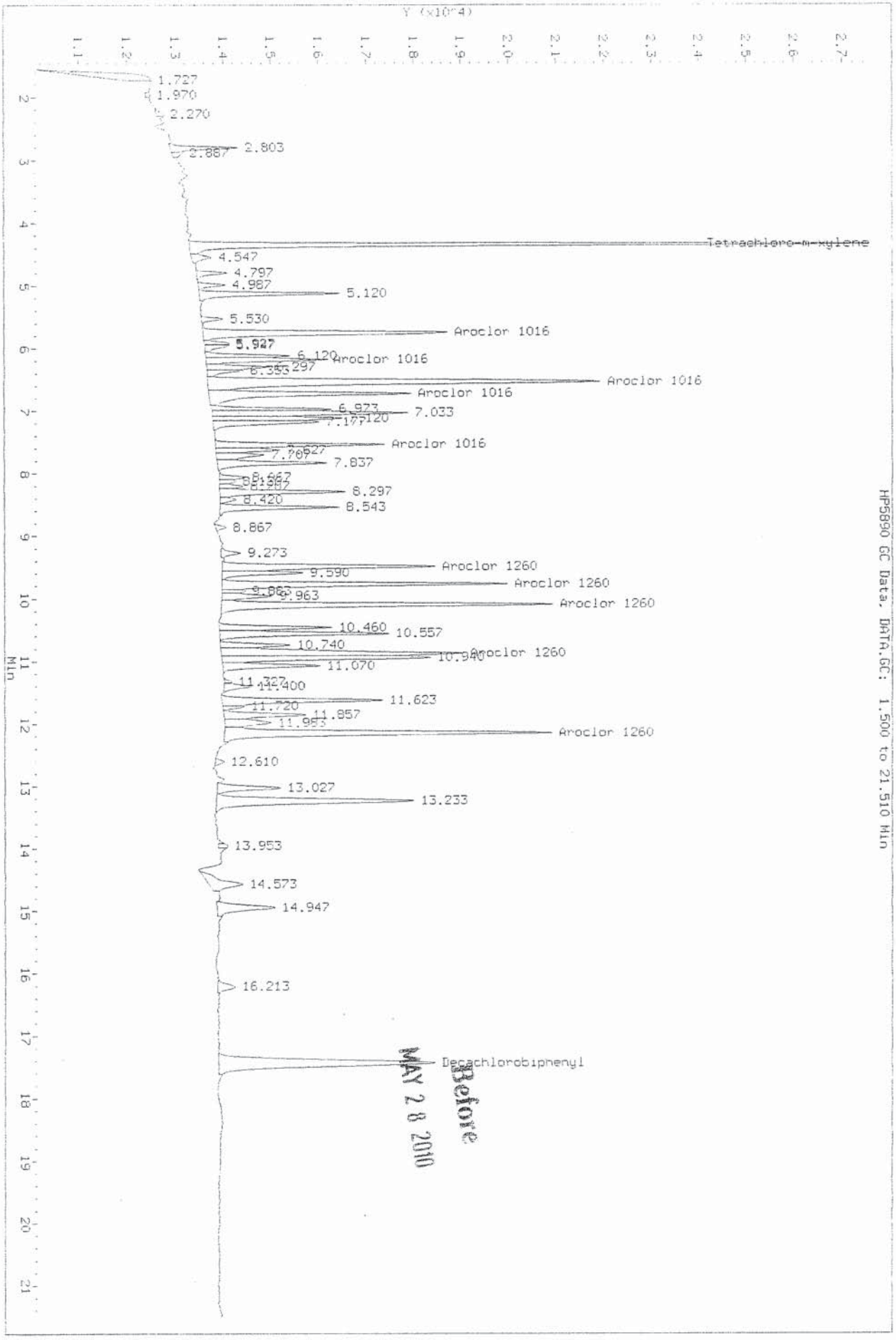
Date : 28-MAY-2010 00:19
Client ID:
Sample Info: 1016/1260 @ 50ppb | PCB5-55G

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

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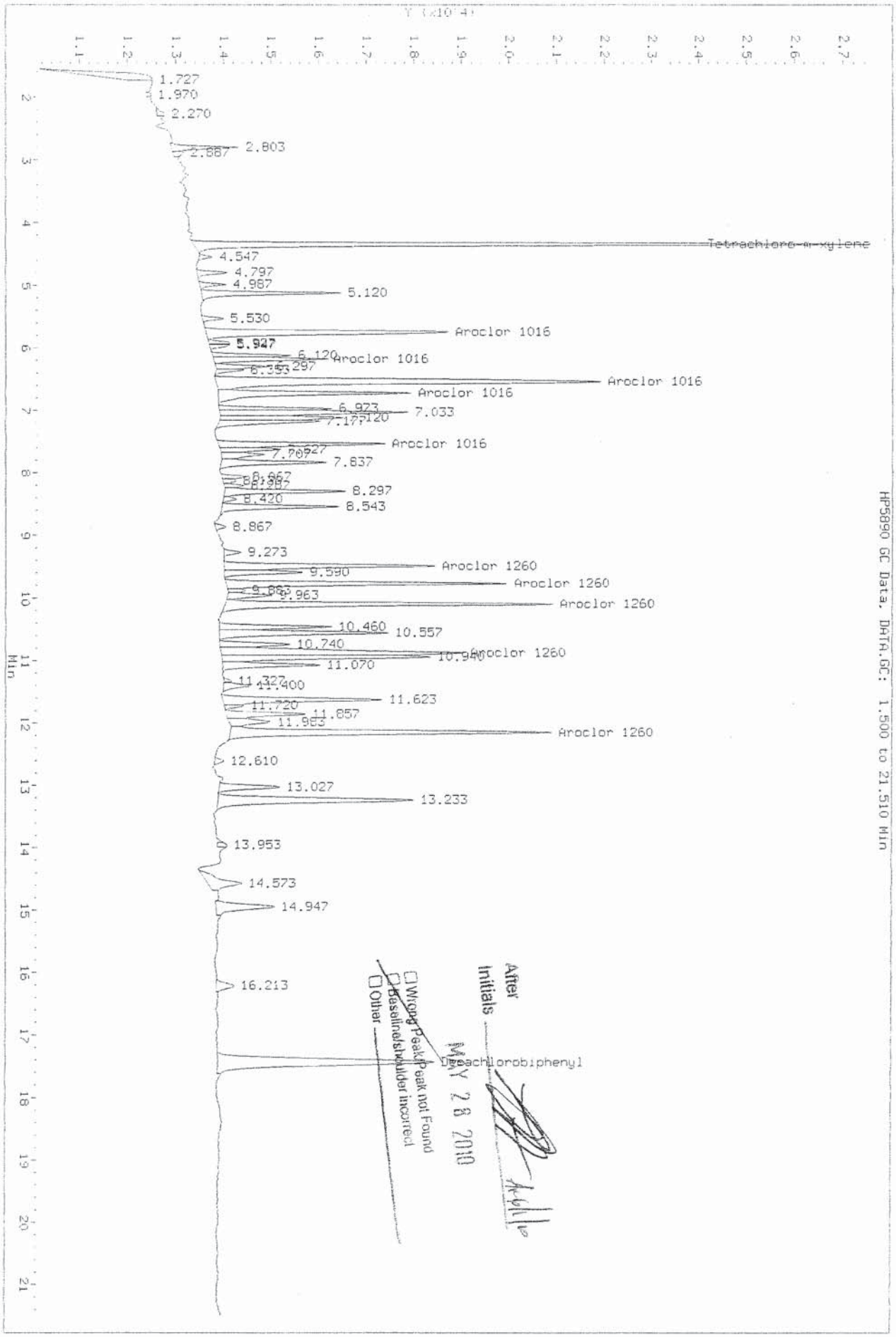
Injection Date: 28-MAY-2010 00:19
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.510 Min
Y (x10^4)
Min
1.727
1.970
2.270
2.803
Tetrachloro-m-xylene
4.547
4.797
4.987
5.120
5.530
Aroclor 1016
5.927
6.120
Aroclor 1016
Aroclor 1016
Aroclor 1016
6.973
7.033
Aroclor 1016
7.837
8.297
8.420
8.543
8.867
9.273
Aroclor 1260
9.590
Aroclor 1260
Aroclor 1260
10.460
10.557
10.740
Aroclor 1260
11.070
11.400
11.623
11.720
11.857
Aroclor 1260
12.610
13.027
13.233
13.953
14.573
14.947
16.213
Decachlorobiphenyl
Before
MAY 28 2010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Injection Date: 28-MAY-2010 00:19
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.510 Min
Y (x10^4)
Min
1.727
1.970
2.270
2.803
Tetrachloro-m-xylene
4.547
4.797
4.967
5.120
5.530
Aroclor 1016
5.927
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.033
Aroclor 1016
7.837
8.297
8.543
8.867
9.273
Aroclor 1260
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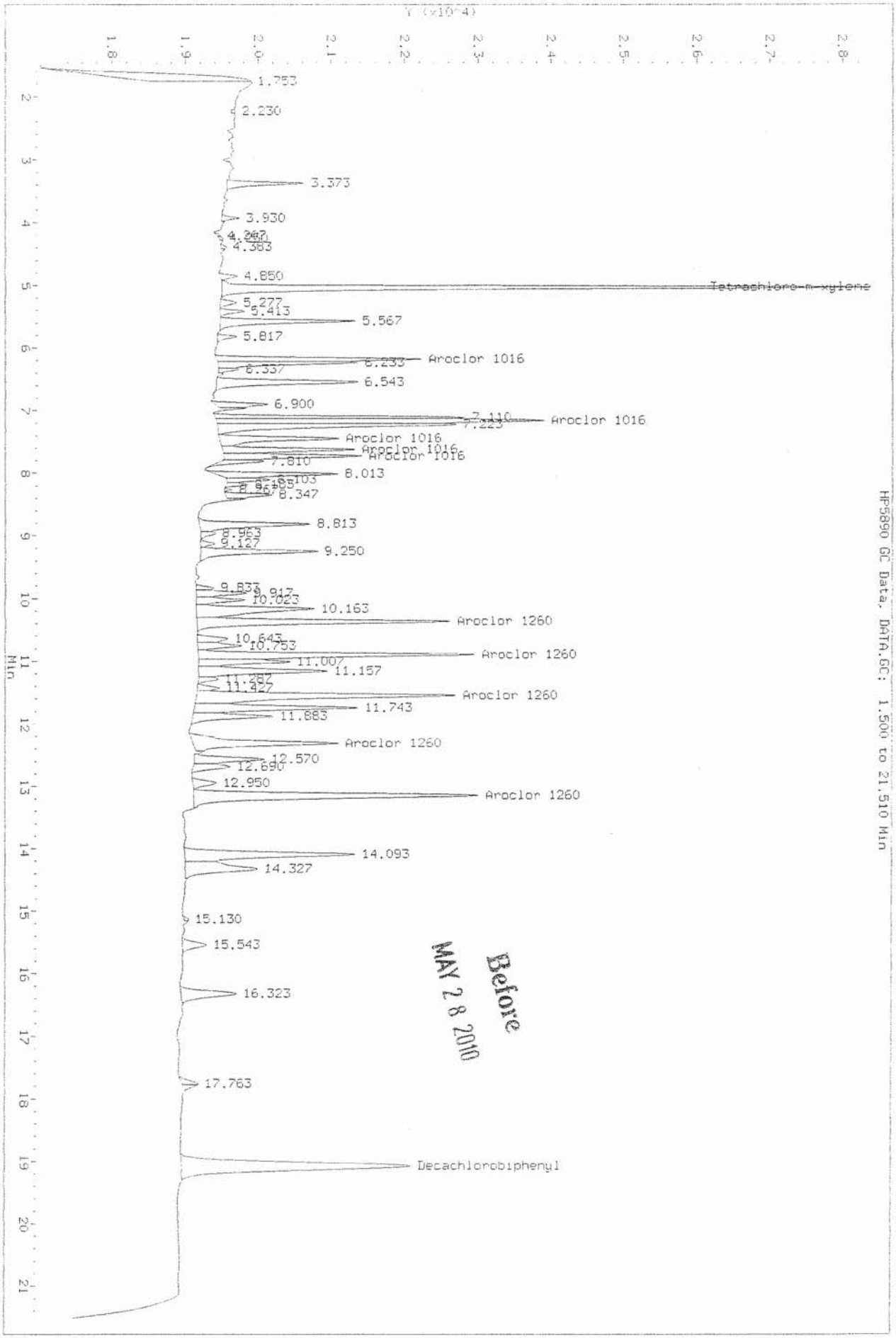
9.963
Aroclor 1260
10.460
10.557
10.740
Aroclor 1260
11.070
11.400
11.623
11.720
11.857
Aroclor 1260
12.610
13.027
13.233
13.953
14.573
14.947
16.213
Decachlorobiphenyl
After
Initials
MAY 28 2010
Wrong Peak
Peak not Found
Baseline/shoulder incorrect
Other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Injection Date: 28-MAY-2010 00:19
Instrument: GC09.i
Client Sample ID:
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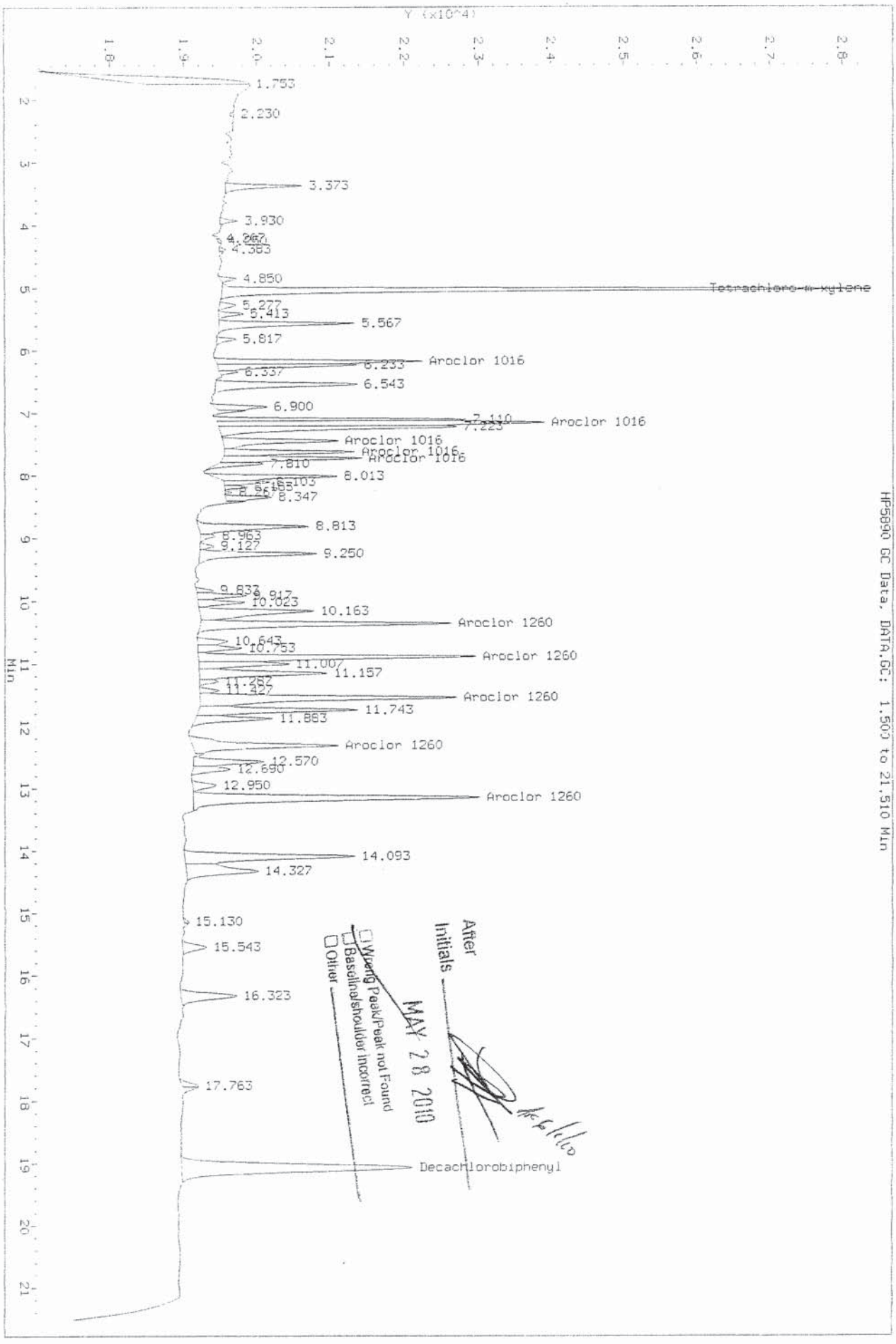
Y (x10^4)
Min
1.753
2.230
3.373
3.930
4.207
4.383
4.850
Tetrachloro-m-xylene
5.277
5.413
5.567
5.817
6.233
Aroclor 1016
6.337
6.543
6.900
7.110
Aroclor 1016
7.223
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.810
8.013
8.103
8.347
8.813
8.963
9.127
9.250
9.833
9.917
10.023
10.163
Aroclor 1260
10.643
10.753
Aroclor 1260
11.007
11.157
11.387
11.427
Aroclor 1260
11.743
11.883
Aroclor 1260
12.570
12.690
12.950
Aroclor 1260
14.093
14.327
15.130
15.543
16.323
17.763
Decachlorobiphenyl
Before
MAY 28 2010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Injection Date: 28-MAY-2010 00:19
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Y (x10^4)
Min
1.753
2.230
3.373
3.930
4.267
4.383
4.850
Tetrachloro-m-xylene
5.277
5.413
5.567
5.817
6.233
Aroclor 1016
6.337
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7.223
Aroclor 1016
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.810
8.013
8.105
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Aroclor 1260
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10.753
Aroclor 1260
11.007
11.157
11.282
11.427
Aroclor 1260
11.743
11.883
Aroclor 1260
12.570
12.690
12.950
Aroclor 1260
14.093
14.327
15.130
15.543
16.323
17.763
Decachlorobiphenyl
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F024.D
Report Date: 28-May-2010 17:28

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F024.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R024.D
Inj Date : 28-MAY-2010 00:46
Sample Info: 1016/1260 @ 500ppb | PCB5-55H
Misc Info :
Cal Date : 28-MAY-2010 13:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.343	5.030	358927	302772	50.5	51.4		100.00
Aroclor 1016	6.180	6.543	76152	95206	340	559	80.00- 120.00	100.00
	6.723	7.163	156129	161123	715	498	162.91- 244.36	205.02
	7.030	7.447	130532	86747	304	549	145.03- 217.55	171.41
	7.540	7.623	106227	70150	427	537	115.01- 172.51	139.49
	7.837	7.723	81220	79490	467	539	76.85- 115.27	106.66
			Average of Peak Amounts =		451	536		
Aroclor 1260	9.487	10.363	154171	145100	510	523	80.00- 120.00	100.00
	9.770	10.893	186470	153968	504	519	99.64- 149.46	120.95
	10.093	11.547	227237	191734	502	515	122.12- 183.18	147.39
	10.873	12.310	226727	126592	523	538	109.38- 164.06	147.06
	12.143	13.147	300748	224610	527	511	135.68- 203.52	195.07
			Average of Peak Amounts =		513	521		
Decachlorobiphenyl	17.433	19.077	301294	264364	51.0	51.5		100.00

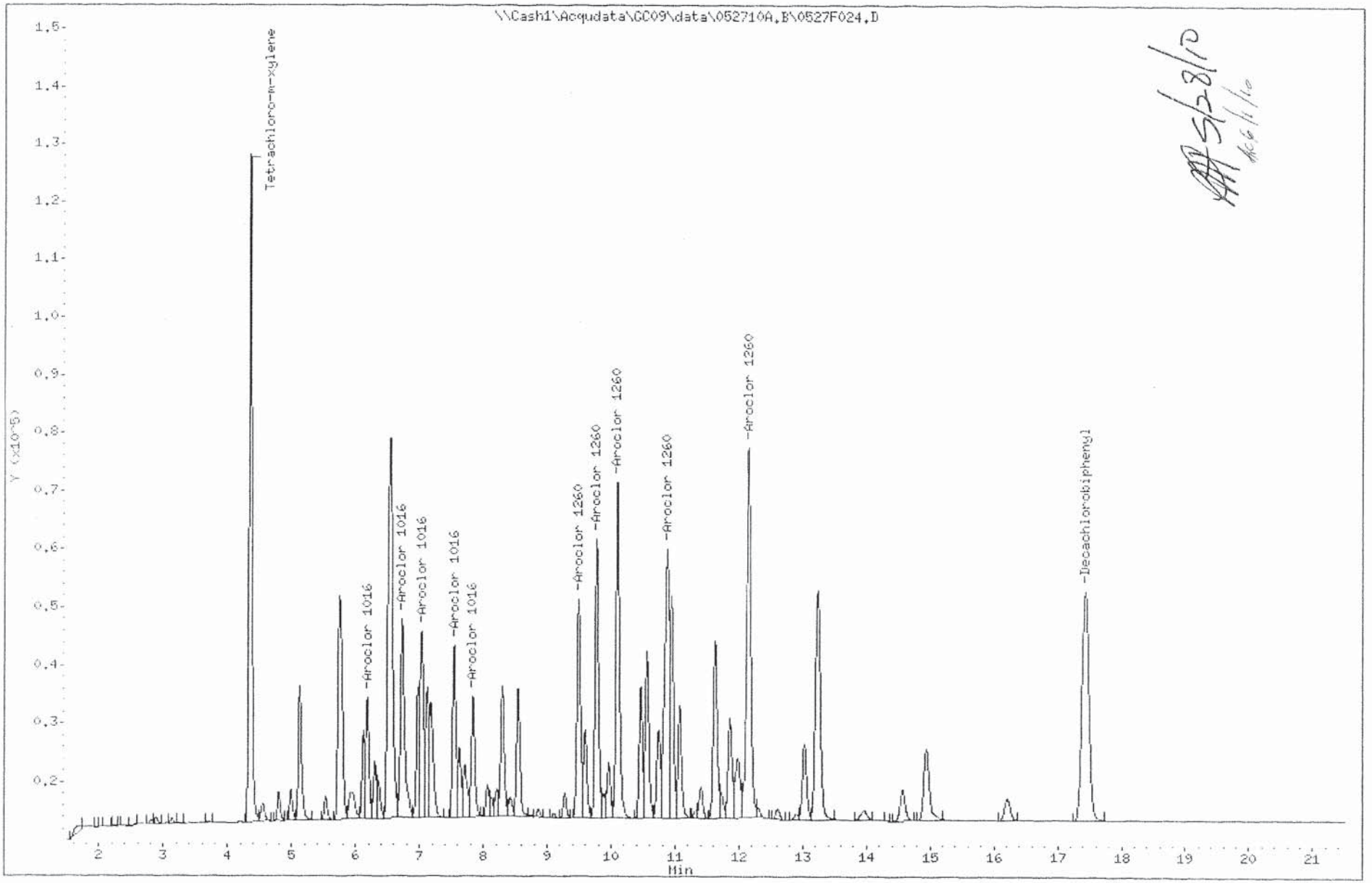
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F024.D
Date : 28-MAY-2010 00:46
Client ID:
Sample Info: 1016/1260 @ 500ppb | PCB5-55H

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

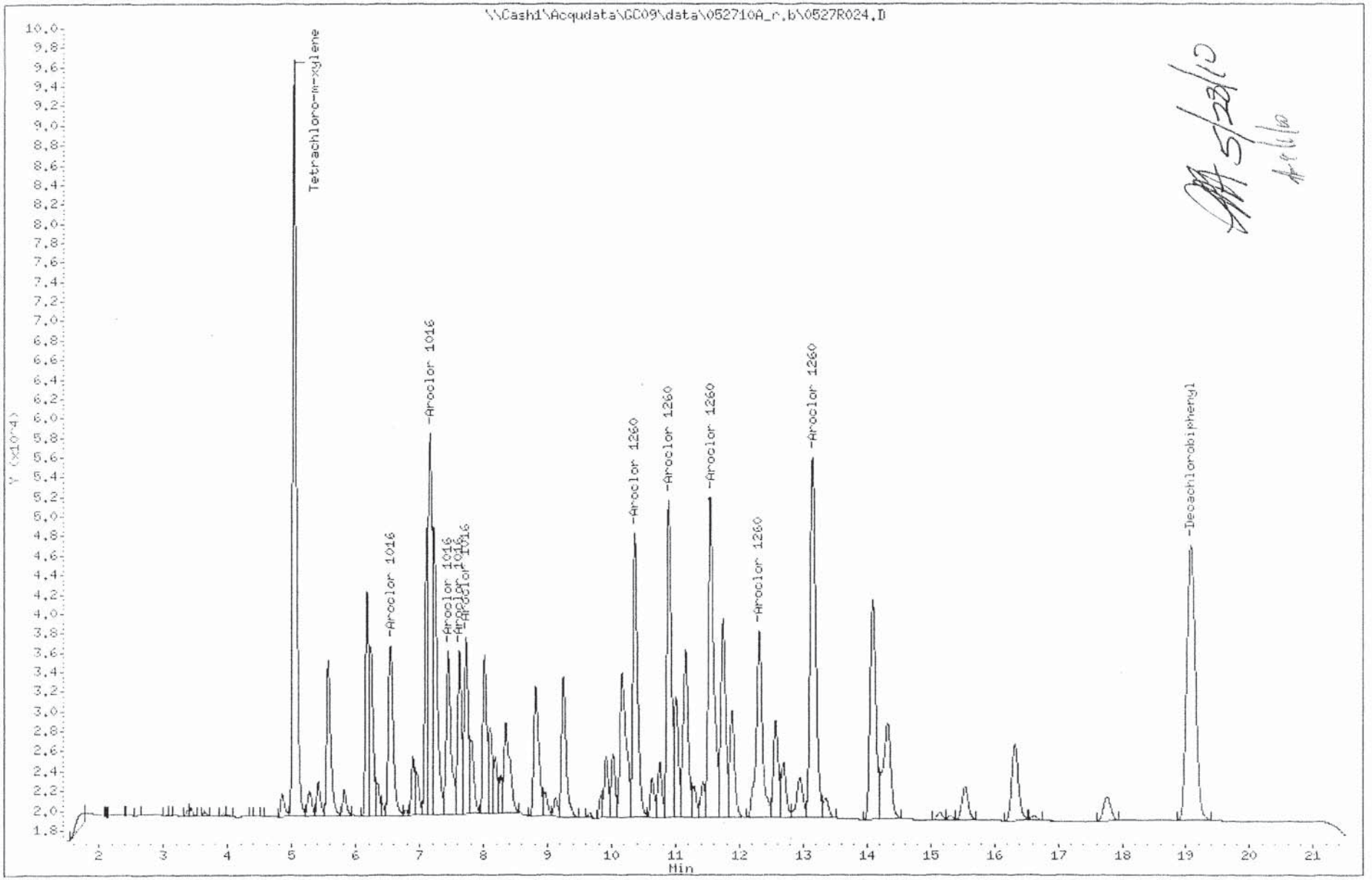
Column phase: DB-35MS

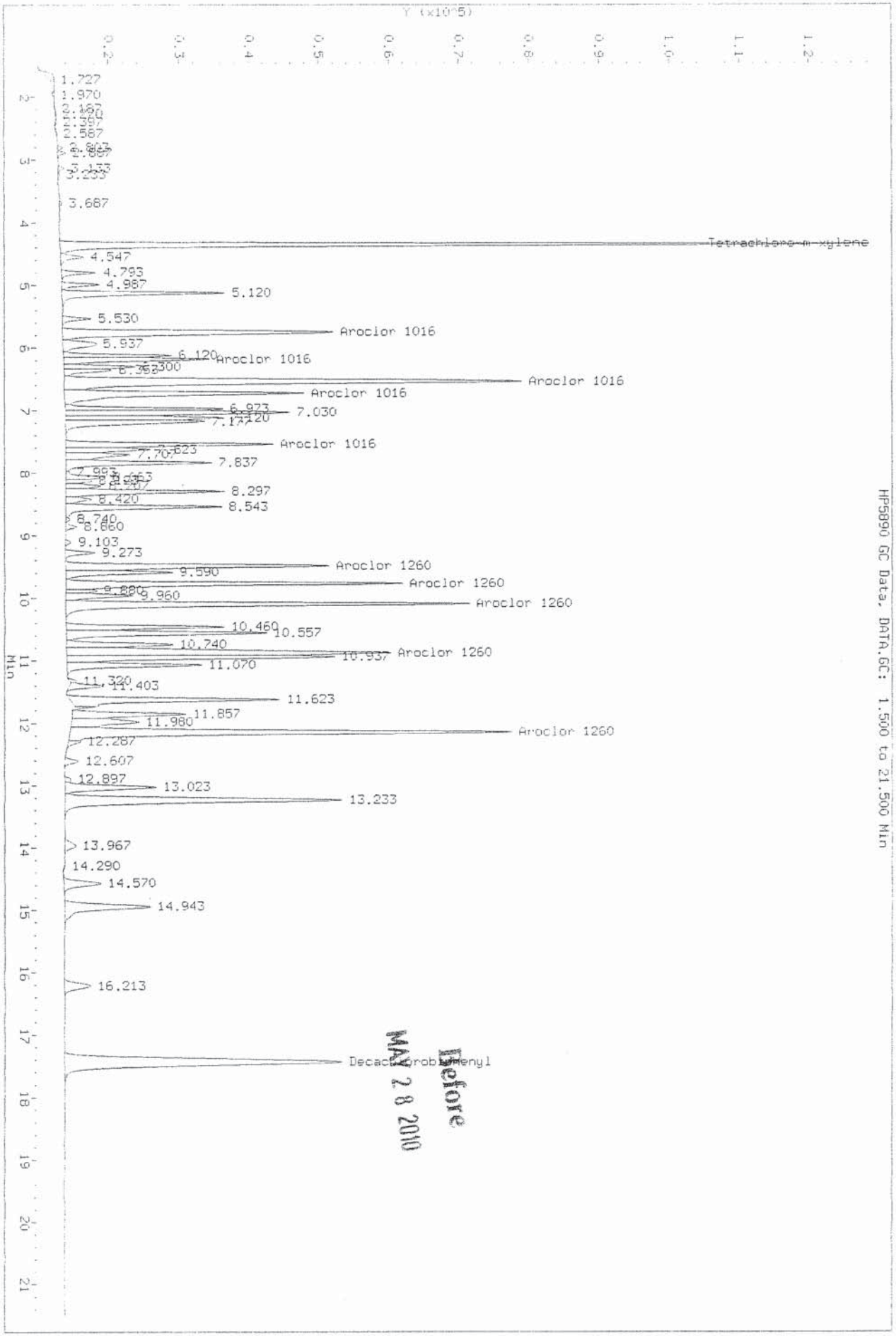


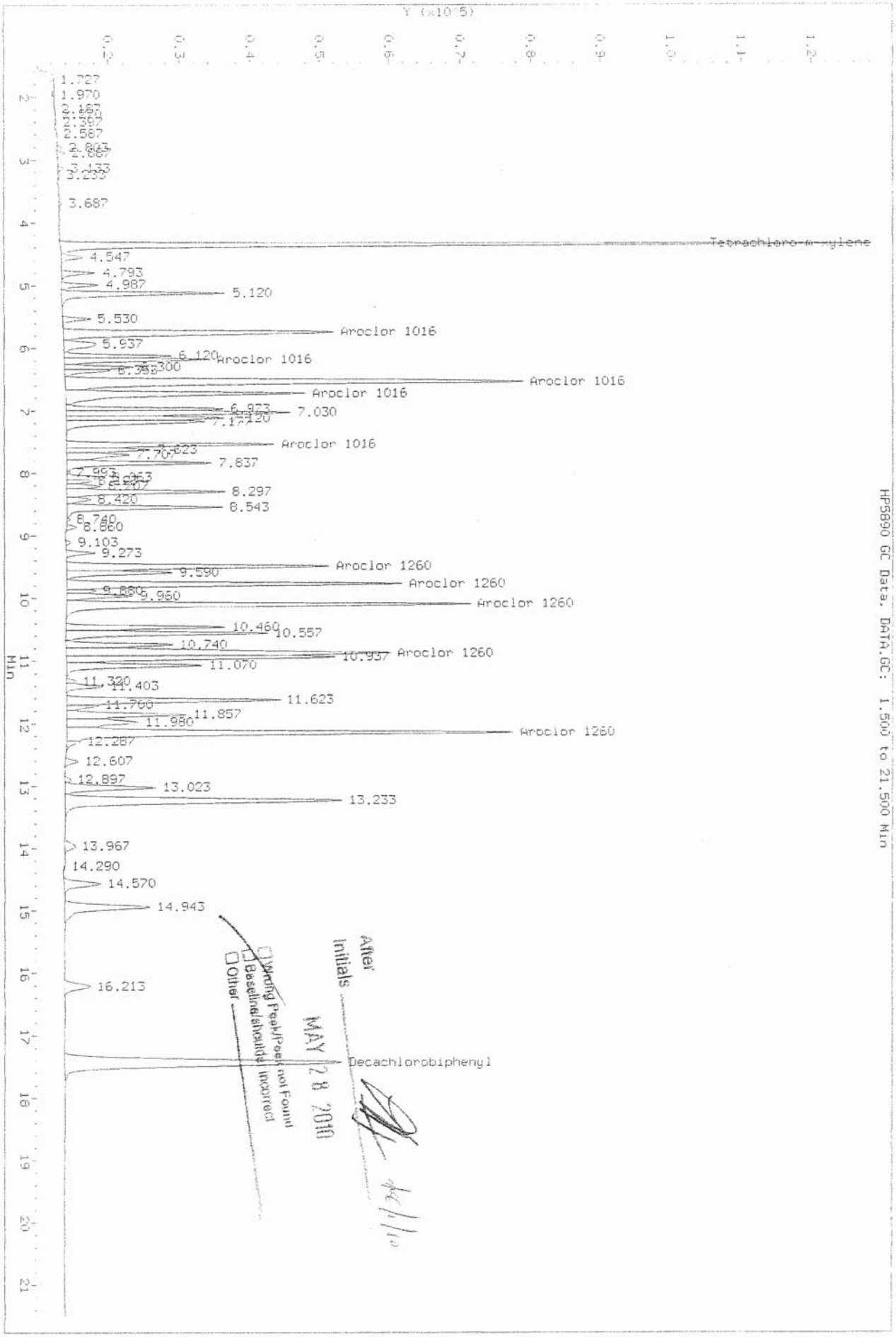
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R024.D
Date : 28-MAY-2010 00:46
Client ID:
Sample Info: 1016/1260 @ 500ppb | PCB5-55H

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F024.D
Injection Date: 28-MAY-2010 00:46
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.500 Min
Y (x10^5)
Min
Tetrachloro-m-xylene
Aroclor 1016
Aroclor 1260
Decachlorobiphenyl
MAY 28 2010
Before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Injection Date: 28-MAY-2010 00:46
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.500 Min
Y (x10^5)
Min
Tetrachloro-m-xylene
Aroclor 1016
Aroclor 1260
Decachlorobiphenyl
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline should be
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F025.D
Report Date: 28-May-2010 17:28

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F025.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R025.D
Inj Date : 28-MAY-2010 01:12
Sample Info: 1016/1260 @ 1000ppb | PCB5-55I
Misc Info :
Cal Date : 28-MAY-2010 13:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

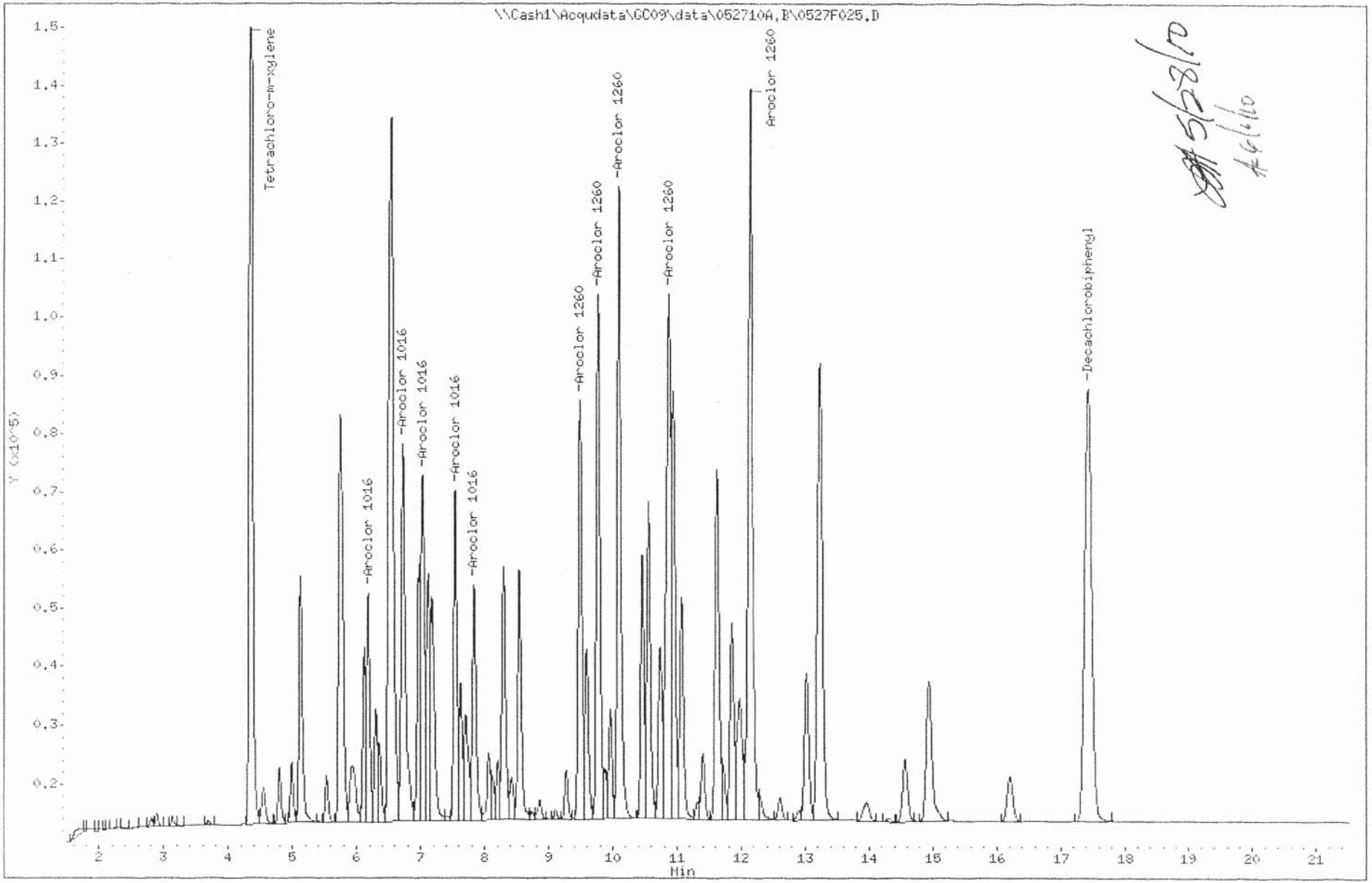
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.343	5.030	679365	602860	95.6	102		100.00
Aroclor 1016	6.180	6.543	143205	187924	747	1100	80.00- 120.00	100.00
	6.723	7.160	299231	319993	1220	989	162.91- 244.36	208.95
	7.030	7.447	234428	175921	649	1110	145.03- 217.55	163.70
	7.540	7.620	204741	136724	880	1050	115.01- 172.51	142.97
	7.837	7.723	162641	156483	975	1060	76.85- 115.27	113.57
			Average of Peak Amounts =		894	1060		
Aroclor 1260	9.487	10.363	291968	280711	966	1010	80.00- 120.00	100.00
	9.767	10.890	350576	296240	948	998	99.64- 149.46	120.07
	10.093	11.543	432194	381715	954	1020	122.12- 183.18	148.03
	10.870	12.310	439578	255621	1010	1080	109.38- 164.06	150.56
	12.143	13.143	592535	449818	1040	1020	135.68- 203.52	202.95
			Average of Peak Amounts =		984	1020		
Decachlorobiphenyl	17.433	19.077	579519	518864	98.2	101		100.00

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F025.D
Date : 28-MAY-2010 01:12
Client ID:
Sample Info: 1016/1260 @ 1000ppb | PCB5-55I

Instrument: GC09.i

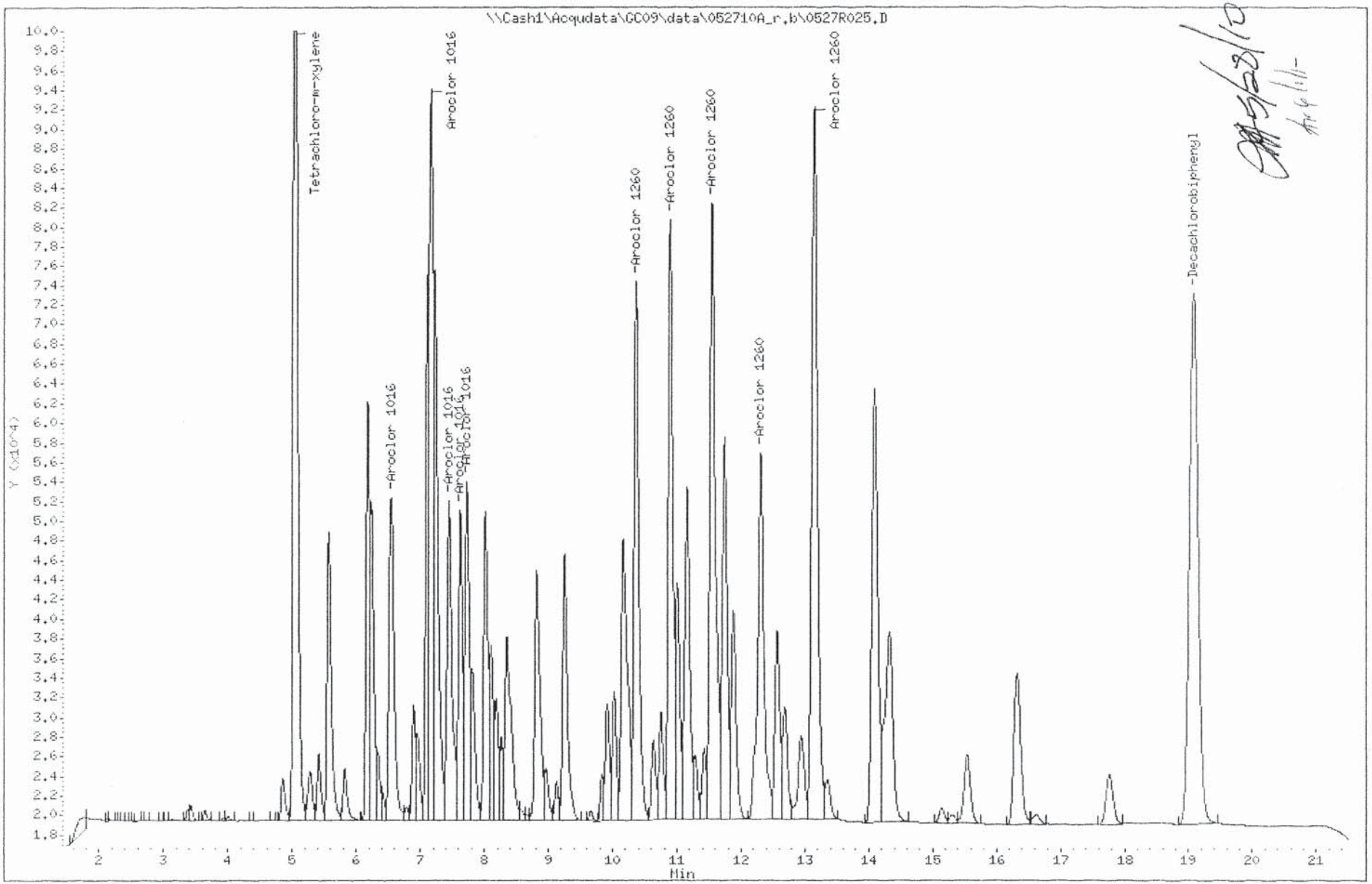
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R025.D
Date : 28-MAY-2010 01:12
Client ID:
Sample Info: 1016/1260 @ 1000ppb | PCB5-55I

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



742

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F026.D
Report Date: 28-May-2010 17:28

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F026.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R026.D
Inj Date : 28-MAY-2010 01:38
Sample Info: 1016/1260 @ 2000ppb | PCB5-55J
Misc Info :
Cal Date : 28-MAY-2010 13:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.343	5.030	1221823	1124897	172	191		100.00
Aroclor 1016	6.177	6.543	255219	346420	1560	1960	80.00- 120.00	100.00
	6.720	7.160	530222	581827	1980	1800	162.91- 244.36	207.75
	7.030	7.443	413223	325655	1370	2060	145.03- 217.55	161.91
	7.540	7.620	370775	248279	1690	1900	115.01- 172.51	145.28
	7.833	7.723	298915	278972	1860	1890	76.85- 115.27	117.12
			Average of Peak Amounts =		1690	1920		
Aroclor 1260	9.487	10.360	534792	514161	1770	1850	80.00- 120.00	100.00
	9.767	10.890	633380	539881	1710	1820	99.64- 149.46	118.43
	10.093	11.543	786956	716222	1740	1920	122.12- 183.18	147.15
	10.870	12.307	825913	489934	1910	2080	109.38- 164.06	154.44
	12.140	13.143	1092934	855710	1910	1950	135.68- 203.52	204.37
			Average of Peak Amounts =		1810	1920		
Decachlorobiphenyl	17.433	19.077	1052797	957927	178	187		100.00

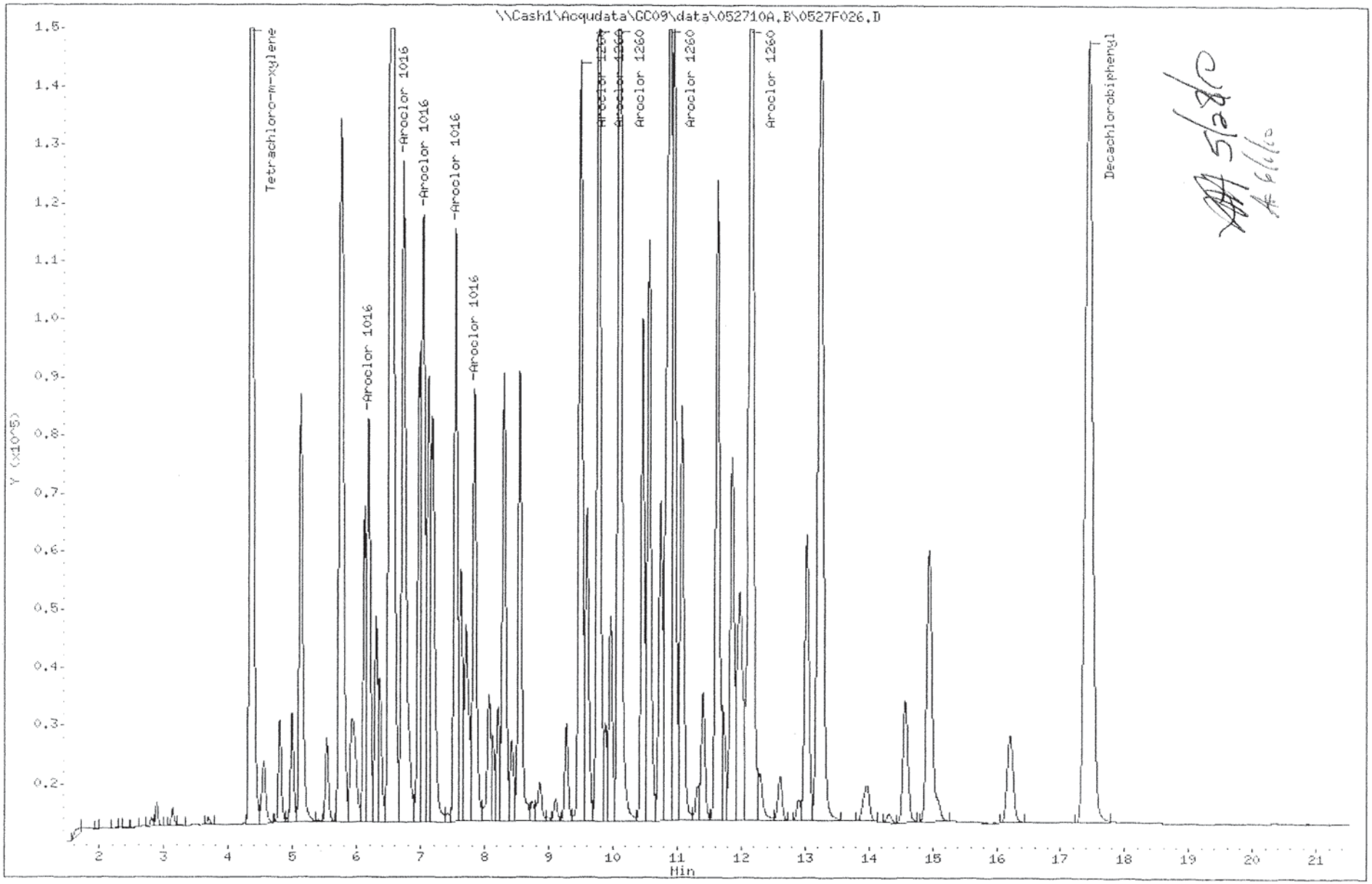
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F026.D
Date : 28-MAY-2010 01:38
Client ID:
Sample Info: 1016/1260 @ 2000ppb | PCB5-55J

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-35MS

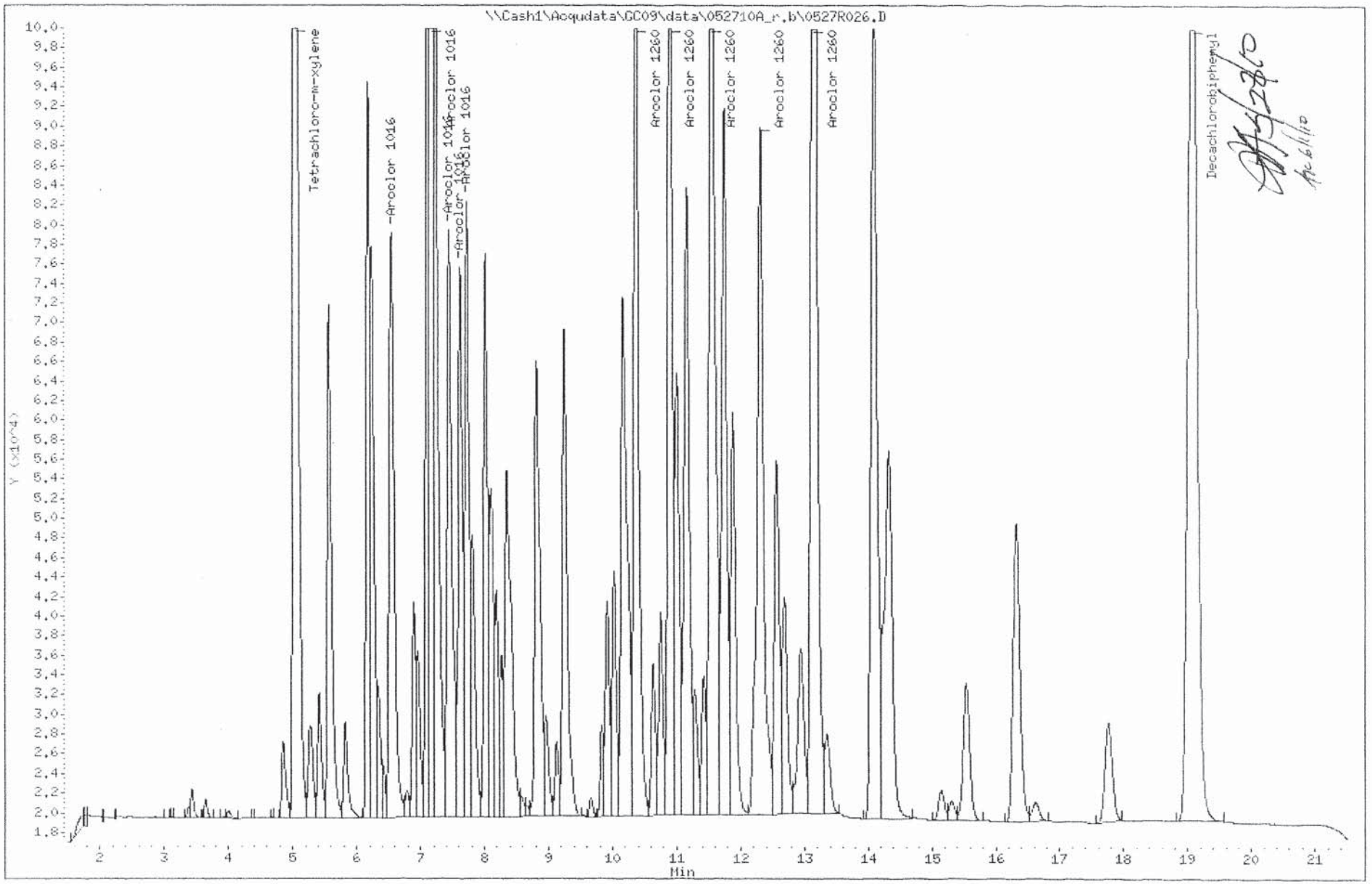


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R026.D
Date : 28-MAY-2010 01:38
Client ID:
Sample Info: 1016/1260 @ 2000ppb | PCB5-55J

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F027.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F027.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R027.D
Inj Date : 28-MAY-2010 02:04
Sample Info: 1016/1260 @ 5000ppb | PCB5-55K
Misc Info :
Cal Date : 28-MAY-2010 13:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

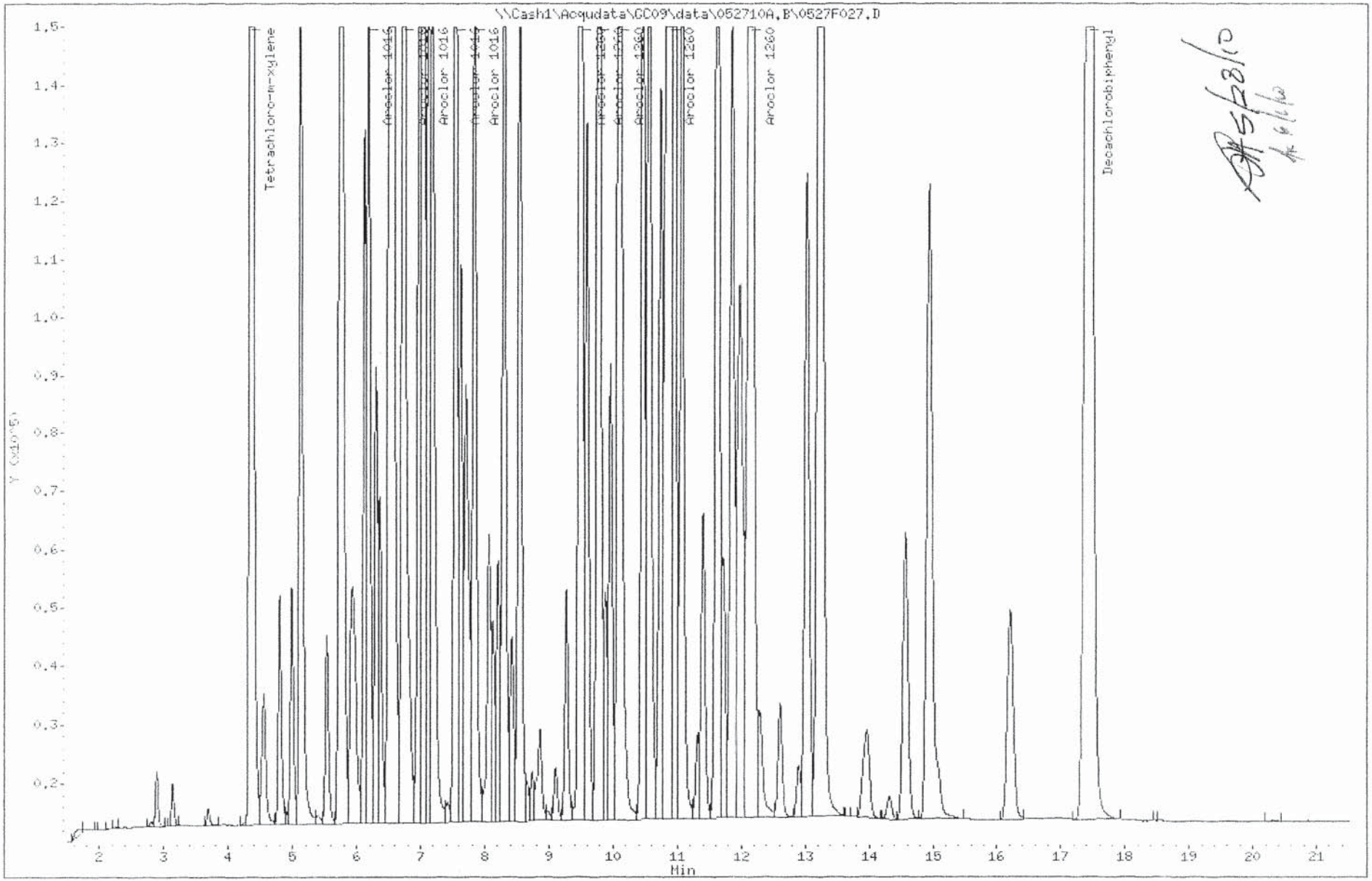
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.343	5.030	2611859	2519566	368	428		100.00
Aroclor 1016	6.177	6.543	543843	772733	3840	4220	80.00- 120.00	100.00
	6.720	7.160	1121268	1276275	3920	3950	162.91- 244.36	206.17
	7.030	7.447	878662	742029	3520	4700	145.03- 217.55	161.57
	7.537	7.620	792247	525888	3800	4020	115.01- 172.51	145.68
	7.833	7.723	658716	625893	4210	4240	76.85- 115.27	121.12
			Average of Peak Amounts =		3860	4230		
Aroclor 1260	9.483	10.360	1177390	1159016	3890	4180	80.00- 120.00	100.00
	9.767	10.890	1370255	1216528	3700	4100	99.64- 149.46	116.38
	10.093	11.543	1708055	1670009	3770	4490	122.12- 183.18	145.07
	10.870	12.307	1867377	1167982	4310	4960	109.38- 164.06	158.60
	12.140	13.143	2477559	2003451	4340	4560	135.68- 203.52	210.43
			Average of Peak Amounts =		4000	4460		
Decachlorobiphenyl	17.427	19.073	2333876	2128229	395	415		100.00

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F027.D
Date : 28-MAY-2010 02:04
Client ID:
Sample Info: 1016/1260 @ 5000ppb | PCB5-55K

Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

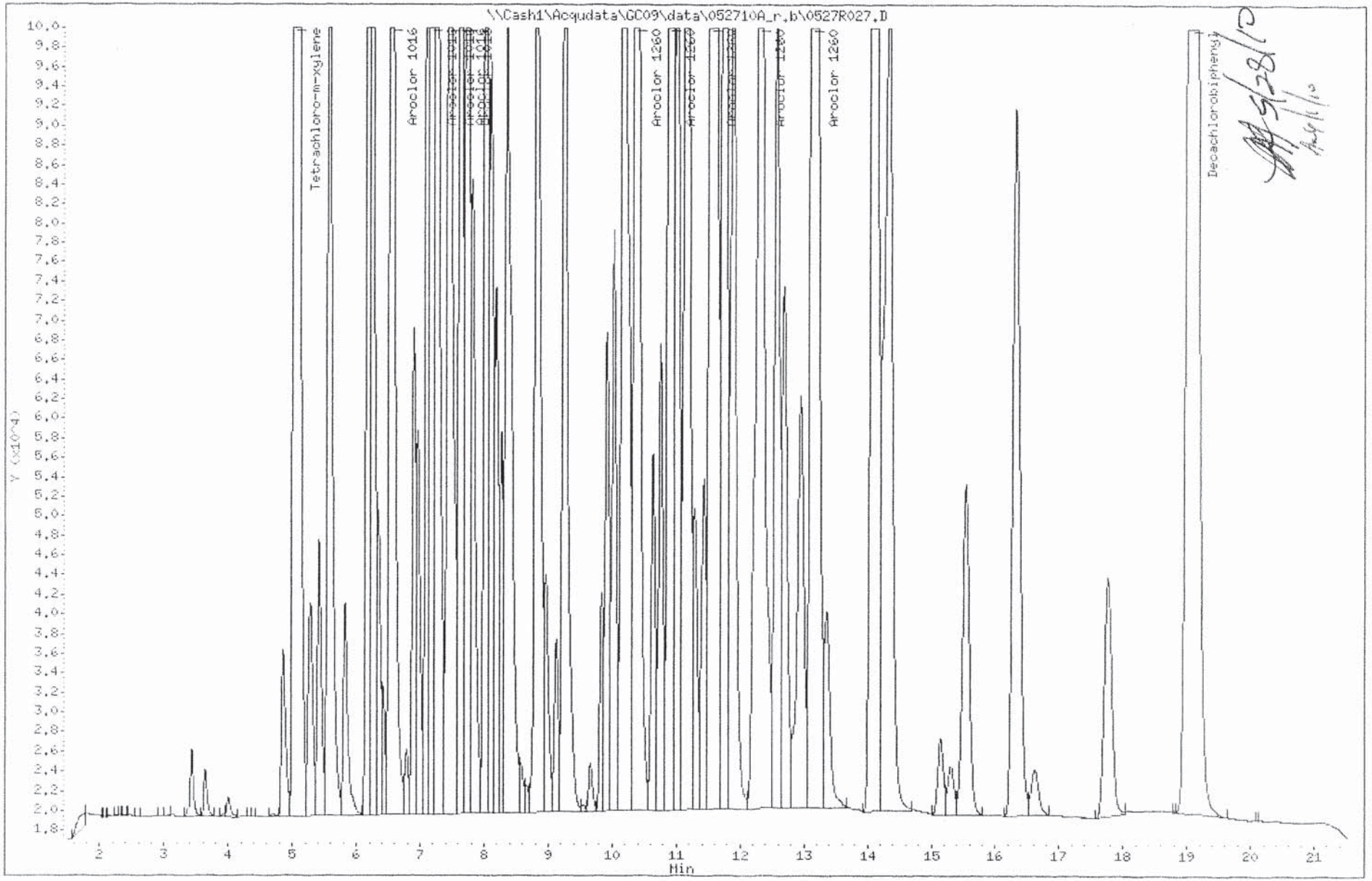


Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R027.D
Date : 28-MAY-2010 02:04
Client ID:
Sample Info: 1016/1260 @ 5000ppb | PCB5-55K

Instrument: GC09,i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F028.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F028.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R028.D
Inj Date : 28-MAY-2010 02:30
Sample Info: 1221/1254 @ 50/25ppb | PCB5-52D
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1221+1254.sub
Sub List #2 : 1221+1254.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.687	4.687	4988	1123	68.5	22.8	80.00- 120.00	100.00(M)
	4.543	4.853	2363	1715	57.9	48.6	51.26- 76.90	47.39(M)
	4.797	5.283	6838	4308	64.9	53.8	133.68- 200.52	137.09(M)
	4.983	5.410	4212	2937	62.3	52.4	85.57- 128.36	84.44(M)
			Average of Peak Amounts =		63.4	44.4		
Aroclor 1254	8.543	9.250	11427	11581	27.9	30.1	80.00- 120.00	100.00(M)
	9.270	9.833	7602	4384	27.1	24.9	61.91- 92.86	66.53(M)
	9.477	10.027	16454	9244	28.7	25.8	116.61- 174.92	143.99(M)
	9.887	10.363	12788	3643	27.4	22.6	101.15- 151.73	111.90(M)
	10.097	10.640	6746	8716	27.0	27.7	53.30- 79.94	59.03(M)
			Average of Peak Amounts =		27.6	26.2		

QC Flag Legend

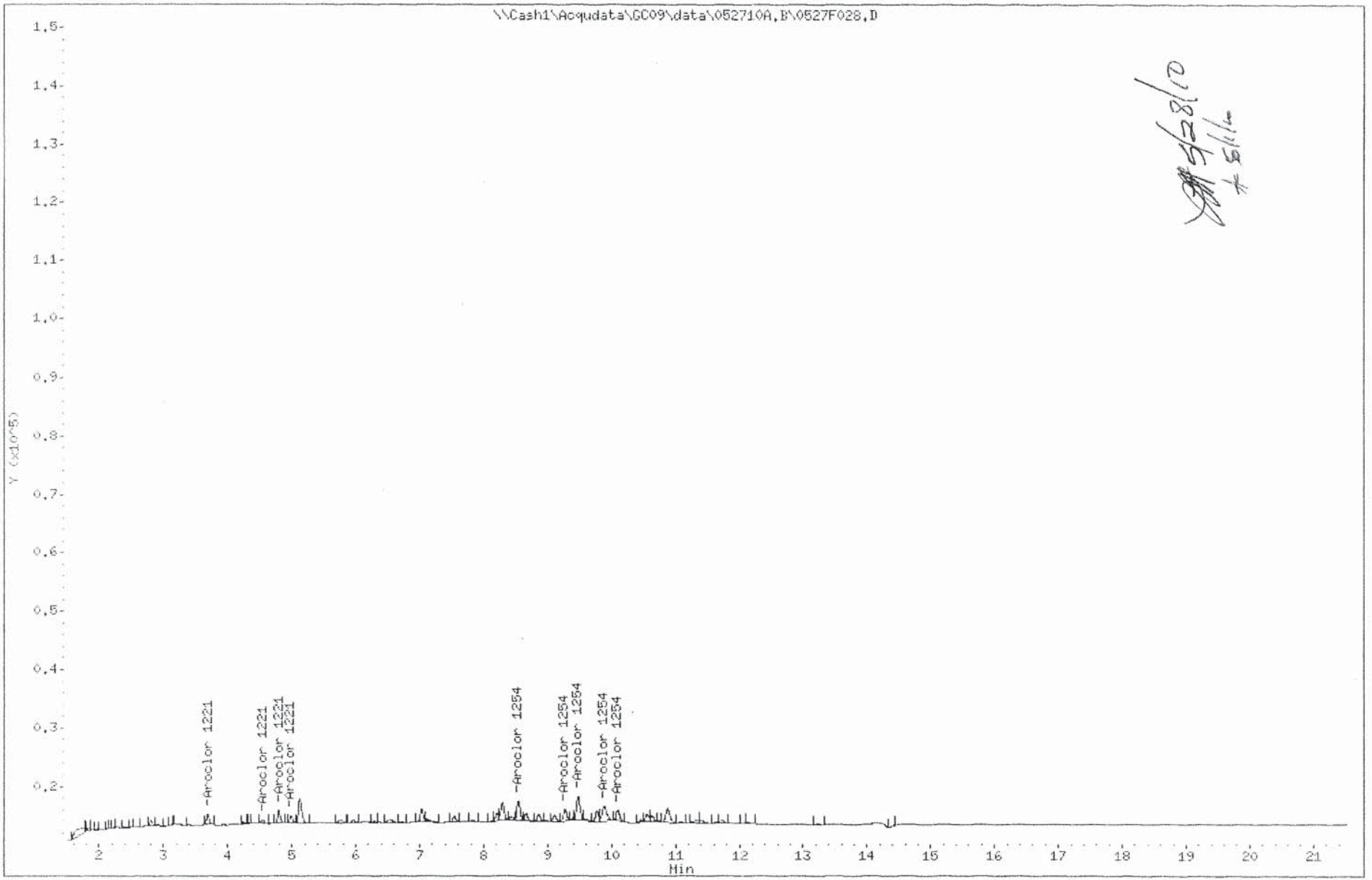
M - Compound response manually integrated.

5/28/10

4/6/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F028.D
Date : 28-MAY-2010 02:30
Client ID:
Sample Info: 1221/1254 @ 50/25ppb | PCB5-52D

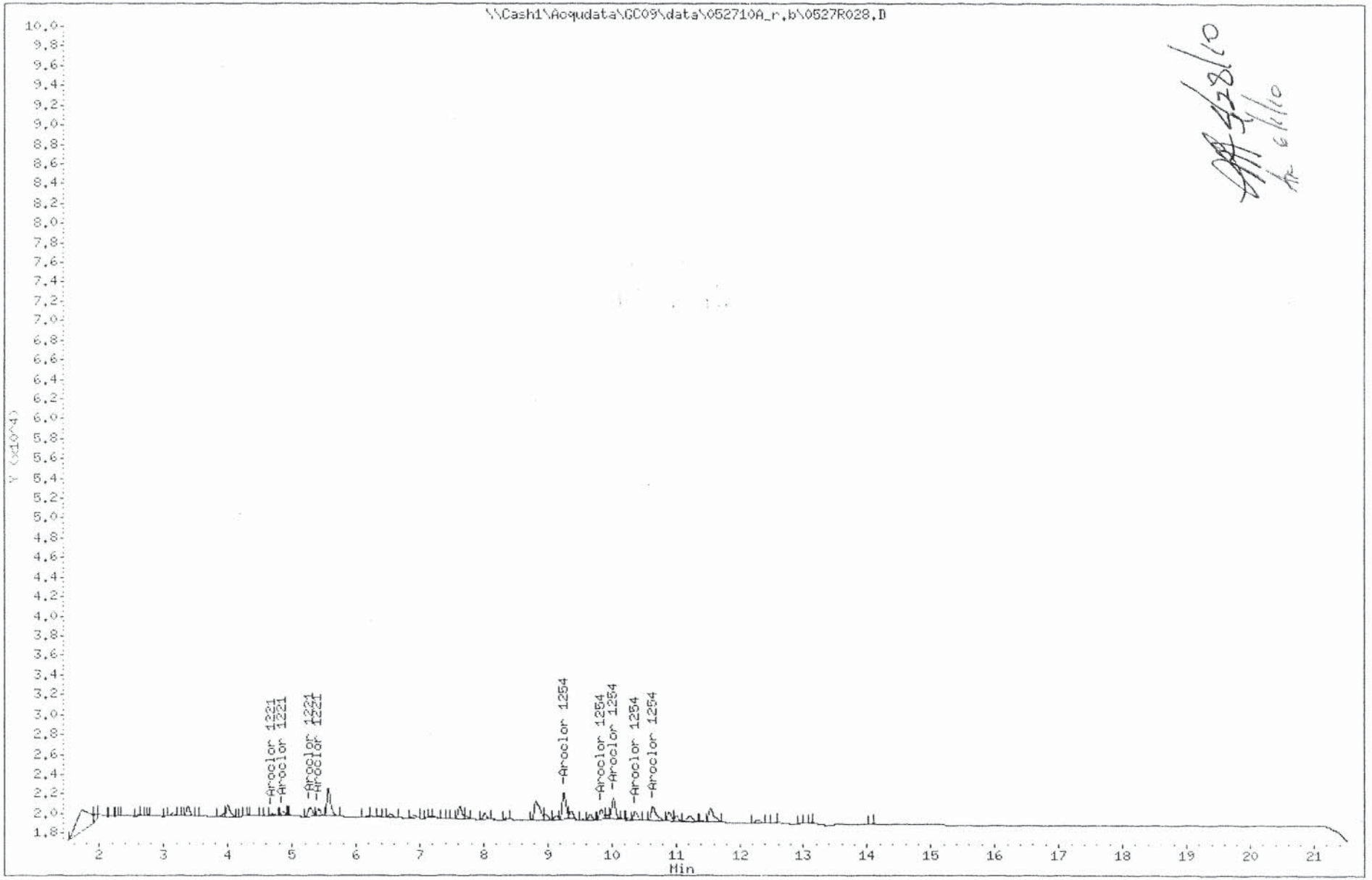
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

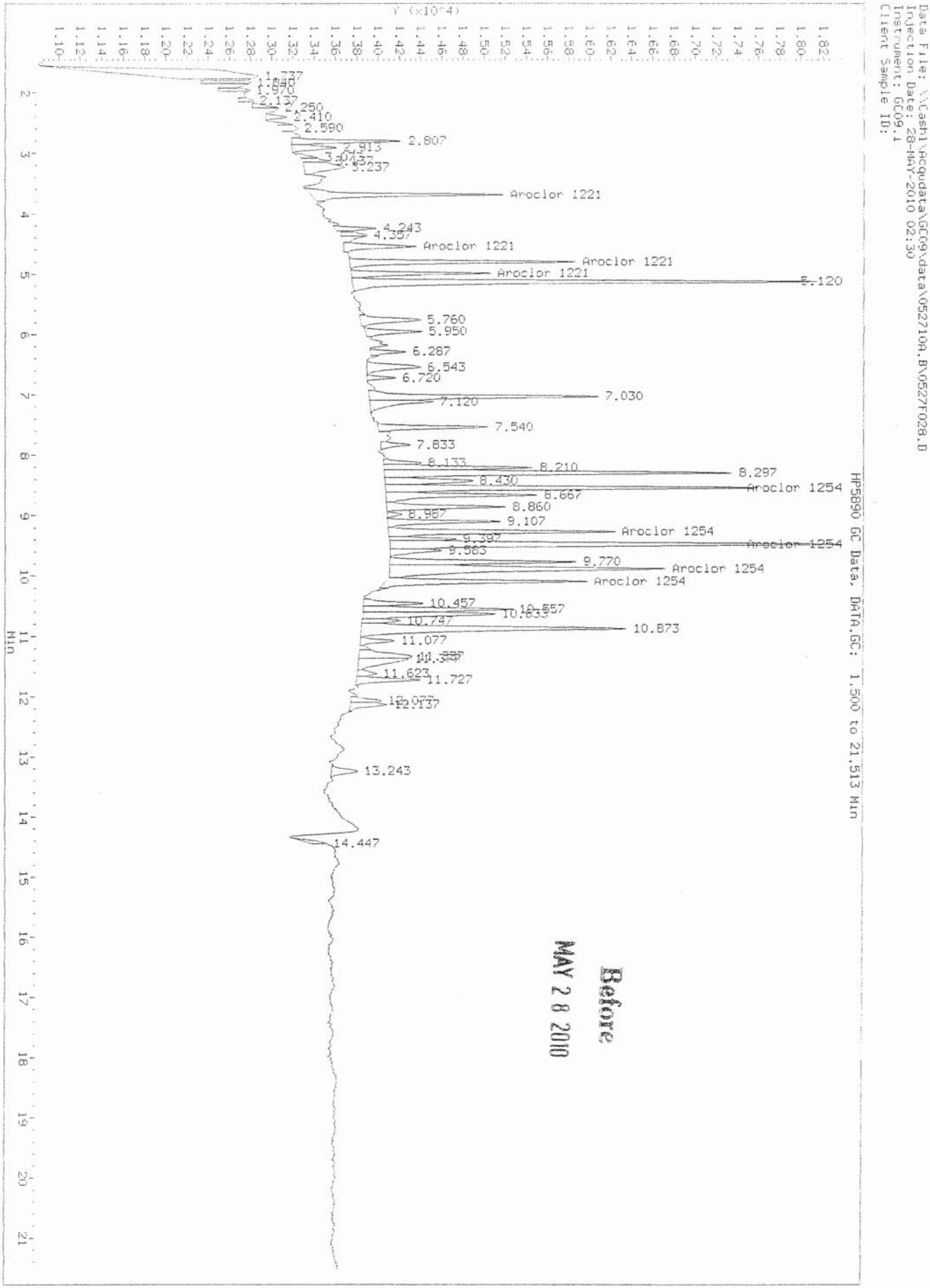


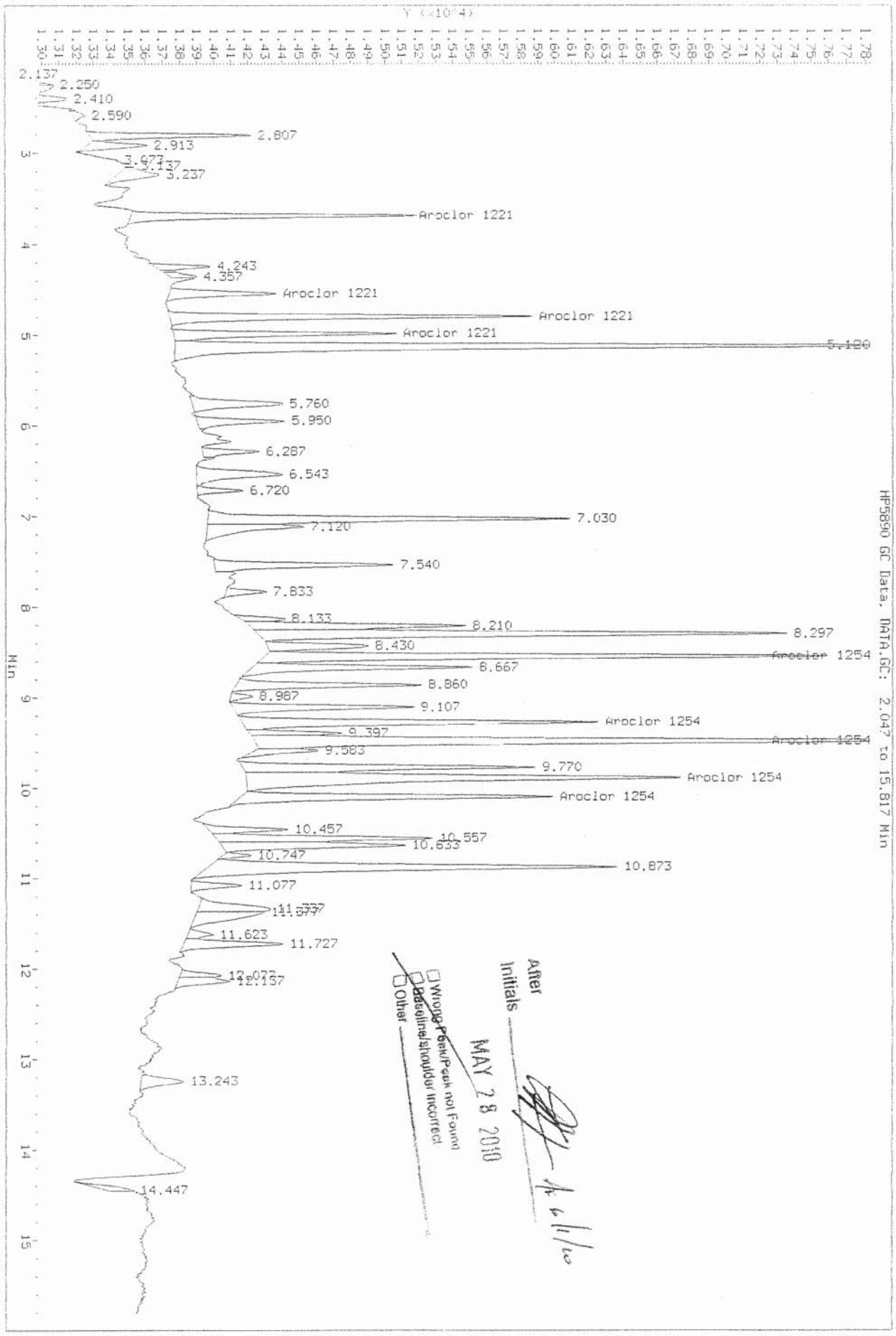
Data File: \\Cash1\Aoqudata\GC09\data\052710A_r.b\0527R028.D
Date : 28-MAY-2010 02:30
Client ID:
Sample Info: 1221/1254 @ 50/25ppb | PCB5-52D

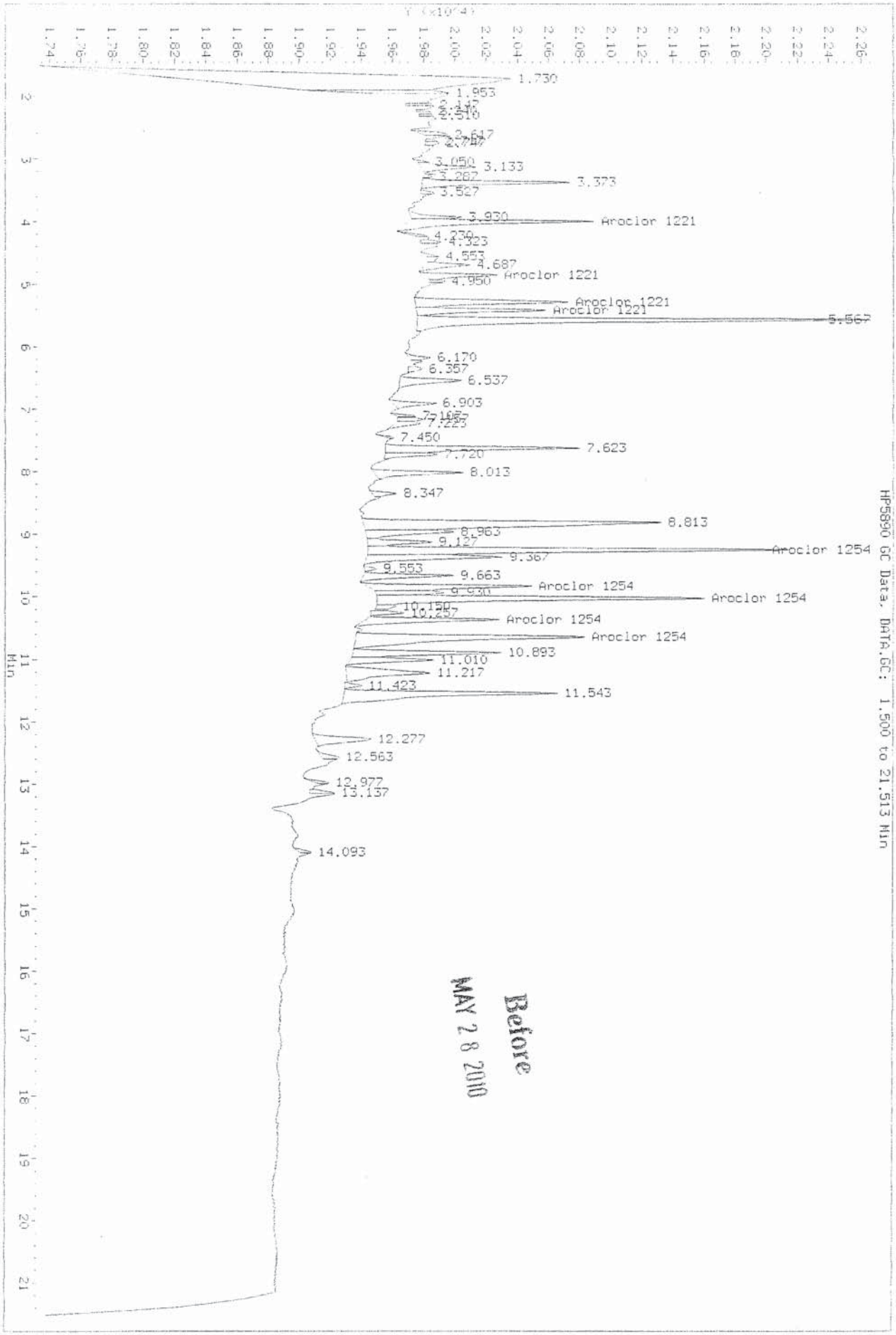
Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

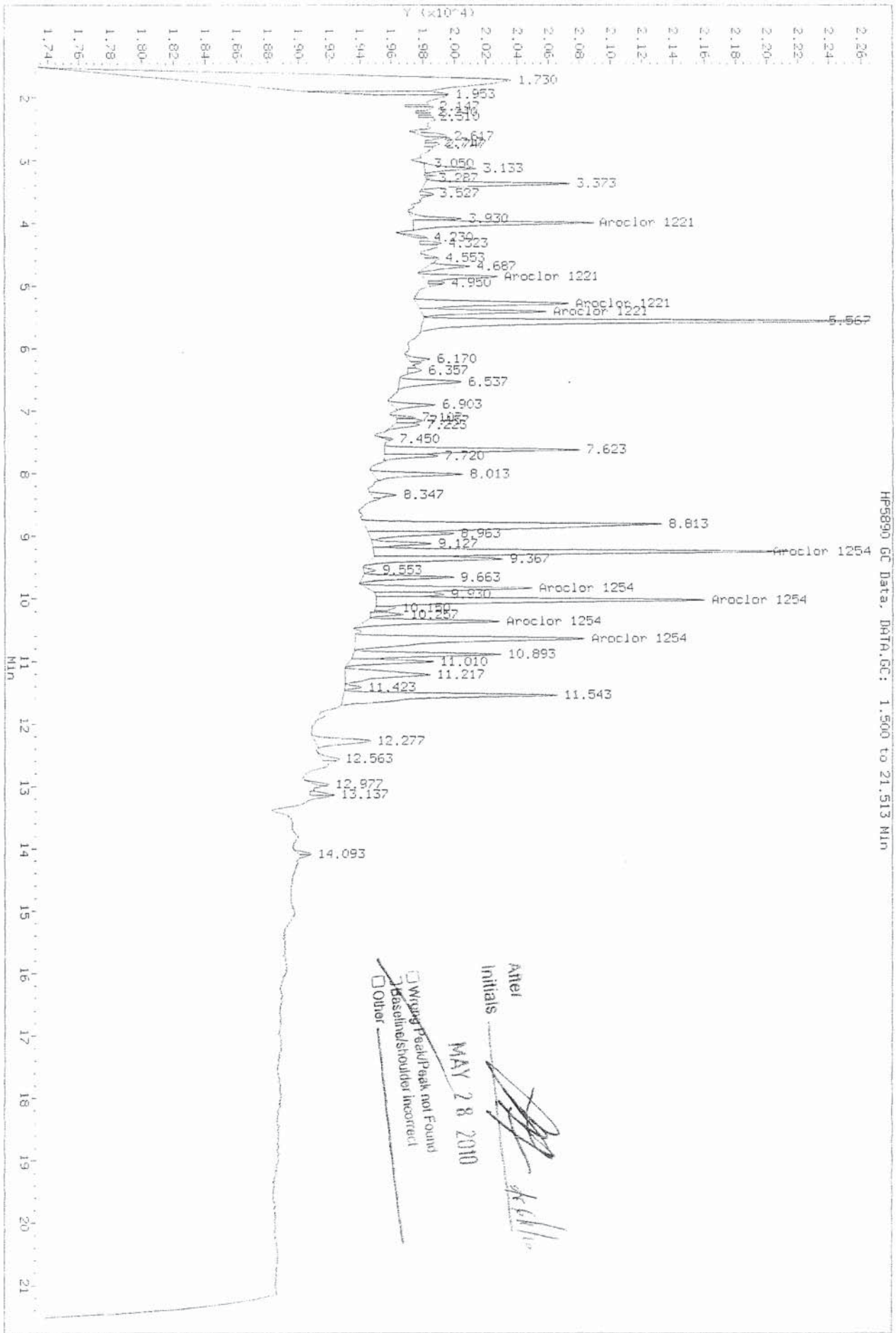



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F028.D
Injection Date: 28-MAY-2010 02:30
Instrument: GC09.1
Client Sample ID:
Y (x10^4)
Min
HP5890 GC Data, DATA.GC: 1.500 to 21.513 Min
2.807
Aroclor 1221
4.243
4.357
Aroclor 1221
Aroclor 1221
Aroclor 1221
5.120
5.760
5.950
6.287
6.543
6.720
7.030
7.120
7.540
7.833
8.133
8.210
8.297
8.430
Aroclor 1254
8.667
8.860
8.987
9.107
Aroclor 1254
9.397
Aroclor 1254
9.563
9.770
Aroclor 1254
Aroclor 1254
10.457
10.557
10.747
10.873
11.077
11.623
11.727
13.243
14.447
Before
MAY 2 8 2010


Data File: \\Cashl\Acqudata\GC09\data\052710A.B\0527F028.D
Injection Date: 28-MAY-2010 02:30
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data, DATA.GC: 2.047 to 15.817 Min
Min
2.135
2.250
2.410
2.590
2.807
2.913
3.077
3.137
3.237
Aroclor 1221
4.243
4.357
Aroclor 1221
Aroclor 1221
Aroclor 1221
5.120
5.760
5.950
6.267
6.543
6.720
7.030
7.120
7.540
7.833
8.133
8.210
8.297
8.430
Aroclor 1254
8.667
8.860
8.987
9.107
Aroclor 1254
9.397
Aroclor 1254
9.583
9.770
Aroclor 1254
Aroclor 1254
10.457
10.557
10.633
10.747
10.873
11.077
11.377
11.623
11.727
12.077
12.137
13.243
14.447
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other
After
Initials
MAY 28 2010



Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R028.D
Injection Date: 28-MAY-2010 02:30
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.513 Min
Y (x10^4)
Min
1.730
1.953
3.050
3.133
3.373
3.930
Aroclor 1221
4.553
4.687
Aroclor 1221
4.950
Aroclor 1221
Aroclor 1221
5.557
6.170
6.357
6.537
6.903
7.450
7.623
7.720
8.013
8.347
8.813
8.963
9.127
Aroclor 1254
9.367
9.553
9.663
Aroclor 1254
9.930
Aroclor 1254
Aroclor 1254
Aroclor 1254
10.893
11.010
11.217
11.423
11.543
12.277
12.563
12.977
13.137
14.093
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F029.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F029.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R029.D
Inj Date : 28-MAY-2010 02:56
Sample Info: 1221/1254 @ 100/50ppb | PCB5-52E
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1221+1254.sub
Sub List #2 : 1221+1254.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.687	4.690	9654	2121	136	55.2	80.00- 120.00	100.00(M)
	4.547	4.850	4981	4103	123	116	51.26- 76.90	51.59(M)
	4.797	5.283	13025	8959	124	112	133.68- 200.52	134.92(M)
	4.987	5.410	7707	6218	116	111	85.57- 128.36	79.84(M)
			Average of Peak Amounts =		125	98.6		
Aroclor 1254	8.543	9.250	24280	22371	60.2	58.2	80.00- 120.00	100.00(M)
	9.273	9.833	15302	8235	54.6	46.8	61.91- 92.86	63.02(M)
	9.477	10.027	33727	17080	58.9	47.8	116.61- 174.92	138.91(M)
	9.887	10.363	24834	7183	53.5	44.7	101.15- 151.73	102.28(M)
	10.097	10.640	13034	15498	52.7	49.3	53.30- 79.94	53.68(M)
			Average of Peak Amounts =		56.0	49.4		

QC Flag Legend

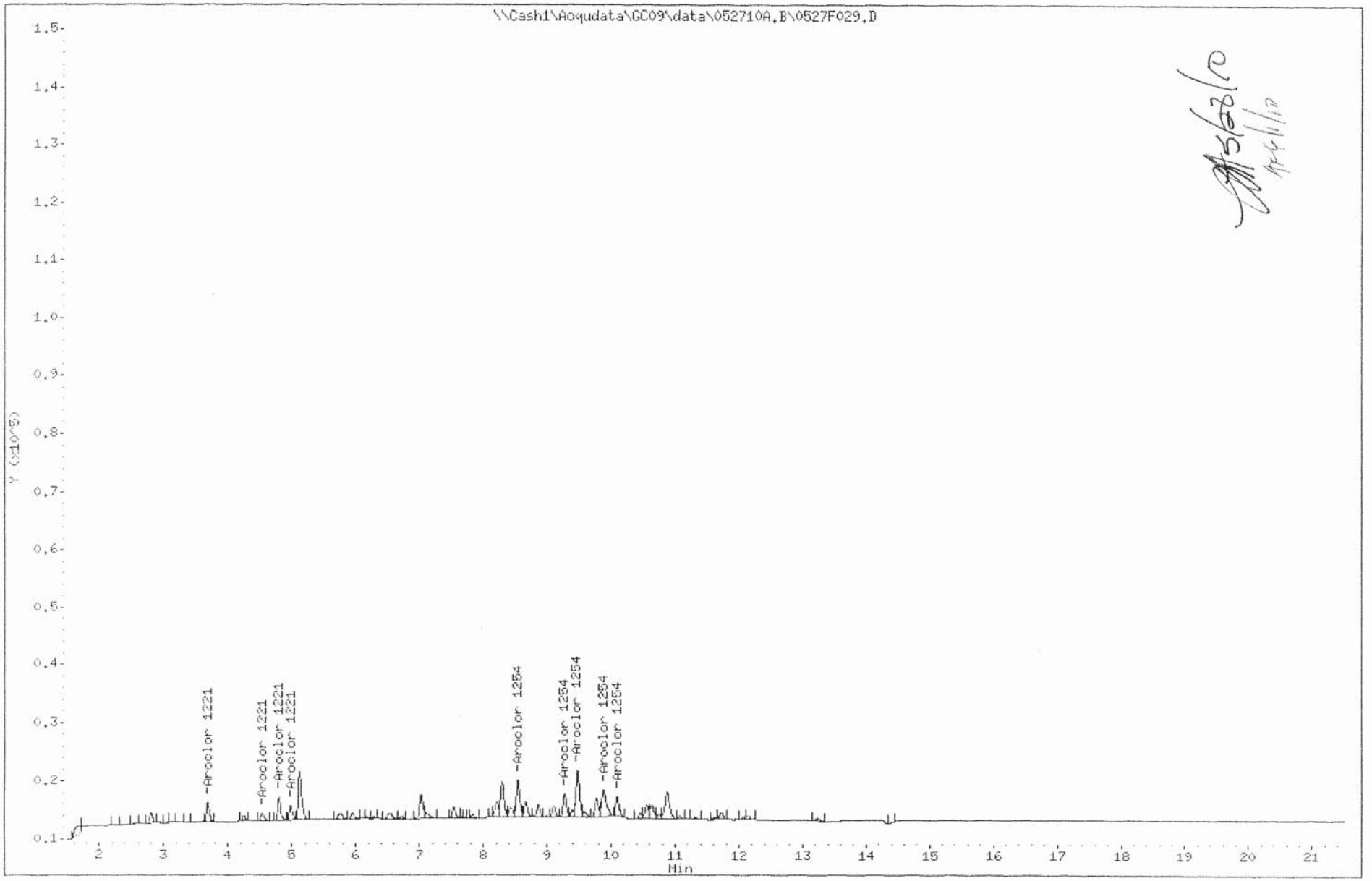
M - Compound response manually integrated.

5/28/10
6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F029,D
Date : 28-MAY-2010 02:56
Client ID:
Sample Info: 1221/1254 @ 100/50ppb | PCB5-52E

Instrument: GC09.i

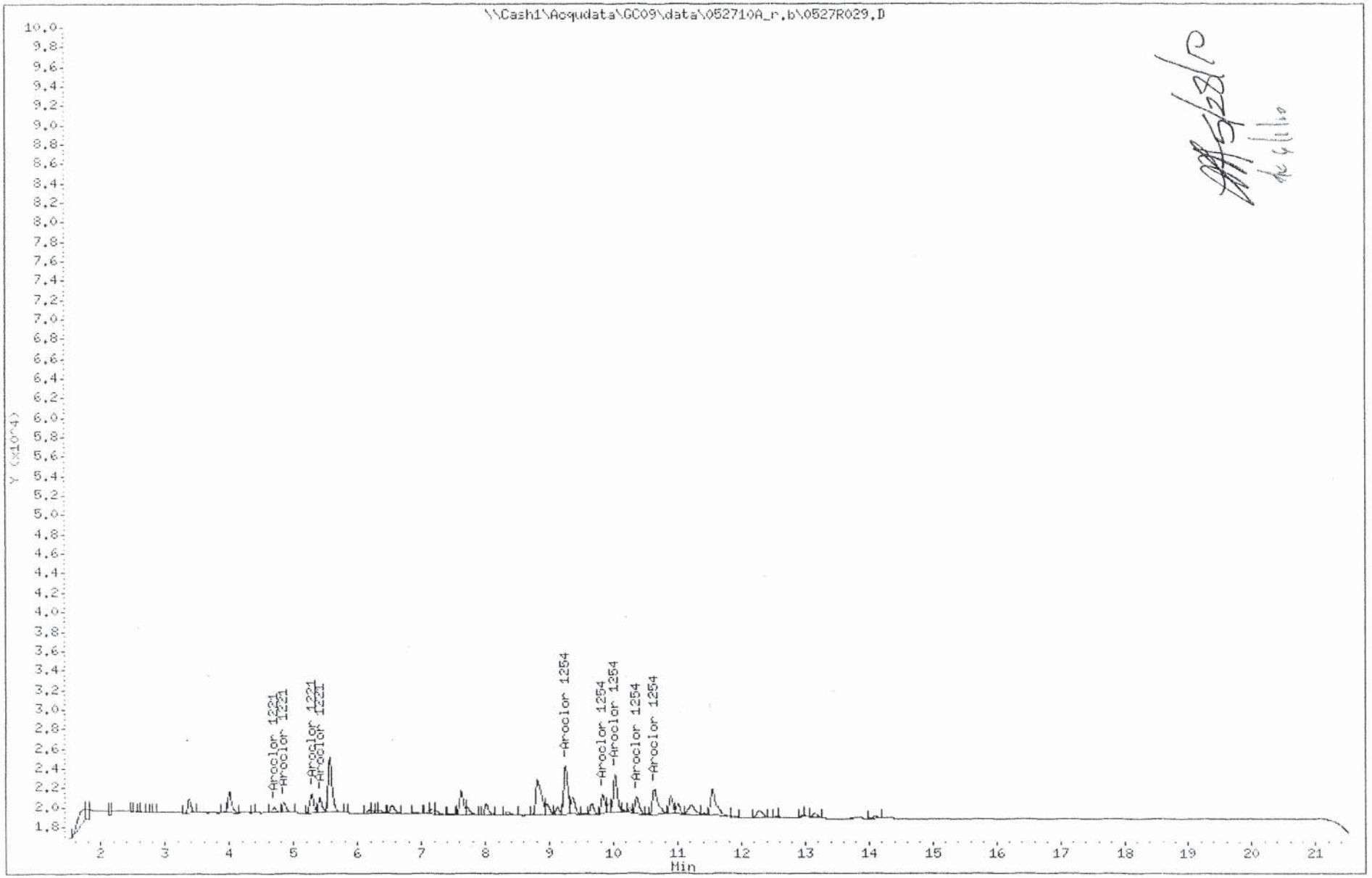
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

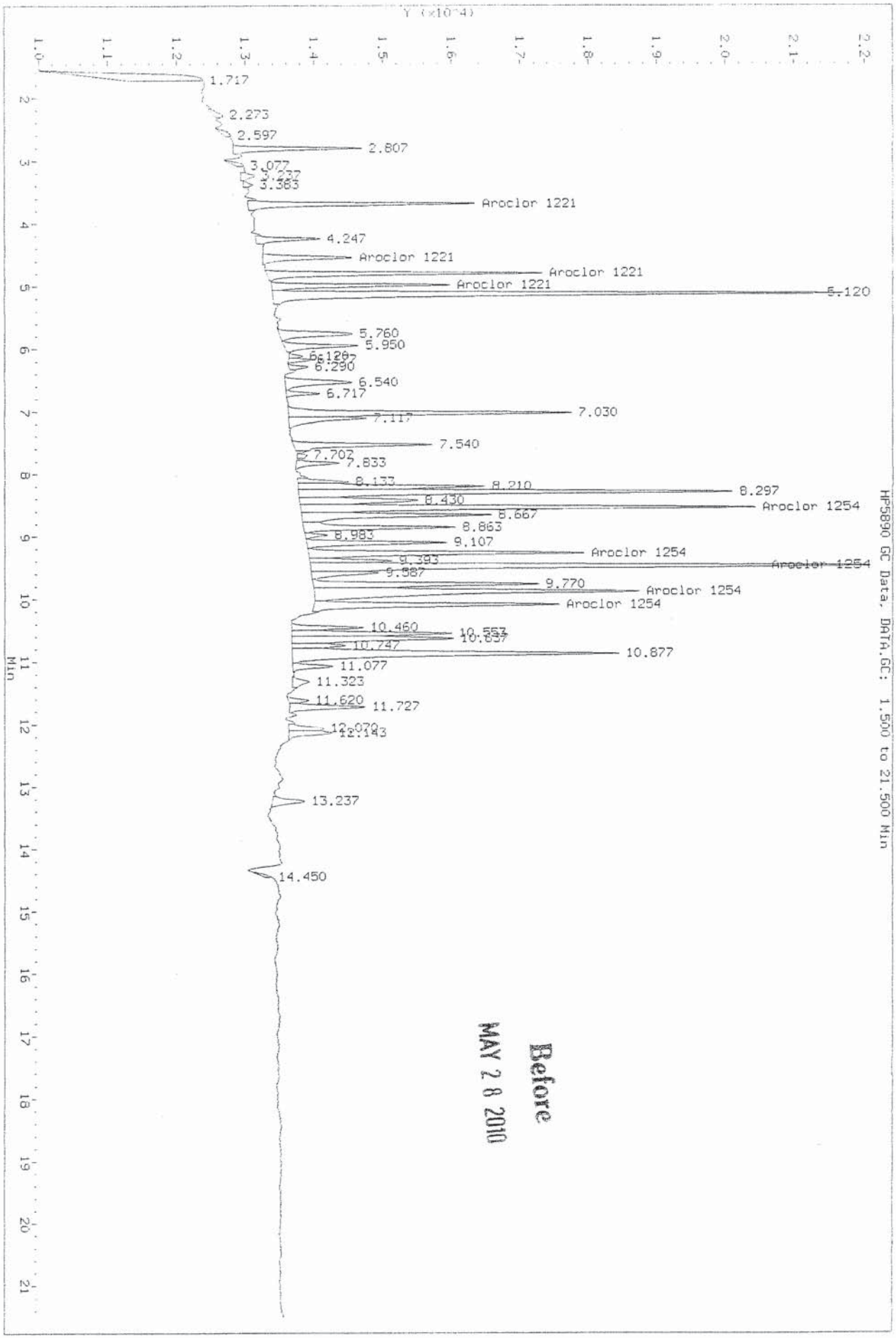


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R029.D
Date : 28-MAY-2010 02:56
Client ID:
Sample Info: 1221/1254 @ 100/50ppb | PCB5-52E

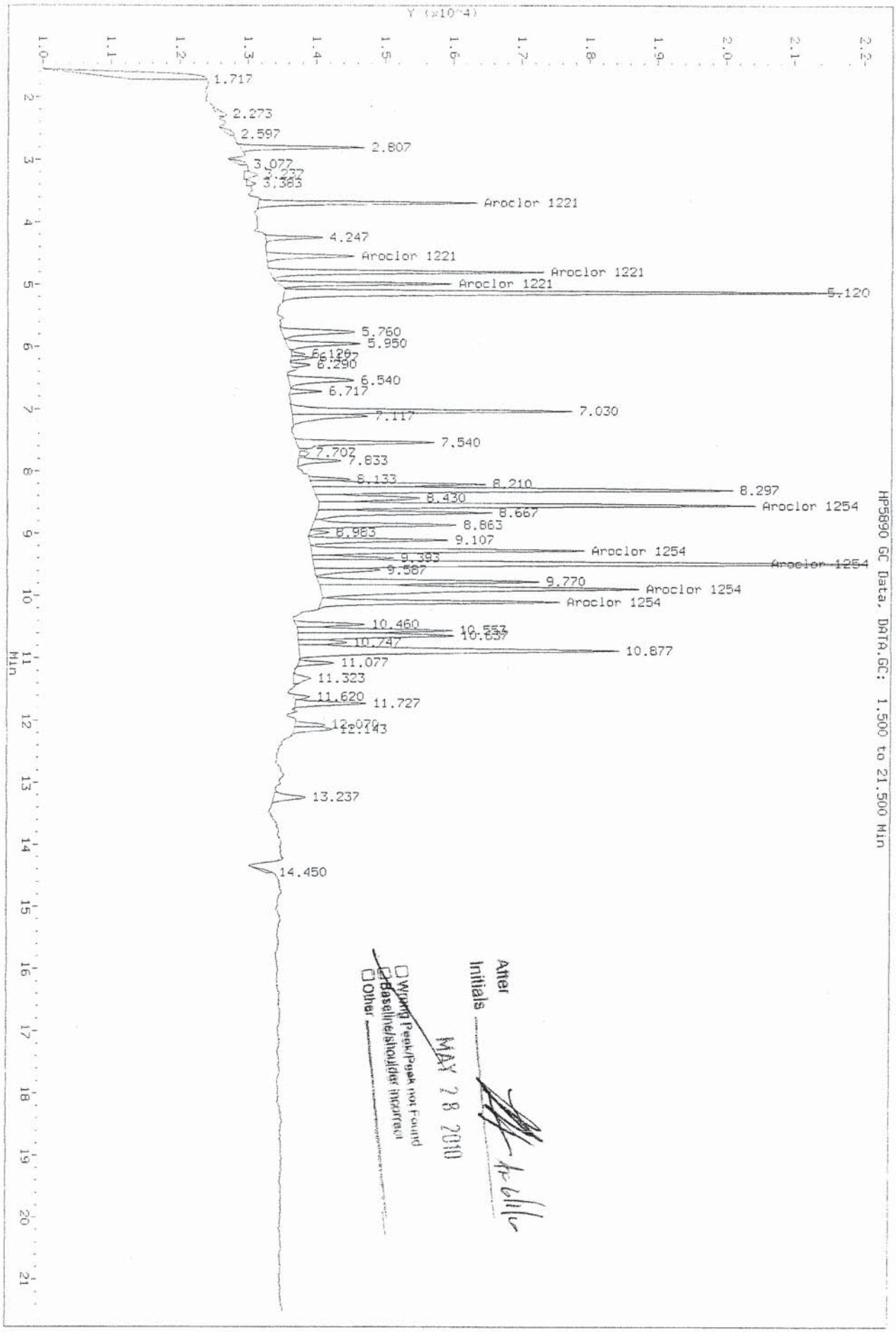
Instrument: GC09.i
Operator: LHarris
Column diameter: 0,53

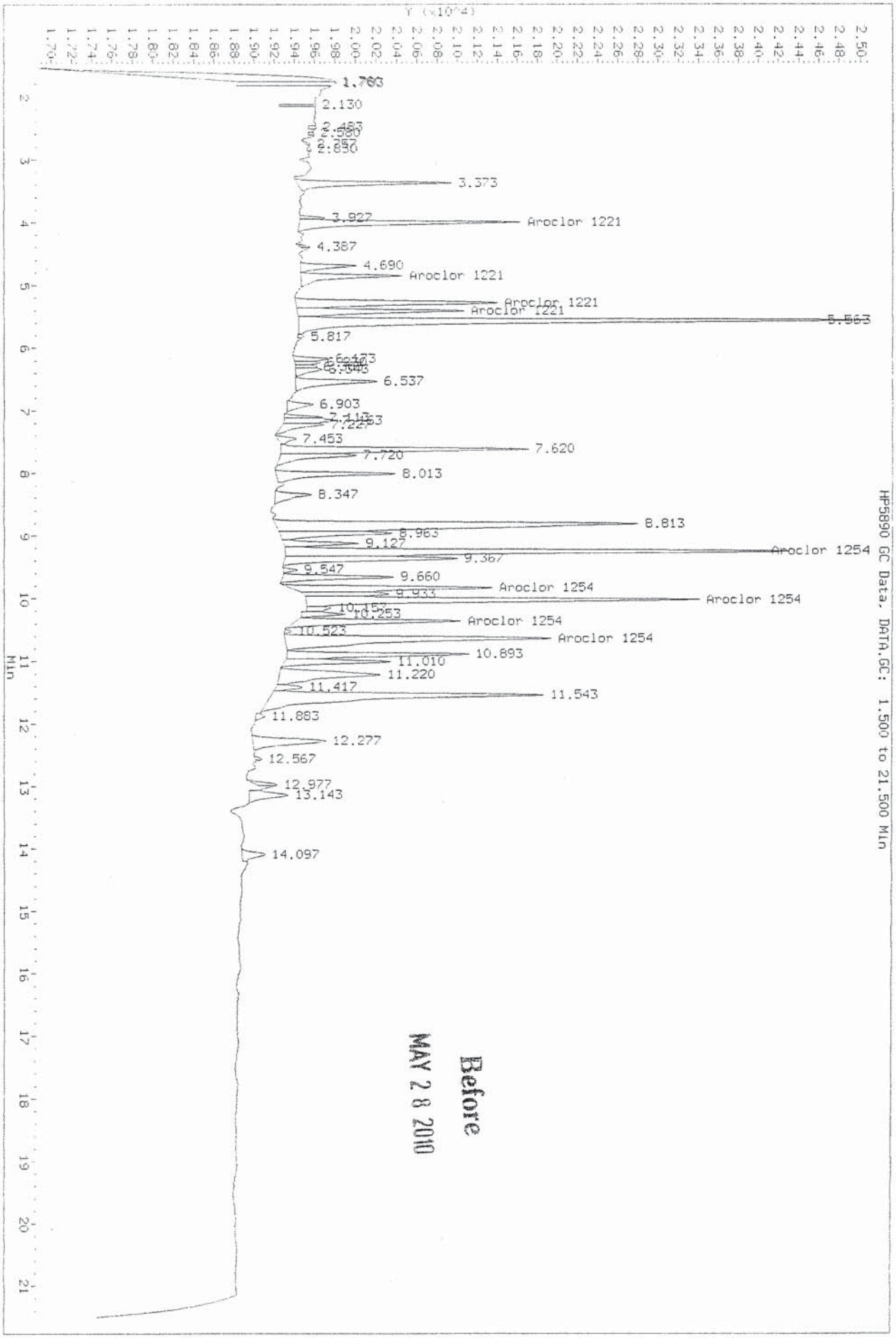
Column phase: DB-XLB





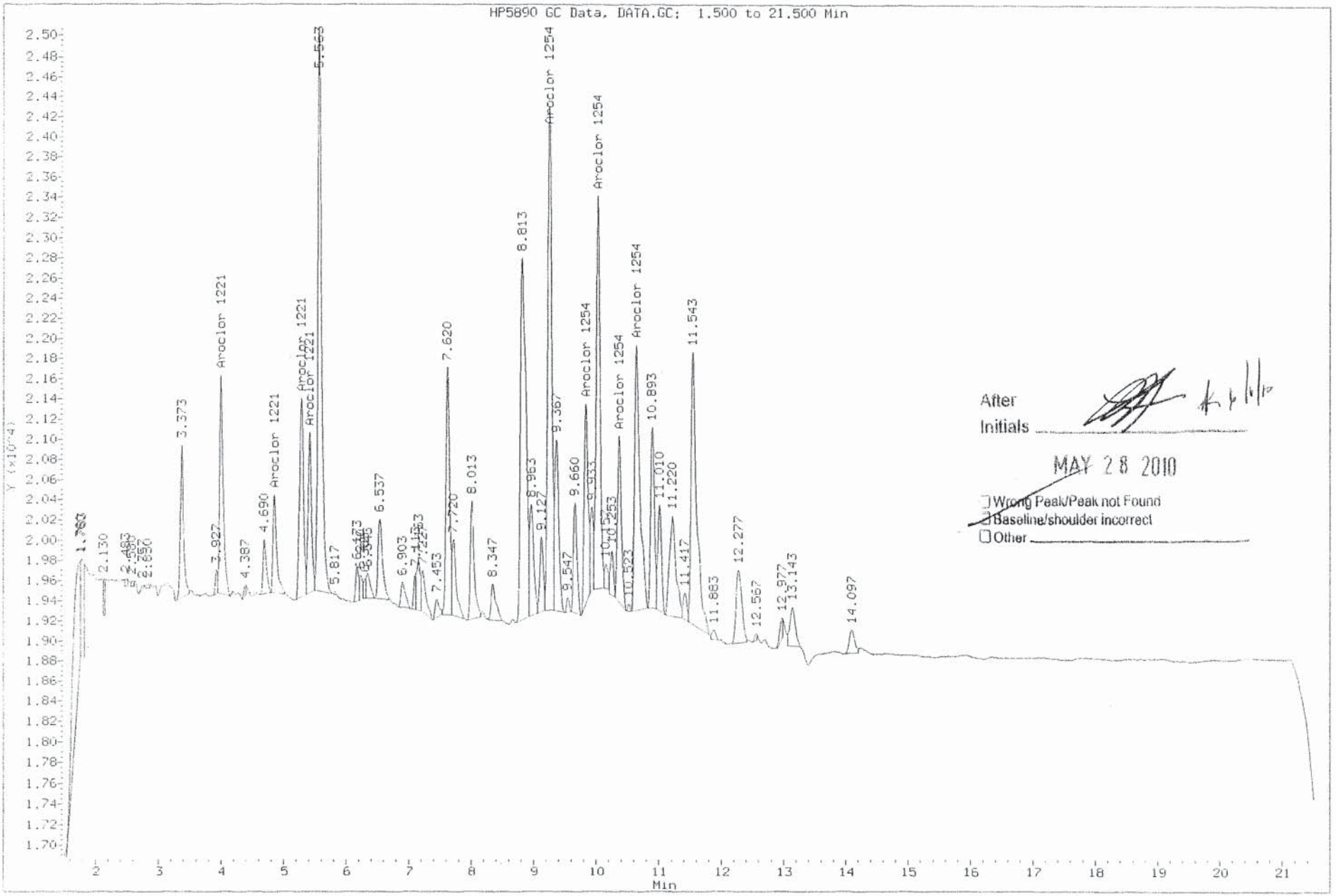
Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F029.D
Injection Date: 28-MAY-2010 02:56
Instrument: GC09.i
Client Sample ID:
Y (x10^4)
1.717
2.273
2.597
2.807
3.077
3.237
3.383
Aroclor 1221
4.247
Aroclor 1221
Aroclor 1221
Aroclor 1221
5.120
5.760
5.950
6.129
6.290
6.540
6.717
7.030
7.117
7.540
7.702
7.833
8.133
8.210
8.297
8.430
Aroclor 1254
8.667
8.863
8.983
9.107
Aroclor 1254
9.393
Aroclor 1254
9.587
9.770
Aroclor 1254
Aroclor 1254
10.460
10.553
10.637
10.747
10.877
11.077
11.323
11.620
11.727
12.070
12.143
13.237
14.450
HP5890 GC Data, DATA.GC; 1.500 to 21.500 Min
Min
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other



Before
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R029.D
Injection Date: 28-MAY-2010 02:56
Instrument: GC09.i
Client Sample ID:



After
Initials

MAY 28 2010

☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder incorrect
☐ Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F030.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F030.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R030.D
Inj Date : 28-MAY-2010 03:22
Sample Info: 1221/1254 @ 1000/500ppb | PCB5-52F
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1221+1254.sub
Sub List #2 : 1221+1254.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.687	4.687	71928	22659	1010	590	80.00- 120.00	100.00(M)
	4.543	4.850	43264	40044	1070	1140	51.26- 76.90	60.15(M)
	4.797	5.283	108157	86110	1030	1080	133.68- 200.52	150.37(M)
	4.983	5.410	71992	61777	1080	1100	85.57- 128.36	100.09(M)
			Average of Peak Amounts =		1050	978		
Aroclor 1254	8.540	9.250	219677	202310	545	526	80.00- 120.00	100.00(M)
	9.270	9.833	149800	96832	536	550	61.91- 92.86	68.19(M)
	9.473	10.023	304398	197264	532	552	116.61- 174.92	138.57(M)
	9.880	10.363	252087	93897	545	584	101.15- 151.73	114.75(M)
	10.093	10.637	136255	169734	554	540	53.30- 79.94	62.03(M)
			Average of Peak Amounts =		542	550		

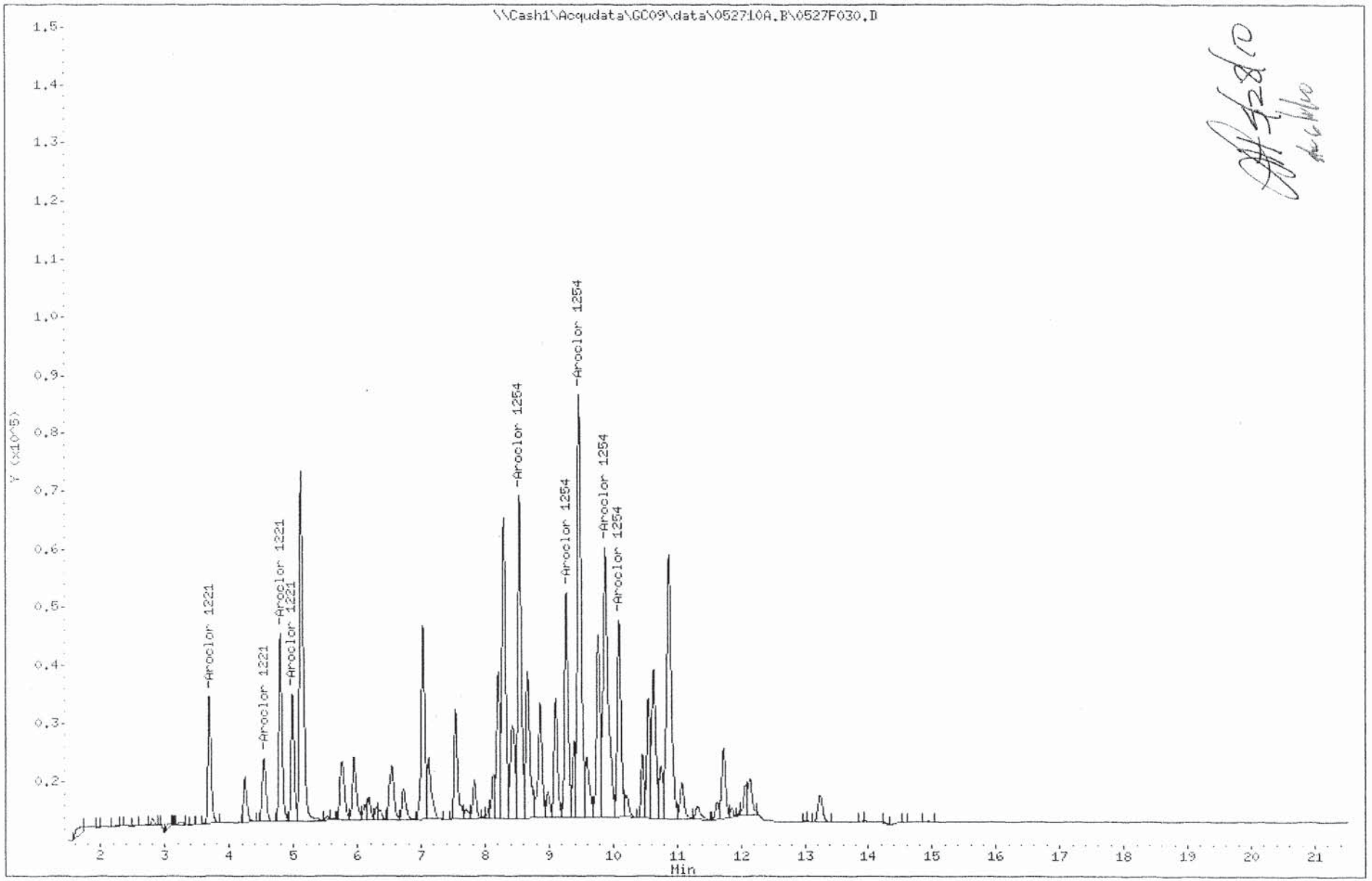
QC Flag Legend

M - Compound response manually integrated.

5/28/10

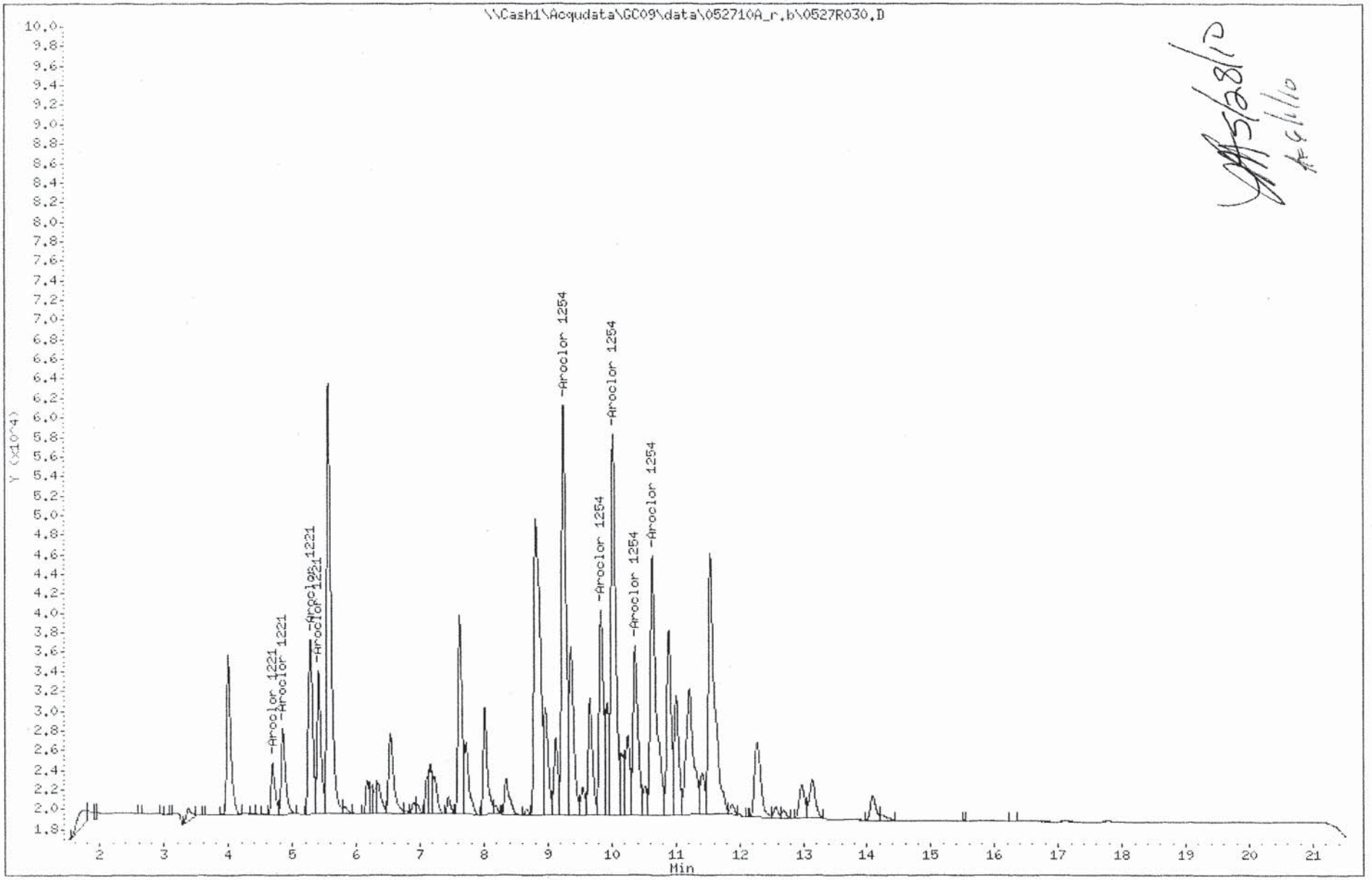
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F030.D
Date : 28-MAY-2010 03:22
Client ID:
Sample Info: 1221/1254 @ 1000/500ppb | PCB5-52F

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

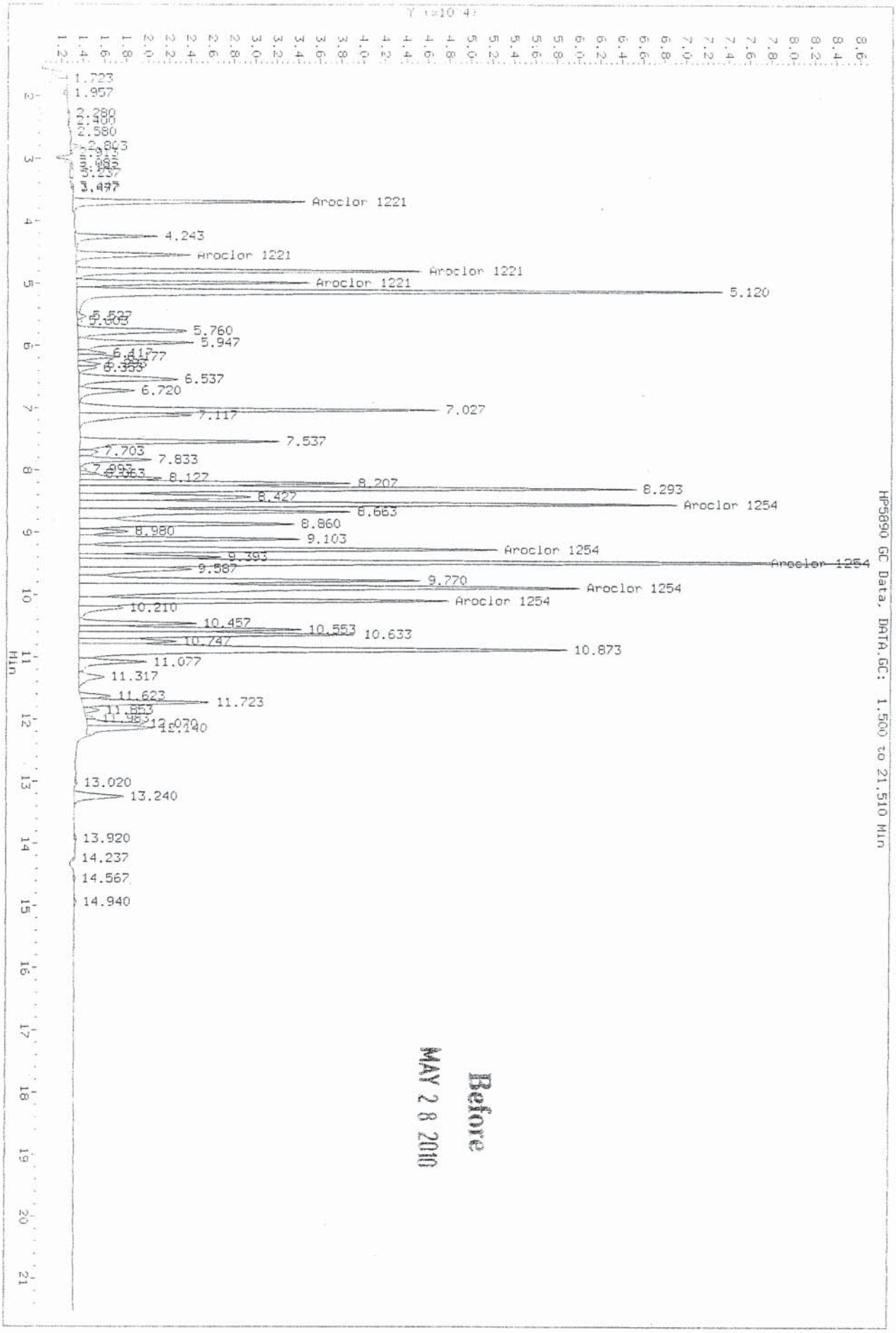


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R030.D
Date : 28-MAY-2010 03:22
Client ID:
Sample Info: 1221/1254 @ 1000/500ppb | PCB5-52F

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

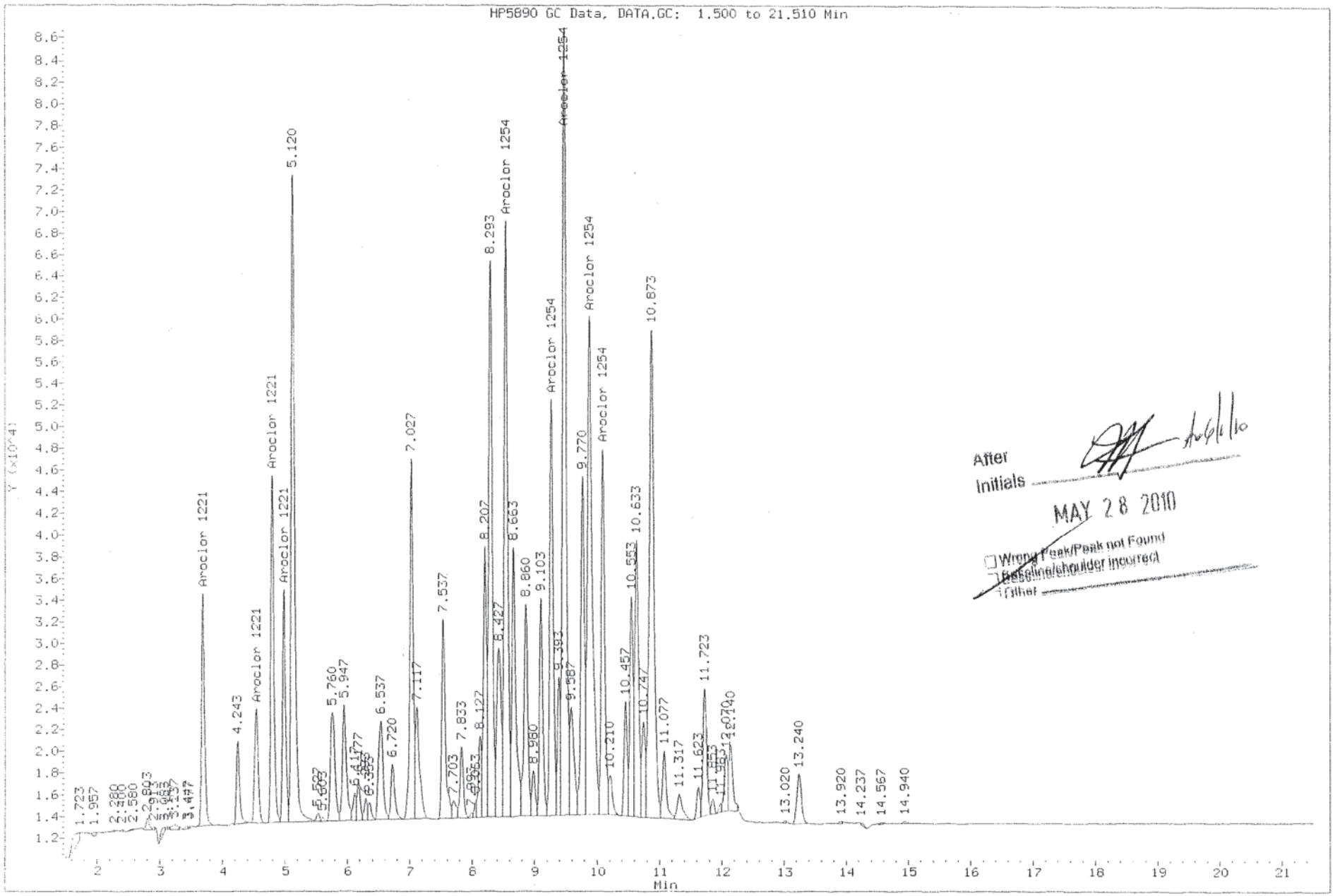


765



Before
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F030.D
Injection Date: 28-MAY-2010 03:22
Instrument: GC09.i
Client Sample ID:



After
Initials
MAY 28 2010
☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder incorrect
☐ Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F031.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F031.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R031.D
Inj Date : 28-MAY-2010 03:48
Sample Info: 1221/1254 @ 2000/1000ppb | PCB5-52G
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1221+1254.sub
Sub List #2 : 1221+1254.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.687	4.687	122122	41167	1720	1310	80.00- 120.00	100.00(M)
	4.543	4.850	77462	72656	1930	2060	51.26- 76.90	63.43(M)
	4.797	5.283	193426	159066	1840	1990	133.68- 200.52	158.39(M)
	4.983	5.410	127548	113283	1910	2020	85.57- 128.36	104.44(M)
			Average of Peak Amounts =		1850	1840		
Aroclor 1254	8.540	9.247	390024	362273	965	942	80.00- 120.00	100.00(M)
	9.270	9.830	278332	180264	992	1020	61.91- 92.86	71.36(M)
	9.473	10.023	555094	365880	967	1020	116.61- 174.92	142.32(M)
	9.880	10.360	462920	171964	996	1070	101.15- 151.73	118.69(M)
	10.093	10.633	250662	312572	1010	994	53.30- 79.94	64.27(M)
			Average of Peak Amounts =		986	1010		

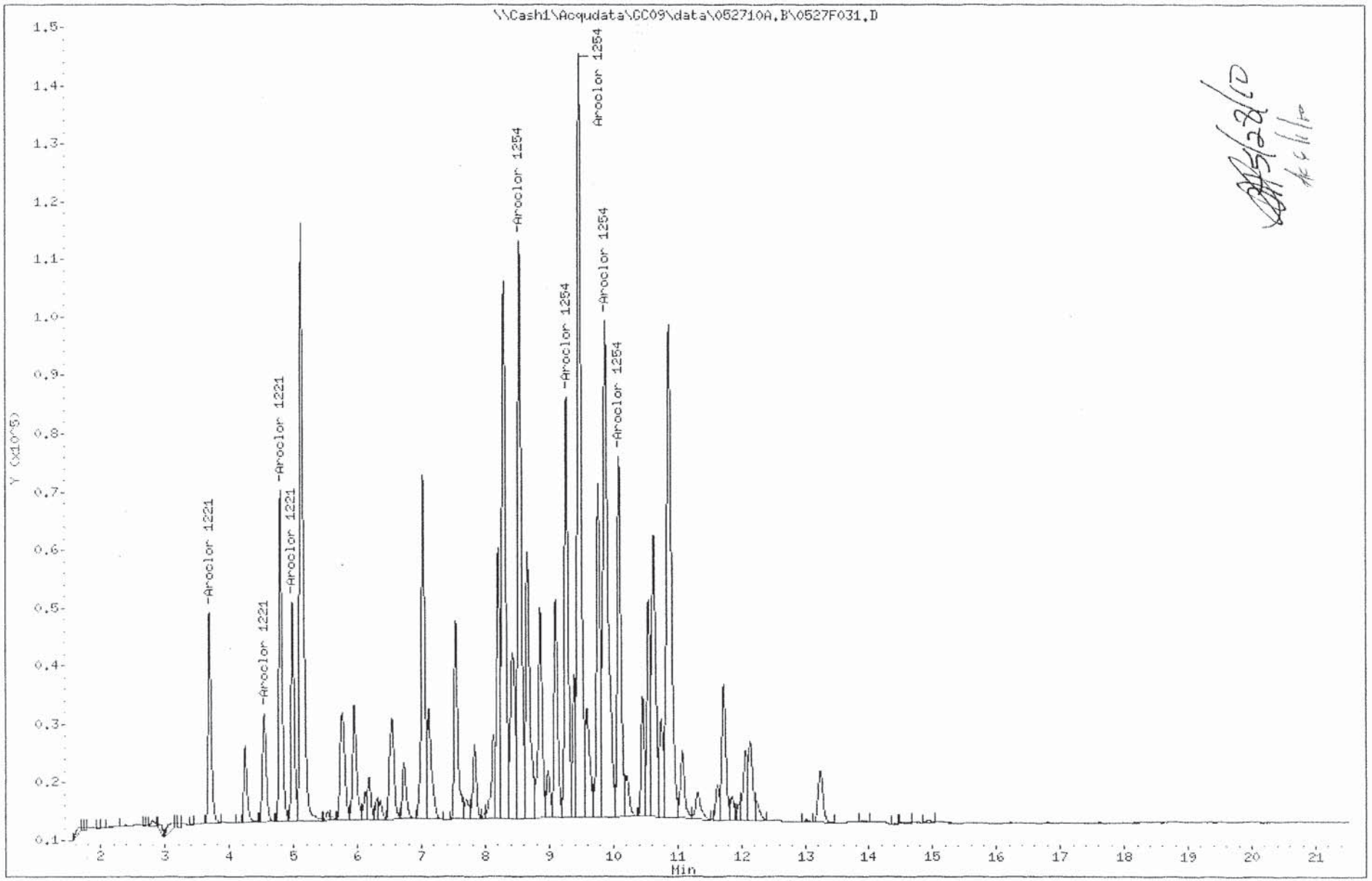
QC Flag Legend

M - Compound response manually integrated.

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F031.D
Date : 28-MAY-2010 03:48
Client ID:
Sample Info: 1221/1254 @ 2000/1000ppb | PCB5-52G

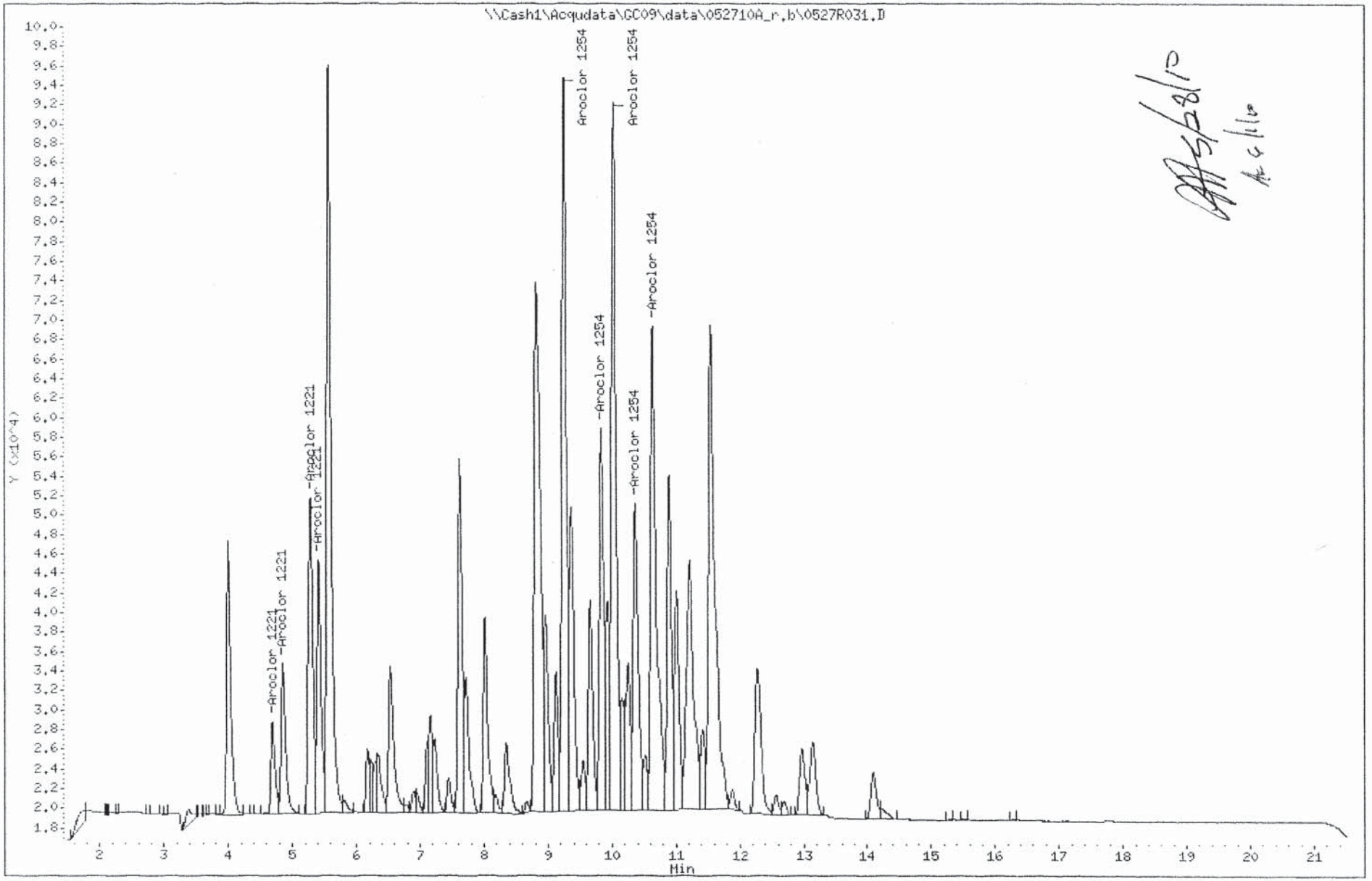
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

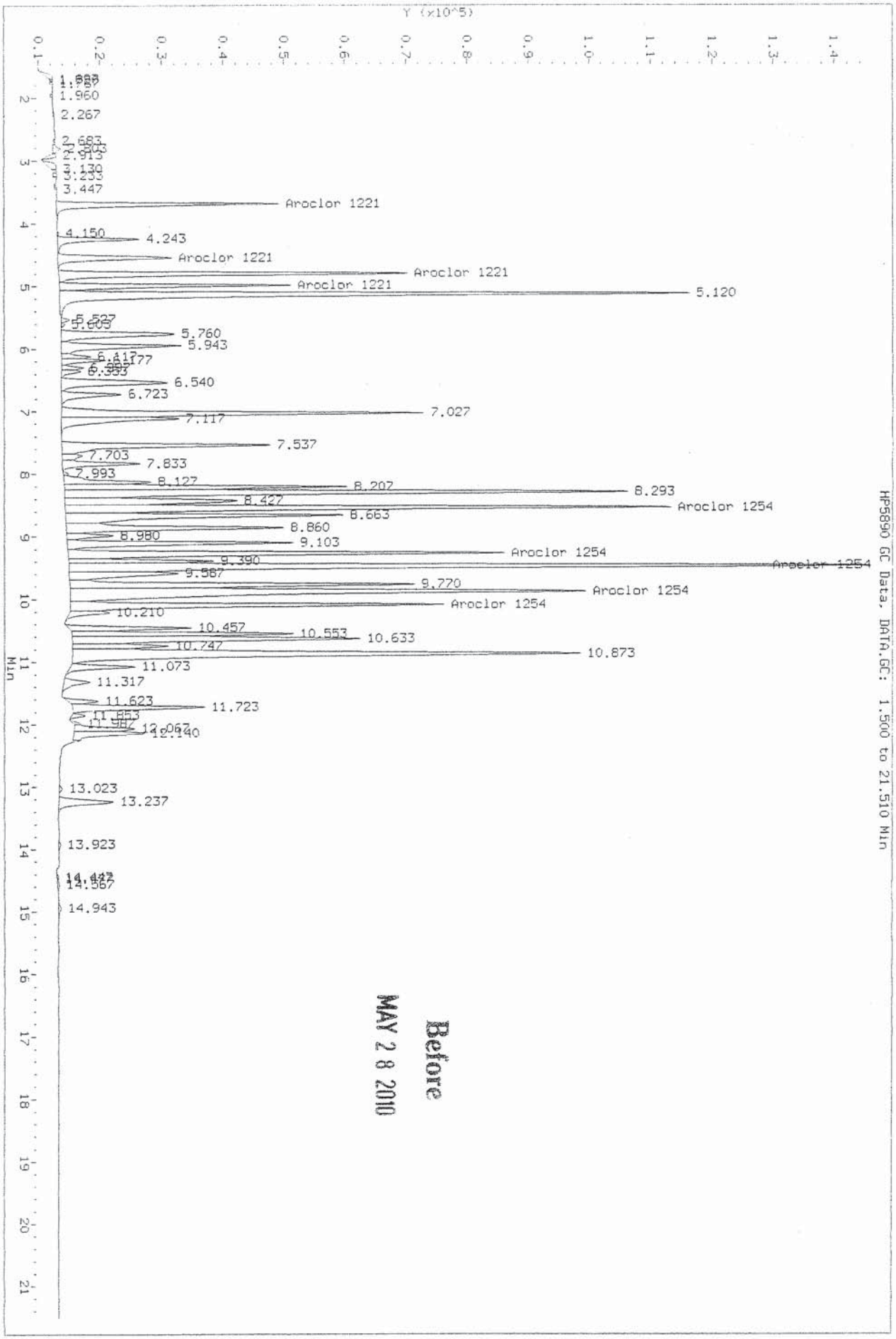


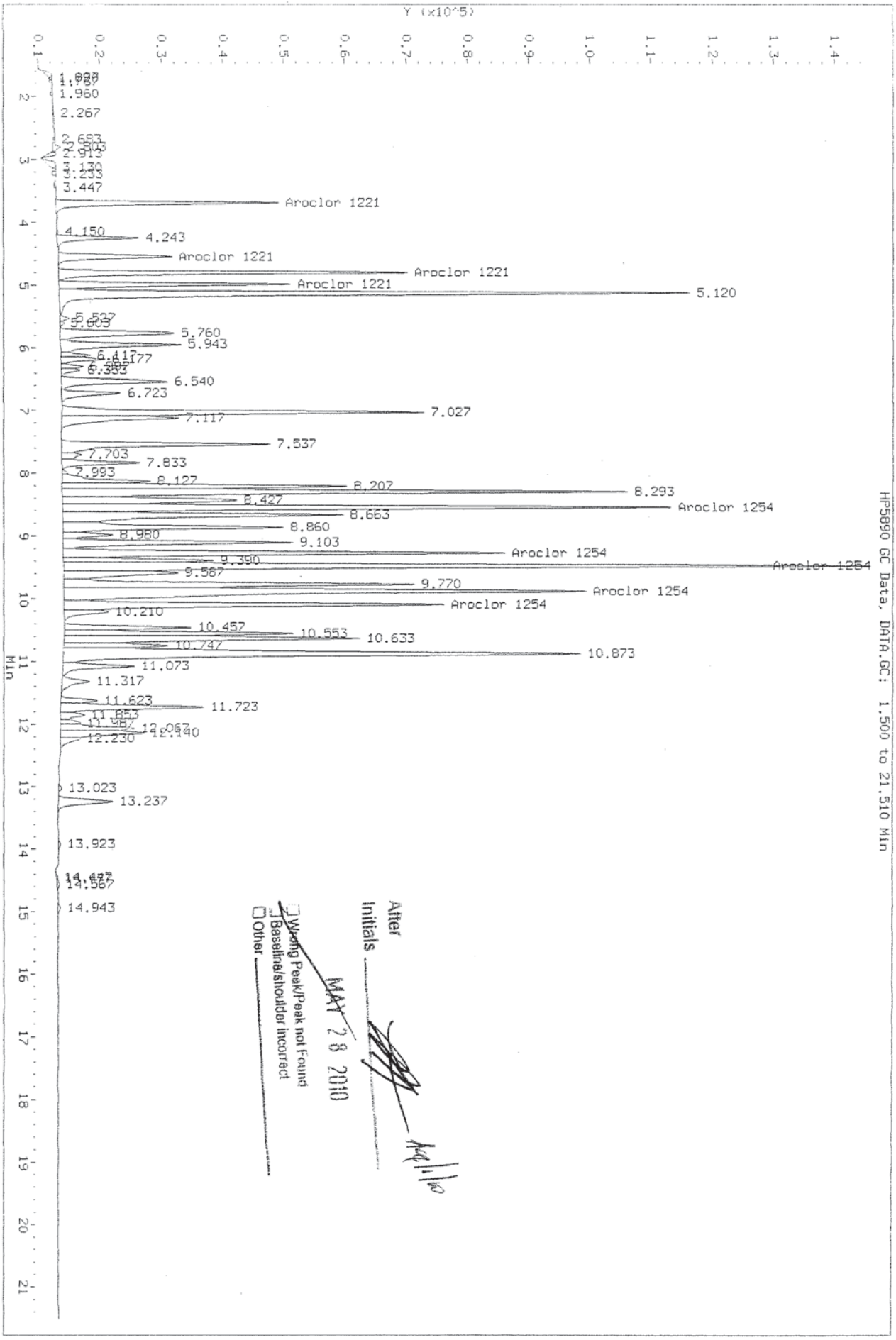
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R031.D
Date : 28-MAY-2010 03:48
Client ID:
Sample Info: 1221/1254 @ 2000/1000ppb | PCB5-52G

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F031.D
Injection Date: 28-MAY-2010 03:48
Instrument: GC09.1
Client Sample ID:
Y (x10^5)
Min
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Aroclor 1221
Aroclor 1254
Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F031.D
Injection Date: 28-MAY-2010 03:48
Instrument: GC09.i
Client Sample ID:
Y (x10^5)
HP5890 GC Data, DATA.GC; 1.500 to 21.510 Min
Min
Aroclor 1221
Aroclor 1254
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other
After
Initials
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F032.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F032.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R032.D
Inj Date : 28-MAY-2010 04:14
Sample Info: 1221/1254 @ 4000/2000ppb | PCB5-52H
Misc Info :
Cal Date : 28-MAY-2010 13:16
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1221+1254.sub
Sub List #2 : 1221+1254.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.683	4.687	211846	74862	2990	2950	80.00- 120.00	100.00
	4.540	4.850	135747	130962	3380	3710	51.26- 76.90	64.08
	4.793	5.283	345748	295862	3290	3690	133.68- 200.52	163.21
	4.983	5.410	225277	208386	3380	3720	85.57- 128.36	106.34
			Average of Peak Amounts =		3260	3520		
Aroclor 1254	8.540	9.247	701844	668933	1740	1740	80.00- 120.00	100.00
	9.270	9.830	523384	357253	1860	2030	61.91- 92.86	74.57
	9.470	10.023	1017858	707364	1770	1980	116.61- 174.92	145.03
	9.880	10.360	866454	334601	1860	2080	101.15- 151.73	123.45
	10.093	10.630	462993	604402	1870	1920	53.30- 79.94	65.97
			Average of Peak Amounts =		1820	1950		

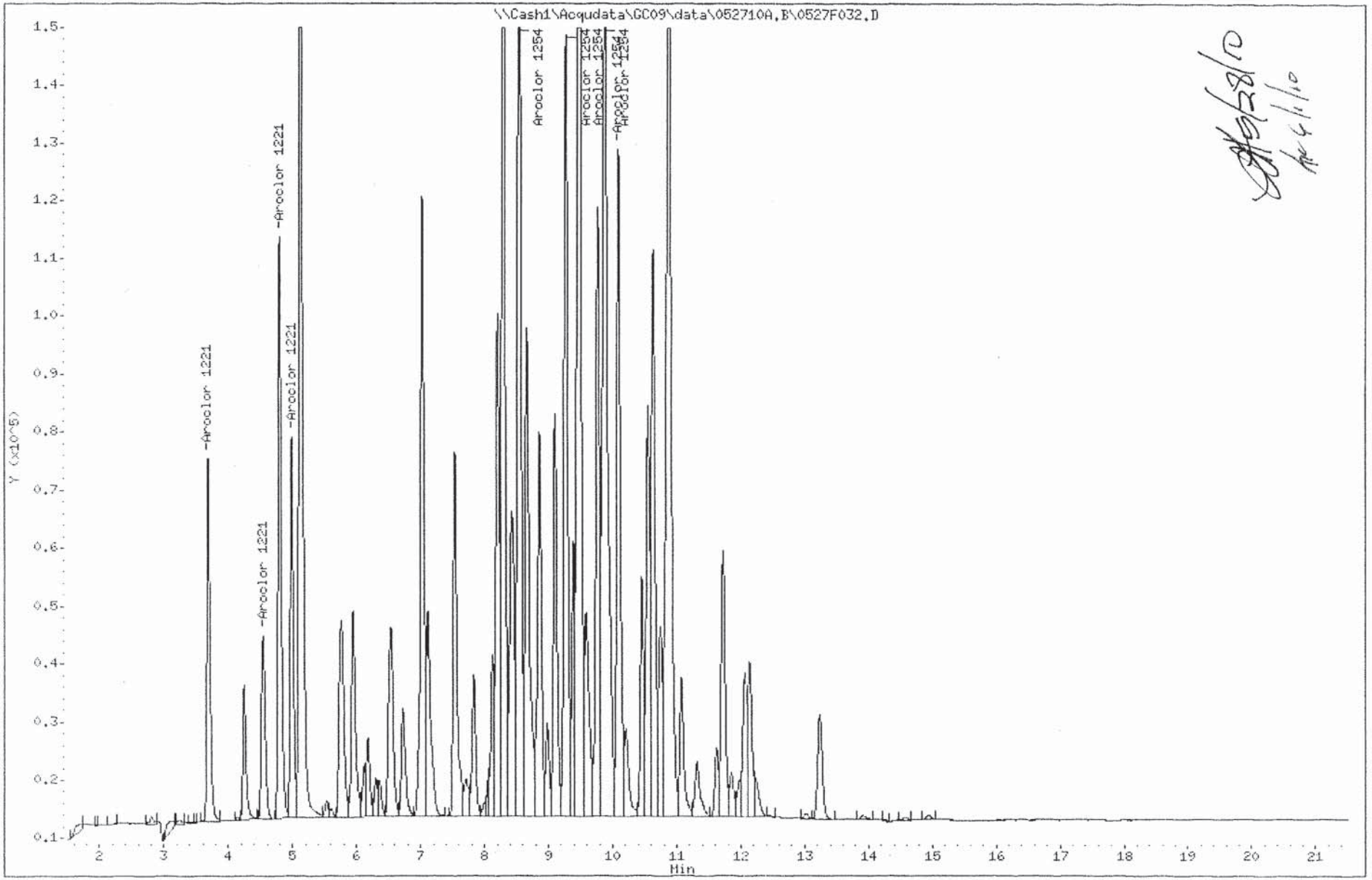
5/28/10

6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F032.D
Date : 28-MAY-2010 04:14
Client ID:
Sample Info: 1221/1254 @ 4000/2000ppb | PCB5-52H

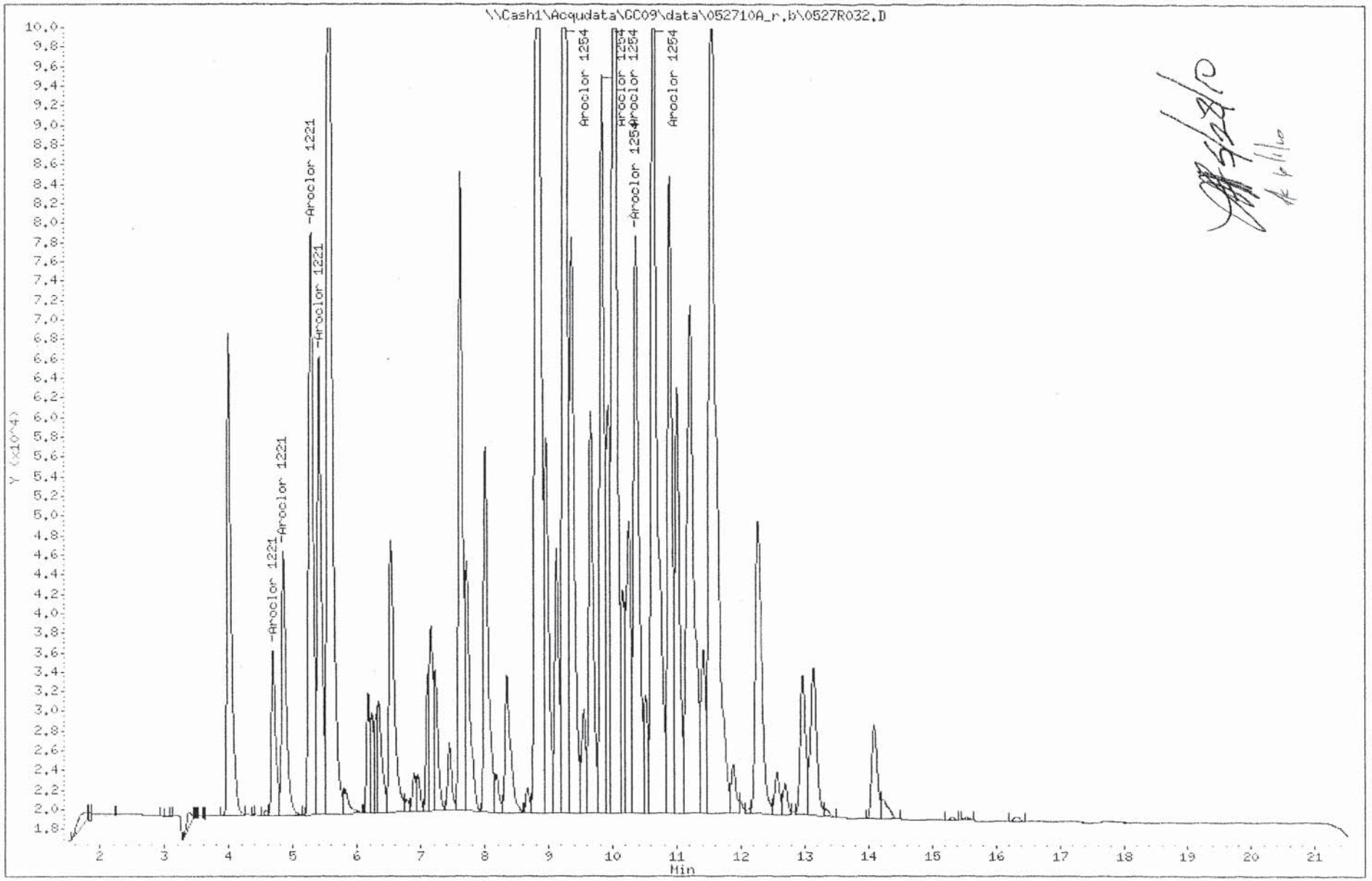
Instrument: GC09.i

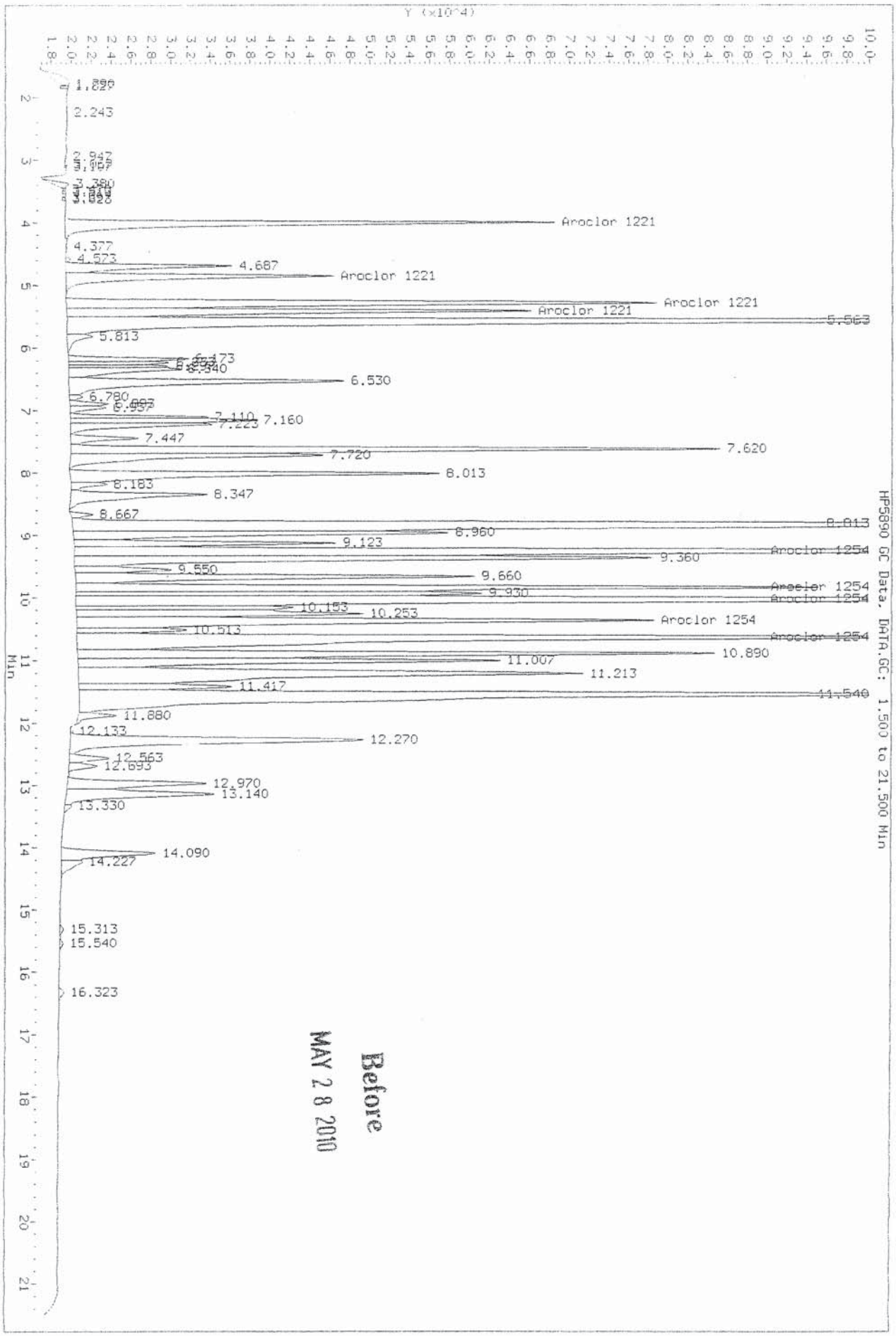
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R032.D
Date : 28-MAY-2010 04:14
Client ID:
Sample Info: 1221/1254 @ 4000/2000ppb | PCB5-52H

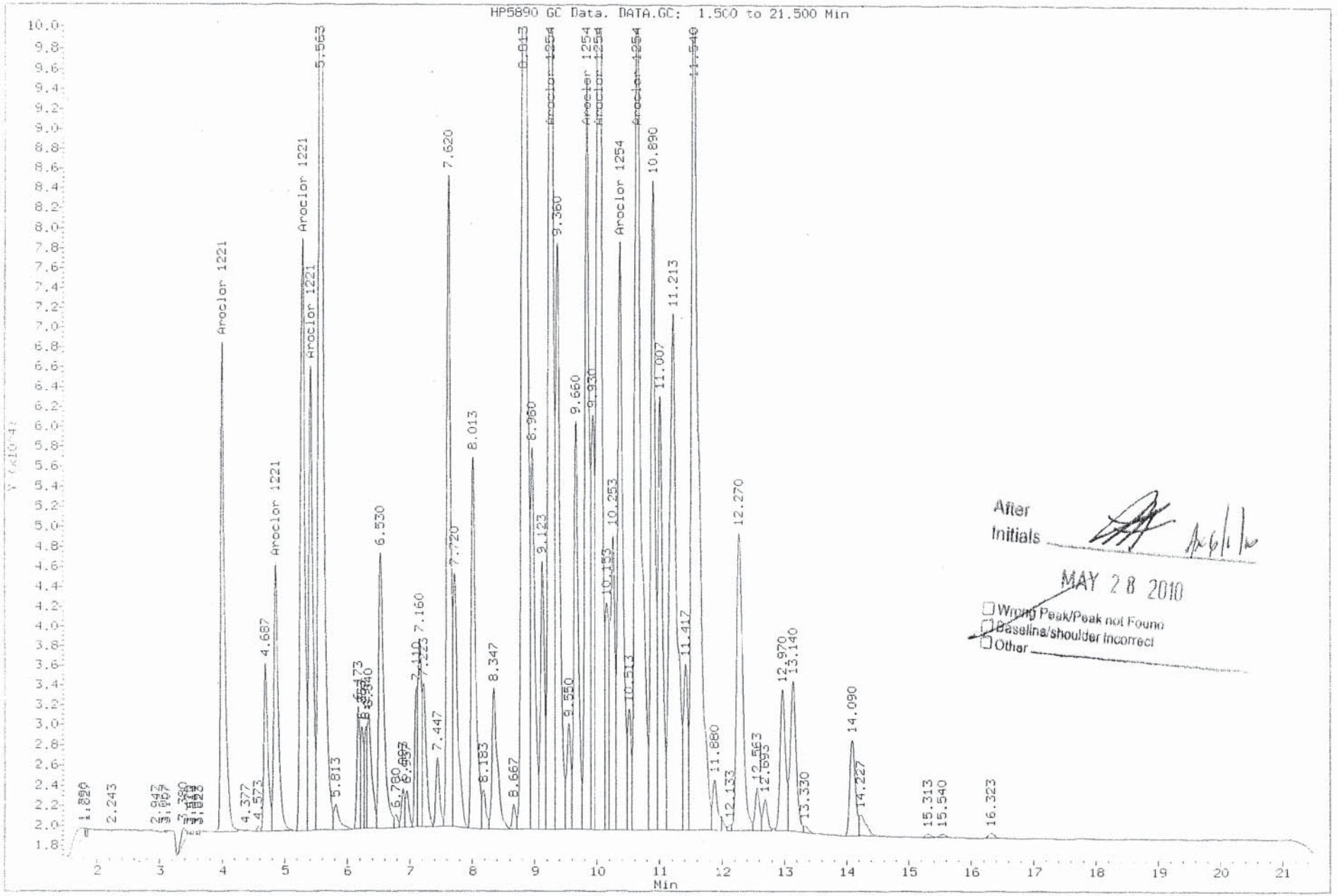
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53





Before
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R032.D
Injection Date: 28-MAY-2010 04:14
Instrument: GC09.i
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F033.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F033.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R033.D
Inj Date : 28-MAY-2010 04:40
Sample Info: 1221/1254 @ 10000/5000ppb | PCB5-52I
Misc Info :
Cal Date : 28-MAY-2010 13:16
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1221+1254.sub
Sub List #2 : 1221+1254.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.687	4.687	432503	159348	6100	7860	80.00- 120.00	100.00(A)
	4.540	4.850	277157	271312	6910	7690	51.26- 76.90	64.08(A)
	4.797	5.283	722722	651571	6880	8140	133.68- 200.52	167.10(A)
	4.983	5.410	462618	445929	6940	7960	85.57- 128.36	106.96(A)
			Average of Peak Amounts =		6710	7910		
Aroclor 1254	8.540	9.247	1508202	1471676	3730	3830	80.00- 120.00	100.00
	9.270	9.830	1167135	819224	4160	4650	61.91- 92.86	77.39
	9.470	10.020	2198476	1601678	3830	4480	116.61- 174.92	145.77
	9.877	10.360	1907007	742950	4100	4620	101.15- 151.73	126.44
	10.090	10.630	1004763	1364075	4060	4340	53.30- 79.94	66.62
			Average of Peak Amounts =		3980	4380		

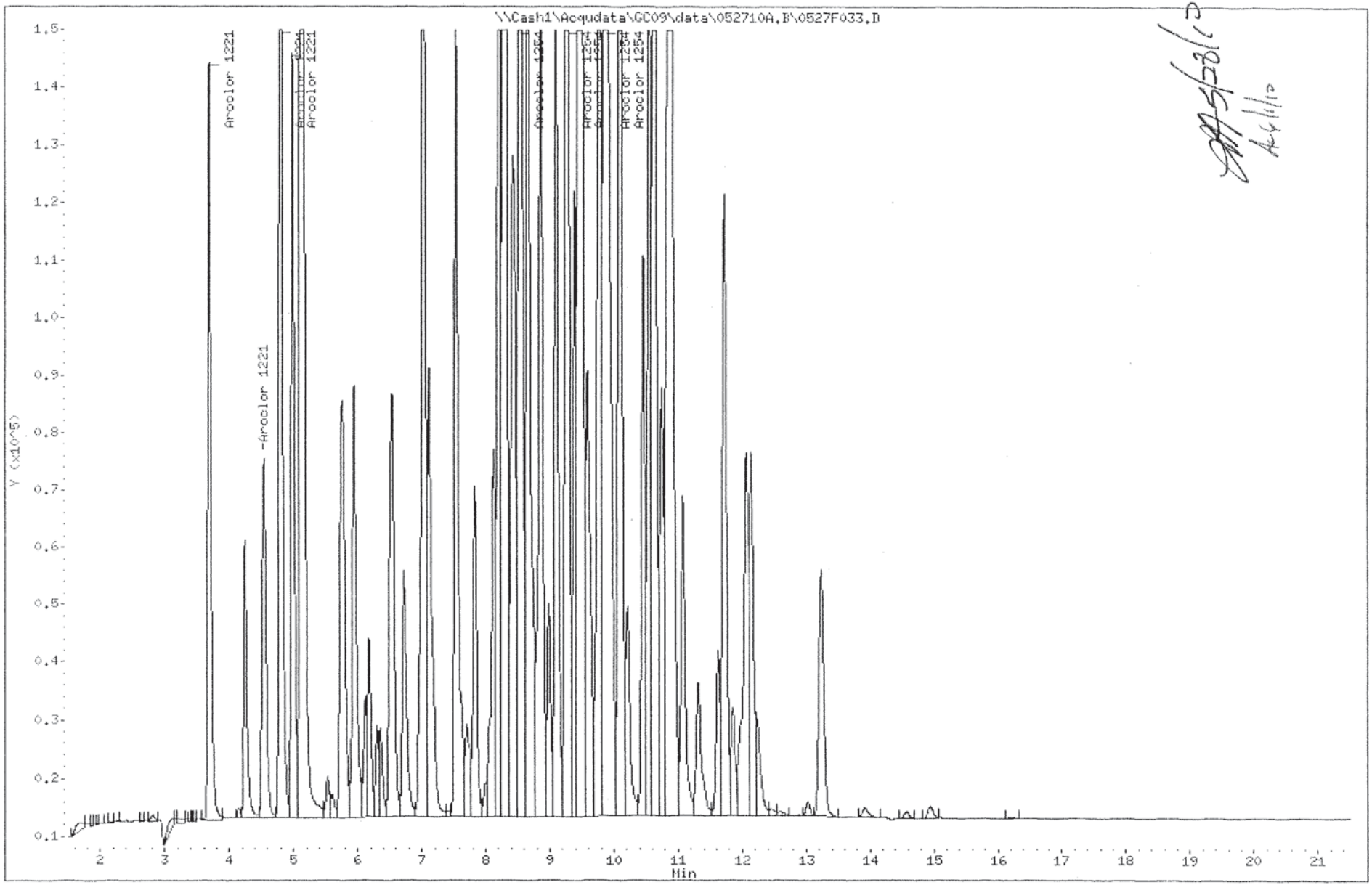
QC Flag Legend

A - Target compound detected but, quantitated amount exceeded maximum amount.

5/28/10
5/1/10

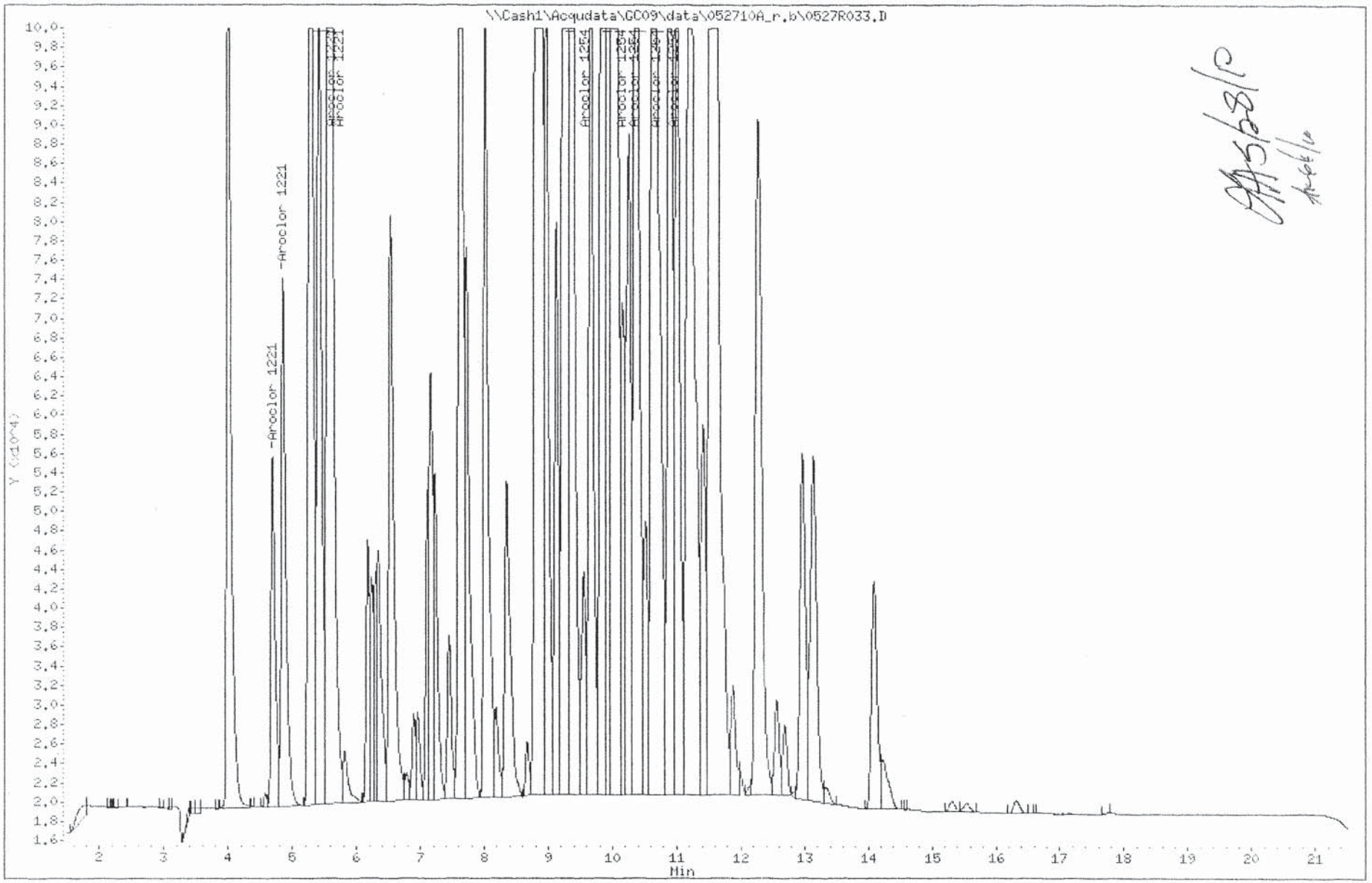
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F033.D
Date : 28-MAY-2010 04:40
Client ID:
Sample Info: 1221/1254 @ 10000/5000ppb | PCB5-52I

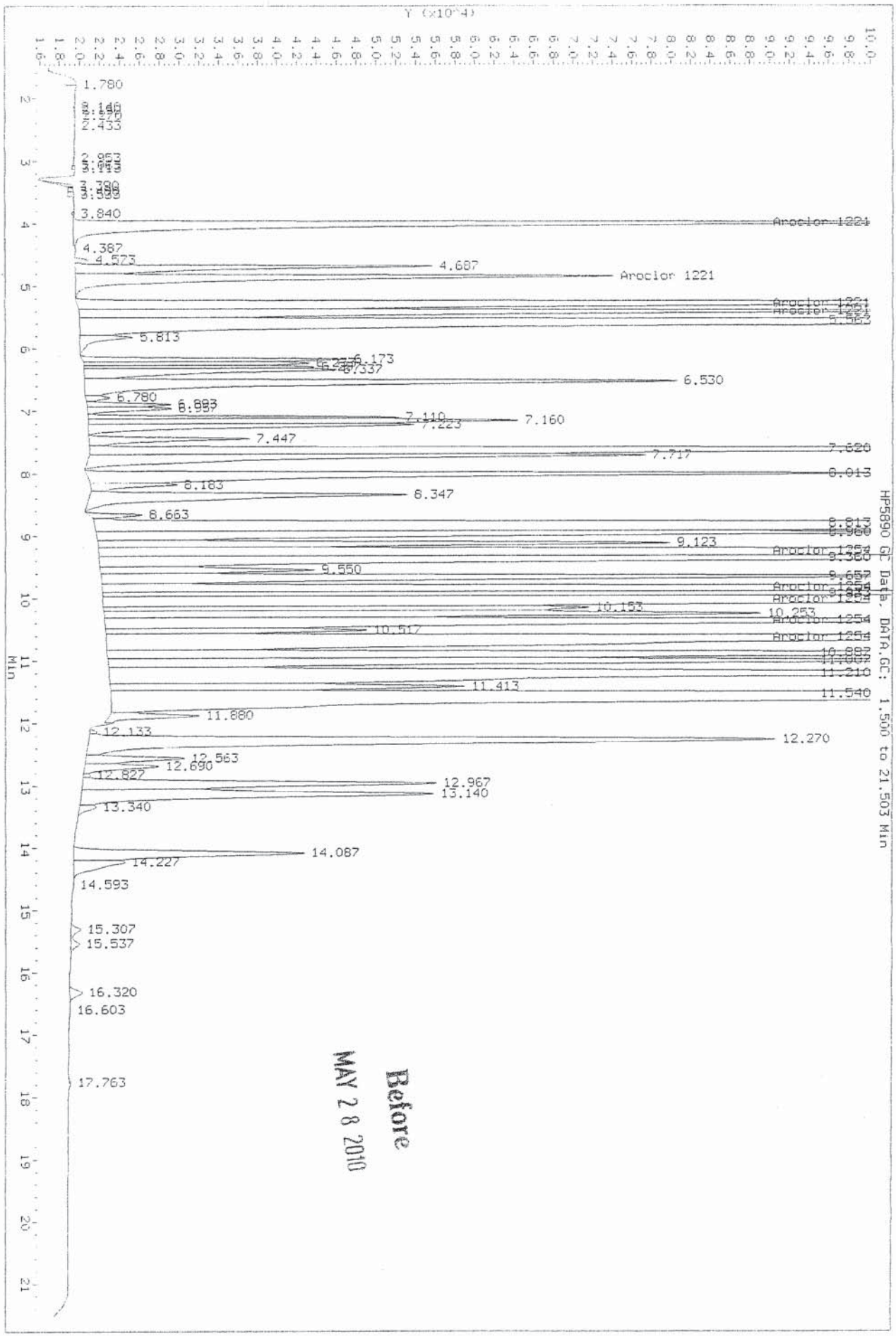
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



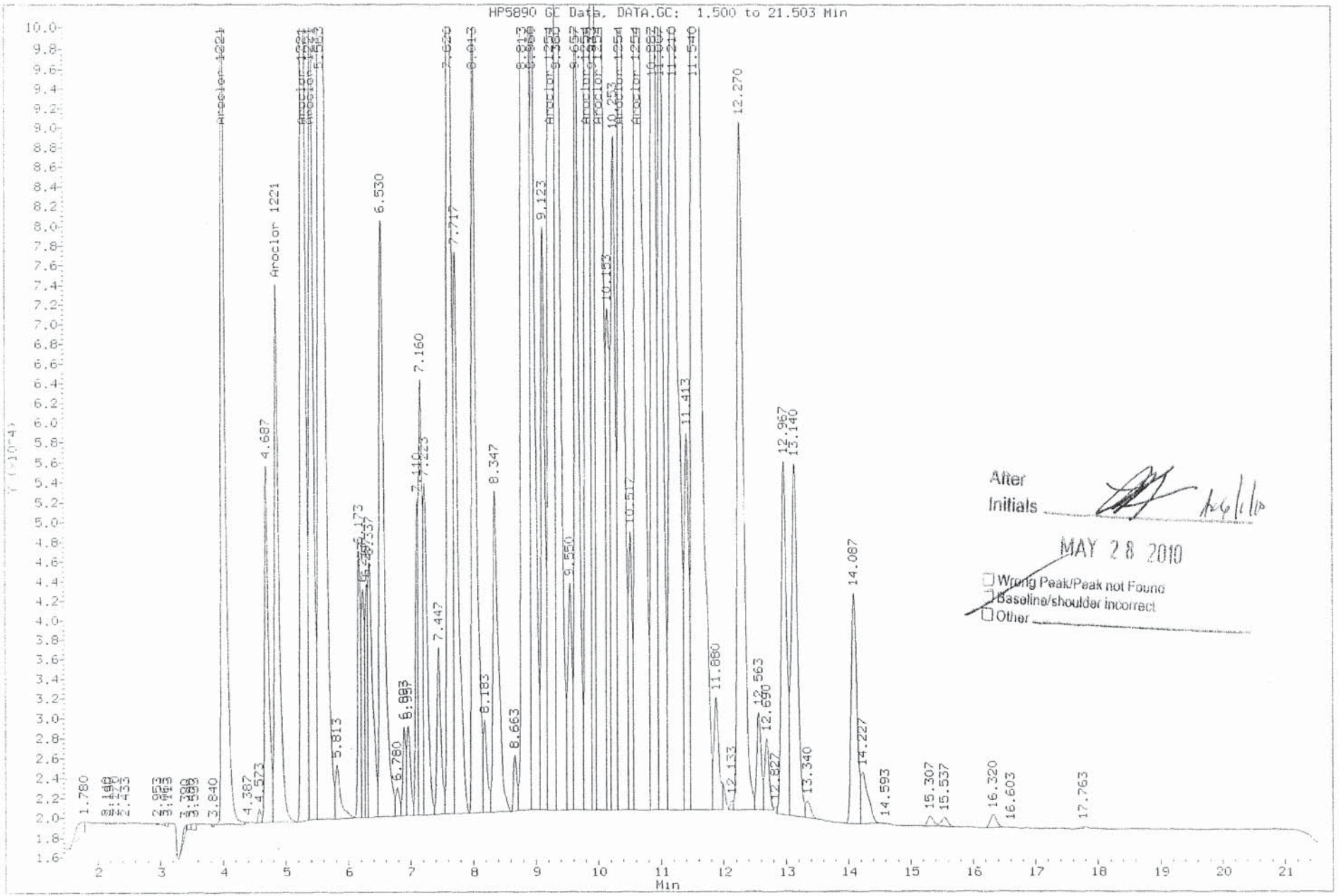
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R033.D
Date : 28-MAY-2010 04:40
Client ID:
Sample Info: 1221/1254 @ 10000/5000ppb | PCB5-52I

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cashl\Acqudata\6C09\data\052710A_r.b\0527R033.D
Injection Date: 28-MAY-2010 04:40
Instrument: GC09.i
Client Sample ID:
Y (x10^4)
HP6890 GC Data, DATA.6C; 1.500 to 21.503 Min
Min
1.780
2.433
3.840
Aroclor 1221
4.387
4.523
4.687
5.813
6.530
6.780
7.160
7.447
7.717
8.183
8.347
8.663
9.123
9.550
10.163
10.517
11.413
11.880
12.133
12.270
12.563
12.690
12.827
12.967
13.140
13.340
14.087
14.227
14.593
15.307
15.537
16.320
16.603
17.763
Before
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R033.D
Injection Date: 28-MAY-2010 04:40
Instrument: GC09.1
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F034.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F034.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R034.D
Inj Date : 28-MAY-2010 05:06
Sample Info: 1232/1262 @ 25ppb | PCB5-52J
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1232+1262.sub
Sub List #2 : 1232+1262.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.567	6575	4758	29.8	26.8	80.00- 120.00	100.00(M)
	5.743	6.230	5934	1365	29.3	21.5	73.60- 110.40	90.24(M)
	6.117	7.223	1118	3538	24.8	27.1	17.45- 26.18	17.01(M)
	6.540	7.443	9911	1399	27.8	19.1	131.13- 196.70	150.72(M)
			Average of Peak Amounts =		27.9	23.6		
Aroclor 1262	10.560	10.893	7377	5679	30.2	23.9	80.00- 120.00	100.00(M)
	10.943	11.550	13889	5042	28.1	24.1	166.27- 249.41	188.28(M)
	11.623	11.743	10597	9064	27.2	26.7	129.39- 194.09	143.66(M)
	12.147	12.313	18661	5818	25.2	22.6	249.74- 374.61	252.96(M)
	13.230	13.150	13311	13769	24.4	24.9	188.68- 283.02	180.45(M)
			Average of Peak Amounts =		27.0	24.4		

QC Flag Legend

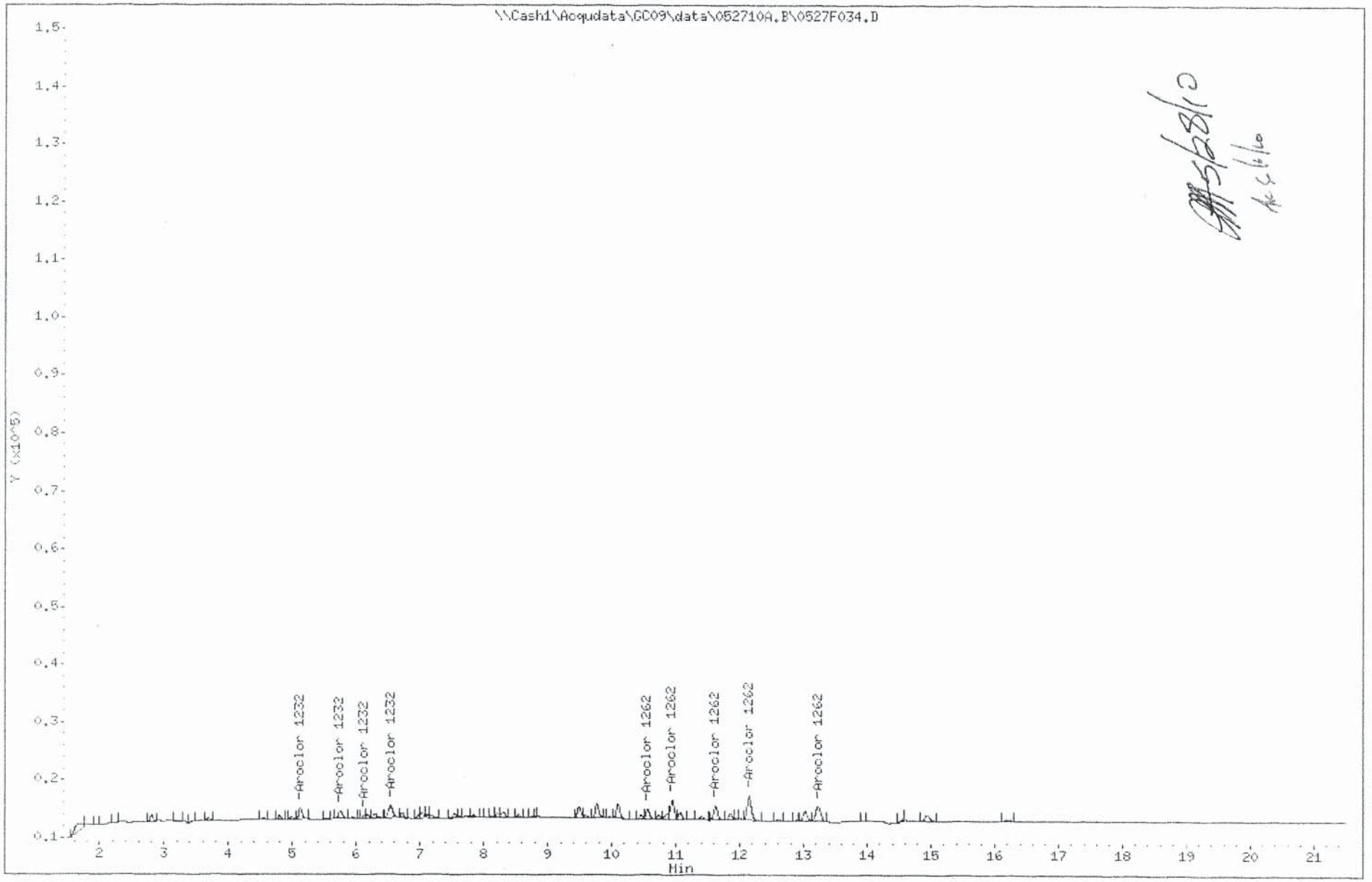
M - Compound response manually integrated.

5/28/10
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F034.D
Date : 28-MAY-2010 05:06
Client ID:
Sample Info: 1232/1262 @ 25ppb | PCB5-62J

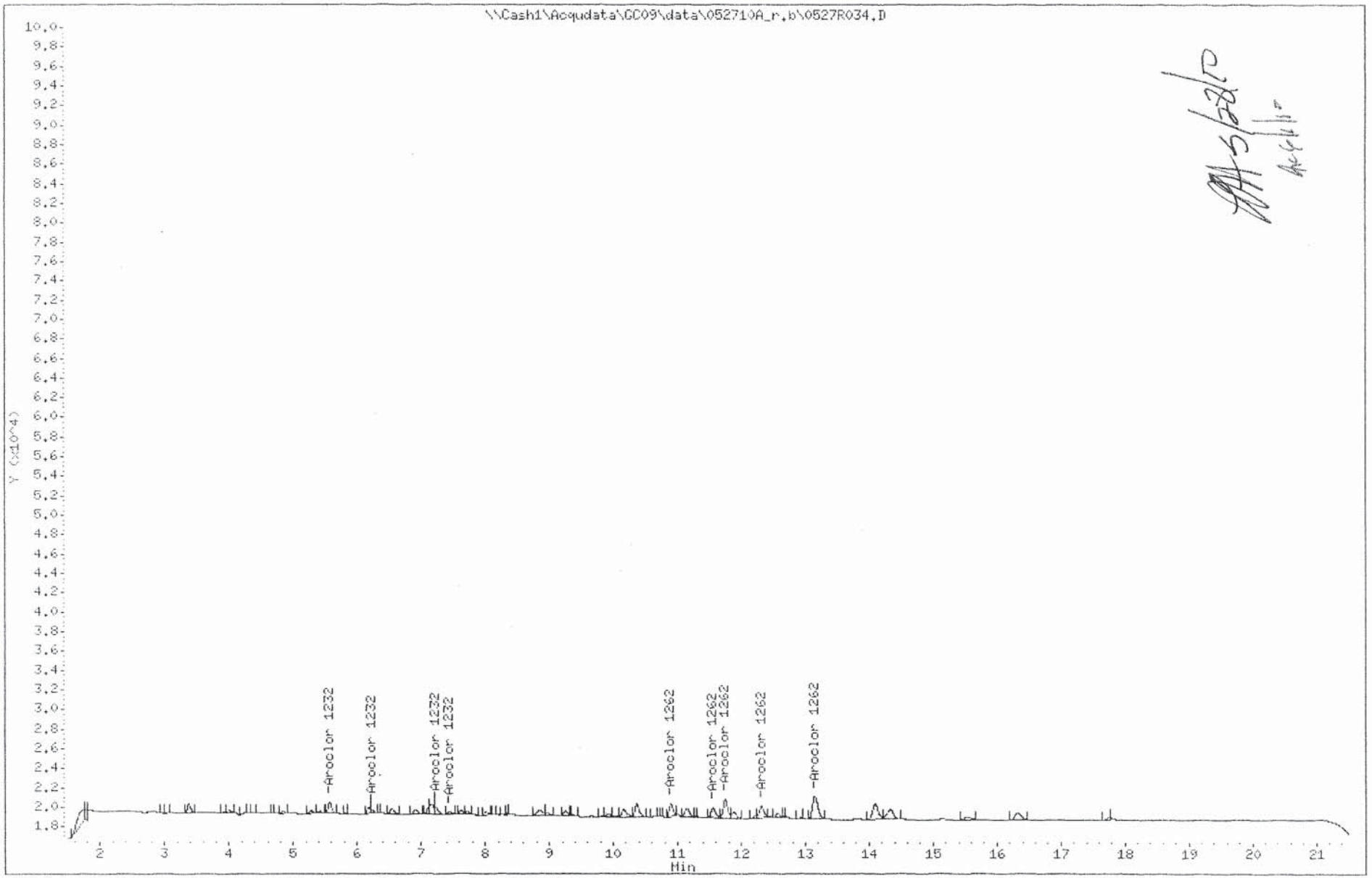
Instrument: GC09.i

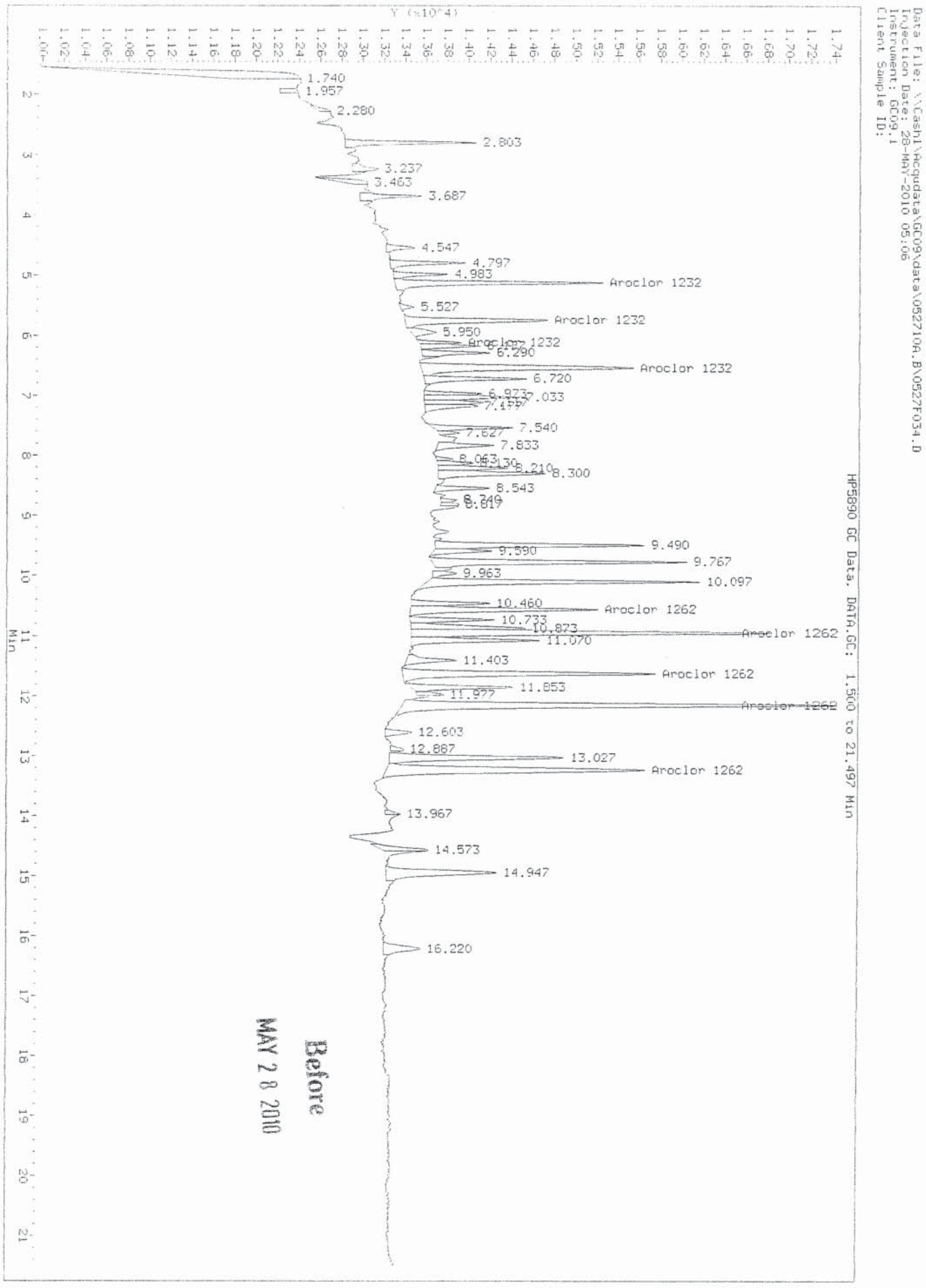
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

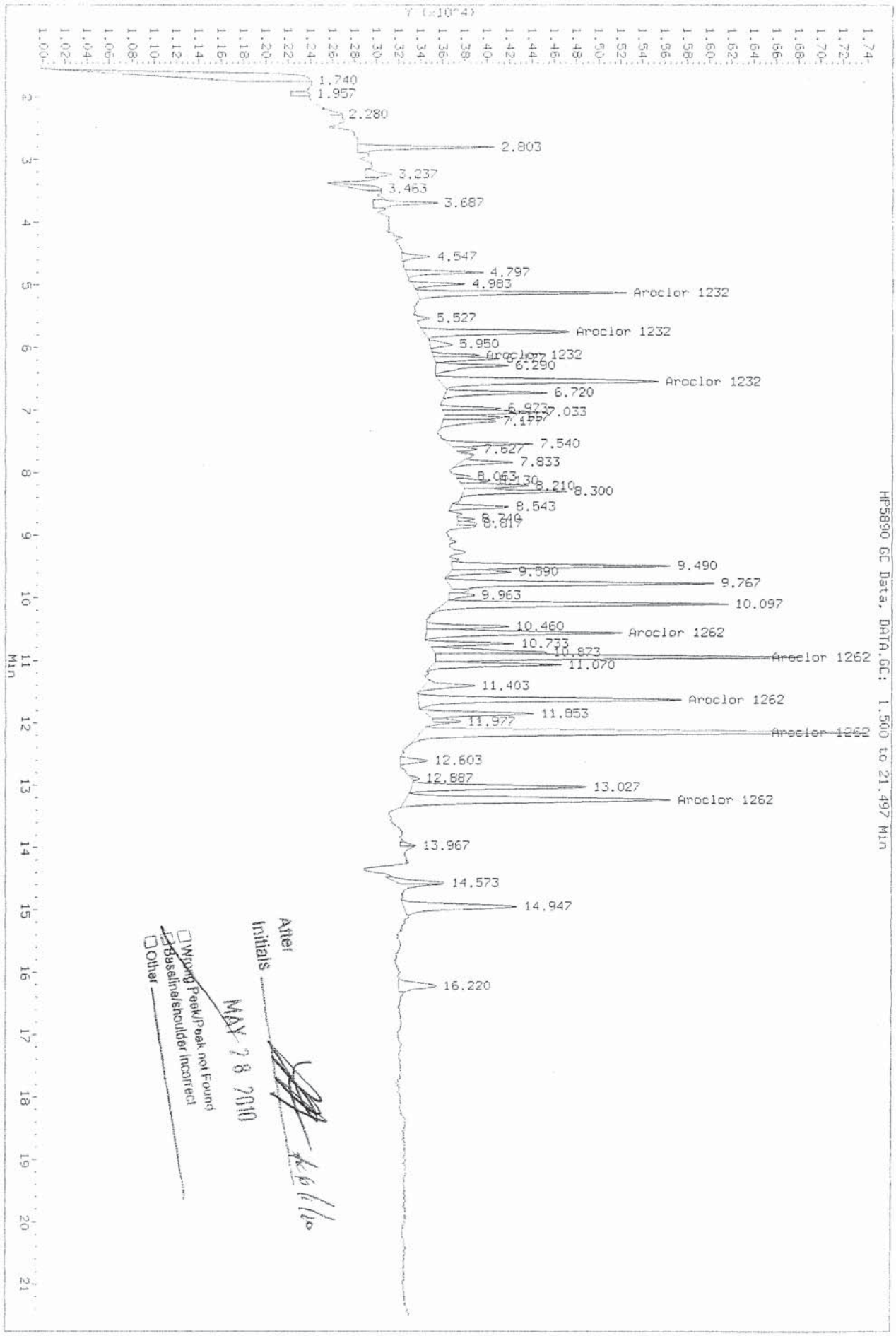


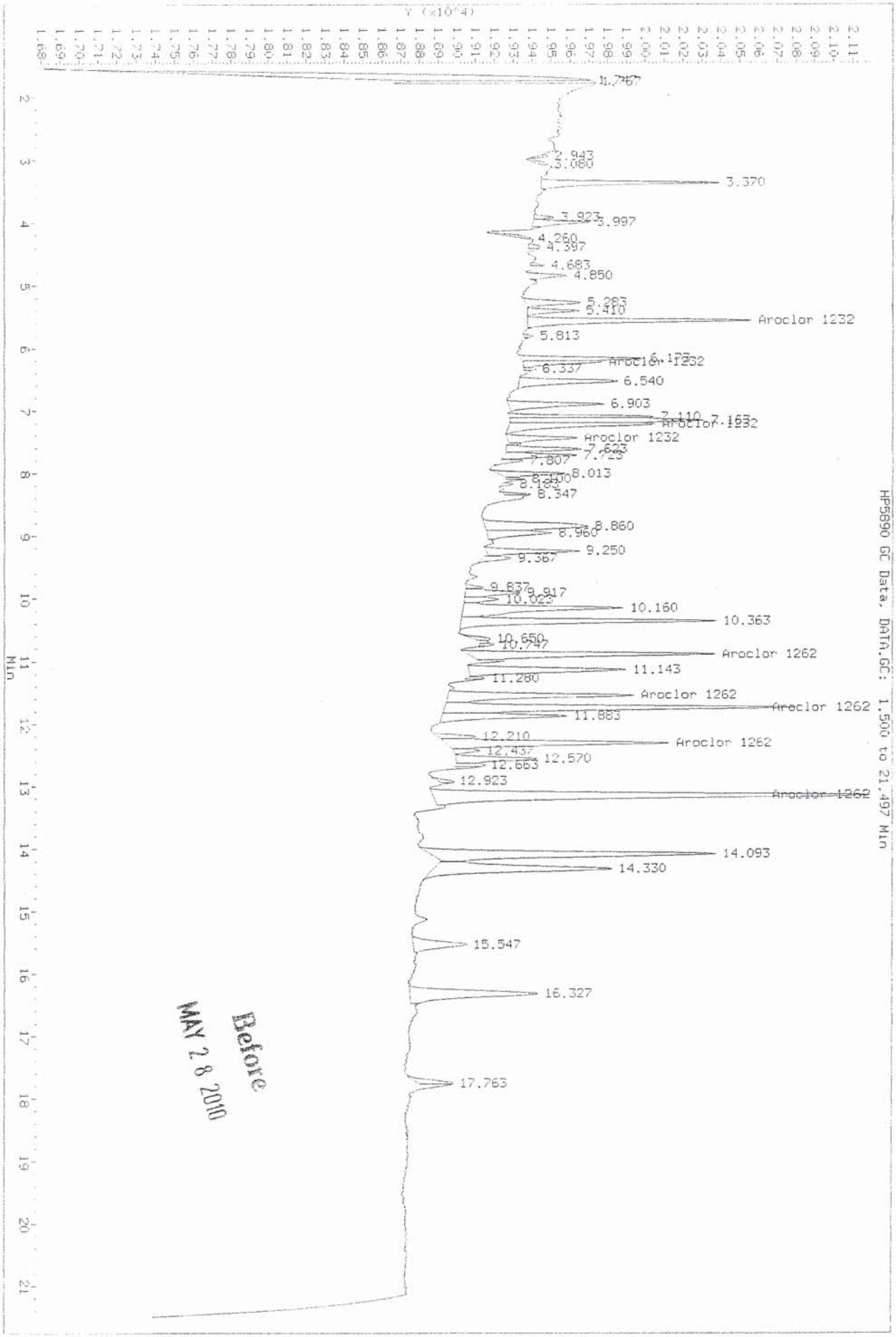
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R034.D
Date : 28-MAY-2010 05:06
Client ID:
Sample Info: 1232/1262 @ 25ppb | PCB5-52J

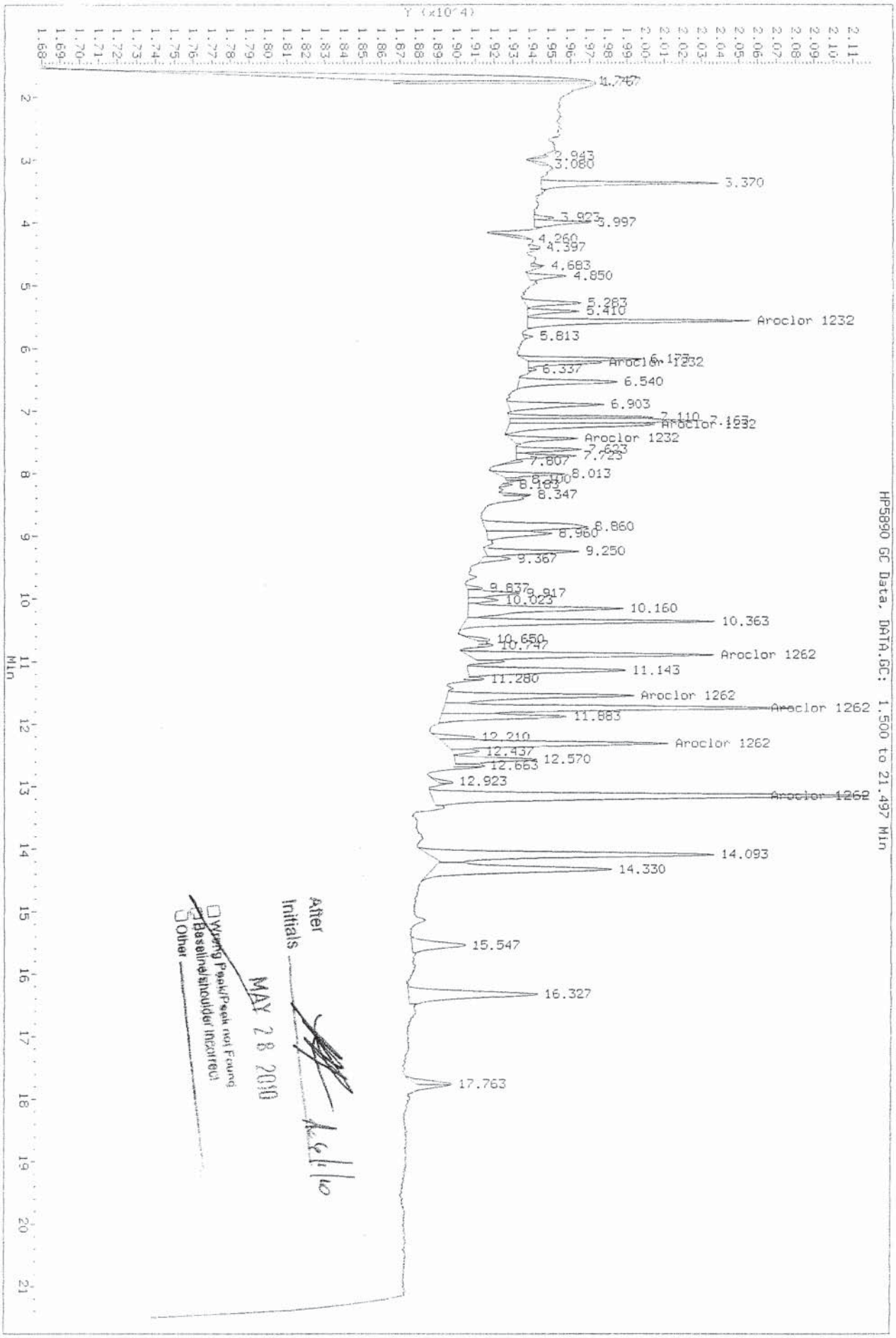
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F034.D
Injection Date: 28-MAY-2010 05:06
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data. DATA.GC; 1.500 to 21.497 Min
Y (x10^4)
Min
1.740
1.957
2.280
2.803
3.237
3.463
3.687
4.547
4.797
4.983
Aroclor 1232
5.527
Aroclor 1232
5.950
Aroclor 1232
6.290
Aroclor 1232
6.720
7.033
7.540
7.833
8.210
8.300
8.543
9.490
9.590
9.767
9.963
10.097
10.460
Aroclor 1262
10.733
10.873
Aroclor 1262
11.070
11.403
Aroclor 1262
11.853
Aroclor 1262
12.603
12.887
13.027
Aroclor 1262
13.967
14.573
14.947
16.220
Before
MAY 2 8 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F034.D
Injection Date: 28-MAY-2010 05:06
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.497 Min
Min
1.740
1.957
2.280
2.803
3.237
3.463
3.687
4.547
4.797
4.983
Aroclor 1232
5.527
Aroclor 1232
5.950
Aroclor 1232
6.290
Aroclor 1232
6.720
7.033
7.540
7.627
7.833
8.210
8.300
8.543
9.490
9.590
9.767
9.963
10.097
10.460
Aroclor 1262
10.733
Aroclor 1262
11.070
11.403
Aroclor 1262
11.853
11.977
Aroclor 1262
12.603
12.887
13.027
Aroclor 1262
13.967
14.573
14.947
16.220
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other


Data File: \\Cashl\Acqudata\GC09\data\052710A_r.b\0527R034.D
Injection Date: 28-MAY-2010 05:06
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.497 Min
Y (x10^4)
Min
11.767
2.943
3.080
3.370
3.923
3.997
4.260
4.397
4.683
4.850
5.283
5.410
Aroclor 1232
5.813
6.337
6.540
6.903
Aroclor 1232
7.623
7.807
8.013
8.347
8.860
8.960
9.250
9.367
10.160
10.363
Aroclor 1262
11.143
11.280
Aroclor 1262
Aroclor 1262
11.883
12.210
Aroclor 1262
12.570
12.923
14.093
14.330
15.547
16.327
17.763
Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R034.D
Injection Date: 28-MAY-2010 05:06
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.497 Min
Y (x10^4)
Min
1.747
2.943
3.080
3.370
3.923
3.997
4.260
4.397
4.683
4.850
5.283
5.410
Aroclor 1232
5.813
6.177
Aroclor 1232
6.337
6.540
6.903
7.110
7.167
Aroclor 1232
Aroclor 1232
7.623
7.723
7.807
8.013
8.347
8.860
8.960
9.250
9.367
9.837
9.917
10.023
10.160
10.363
10.650
10.747
Aroclor 1262
11.143
11.280
Aroclor 1262
Aroclor 1262
11.883
12.210
Aroclor 1262
12.437
12.570
12.663
12.923
Aroclor 1262
14.093
14.330
15.547
16.327
17.763
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F035.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F035.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R035.D
Inj Date : 28-MAY-2010 05:32
Sample Info: 1232/1262 @ 50ppb | PCB5-52K
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1232+1262.sub
Sub List #2 : 1232+1262.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.567	13403	9859	61.4	56.0	80.00- 120.00	100.00(M)
	5.743	6.230	11859	3282	59.0	51.3	73.60- 110.40	90.13(M)
	6.117	7.223	2234	7128	49.5	55.0	17.45- 26.18	16.67(M)
	6.540	7.447	21022	3421	58.9	48.5	131.13- 196.70	156.85(M)
			Average of Peak Amounts =		57.2	52.7		
Aroclor 1262	10.560	10.893	13924	13007	57.0	55.1	80.00- 120.00	100.00(M)
	10.943	11.550	29206	10645	59.0	50.8	166.27- 249.41	209.75(M)
	11.623	11.743	21741	18744	56.0	55.1	129.39- 194.09	156.14(M)
	12.147	12.313	40850	13653	55.6	53.1	249.74- 374.61	293.38(M)
	13.230	13.150	28630	28252	52.6	51.2	188.68- 283.02	205.62(M)
			Average of Peak Amounts =		56.0	53.1		

QC Flag Legend

M - Compound response manually integrated.

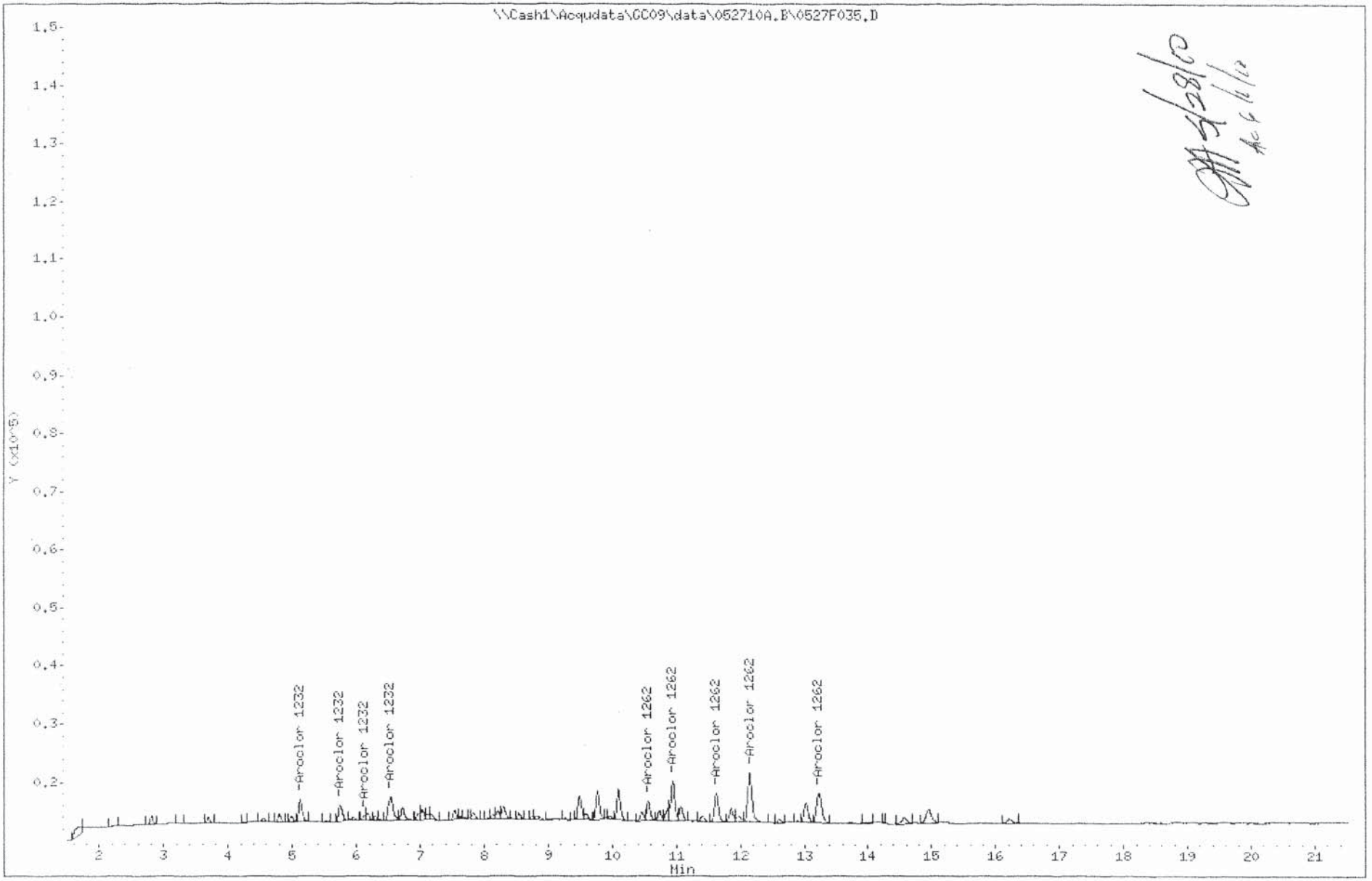
5/28/10

Ac 6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F035.D
Date : 28-MAY-2010 05:32
Client ID:
Sample Info: 1232/1262 @ 50ppb | PCB5-52K

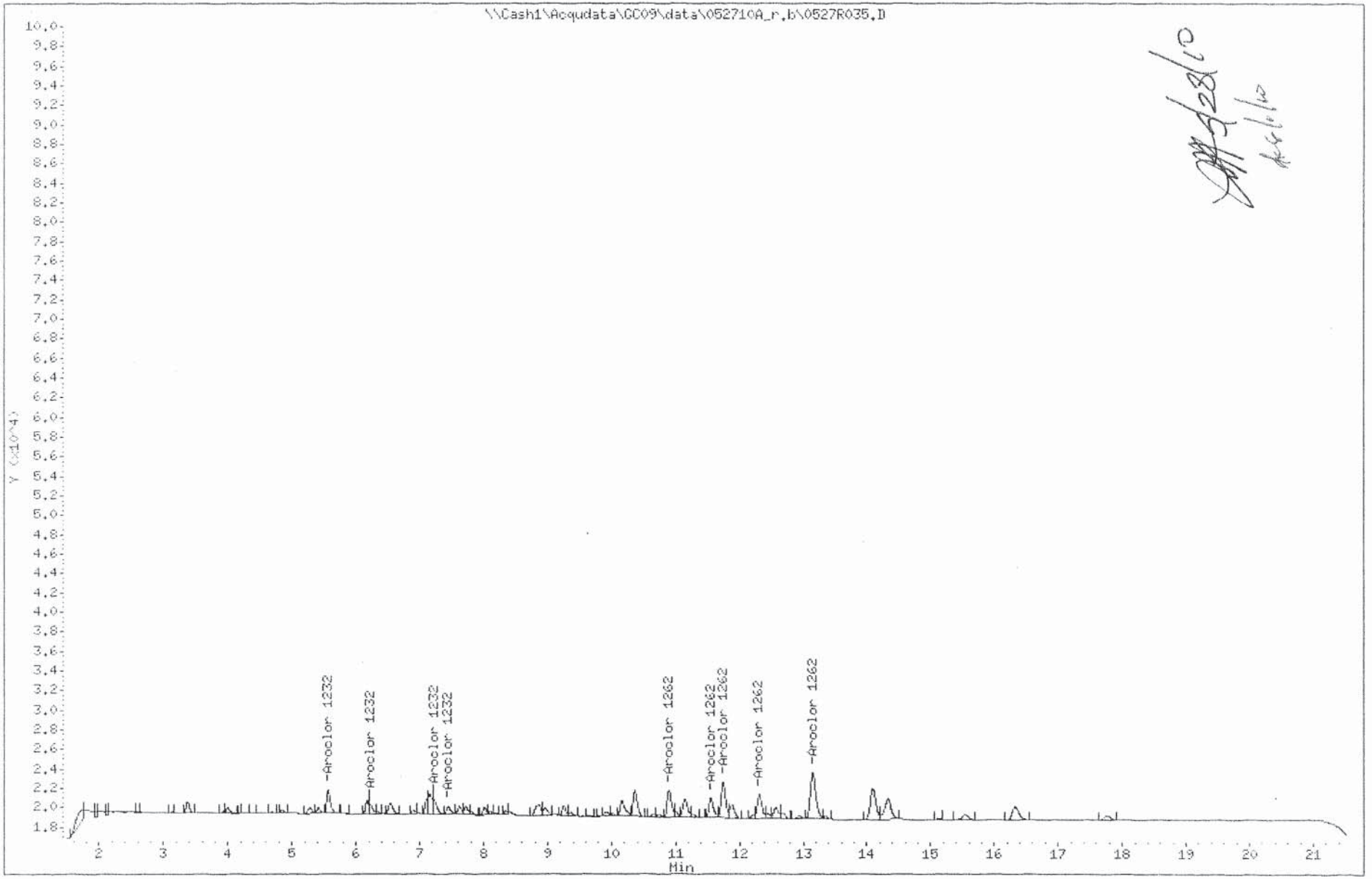
Instrument: GC09.i

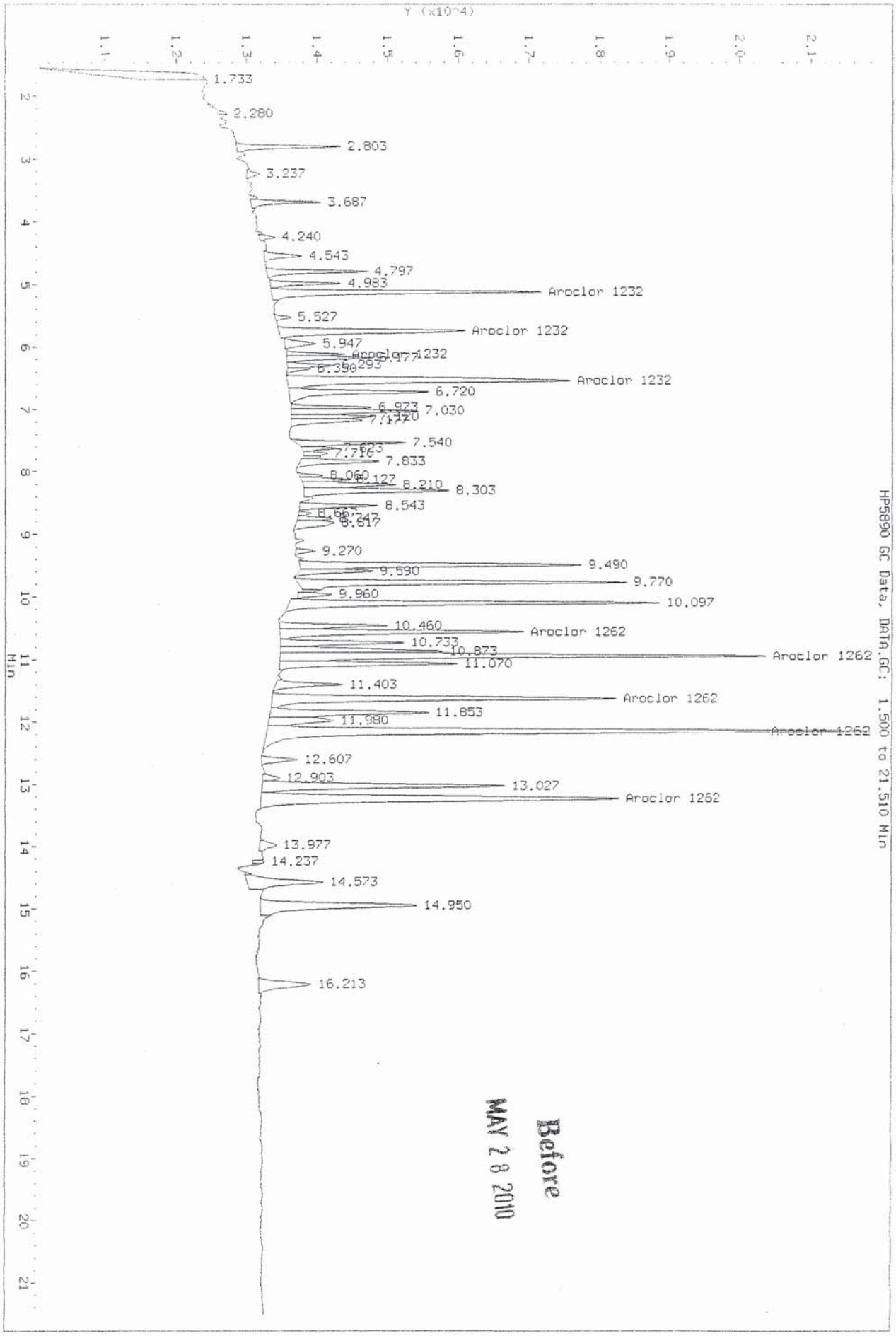
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

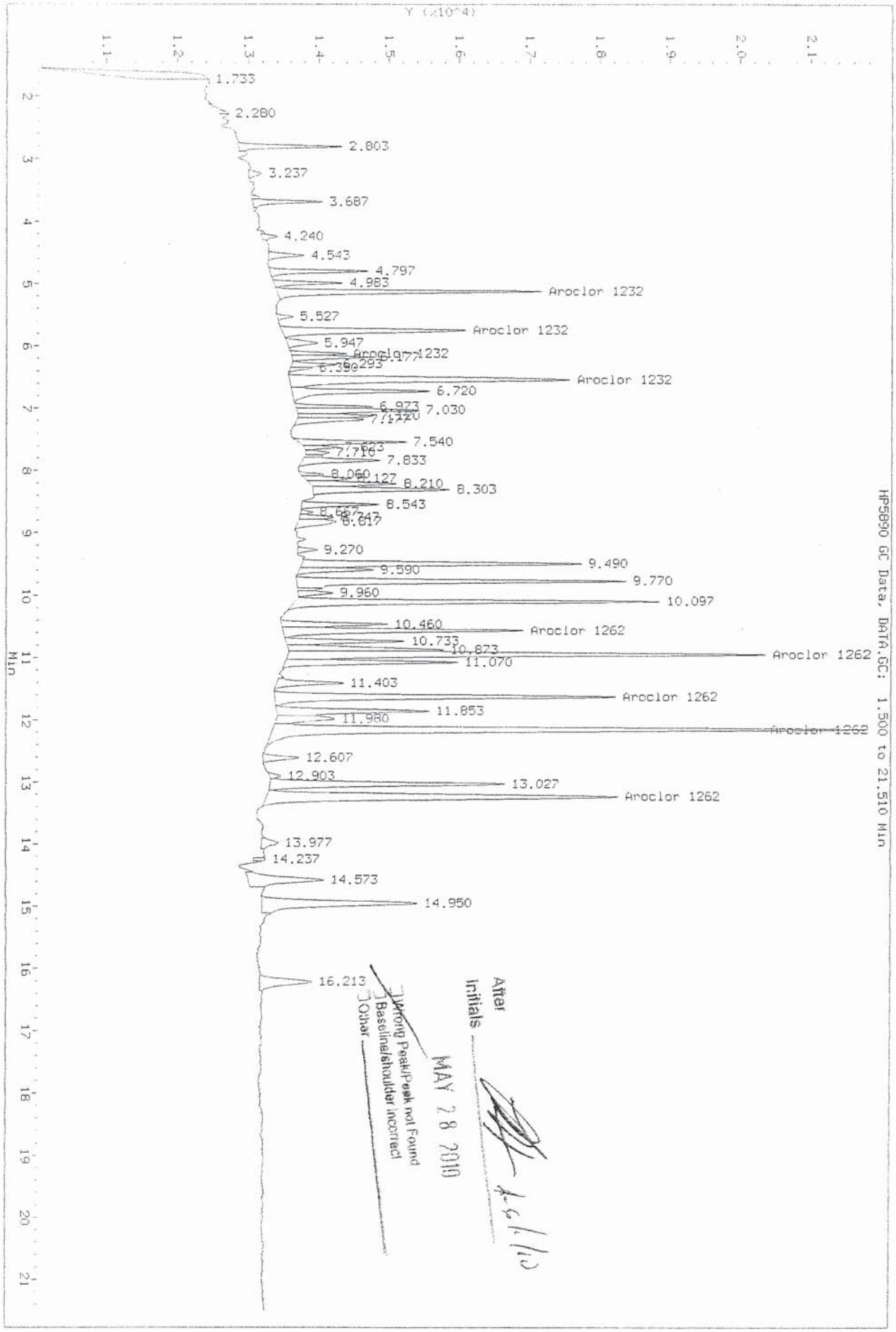


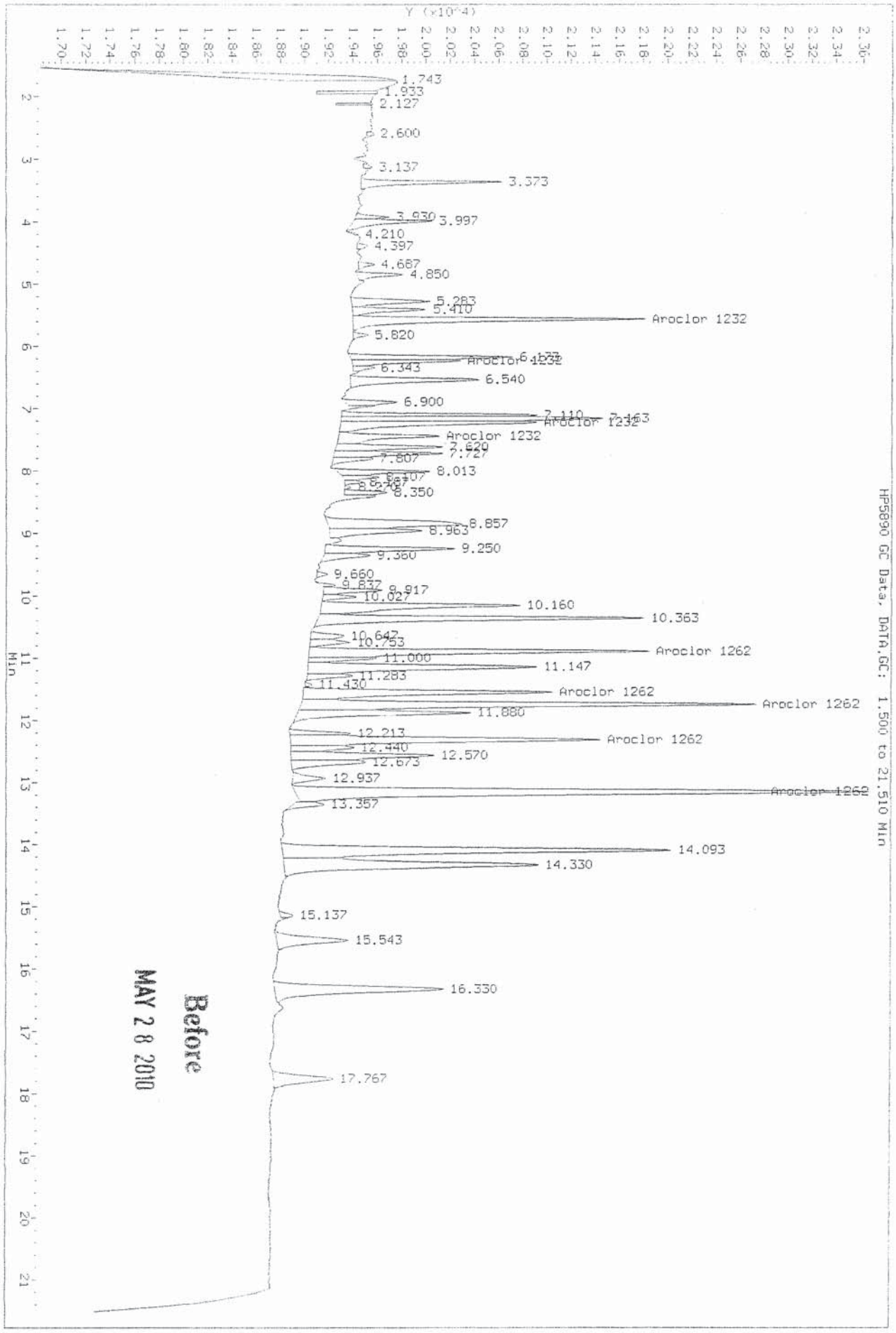
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R035.D
Date : 28-MAY-2010 05:32
Client ID:
Sample Info: 1232/1262 @ 50ppb | PCB5-52K

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F035.D
Injection Date: 28-MAY-2010 05:32
Instrument: GC09.I
Client Sample ID:
Y (x10^4)
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Min
1.733
2.280
2.803
3.237
3.687
4.240
4.543
4.797
4.983
Aroclor 1232
5.527
Aroclor 1232
5.947
Aroclor 1232
Aroclor 1232
6.720
7.030
7.540
7.833
8.210
8.303
8.543
9.270
9.490
9.590
9.770
9.960
10.097
10.460
Aroclor 1262
10.733
10.873
Aroclor 1262
11.070
11.403
Aroclor 1262
11.853
11.980
12.607
12.903
13.027
Aroclor 1262
13.977
14.237
14.573
14.950
16.213
Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F035.D
Injection Date: 28-MAY-2010 05:32
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.510 Min
Y (x10^4)
Min
1.733
2.280
2.803
3.237
3.687
4.240
4.543
4.797
4.983
Aroclor 1232
5.527
Aroclor 1232
5.947
Aroclor 1232
Aroclor 1232
6.720
6.973
7.030
7.540
7.833
8.210
8.303
8.543
9.270
9.490
9.590
9.770
9.960
10.097
10.460
Aroclor 1262
10.733
10.873
Aroclor 1262
11.070
11.403
Aroclor 1262
11.853
11.980
12.607
12.903
13.027
Aroclor 1262
13.977
14.237
14.573
14.950
16.213
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\052RR035.D
Injection Date: 28-MAY-2010 05:32
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.510 Min
Y (x10^4)
Min
1.743
1.933
2.127
2.600
3.137
3.373
3.930
3.997
4.210
4.397
4.687
4.850
5.283
5.410
Aroclor 1232
5.820
Aroclor 1232
6.343
6.540
6.900
7.110
Aroclor 1232
Aroclor 1232
7.629
7.727
7.807
8.013
8.350
8.857
8.963
9.250
9.360
9.660
9.837
8.917
10.027
10.160
10.363
10.647
10.753
Aroclor 1262
11.000
11.147
11.283
11.430
Aroclor 1262
Aroclor 1262
11.880
12.213
Aroclor 1262
12.440
12.570
12.673
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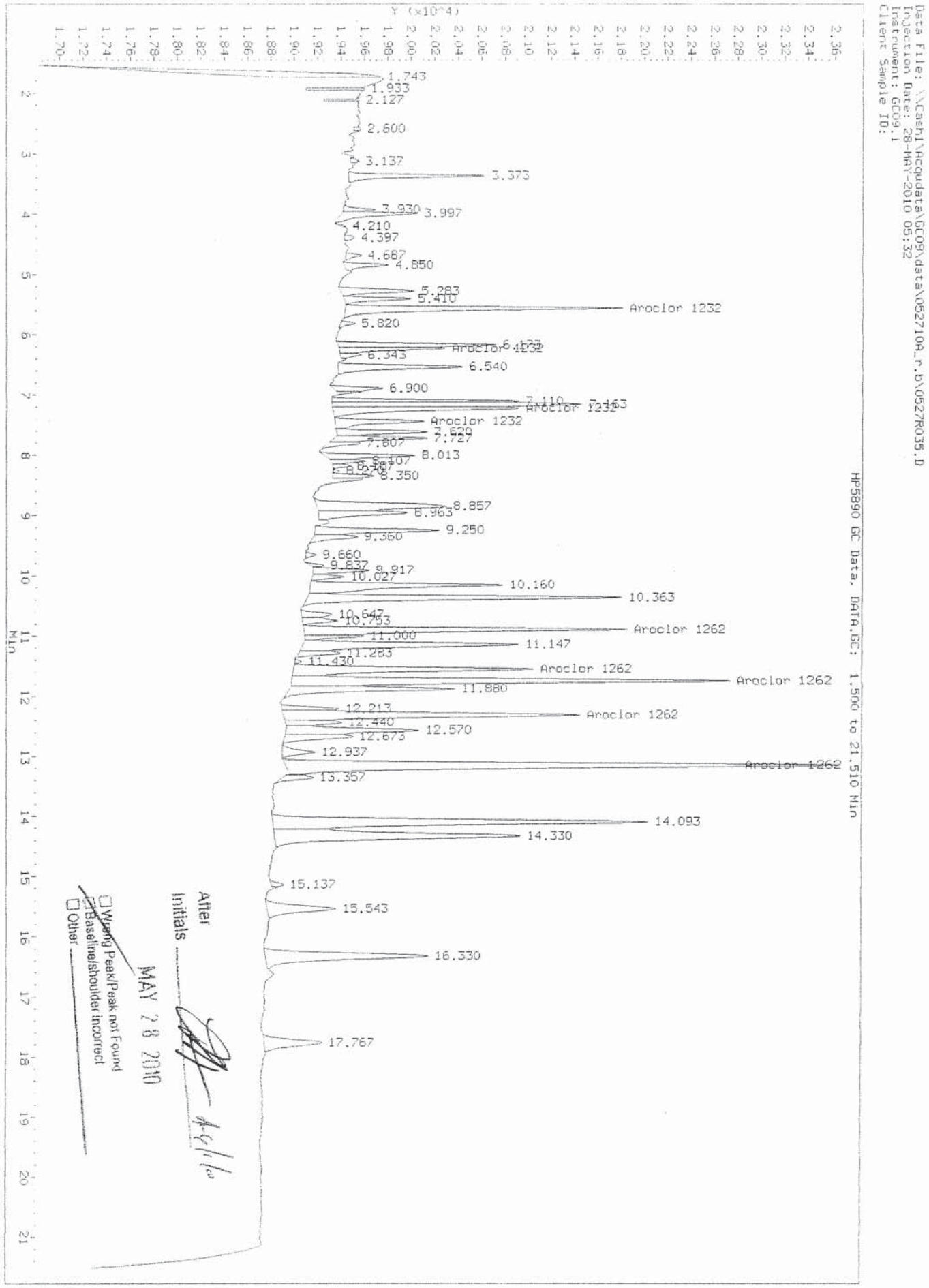
Aroclor 1262
13.357
14.093
14.330
15.137
15.543
16.330
17.767
Before
MAY 2 8 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R035.D
Injection Date: 28-MAY-2010 05:32
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Y (x10^4)
Min
1.743
1.933
2.127
2.600
3.137
3.373
3.930
3.997
4.210
4.397
4.687
4.850
5.283
5.410
Aroclor 1232
5.820
Aroclor 1232
6.345
6.540
6.900
Aroclor 1232
7.807
7.727
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10.027
10.160
10.363
10.753
Aroclor 1262
11.000
11.147
11.283
11.430
Aroclor 1262
Aroclor 1262
11.880
12.213
Aroclor 1262
12.440
12.570
12.675
12.937
Aroclor 1262
13.357
14.093
14.330
15.137
15.543
16.330
17.767
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F036.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F036.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R036.D
Inj Date : 28-MAY-2010 05:58
Sample Info: 1232/1262 @ 500ppb | PCB5-52L
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1232+1262.sub
Sub List #2 : 1232+1262.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.563	121497	100404	556	570	80.00- 120.00	100.00(M)
	5.743	6.230	111779	38905	555	600	73.60- 110.40	92.00(M)
	6.117	7.223	26508	76783	585	573	17.45- 26.18	21.82(M)
	6.537	7.447	199150	47057	558	622	131.13- 196.70	163.91(M)
			Average of Peak Amounts =		564	591		
Aroclor 1262	10.557	10.893	130098	136022	533	575	80.00- 120.00	100.00(M)
	10.943	11.547	270396	117979	546	562	166.27- 249.41	207.84(M)
	11.620	11.743	210418	188626	542	554	129.39- 194.09	161.74(M)
	12.143	12.313	406130	145378	552	564	249.74- 374.61	312.17(M)
	13.227	13.147	306832	305388	564	553	188.68- 283.02	235.85(M)
			Average of Peak Amounts =		547	562		

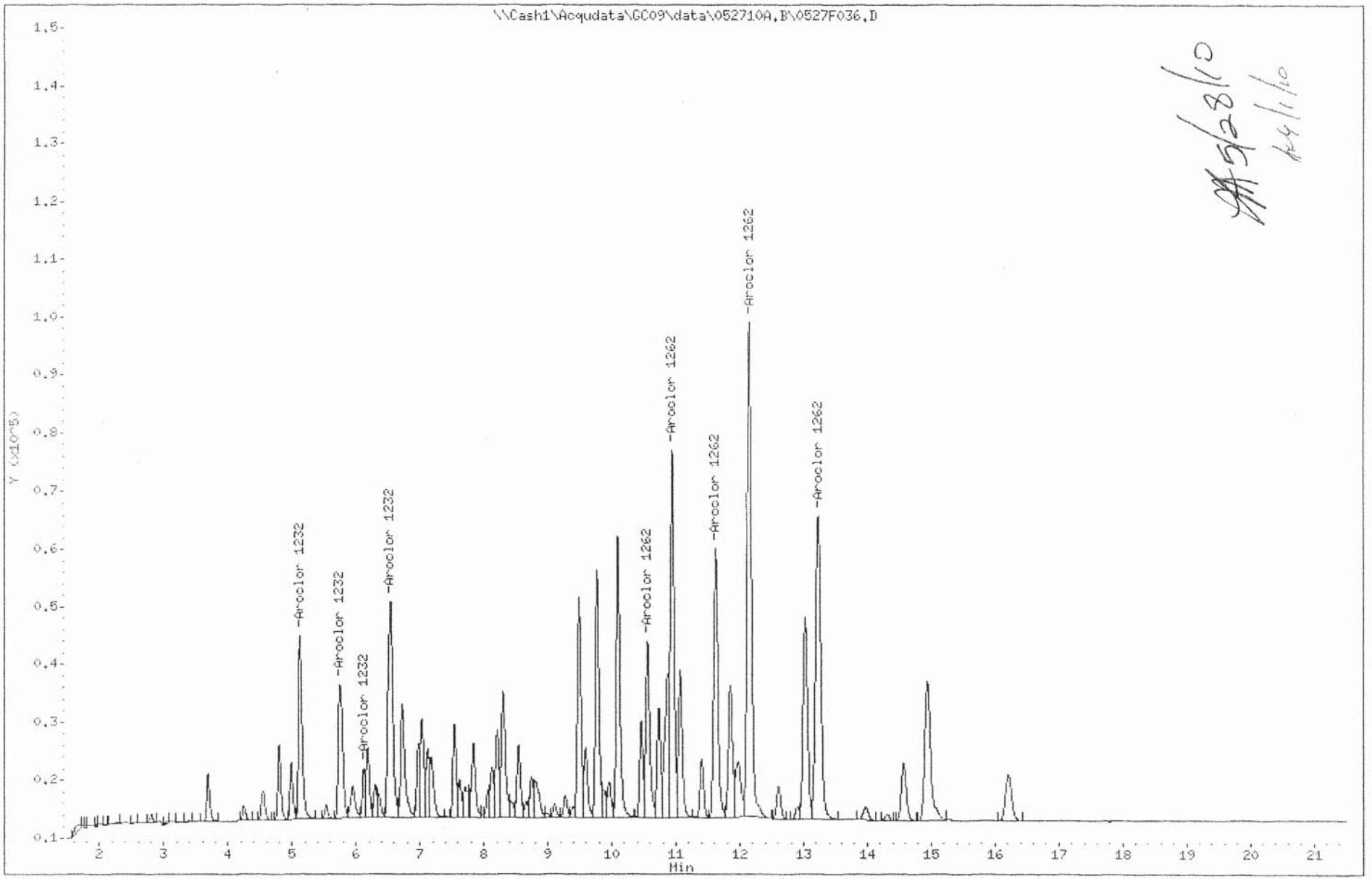
QC Flag Legend

M - Compound response manually integrated.

5/28/10
Acpt 1/10

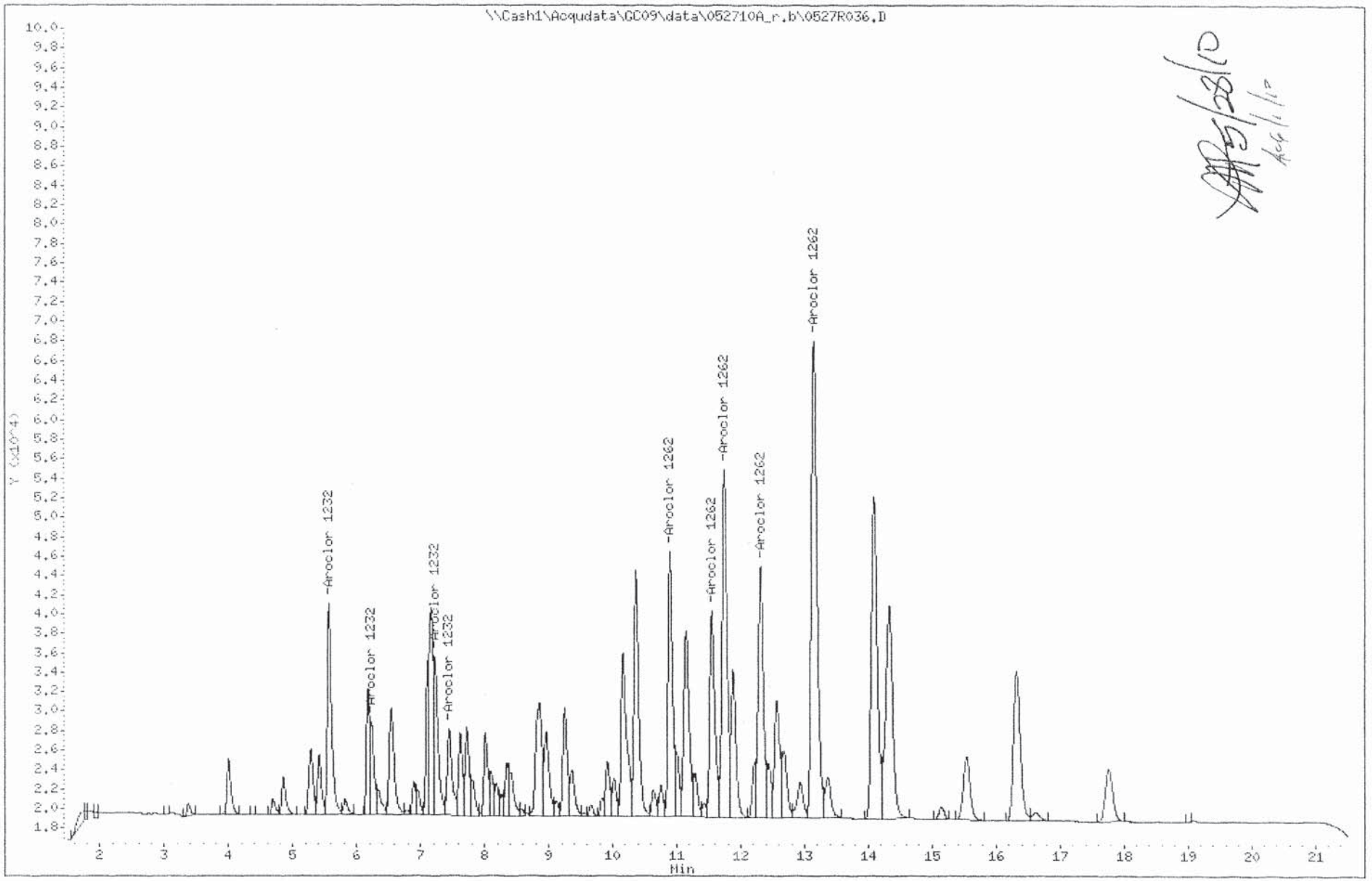
Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F036.D
Date : 28-MAY-2010 05:58
Client ID:
Sample Info: 1232/1262 @ 500ppb | PCB5-52L

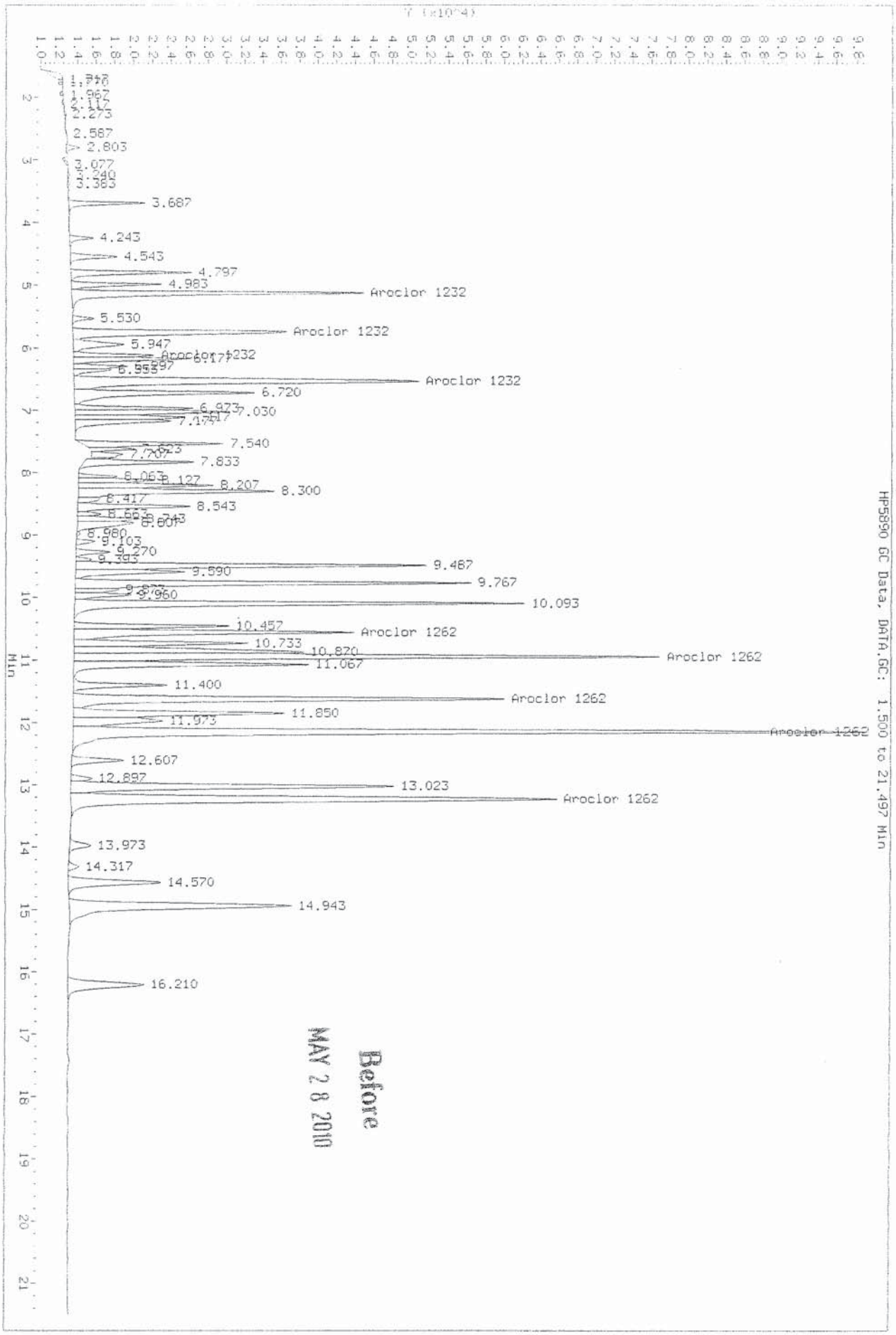
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



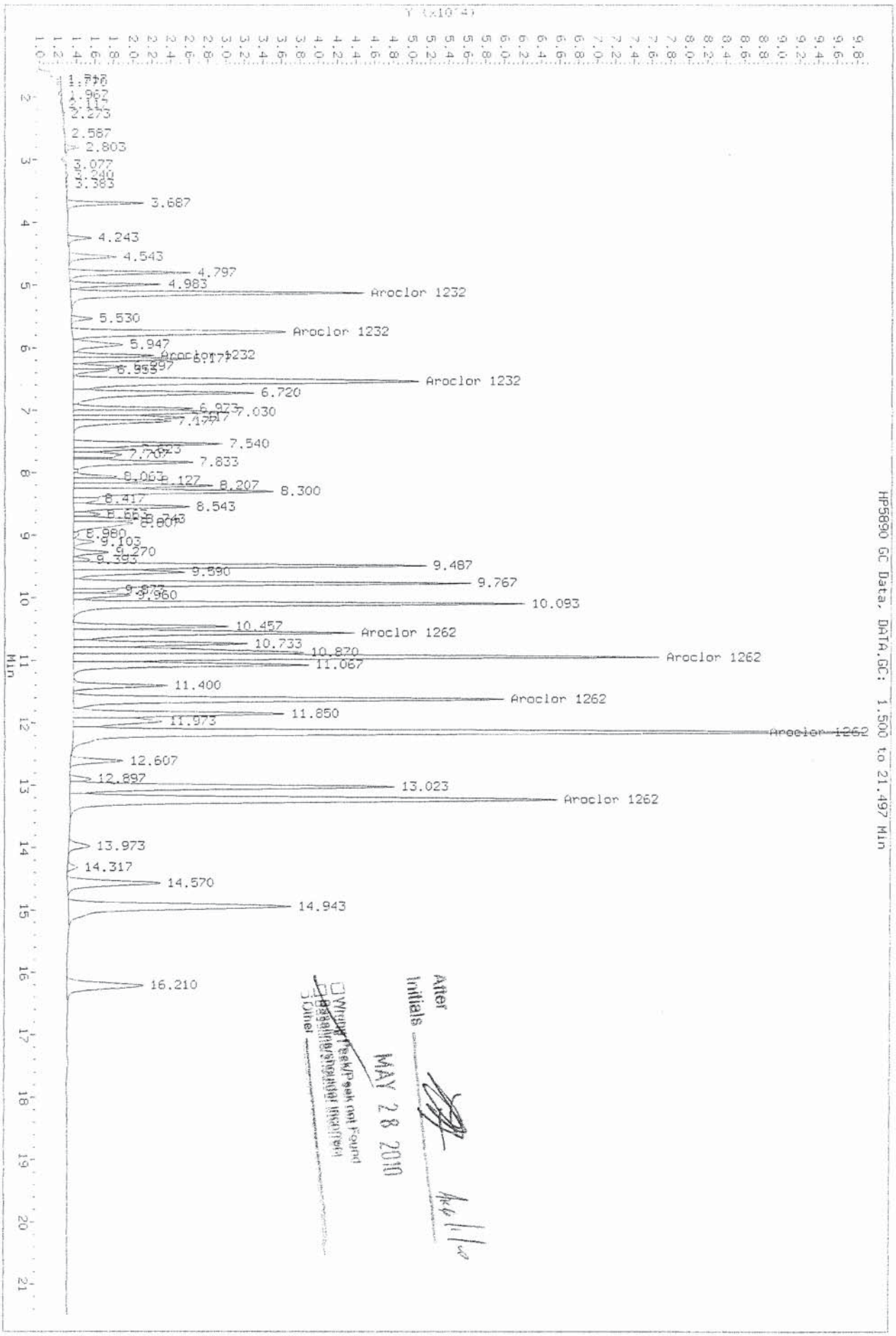
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R036.D
Date : 28-MAY-2010 05:58
Client ID:
Sample Info: 1232/1262 @ 500ppb | PCB5-52L

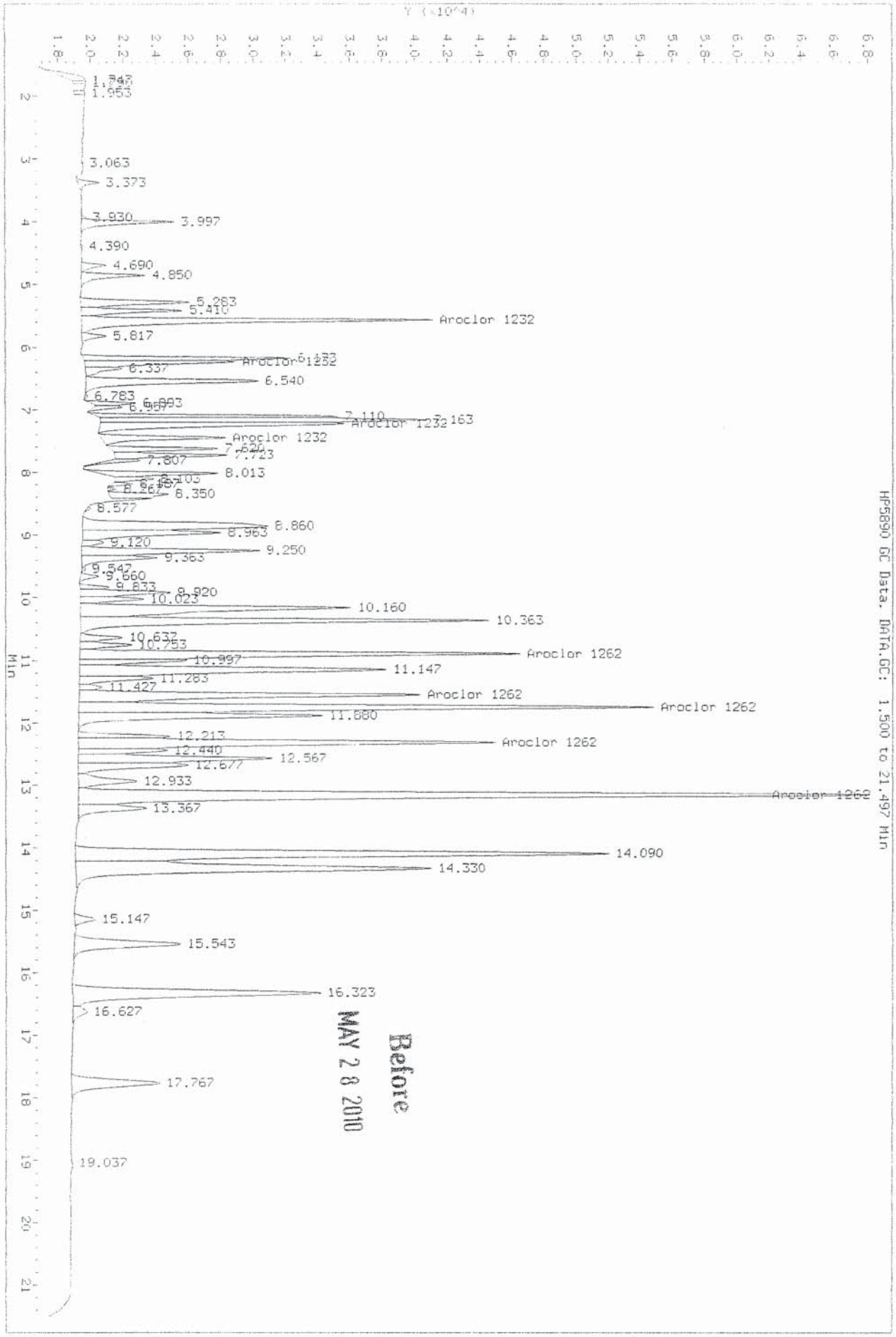
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

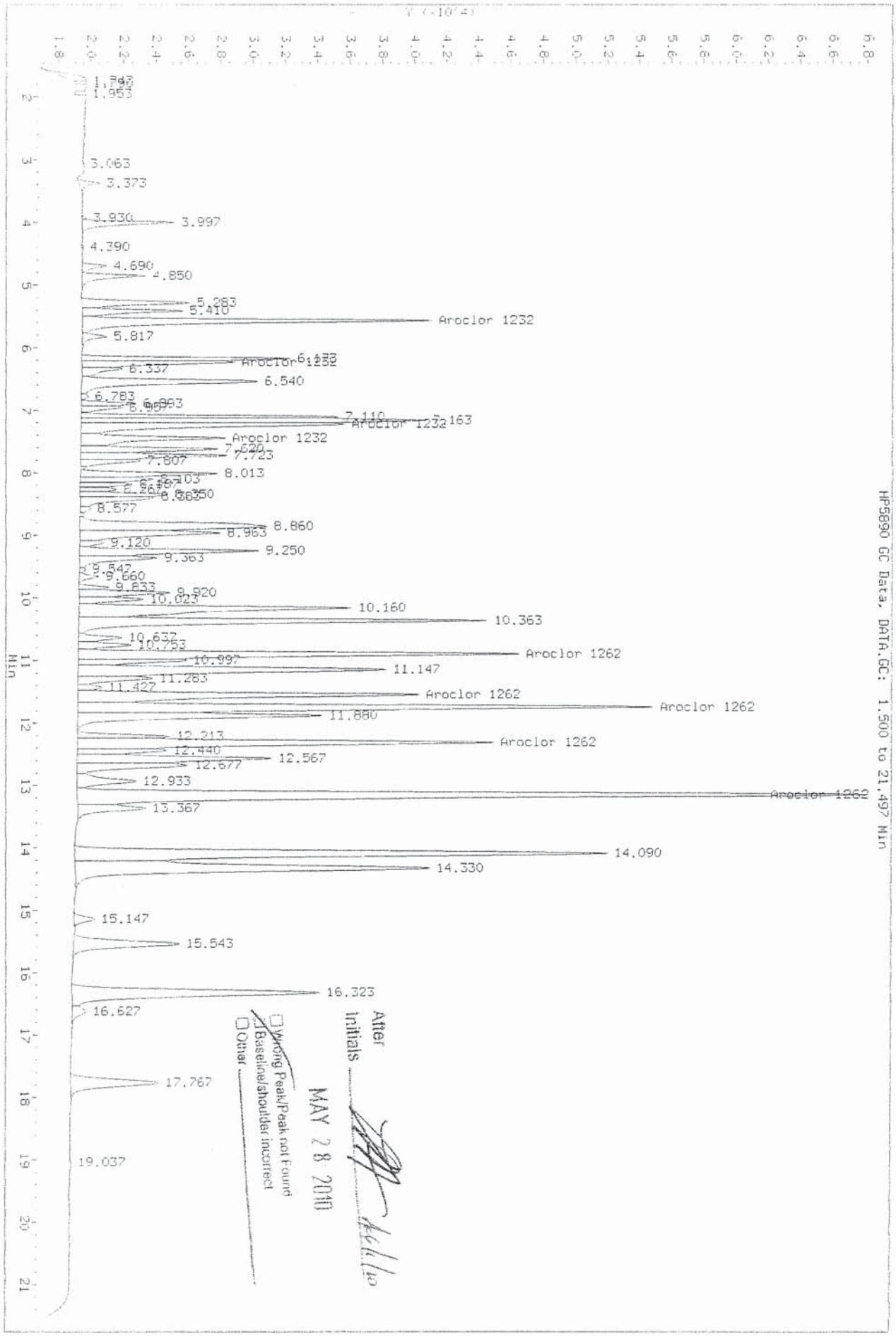




Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F036.D
Injection Date: 28-MAY-2010 05:58
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.497 Min
Y (x10^4)
Min
2.587
2.803
3.687
4.243
4.543
4.797
4.983
Aroclor 1232
5.530
Aroclor 1232
5.947
Aroclor 1232
6.720
7.030
7.540
7.833
8.207
8.300
8.543
9.487
9.767
10.093
10.457
Aroclor 1262
10.733
10.870
Aroclor 1262
11.067
11.400
Aroclor 1262
11.850
11.973
Aroclor 1262
12.607
12.897
13.023
Aroclor 1262
13.973
14.317
14.570
14.943
16.210
After
Initials
MAY 28 2010
Other


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R036.D
Injection Date: 28-MAY-2010 05:58
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.497 Min
Y (x10^4)
Min
1.743
1.953
3.063
3.373
3.930
3.997
4.390
4.690
4.850
5.283
5.410
Aroclor 1232
5.817
6.337
6.540
6.783
7.110
Aroclor 1232
Aroclor 1232
7.807
7.723
8.013
8.350
8.577
8.860
8.963
9.120
9.250
9.363
9.542
9.660
9.833
10.023
10.160
10.363
10.637
10.753
Aroclor 1262
10.997
11.147
11.283
11.427
Aroclor 1262
Aroclor 1262
11.680
12.213
Aroclor 1262
12.440
12.567
12.677
12.933
Aroclor 1262
13.367
14.090
14.330
15.147
15.543
16.323
16.627
17.767
19.037
Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R036.D
Injection Date: 28-MAY-2010 05:58
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.497 Min
Y (x10^4)
Min
1.963
3.063
3.373
3.930
3.997
4.390
4.690
4.850
5.083
5.410
Aroclor 1232
5.817
6.337
6.540
6.783
7.110
Aroclor 1232
Aroclor 1232
7.807
7.923
8.013
8.577
8.860
8.963
9.120
9.250
9.363
9.660
9.833
8.920
10.623
10.160
10.363
10.637
10.753
Aroclor 1262
10.997
11.147
11.283
11.422
Aroclor 1262
Aroclor 1262
11.880
12.213
Aroclor 1262
12.440
12.567
12.677
12.933
Aroclor 1262
13.367
14.090
14.330
15.147
15.543
16.323
16.627
17.767
19.037
After Initials
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F037.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F037.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R037.D
Inj Date : 28-MAY-2010 06:24
Sample Info: 1232/1262 @ 1000ppb | PCB5-52M
Misc Info :
Cal Date : 28-MAY-2010 12:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1232+1262.sub
Sub List #2 : 1232+1262.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.563	208284	176778	954	1000	80.00- 120.00	100.00(M)
	5.747	6.230	194171	68619	962	1060	73.60- 110.40	93.22(M)
	6.120	7.223	47599	133858	1060	999	17.45- 26.18	22.85(M)
	6.540	7.447	349443	83549	979	1100	131.13- 196.70	167.77(M)
			Average of Peak Amounts =		989	1040		
Aroclor 1262	10.560	10.893	231703	240748	949	1020	80.00- 120.00	100.00(M)
	10.943	11.547	481660	213920	973	1020	166.27- 249.41	207.88(M)
	11.623	11.743	383332	337267	987	991	129.39- 194.09	165.44(M)
	12.143	12.313	740624	264534	1010	1030	249.74- 374.61	319.64(M)
	13.230	13.147	558903	563040	1030	1020	188.68- 283.02	241.22(M)
			Average of Peak Amounts =		990	1020		

QC Flag Legend

M - Compound response manually integrated.

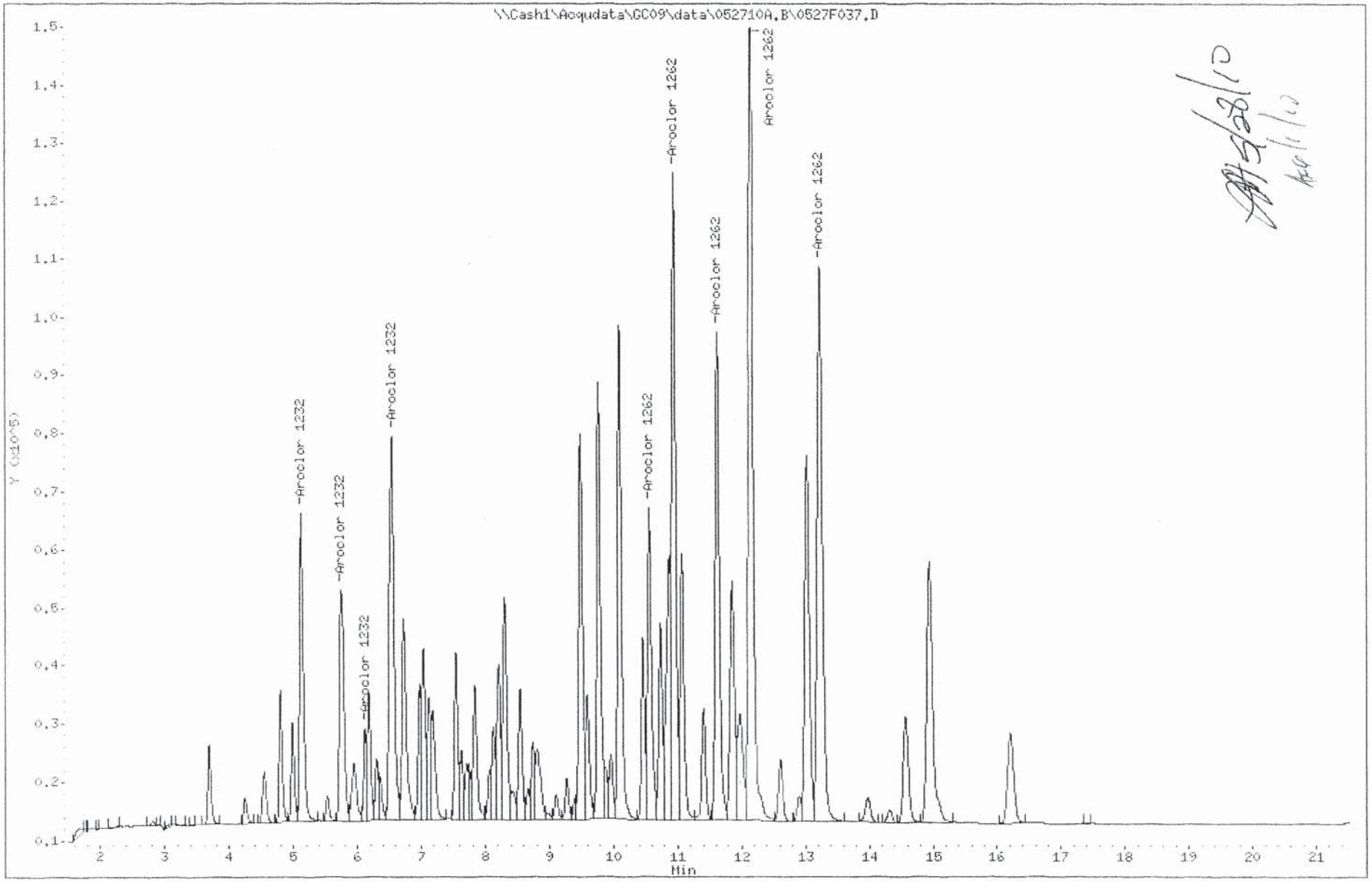
5/28/10
6/4/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F037.D
Date : 28-MAY-2010 06:24
Client ID:
Sample Info: 1232/1262 @ 1000ppb | PCB5-52M

Instrument: GC09.i

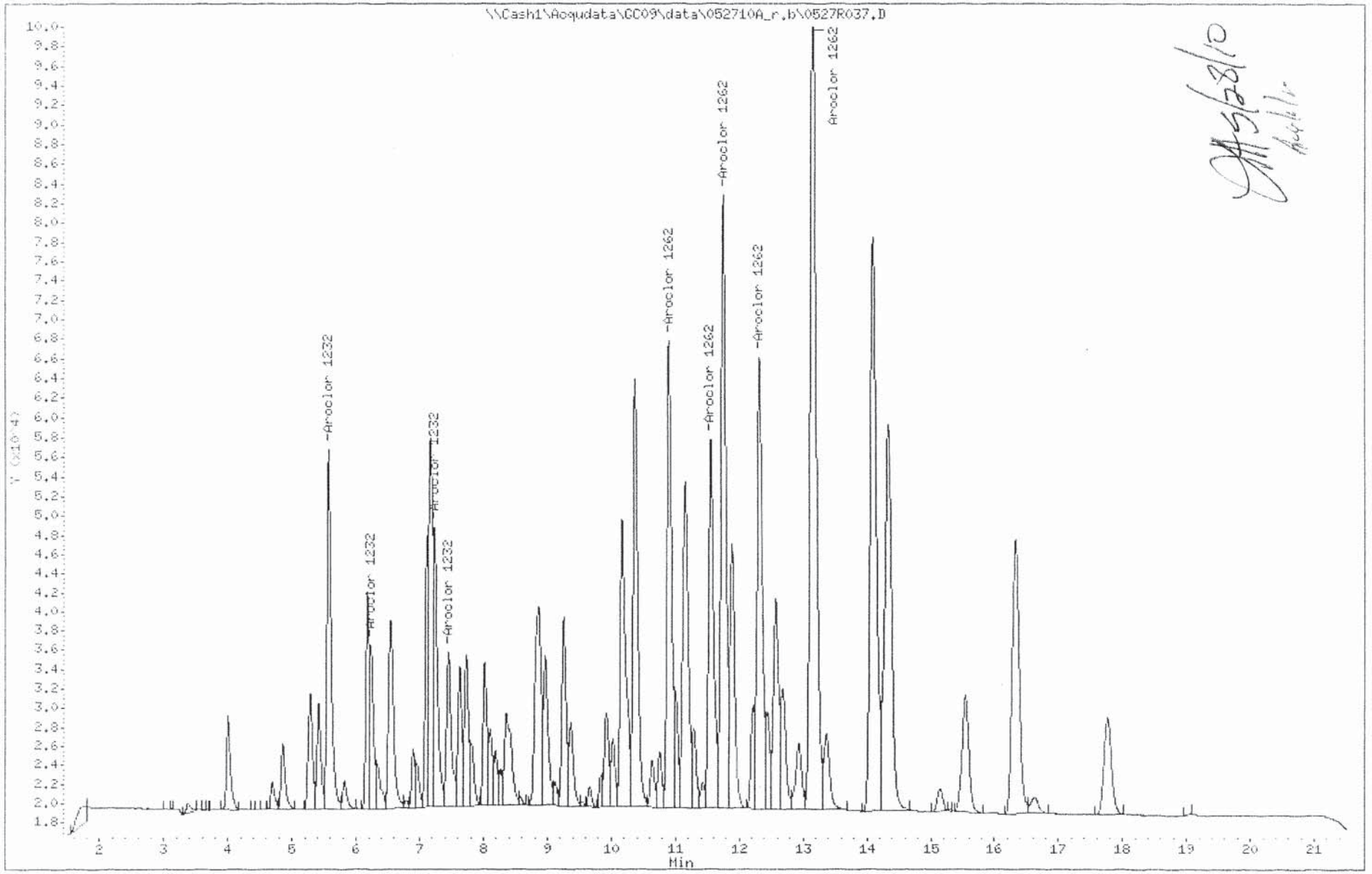
Operator: LHarris
Column diameter: 0.53

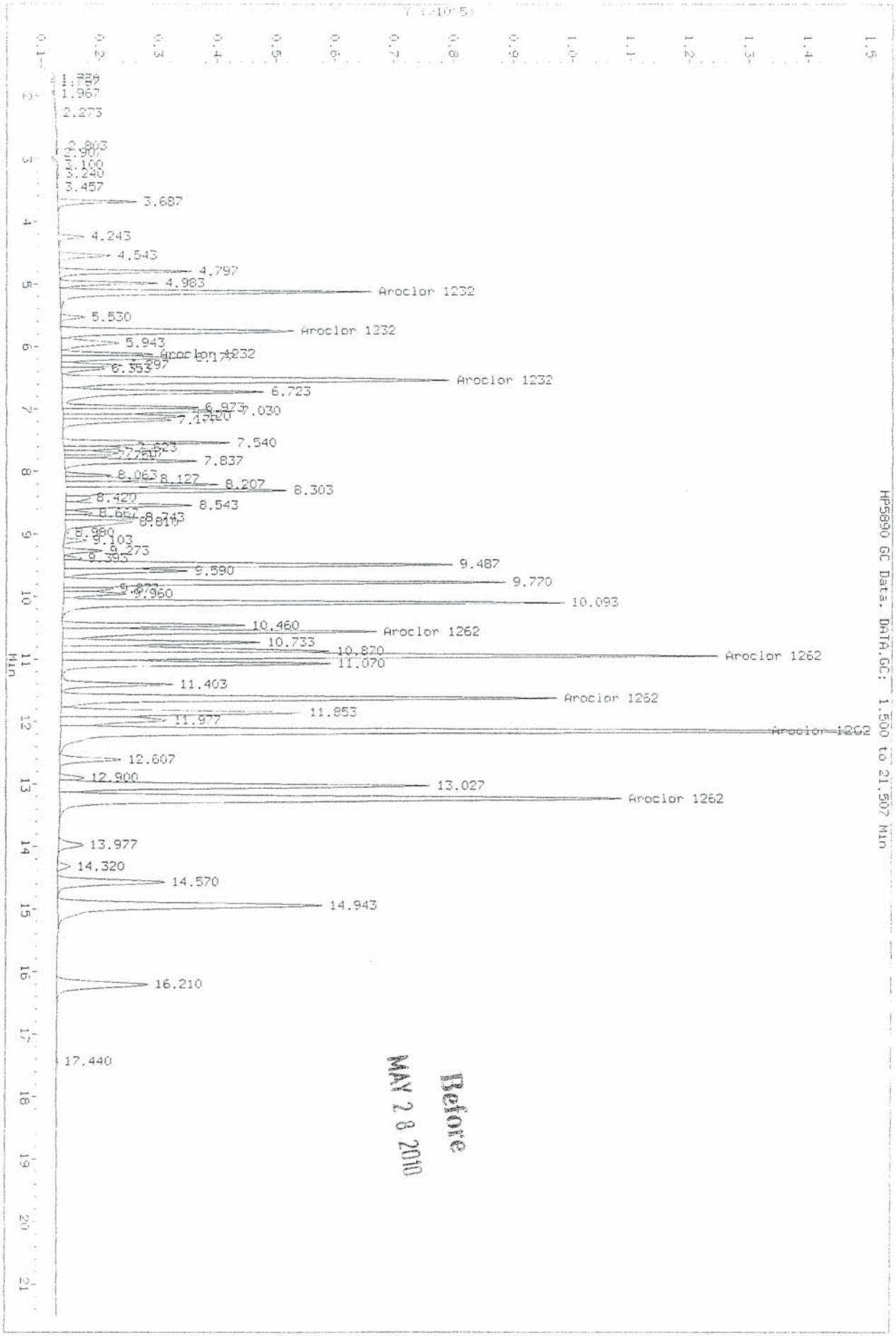
Column phase: DB-35MS



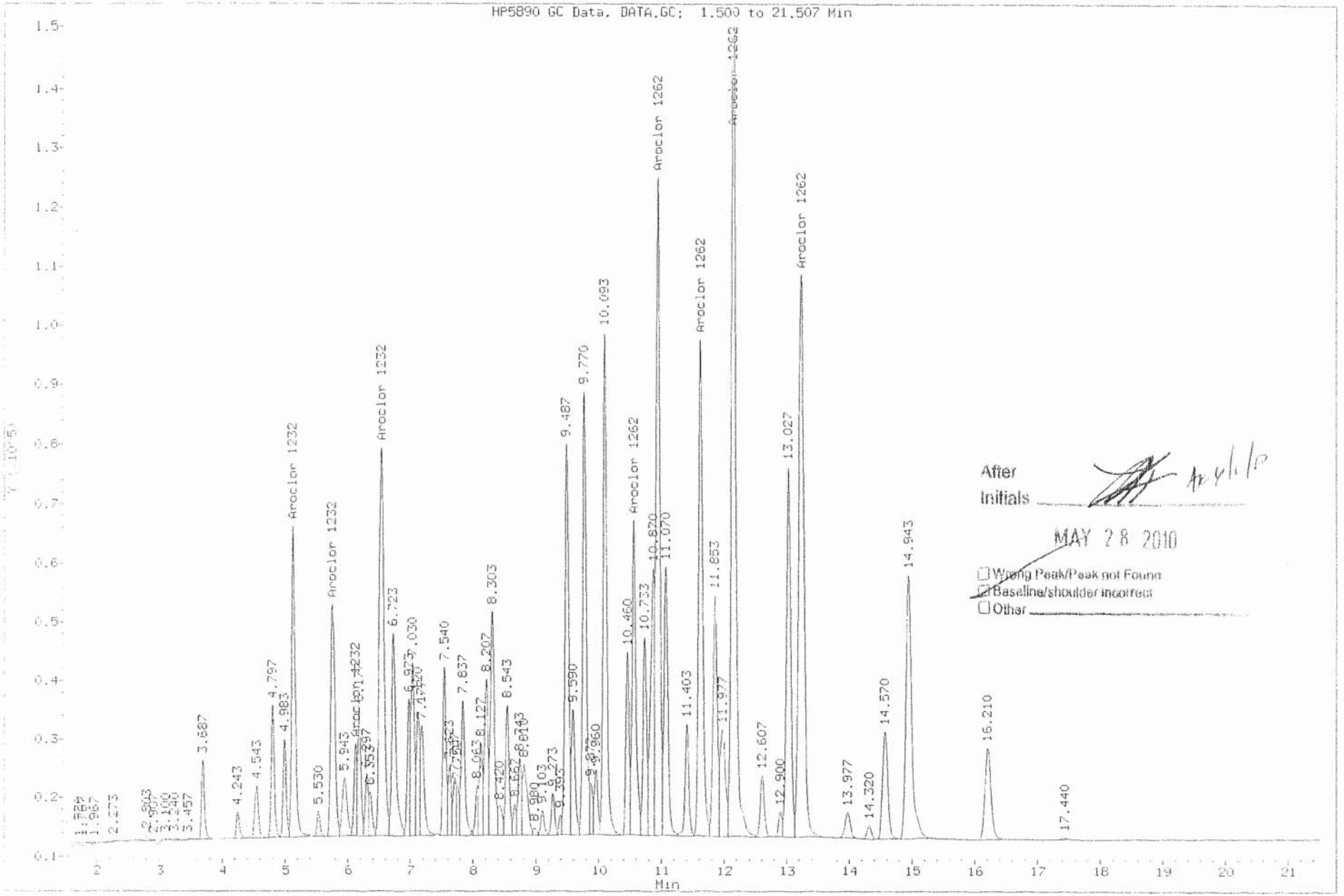
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R037.D
Date : 28-MAY-2010 06:24
Client ID:
Sample Info: 1232/1262 @ 1000ppb | PCB5-52M

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cashl\Acqudata\GC09\data\052710A.B\052F037.D
Injection Date: 28-MAY-2010 06:24
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.507 Min
Min
1.967
2.273
3.687
4.243
4.543
4.797
4.983
Aroclor 1232
5.530
Aroclor 1232
5.943
Aroclor 1232
Aroclor 1232
6.723
7.030
7.540
7.837
8.063
8.127
8.207
8.303
8.420
8.543
8.810
8.960
9.103
9.273
9.395
9.487
9.590
9.770
9.960
10.093
10.460
Aroclor 1262
10.733
10.870
Aroclor 1262
11.070
11.403
Aroclor 1262
11.853
11.977
Aroclor 1262
12.607
12.900
13.027
Aroclor 1262
13.977
14.320
14.570
14.943
16.210
17.440
Before
MAY 28 2010

Data File: \\Cashl\Acqudata\GC09\data\052710A.B\0527F037.D
Injection Date: 28-MAY-2010 06:24
Instrument: GC09.i
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F038.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F038.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R038.D
Inj Date : 28-MAY-2010 06:50
Sample Info: 1232/1262 @ 2000ppb | PCB5-52N
Misc Info :
Cal Date : 28-MAY-2010 13:16
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

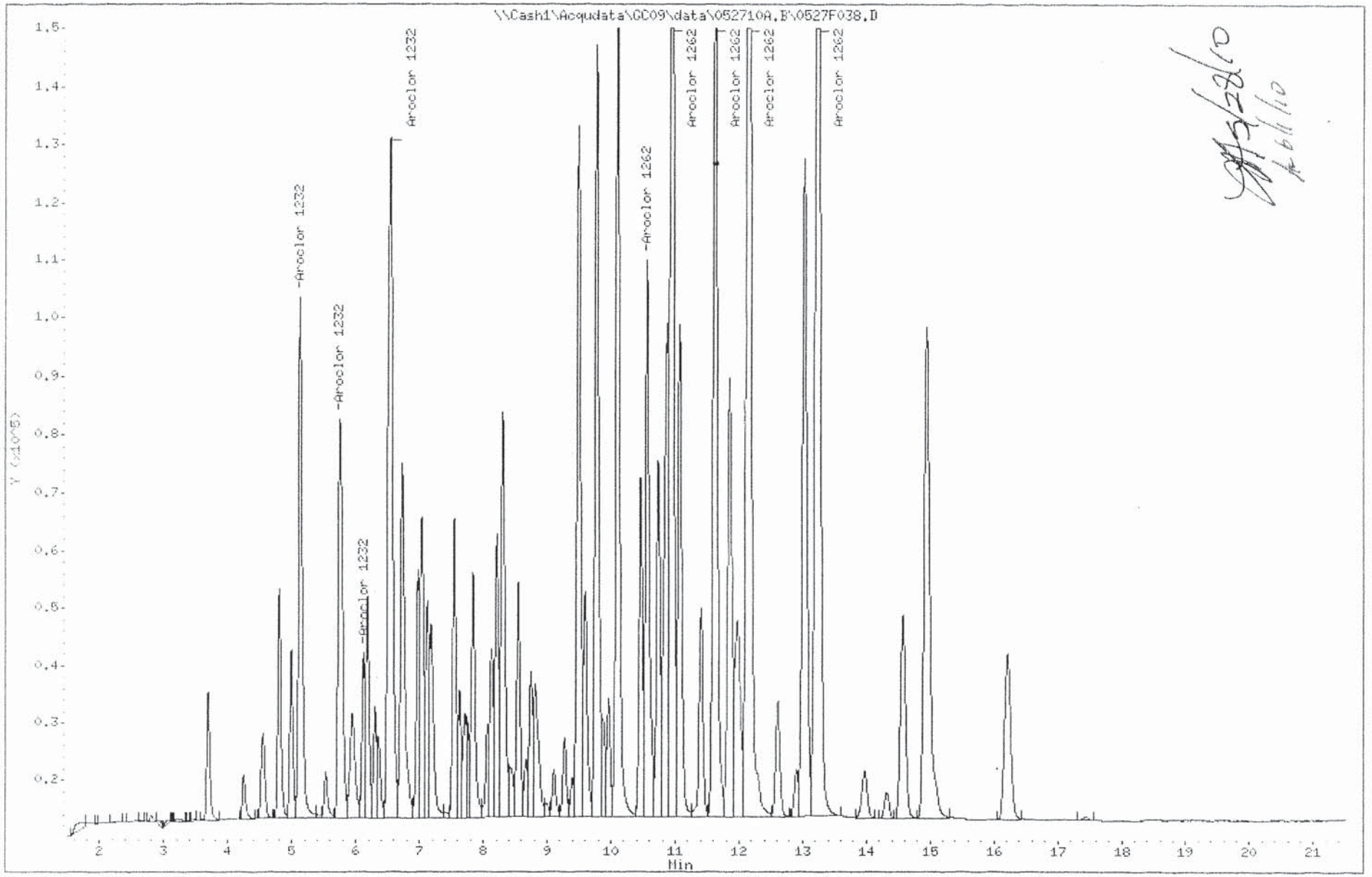
Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1232+1262.sub
Sub List #2 : 1232+1262.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.563	360855	314112	1650	1780	80.00- 120.00	100.00
	5.743	6.230	343131	131899	1700	2020	73.60- 110.40	95.09
	6.120	7.220	86461	253535	1920	1880	17.45- 26.18	23.96
	6.540	7.447	622650	160958	1740	2110	131.13- 196.70	172.55
			Average of Peak Amounts =		1750	1950		
Aroclor 1262	10.557	10.893	420338	446871	1720	1880	80.00- 120.00	100.00
	10.943	11.547	859523	410604	1740	1950	166.27- 249.41	204.48
	11.620	11.743	702866	623192	1810	1830	129.39- 194.09	167.21
	12.143	12.310	1360891	502806	1950	1950	249.74- 374.61	323.76
	13.230	13.143	1031973	1060977	1900	1920	188.68- 283.02	245.51
			Average of Peak Amounts =		1800	1910		

5/28/10
Analyst

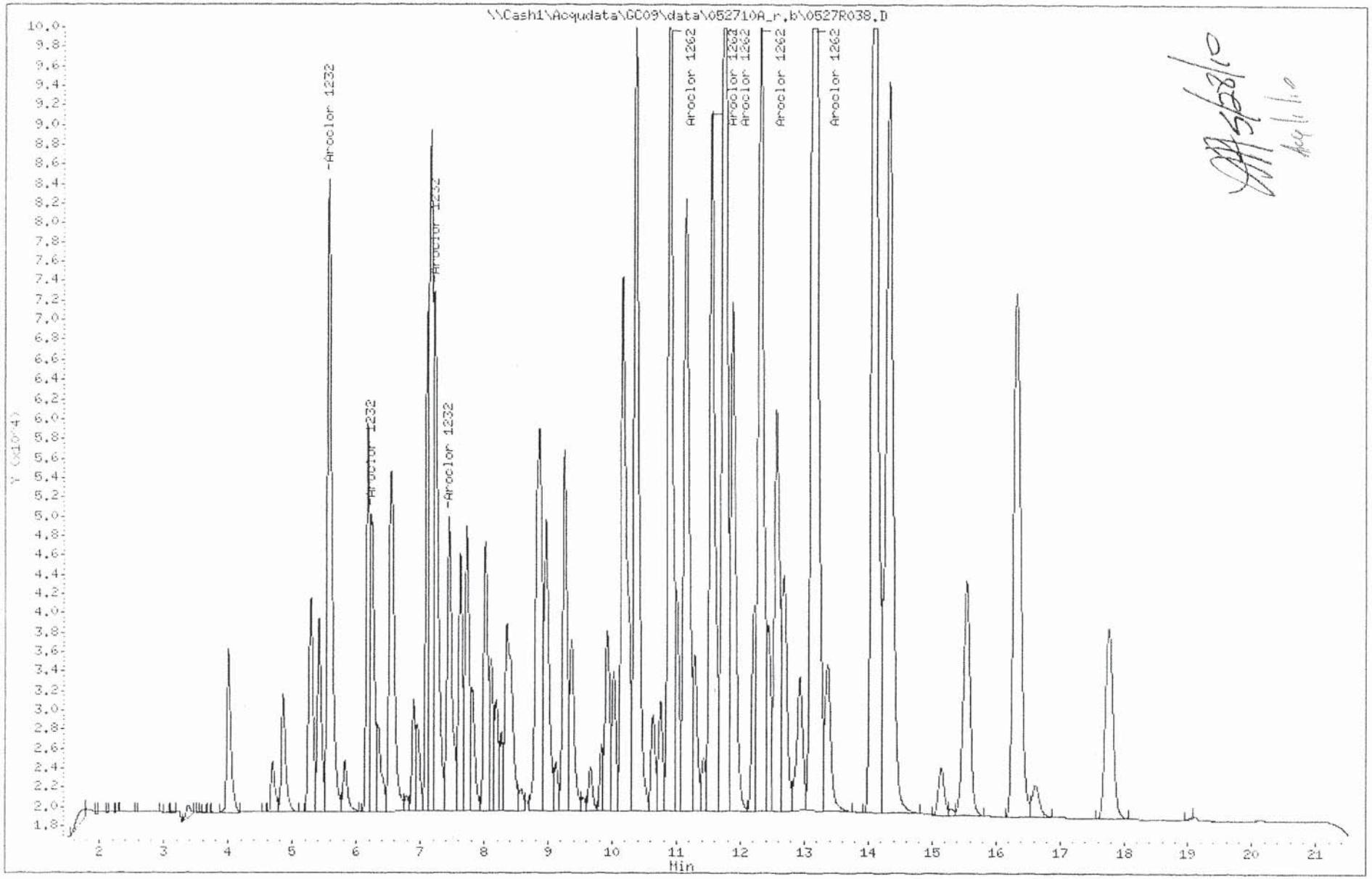
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F038.D
Date : 28-MAY-2010 06:50
Client ID:
Sample Info: 1232/1262 @ 2000ppb | PCB5-52N

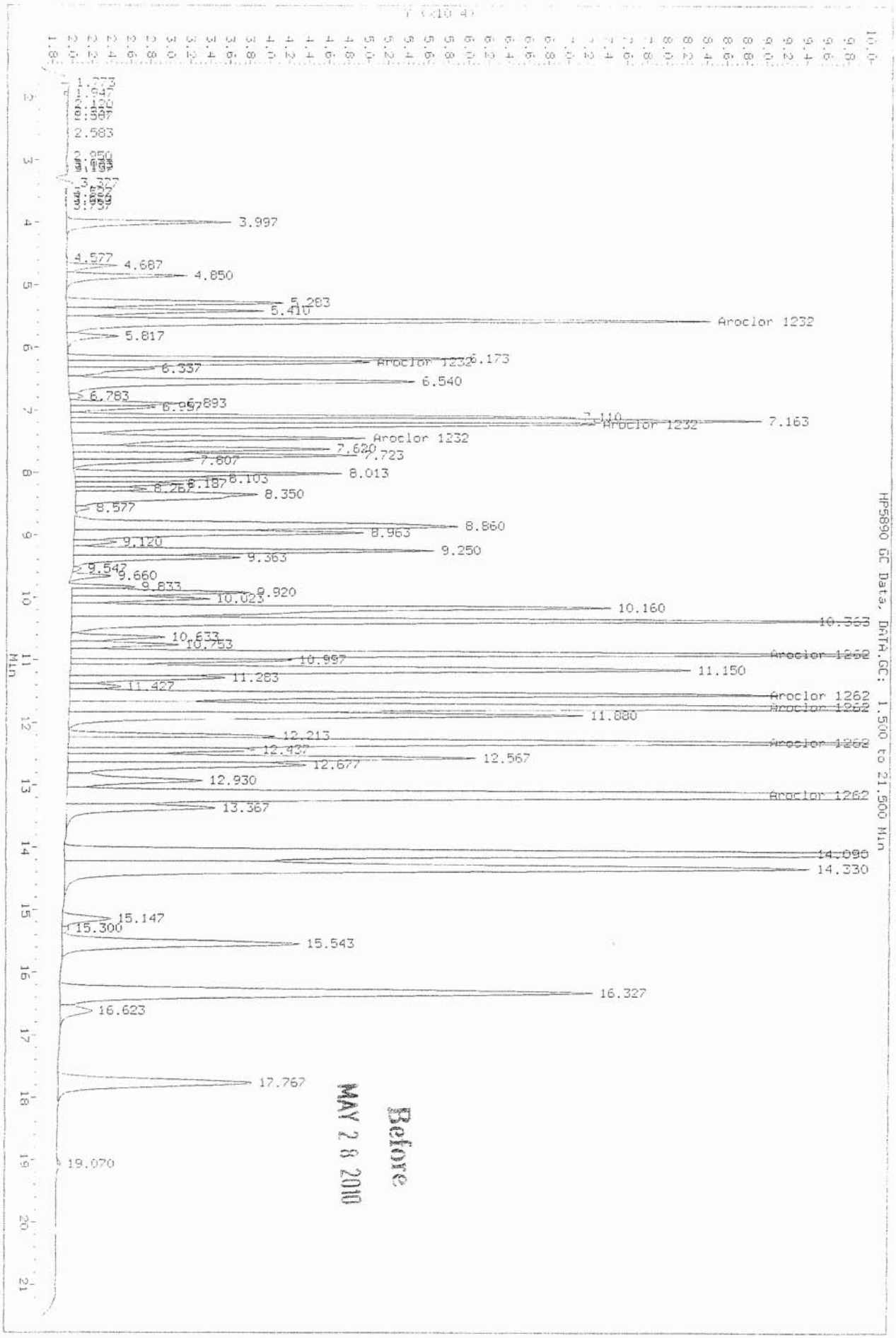
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

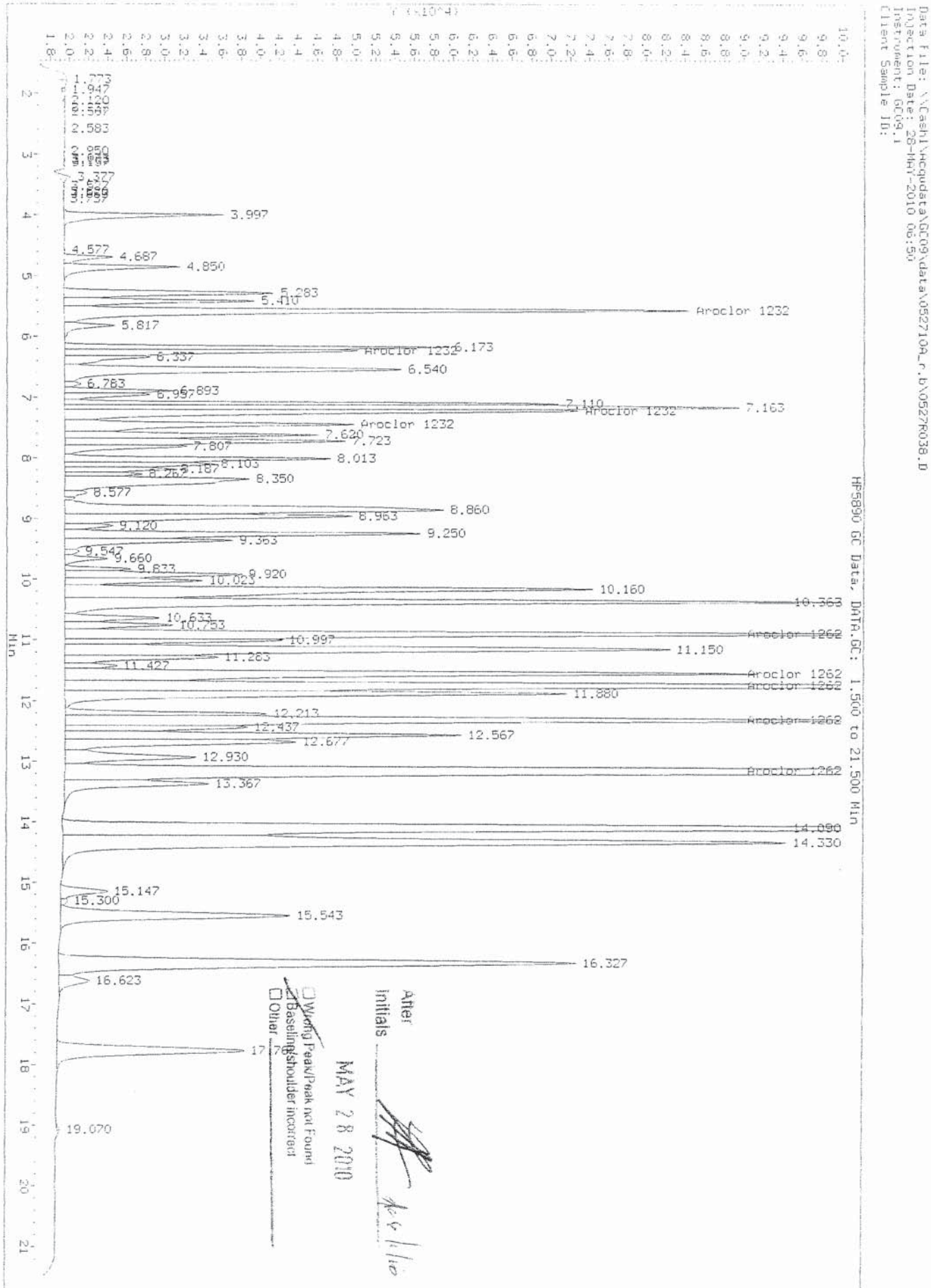


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R038.D
Date : 28-MAY-2010 06:50
Client ID:
Sample Info: 1232/1262 @ 2000ppb | PCB5-52N

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cash1\acqudata\GC09\data\052710A_r.b\0527R038.D
Injection Date: 28-MAY-2010 06:50
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.500 Min
Min
Aroclor 1232
Aroclor 1262
Before
MAY 2 8 2010


Data file: \\Cash\Hcqudata\GC09\data\052710A_r.b\0527R038.D
Injection Date: 28-MAY-2010 06:50
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.500 Min
Min
3.997
4.577
4.687
4.850
5.283
5.410
Aroclor 1232
5.817
Aroclor 1232
6.173
6.337
6.540
6.783
6.997
6.893
7.110
Aroclor 1232
7.163
Aroclor 1232
7.620
7.807
7.723
8.013
8.103
8.187
8.267
8.350
8.577
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9.120
9.250
9.363
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9.660
9.833
9.920
10.023
10.160
10.363
10.633
10.753
Aroclor 1262
10.997
11.150
11.263
11.427
Aroclor 1262
Aroclor 1262
11.880
12.213
Aroclor 1262
12.437
12.567
12.677
12.930
Aroclor 1262
13.367
14.090
14.330
15.147
15.300
15.543
16.327
16.623
17.780
19.070
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other
After initials
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F039.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F039.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R039.D
Inj Date : 28-MAY-2010 07:17
Sample Info: 1232/1262 @ 5000ppb | PCB5-520
Misc Info :
Cal Date : 28-MAY-2010 13:16
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

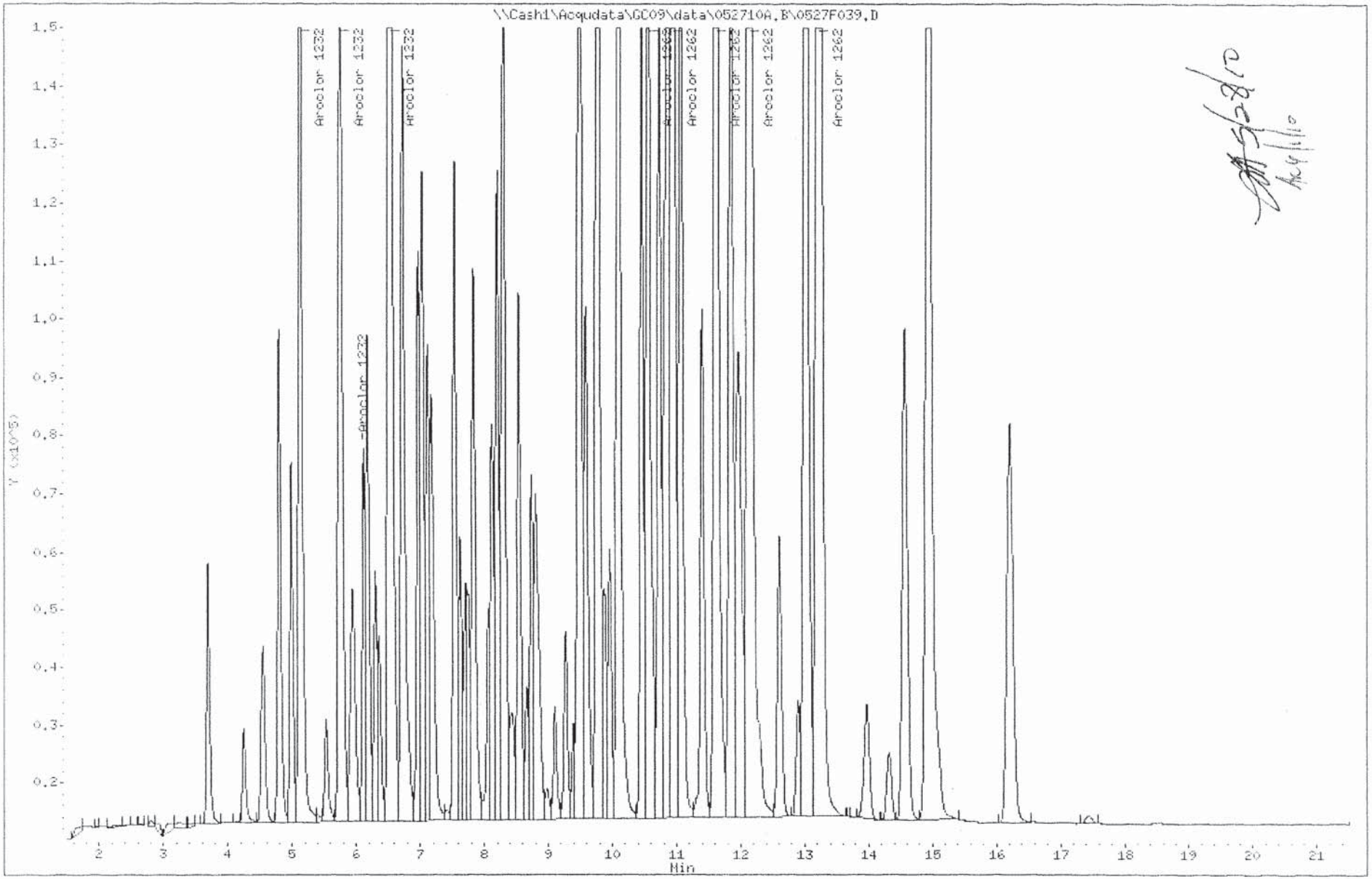
Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1232+1262.sub
Sub List #2 : 1232+1262.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.563	738377	675383	3380	3830	80.00- 120.00	100.00
	5.743	6.230	718670	290868	3560	4460	73.60- 110.40	97.33
	6.117	7.220	192260	553847	4260	4110	17.45- 26.18	26.04
	6.537	7.447	1333496	371430	3740	4880	131.13- 196.70	180.60
			Average of Peak Amounts =		3740	4320		
Aroclor 1262	10.557	10.890	945544	998865	3870	4210	80.00- 120.00	100.00
	10.940	11.547	1893966	959335	3820	4560	166.27- 249.41	200.30
	11.620	11.743	1579879	1411105	4070	4140	129.39- 194.09	167.09
	12.140	12.310	3070079	1178432	4180	4570	249.74- 374.61	324.69
	13.227	13.143	2348218	2476048	4320	4480	188.68- 283.02	248.35
			Average of Peak Amounts =		4050	4390		

5/28/10
AG 6/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F039,D
Date : 28-MAY-2010 07:17
Client ID:
Sample Info: 1232/1262 @ 5000ppb | PCB5-520

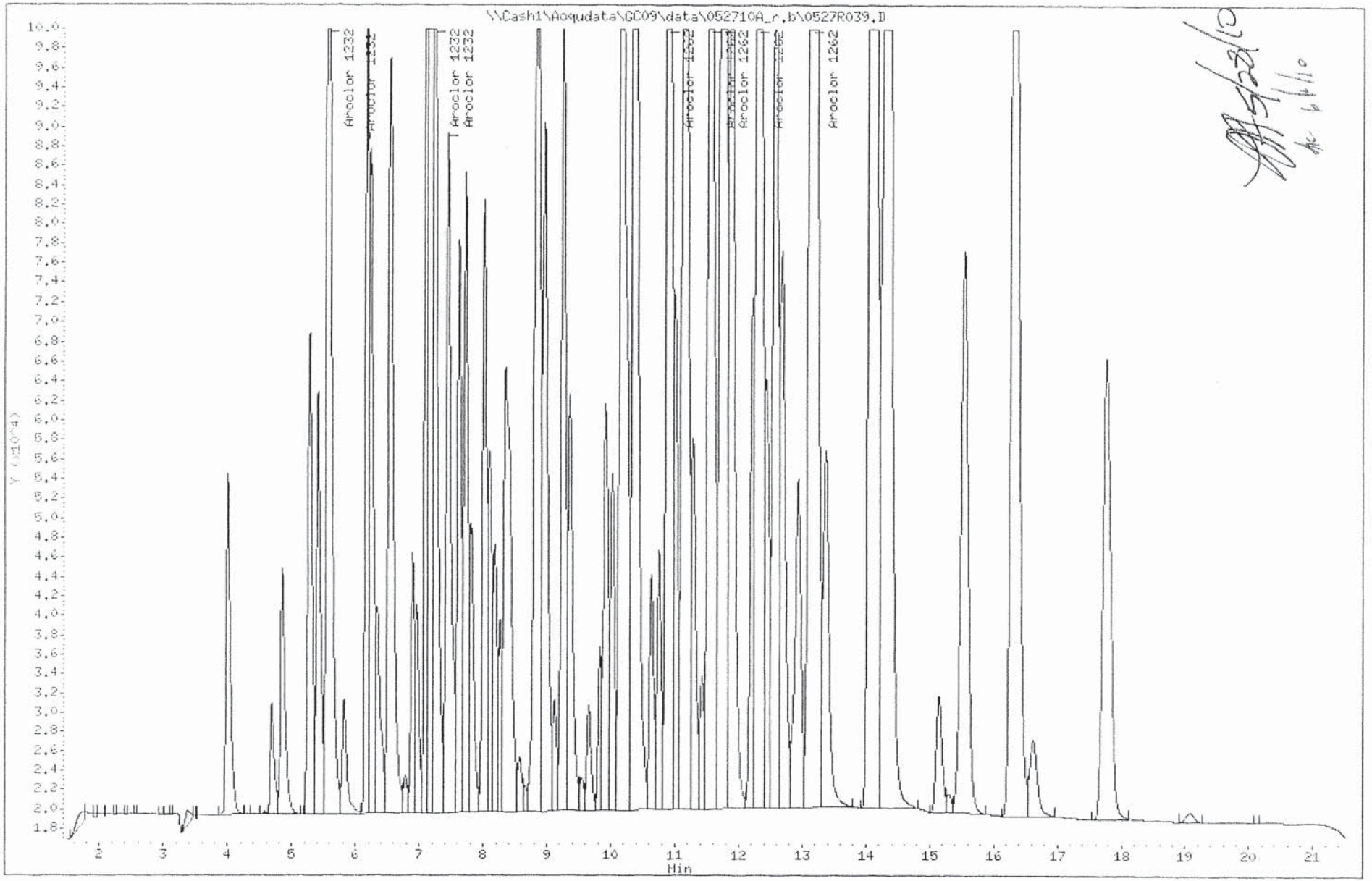
Instrument: GC09,i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R039.D
Date : 28-MAY-2010 07:17
Client ID:
Sample Info: 1232/1262 @ 5000ppb | PCB5-520

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F040.D
Report Date: 28-May-2010 17:29

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F040.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R040.D
Inj Date : 28-MAY-2010 07:43
Sample Info: 1242/1268 @ 25ppb | PCB5-52P
Misc Info :
Cal Date : 28-MAY-2010 12:05
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1242+1268.sub
Sub List #2 : 1242+1268.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.117	6.230	2189	2592	28.6	20.7	80.00- 120.00	100.00(M)
	6.540	6.540	19058	4258	30.9	26.4	727.90-1091.85	870.63(M)
	7.117	7.110	3946	3951	28.8	28.9	150.61- 225.91	180.26(M)
	7.540	7.623	5935	2581	30.0	23.0	230.50- 345.75	271.13(M)
	7.837	7.727	4985	2850	26.7	22.1	198.70- 298.04	227.73(M)
			Average of Peak Amounts =		29.0	24.2		
Aroclor 1268	13.027	14.340	30733	20882	30.3	27.2	80.00- 120.00	100.00(M)
	13.213	15.147	27378	16722	29.3	27.5	70.89- 106.33	89.08(M)
	13.977	15.480	22388	4112	29.7	25.5	58.04- 87.06	72.85(M)
	14.947	16.330	7889	6294	25.5	25.2	20.96- 31.43	25.67(M)
	16.213	17.770	66814	53482	31.2	29.4	172.26- 258.39	217.40(M)
			Average of Peak Amounts =		29.2	27.0		

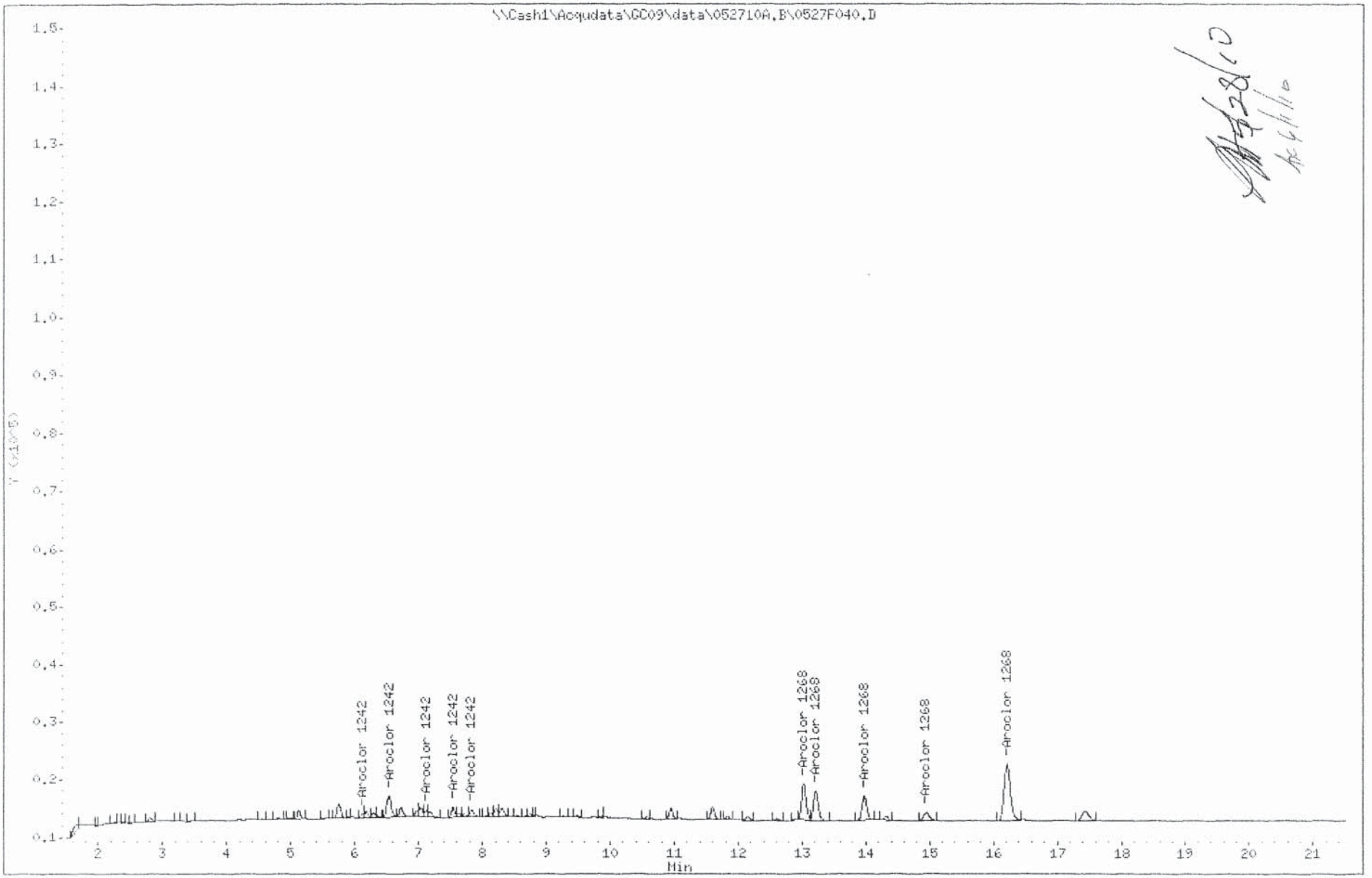
QC Flag Legend

M - Compound response manually integrated.

5/28/10
6/1/10

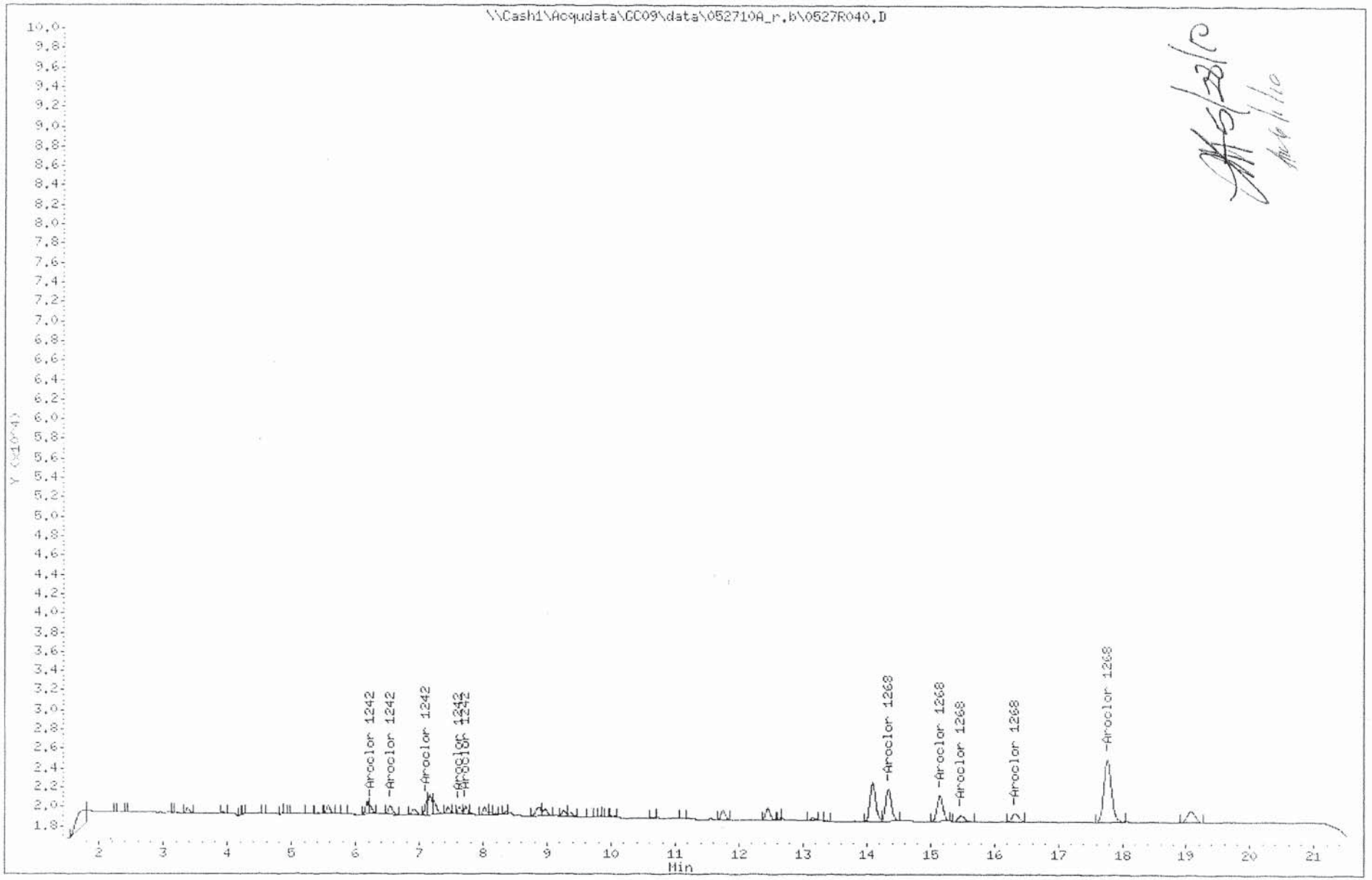
Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F040.D
Date : 28-MAY-2010 07:43
Client ID:
Sample Info: 1242/1268 @ 25ppb | PCB5-52P

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

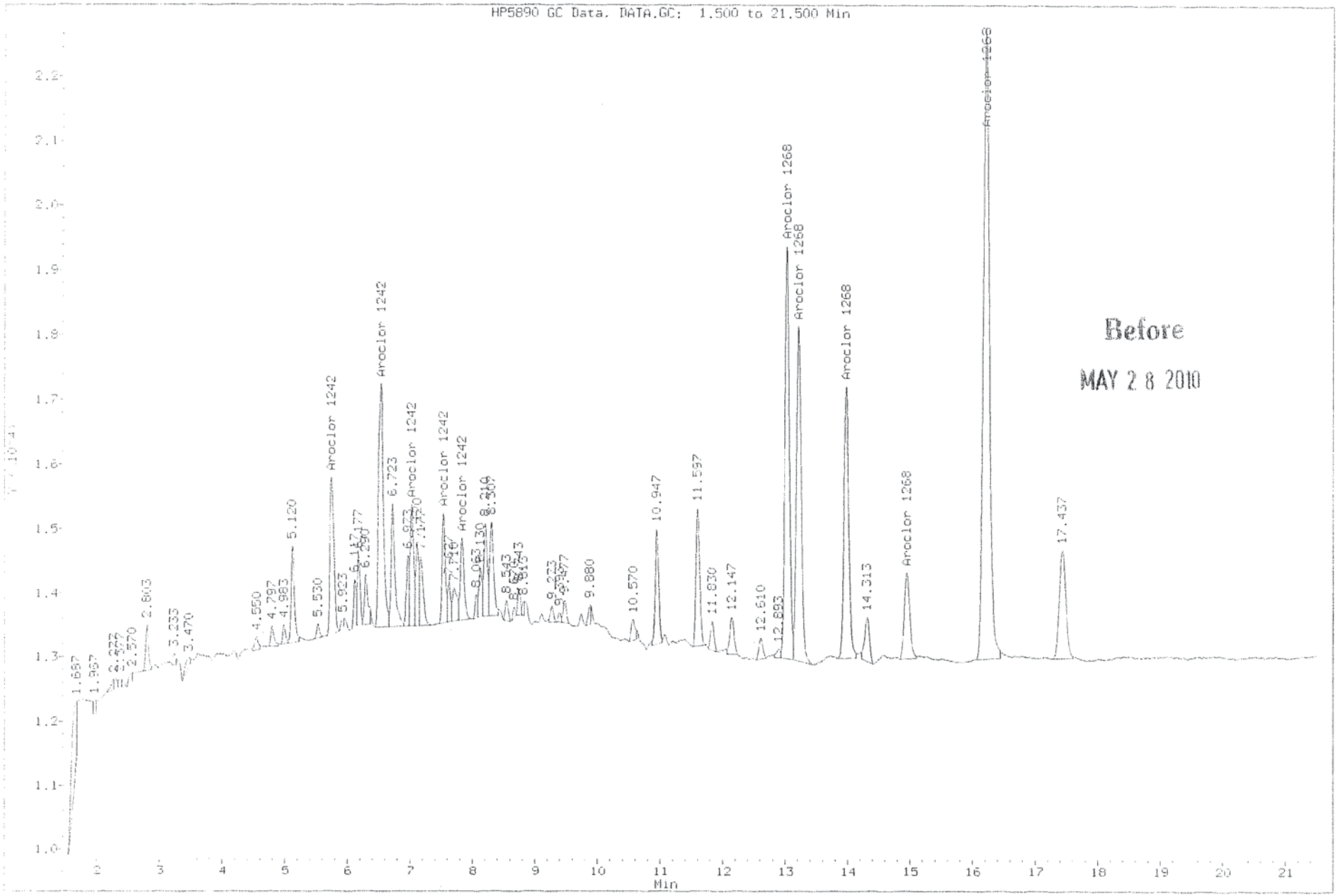


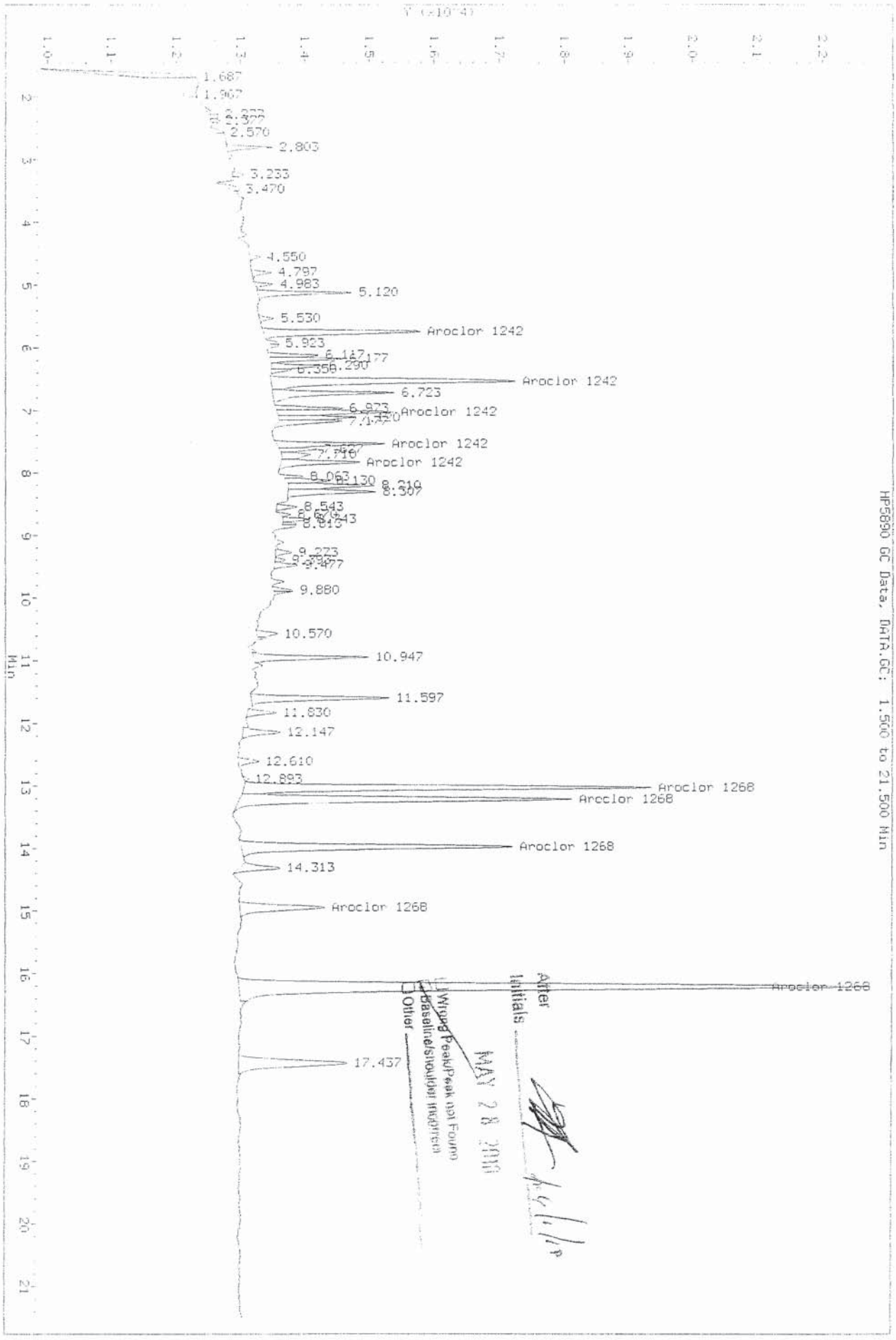
Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R040.D
Date : 28-MAY-2010 07:43
Client ID:
Sample Info: 1242/1268 @ 25ppb I PCB5-52P

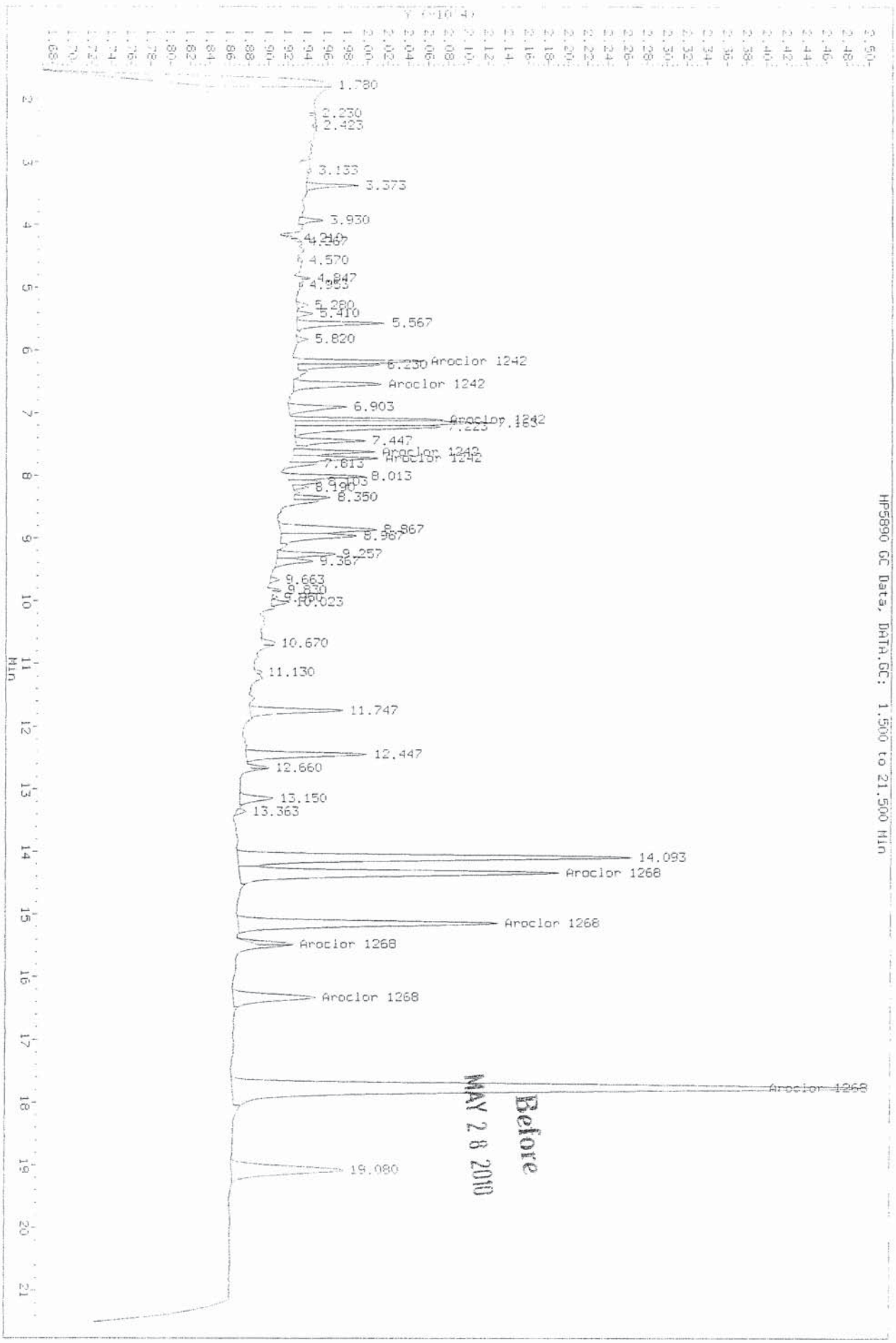
Instrument: GC09,i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



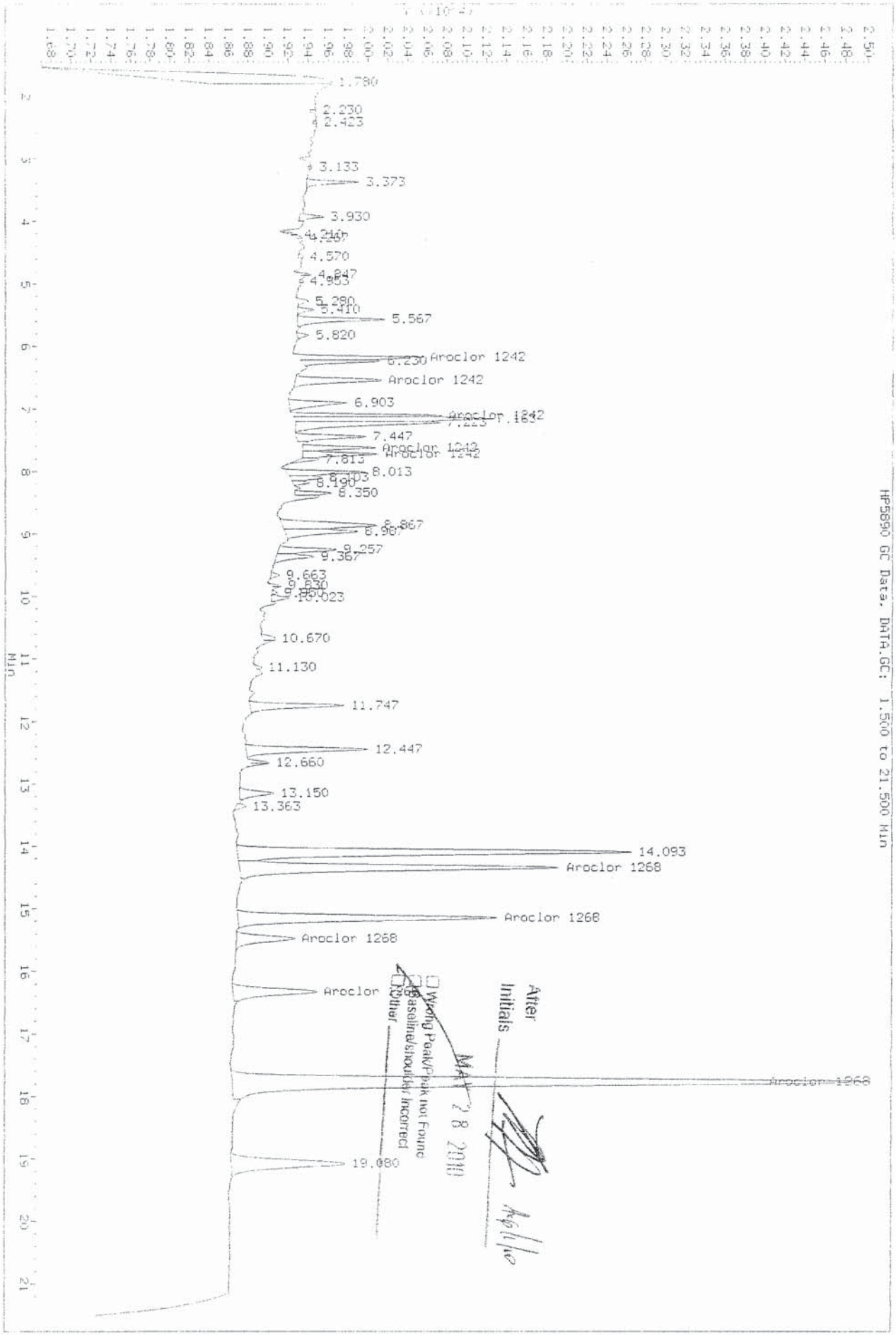
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F040.D
Injection Date: 28-MAY-2010 07:43
Instrument: GC09.1
Client Sample ID:




Data File: \\Cashl\Acqudata\GC09\data\0527100.B\0527F040.D
Injection Date: 28-MAY-2010 07:43
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.500 Min
Min
1.687
1.967
2.570
2.803
3.233
3.470
4.550
4.797
4.983
5.120
5.530
Aroclor 1242
5.923
6.723
Aroclor 1242
Aroclor 1242
Aroclor 1242
8.210
8.309
8.543
9.273
9.880
10.570
10.947
11.597
11.830
12.147
12.610
12.893
Aroclor 1268
Aroclor 1268
Aroclor 1268
14.313
Aroclor 1268
Aroclor 1268
17.437
After
Initials
Wrong Peak/Peak not Found
Baseline/shoulder integrated
Other
MAY 28 2010



Before
MAY 28 2010


Data File: \\Cashj\acqdata\GC09\data\052710A_r.b\0527R040.D
Injection Date: 28-MAY-2010 07:43
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.500 Min
1.780
2.230
2.423
3.133
3.373
3.930
4.570
5.290
5.410
5.567
5.820
6.230
Aroclor 1242
Aroclor 1242
6.903
Aroclor 1242
7.447
Aroclor 1242
Aroclor 1242
7.813
8.013
8.350
8.867
8.967
9.257
9.367
9.663
10.023
10.670
11.130
11.747
12.447
12.660
13.150
13.363
14.093
Aroclor 1268
Aroclor 1268
Aroclor 1268
Aroclor
Aroclor 1268
19.080
Min
After
Initials
Wrong Peak/Peak not Found
Baseline/Shoulder Incorrect
Other
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F041.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F041.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R041.D
Inj Date : 28-MAY-2010 08:09
Sample Info: 1242/1268 @ 50ppb | PCB5-52Q
Misc Info :
Cal Date : 28-MAY-2010 13:56
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1242+1268.sub
Sub List #2 : 1242+1268.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.120	6.233	3866	5771	50.4	49.7	80.00- 120.00	100.00(M)
	6.540	6.543	35175	8272	57.6	51.3	727.90-1091.85	909.88(M)
	7.120	7.113	7278	7955	53.2	58.3	150.61- 225.91	188.26(M)
	7.540	7.623	11139	6126	56.2	54.7	230.50- 345.75	288.13(M)
	7.837	7.727	9602	6766	51.5	52.4	198.70- 298.04	248.37(M)
			Average of Peak Amounts =		53.8	53.3		
Aroclor 1268	13.030	14.343	57560	40523	56.8	52.8	80.00- 120.00	100.00
	13.217	15.150	51003	32054	54.6	52.8	70.89- 106.33	88.61
	13.980	15.480	41761	7540	55.4	46.8	58.04- 87.06	72.55
	14.950	16.330	15078	12275	48.8	49.1	20.96- 31.43	26.20
	16.213	17.773	123942	100149	58.0	55.1	172.26- 258.39	215.33
			Average of Peak Amounts =		54.7	51.3		

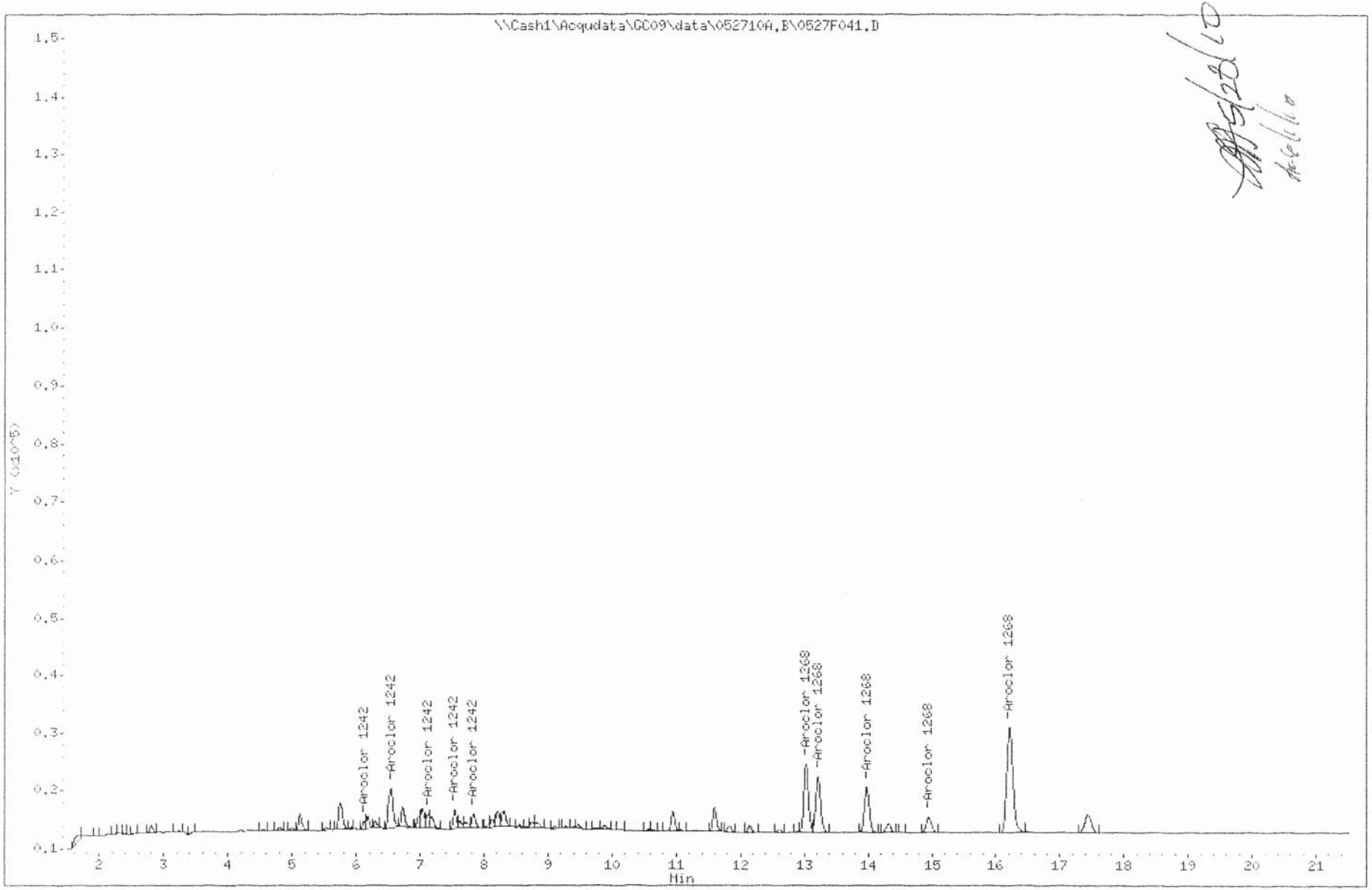
QC Flag Legend

M - Compound response manually integrated.

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F041.D
Date : 28-MAY-2010 08:09
Client ID:
Sample Info: 1242/1268 @ 50ppb | PCB5-52Q

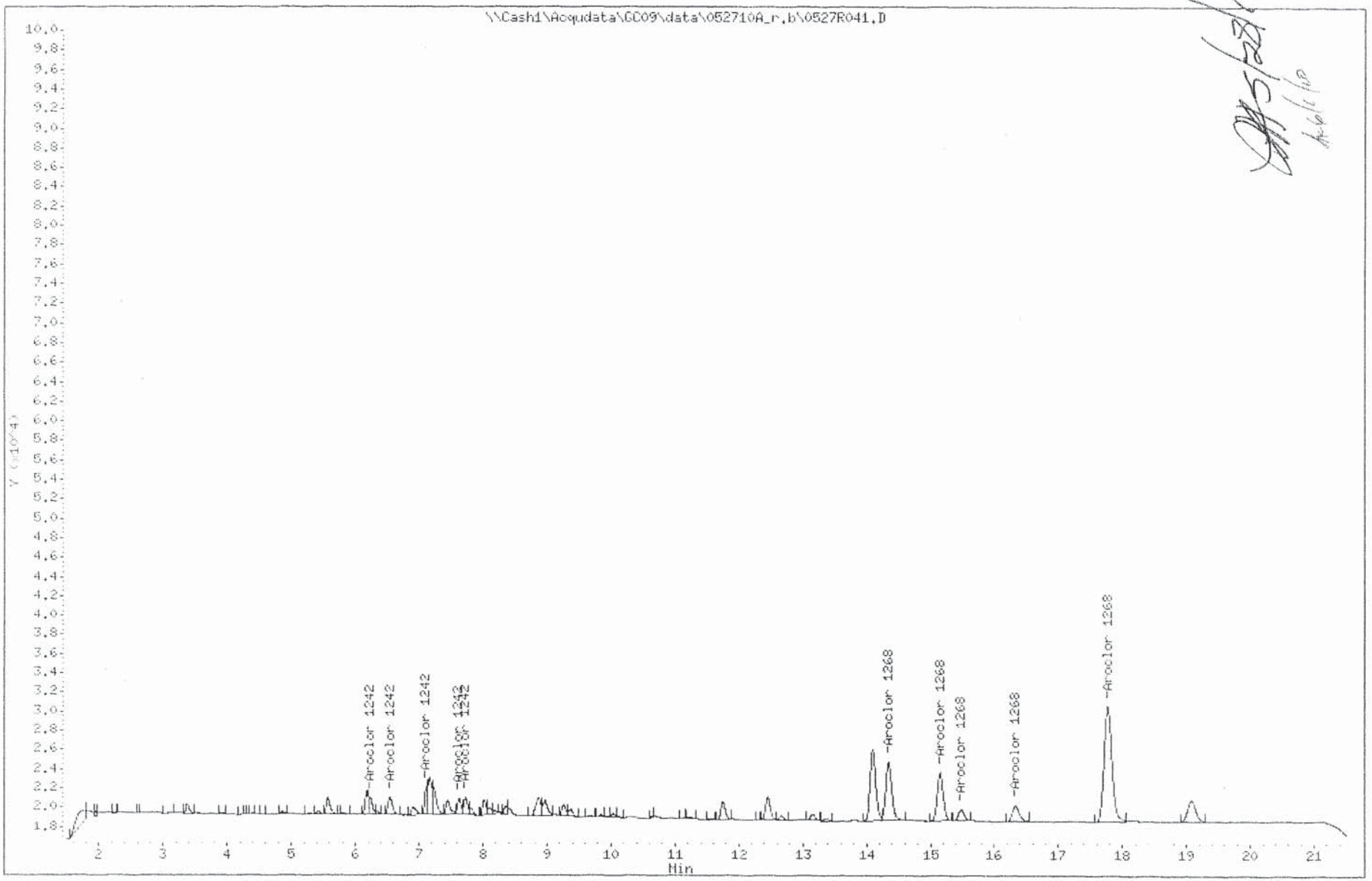
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



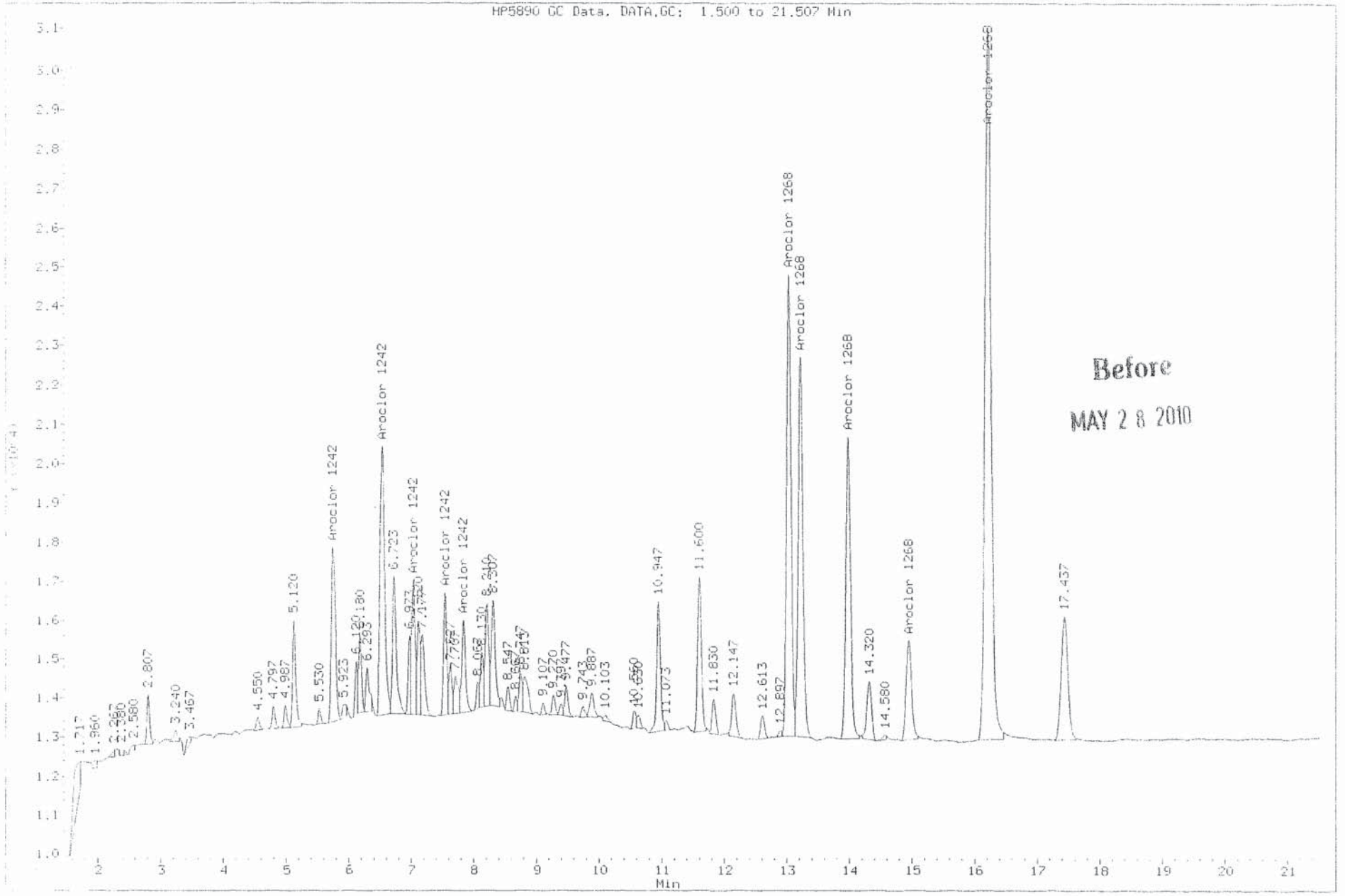
Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R041.D
Date : 28-MAY-2010 08:09
Client ID:
Sample Info: 1242/1268 @ 50ppb | PCB5-52Q

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



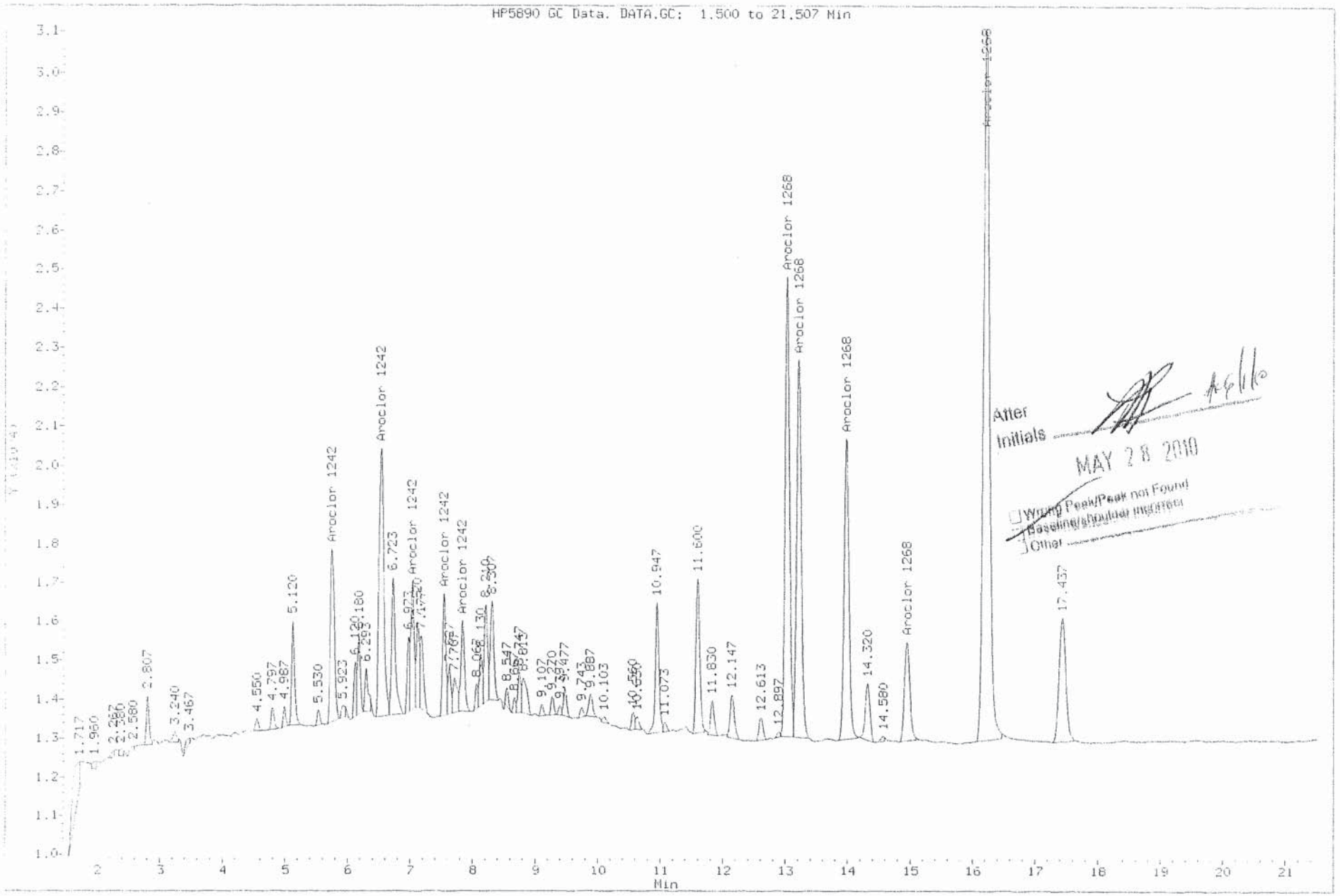
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F041.D
Injection Date: 28-MAY-2010 08:09
Instrument: GC09.i
Client Sample ID:

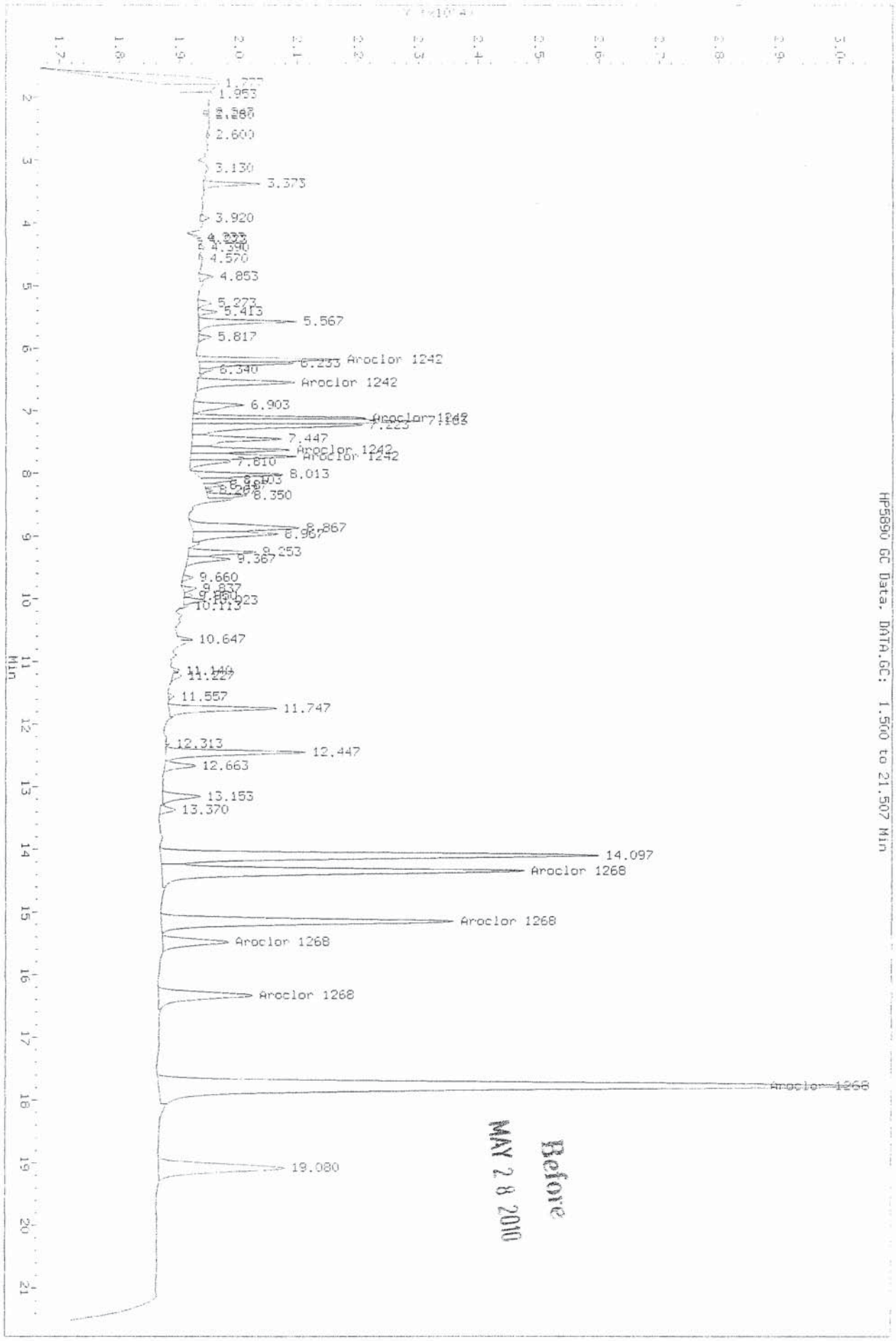


Before

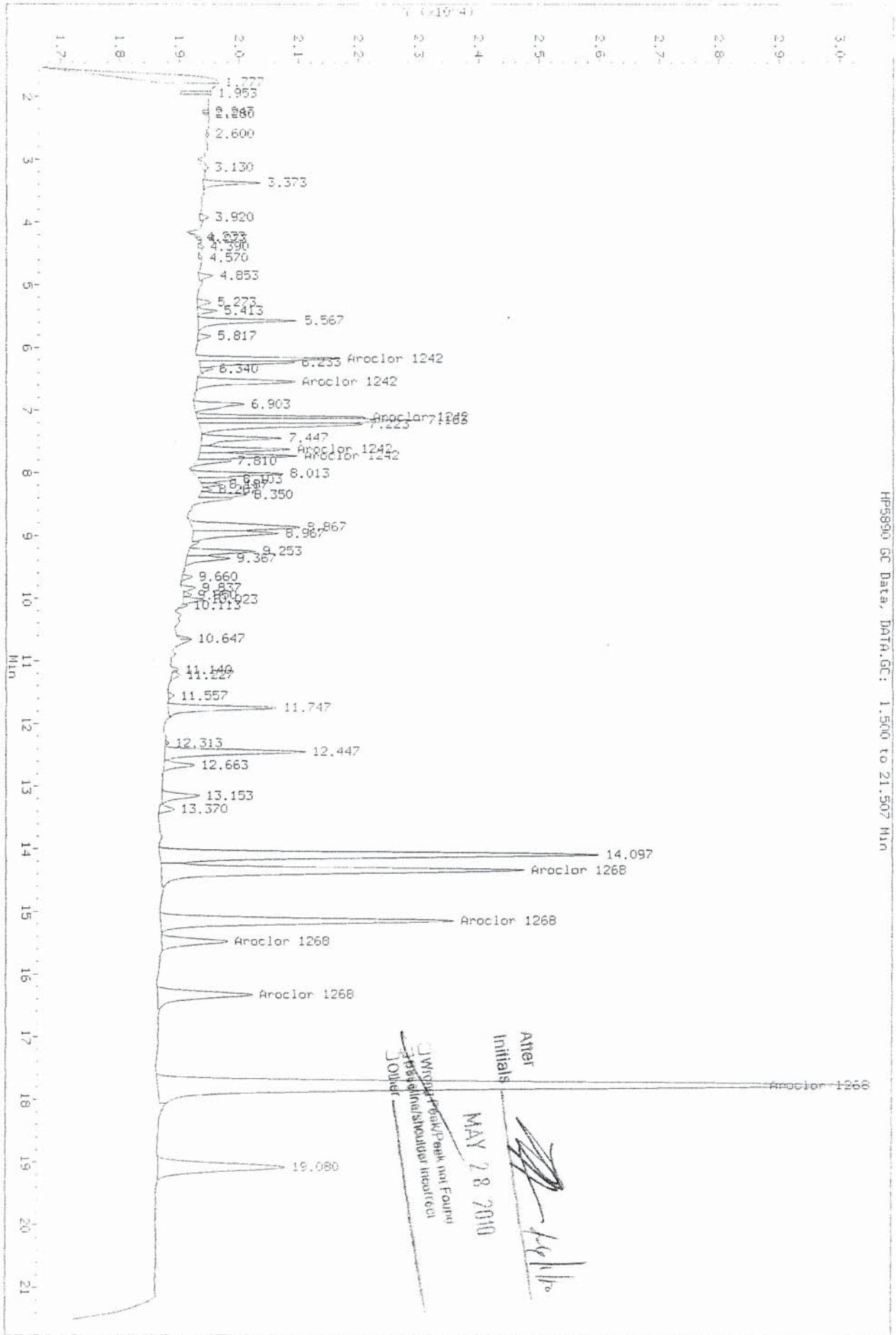
MAY 2 8 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F041.D
Injection Date: 28-MAY-2010 08:09
Instrument: GC09.i
Client Sample ID:





Before
MAY 2 8 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R041.D
Injection Date: 28-MAY-2010 08:09
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.507 Min
Min
1.777
1.953
2.600
3.130
3.373
3.920
4.570
4.853
5.567
5.817
Aroclor 1242
6.340
Aroclor 1242
6.903
7.447
Aroclor 1242
7.810
8.013
8.350
8.867
8.967
9.253
9.387
9.660
10.113
10.647
11.557
11.747
12.313
12.447
12.663
13.153
13.370
14.097
Aroclor 1268
Aroclor 1268
Aroclor 1268
Aroclor 1268
19.080
After
Initials
MAY 28 2010
Outlier

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F042.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F042.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R042.D
Inj Date : 28-MAY-2010 08:35
Sample Info: 1242/1268 @ 500ppb | PCB5-52R
Misc Info :
Cal Date : 28-MAY-2010 13:56
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1242+1268.sub
Sub List #2 : 1242+1268.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

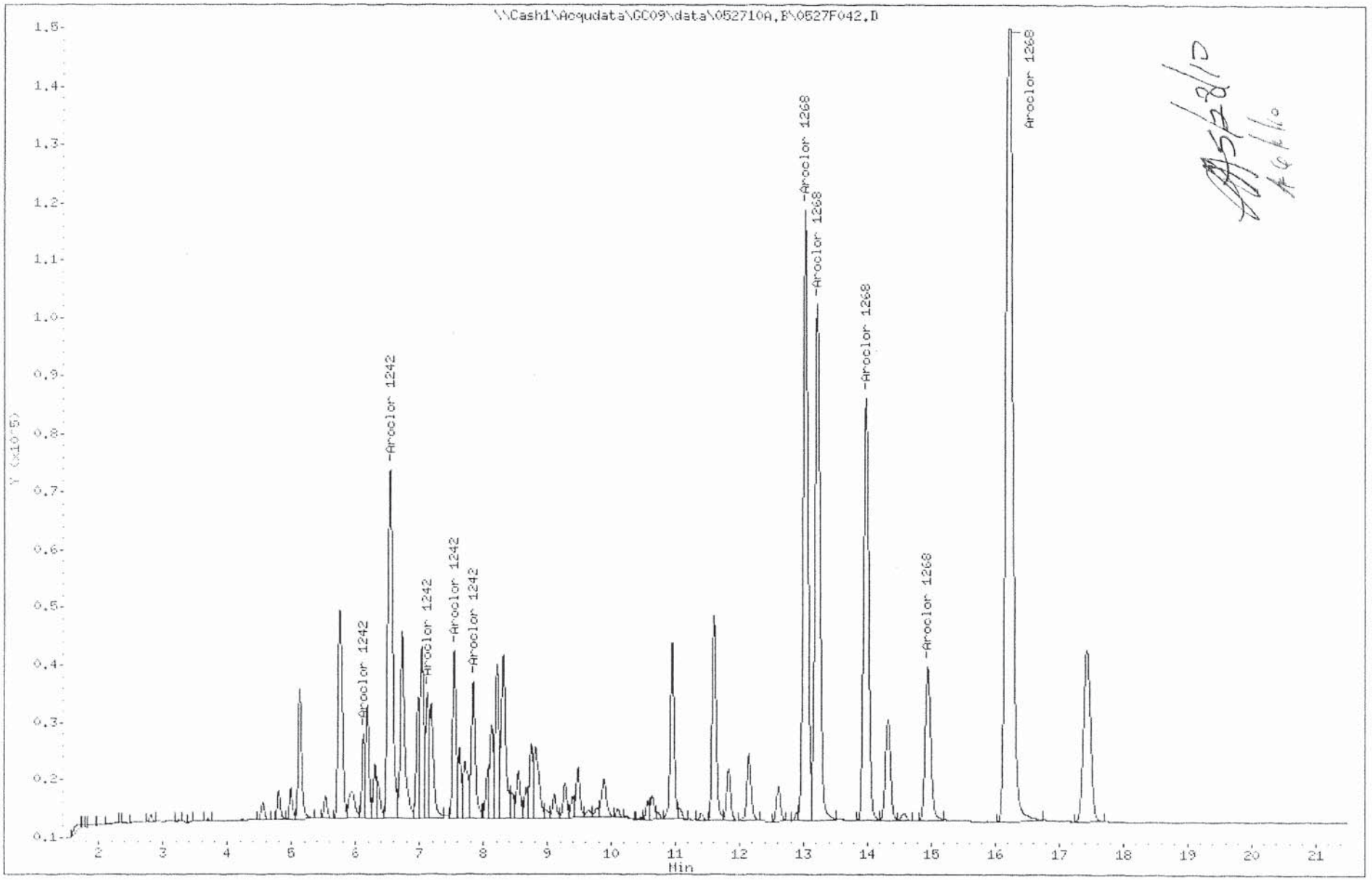
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.120	6.233	42663	60675	243	522	80.00- 120.00	100.00
	6.540	6.543	327749	88798	528	551	727.90-1091.85	768.22
	7.120	7.113	76765	70268	458	515	150.61- 225.91	179.93
	7.540	7.623	106545	65072	538	581	230.50- 345.75	249.74
	7.837	7.727	105831	75230	568	582	198.70- 298.04	248.06
			Average of Peak Amounts =		467	550		
Aroclor 1268	13.027	14.343	530418	411997	524	536	80.00- 120.00	100.00
	13.213	15.147	496919	321936	532	530	70.89- 106.33	93.68
	13.977	15.477	396320	86145	526	535	58.04- 87.06	74.72
	14.947	16.327	166509	110918	539	523	20.96- 31.43	31.39
	16.213	17.770	1118039	966060	523	532	172.26- 258.39	210.78
			Average of Peak Amounts =		529	531		

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F042.D
Date : 28-MAY-2010 08:35
Client ID:
Sample Info: 1242/1268 @ 500ppb | PCB5-52R

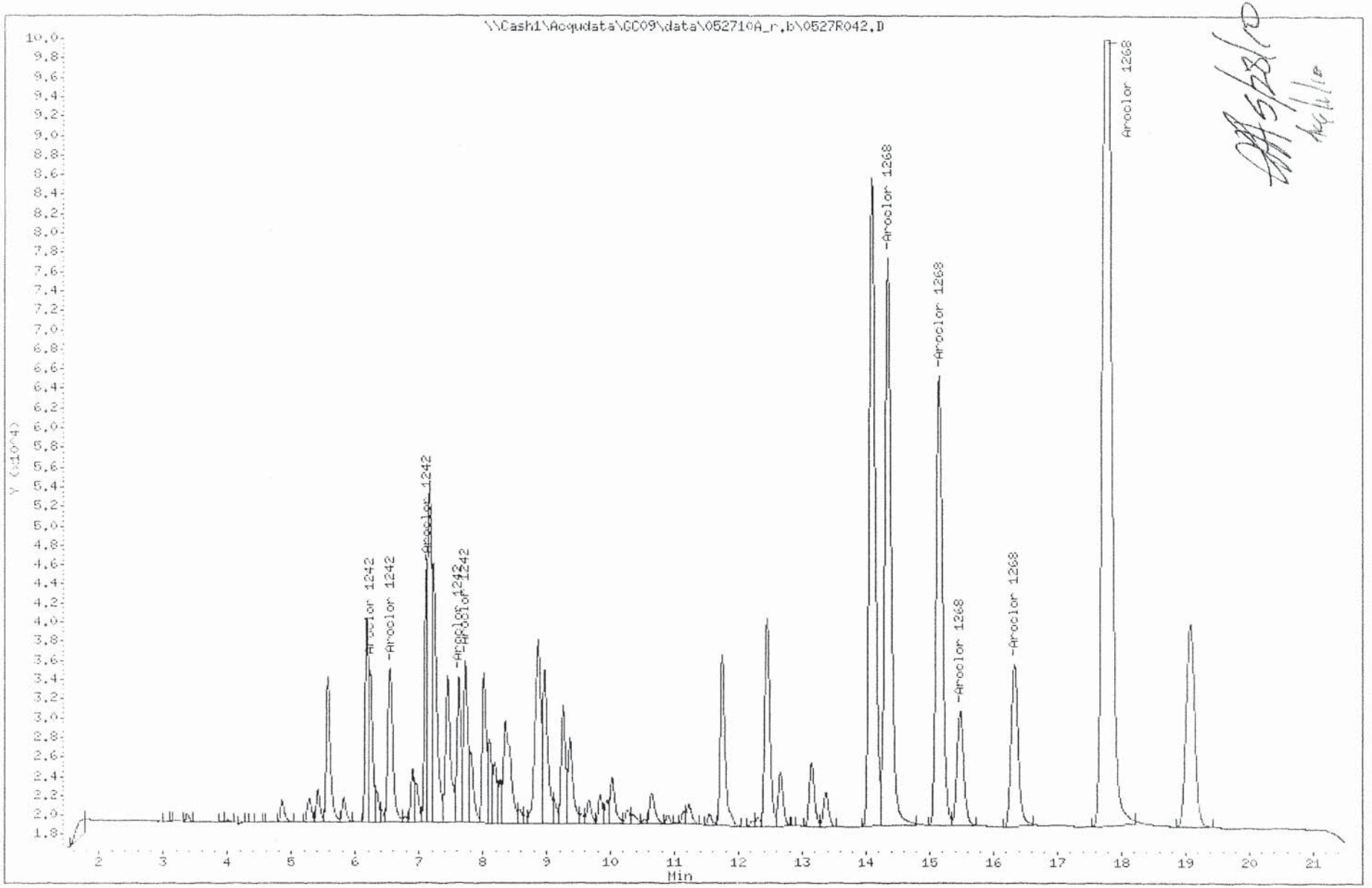
Instrument: GC09,i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0,53

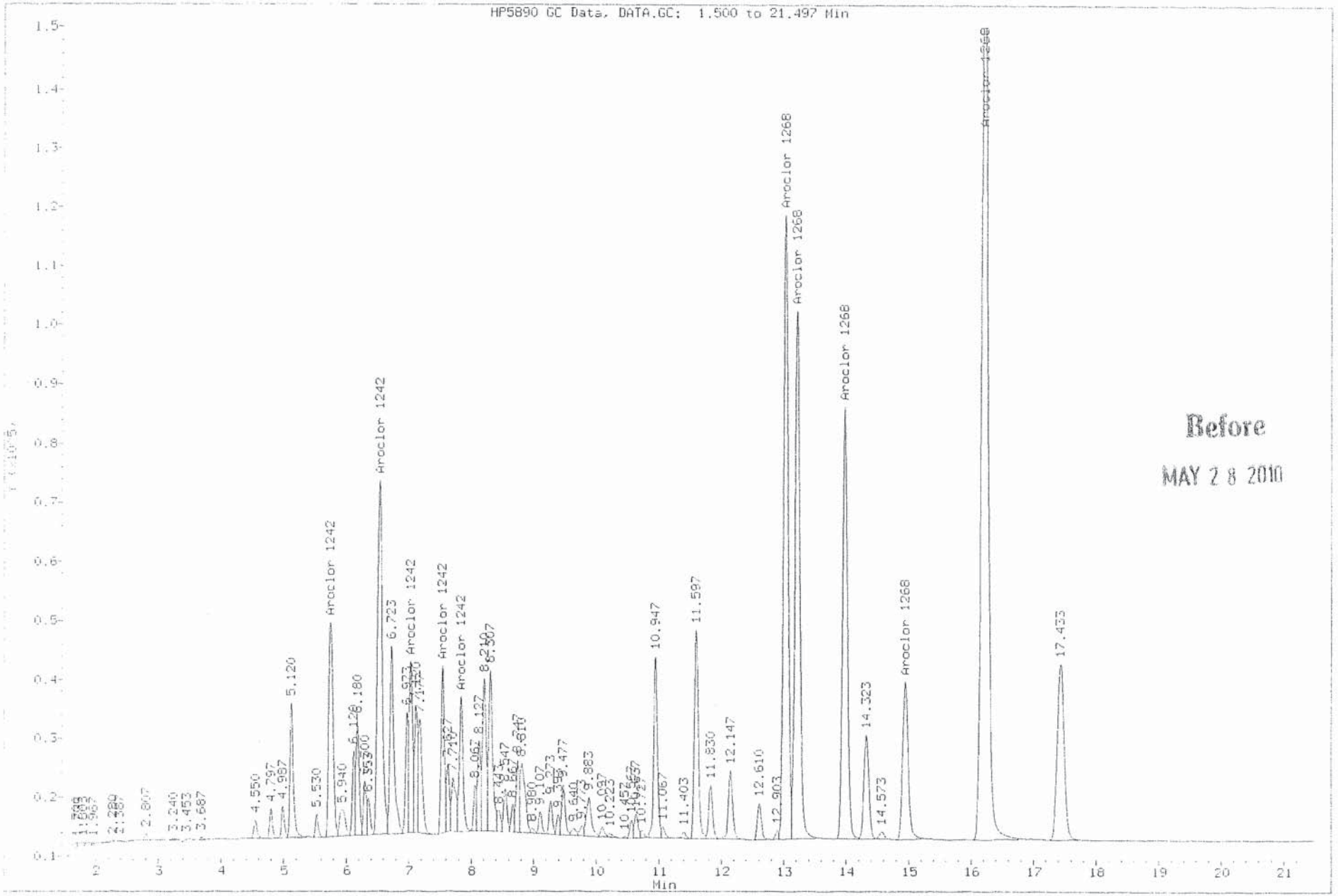


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R042.D
Date : 28-MAY-2010 08:35
Client ID:
Sample Info: 1242/1268 @ 500ppb | PCB5-52R

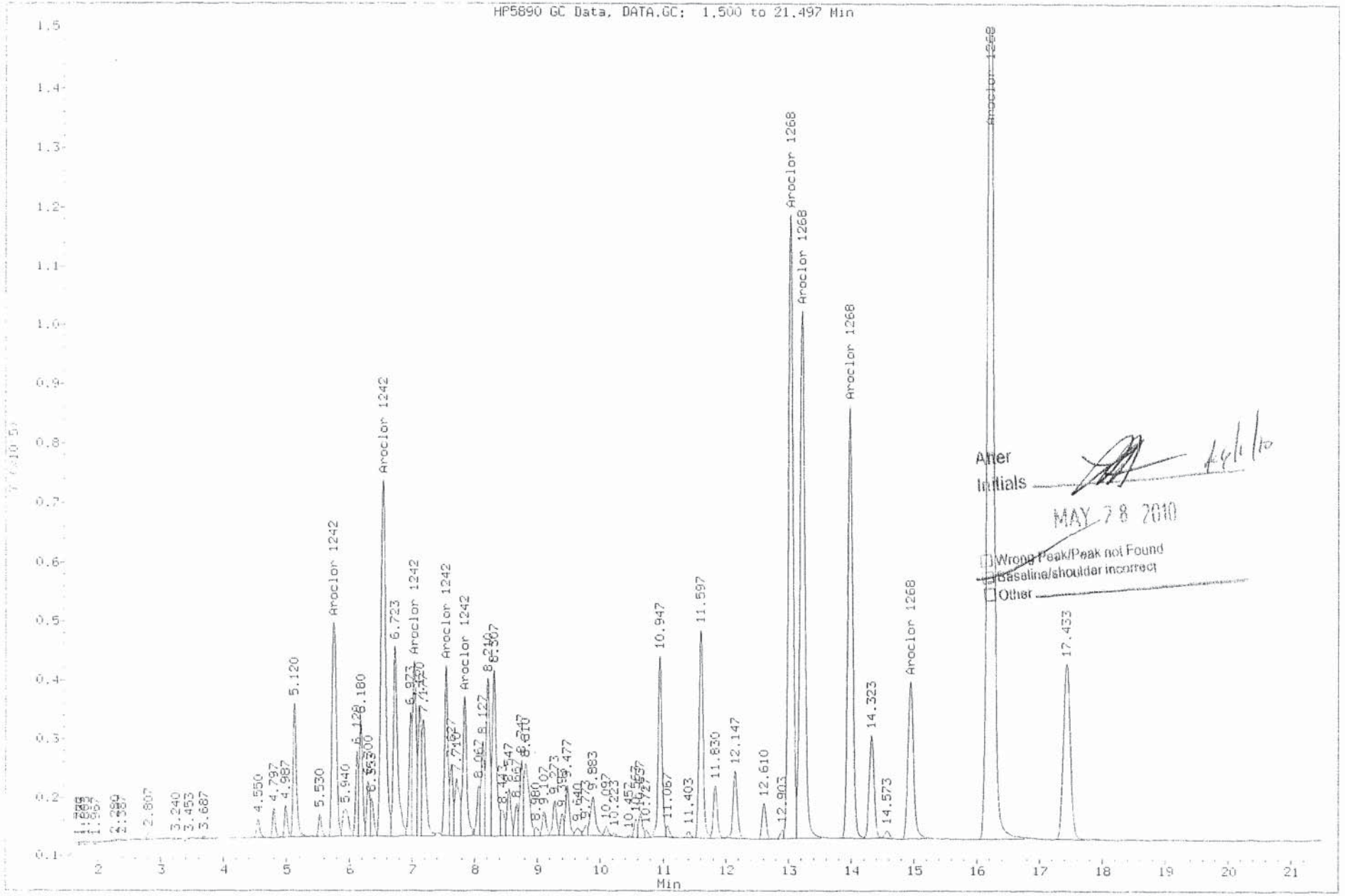
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F042.D
Injection Date: 28-MAY-2010 08:35
Instrument: GC09.I
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F042.D
Injection Date: 28-MAY-2010 08:35
Instrument: GC09.i
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F043.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F043.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R043.D
Inj Date : 28-MAY-2010 09:01
Sample Info: 1242/1268 @ 1000ppb | PCB5-52S
Misc Info :
Cal Date : 28-MAY-2010 13:56
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1242+1268.sub
Sub List #2 : 1242+1268.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.120	6.230	77031	114906	563	1020	80.00- 120.00	100.00
	6.540	6.543	583632	164054	940	1020	727.90-1091.85	757.66
	7.117	7.110	136697	134281	880	984	150.61- 225.91	177.46
	7.540	7.623	192418	118806	971	1060	230.50- 345.75	249.79
	7.837	7.727	191352	139765	1030	1080	198.70- 298.04	248.41
			Average of Peak Amounts =		877	1030		
Aroclor 1268	13.027	14.343	970655	772430	958	1000	80.00- 120.00	100.00
	13.213	15.150	913928	606755	979	999	70.89- 106.33	94.16
	13.977	15.480	730324	166699	969	1030	58.04- 87.06	75.24
	14.943	16.327	321406	256412	1040	1020	20.96- 31.43	33.11
	16.210	17.770	2018454	1782067	944	981	172.26- 258.39	207.95
			Average of Peak Amounts =		978	1010		

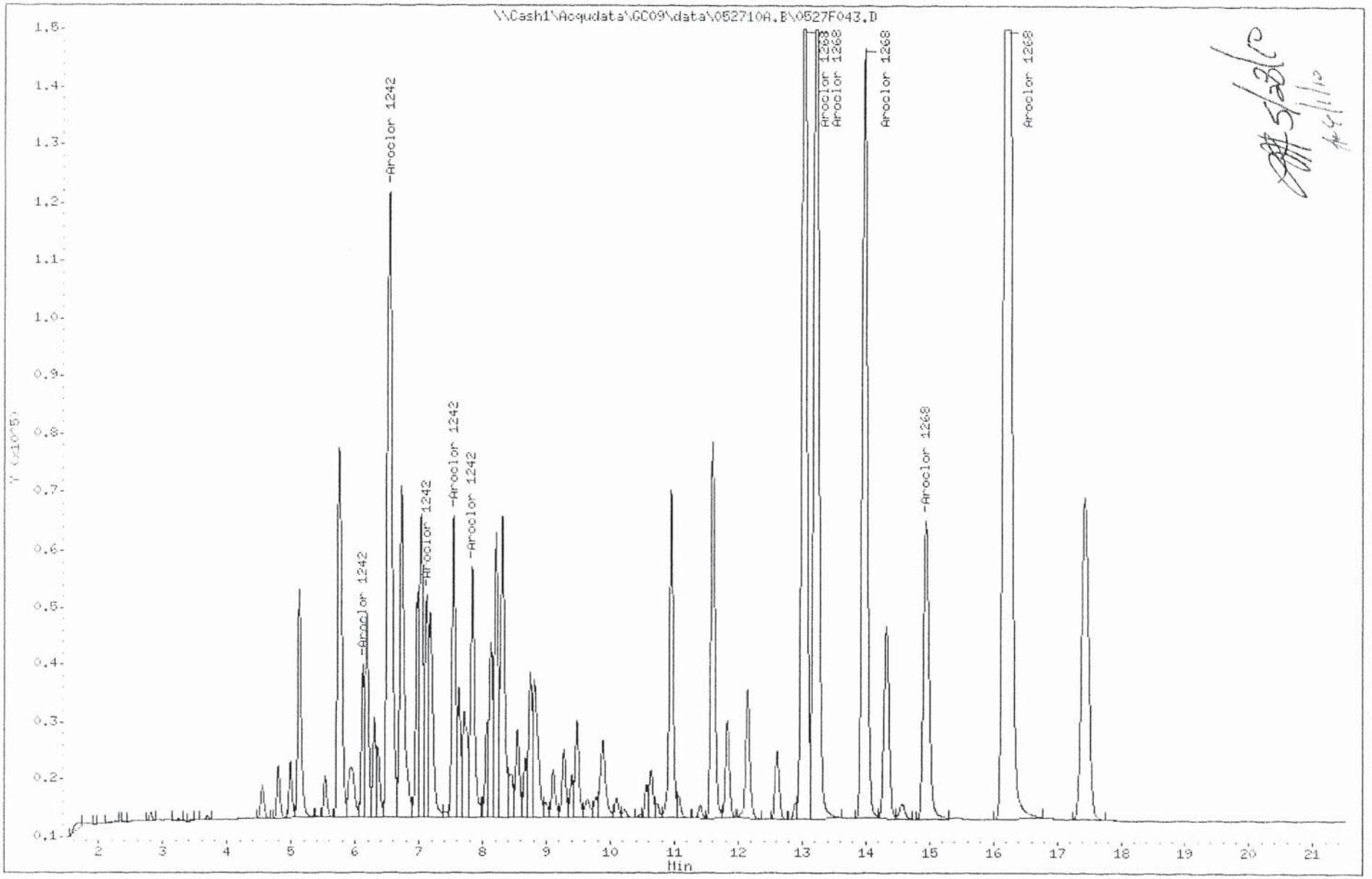
5/28/10
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F043.D
Date : 28-MAY-2010 09:01
Client ID:
Sample Info: 1242/1268 @ 1000ppb | PCB5-52S

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

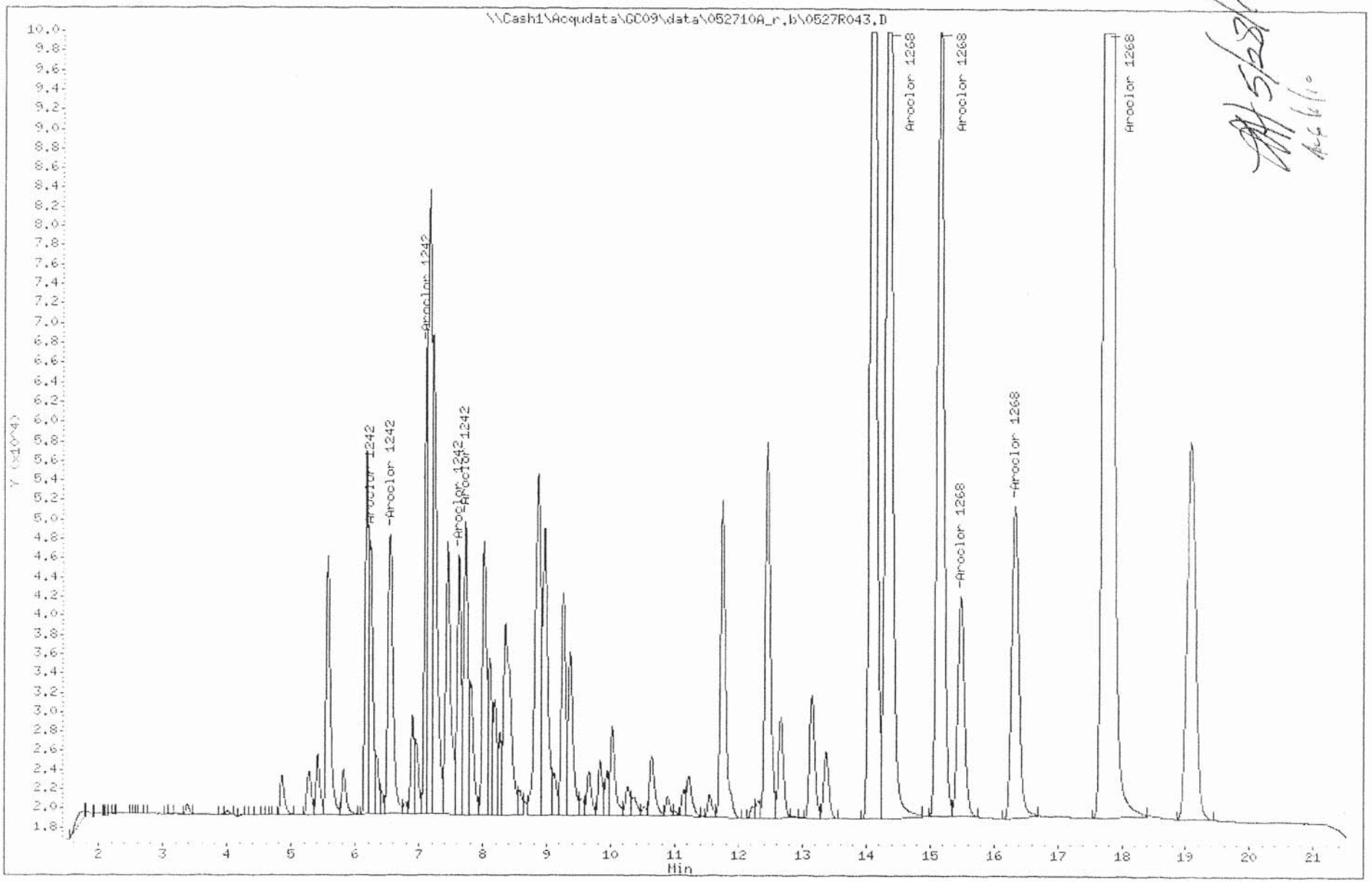
Column phase: DB-35MS



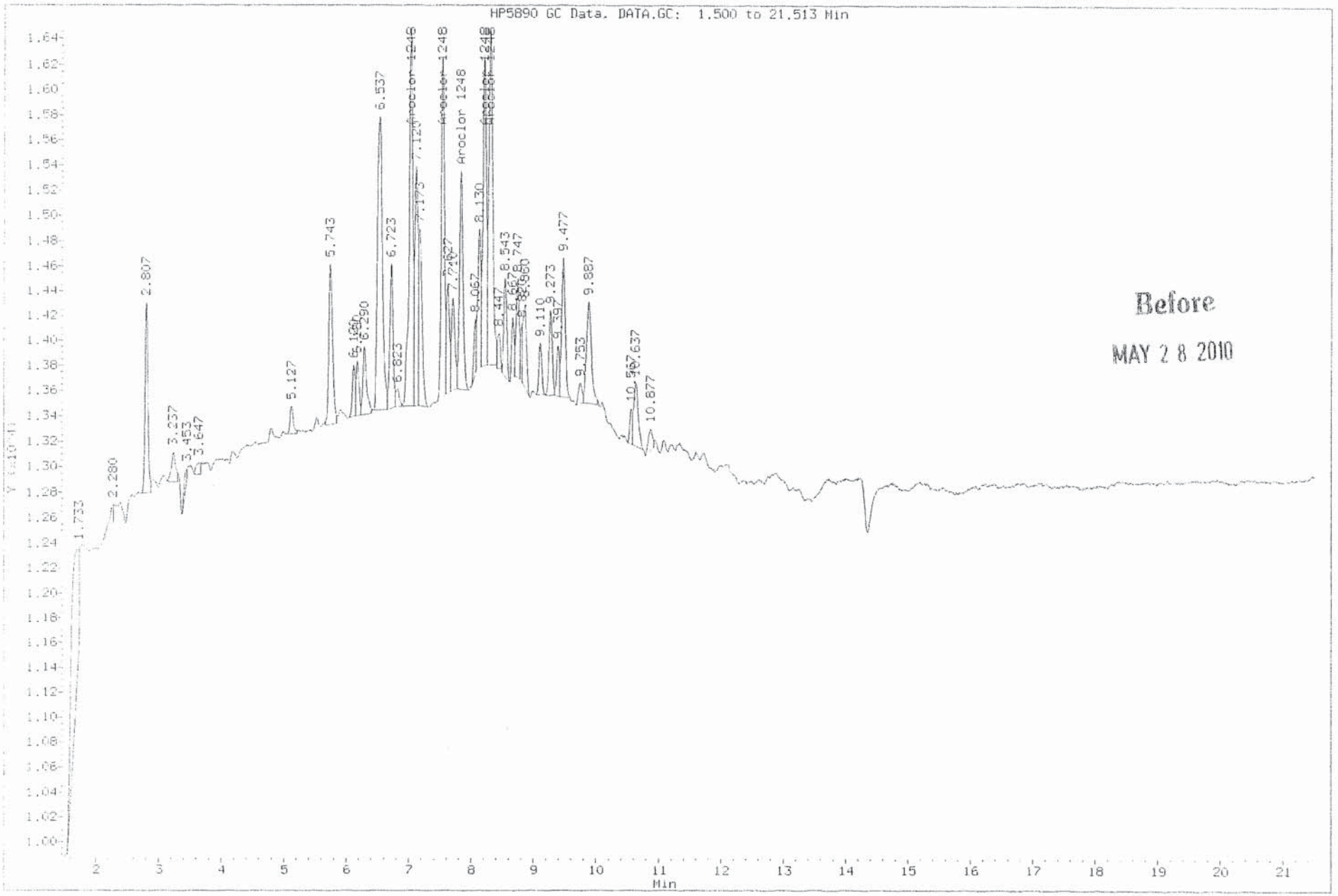
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R043.D
Date : 28-MAY-2010 09:01
Client ID:
Sample Info: 1242/1268 @ 1000ppb | PCB5-52S

Instrument: GC09.i

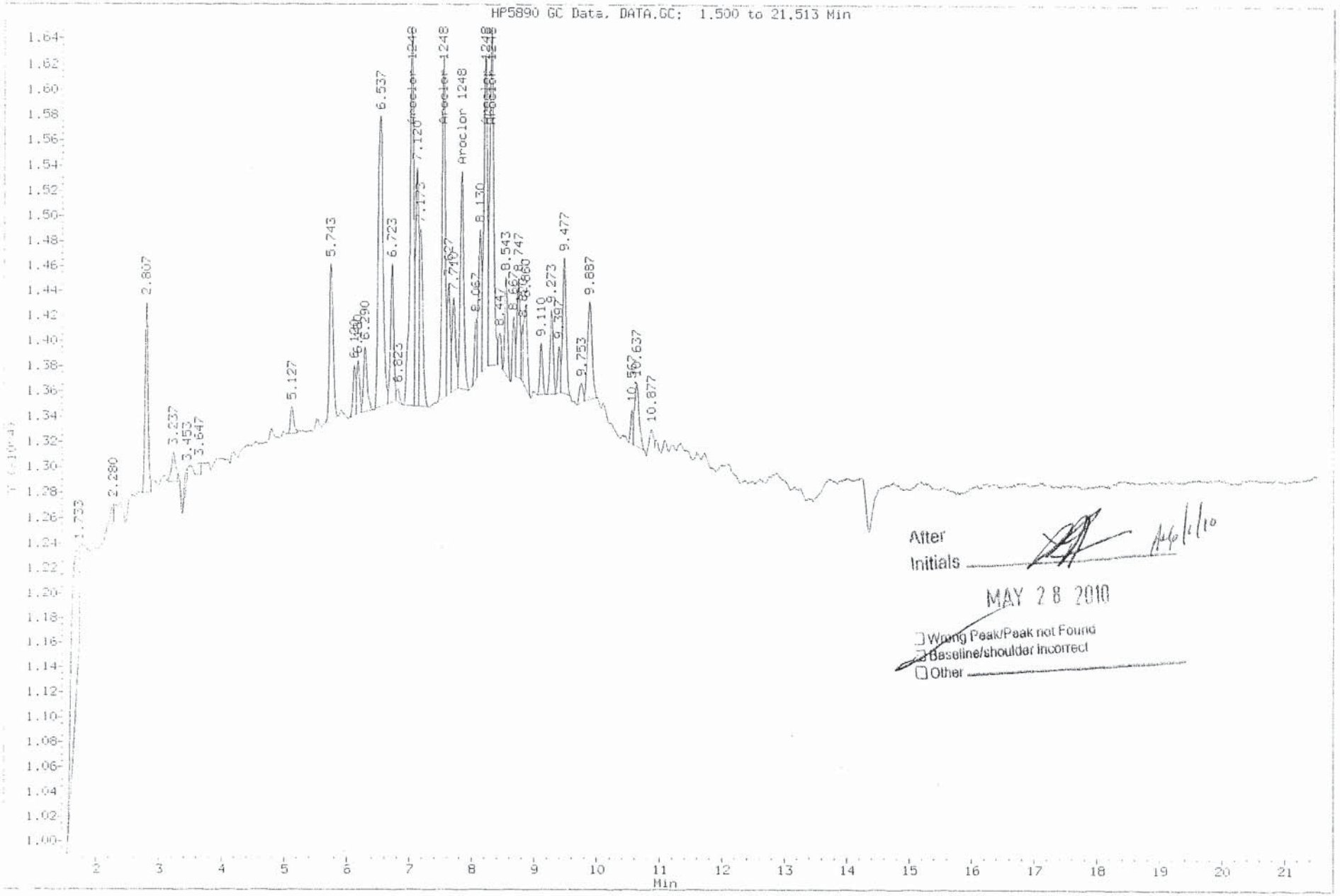
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F046.D
Injection Date: 28-MAY-2010 10:19
Instrument: GC09.i
Client Sample ID:



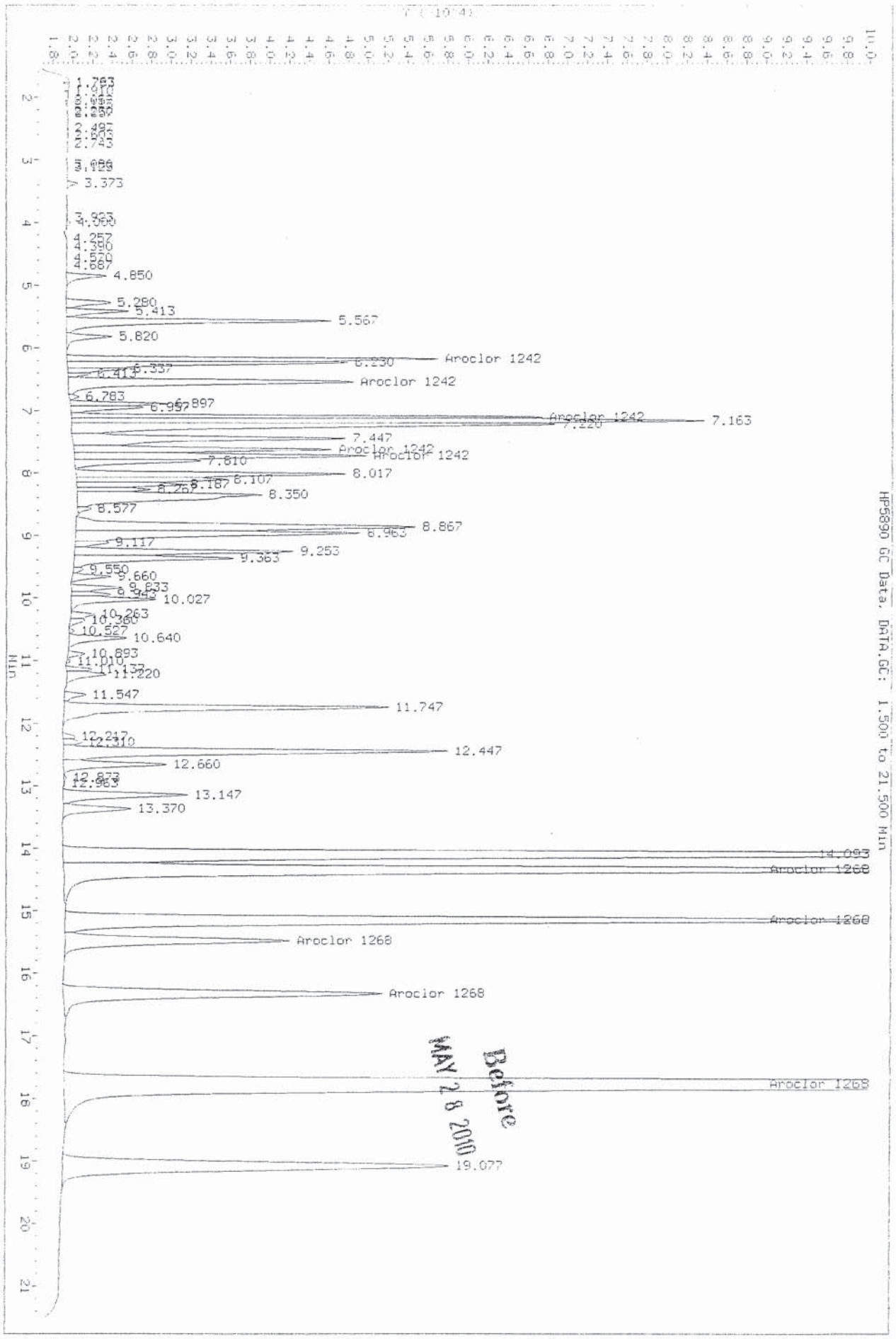
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F046.D
Injection Date: 28-MAY-2010 10:19
Instrument: GC09.i
Client Sample ID:

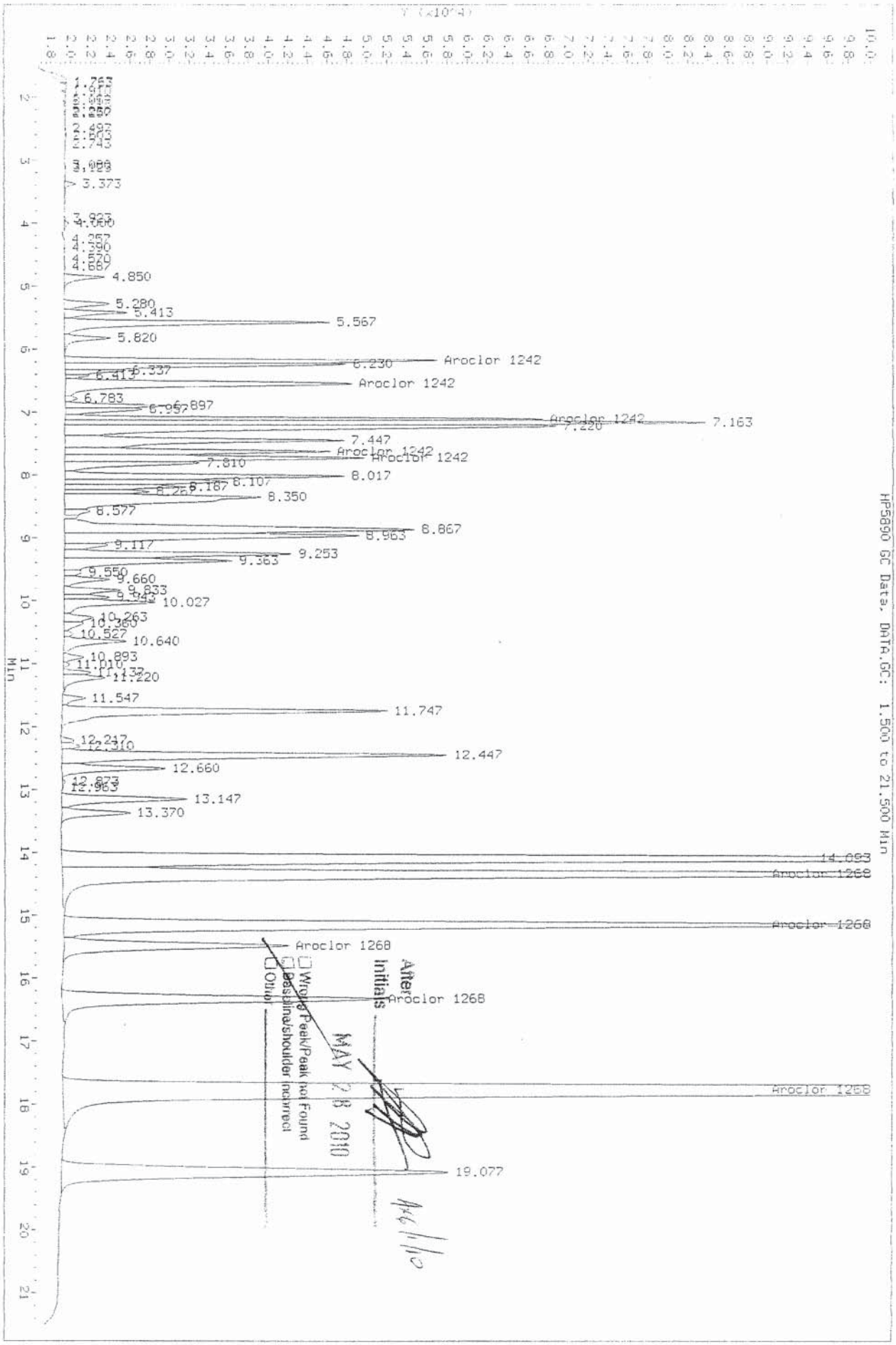


After
Initials _____ 5/28/10

MAY 28 2010

☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder incorrect
☐ Other _____


Data File: \\Cashl\Acqudata\5C09\data\052710A_r.b\052?R043.D
Injection Date: 28-MAY-2010 09:01
Instrument: 6C09.I
Client Sample ID:
HP6890 GC Data, DATA.GC; 1.500 to 21.500 Min
Aroclor 1242
Aroclor 1268
Before
MAY 28 2010
19.077


Data File: \\Cashl\hcqudata\GC09\data\052710A_r.b\052706043.D
Injection Date: 28-MAY-2010 09:01
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.500 Min
Min
Aroclor 1242
Aroclor 1242
Aroclor 1242
Aroclor 1242
Aroclor 1242
Aroclor 1268
Aroclor 1268
Aroclor 1268
Aroclor 1268
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other
Initials
After
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F044.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F044.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R044.D
Inj Date : 28-MAY-2010 09:27
Sample Info: 1242/1268 @ 2000ppb | PCB5-52T
Misc Info :
Cal Date : 28-MAY-2010 13:56
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1242+1268.sub
Sub List #2 : 1242+1268.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.120	6.230	142461	214001	1380	1940	80.00- 120.00	100.00
	6.540	6.543	1040414	306148	1680	1900	727.90-1091.85	730.31
	7.117	7.110	245200	242043	1690	1770	150.61- 225.91	172.12
	7.540	7.623	347926	211354	1760	1890	230.50- 345.75	244.23
	7.837	7.727	348197	250884	1870	1940	198.70- 298.04	244.42
			Average of Peak Amounts =		1680	1890		
Aroclor 1268	13.027	14.340	1769630	1436735	1750	1870	80.00- 120.00	100.00
	13.210	15.150	1673324	1132862	1790	1860	70.89- 106.33	94.56
	13.977	15.480	1342323	321903	1780	2000	58.04- 87.06	75.85
	14.943	16.330	608256	497952	1970	1990	20.96- 31.43	34.37
	16.210	17.770	3655383	3237262	1710	1780	172.26- 258.39	206.56
			Average of Peak Amounts =		1800	1900		

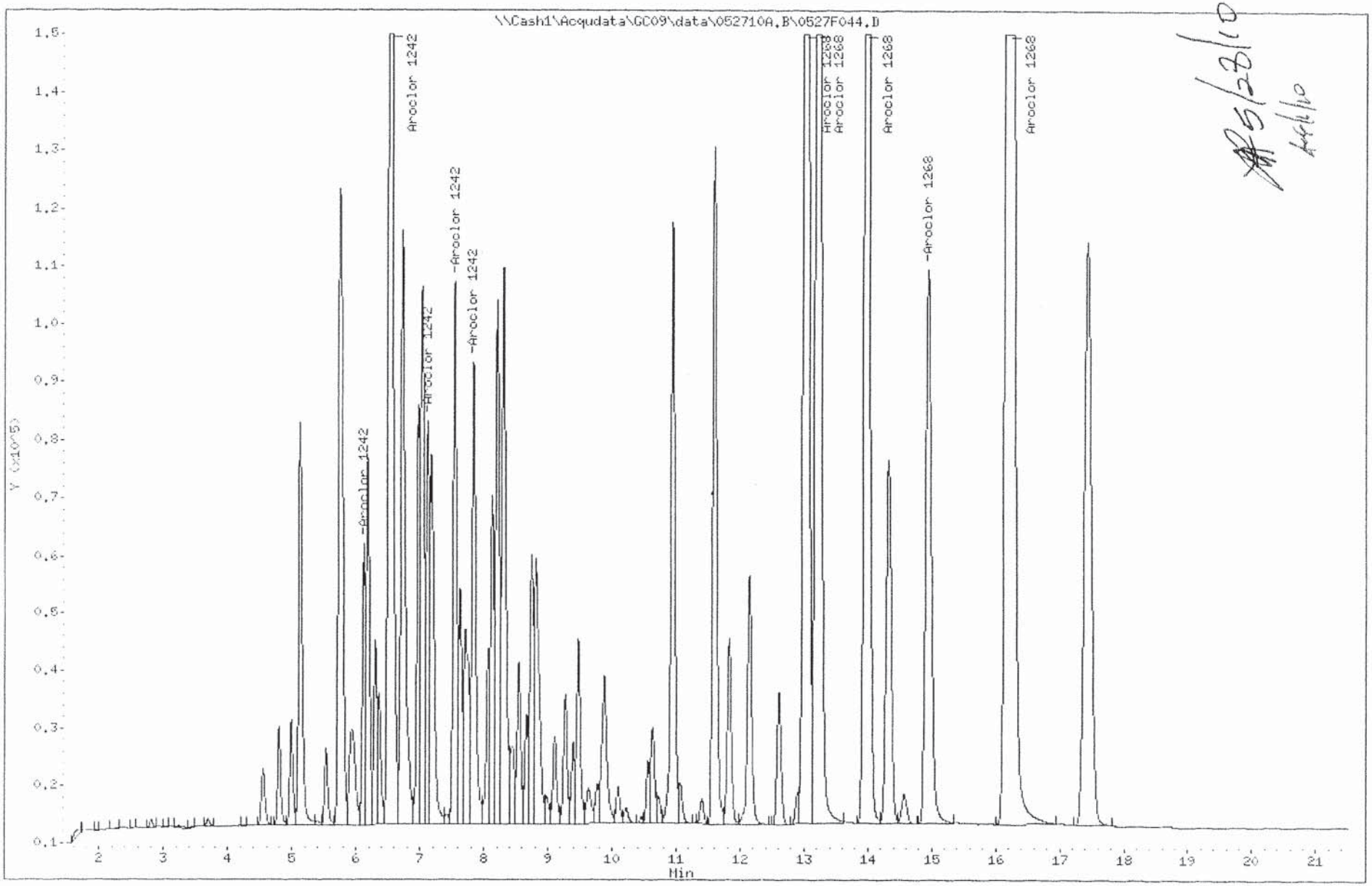
5/28/10
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F044.D
Date : 28-MAY-2010 09:27
Client ID:
Sample Info: 1242/1268 @ 2000ppb | PCB5-52T

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

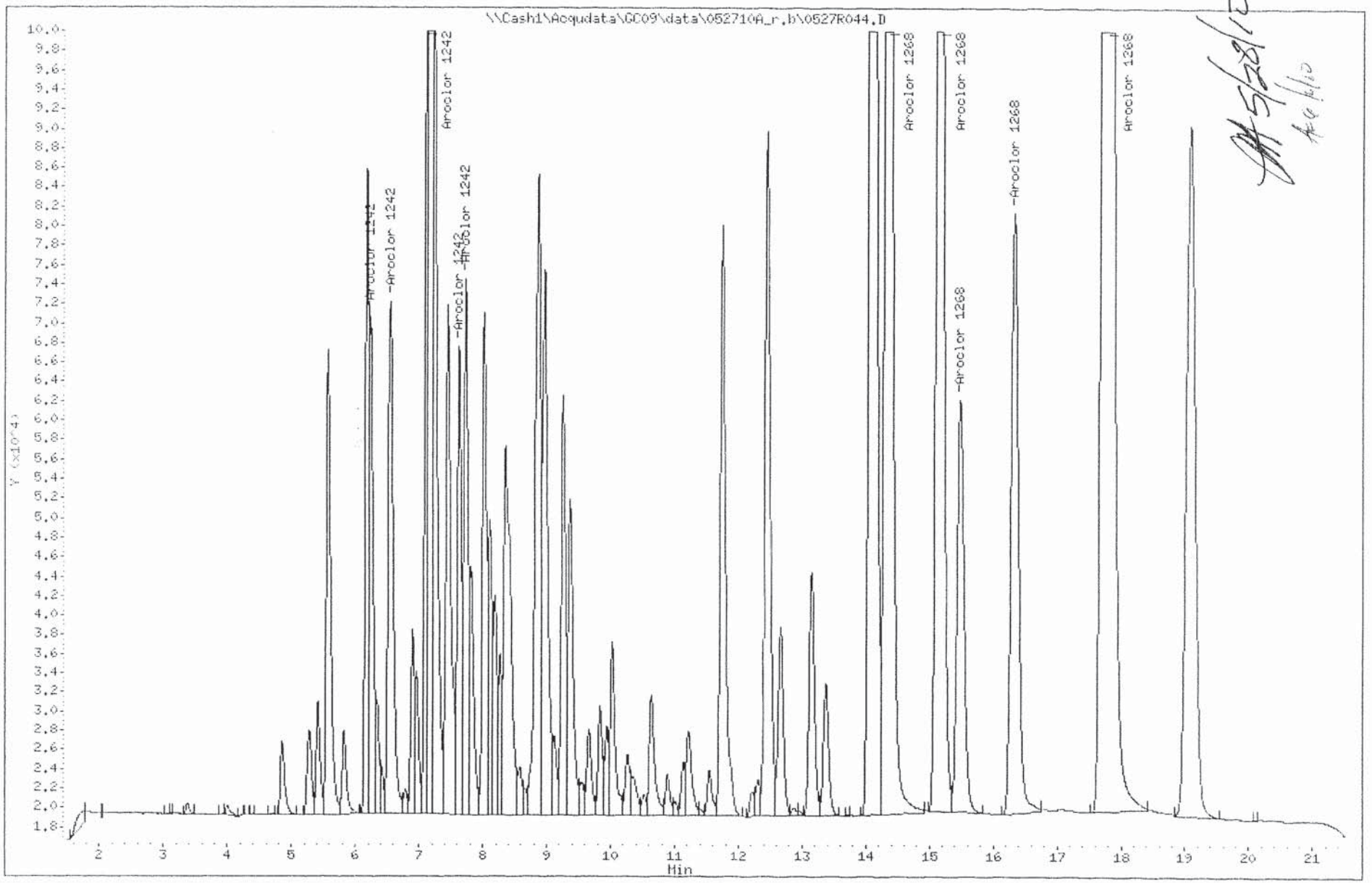
Column phase: DB-35MS



844

Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R044.D
Date : 28-MAY-2010 09:27
Client ID:
Sample Info: 1242/1268 @ 2000ppb | PCB5-52T

Instrument: GC09,i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F045.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F045.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R045.D
Inj Date : 28-MAY-2010 09:53
Sample Info: 1242/1268 @ 5000ppb | PCB5-53A
Misc Info :
Cal Date : 28-MAY-2010 13:56
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : 1242+1268.sub
Sub List #2 : 1242+1268.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.120	6.230	306390	472202	4000	4310	80.00- 120.00	100.00
	6.540	6.547	2176586	681089	3500	4230	727.90-1091.85	710.40
	7.120	7.110	522070	530644	3820	3890	150.61- 225.91	170.39
	7.540	7.623	743958	457086	3760	4080	230.50- 345.75	242.81
	7.837	7.727	751231	550847	4030	4260	198.70- 298.04	245.19
			Average of Peak Amounts =		3820	4150		
Aroclor 1268	13.027	14.340	3909122	3239299	3860	4220	80.00- 120.00	100.00
	13.210	15.150	3716596	2582867	3980	4250	70.89- 106.33	95.07
	13.977	15.480	2983479	757129	3960	4700	58.04- 87.06	76.32
	14.943	16.327	1395840	1181726	4520	4720	20.96- 31.43	35.71
	16.213	17.770	7966697	7114555	3730	3920	172.26- 258.39	203.80
			Average of Peak Amounts =		4010	4360		

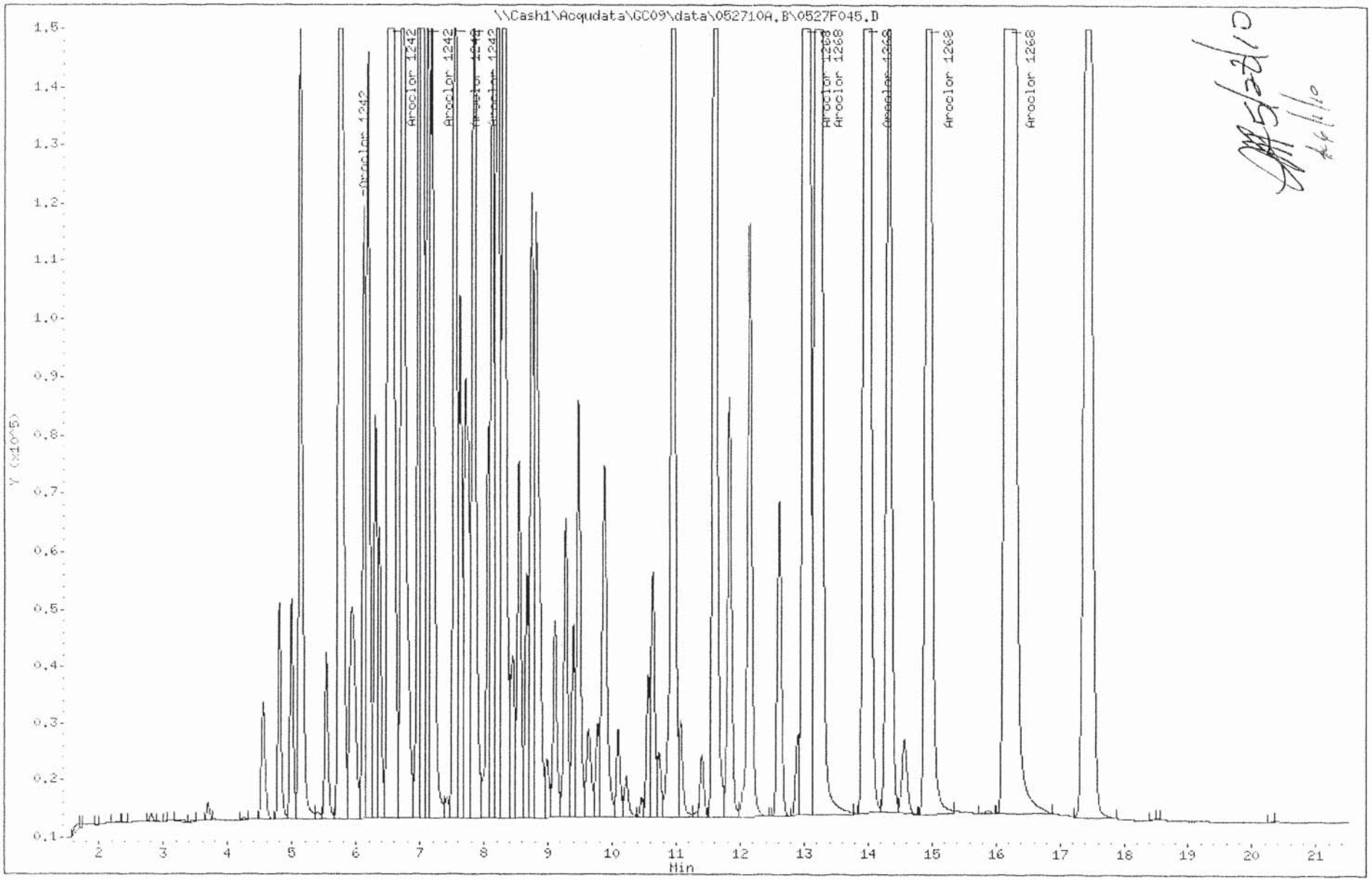
5/28/10
for 6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F045.D
Date : 28-MAY-2010 09:53
Client ID:
Sample Info: 1242/1268 @ 5000ppb | PCB5-53A

Instrument: GC09.i

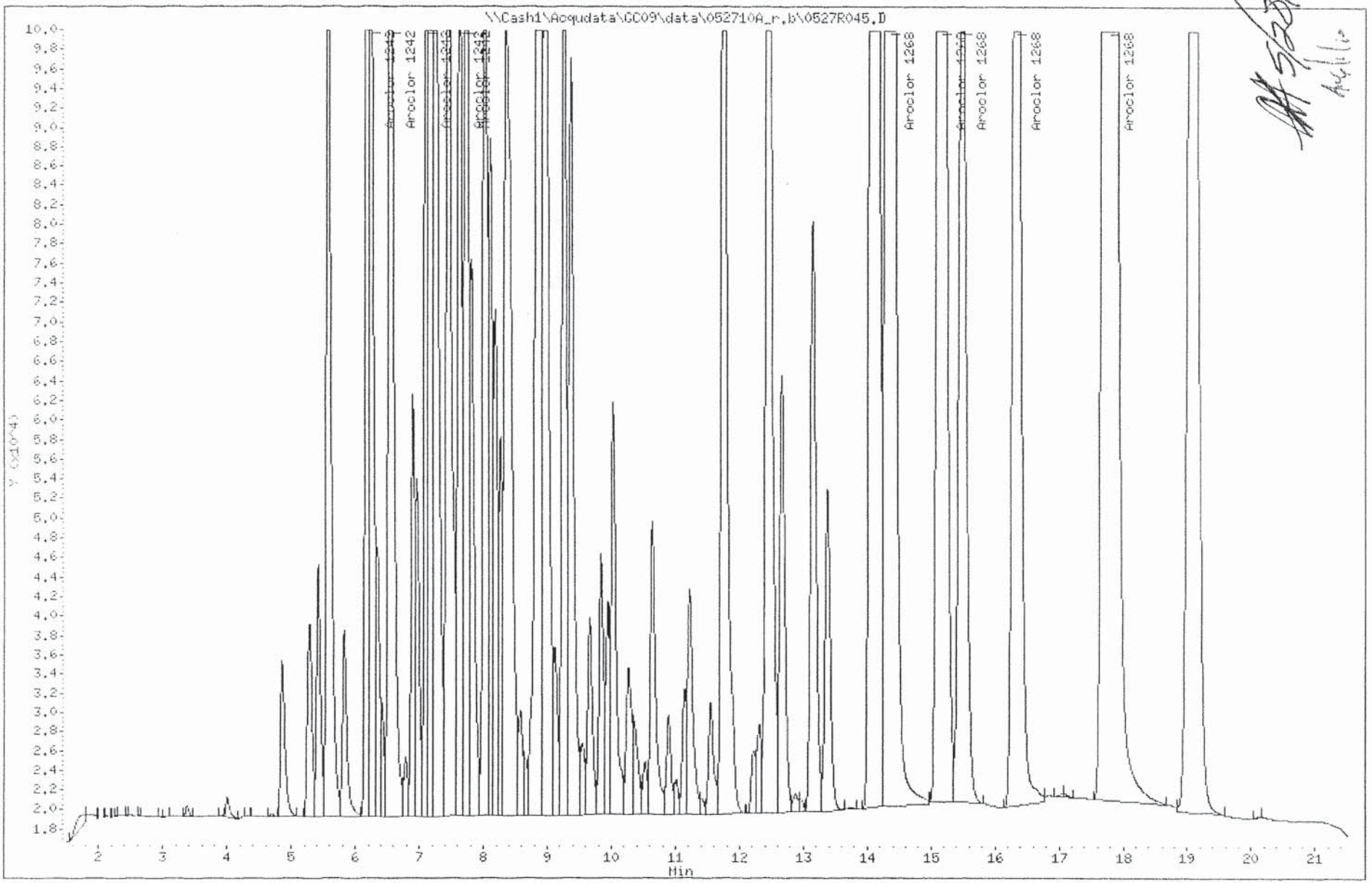
Operator: LHarris
Column diameter: 0.53

Column phase: DB-35MS



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R045.D
Date : 28-MAY-2010 09:53
Client ID:
Sample Info: 1242/1268 @ 5000ppb | PCB5-53A

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F046.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F046.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R046.D
Inj Date : 28-MAY-2010 10:19
Sample Info: 1248 @ 25ppb | PCB5-53B
Misc Info :
Cal Date : 28-MAY-2010 14:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.537	7.623	12314	5847	29.8	30.1	80.00- 120.00	100.00(M)
	7.173	7.727	5168	5459	24.6	28.3	39.24- 58.85	41.97(M)
	7.837	8.017	6772	5490	27.0	28.2	43.38- 65.07	54.99(M)
	8.210	8.867	9986	10389	28.7	27.4	66.78- 100.17	81.09(M)
	8.307	8.967	11851	7209	27.9	26.2	84.10- 126.14	96.24(M)
			Average of Peak Amounts =		27.6	28.0		

QC Flag Legend

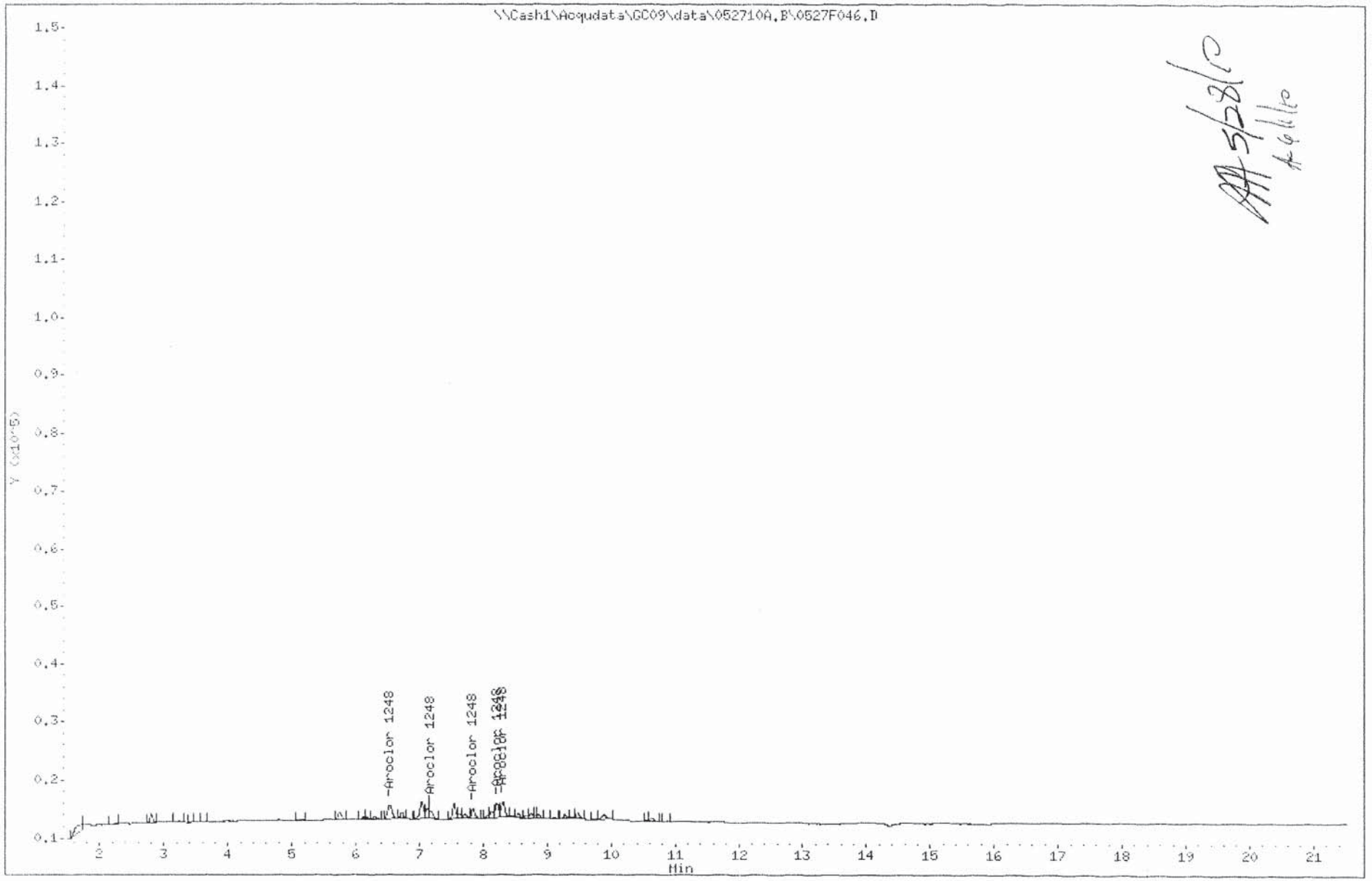
M - Compound response manually integrated.

5/28/10
6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F046.D
Date : 28-MAY-2010 10:19
Client ID:
Sample Info: 1248 @ 25ppb | PCB5-53B

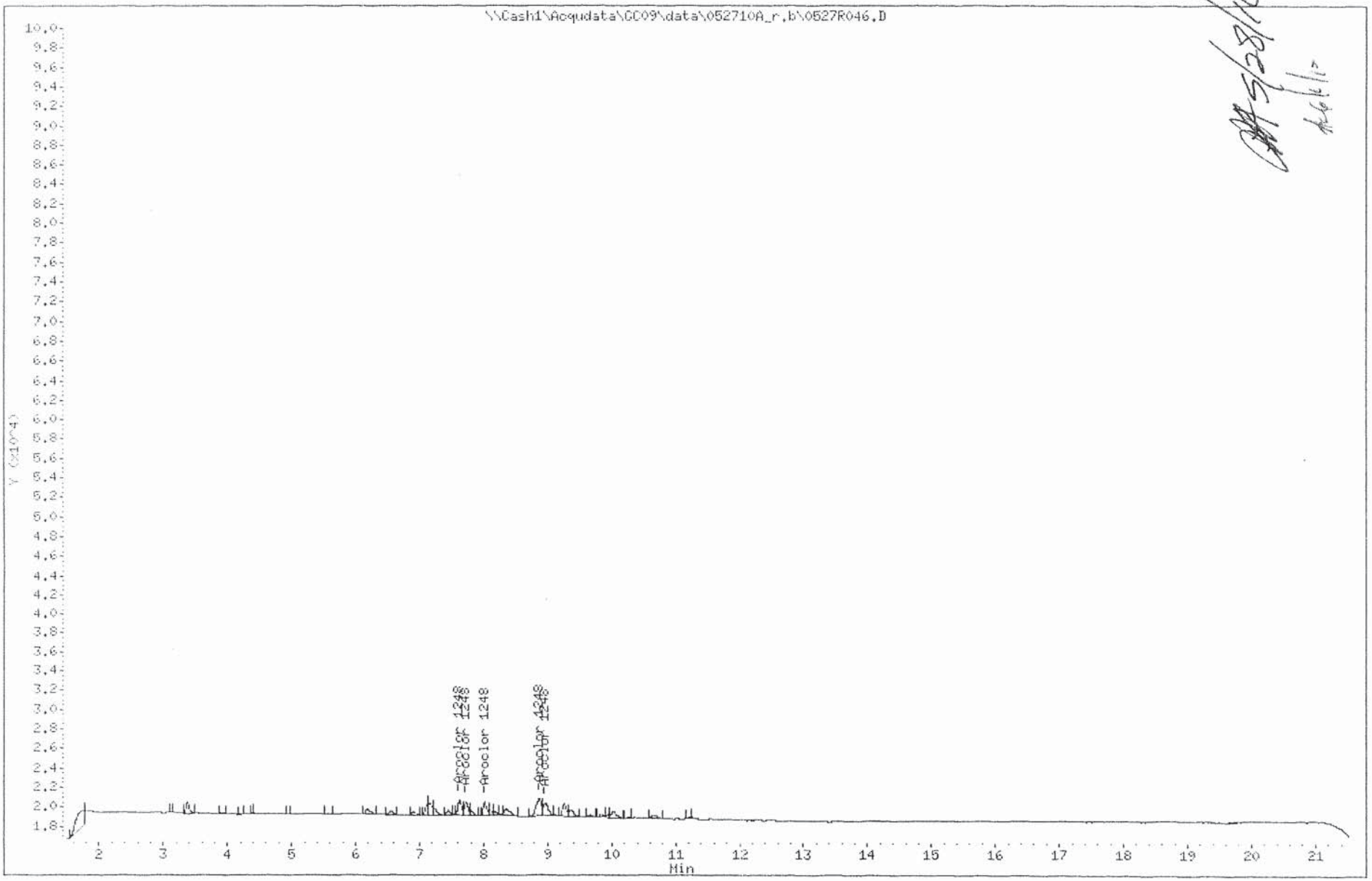
Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

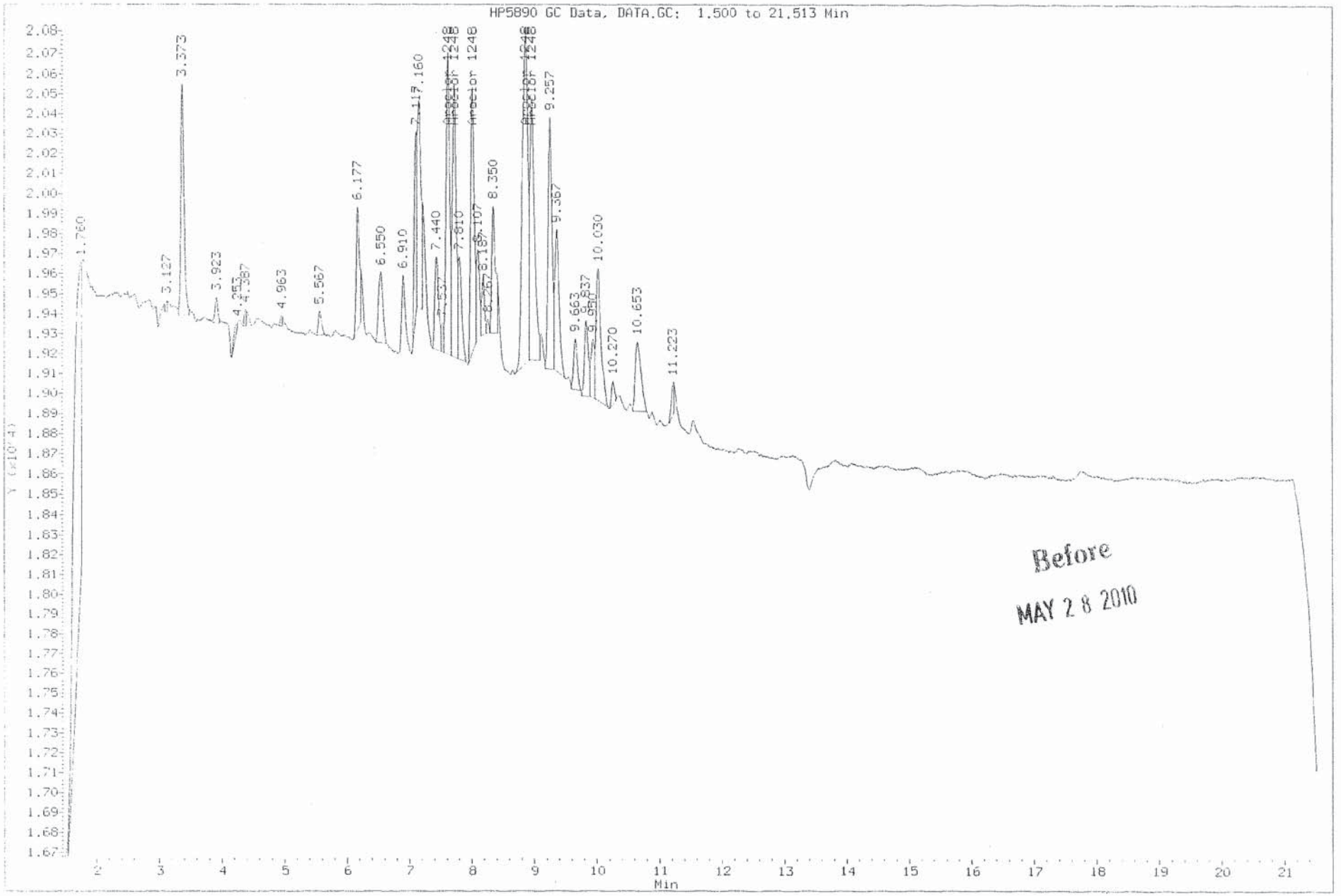


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R046.D
Date : 28-MAY-2010 10:19
Client ID:
Sample Info: 1248 @ 25ppb | PCB5-53B

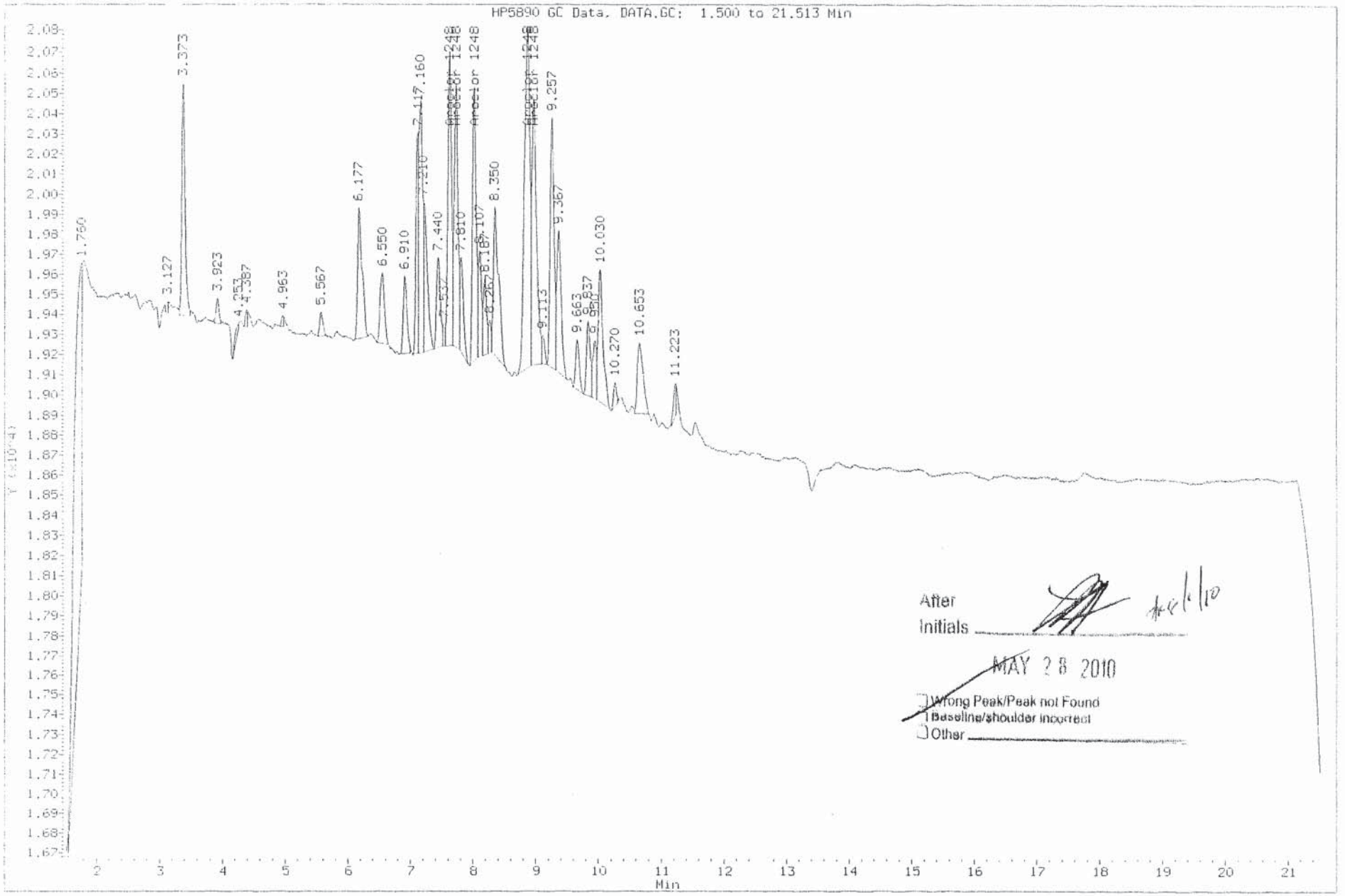
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R046.D
Injection Date: 28-MAY-2010 10:19
Instrument: GC09.i
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R046.D
Injection Date: 28-MAY-2010 10:19
Instrument: GC09.i
Client Sample ID:



After
Initials _____ 5/28/10

MAY 28 2010

☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder incorrect
☐ Other _____

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F047.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F047.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R047.D
Inj Date : 28-MAY-2010 10:45
Sample Info: 1248 @ 50ppb | PCB5-53C
Misc Info :
Cal Date : 28-MAY-2010 14:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.537	7.623	24649	12105	59.8	60.7	80.00- 120.00	100.00(M)
	7.177	7.727	12089	11359	57.6	57.4	39.24- 58.85	49.04(M)
	7.837	8.017	13367	11236	53.2	56.3	43.38- 65.07	54.23(M)
	8.210	8.867	20576	20900	59.2	55.2	66.78- 100.17	83.48(M)
	8.303	8.967	25911	15312	61.1	55.9	84.10- 126.14	105.12(M)
			Average of Peak Amounts =		58.2	57.1		

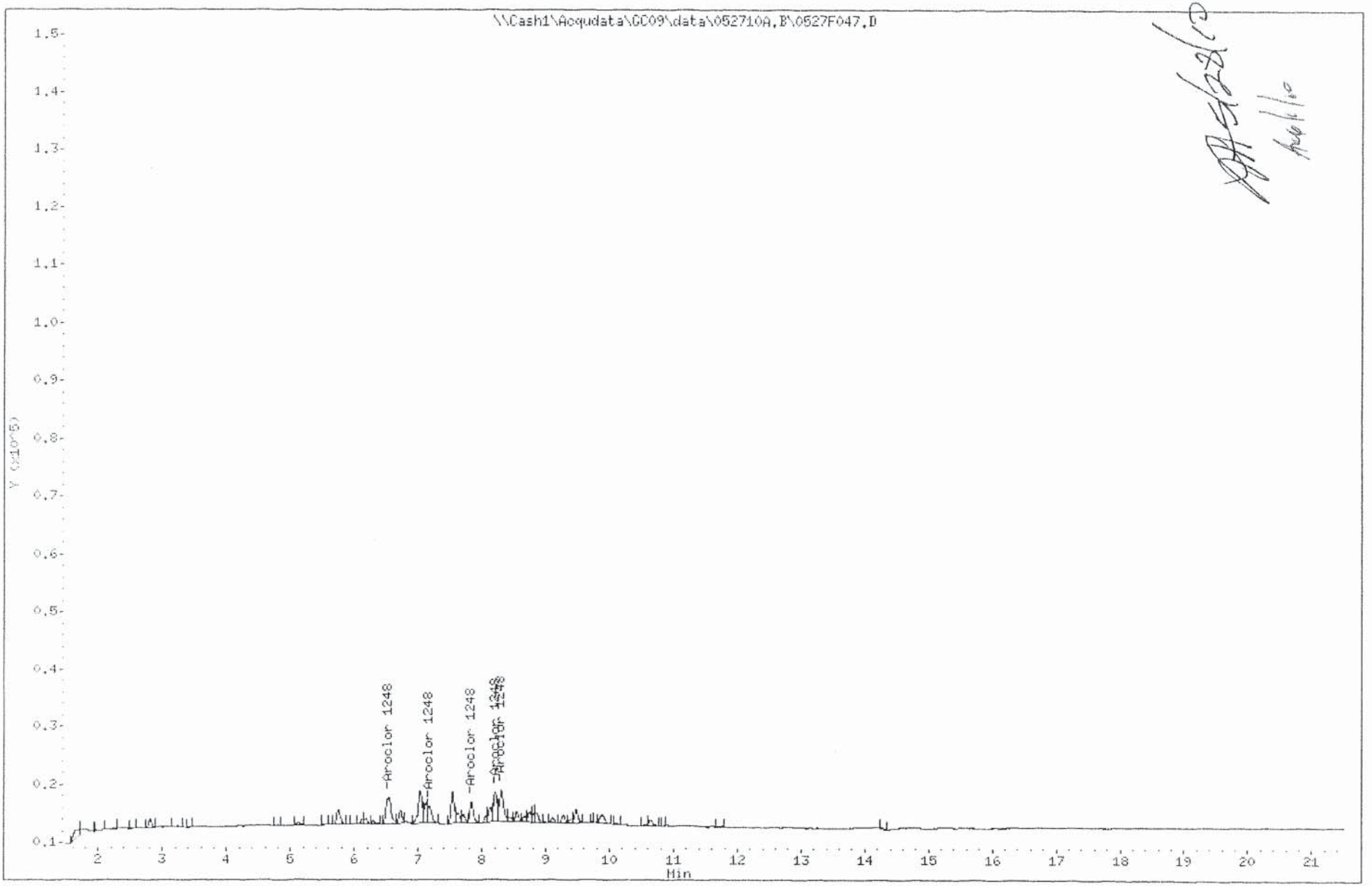
QC Flag Legend

M - Compound response manually integrated.

5/28/10

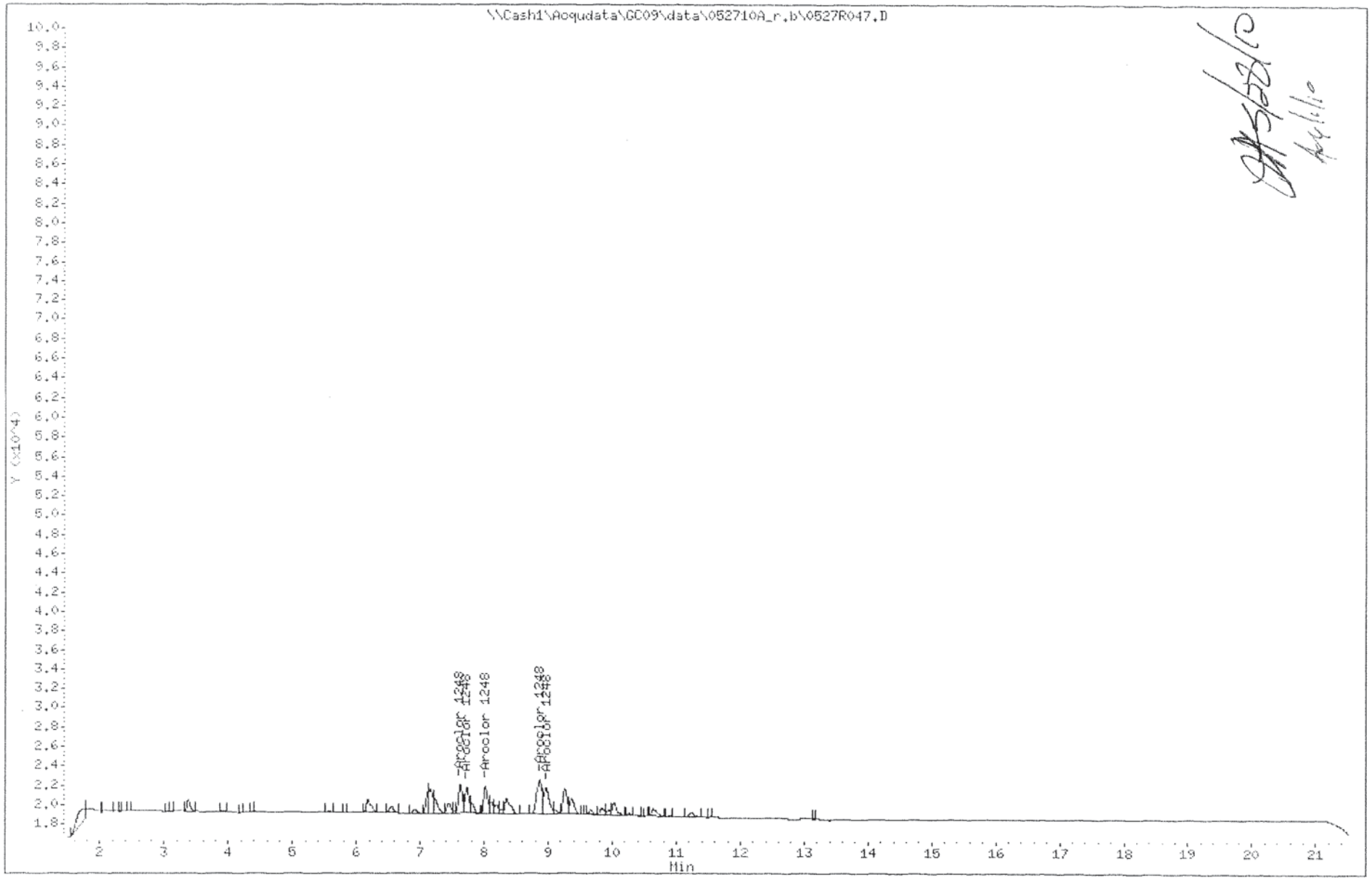
Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F047.D
Date : 28-MAY-2010 10:45
Client ID:
Sample Info: 1248 @ 50ppb | PCB5-53C

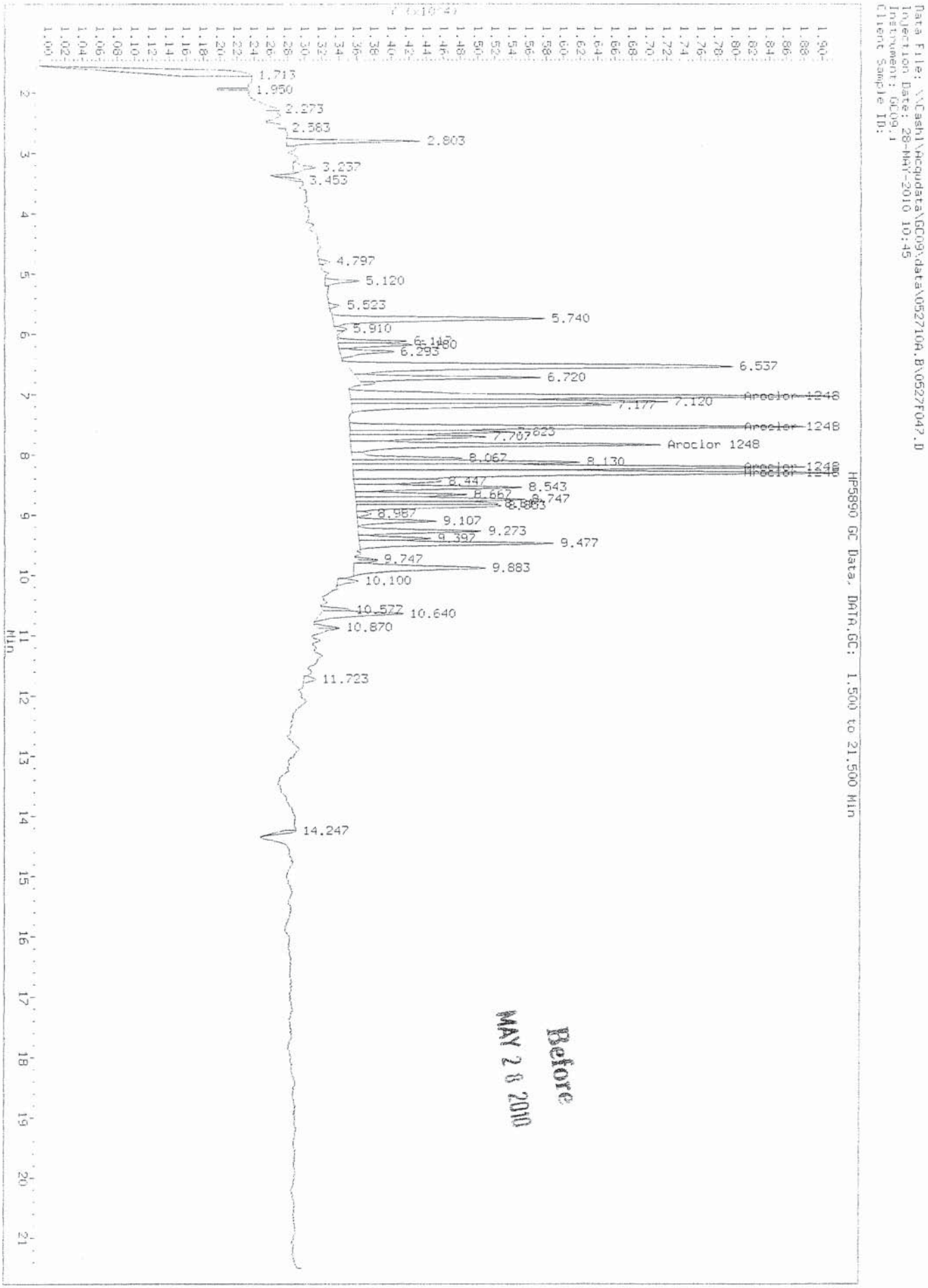
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



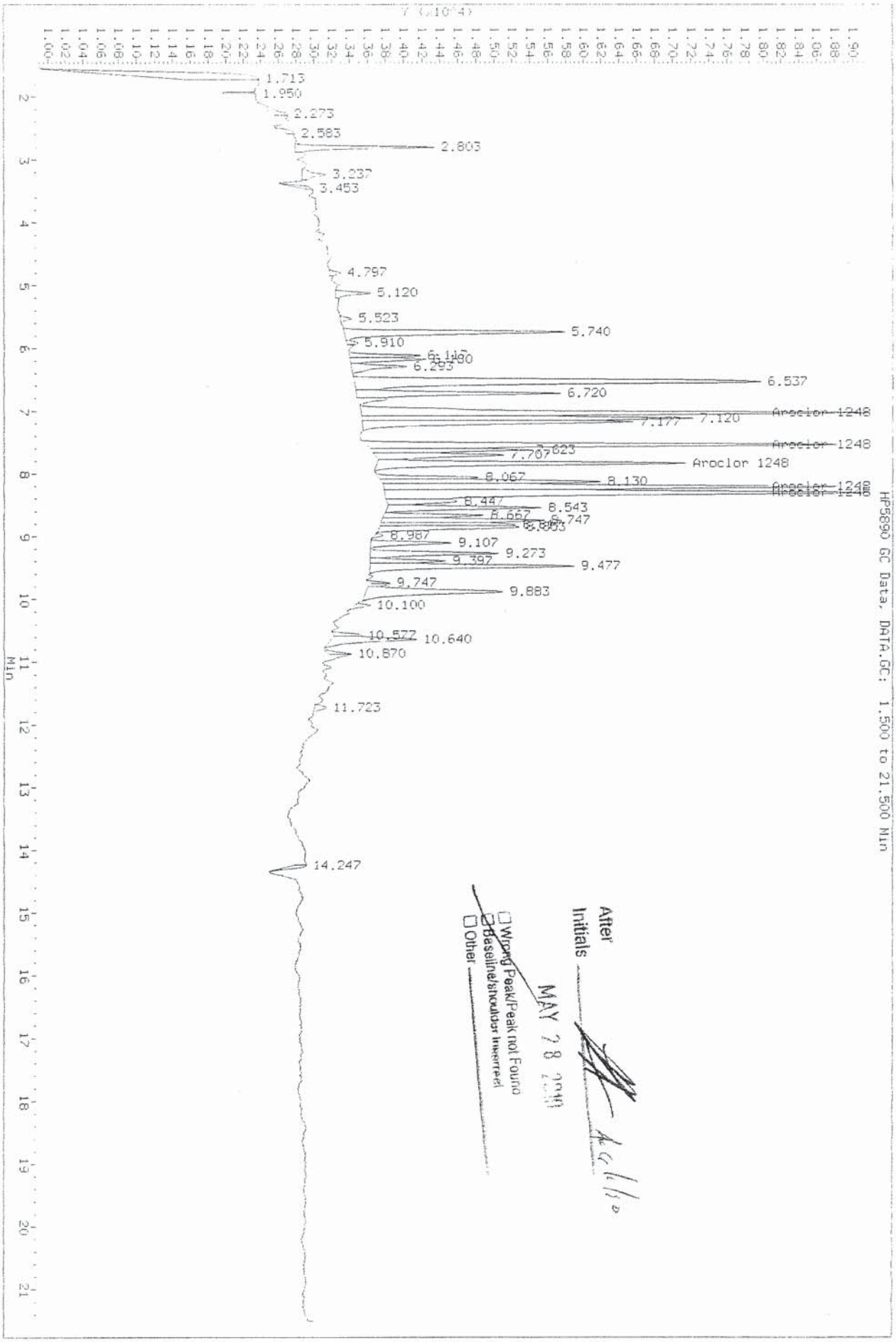
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R047.D
Date : 28-MAY-2010 10:45
Client ID:
Sample Info: 1248 @ 50ppb | PCB5-53C

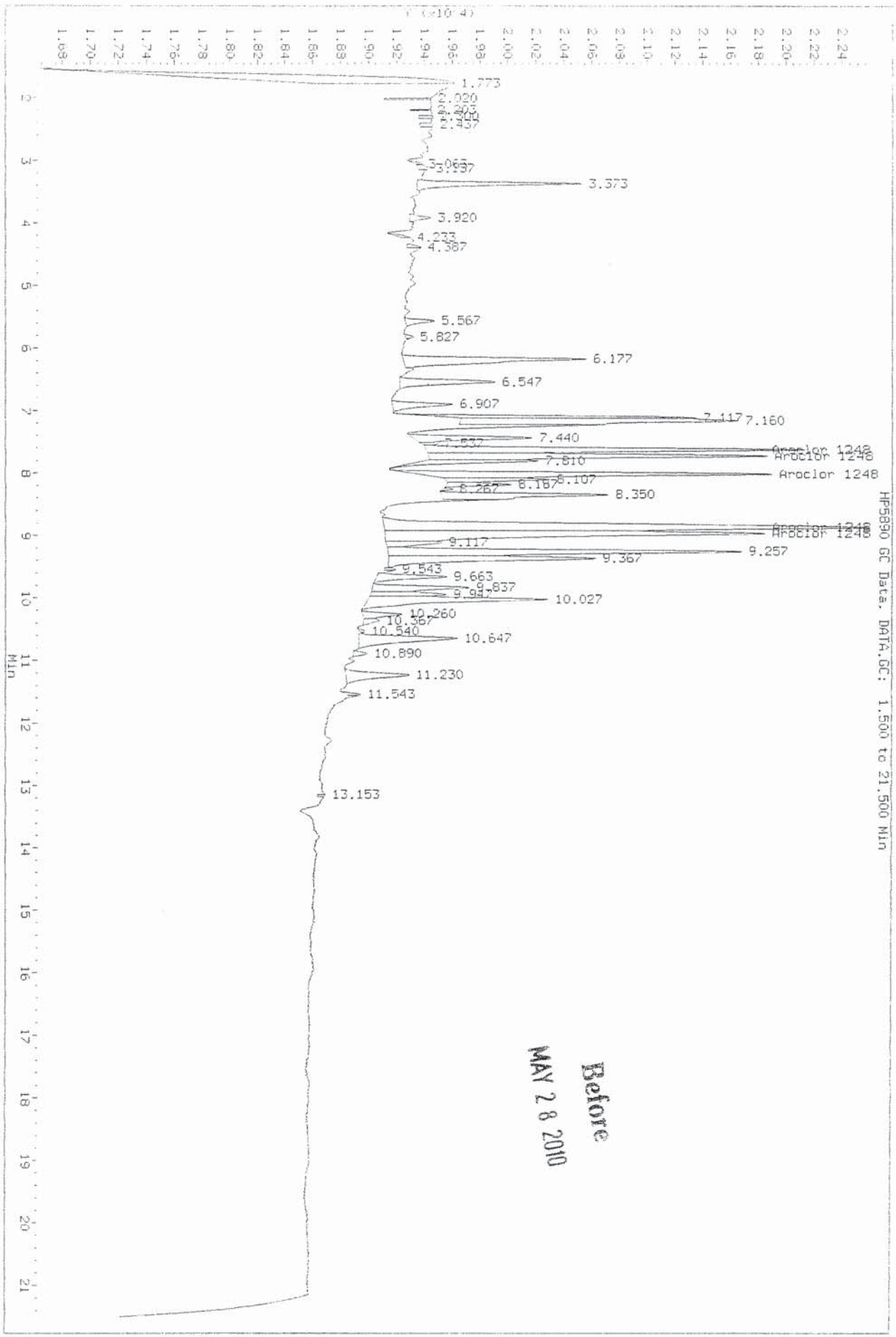
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53





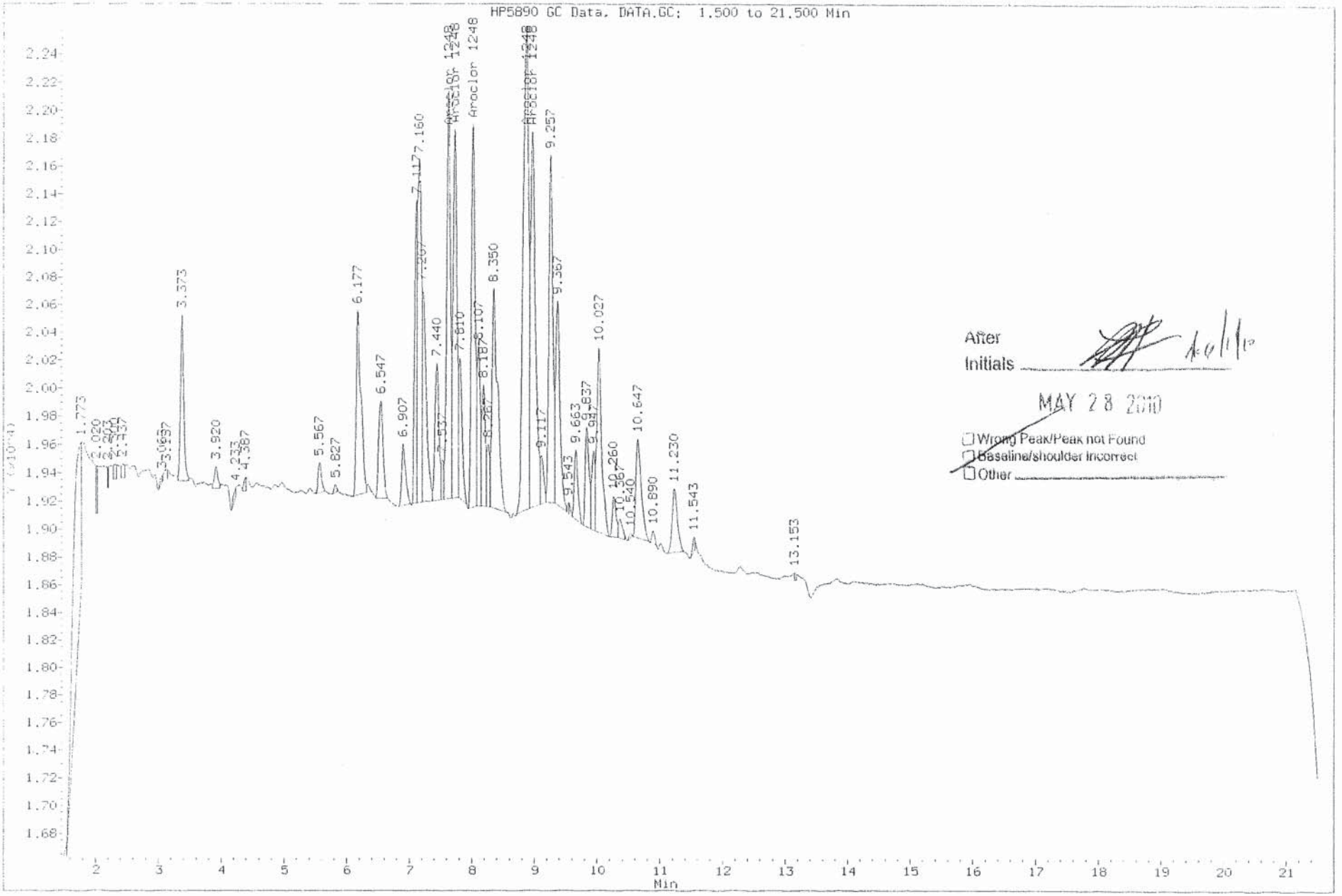
Before
MAY 2 6 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\052FG47.D
Injection Date: 28-MAY-2010 10:45
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.500 Min
Min
1.713
1.950
2.273
2.583
2.803
3.237
3.453
4.797
5.120
5.523
5.740
5.910
6.293
6.537
6.720
Aroclor 1248
7.177
7.120
Aroclor 1248
7.707
Aroclor 1248
8.067
8.130
Aroclor 1248
8.447
8.543
8.747
8.987
9.107
9.273
9.397
9.477
9.747
9.883
10.100
10.572
10.640
10.870
11.723
14.247
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/Integrator Incorrect
Other



Before
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R047.D
Injection Date: 28-MAY-2010 10:45
Instrument: GC09.1
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F048.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F048.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R048.D
Inj Date : 28-MAY-2010 11:11
Sample Info: 1248 @ 500ppb | PCB5-53D
Misc Info :
Cal Date : 28-MAY-2010 14:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

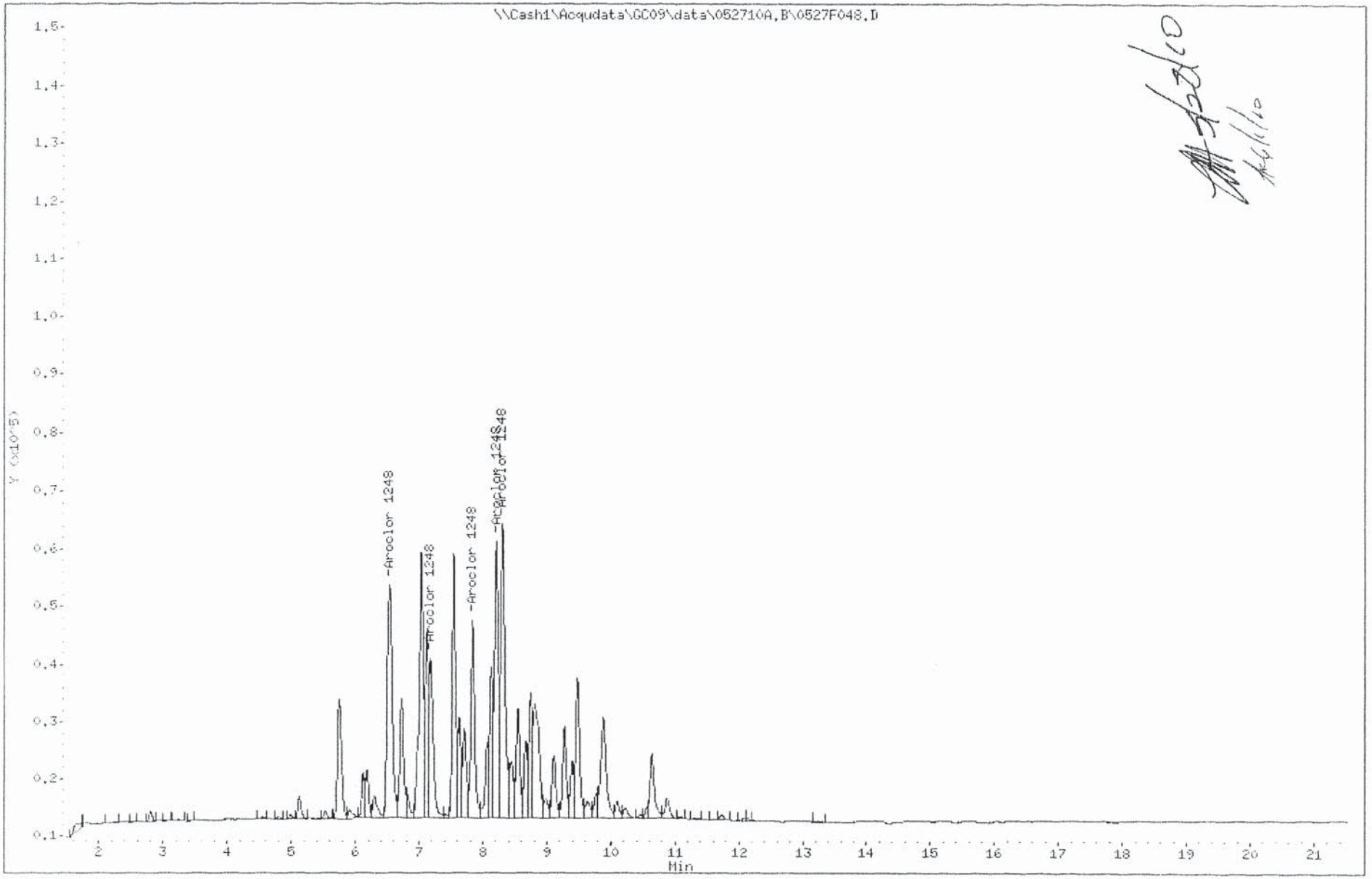
Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.537	7.623	226287	108687	691	545	80.00- 120.00	100.00
	7.177	7.723	121416	111692	578	564	39.24- 58.85	53.66
	7.837	8.013	144855	110702	577	555	43.38- 65.07	64.01
	8.207	8.867	192923	208975	555	552	66.78- 100.17	85.26
	8.303	8.963	240006	152745	566	558	84.10- 126.14	106.06
			Average of Peak Amounts =		593	555		

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F048.D
Date : 28-MAY-2010 11:11
Client ID:
Sample Info: 1248 @ 500ppb | PCB5-53D

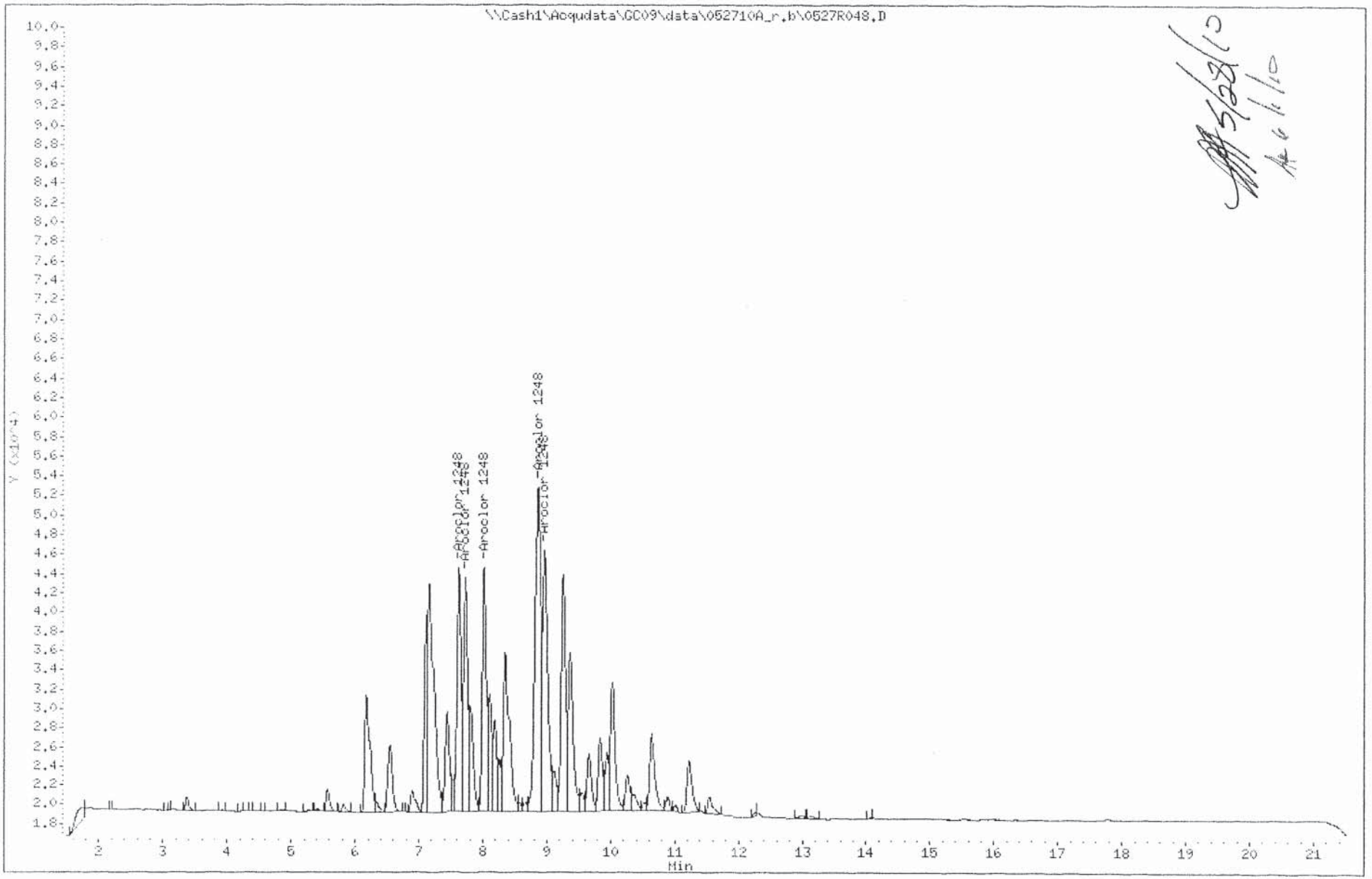
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R048.D
Date : 28-MAY-2010 11:11
Client ID:
Sample Info: 1248 @ 500ppb | PCB5-53D

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0,53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F049.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F049.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R049.D
Inj Date : 28-MAY-2010 11:37
Sample Info: 1248 @ 1000ppb | PCB5-53E
Misc Info :
Cal Date : 28-MAY-2010 14:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

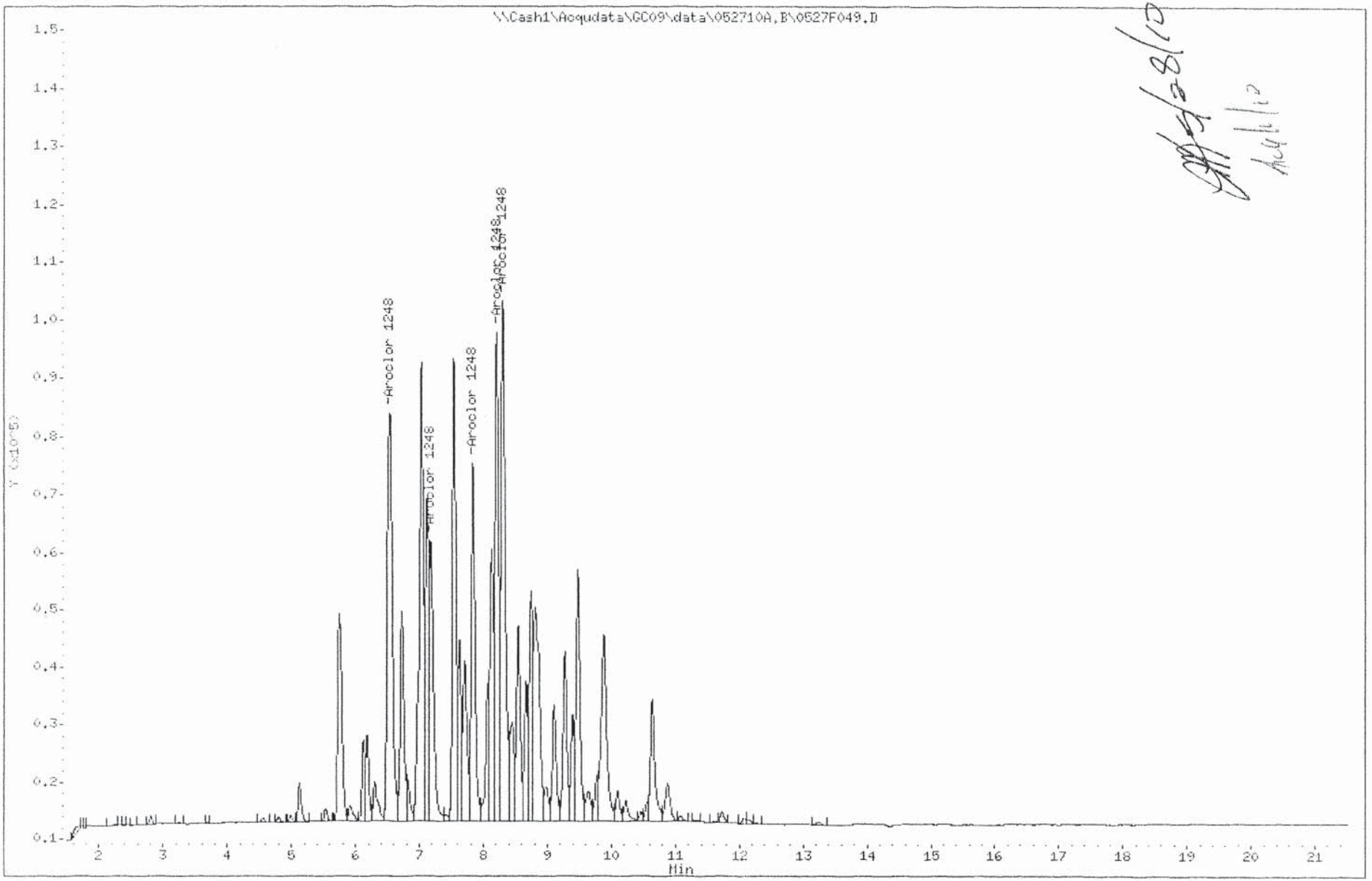
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.537	7.623	394337	191702	1090	962	80.00- 120.00	100.00
	7.177	7.723	219634	193574	1050	978	39.24- 58.85	55.70
	7.837	8.013	256124	196694	1020	986	43.38- 65.07	64.95
	8.207	8.863	337925	374966	973	991	66.78- 100.17	85.69
	8.303	8.963	412366	271790	972	993	84.10- 126.14	104.57
			Average of Peak Amounts =		1020	982		

5/28/10
Acq 5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F049.D
Date : 28-MAY-2010 11:37
Client ID:
Sample Info: 1248 @ 1000ppb | PCB5-53E

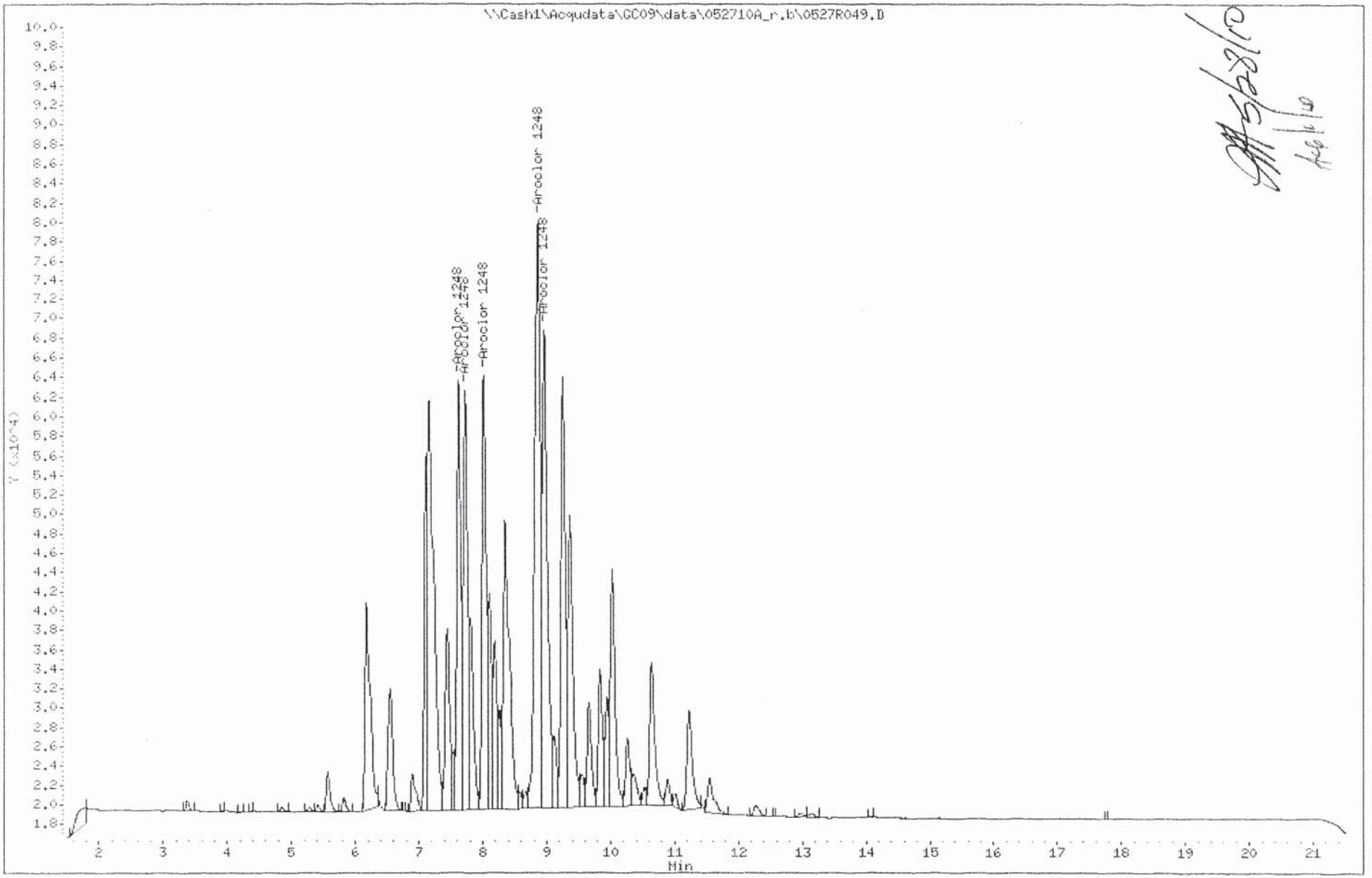
Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R049.D
Date : 28-MAY-2010 11:37
Client ID:
Sample Info: 1248 @ 1000ppb | PCB5-53E

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F050.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F050.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R050.D
Inj Date : 28-MAY-2010 12:03
Sample Info: 1248 @ 2000ppb | PCB5-53F
Misc Info :
Cal Date : 28-MAY-2010 14:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

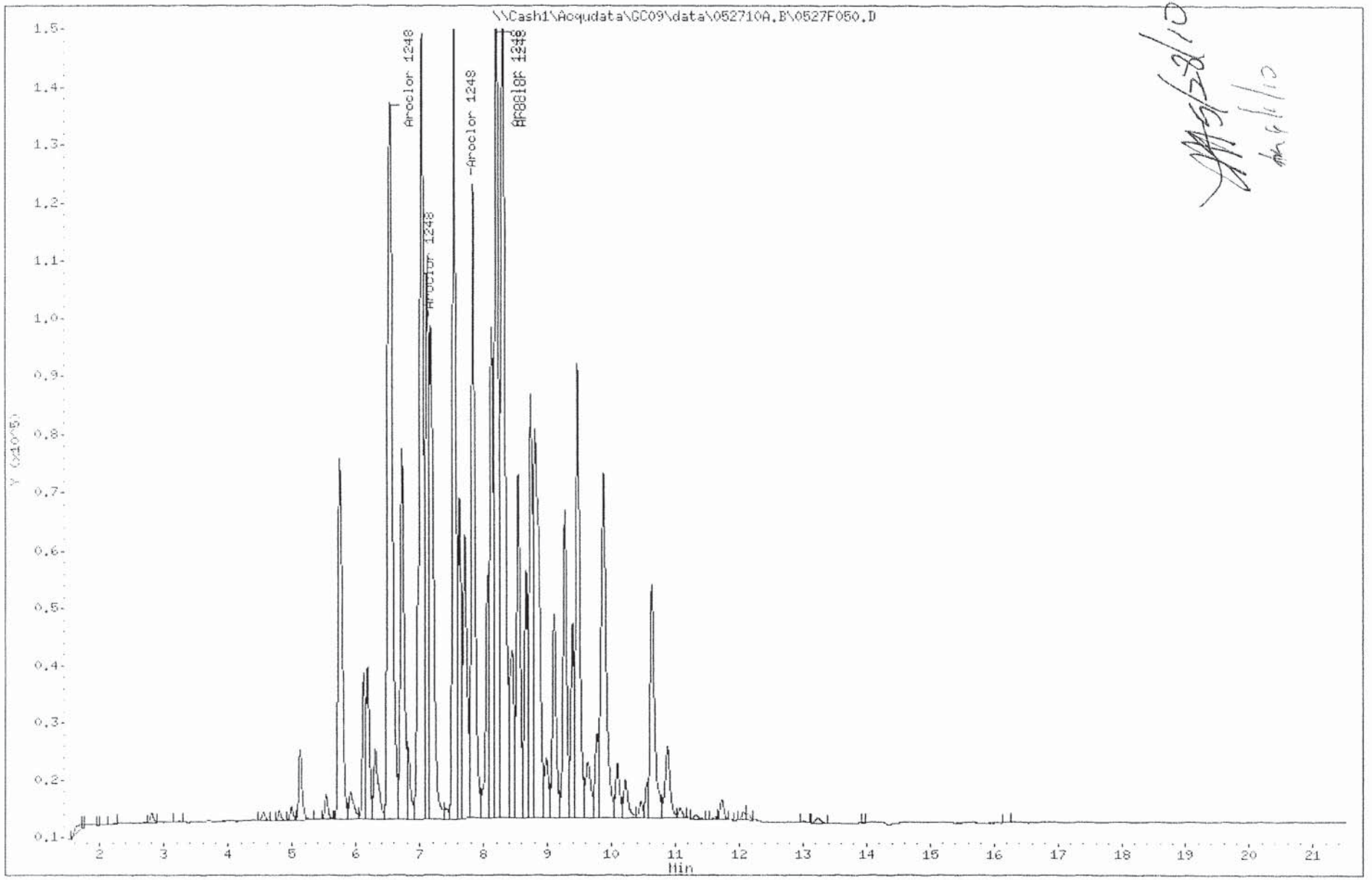
Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.537	7.620	688597	336241	1770	1680	80.00- 120.00	100.00
	7.177	7.723	376350	354966	1790	1790	39.24- 58.85	54.65
	7.833	8.013	449526	356802	1790	1780	43.38- 65.07	65.28
	8.207	8.863	590433	676997	1700	1780	66.78- 100.17	85.74
	8.303	8.963	706812	504603	1670	1840	84.10- 126.14	102.65
			Average of Peak Amounts =		1740	1770		

5/28/10
Acc 6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F050.D
Date : 28-MAY-2010 12:03
Client ID:
Sample Info: 1248 @ 2000ppb I PCB5-53F

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R050.D

Date : 28-MAY-2010 12:03

Client ID:

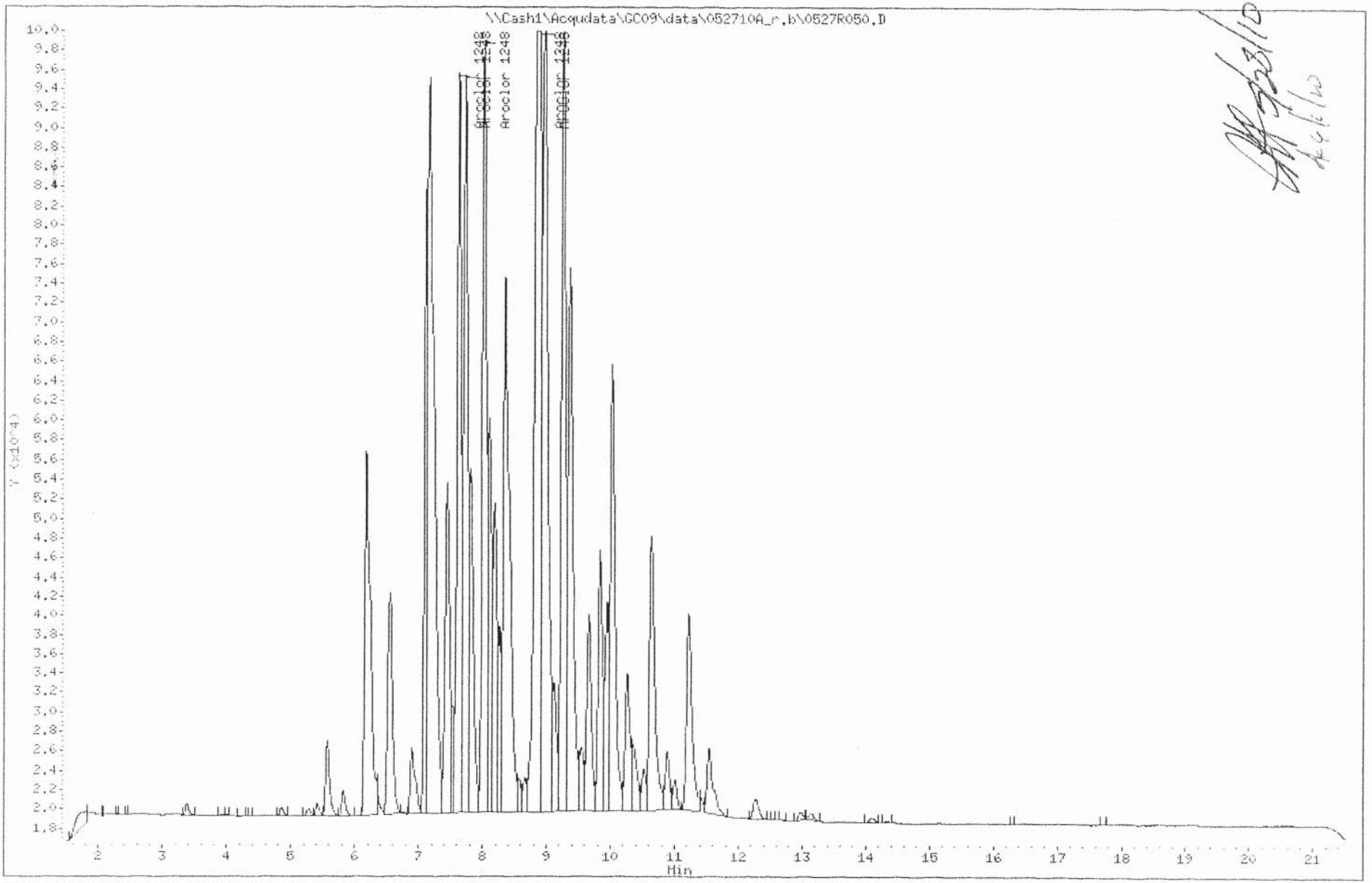
Sample Info: 1248 @ 2000ppb | PCB5-53F

Instrument: GC09.i

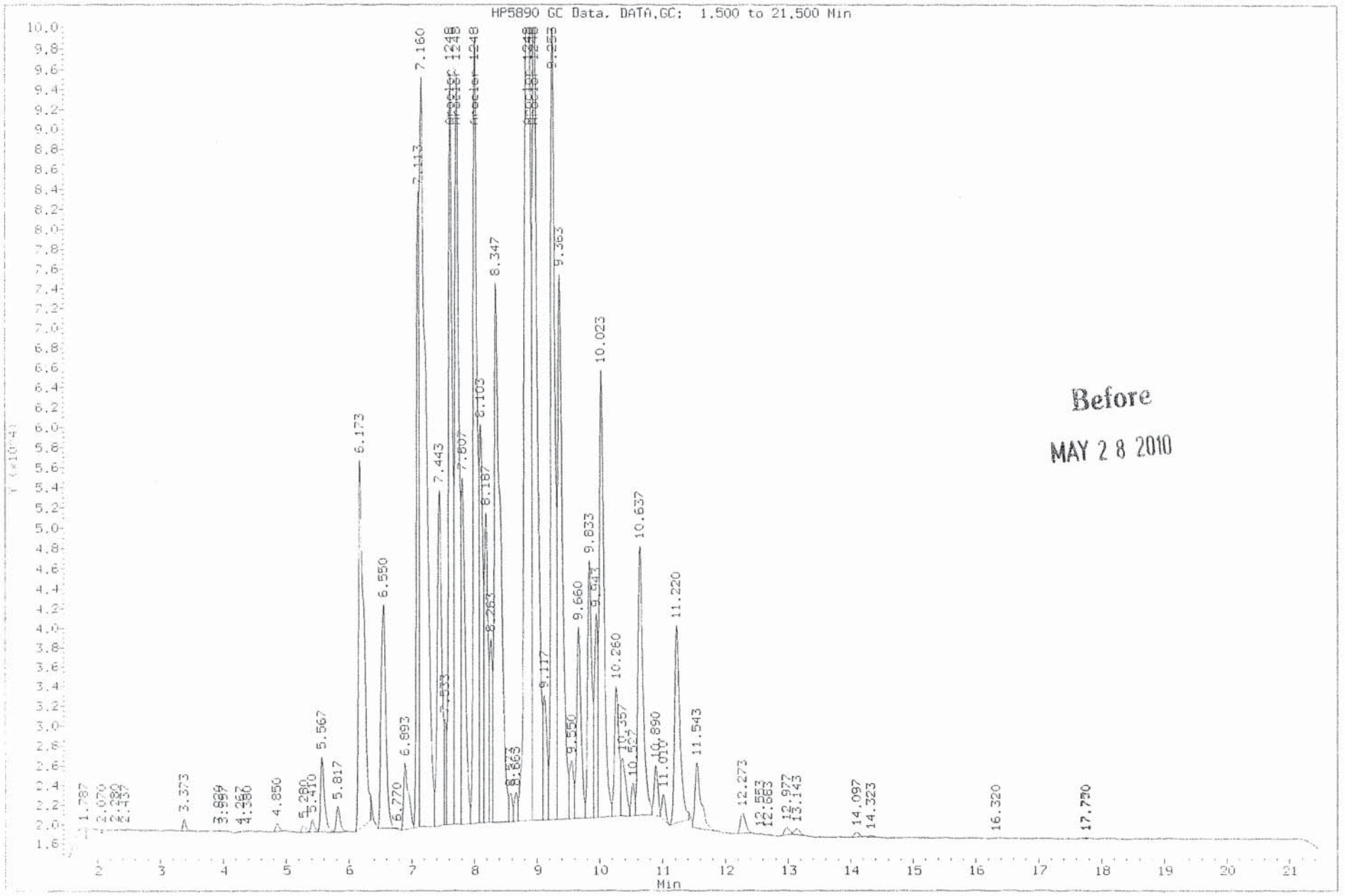
Operator: LHarris

Column phase: DB-XLB

Column diameter: 0.53



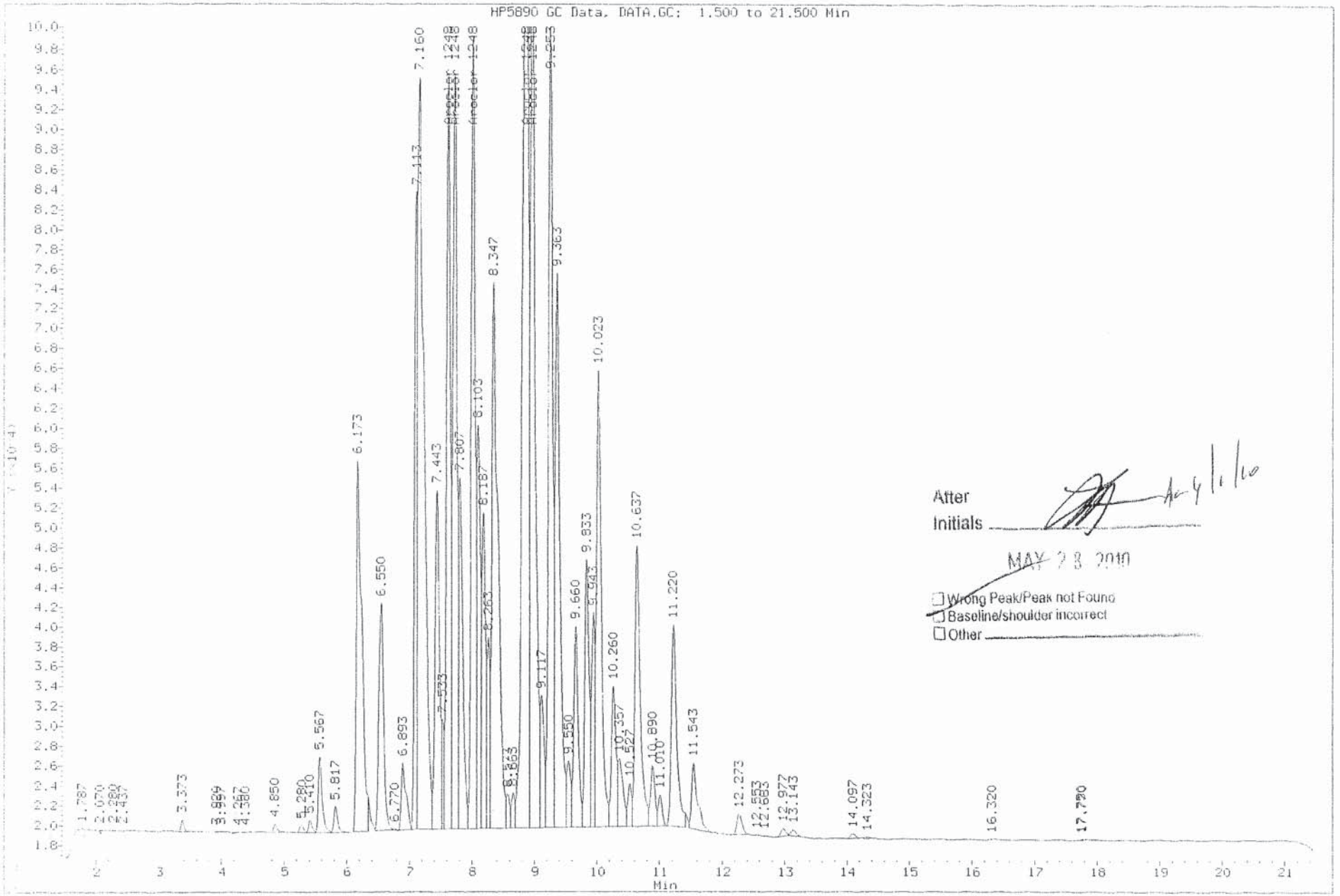
Data File: \\Cash1\Acqudata\GC09\data\0527104_r.b\0527R050.D
Injection Date: 28-MAY-2010 12:03
Instrument: GC09.i
Client Sample ID:



Before

MAY 2 8 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R050.D
Injection Date: 28-MAY-2010 12:03
Instrument: GC09.1
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F051.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F051.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R051.D
Inj Date : 28-MAY-2010 12:29
Sample Info: 1248 @ 5000ppb | PCB5-53G
Misc Info :
Cal Date : 28-MAY-2010 14:04
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

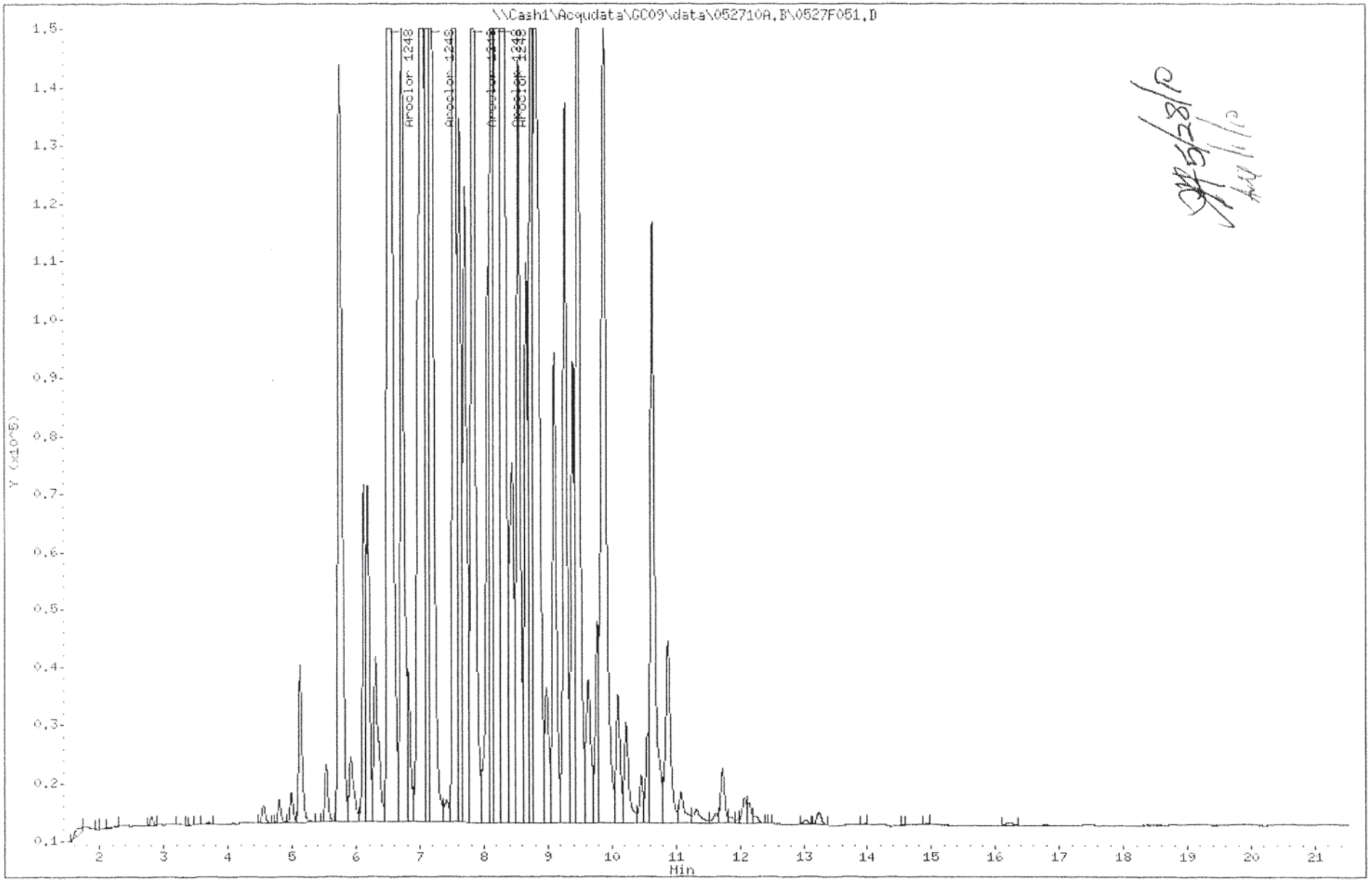
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.533	7.620	1481123	732014	3570	3660	80.00- 120.00	100.00
	7.177	7.723	799924	762186	3810	3840	39.24- 58.85	54.01
	7.833	8.013	988197	798492	3940	3990	43.38- 65.07	66.72
	8.207	8.860	1272422	1558356	3660	4110	66.78- 100.17	85.91
	8.303	8.960	1534313	1125026	3620	4090	84.10- 126.14	103.59
			Average of Peak Amounts =		3720	3940		

5/28/10
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F051.D
Date : 28-MAY-2010 12:29
Client ID:
Sample Info: 1248 @ 5000ppb | PCB5-53G

Instrument: GC09,i

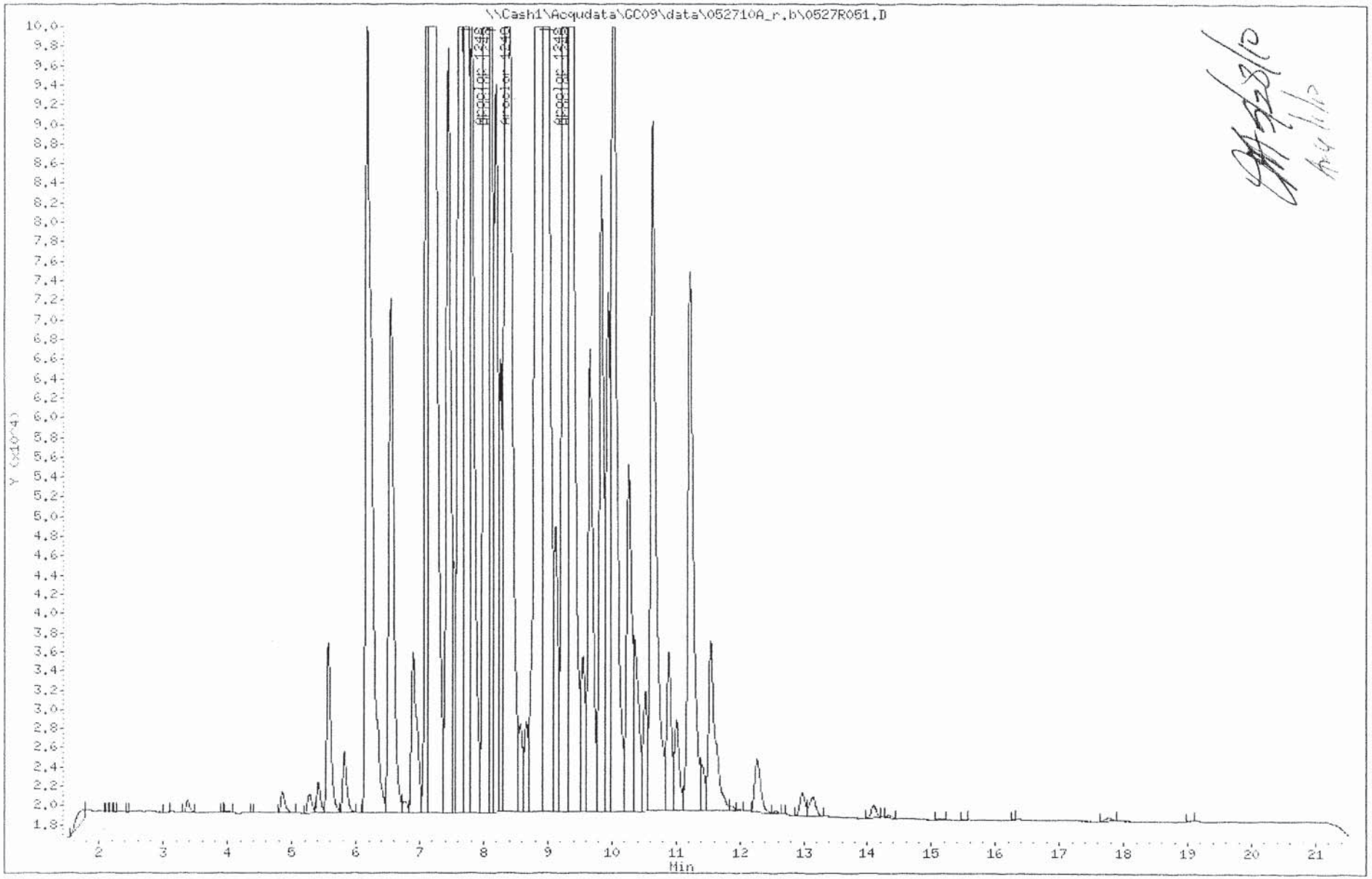
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R051.D
Date : 28-MAY-2010 12:29
Client ID:
Sample Info: 1248 @ 5000ppb | PCB5-53C

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F052.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F052.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R052.D
Inj Date : 28-MAY-2010 12:55
Sample Info: 1016 @ 1000ppb | PCB5-49E
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1016.sub
Sub List #2 : AR1016.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1016	6.180	6.543	160050	194613	1130	1060	80.00- 120.00	100.00(M)
	6.723	7.163	333864	326260	1160	1010	162.91- 244.36	208.60(M)
	7.030	7.447	256998	181463	1060	1150	145.03- 217.55	160.57(M)
	7.540	7.623	224481	137090	1080	1050	115.01- 172.51	140.26(M)
	7.837	7.723	198343	162665	1290	1100	76.85- 115.27	123.93(M)
			Average of Peak Amounts =		1140	1070		

QC Flag Legend

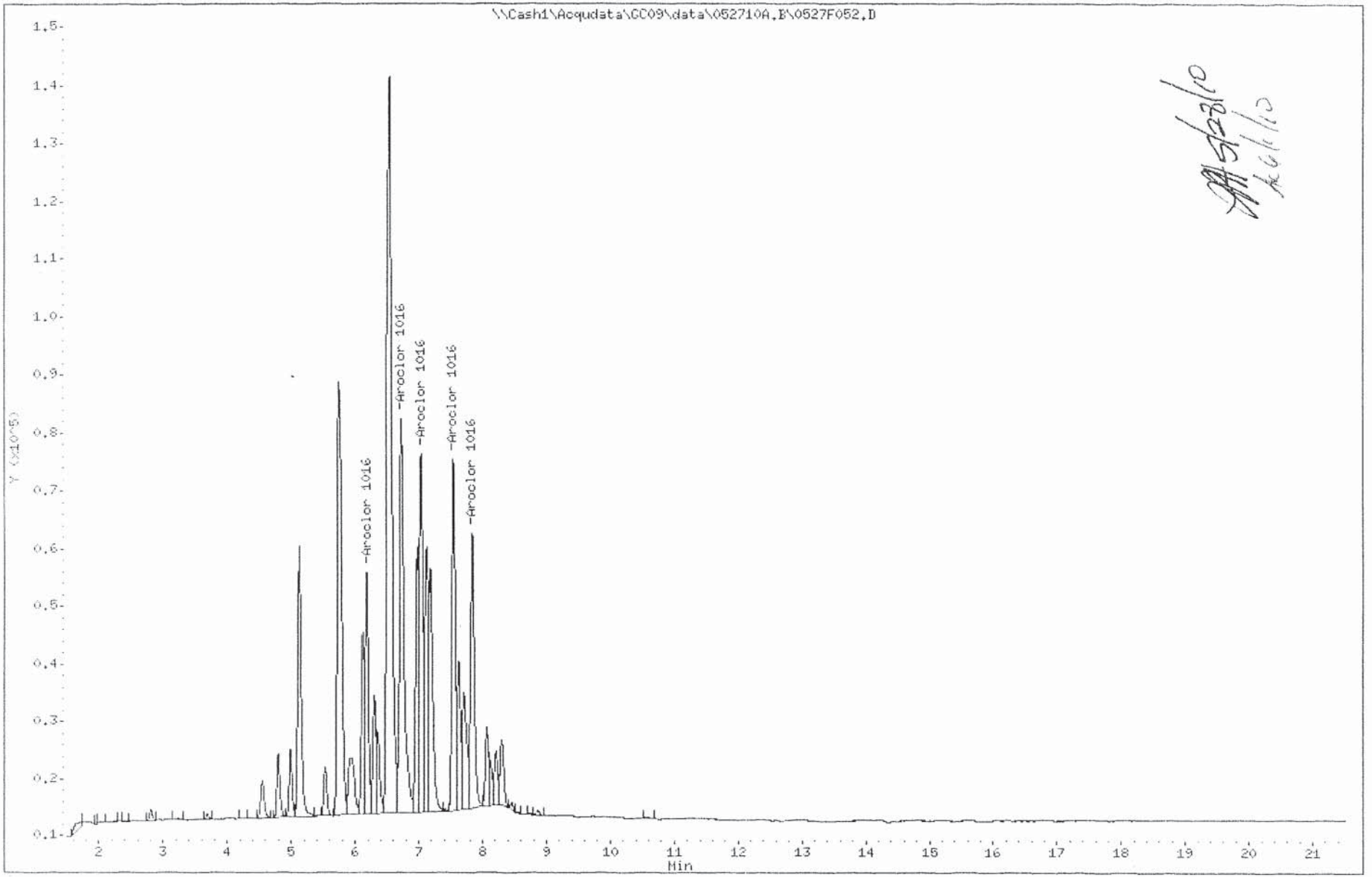
M - Compound response manually integrated.

5/28/10
5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F052.D
Date : 28-MAY-2010 12:55
Client ID:
Sample Info: 1016 @ 1000ppb | PCB5-49E

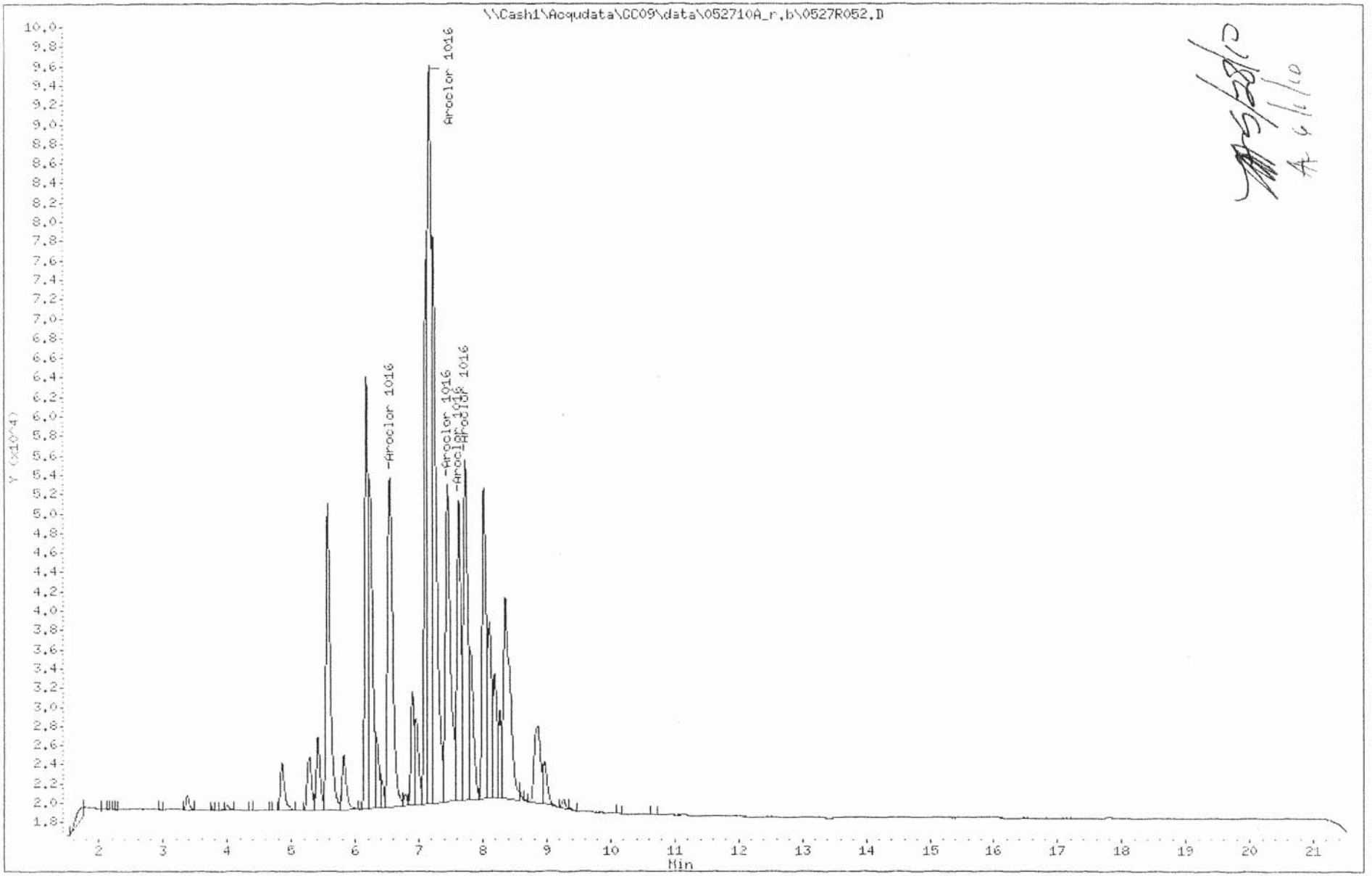
Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

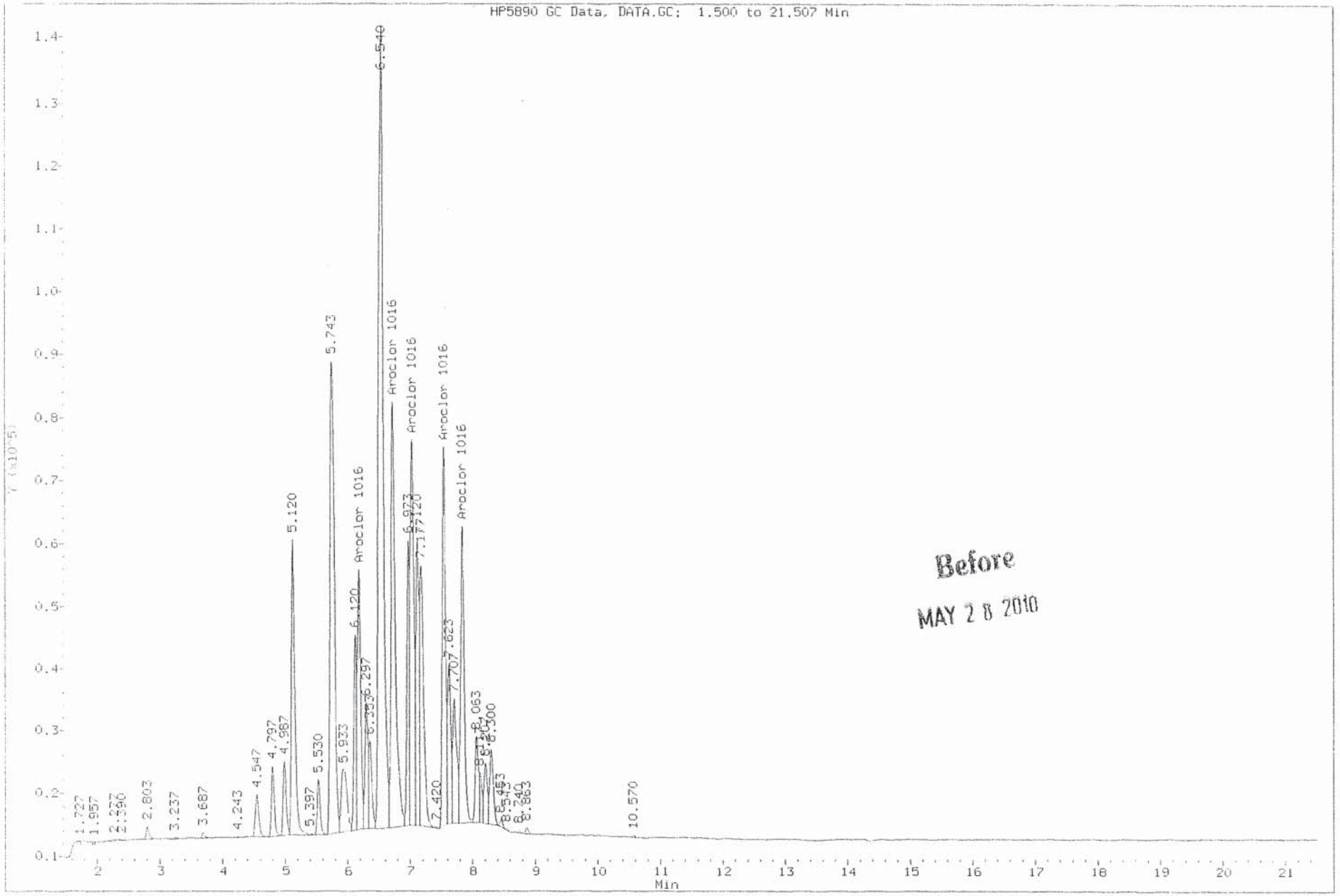


Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R052.D
Date : 28-MAY-2010 12:55
Client ID:
Sample Info: 1016 @ 1000ppb | PCB5-49E

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

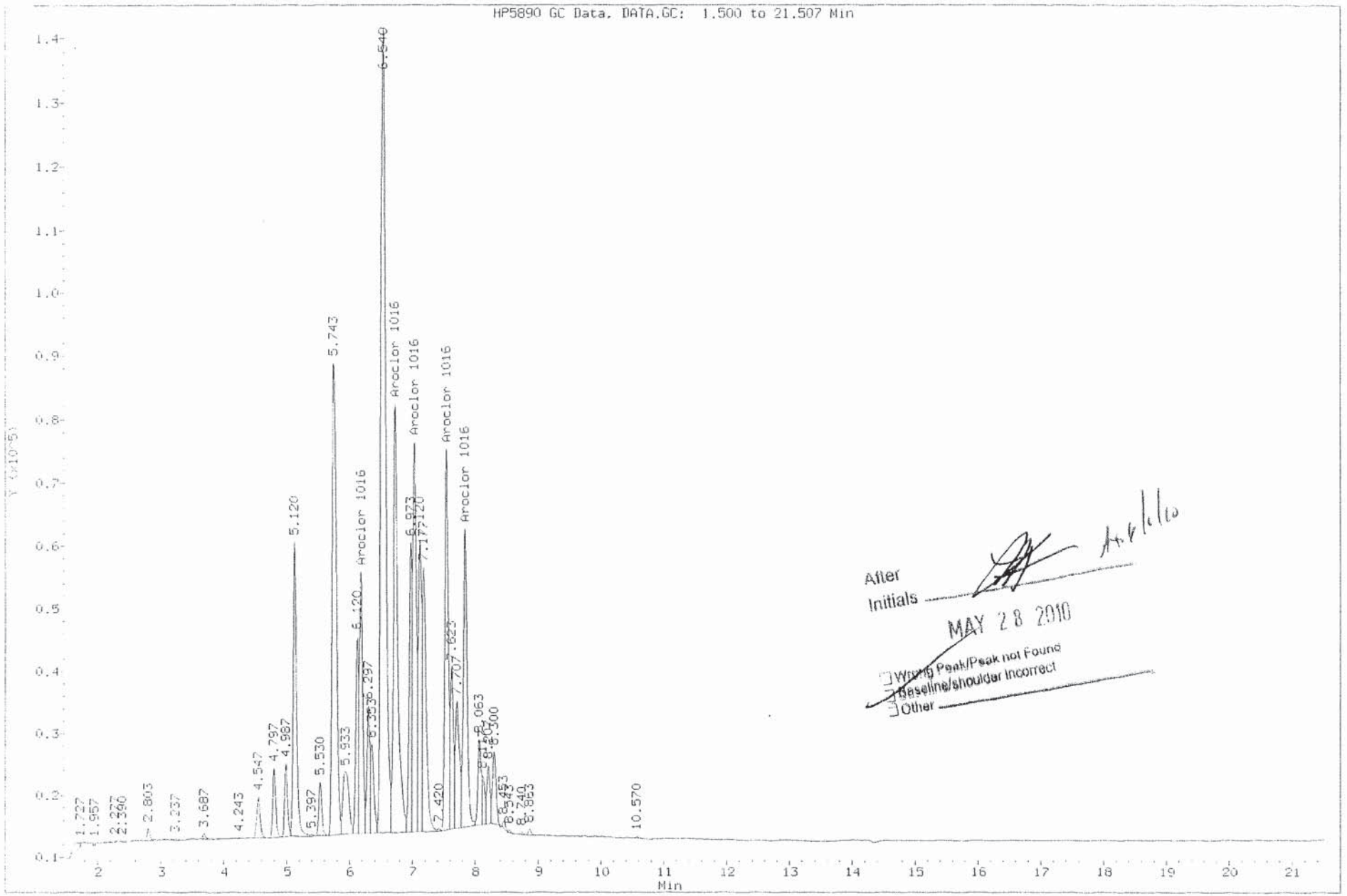


Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F052.D
Injection Date: 28-MAY-2010 12:55
Instrument: GC09.i
Client Sample ID:



Before
MAY 28 2010

Data file: \\Cash1\Acqudata\GC09\data\052710A.B\0527F052.D
Injection Date: 28-MAY-2010 12:55
Instrument: GC09.i
Client Sample ID:

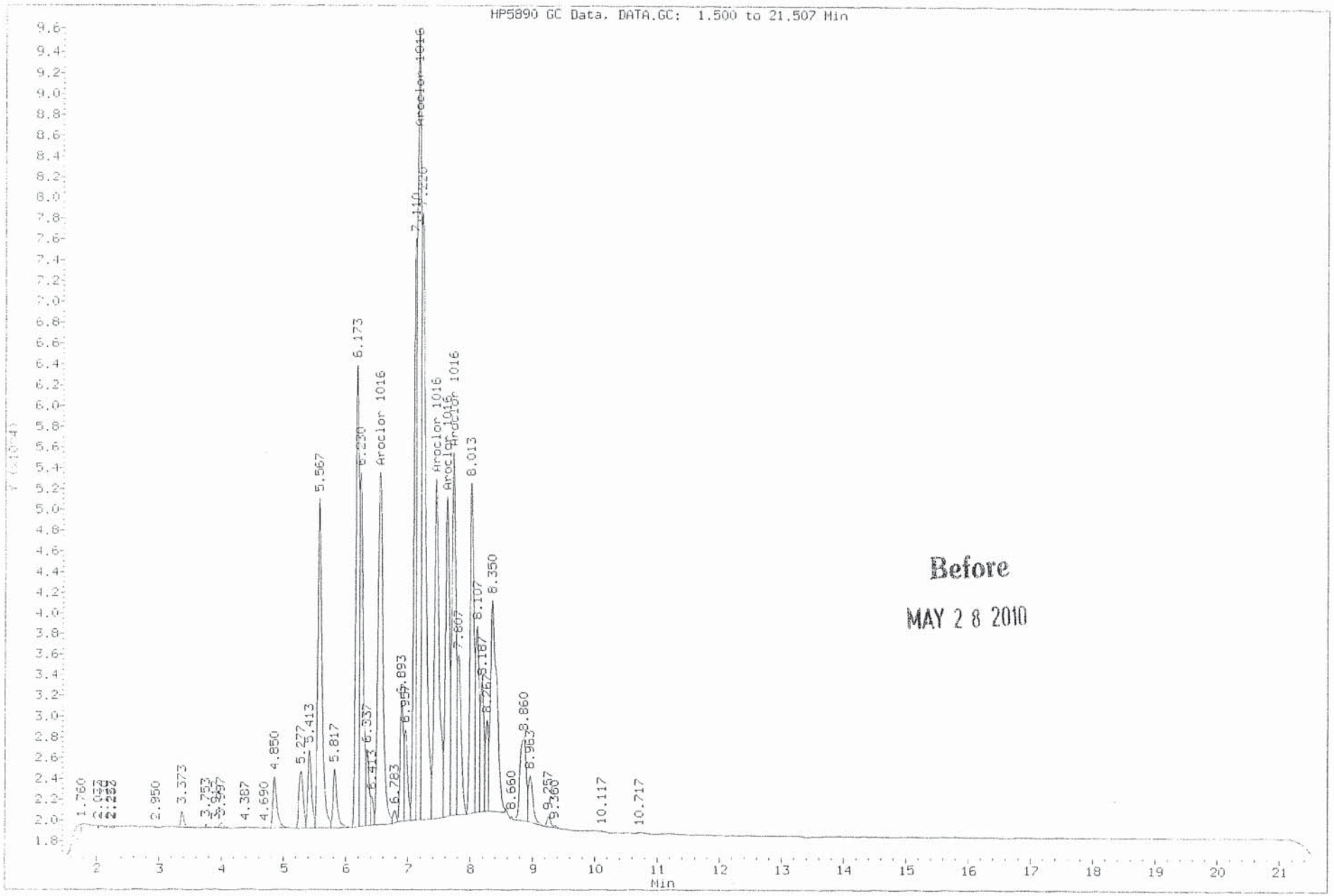


After
Initials

MAY 28 2010

Wrong Peak/Peak not Found
Baseline/shoulder Incorrect
Other

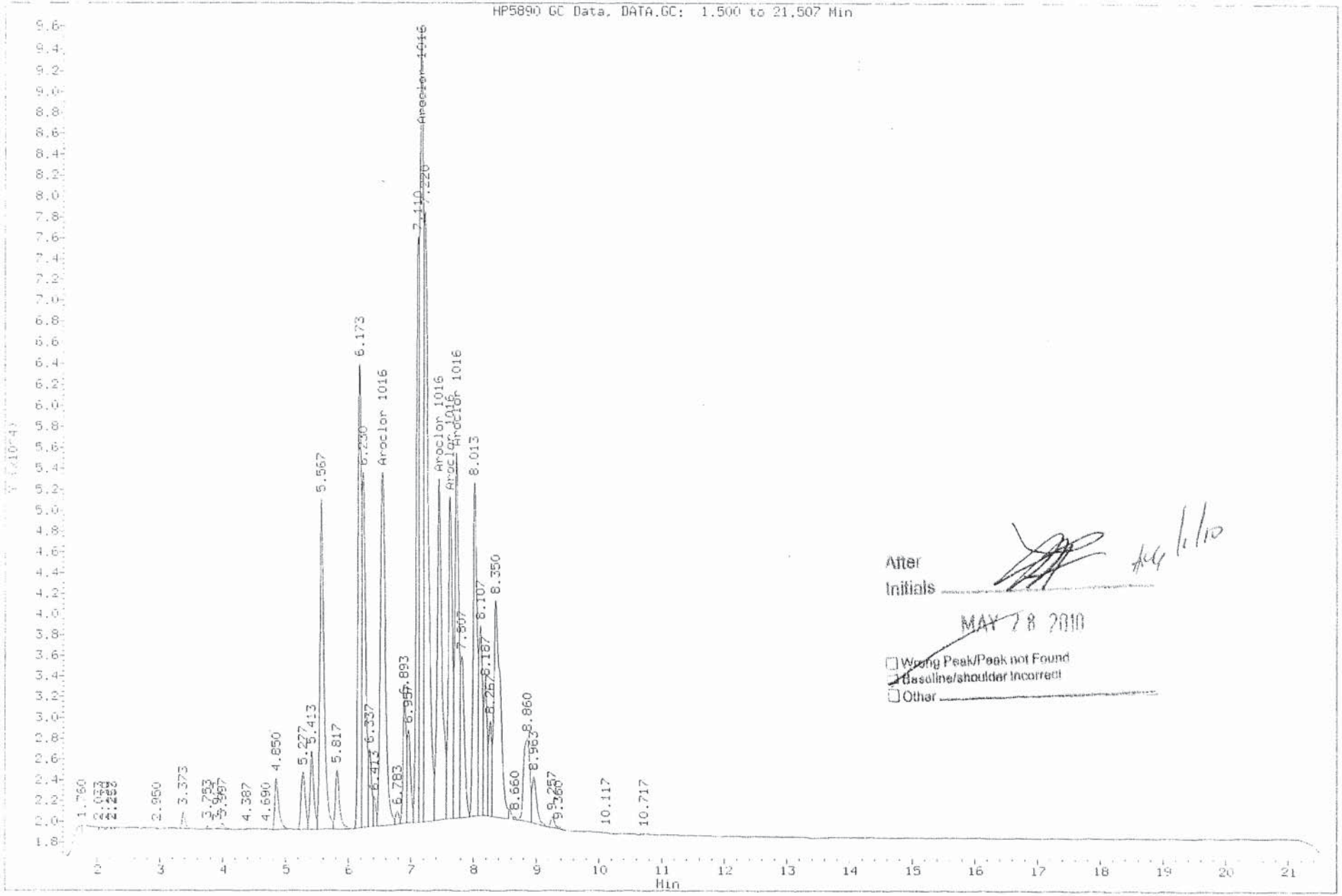
Data File: \\Cash1\acqudata\GC09\data\052710A_r.b\0527R052.D
Injection Date: 28-MAY-2010 12:55
Instrument: GC09.1
Client Sample ID:



Before

MAY 2 8 2010

Data File: \\Cashl\Acqudata\GC09\data\052710A_r.b\0527R052.D
Injection Date: 28-MAY-2010 12:55
Instrument: GC09.1
Client Sample ID:



After
Initials _____ 5/11/10

MAY 28 2010

☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder Incorrect
☐ Other _____

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F053.D
Report Date: 28-May-2010 18:05

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F053.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R053.D
Inj Date : 28-MAY-2010 13:21
Sample Info: 1221 @ 1000ppb | PCB5-53J
Misc Info :
Cal Date : 28-MAY-2010 18:00
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1221.SUB
Sub List #2 : AR1221.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1221	3.687	4.690	66551	19366	925	956	80.00- 120.00	100.00(M)
	4.543	4.850	41087	35793	950	1020	51.26- 76.90	61.74(M)
	4.797	5.287	101432	77078	938	962	133.68- 200.52	152.41(M)
	4.987	5.413	67723	55532	941	992	85.57- 128.35	101.76(M)
			Average of Peak Amounts =		938	982		

QC Flag Legend

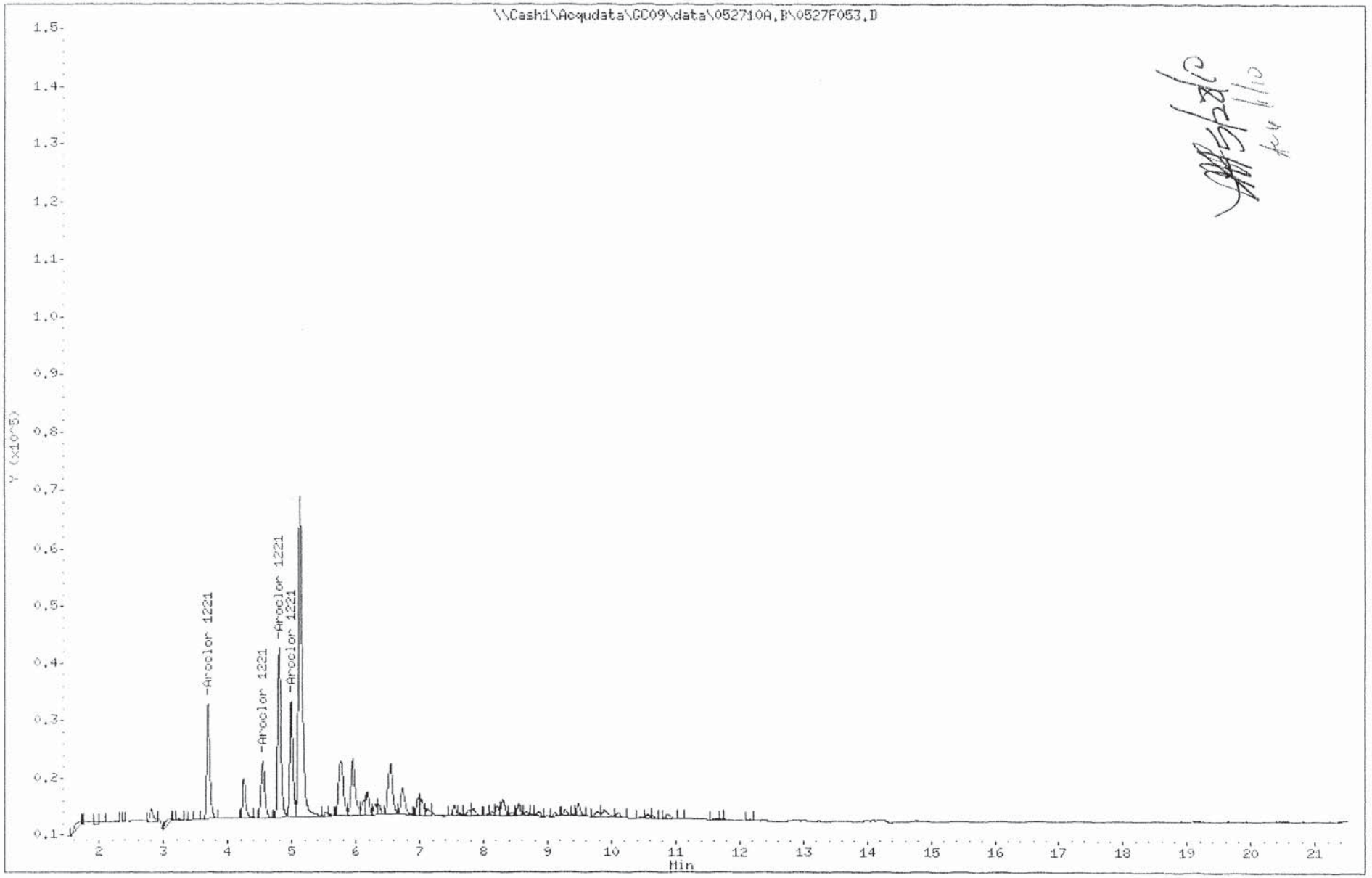
M - Compound response manually integrated.

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F053.D
Date : 28-MAY-2010 13:21
Client ID:
Sample Info: 1221 @ 1000ppb | PCB5-53J

Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



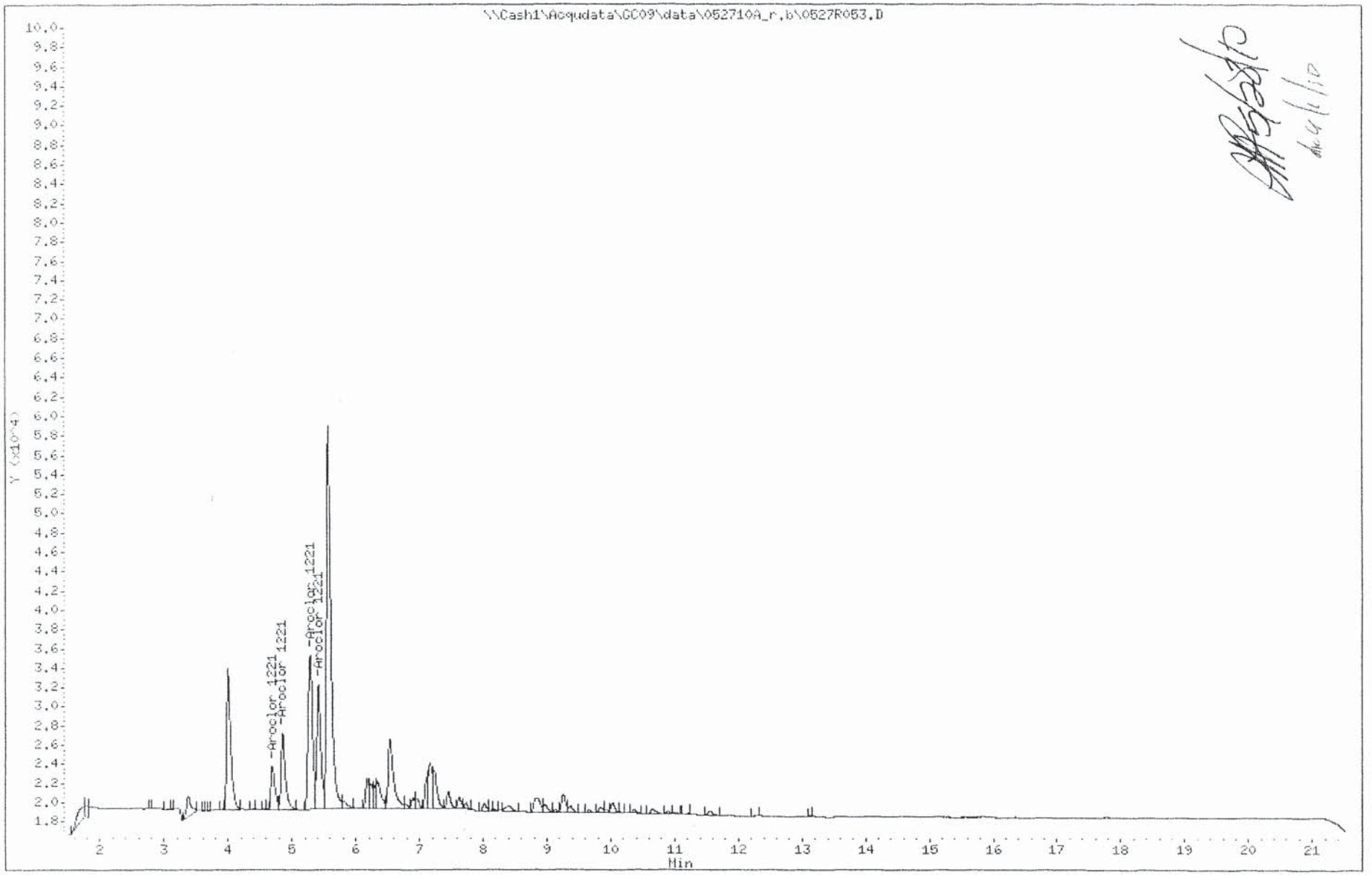
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R053.D
Date : 28-MAY-2010 13:21
Client ID:
Sample Info: 1221 @ 1000ppb | PCB5-53J

Instrument: GC09.i

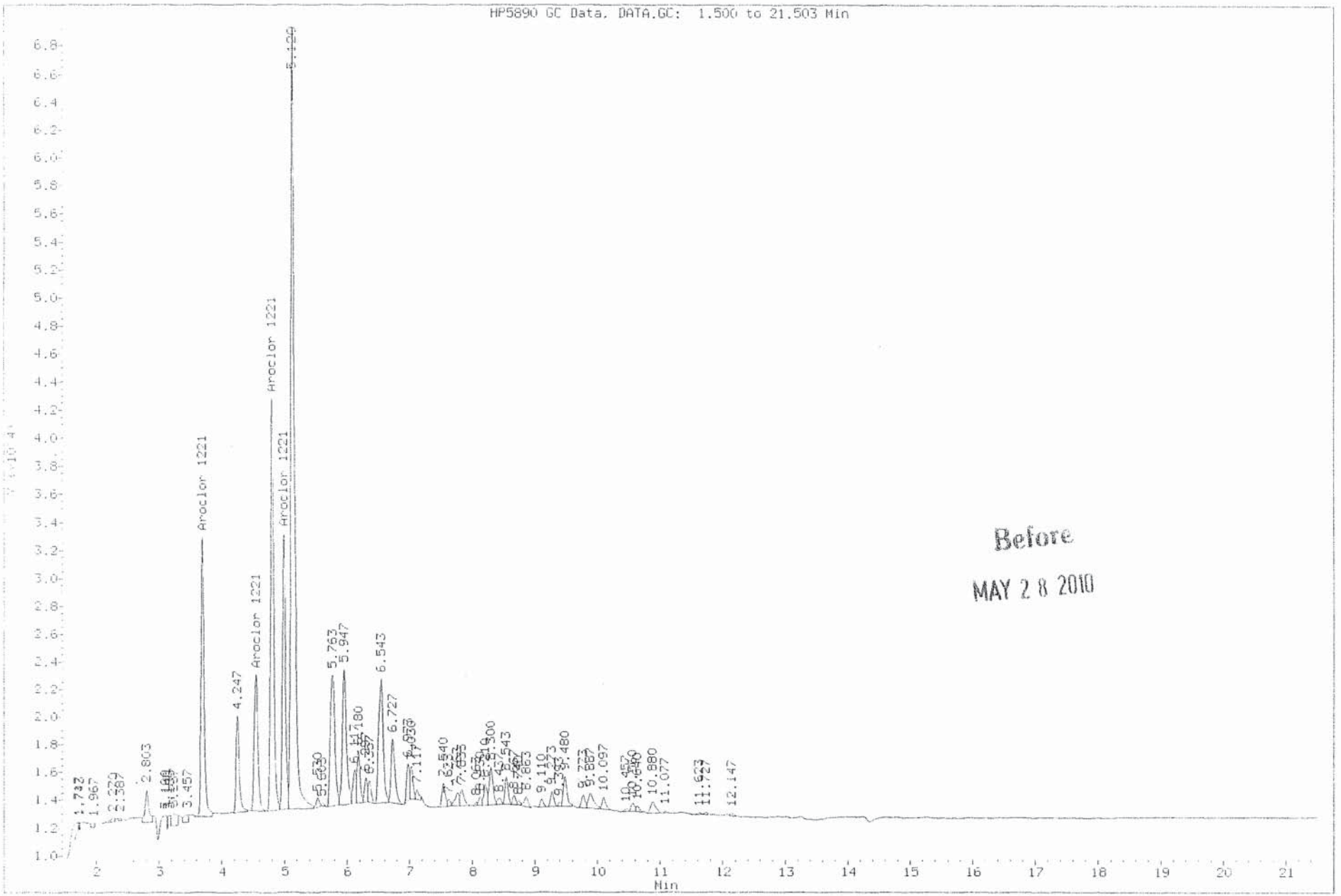
Operator: LHarris

Column phase: DB-XLB

Column diameter: 0.53



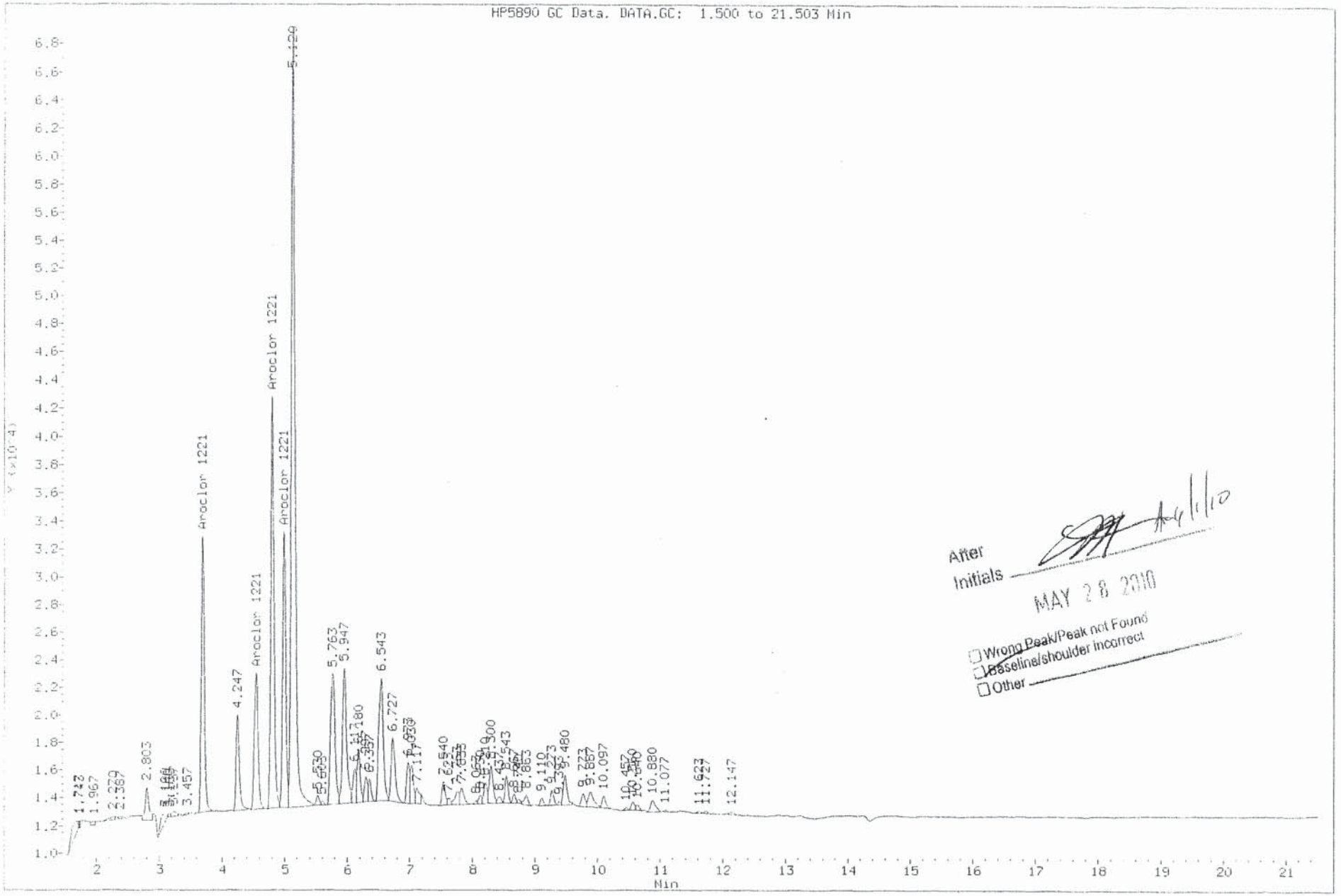
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F053.D
Injection Date: 28-MAY-2010 13:21
Instrument: GC09.1
Client Sample ID:



Before

MAY 2 8 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F053.D
Injection Date: 28-MAY-2010 13:21
Instrument: GC09.1
Client Sample ID:



After
Initials
MAY 28 2010
☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder incorrect
☐ Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F054.D
Report Date: 28-May-2010 18:05

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F054.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R054.D
Inj Date : 28-MAY-2010 13:47
Sample Info: 1232 @ 1000ppb | PCB5-49F
Misc Info :
Cal Date : 28-MAY-2010 18:00
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1232.SUB
Sub List #2 : AR1232.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1232	5.120	5.567	184400	151682	793	861	80.00- 120.00	100.00(M)
	5.747	6.233	178916	60929	889	935	73.60- 110.40	97.03(M)
	6.120	7.223	43415	125482	959	930	17.45- 26.18	23.54(M)
	6.540	7.447	321246	74165	900	974	131.13- 196.70	174.21(M)
			Average of Peak Amounts =		885	925		

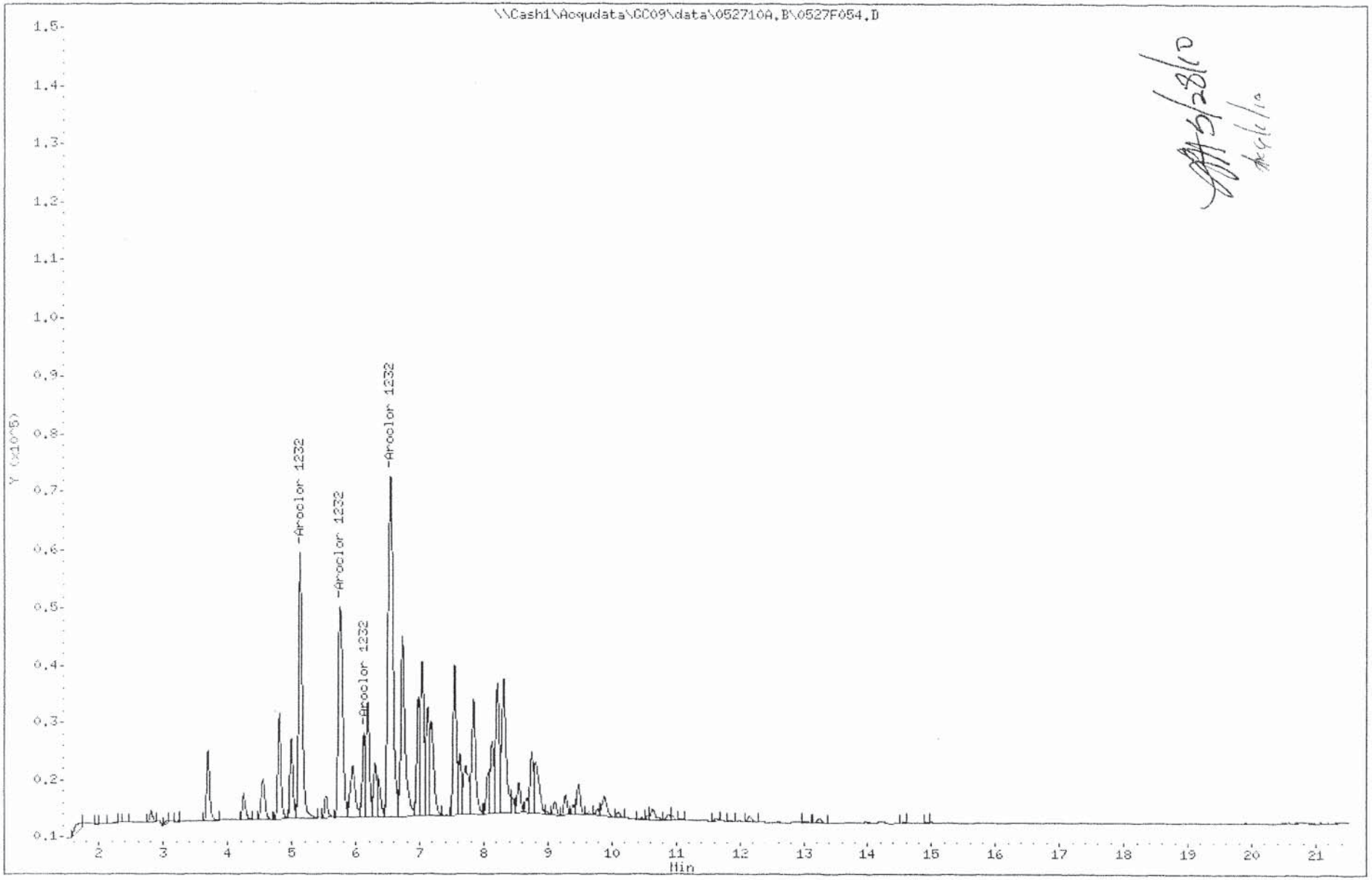
QC Flag Legend

M - Compound response manually integrated.

5/28/10
6/1/10

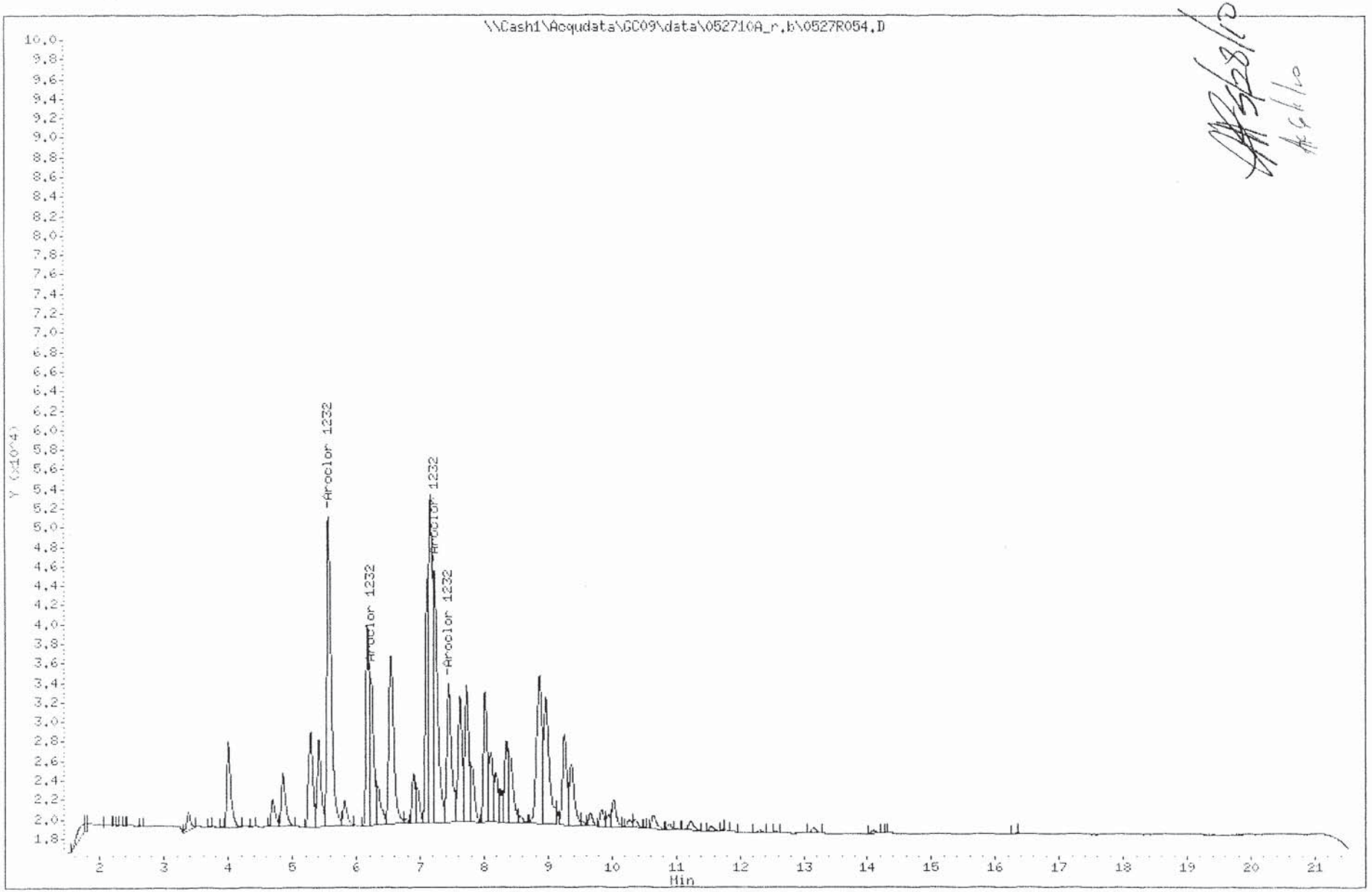
Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F054.D
Date : 28-MAY-2010 13:47
Client ID:
Sample Info: 1232 @ 1000ppb | PCB6-49F

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



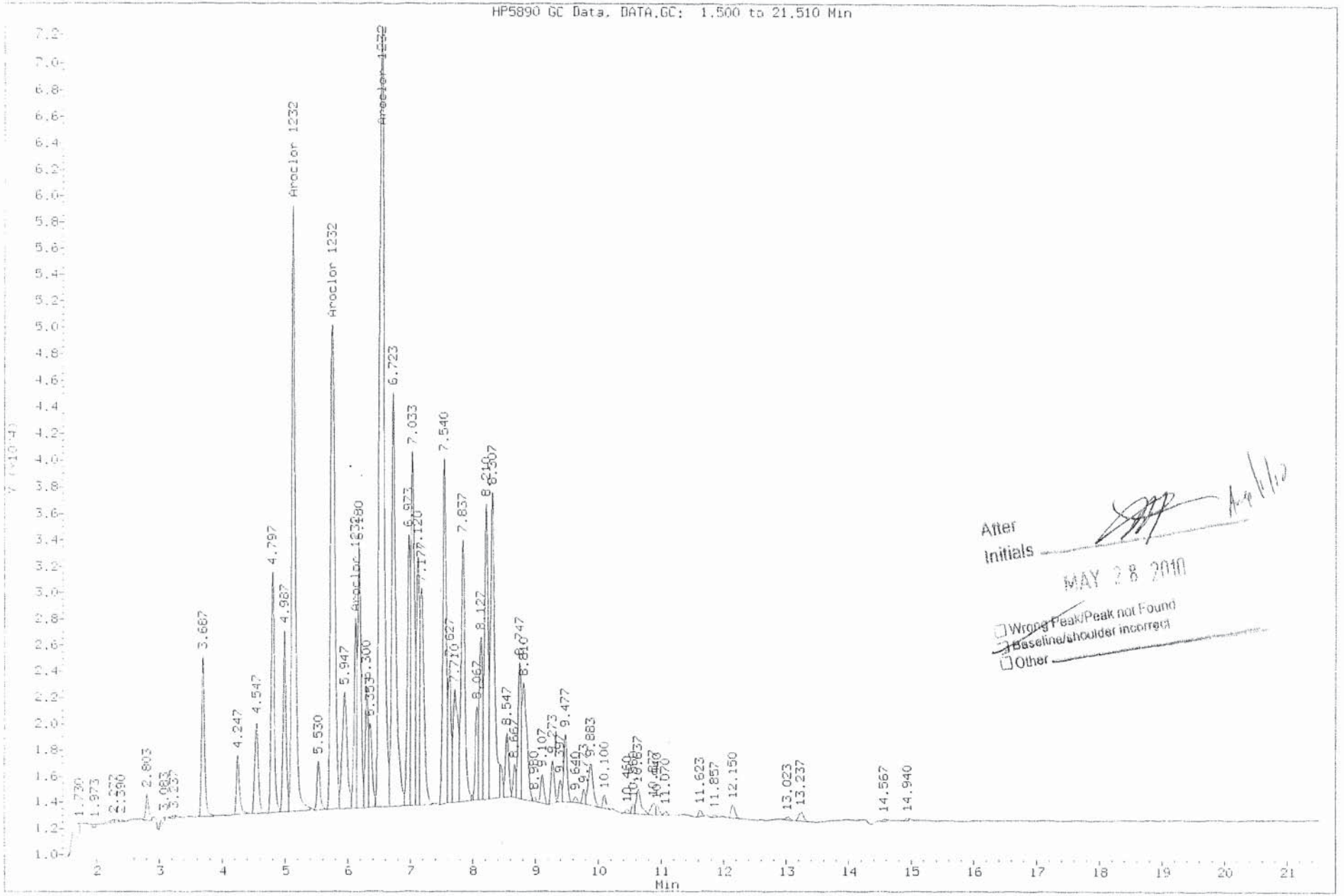
Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R054.D
Date : 28-MAY-2010 13:47
Client ID:
Sample Info: 1232 @ 1000ppb | PCB5-49F

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

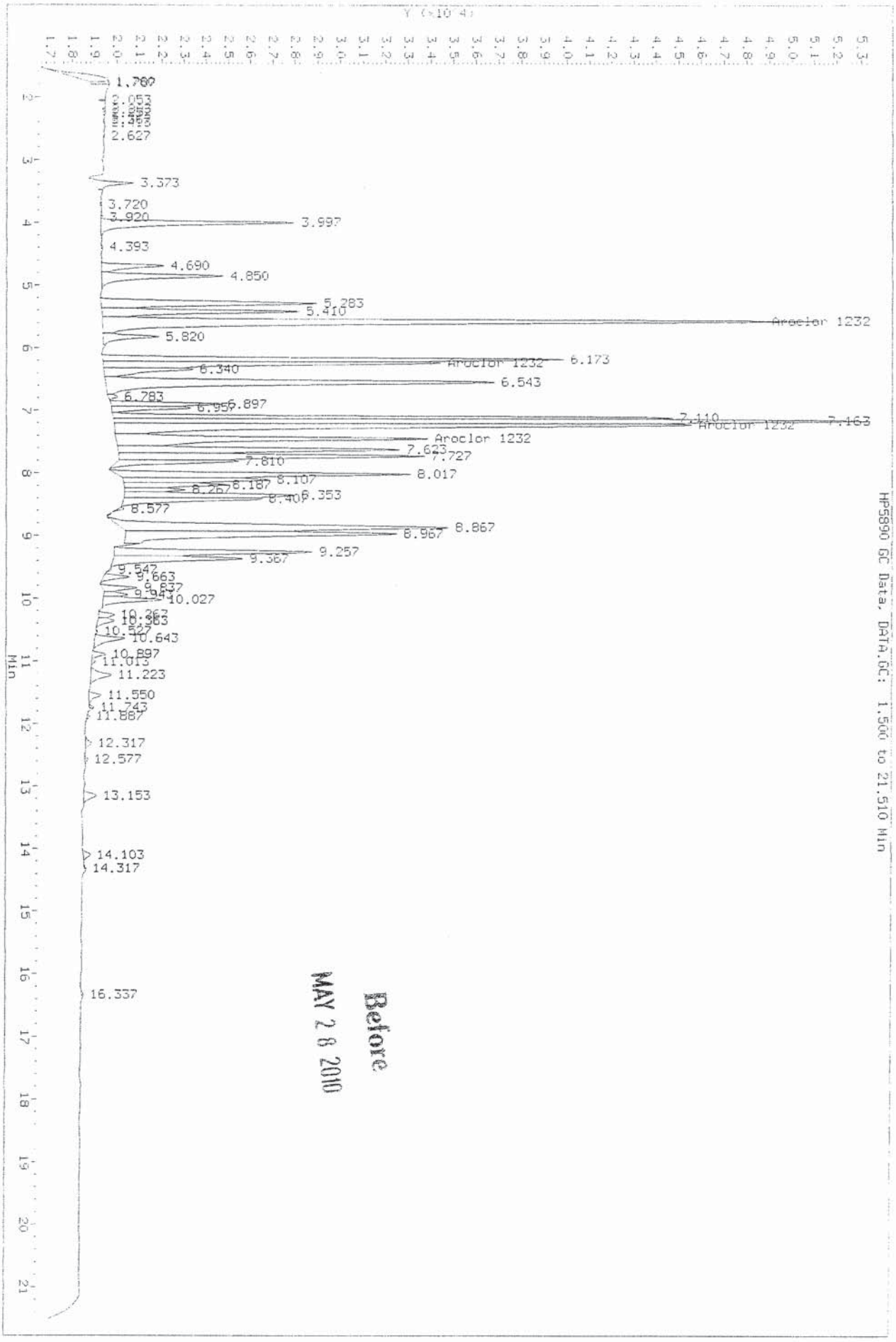


Data file: \\Esh1\hcqudata\GC09\data\052710A.B\0527F054.D
Injection Date: 28-MAY-2010 13:47
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Min
1.730
1.973
2.390
2.803
3.083
3.237
3.687
4.247
4.547
4.797
4.987
Aroclor 1232
5.530
Aroclor 1232
5.947
Aroclor 1
6.480
6.725
6.973
7.033
7.540
7.627
7.837
8.067
8.127
8.210
8.307
8.547
8.662
8.747
8.960
9.107
9.273
9.477
9.640
9.883
10.100
10.460
10.837
11.070
11.623
11.857
12.150
13.023
13.237
14.567
14.940
Before
MAY 28 2010

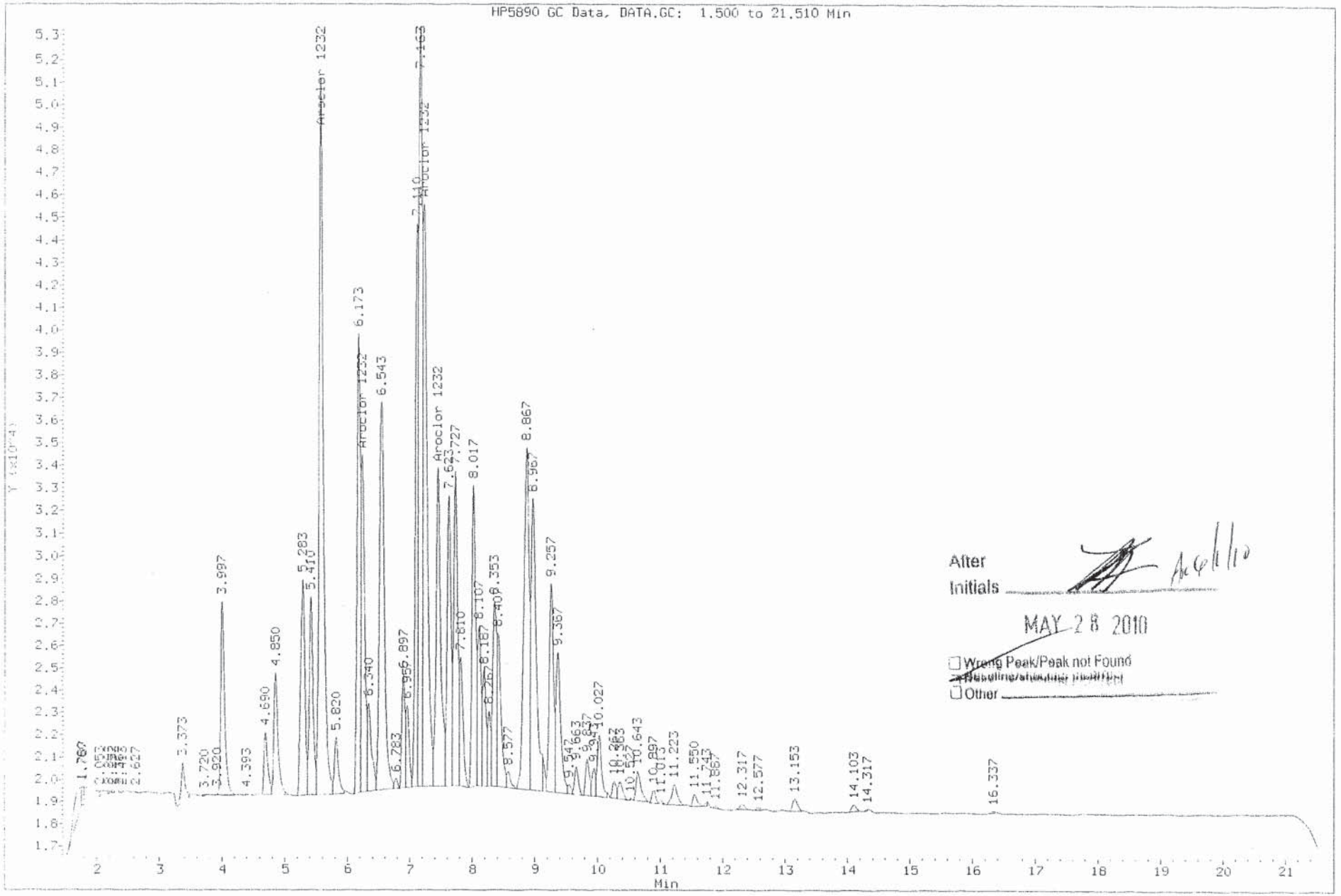
Data File: \\Cashl\Acqudata\GC09\data\052710A.B\0527F054.D
Injection Date: 28-MAY-2010 13:47
Instrument: GC09.i
Client Sample ID:



After
Initials
MAY 28 2010
☐ Wrong Peak/Peak not Found
☐ Baseline/shoulder incorrect
☐ Other


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R054.D
Injection Date: 28-MAY-2010 13:47
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.510 Min
Y (x10^4)
Min
1.780
2.627
3.373
3.720
3.920
3.997
4.393
4.690
4.850
5.283
5.410
Aroclor 1232
5.820
Aroclor 1232
6.173
6.340
6.543
6.783
6.956
6.897
7.110
Aroclor 1232
7.163
Aroclor 1232
7.623
7.727
7.810
8.017
8.107
8.187
8.267
8.353
8.577
8.867
8.967
9.257
9.367
9.542
9.663
9.837
10.027
10.643
10.897
11.015
11.223
11.550
11.743
11.887
12.317
12.577
13.153
14.103
14.317
16.337
Before
MAY 28 2010

Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R054.D
Injection Date: 28-MAY-2010 13:47
Instrument: GC09.i
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F055.D
Report Date: 28-May-2010 17:30

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F055.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R055.D
Inj Date : 28-MAY-2010 14:13
Sample Info: 1242 @ 1000ppb | PCB5-49G
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1242.SUB
Sub List #2 : AR1242.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1242	6.120	6.230	73732	104456	962	954	80.00- 120.00	100.00(M)
	6.540	6.543	550636	153844	902	955	727.90-1091.85	746.80(M)
	7.120	7.110	132541	121334	969	889	150.61- 225.91	179.76(M)
	7.540	7.623	181946	116852	918	1040	230.50- 345.75	246.77(M)
	7.837	7.727	174489	134745	936	1040	198.70- 298.04	236.65(M)
			Average of Peak Amounts =		937	976		

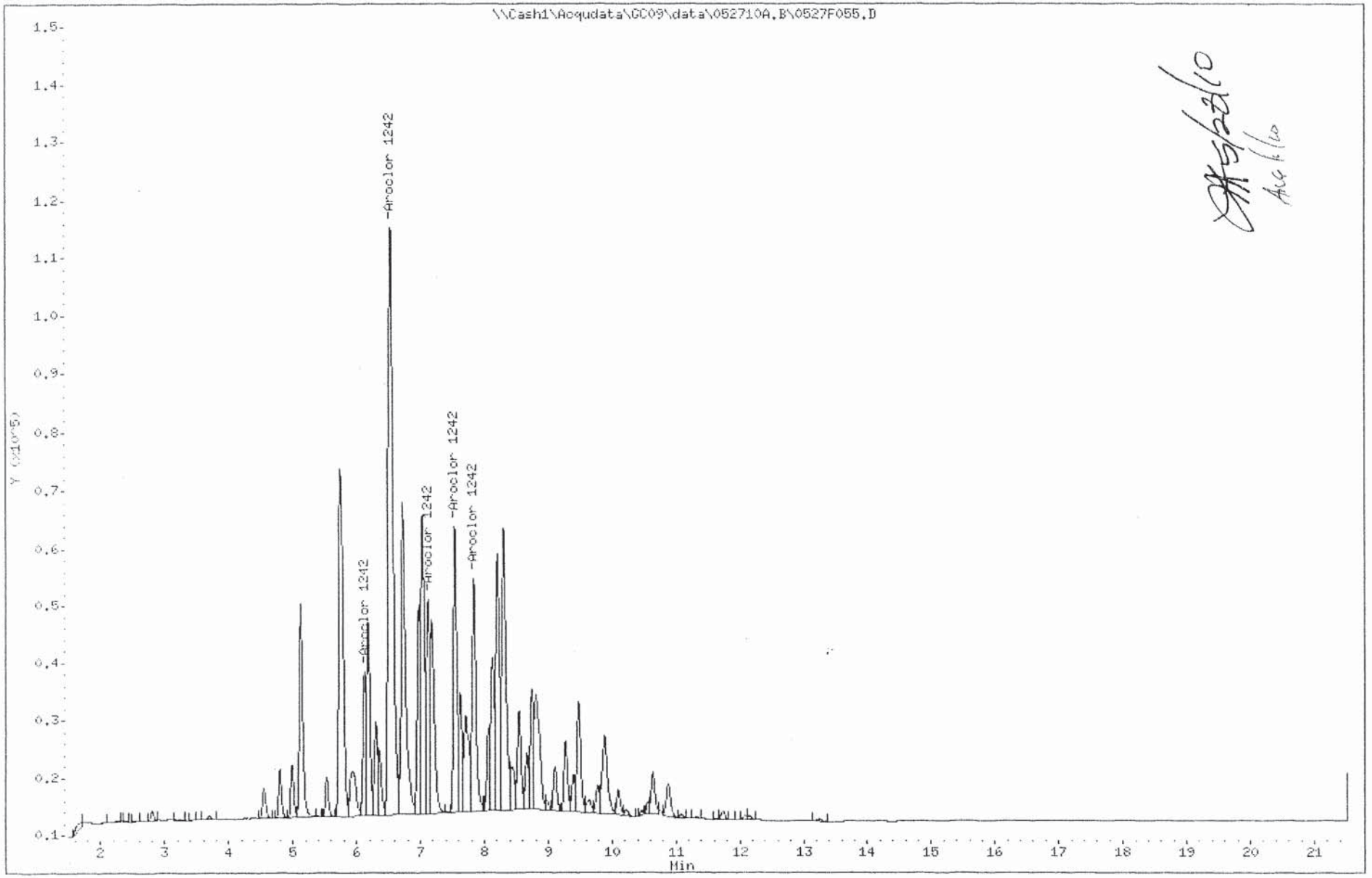
QC Flag Legend

M - Compound response manually integrated.

5/28/10
5/28/10

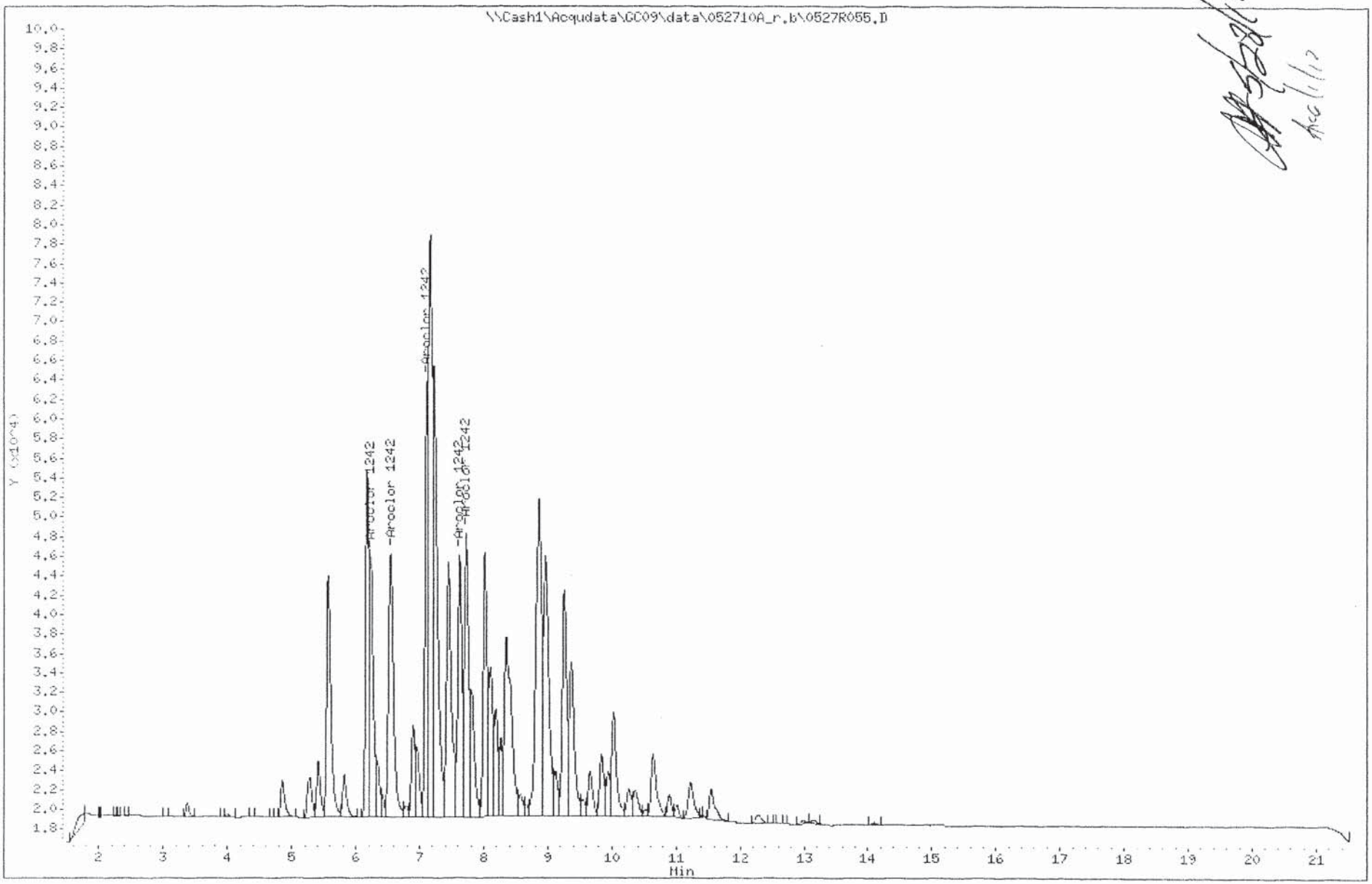
Data File: \\Cash1\Aoqudata\GC09\data\052710A.B\0527F055.D
Date : 28-MAY-2010 14:13
Client ID:
Sample Info: 1242 @ 1000ppb I PCB5-49G

Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

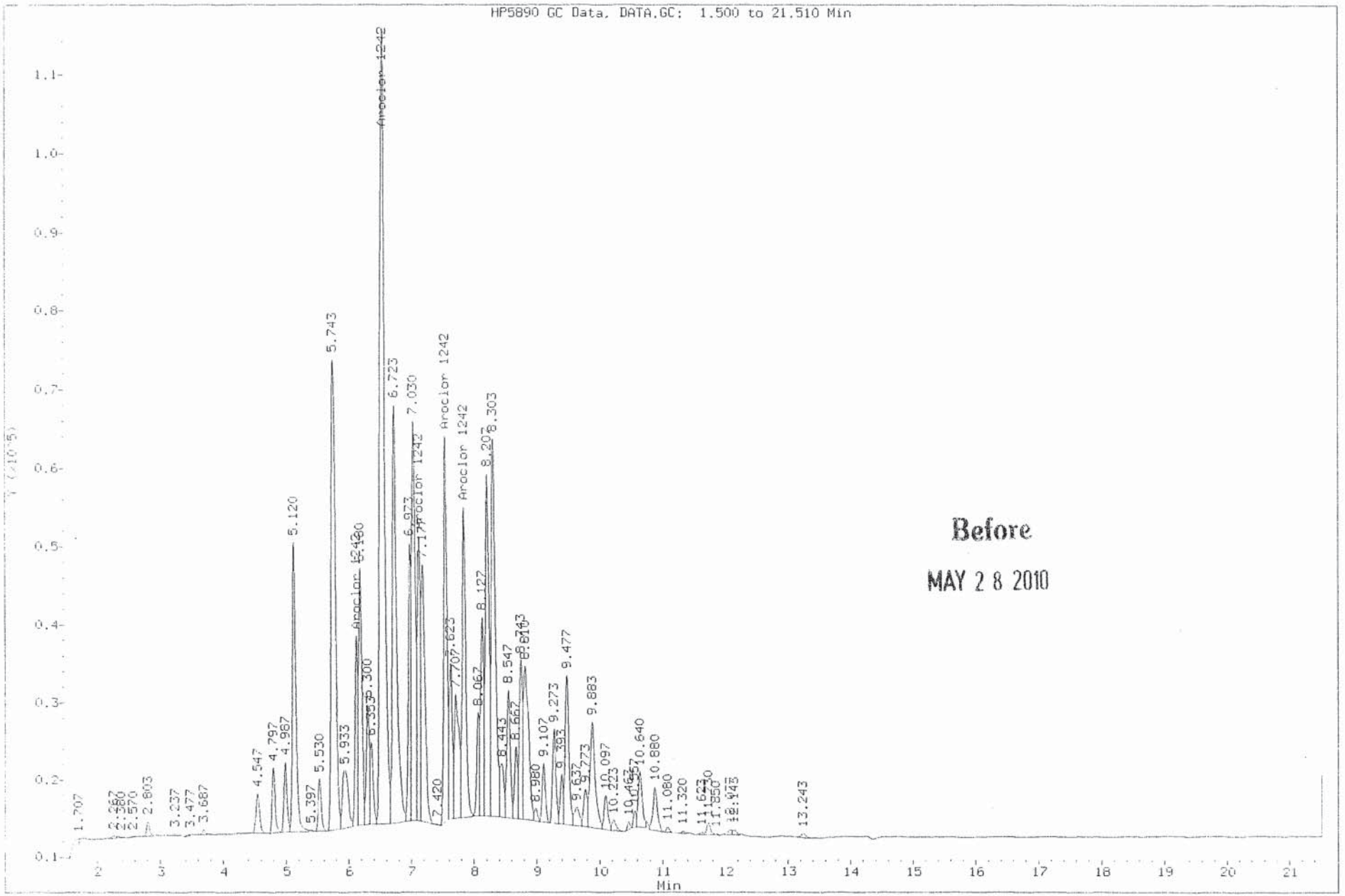


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R055.D
Date : 28-MAY-2010 14:13
Client ID:
Sample Info: 1242 @ 1000ppb | PCB5-49G

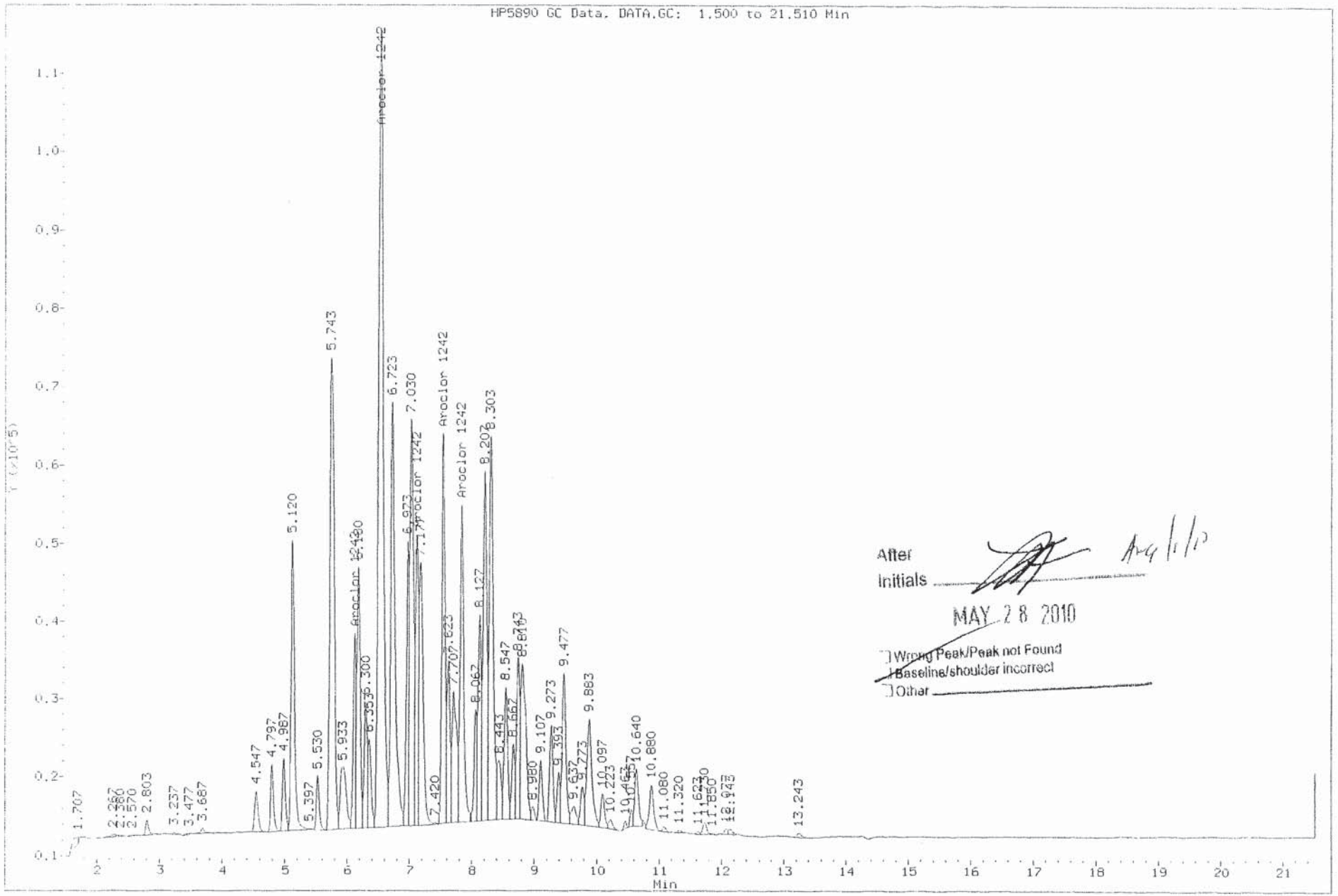
Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F055.D
Injection Date: 28-MAY-2010 14:13
Instrument: GC09.i
Client Sample ID:



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F055.D
Injection Date: 28-MAY-2010 14:13
Instrument: GC09.i
Client Sample ID:



After
Initials

MAY 28 2010

[ ] Wrong Peak/Peak not Found
[ ] Baseline/shoulder incorrect
[ ] Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F056.D
Report Date: 28-May-2010 17:31

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F056.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R056.D
Inj Date : 28-MAY-2010 14:39
Sample Info: 1248 @ 1000ppb | PCB5-49H
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1248.SUB
Sub List #2 : AR1248.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1248	6.537	7.623	373368	178975	906	895	80.00- 120.00	100.00(M)
	7.177	7.727	191888	181363	914	912	39.24- 58.85	51.39(M)
	7.837	8.013	227914	181597	908	908	43.38- 65.07	61.04(M)
	8.207	8.863	302327	331516	870	874	66.78- 100.17	80.97(M)
	8.303	8.963	373583	242584	881	882	84.10- 126.14	100.06(M)
			Average of Peak Amounts =		896	894		

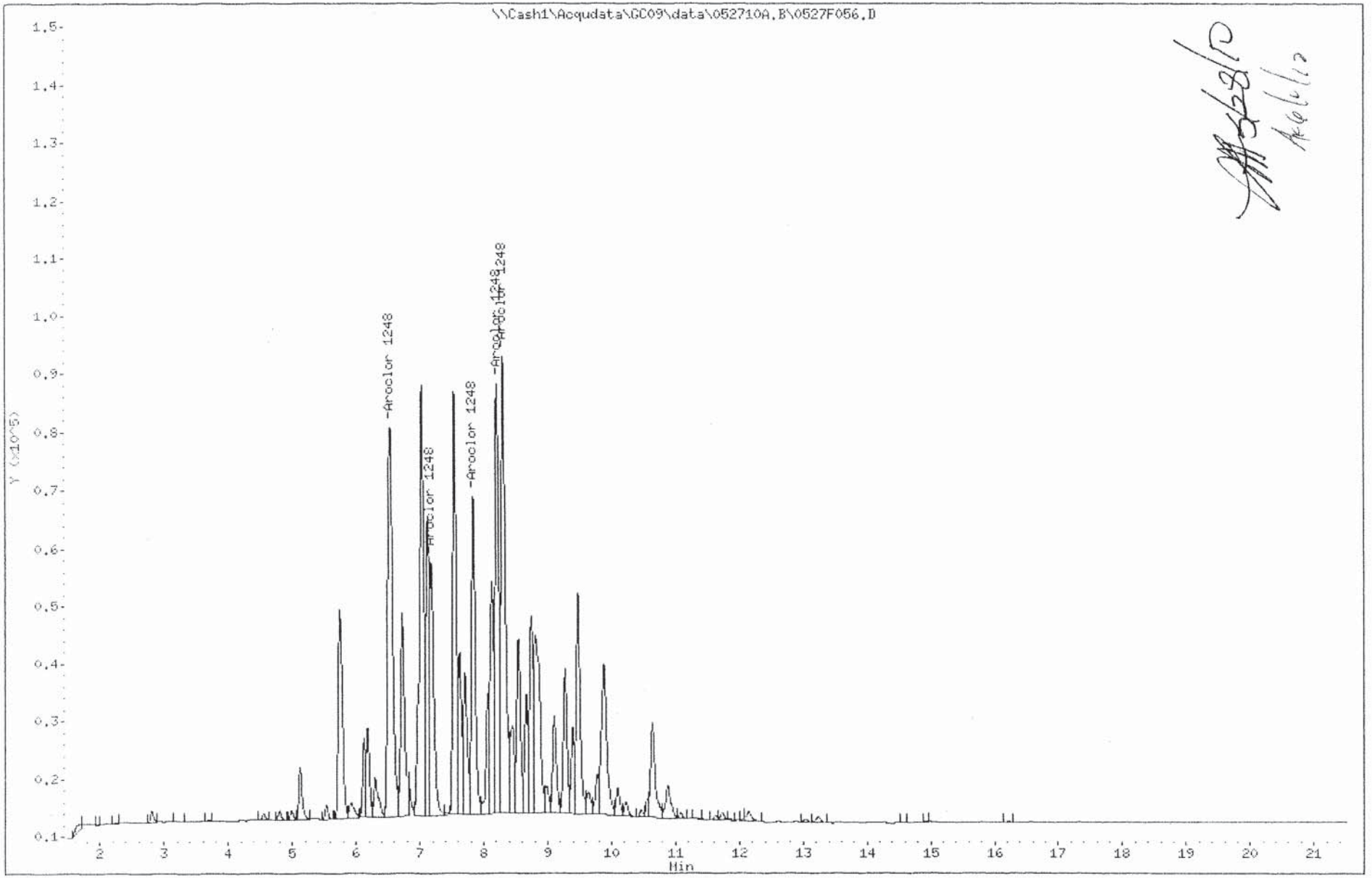
QC Flag Legend

M - Compound response manually integrated.

5/28/10

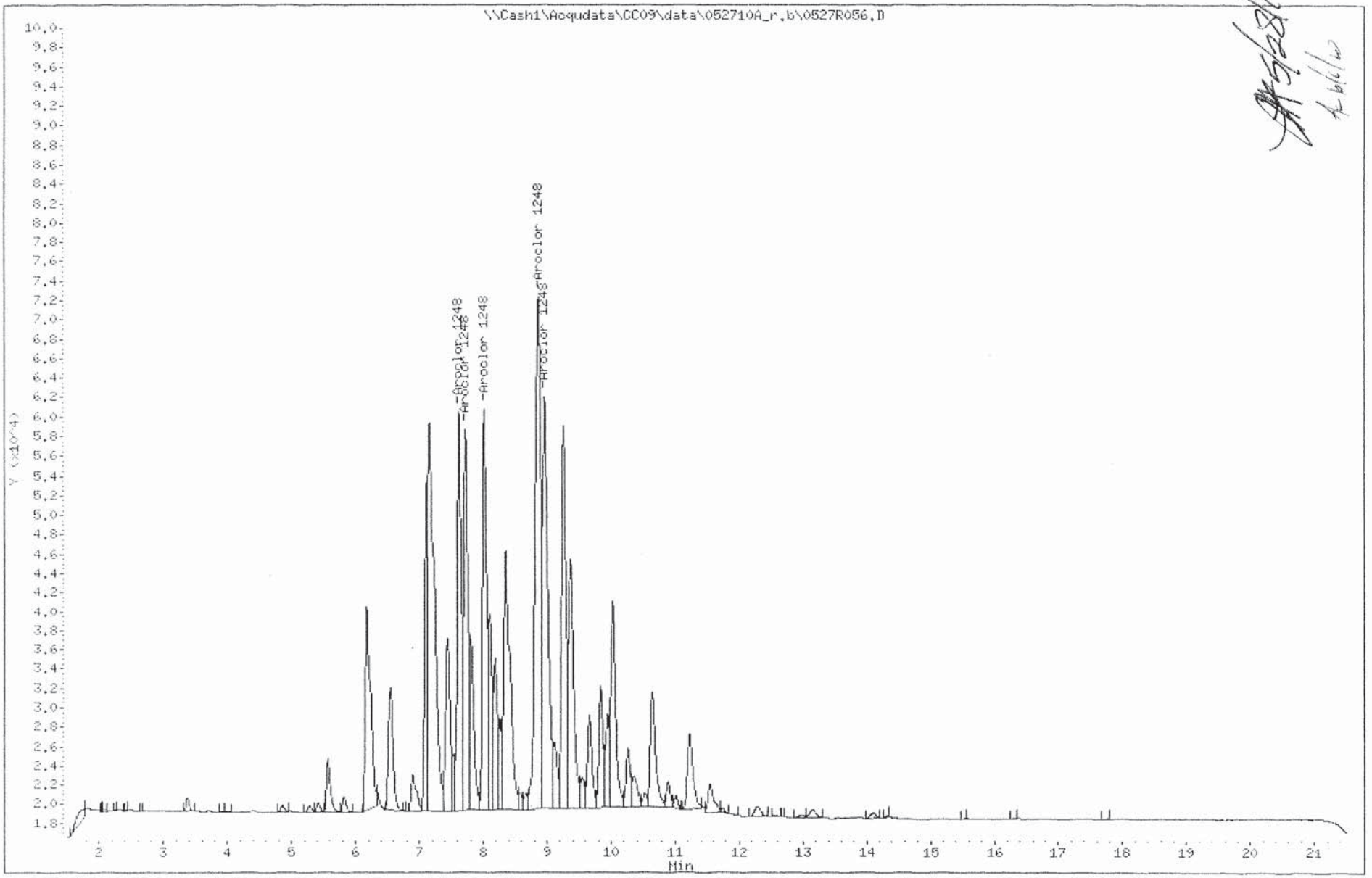
Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F056.D
Date : 28-MAY-2010 14:39
Client ID:
Sample Info: 1248 @ 1000ppb | PCB5-49H

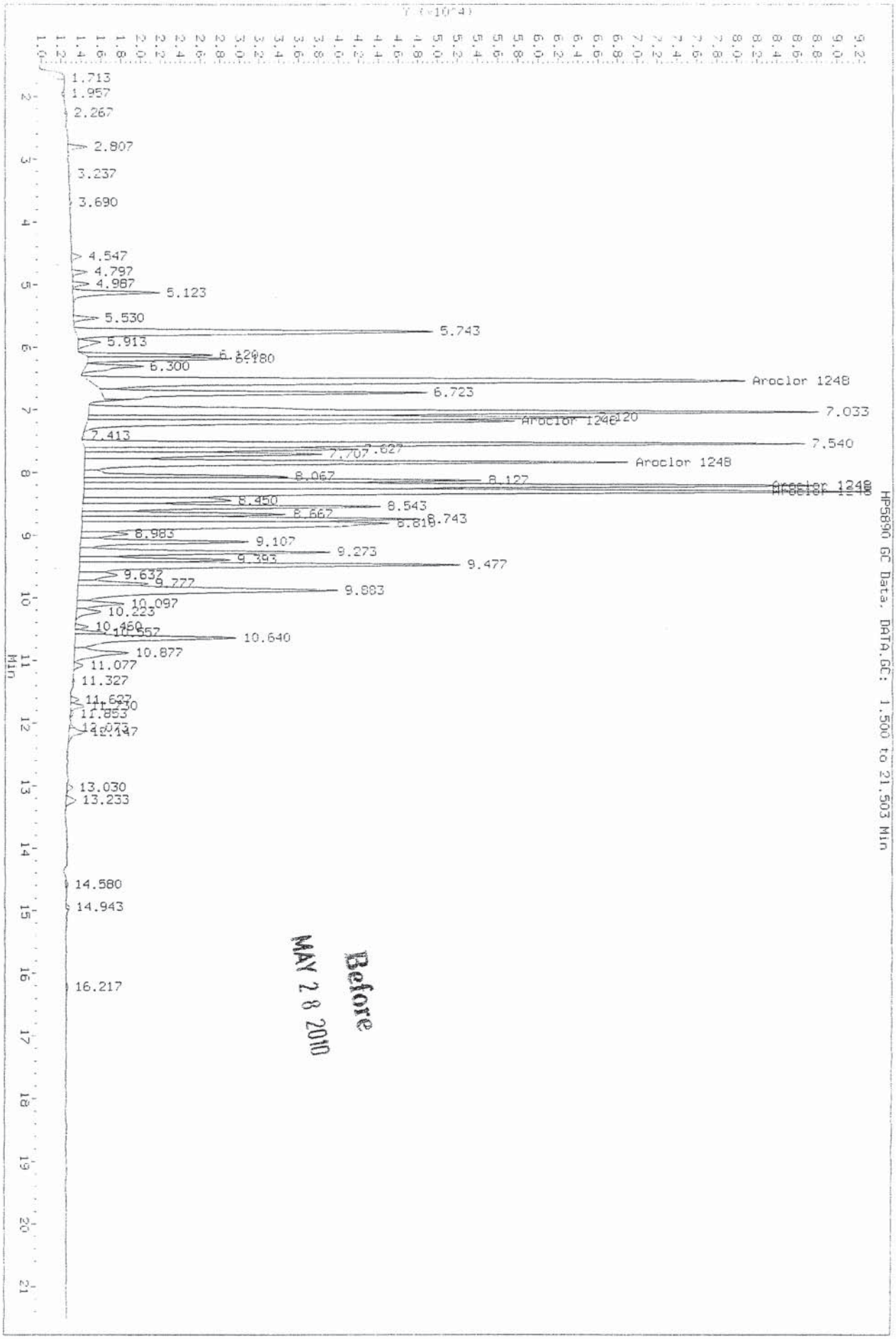
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



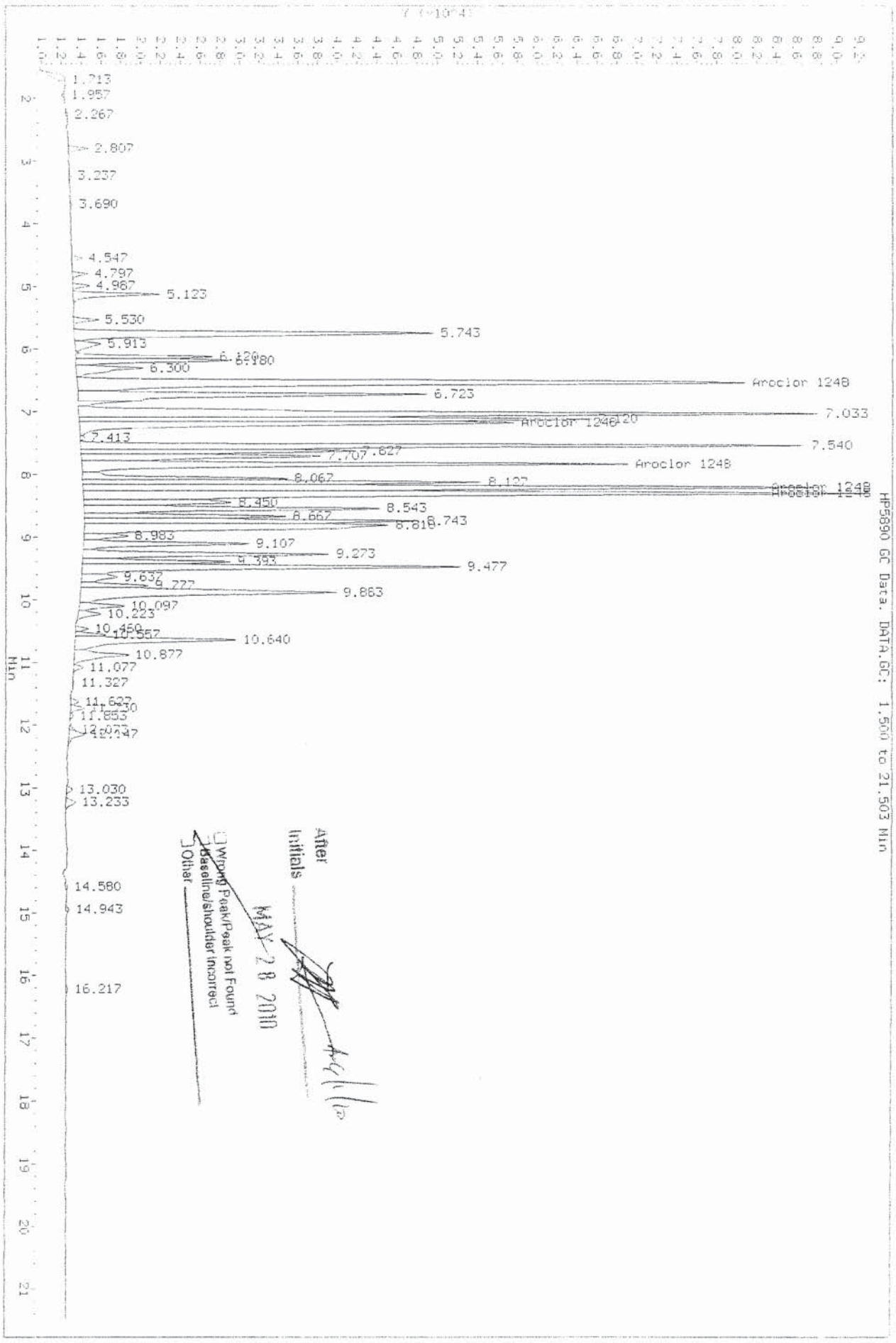
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R056.D
Date : 28-MAY-2010 14:39
Client ID:
Sample Info: 1248 @ 1000ppb | PCB5-49H

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53





Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052704.B\052706.D
Injection Date: 28-MAY-2010 14:39
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.503 Min
Min
1.713
1.957
2.267
2.807
3.237
3.690
4.547
4.797
4.987
5.123
5.530
5.743
5.913
6.120
6.180
6.300
Aroclor 1248
6.723
7.033
Aroclor 1248-20
2.413
7.540
7.707
7.627
Aroclor 1248
8.067
8.127
Aroclor 1248
8.450
8.543
8.667
8.810
8.743
8.983
9.107
9.273
9.393
9.477
9.637
9.727
9.863
10.097
10.223
10.450
10.557
10.640
10.877
11.077
11.327
11.627
11.730
11.853
12.073
12.147
13.030
13.233
14.580
14.943
16.217
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F057.D
Report Date: 28-May-2010 17:31

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F057.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R057.D
Inj Date : 28-MAY-2010 15:05
Sample Info: 1254 @ 1000ppb | PCB5-53K
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

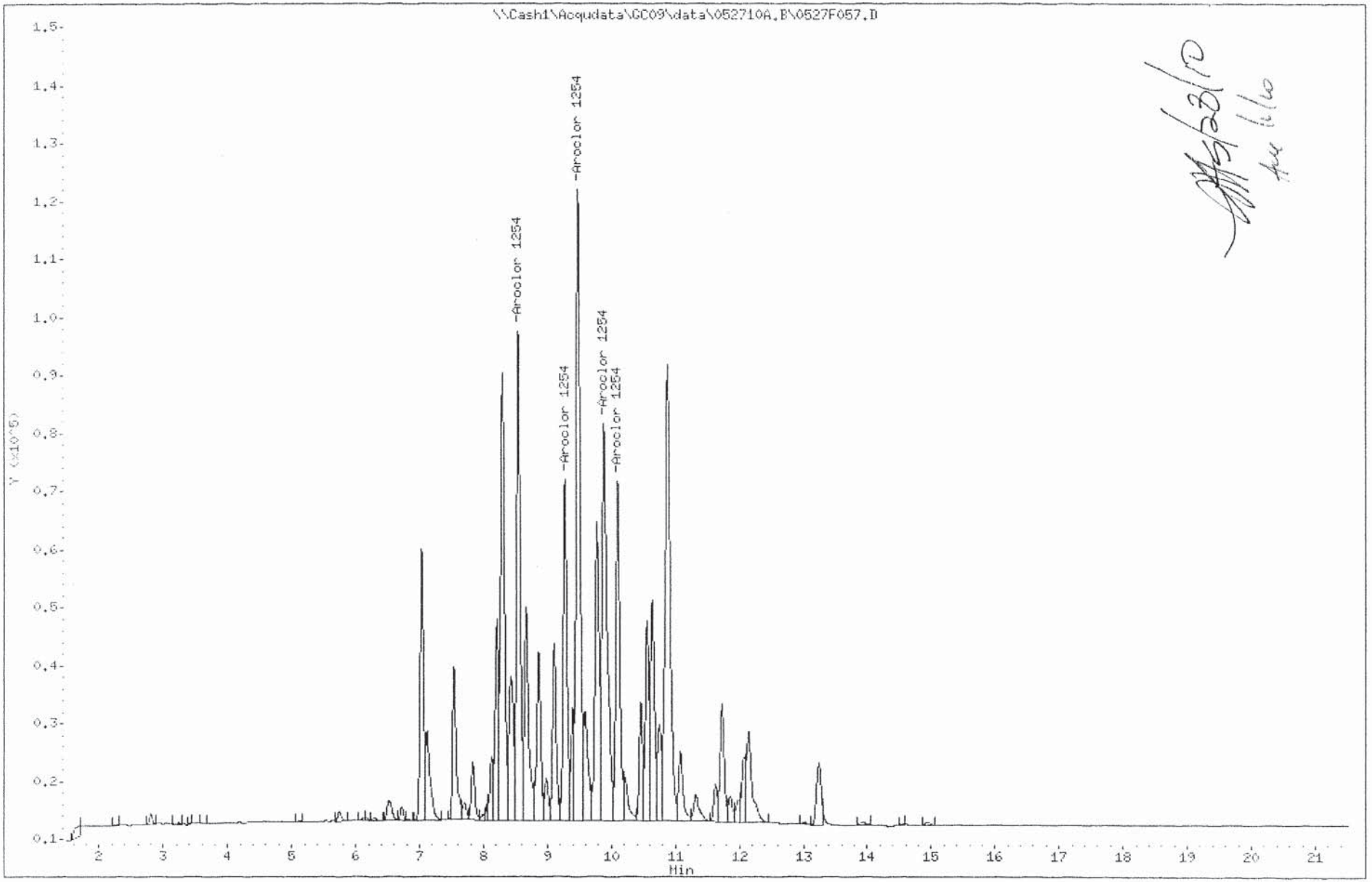
Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1254.SUB
Sub List #2 : AR1254.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1254	8.543	9.250	338913	303655	839	790	80.00- 120.00	100.00
	9.273	9.833	238260	146943	849	834	61.91- 92.86	70.30
	9.477	10.023	473764	295305	825	826	116.61- 174.92	139.79
	9.883	10.367	399146	159259	858	990	101.15- 151.73	117.77
	10.097	10.637	248633	259604	1000	826	53.30- 79.94	73.36
			Average of Peak Amounts =		874	853		

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F057.D
Date : 28-MAY-2010 15:05
Client ID:
Sample Info: 1254 @ 1000ppb | PCB5-53K

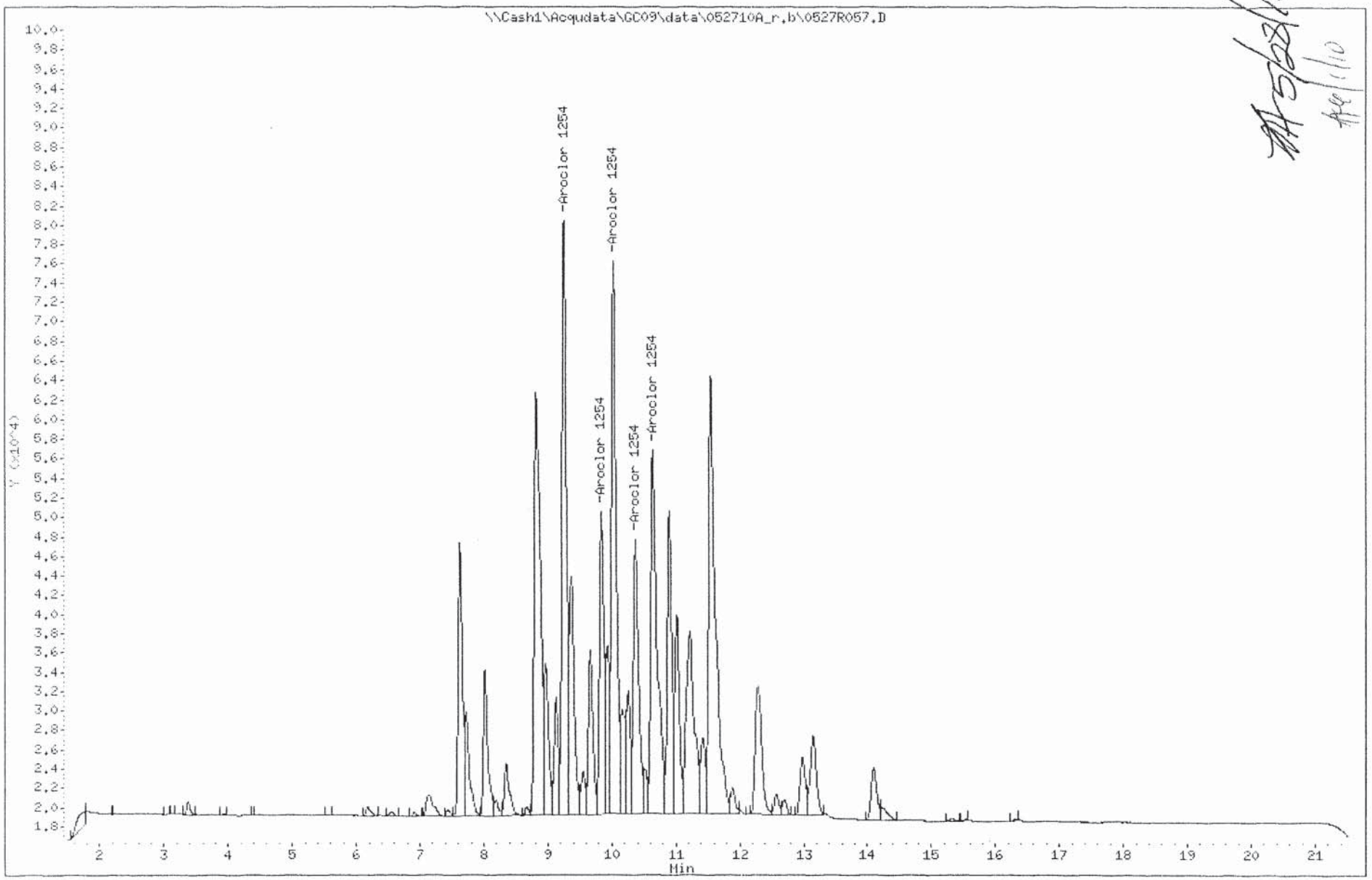
Instrument: GC09.i

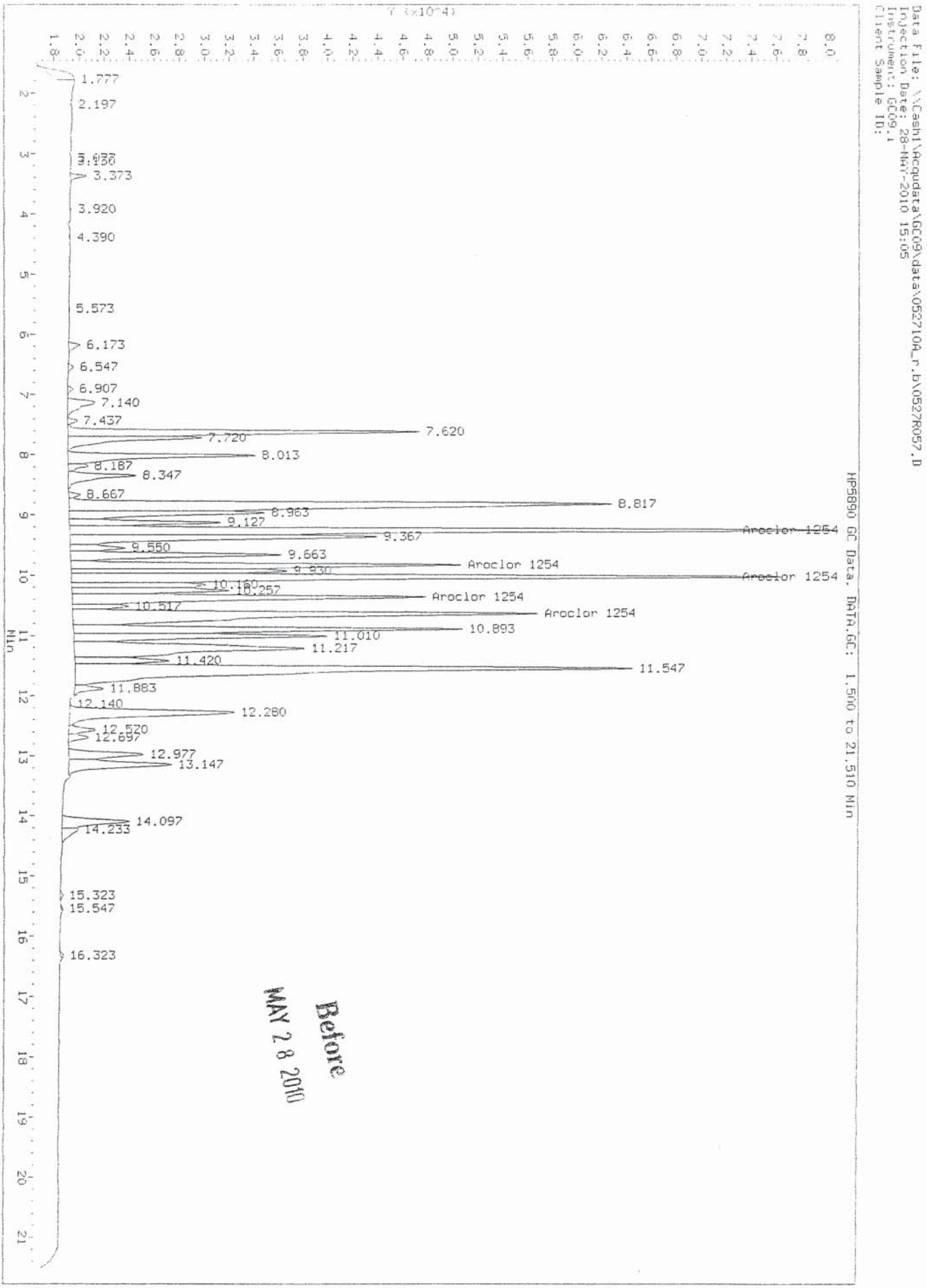
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



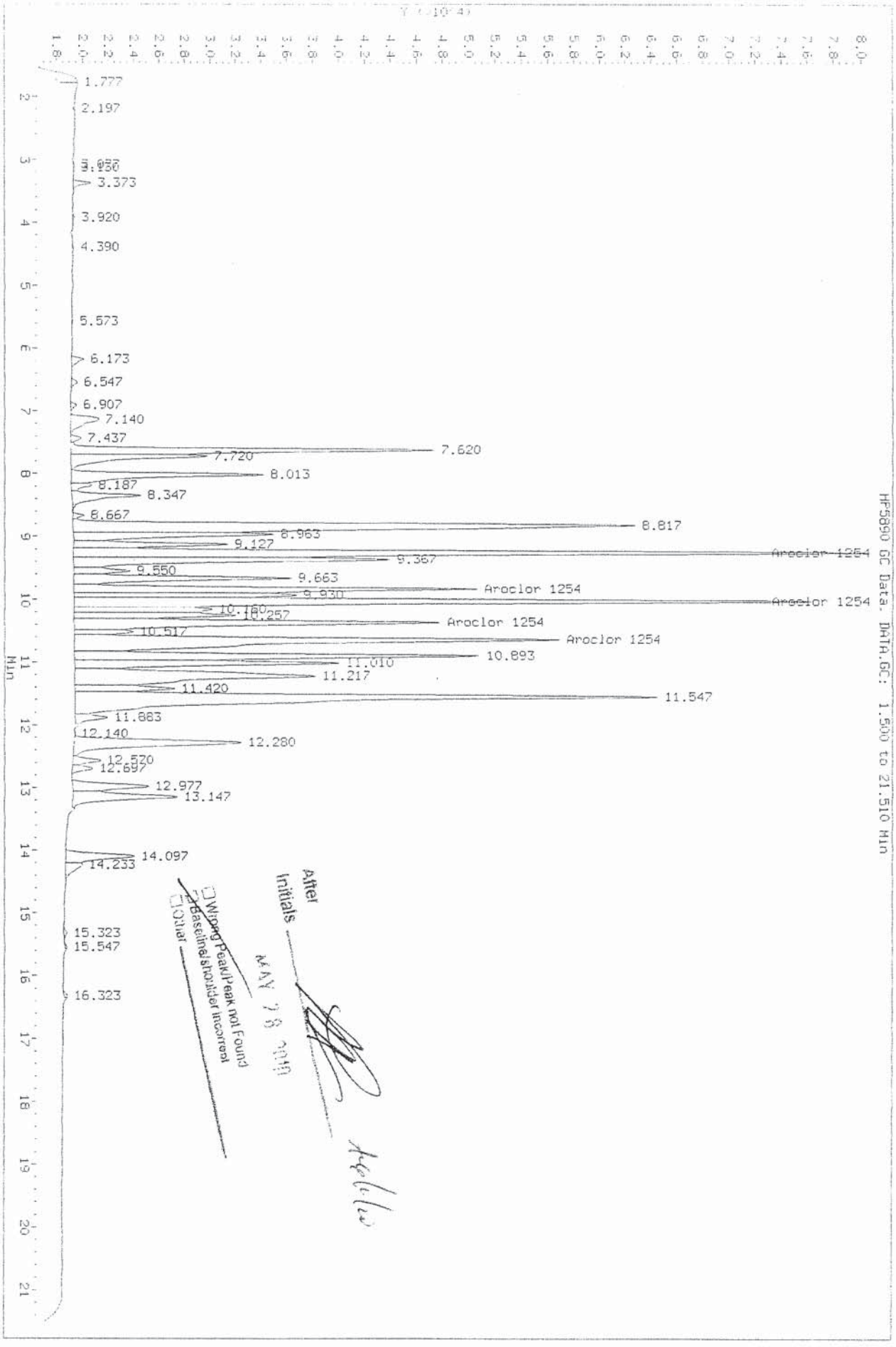
Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R057.D
Date : 28-MAY-2010 15:05
Client ID:
Sample Info: 1254 @ 1000ppb | PCB5-53K

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53





Before
MAY 28 2010


Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R057.D
Injection Date: 28-MAY-2010 15:05
Instrument: GC09.I
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.510 Min
Min
1.777
2.197
3.373
3.920
4.390
5.573
6.173
6.547
6.907
7.140
7.437
7.620
7.720
8.013
8.187
8.347
8.667
8.817
8.963
9.127
Aroclor 1254
9.367
9.550
9.663
Aroclor 1254
9.930
Aroclor 1254
10.257
Aroclor 1254
10.517
Aroclor 1254
10.893
11.010
11.217
11.420
11.547
11.883
12.140
12.280
12.530
12.697
12.977
13.147
14.097
14.253
15.323
15.547
16.323
After
Initials
MAY 28 2010
Wrong Peak/Peak not Found
Baseline/shoulder incorrect
Other

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F058.D
Report Date: 28-May-2010 17:31

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F058.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R058.D
Inj Date : 28-MAY-2010 15:31
Sample Info: 1260 @ 1000ppb | PCB5-49I
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1260.sub
Sub List #2 : AR1260.sub
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1260	9.490	10.363	286956	280458	949	1010	80.00- 120.00	100.00(M)
	9.770	10.893	376991	291710	1020	983	99.64- 149.46	131.38(M)
	10.093	11.547	450202	338687	994	910	122.12- 183.18	156.89(M)
	10.873	12.313	374032	281950	863	1200	109.38- 164.06	130.34(M)
	12.143	13.150	671293	505893	1180	1150	135.68- 203.52	243.84(M)
			Average of Peak Amounts =		1000	1050		

QC Flag Legend

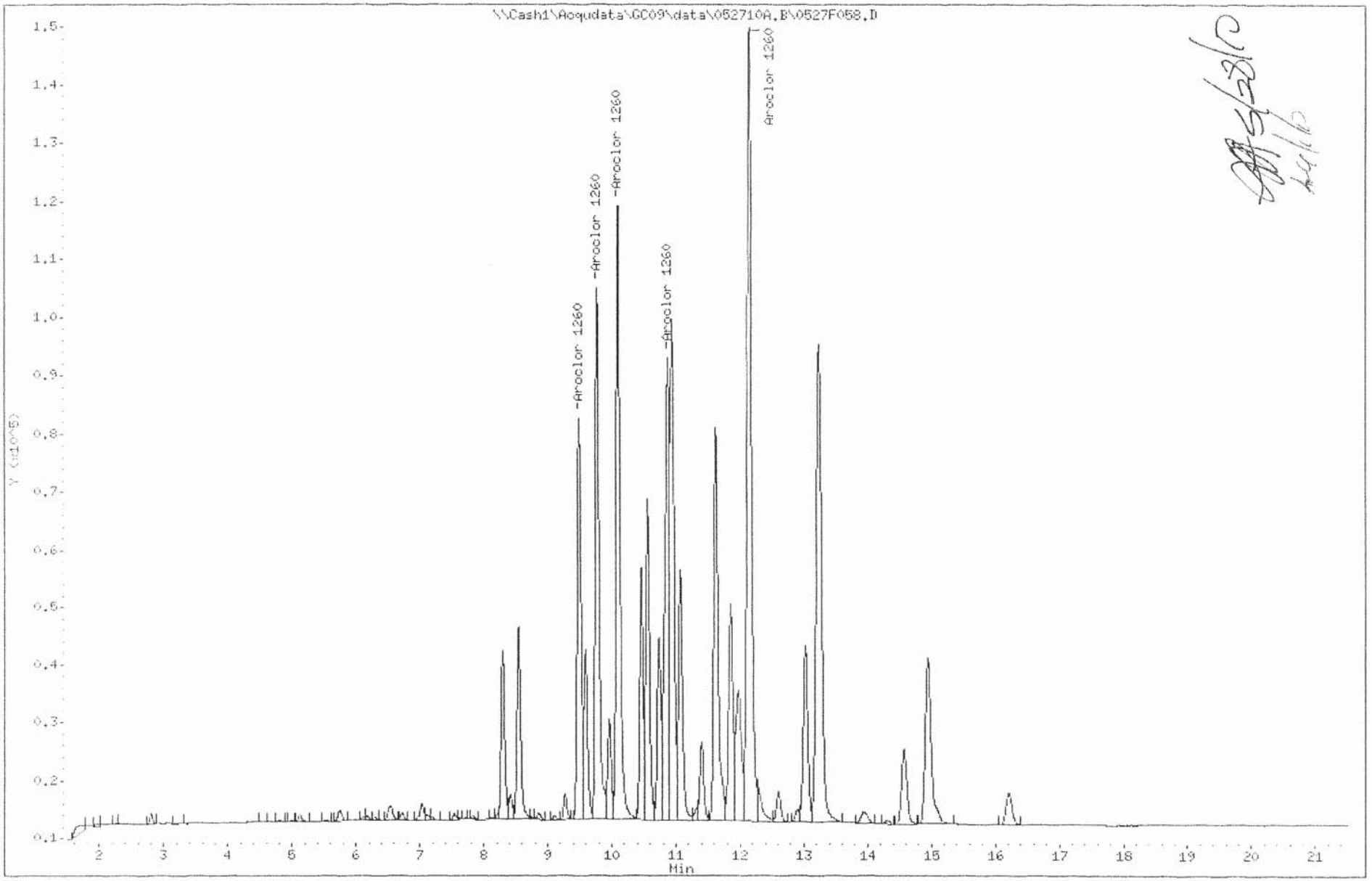
M - Compound response manually integrated.

5/28/10

6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F058.D
Date : 28-MAY-2010 15:31
Client ID:
Sample Info: 1260 @ 1000ppb | PCB5-49I

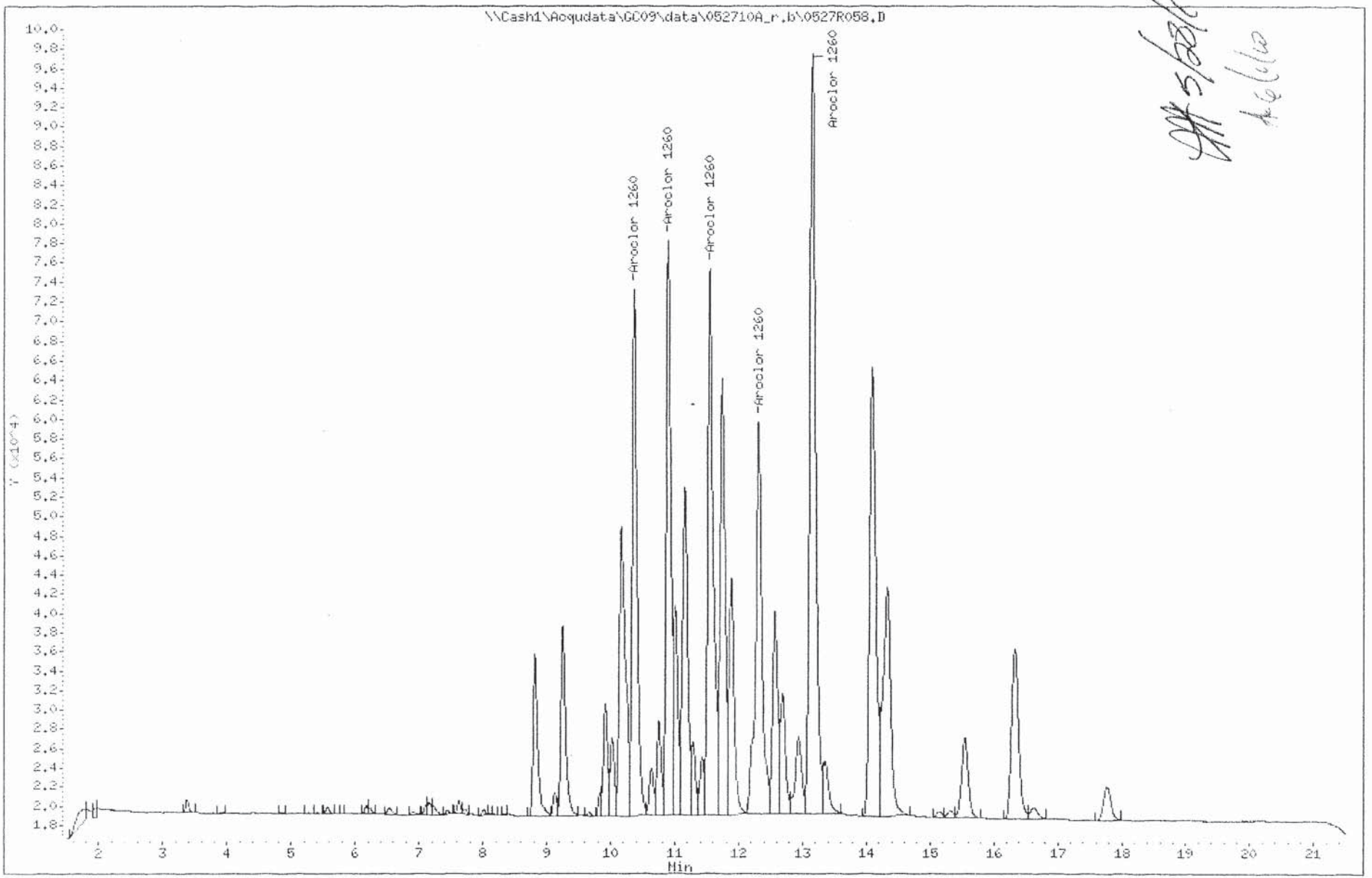
Instrument: GC09.i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R058.D
Date : 28-MAY-2010 15:31
Client ID:
Sample Info: 1260 @ 1000ppb | PCB5-49I

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F059.D
Report Date: 28-May-2010 17:31

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F059.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R059.D
Inj Date : 28-MAY-2010 15:57
Sample Info: 1262 @ 1000ppb | PCB5-49J
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1262.SUB
Sub List #2 : AR1262.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

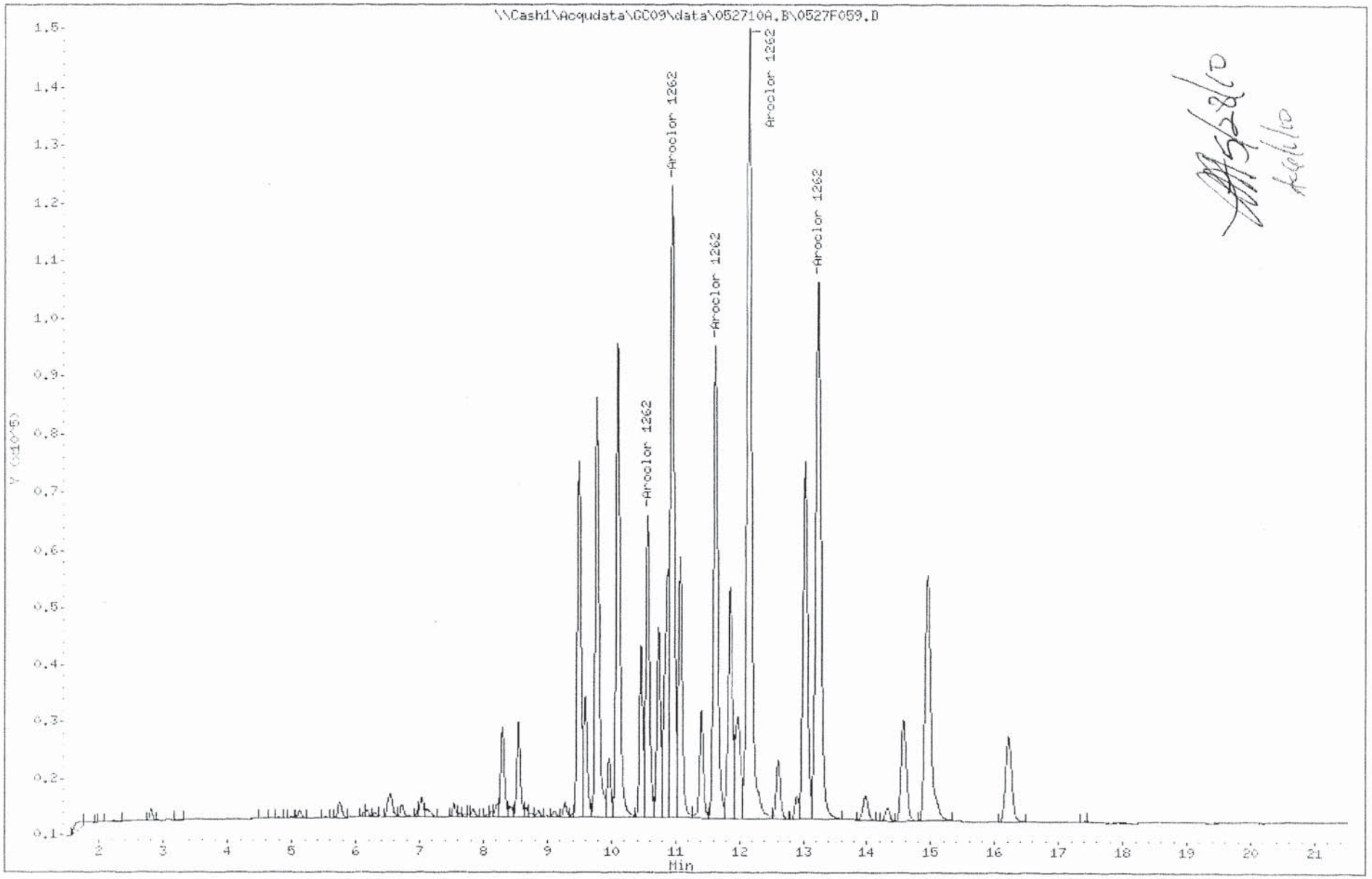
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1262	10.560	10.893	222102	234302	910	988	80.00- 120.00	100.00
	10.943	11.550	481487	207257	972	986	166.27- 249.41	216.79
	11.623	11.743	381965	333659	983	978	129.39- 194.09	171.98
	12.147	12.313	735672	259339	1000	1000	249.74- 374.61	331.23
	13.230	13.150	553496	546094	1020	988	188.68- 283.02	249.21
			Average of Peak Amounts =		977	988		

5/28/10

Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F059.D
Date : 28-MAY-2010 15:57
Client ID:
Sample Info: 1262 @ 1000ppb | PCB5-49J

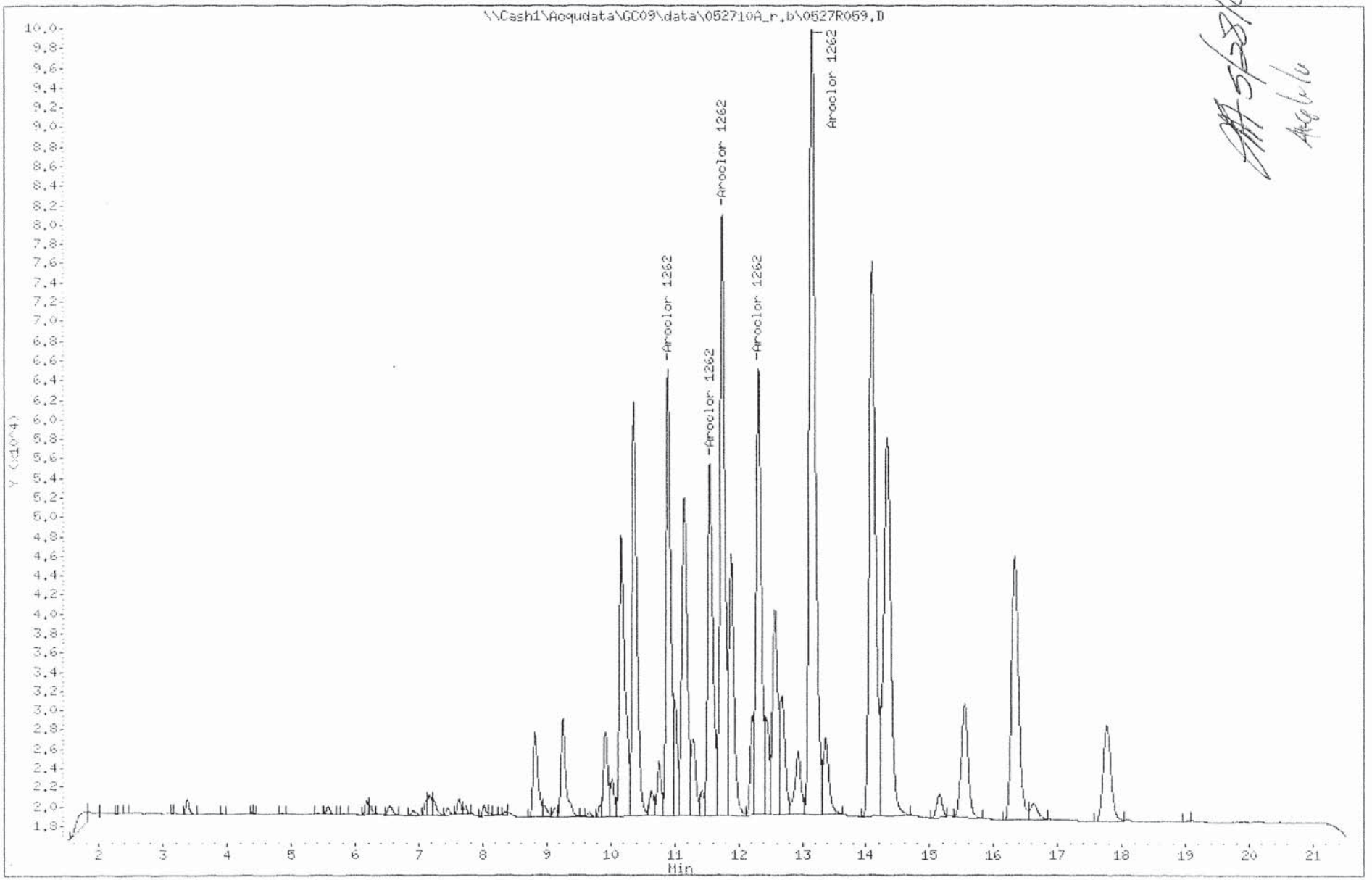
Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R059.D
Date : 28-MAY-2010 15:57
Client ID:
Sample Info: 1262 @ 1000ppb | PCB5-49J

Instrument: GC09.i
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Data File: \\Cash1\Acqudata\GC09\data\052710A.B\0527F060.D
Report Date: 28-May-2010 17:31

Columbia Analytical Services

Sample #1 : \\Cash1\Acqudata\GC09\data\052710A.B\0527F060.D
Sample #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\0527R060.D
Inj Date : 28-MAY-2010 16:23
Sample Info: 1268 @ 1000ppb | PCB5-49K
Misc Info :
Cal Date : 28-MAY-2010 14:15
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

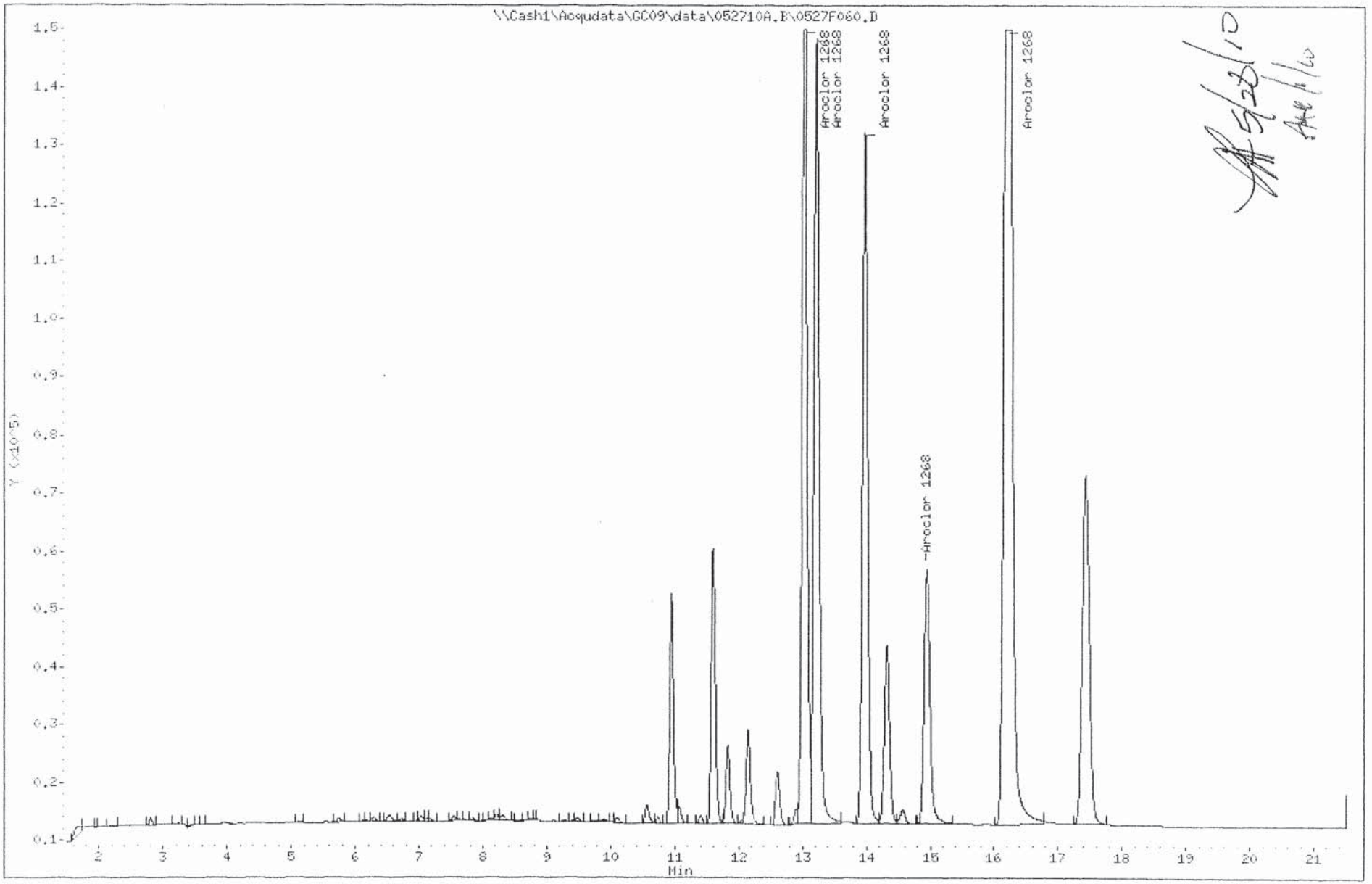
Method #1 : \\Cash1\Acqudata\GC09\data\052710A.B\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\052710A_r.b\052710A_r.m
Sub List #1 : AR1268.SUB
Sub List #2 : AR1268.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Aroclor 1268	13.023	14.340	784164	632174	774	823	80.00- 120.00	100.00
	13.213	15.147	763165	542902	817	894	70.89- 106.33	97.32
	13.977	15.477	660290	153221	876	951	58.04- 87.06	84.20
	14.947	16.327	284662	225700	921	902	20.96- 31.43	36.30
	16.210	17.770	1934553	1692617	905	932	172.26- 258.39	246.70
			Average of Peak Amounts =		859	900		

5/28/10
6/1/10

Data File: \\Cash1\Acqudata\GC09\data\052710A,B\0527F060.D
Date : 28-MAY-2010 16:23
Client ID:
Sample Info: 1268 @ 1000ppb | PCB5-49K

Instrument: GC09,i
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0,53

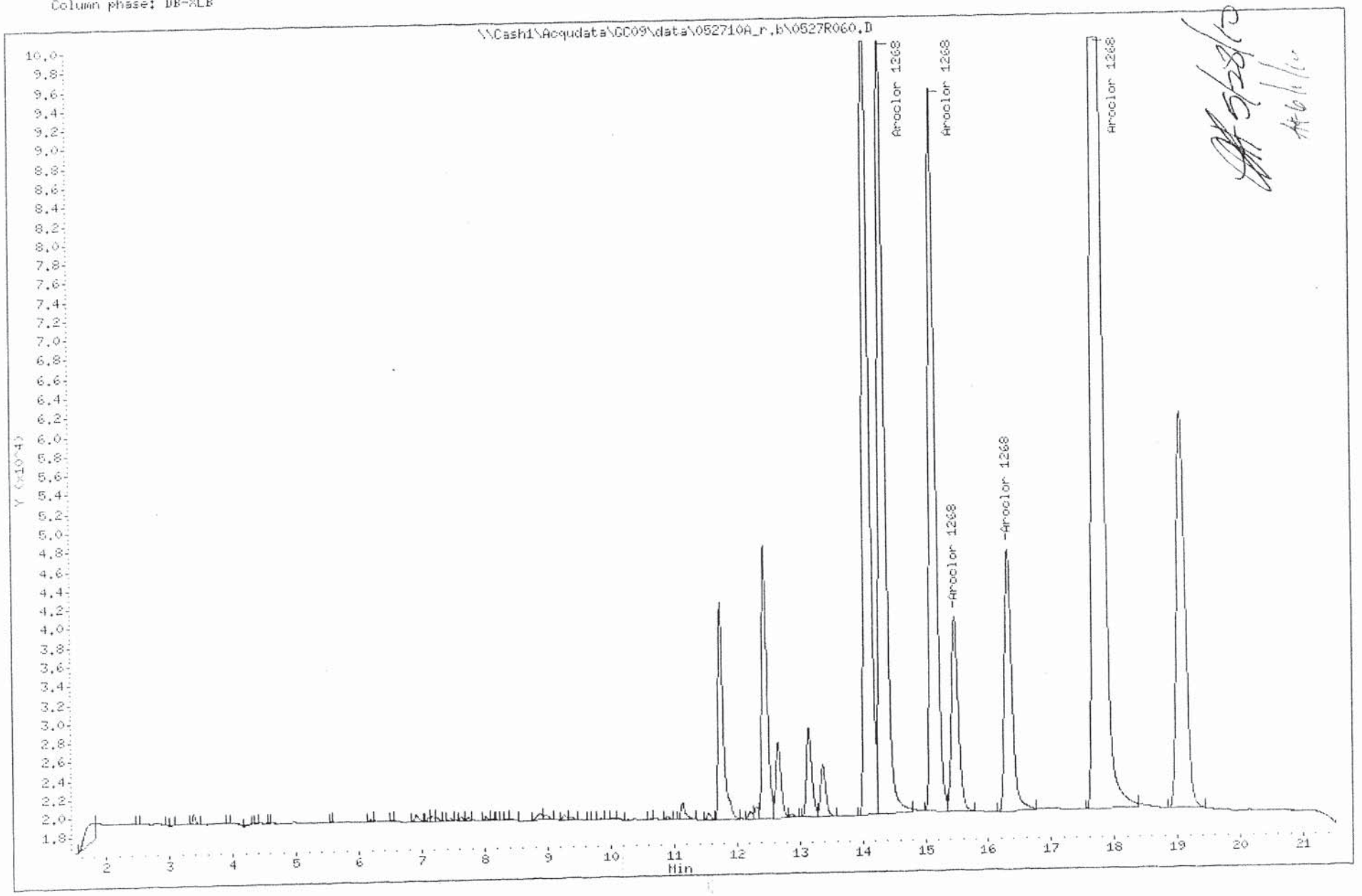


Data File: \\Cash1\Acqudata\GC09\data\052710A_r,b\0527R060.D
Date : 28-MAY-2010 16:23
Client ID:
Sample Info: 1268 @ 1000ppb | PCB5-49K

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/08/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F003.D

Column ID: DB-35MS

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	97	5900	5740	-3	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	1000	142	146	3	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1000	286	300	5	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1000	243	245	1	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	970	207	201	-3	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	154	158	2	NA	± 100 %	AverageRF
Aroclor 1016	1000	1000	NA	NA	NA	2	± 15 %	NA
Aroclor 1260 {1}	1000	950	302	288	-5	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	930	370	346	-7	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	940	453	425	-6	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	990	433	429	-1	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	990	571	567	-1	NA	± 100 %	AverageRF
Aroclor 1260	1000	960	NA	NA	NA	-4	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/08/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R003.D

Column ID: DB-XLB

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	100	5130	5210	1	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	990	183	181	-1	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	960	323	311	-4	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	158	172	9	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1000	131	135	3	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	147	154	5	NA	± 100 %	AverageRF
Aroclor 1016	1000	1000	NA	NA	NA	2	± 15 %	NA
Aroclor 1260 {1}	1000	990	277	275	-1	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	980	297	290	-2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	1000	372	374	1	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	1100	235	252	7	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1000	440	442	1	NA	± 100 %	AverageRF
Aroclor 1260	1000	1000	NA	NA	NA	1	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F003.D	**Date Acquired:**	06/08/2010 18:27
Lab ID:	KWG1005547-2	**Date Quantitated:**	06/09/2010 10:28
RunType:	CCV	**Batch ID:**	KWG1005547
Matrix:	NOT APPLICABLE	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: MM 6/9/10

Secondary Review: 6/10/10

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R003.D	Date Acquired:	06/08/2010 18:27
Lab ID:	KWG1005547-2	Date Quantitated:	06/09/2010 10:29
RunType:	CCV	Batch ID:	KWG1005547
Matrix:	NOT APPLICABLE	Analysis Method:	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: MBQD

Secondary Review: [illegible]

Quantitation Report

Bottle ID:		Tier:		Matrix:	NOT APPLICABL
Prod Code:	8082 PCB	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:		Report Group:	
Analysis Method:	8082	Prep Method:			
Prep Ref:		Prep Date:			

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:		Method ID:	MJ579
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F003.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R003.D			Vial:	96
Acqu Date:	06/08/2010 18:27	Quant Date:	06/09/2010 10:28	Dilution:	1.0
Run Type:	CCV			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005547-2				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	4.33	5.02	692263m	601402m	97.43	102.13			NA
			%Recovery =		NA	NA	Limits =	21-114	
Decachlorobiphenyl	17.39	19.02	574224	520631	97.27	101.43			NA
			%Recovery =		NA	NA	Limits =	36-113	

Target Compounds

Final Conc. Units: ug/L

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/L #1	ug/L #2	Rpt
Aroclor 1016			0	0	1,016	1,024			
Aroclor 1016 {1}	6.17	6.53	145881m	181457m	1,030	990.64			
Aroclor 1016 {2}	6.71	7.15	300418m	310938m	1,049	961.49			
Aroclor 1016 {3}	7.02	7.43	245498m	171944m	1,009	1,089			
Aroclor 1016 {4}	7.52	7.60	200913m	134940m	969.88	1,033			
Aroclor 1016 {5}	7.82	7.71	157560m	154205m	1,022	1,046			
Aroclor 1260			0	0	961.98	1,010			
Aroclor 1260 {1}	9.47	10.34	288291m	275321m	953.54	992.38			
Aroclor 1260 {2}	9.75	10.87	345727m	289848m	934.97	976.63			
Aroclor 1260 {3}	10.08	11.52	425217m	374159m	939.01	1,006			
Aroclor 1260 {4}	10.86	12.29	428795m	251661m	989.82	1,069			
Aroclor 1260 {5}	12.12	13.12	566750m	441958m	992.59	1,005			

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: If at/above MRL, also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Calibration Verification Report

Calibration ID: CAL9510
Method ID: MJ579
DataFile: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F003.D

Parameter Name	Type	PARM Type	Curve Fit	Method Criteria	Min RF	ICAL RF	CCV RF	%Diff	Sol'n Conc.	True Value	% Drift
Aroclor 1016		MS	NA	15					1,016	1,000	1.6
Aroclor 1260		MS	NA	15					962.0	1,000	-3.8
Tetrachloro-m-xylene		SURR	AverageRF	15		7.1E+3	6.9E+3	-2.6			
Aroclor 1016 {1}		MULTI	AverageRF	100		1.4E+2	1.5E+2	3.0			
Aroclor 1016 {2}		MULTI	AverageRF	100		2.9E+2	3.0E+2	4.9			
Aroclor 1016 {3}		MULTI	AverageRF	100		2.4E+2	2.5E+2	0.9			
Aroclor 1016 {4}		MULTI	AverageRF	100		2.1E+2	2.0E+2	-3.0			
Aroclor 1016 {5}		MULTI	AverageRF	100		1.5E+2	1.6E+2	2.2			
Aroclor 1260 {1}		MULTI	AverageRF	100		3.0E+2	2.9E+2	-4.6			
Aroclor 1260 {2}		MULTI	AverageRF	100		3.7E+2	3.5E+2	-6.5			
Aroclor 1260 {3}		MULTI	AverageRF	100		4.5E+2	4.3E+2	-6.1			
Aroclor 1260 {4}		MULTI	AverageRF	100		4.3E+2	4.3E+2	-1.0			
Aroclor 1260 {5}		MULTI	AverageRF	100		5.7E+2	5.7E+2	-0.7			
Decachlorobiphenyl		SURR	AverageRF	15		5.9E+3	5.7E+3	-2.7			

Alternate Calibration Evaluation Summary

Maximum Allowable Average %D =	15.0
Calculated Average %D =	3.2

Calibration Verification Report

Calibration ID: CAL9510
Method ID: MJ579
DataFile: \\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R003.D

Parameter Name	Type	PARM Type	Curve Fit	Method Criteria	Min RF	ICAL RF	CCV RF	%Diff	Sol'n Conc.	True Value	% Drift
Aroclor 1016		MS	NA	15					1,024	1,000	2.4
Aroclor 1260		MS	NA	15					1,010	1,000	1.0
Tetrachloro-m-xylene		SURR	AverageRF	15		5.9E+3	6.0E+3	2.1			
Aroclor 1016 {1}		MULTI	AverageRF	100		1.8E+2	1.8E+2	-0.9			
Aroclor 1016 {2}		MULTI	AverageRF	100		3.2E+2	3.1E+2	-3.9			
Aroclor 1016 {3}		MULTI	AverageRF	100		1.6E+2	1.7E+2	8.9			
Aroclor 1016 {4}		MULTI	AverageRF	100		1.3E+2	1.3E+2	3.3			
Aroclor 1016 {5}		MULTI	AverageRF	100		1.5E+2	1.5E+2	4.6			
Aroclor 1260 {1}		MULTI	AverageRF	100		2.8E+2	2.8E+2	-0.8			
Aroclor 1260 {2}		MULTI	AverageRF	100		3.0E+2	2.9E+2	-2.3			
Aroclor 1260 {3}		MULTI	AverageRF	100		3.7E+2	3.7E+2	0.5			
Aroclor 1260 {4}		MULTI	AverageRF	100		2.4E+2	2.5E+2	6.9			
Aroclor 1260 {5}		MULTI	AverageRF	100		4.4E+2	4.4E+2	0.5			
Decachlorobiphenyl		SURR	AverageRF	15		5.1E+3	5.2E+3	1.4			

Alternate Calibration Evaluation Summary

Maximum Allowable Average %D =	15.0
Calculated Average %D =	3.0

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F003.D
Report Date: 09-Jun-2010 10:28

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F003.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R003.D
Inj Date : 08-JUN-2010 18:27
Sample Info: 1660 @ 1000ppb | PCB5-55I | KWG1005547-2
Misc Info : SEMIVOA GC\W1005547\2-CCV.H
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.333	5.017	692263	601402	97.4	102		100.00(M)
Aroclor 1016	6.167	6.527	145881	181457	1030	991	80.00- 120.00	100.00(M)
	6.707	7.147	300418	310938	1050	961	164.25- 246.38	205.93(M)
	7.017	7.430	245498	171944	1010	1090	130.18- 195.27	168.29(M)
	7.523	7.603	200913	134940	970	1030	112.90- 169.35	137.72(M)
	7.820	7.707	157560	154205	1020	1040	87.15- 130.73	108.01(M)
			Average of Peak Amounts =		1020	1020		
Aroclor 1260	9.470	10.343	288291	275321	954	992	80.00- 120.00	100.00(M)
	9.753	10.873	345727	289848	935	977	99.43- 149.15	119.92(M)
	10.077	11.523	425217	374159	939	1000	120.99- 181.49	147.50(M)
	10.857	12.287	428795	251661	990	1070	123.90- 185.85	148.74(M)
	12.123	13.117	566750	441958	992	1000	165.84- 248.76	196.59(M)
			Average of Peak Amounts =		962	1010		
Decachlorobiphenyl	17.390	19.023	574224	520631	97.3	101		100.00

QC Flag Legend

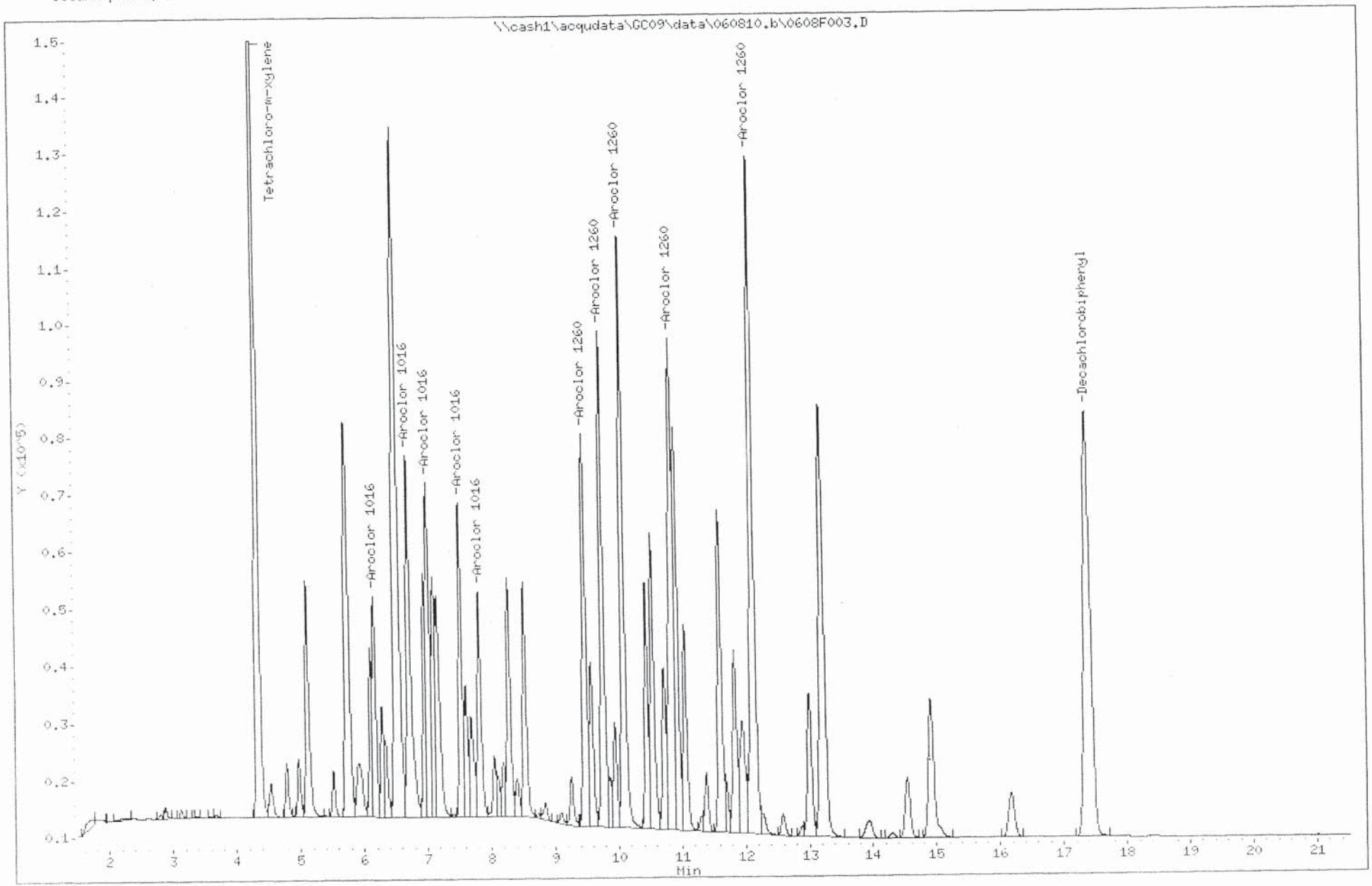
M - Compound response manually integrated.

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F003.D
Date : 08-JUN-2010 18:27
Client ID:
Sample Info: 1660 @ 1000ppb | PCB5-55I | KWG1005547-2

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

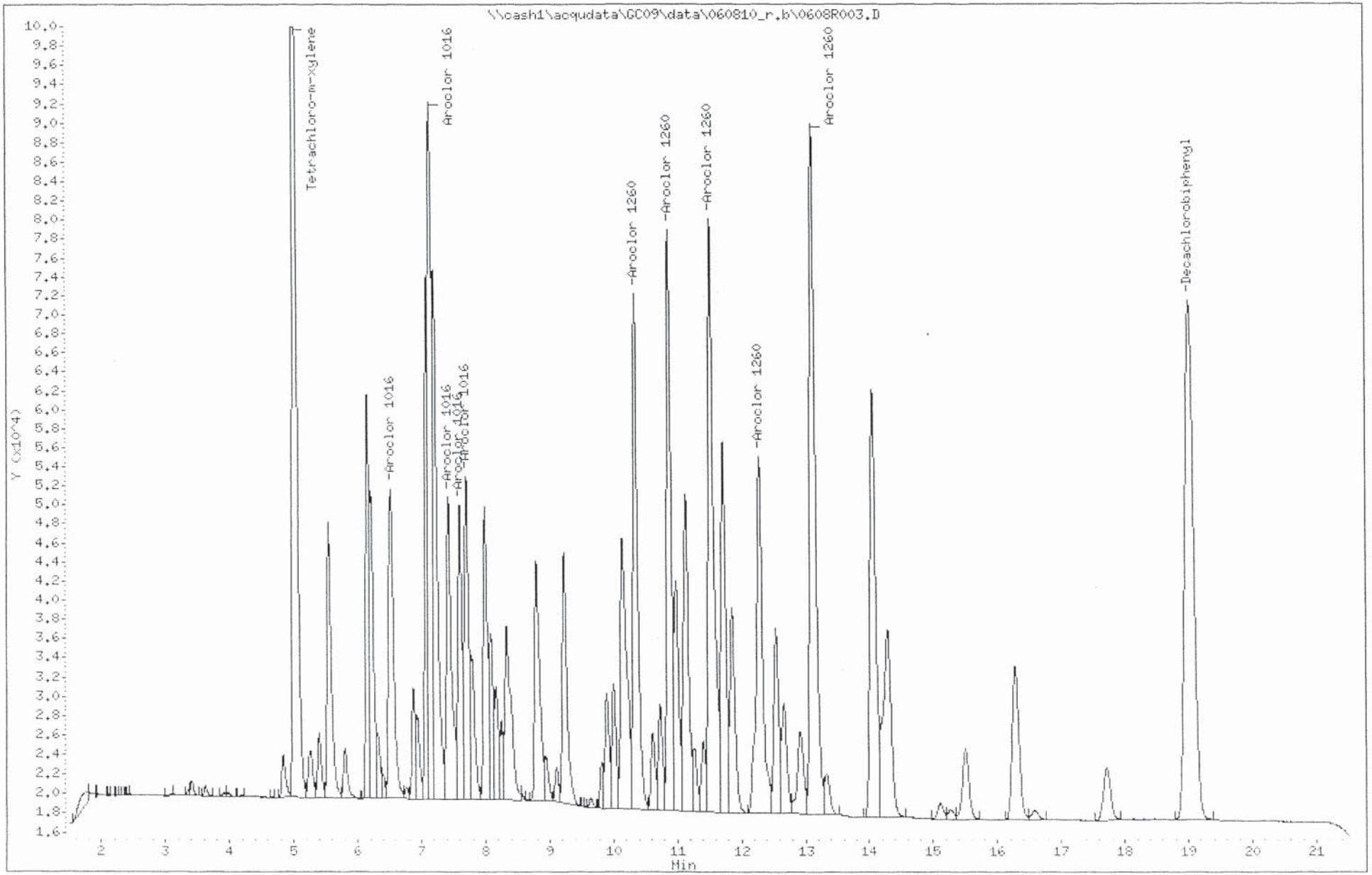
Column phase: DB-35MS

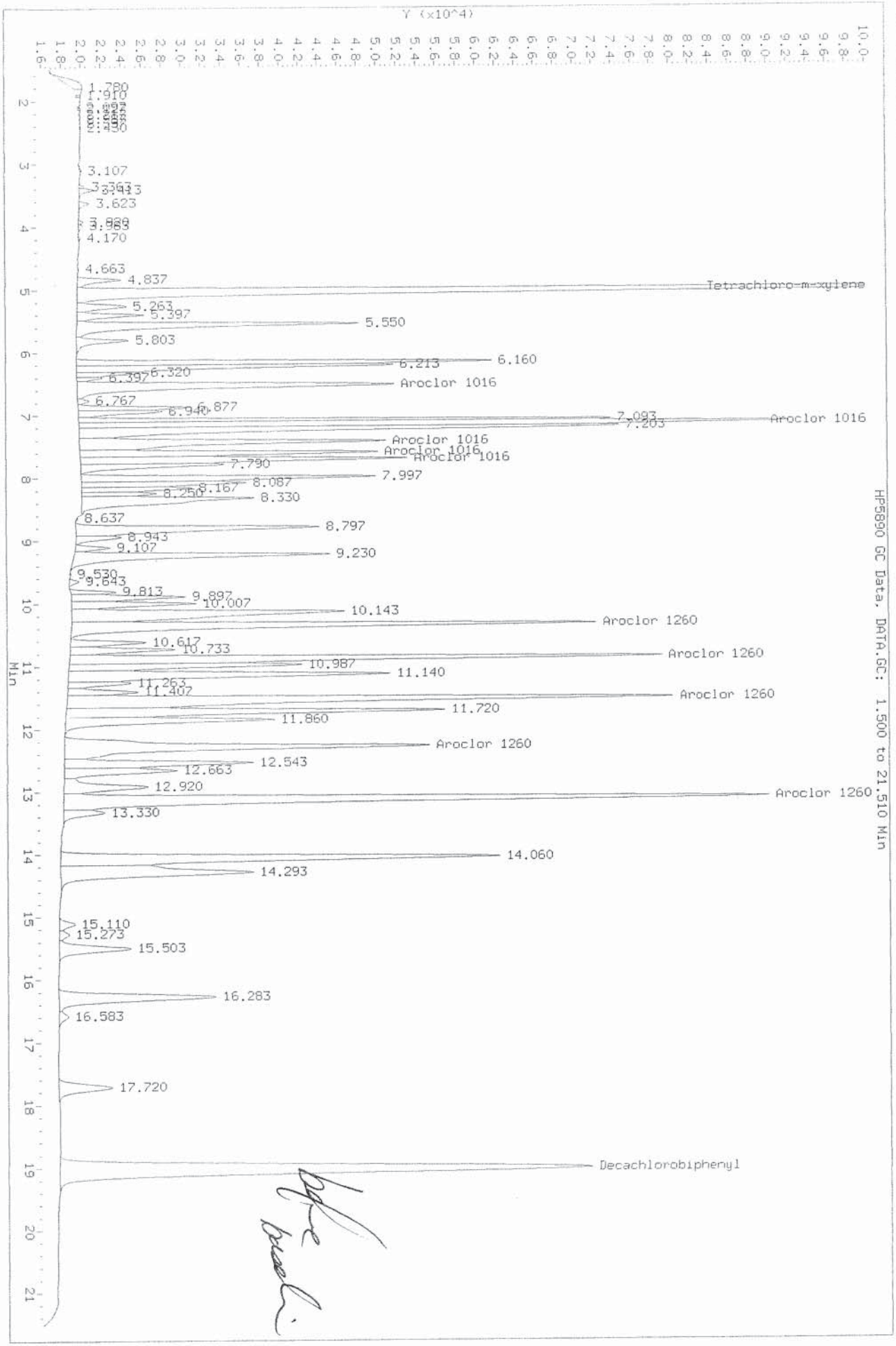


Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R003.D
Date : 08-JUN-2010 18:27
Client ID:
Sample Info: 1660 @ 1000ppb | PCB5-55I | KWG1005547-2

Instrument: GC09.i

Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53

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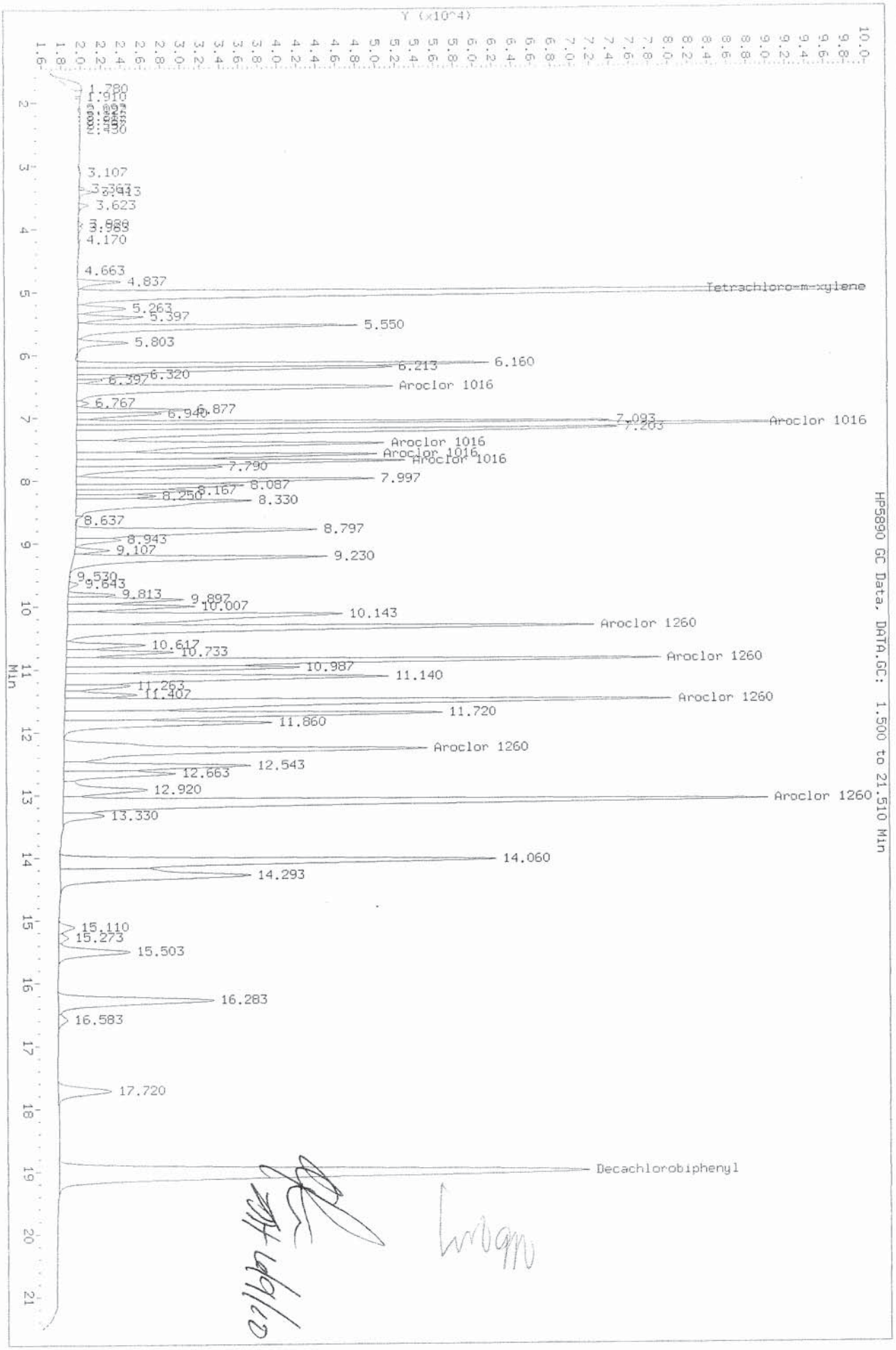
Injection Date: 08-JUN-2010 18:27
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC; 1.500 to 21.510 Min
Y (x10^4)
Min
1.780
1.910
3.107
3.623
4.170
4.663
4.837
Tetrachloro-m-xylene
5.263
5.397
5.550
5.803
6.213
6.160
6.320
6.397
Aroclor 1016
6.767
6.940
6.877
7.093
7.203
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.790
7.997
8.167
8.087
8.250
8.330
8.637
8.797
8.943
9.107
9.230
9.530
9.643
9.813
9.897
10.007
10.143
Aroclor 1260
10.617
10.733
Aroclor 1260
10.987
11.140
11.263
11.407
Aroclor 1260
11.720
11.860
Aroclor 1260
12.543
12.663
12.920
Aroclor 1260
13.330
14.060
14.293
15.110
15.273
15.503
16.283
16.583
17.720
Decachlorobiphenyl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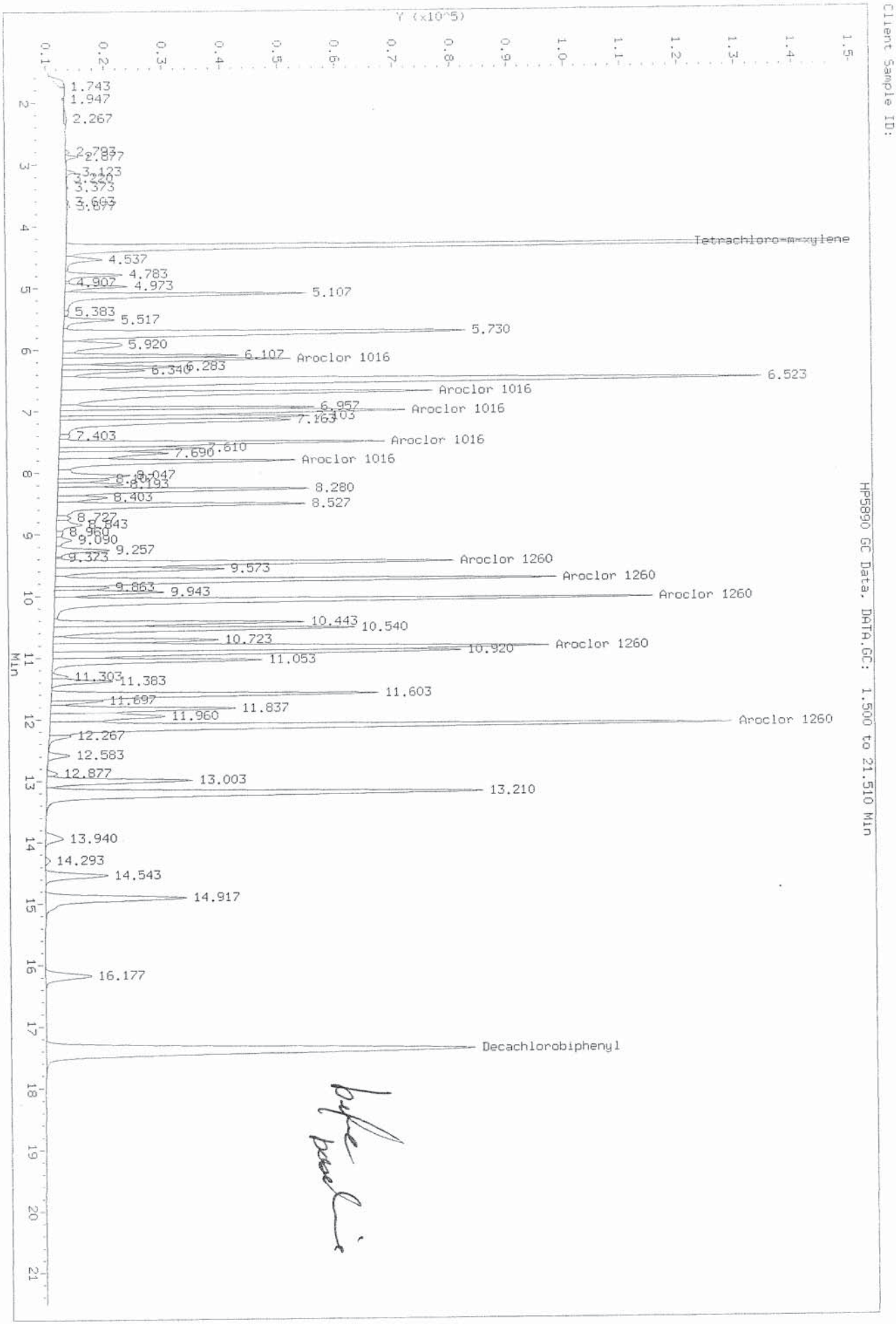
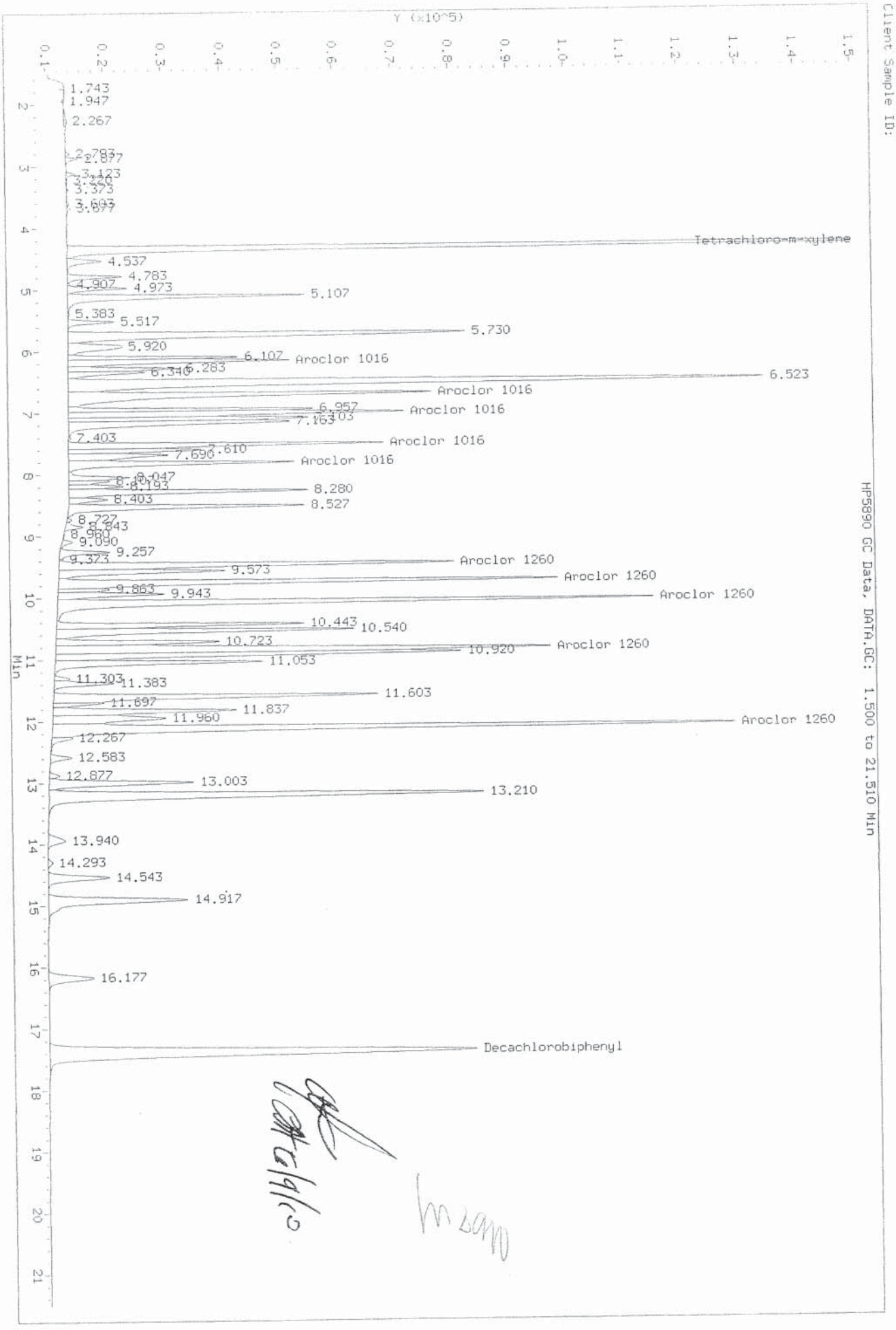
Injection Date: 08-JUN-2010 18:27
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Y (x10^4)
Min
1.780
1.910
3.107
3.413
3.623
3.988
4.170
4.663
4.837
Tetrachloro-m-xylene
5.263
5.397
5.550
5.803
6.160
6.213
6.320
6.397
Aroclor 1016
6.767
6.877
6.940
7.093
7.203
Aroclor 1016
Aroclor 1016
Aroclor 1016
Aroclor 1016
7.790
7.997
8.087
8.167
8.250
8.330
8.637
8.797
8.943
9.107
9.230
9.530
9.643
9.813
9.897
10.007
10.143
Aroclor 1260
10.617
10.733
Aroclor 1260
10.987
11.140
11.263
11.407
Aroclor 1260
11.720
11.860
Aroclor 1260
12.543
12.663
12.920
Aroclor 1260
13.330
14.060
14.293
15.110
15.273
15.503
16.283
16.583
17.720
Decachlorobiphenyl


Data File: \\Cash1\Acqudata\GC09\data\060810.b\0608F003.D
Injection Date: 08-JUN-2010 18:27
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data. DATA.GC: 1.500 to 21.510 Min
Y (x10^5)
Min
1.743
1.947
2.267
2.783
2.877
3.123
3.320
3.373
3.603
3.677
Tetrachloro-m-xylene
4.537
4.783
4.902
4.973
5.107
5.383
5.517
5.730
5.920
6.107
Aroclor 1016
6.340
6.283
6.523
Aroclor 1016
6.957
Aroclor 1016
7.103
7.163
7.403
Aroclor 1016
7.610
7.690
Aroclor 1016
8.047
8.193
8.280
8.403
8.527
8.727
8.843
8.960
9.090
9.257
9.323
Aroclor 1260
9.573
Aroclor 1260
9.863
9.943
Aroclor 1260
10.443
10.540
10.723
10.920
Aroclor 1260
11.053
11.303
11.383
11.603
11.697
11.837
11.960
Aroclor 1260
12.267
12.583
12.877
13.003
13.210
13.940
14.293
14.543
14.917
16.177
Decachlorobiphenyl


Data File: \\Cashl\Acqudata\GC09\data\060810.b\0608F003.D
Injection Date: 08-JUN-2010 18:27
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.510 Min
Y (x10^5)
Min
1.743
1.947
2.267
2.787
2.877
3.123
3.220
3.373
3.693
3.877
Tetrachloro-m-xylene
4.537
4.783
4.907
4.973
5.107
5.383
5.517
5.730
5.920
6.107
Aroclor 1016
6.283
6.340
6.523
Aroclor 1016
6.957
Aroclor 1016
7.103
7.163
7.403
Aroclor 1016
7.610
7.690
Aroclor 1016
8.047
8.193
8.280
8.403
8.527
8.727
8.843
8.960
9.090
9.257
9.323
Aroclor 1260
9.573
Aroclor 1260
9.863
9.943
Aroclor 1260
10.443
10.540
10.723
10.920
Aroclor 1260
11.053
11.303
11.383
11.603
11.697
11.837
11.960
Aroclor 1260
12.267
12.583
12.877
13.003
13.210
13.940
14.293
14.543
14.917
16.177
Decachlorobiphenyl

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/09/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F019.D

Column ID: DB-35MS

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	98	5900	5760	-2	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	1000	142	143	1	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	1000	286	295	3	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	960	243	232	-4	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	970	207	201	-3	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	154	156	1	NA	± 100 %	AverageRF
Aroclor 1016	1000	990	NA	NA	NA	-1	± 15 %	NA
Aroclor 1260 {1}	1000	940	302	284	-6	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	930	370	344	-7	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	930	453	421	-7	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	990	433	427	-1	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1000	571	571	0	NA	± 100 %	AverageRF
Aroclor 1260	1000	960	NA	NA	NA	-4	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 06/09/2010

Continuing Calibration Verification Summary
Polychlorinated Biphenyls (PCBs)

Calibration Type: External Standard
Analysis Method: 8082
CCV Standard ID: PCB5-55I

Calibration Date: 05/27/2010
Calibration ID: CAL9510
Analysis Lot: KWG1005547
Units: ng/mL

File ID: \\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R019.D

Column ID: DB-XLB

Analyte Name	Expected	Result	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Decachlorobiphenyl	100	100	5130	5190	1	NA	± 15 %	AverageRF
Aroclor 1016 {1}	1000	970	183	178	-3	NA	± 100 %	AverageRF
Aroclor 1016 {2}	1000	950	323	306	-5	NA	± 100 %	AverageRF
Aroclor 1016 {3}	1000	1100	158	170	7	NA	± 100 %	AverageRF
Aroclor 1016 {4}	1000	1000	131	133	2	NA	± 100 %	AverageRF
Aroclor 1016 {5}	1000	1000	147	152	3	NA	± 100 %	AverageRF
Aroclor 1016	1000	1000	NA	NA	NA	1	± 15 %	NA
Aroclor 1260 {1}	1000	990	277	274	-1	NA	± 100 %	AverageRF
Aroclor 1260 {2}	1000	980	297	292	-2	NA	± 100 %	AverageRF
Aroclor 1260 {3}	1000	1000	372	374	0	NA	± 100 %	AverageRF
Aroclor 1260 {4}	1000	1100	235	251	7	NA	± 100 %	AverageRF
Aroclor 1260 {5}	1000	1000	440	443	1	NA	± 100 %	AverageRF
Aroclor 1260	1000	1000	NA	NA	NA	1	± 15 %	NA

Results flagged with an asterisk (*) indicate values outside control criteria.

Exception Report

Data File: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F019.D
Lab ID: KWG1005547-4
RunType: CCV
Matrix: NOT APPLICABLE

Date Acquired: 06/09/2010 01:25
Date Quantitated: 06/09/2010 10:29
Batch ID: KWG1005547
Analysis Method: 8082
MethodJoinID: MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature] 6/10/10

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R019.D	**Date Acquired:**	06/09/2010 01:25
Lab ID:	KWG1005547-4	**Date Quantitated:**	06/09/2010 10:30
RunType:	CCV	**Batch ID:**	KWG1005547
Matrix:	NOT APPLICABLE	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: MBGN

Secondary Review: 6/10/2010

Quantitation Report

Bottle ID:		Tier:		Matrix:	NOT APPLICABL
Prod Code:	8082 PCB	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:		Report Group:	
Analysis Method:	8082	Prep Method:			
Prep Ref:		Prep Date:			

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:		Method ID:	MJ579
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F019.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R019.D			Vial:	96
Acqu Date:	06/09/2010 01:25	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	CCV			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005547-4				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	4.33	5.02	675384	588830m	95.05	99.99			NA
			%Recovery =		NA	NA	Limits =	21-114	
Decachlorobiphenyl	17.39	19.02	576179	519487	97.60	101.20			NA
			%Recovery =		NA	NA	Limits =	36-113	

Target Compounds

Final Conc. Units: ug/L

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/L #1	ug/L #2	Rpt
Aroclor 1016			0	0	994.26	1,008			
Aroclor 1016 {1}	6.17	6.53	142527m	177625m	1,007	969.72			
Aroclor 1016 {2}	6.71	7.15	295434m	305811m	1,032	945.64			
Aroclor 1016 {3}	7.02	7.43	232366m	169767m	955.08	1,075			
Aroclor 1016 {4}	7.53	7.60	200576m	133314m	968.26	1,020			
Aroclor 1016 {5}	7.82	7.71	155701m	151923m	1,010	1,030			
Aroclor 1260			0	0	956.86	1,010			
Aroclor 1260 {1}	9.47	10.34	284077m	274409m	939.61	989.09			
Aroclor 1260 {2}	9.75	10.87	343579m	291607m	929.16	982.56			
Aroclor 1260 {3}	10.08	11.52	421480m	373679m	930.76	1,004			
Aroclor 1260 {4}	10.85	12.29	426925m	251121m	985.50	1,066			
Aroclor 1260 {5}	12.12	13.12	570582m	443020m	999.30	1,008			

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Dtt above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Calibration Verification Report

Calibration ID: CAL9510
Method ID: MJ579
DataFile: \\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F019.D

Parameter Name	Type	PARM Type	Curve Fit	Method Criteria	Min RF	ICAL RF	CCV RF	%Diff	Sol'n Conc.	True Value	% Drift
Aroclor 1016		MS	NA	15					994.3	1,000	-0.6
Aroclor 1260		MS	NA	15					956.9	1,000	-4.3
Tetrachloro-m-xylene		SURR	AverageRF	15		7.1E+3	6.8E+3	-5.0			
Aroclor 1016 {1}		MULTI	AverageRF	100		1.4E+2	1.4E+2	0.7			
Aroclor 1016 {2}		MULTI	AverageRF	100		2.9E+2	3.0E+2	3.2			
Aroclor 1016 {3}		MULTI	AverageRF	100		2.4E+2	2.3E+2	-4.5			
Aroclor 1016 {4}		MULTI	AverageRF	100		2.1E+2	2.0E+2	-3.2			
Aroclor 1016 {5}		MULTI	AverageRF	100		1.5E+2	1.6E+2	1.0			
Aroclor 1260 {1}		MULTI	AverageRF	100		3.0E+2	2.8E+2	-6.0			
Aroclor 1260 {2}		MULTI	AverageRF	100		3.7E+2	3.4E+2	-7.1			
Aroclor 1260 {3}		MULTI	AverageRF	100		4.5E+2	4.2E+2	-6.9			
Aroclor 1260 {4}		MULTI	AverageRF	100		4.3E+2	4.3E+2	-1.5			
Aroclor 1260 {5}		MULTI	AverageRF	100		5.7E+2	5.7E+2	-0.1			
Decachlorobiphenyl		SURR	AverageRF	15		5.9E+3	5.8E+3	-2.4			

Alternate Calibration Evaluation Summary

Maximum Allowable Average %D =	15.0
Calculated Average %D =	3.4

Calibration Verification Report

Calibration ID: CAL9510
Method ID: MJ579

DataFile: \\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R019.D

Parameter Name	Type	PARM Type	Curve Fit	Method Criteria	Min RF	ICAL RF	CCV RF	%Diff	Sol'n Conc.	True Value	% Drift
Aroclor 1016		MS	NA	15					1,008	1,000	0.8
Aroclor 1260		MS	NA	15					1,010	1,000	1.0
Tetrachloro-m-xylene		SURR	AverageRF	15		5.9E+3	5.9E+3	0.0			
Aroclor 1016 {1}		MULTI	AverageRF	100		1.8E+2	1.8E+2	-3.0			
Aroclor 1016 {2}		MULTI	AverageRF	100		3.2E+2	3.1E+2	-5.4			
Aroclor 1016 {3}		MULTI	AverageRF	100		1.6E+2	1.7E+2	7.5			
Aroclor 1016 {4}		MULTI	AverageRF	100		1.3E+2	1.3E+2	2.0			
Aroclor 1016 {5}		MULTI	AverageRF	100		1.5E+2	1.5E+2	3.0			
Aroclor 1260 {1}		MULTI	AverageRF	100		2.8E+2	2.7E+2	-1.1			
Aroclor 1260 {2}		MULTI	AverageRF	100		3.0E+2	2.9E+2	-1.7			
Aroclor 1260 {3}		MULTI	AverageRF	100		3.7E+2	3.7E+2	0.4			
Aroclor 1260 {4}		MULTI	AverageRF	100		2.4E+2	2.5E+2	6.6			
Aroclor 1260 {5}		MULTI	AverageRF	100		4.4E+2	4.4E+2	0.8			
Decachlorobiphenyl		SURR	AverageRF	15		5.1E+3	5.2E+3	1.2			

Alternate Calibration Evaluation Summary

Maximum Allowable Average %D =	15.0
Calculated Average %D =	2.7

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F019.D
Report Date: 09-Jun-2010 10:29

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F019.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R019.D
Inj Date : 09-JUN-2010 01:25
Sample Info: 1660 @ 1000ppb | PCB5-55I | KWG1005547-4
Misc Info : SEMIVOA GC\W1005547\4-CCV.H
Cal Date : 09-JUN-2010 08:49
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : AR1660.SUB
Sub List #2 : AR1660.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio
Tetrachloro-m-xylene	4.333	5.017	675384	588830	95.0	100		100.00
Aroclor 1016	6.167	6.527	142527	177625	1010	970	80.00- 120.00	100.00(M)
	6.710	7.147	295434	305811	1030	946	164.25- 246.38	207.28(M)
	7.017	7.430	232366	169767	955	1070	130.18- 195.27	163.03(M)
	7.527	7.603	200576	133314	968	1020	112.90- 169.35	140.73(M)
	7.820	7.707	155701	151923	1010	1030	87.15- 130.73	109.24(M)
			Average of Peak Amounts =		995	1010		
Aroclor 1260	9.470	10.343	284077	274409	940	989	80.00- 120.00	100.00(M)
	9.753	10.873	343579	291607	929	982	99.43- 149.15	120.95(M)
	10.080	11.523	421480	373679	931	1000	120.99- 181.49	148.37(M)
	10.853	12.287	426925	251121	985	1070	123.90- 185.85	150.28(M)
	12.123	13.117	570582	443020	999	1010	165.84- 248.76	200.85(M)
			Average of Peak Amounts =		957	1010		
Decachlorobiphenyl	17.393	19.023	576179	519487	97.6	101		100.00

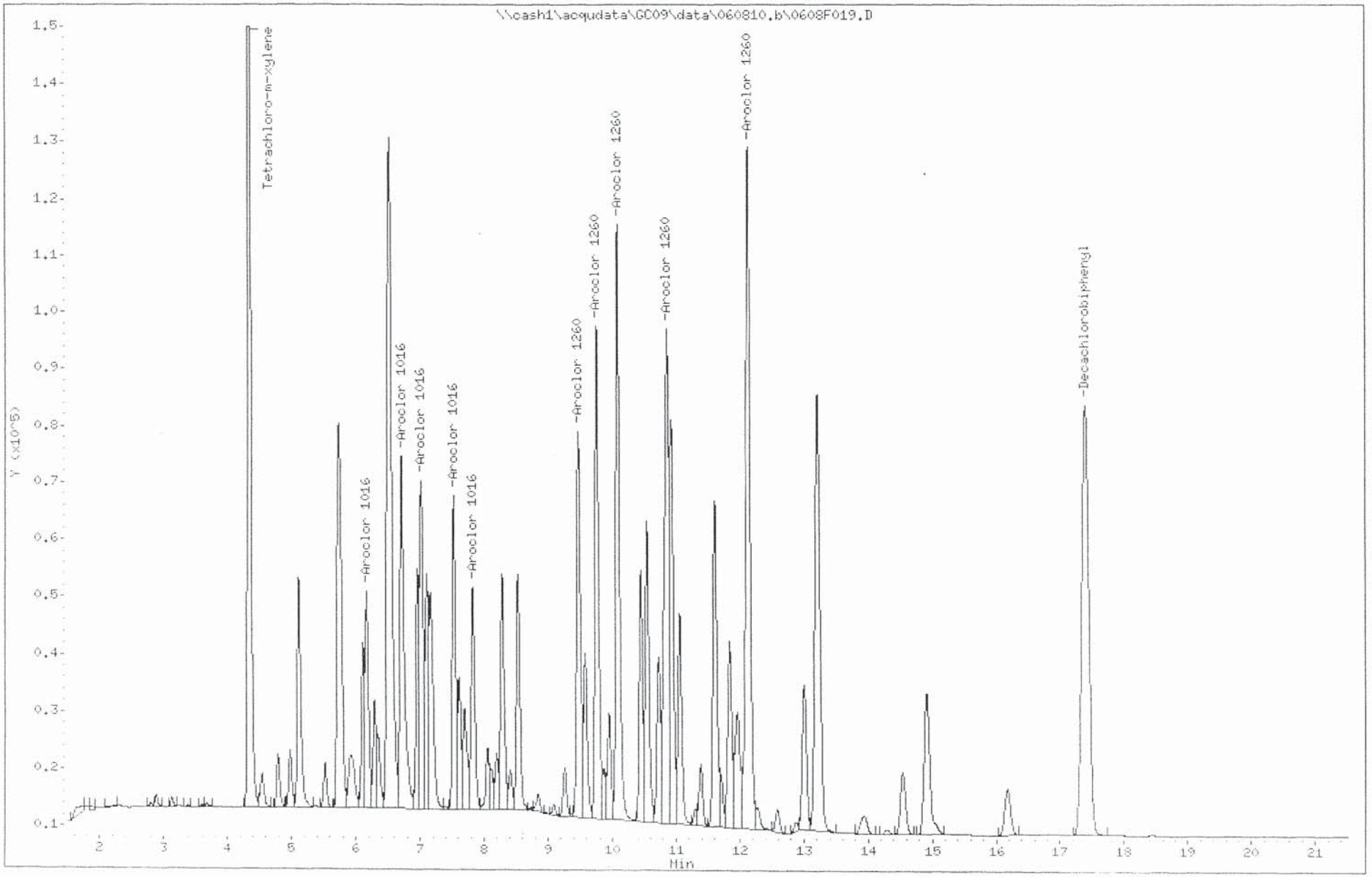
QC Flag Legend

M - Compound response manually integrated.

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F019.D
Date : 09-JUN-2010 01:25
Client ID:
Sample Info: 1660 @ 1000ppb | PCB5-55I | KWG1005547-4

Instrument: GC09.i

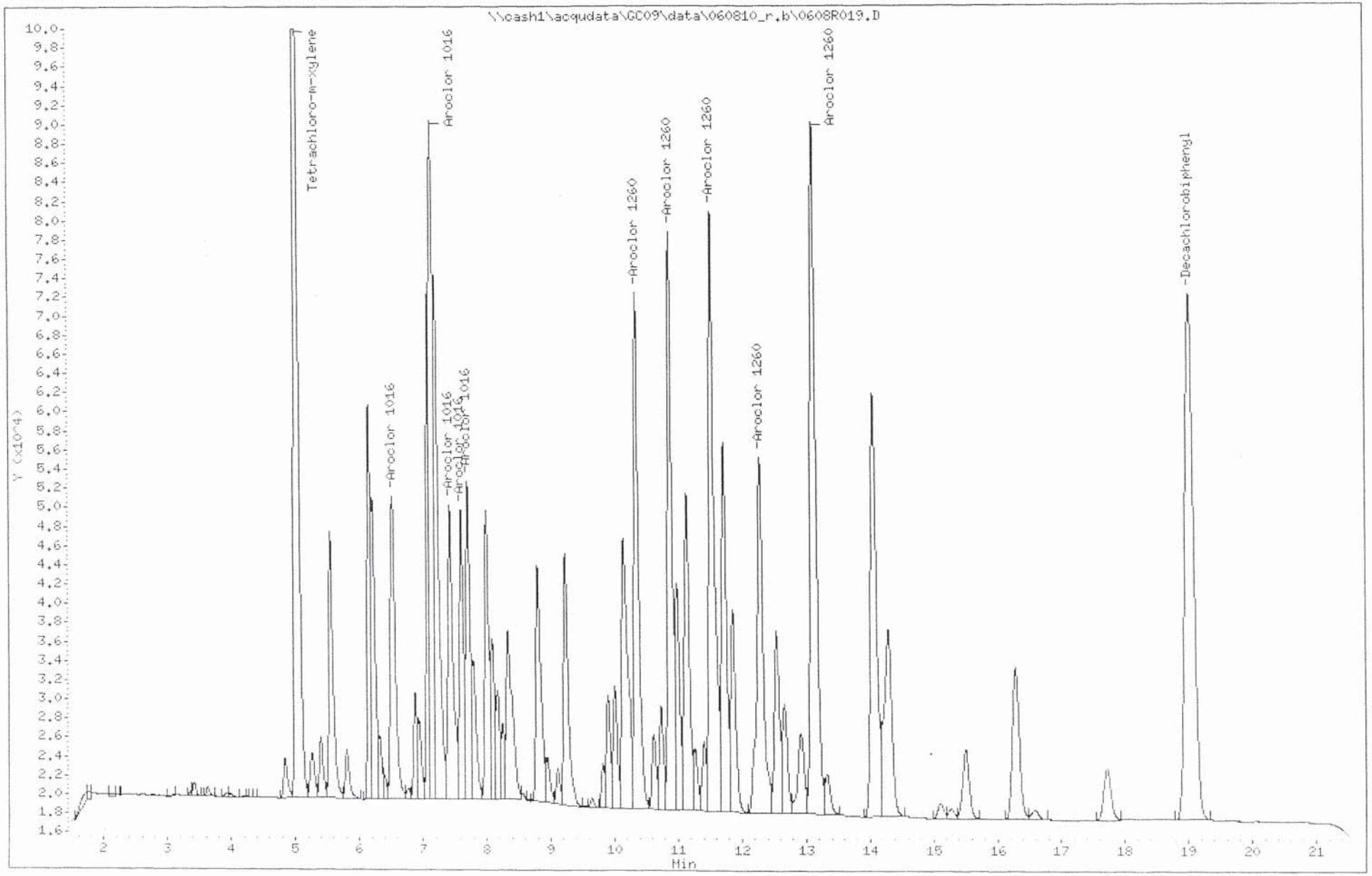
Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53

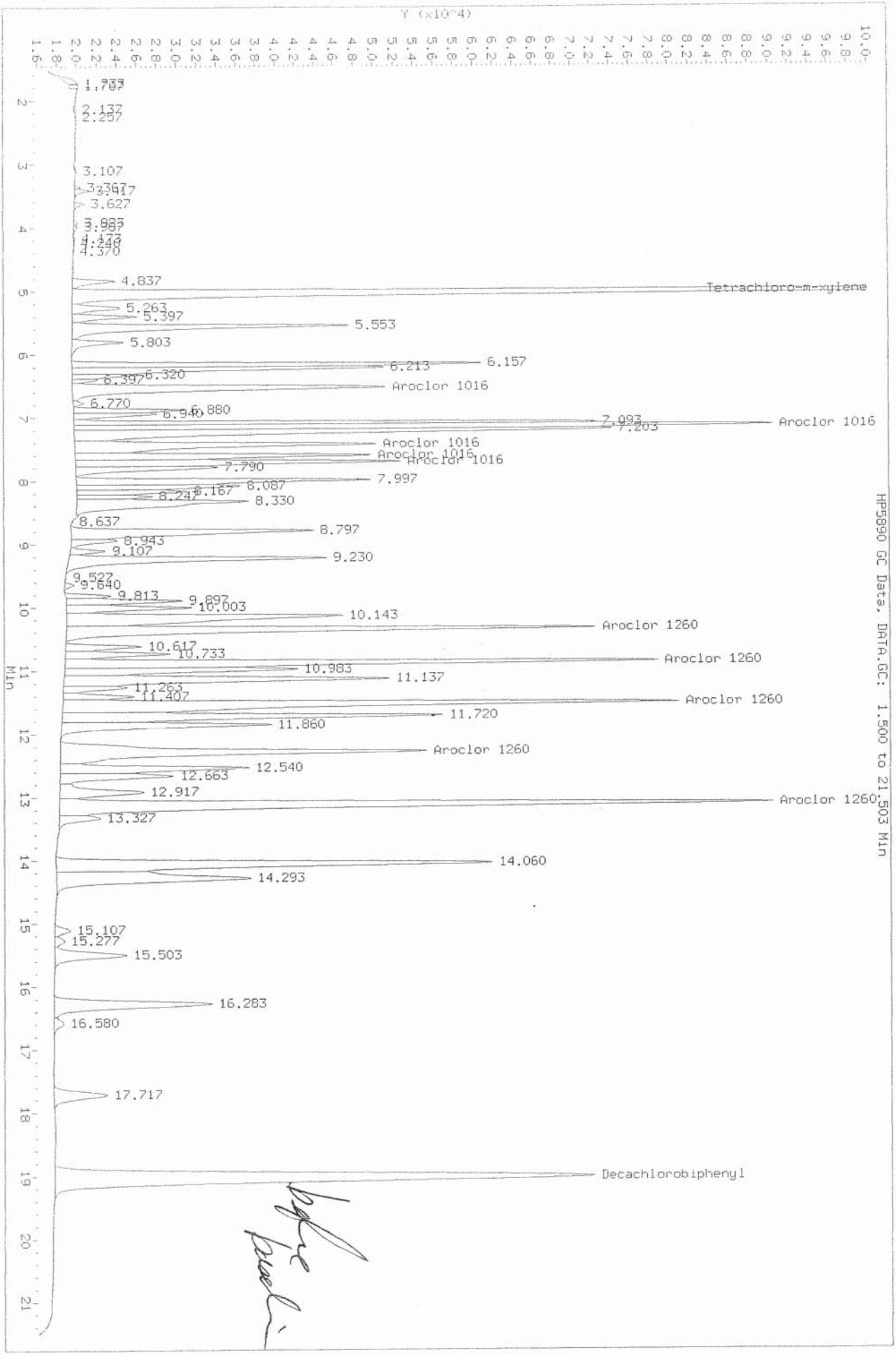


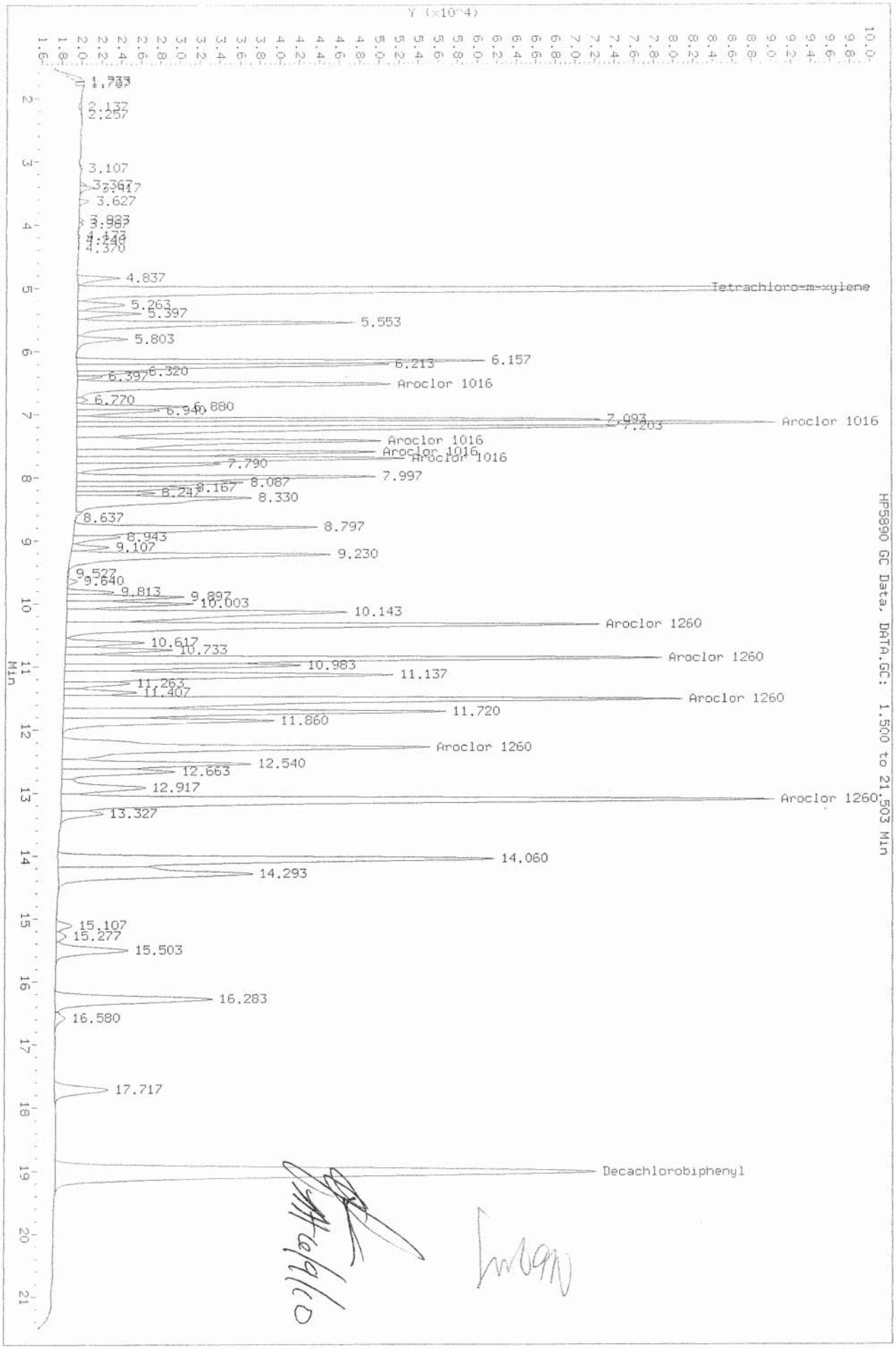
Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R019.D
Date : 09-JUN-2010 01:25
Client ID:
Sample Info: 1660 @ 1000ppb | PCB5-55I | KWG1005547-4

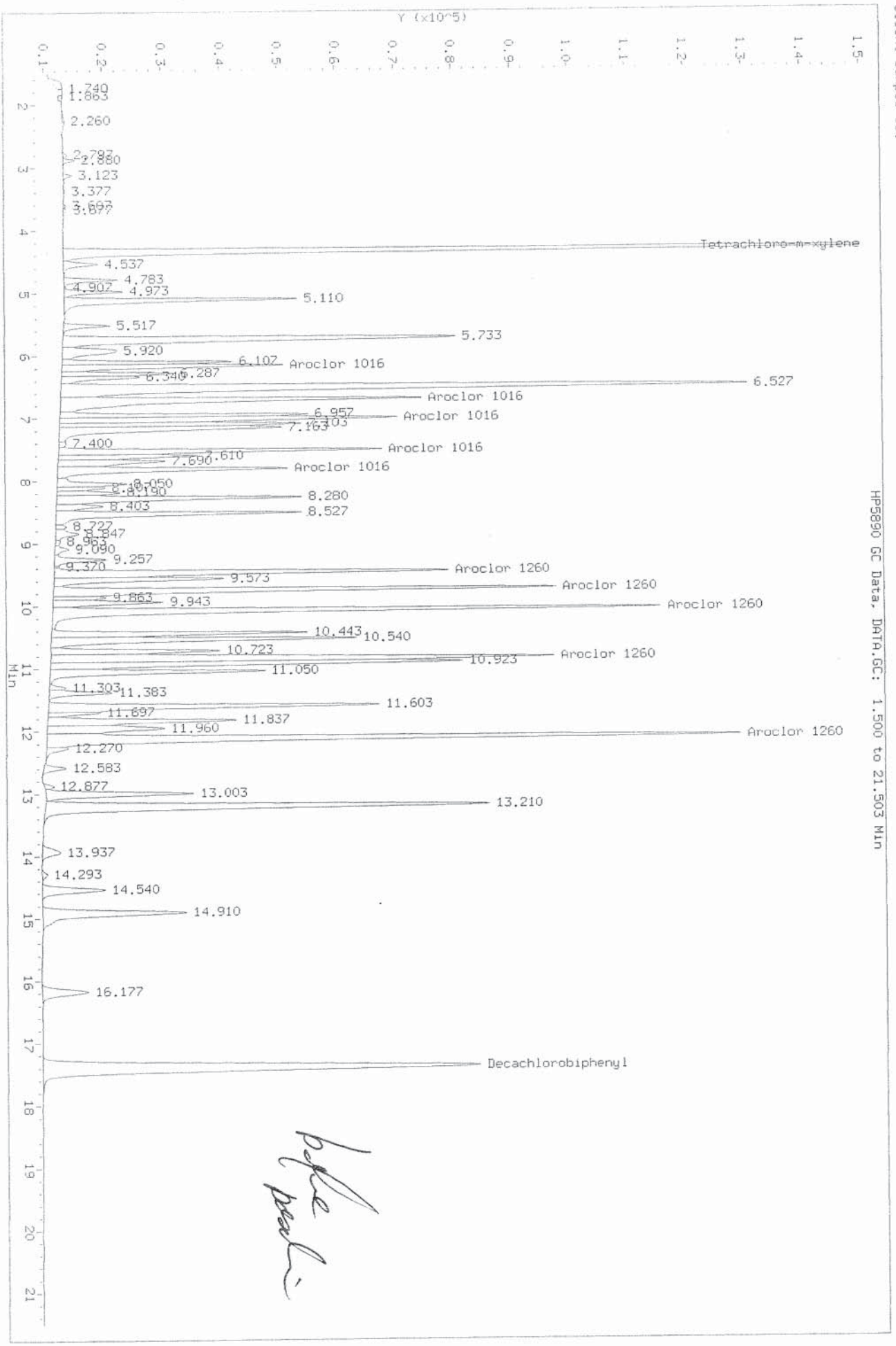
Instrument: GC09.i

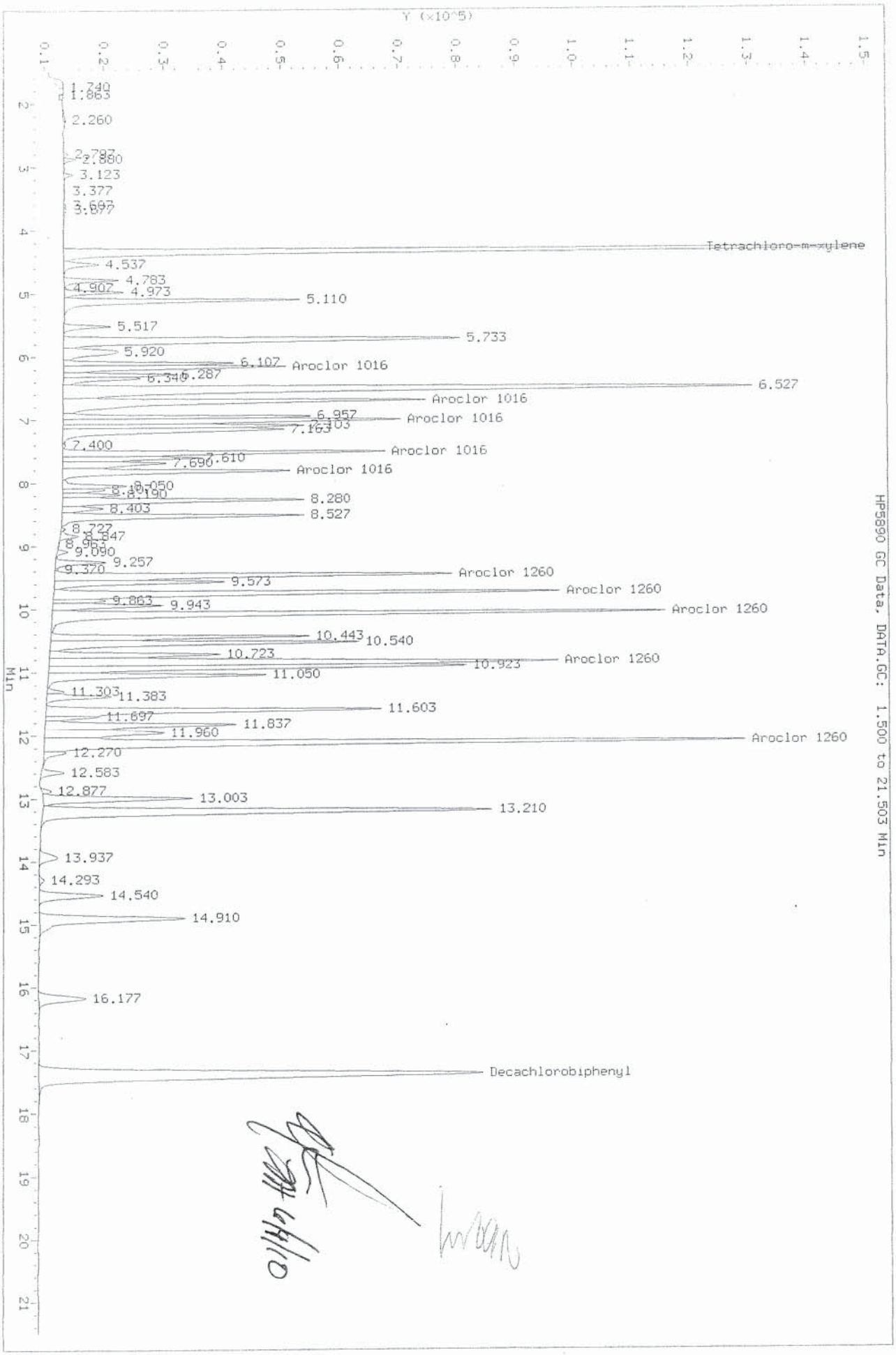
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53




Data File: \\Cash1\Acqudata\GC09\data\060810_r.b\0608R019.D
Injection Date: 09-JUN-2010 01:25
Instrument: GC09.1
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.503 Min
Y (x10^4)
Min
Tetrachloro-m-xylene
Aroclor 1016
Aroclor 1260
Decachlorobiphenyl


Data File: \\Cash1\Acqudata\GC09\data\060810_r.b\0608R019.D
Injection Date: 09-JUN-2010 01:25
Instrument: GC09.i
Client Sample ID:
HP5890 GC Data, DATA.GC: 1.500 to 21.503 Min
Y (x10^4)
Min
Tetrachloro-m-xylene
Aroclor 1016
Aroclor 1260
Decachlorobiphenyl


Y (x10^5)
Min
1.740
1.863
2.260
2.797
2.880
3.123
3.377
3.697
3.877
Tetrachloro-m-xylene
4.537
4.783
4.902
4.973
5.110
5.517
5.733
5.920
6.102
Aroclor 1016
6.287
6.340
6.527
Aroclor 1016
6.957
Aroclor 1016
7.103
7.163
7.400
Aroclor 1016
7.610
7.690
Aroclor 1016
8.050
8.190
8.280
8.403
8.527
8.727
8.847
8.963
9.090
9.257
9.320
Aroclor 1260
9.573
Aroclor 1260
9.863
9.943
Aroclor 1260
10.443
10.540
10.723
10.923
Aroclor 1260
11.050
11.303
11.383
11.603
11.697
11.837
11.960
Aroclor 1260
12.270
12.583
12.877
13.003
13.210
13.937
14.293
14.540
14.910
16.177
Decachlorobiphenyl
Data File: \\Cashl\Acqudata\GC09\data\060810.b\0608F019.D
Injection Date: 09-JUN-2010 01:25
Instrument: GC09.i
Client Sample ID:
HP6890 GC Data. DATA.GC: 1.500 to 21.503 Min


Y (x10^5)
0.1
0.2
0.3
0.4
0.5
0.6
0.7
0.8
0.9
1.0
1.1
1.2
1.3
1.4
1.5
Min
1
2
3
4
5
6
7
8
9
10
11
12
13
14
15
16
17
18
19
20
21
1.740
1.863
2.260
2.787
2.880
3.123
3.377
3.697
3.877
Tetrachloro-m-xylene
4.537
4.783
4.902
4.973
5.110
5.517
5.733
5.920
6.107
Aroclor 1016
6.287
6.340
6.527
Aroclor 1016
6.957
Aroclor 1016
7.163
7.303
7.400
Aroclor 1016
7.610
7.690
Aroclor 1016
8.050
8.190
8.280
8.403
8.527
8.722
8.947
9.090
9.257
9.370
Aroclor 1260
9.573
Aroclor 1260
9.863
9.943
Aroclor 1260
10.443
10.540
10.723
Aroclor 1260
10.923
11.050
11.303
11.383
11.603
11.697
11.837
11.960
Aroclor 1260
12.270
12.583
12.877
13.003
13.210
13.937
14.293
14.540
14.910
16.177
Decachlorobiphenyl
HP5890 GC Data, DATA.GC: 1.500 to 21.503 Min
Data File: \\Cash1\Acqudata\GC09\data\060810.b\0608F019.D
Injection Date: 09-JUN-2010 01:25
Instrument: GC09.i
Client Sample ID:

Organic Analysis:
Polychlorinated Biphenyls (PCBs)

Validation Package

Sample Prep and Screen Data

Preparation Information

Group ID:	KWG1005541	**Prep Method:**	EPA 3541	**Prep Date:**	06/02/10 12:00
Department:	Semivoa GC				

Lab Code	Client ID	Product	Matrix	Amt. Ext.	Final Vol.	Solids
K1005244-003	D-4-16	8082 PCB_LL	MISC. SOL	22.55g	4mL	
KWG1005541-1	Lab Control Sample	8082 PCB_LL	MISC. SOL	20.00g	4mL	
KWG1005541-2	Method Blank	8082 PCB_LL	MISC. SOL	22.78g	4mL	
KWG1005541-3	Matrix Spike	8082 PCB_LL	MISC. SOL	22.78g	4mL	
KWG1005541-4	Duplicate Matrix Spike	8082 PCB_LL	MISC. SOL	22.33g	4mL	

Lab Code	Parent Lab Code	Comments
KWG1005541-1		KQ1005123-01
KWG1005541-2		KQ1005123-03
KWG1005541-3	K1005244-003	KQ1005123-04
KWG1005541-4	K1005244-003	KQ1005123-05

Lab Code	Prep Event ID	Surrogate Solution ID	Amount Added	Spike Solution ID	Amount Added	Witness
K1005244-003	917270					SHeflin
KWG1005541-1	917271					SHeflin
KWG1005541-2	917272					SHeflin
KWG1005541-3	917273					SHeflin
KWG1005541-4	917274					SHeflin

Comments:

Started By: DWood Assisted By:

Completed By: km Assisted By:

Training: Yes / No (Yes) — Yes / No

Reviewed By: Date: 6/4/10 Storage:

Chain of Custody

Relinquished By: Date: 6-8-10

Received By: Date: 6/8/10

Extracts Examined: (Yes) / No

Columbia Analytical Services

Preparation Information Benchsheet

Prep Run: 112728 **Prep Workflow:** OrgExtS (14) **Status:** Prepped **Prep Date:** 06/02/2010 12:00
Team: Semivoa GC **Prep Method:** EPA 3541 **Current Step:** Final Volume **Due Date:** 06/01/2010
Analyst: DWood **Rush/NPDES:** N/A

Lab Code	Client ID	Bottle #	Target Amt	Initial Amt	Final Volume	TestNo List	Comments
K1005244-003	D-4-16	.03	22.05 g	22.55 g	4 mL	PCB_LL	
KQ1005123-01	Lab Control Sample		20.00 g	20.00 g	4 mL	PCB_LL	
KQ1005123-03	Method Blank		20.00 g	22.78 g	4 mL	PCB_LL	
K1005244-003: KQ1005123-04	Matrix Spike	.03	20.00 g	22.78 g	4 mL	PCB_LL	
K1005244-003: KQ1005123-05	Duplicate Matrix Spike	.03	20.00 g	22.33 g	4 mL	PCB_LL	

5 Total Samples consisting of 1 Client Sample, 2 Client QC Samples, 2 Batch QC Samples associated with the current Prep Run.

Spiking Solutions

Witness: SHeflin

Preparation Steps

Step	Started	Finished	By	Assisted By	Training?	Comments
Extraction	02-JUN-10 12:00	02-JUN-10 12:00	DWood		N	
Final Volume	08-JUN-10 12:00	08-JUN-10 12:00	km		N	

Comments

Review

Reviewed by: _____ Date: _____

Chain of Custody

Relinquished By: _____ Date: 6-8-10

Received By: _____ Date: 6/8/10

Extracts/Digestions Examined: Yes (circled) / No

Columbia Analytical Services™ **Preparation Information Benchsheet**

Prep Run: 112728 **Prep Workflow:** OrgExtS (14) **Status:** Draft **Prep Date:** 06/02/2010 05:46
Team: Semivoa GC **Prep Method:** EPA 3541 **Current Step:** Extraction **Due Date:** 06/01/2010
Analyst: DWOOD **Rush/NPDES:** N/A

Lab Code	Client ID	Bottle #	✓	Target Amt	Initial Amount	Inter. Volume	Final Volume	Surr Amt	Spike Amt	TestNo List
K1005244-003	D-4-16	.03		22.05 g	22.55	NA	4ml	200ul	NA	PCB_LL
KQ1005123-01	Lab Control Sample			20.00 g	20.00	↓	↓	↓	100ul	PCB_LL
KQ1005123-03	Method Blank			20.00 g	~~22.78~~	↓	↓	↓	NA	PCB_LL
K1005244-003: KQ1005123-04	Matrix Spike	.03		20.00 g	22.78	↓	↓	↓	100ul	PCB_LL
K1005244-003: KQ1005123-05	Duplicate Matrix Spike	.03		20.00 g	22.33	↓	↓	↓	↓	PCB_LL

5 Total Samples consisting of 1 Client Sample, 2 Client QC Samples, 2 Batch QC Samples associated with the current Prep Run.

Spiking Solutions **Witness:** SH 6-2-10

PCB5-50L 2ug/mL 200ul exp 11-26-10
PCB5-49L 40ug/mL 100ul exp 7-27-10

Preparation Steps

Step	Started	Finished	By	Assisted By	Training?	Comments
Extraction	6-2-10	6-8-10	sm for DW			
Final Volume	6-8-10	K				

Comments

Additional Prep Information for Pest/PCB in Soil by EPA 3541

Service Request # K1005244 Work Group # Pest: _____

PCB: KQ1005123

Extraction (3541):

Sulfate Lot # G2010025P Matrix Sand Lot # 1-14-10

DCM Lot # DA328

Soxtherm Start (Time/Date/Initial): ~~0928~~ 0900 6-2-10 DW SH

Soxtherm Stop (Time/Date/Initial): _____

Clean-ups:

GPC Clean-up (3640): — (Initial/Date)

Solvent Exchanged To Hexane (Initial/Date): 6-4-10 SM S-Evap Temp: —

Hexane Lot # 49343 N-Evap Temp: 35°C

Carbon Clean-up: — 4:1 Hexane/DCM Lot # _____

Florisil Clean-up (3620): — Florisil Lot # _____

1:1 Hexane/Acetone Lot # — 9:1 Hexane/Acetone Lot # _____

Sulfuric Acid Clean-up (3665): 6-4-10 SM Sulfuric Acid Lot # 47264

Sulfur Clean-up (3660M): 6-8-10 K Mercury Lot # 48219908

Pesticide Vials: — Vial Storage: _____

PCB Vials: Green Vial Storage: Group E1-5

Archived Extract Storage: Archived

Comments/Observations: _____

Bench Sheet Review Check List	
☐	Hold Times Met (if no, Reason: _____)
☐	Prep date, dept, method, product code correct in stealth
☐	Spike Information correct
☐	Weights/Volumes and units correct on raw and final bench sheets
☐	Sample IDs have been checked—Bottle numbers appended if required
☐	Names present for: Started by, Completed by, relinquished by, and witnessed by.
☐	Training has been circled
☐	Extract Storage recorded
☐	Additional Prep Sheet completely filled out (NA or line out Blanks)
☐	All clean-ups have been noted on additional prep sheet
☐	Signed service request with Form V, if applicable, has been attached

Sequence Name: D:\GC09\SEQUENCE\060810.S
Comment: PCB Aroclors by EPA 8082
Operator: LHarris
Data Path: D:\GC09\DATA\060810.B\
Pre-Seq Cmd:
Post-Seq Cmd:

Method Sections To Run
(X) Full Method
( ) Reprocessing Only

On A Barcode Mismatch
(X) Inject Anyway
( ) Don't Inject

10AL 9510

Run # 204050

Line	Type	Vial	DataFile	Method	Sample Name	
1	DeleteGC					
2	MaskName	-	----r---			
3	IB	1	0608F001	PCB_REG	IB	
4	IB	2	0608F002	PCB_REG	KWG1005547-1	IB | 8082 PCB
OK, OK 5	CCV	96	0608F003	PCB_REG	1660 @ 1000ppb	PCB5-55I | KWG100554
6	SMPL	3	0608F004	PCB_REG	K1005477-002	Influent 5/25
7	MB	4	0608F005	PCB_REG	KWG1005541-2	MB | 8082 PCB_LL | MIS
8	LCS	5	0608F006	PCB_REG	KWG1005541-1	LCS | 8082 PCB_LL | MI
9	SMPL	6	0608F007	PCB_REG	K1005244-003	D-4-16
10	MS	7	0608F008	PCB_REG	KWG1005541-3	K1005244-003MS | 8082
11	DMS	8	0608F009	PCB_REG	KWG1005541-4	K1005244-003DMS | 8082
12	MB	9	0608F010	PCB_REG	KWG1005543-4	MB | 8082 PCB | SEDIME
13	LCS	10	0608F011	PCB_REG	KWG1005543-3	LCS | 8082 PCB | SEDIM
14	SMPL	11	0608F012	PCB_REG	K1005351-010	SJNE 050-GR1
15	MS	12	0608F013	PCB_REG	KWG1005543-1	K1005351-010MS | 8082
16	DMS	13	0608F014	PCB_REG	KWG1005543-2	K1005351-010DMS | 8082
17	SMPL	14	0608F015	PCB_REG	K1005351-015	SJNE 067-GR1
18	SMPL	15	0608F016	PCB_REG	K1005351-016	SJNE 069-GR1
19	IB	1	0608F017	PCB_REG	IB	
20	IB	2	0608F018	PCB_REG	KWG1005547-3	IB | 8082 PCB
X, OK 21	CCV	96	0608F019	PCB_REG	1660 @ 1000ppb	PCB5-55I | KWG100554
22	SMPL	16	0608F020	PCB_REG	K1005351-017	SJNE 070-GR1
23	SMPL	17	0608F021	PCB_REG	K1005351-021	SJNE 066-GR1
24	SMPL	18	0608F022	PCB_REG	K1005353-004	SJGB 013 S4
25	SMPL	19	0608F023	PCB_REG	K1005353-006	SJGB 014 S1 NR, need 20X
26	SMPL	20	0608F024	PCB_REG	K1005353-010	SJGB 014 S5
27	SMPL	21	0608F025	PCB_REG	K1005353-015	SJGB 015 S5
28	SMPL	22	0608F026	PCB_REG	K1005353-018	SJGB 016 S3
29	SMPL	23	0608F027	PCB_REG	K1005353-024	SJGB 017 S5
30	SMPL	24	0608F028	PCB_REG	K1005423-001	ACS-6
31	SMPL	25	0608F029	PCB_REG	K1005423-002	ACS-7
32	IB	1	0608F030	PCB_REG	IB	
33	IB	2	0608F031	PCB_REG	KWG1005547-5	IB | 8082 PCB
C, OK 34	CCV	96	0608F032	PCB_REG	1660 @ 1000ppb	PCB5-55I | KWG100554

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F002.D	**Date Acquired:**	06/08/2010 18:01
Lab ID:	KWG1005547-1	**Date Quantitated:**	06/09/2010 10:28
RunType:	IB	**Batch ID:**	KWG1005547
Matrix:	NOT APPLICABLE	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: MG 6/9/10

Secondary Review: 6/10/10

Printed: 06/09/2010 11:00:06

u:\Stealth\Crystal.rpt\except2.rpt

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R002.D	Date Acquired:	06/08/2010 18:01
Lab ID:	KWG1005547-1	Date Quantitated:	06/09/2010 10:29
RunType:	IB	Batch ID:	KWG1005547
Matrix:	NOT APPLICABLE	Analysis Method:	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: MJM

Secondary Review: 6/10/2010

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	NOT APPLICABL
Prod Code:	8082 PCB	**Collect Date:**		**Receive Date:**	06/08/2010

Analysis Lot:	KWG1005547	**Prep Lot:**		**Report Group:**	
Analysis Method:	8082	**Prep Method:**			
Prep Ref:		**Prep Date:**			

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	**Calibration ID:**	CAL9510
Title:			
MB Ref:		**Method ID:**	MJ731
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F002.D			**Instrument:**	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R002.D			**Vial:**	2
Acqu Date:	06/08/2010 18:01	**Quant Date:**	06/09/2010 10:28	**Dilution:**	1.0
Run Type:	IB			**Soln Conc. Units:**	ng/mL
Lab ID:	KWG1005547-1				
Signal #1:	DB-35MS	**Signal #2:**	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	0.00		0d	0		0.0000			NA
			%Recovery =		NA	NA	Limits =	10-135	
Decachlorobiphenyl	0.00		0	0		0.0000			NA
			%Recovery =		NA	NA	Limits =	35-133	

Target Compounds

Final Conc. Units:

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	mg/Kg #1	mg/Kg #2	Rpt
Aroclor 1016			0	0	0.0000	0.0000			
Aroclor 1016 {1}			0	0	0.0000	0.0000			
Aroclor 1016 {2}			0	0	0.0000	0.0000			
Aroclor 1016 {3}			0	0	0.0000	0.0000			
Aroclor 1016 {4}			0	0	0.0000	0.0000			
Aroclor 1016 {5}			0	0	0.0000	0.0000			
Aroclor 1221			0	0	0.0000	0.0000			
Aroclor 1221 {1}			0	0	0.0000	0.0000			
Aroclor 1221 {2}			0	0	0.0000	0.0000			
Aroclor 1221 {3}			0	0	0.0000	0.0000			
Aroclor 1221 {4}			0	0	0.0000	0.0000			
Aroclor 1232			0	0	0.0000	0.0000			
Aroclor 1232 {1}			0	0	0.0000	0.0000			
Aroclor 1232 {2}			0	0	0.0000	0.0000			
Aroclor 1232 {3}			0	0	0.0000	0.0000			
Aroclor 1232 {4}			0	0	0.0000	0.0000			
Aroclor 1242			0	0	0.0000	0.0000			
Aroclor 1242 {1}			0	0	0.0000	0.0000			

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F002.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R002.D			Vial:	2
Acqu Date:	06/08/2010 18:01	Quant Date:	06/09/2010 10:28	Dilution:	1.0
Run Type:	IB			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005547-1				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds — Final Conc. Units:

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	mg/Kg #1	mg/Kg #2	Rpt
Aroclor 1242 {2}			0	0	0.0000	0.0000			
Aroclor 1242 {3}			0	0	0.0000	0.0000			
Aroclor 1242 {4}			0	0	0.0000	0.0000			
Aroclor 1242 {5}			0	0	0.0000	0.0000			
Aroclor 1248			0	0	0.0000	0.0000			
Aroclor 1248 {1}			0	0	0.0000	0.0000			
Aroclor 1248 {2}			0	0	0.0000	0.0000			
Aroclor 1248 {3}			0	0	0.0000	0.0000			
Aroclor 1248 {4}			0	0	0.0000	0.0000			
Aroclor 1248 {5}			0	0	0.0000	0.0000			
Aroclor 1254			0	0	0.0000	0.0000			
Aroclor 1254 {1}			0	0	0.0000	0.0000			
Aroclor 1254 {2}			0	0	0.0000	0.0000			
Aroclor 1254 {3}			0	0	0.0000	0.0000			
Aroclor 1254 {4}			0	0	0.0000	0.0000			
Aroclor 1254 {5}			0	0	0.0000	0.0000			
Aroclor 1260			0	0	0.0000	0.0000			
Aroclor 1260 {1}			0	0	0.0000	0.0000			
Aroclor 1260 {2}			0	0	0.0000	0.0000			
Aroclor 1260 {3}			0	0	0.0000	0.0000			
Aroclor 1260 {4}			0	0	0.0000	0.0000			
Aroclor 1260 {5}			0	0	0.0000	0.0000			
Aroclor 1262			0	0	0.0000	0.0000			
Aroclor 1262 {1}			0	0	0.0000	0.0000			
Aroclor 1262 {2}			0	0	0.0000	0.0000			
Aroclor 1262 {3}			0	0	0.0000	0.0000			
Aroclor 1262 {4}			0	0	0.0000	0.0000			
Aroclor 1262 {5}			0	0	0.0000	0.0000			
Aroclor 1268			0	0	0.0000	0.0000			
Aroclor 1268 {1}			0	0	0.0000	0.0000			
Aroclor 1268 {2}			0	0	0.0000	0.0000			
Aroclor 1268 {3}			0	0	0.0000	0.0000			
Aroclor 1268 {4}			0	0	0.0000	0.0000			
Aroclor 1268 {5}			0	0	0.0000	0.0000			

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F002.D
Report Date: 09-Jun-2010 10:28

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F002.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R002.D
Inj Date : 08-JUN-2010 18:01
Sample Info: KWG1005547-1 | IB | 8082 PCB
Misc Info : SEMIVOA GC\W1005547\1-IB.H
Cal Date : 08-JUN-2010 12:20
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

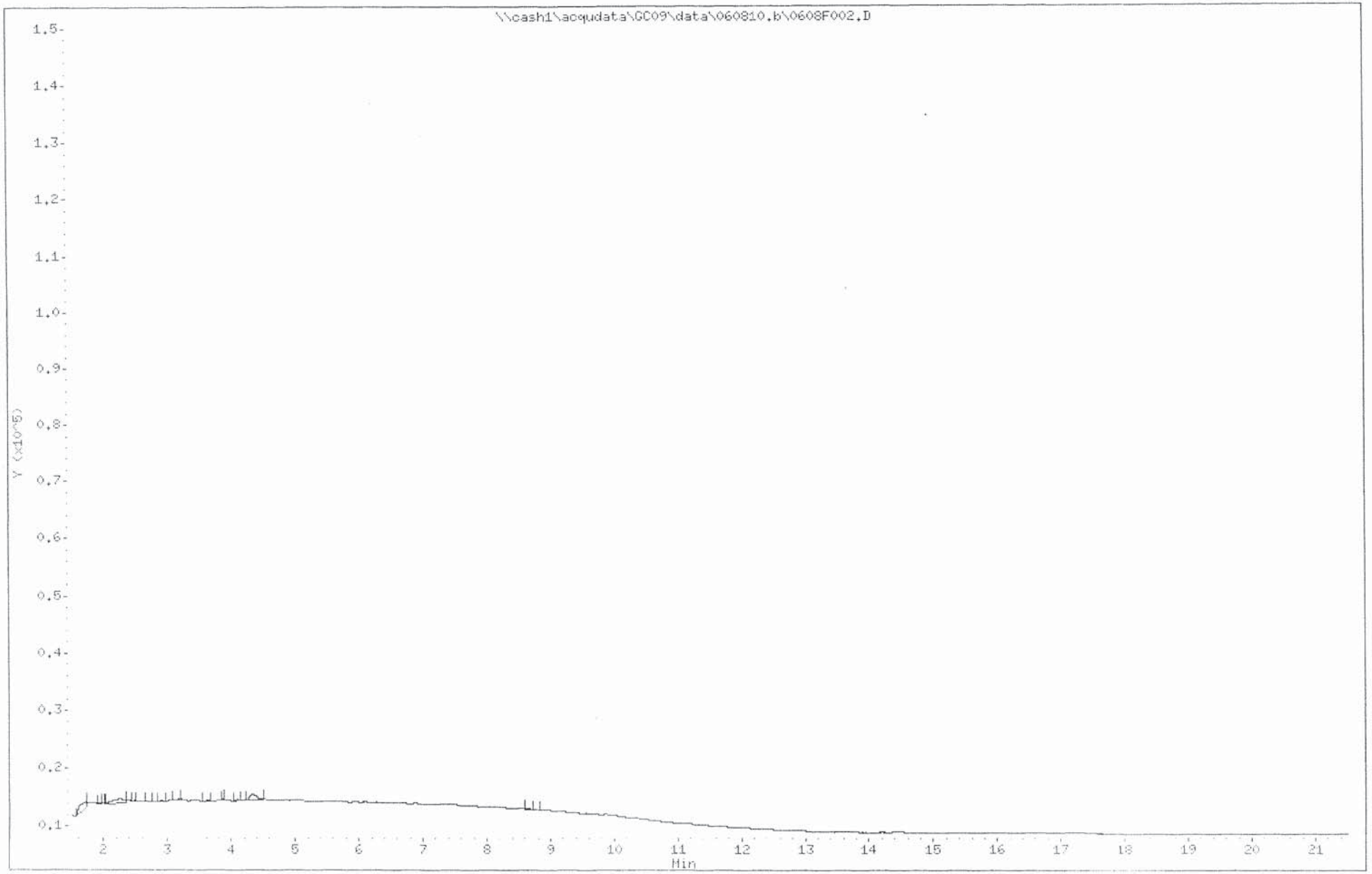
Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F002.D
Date : 08-JUN-2010 18:01
Client ID:
Sample Info: KWG1005547-1 | IB | 8082 PCB

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-35MS

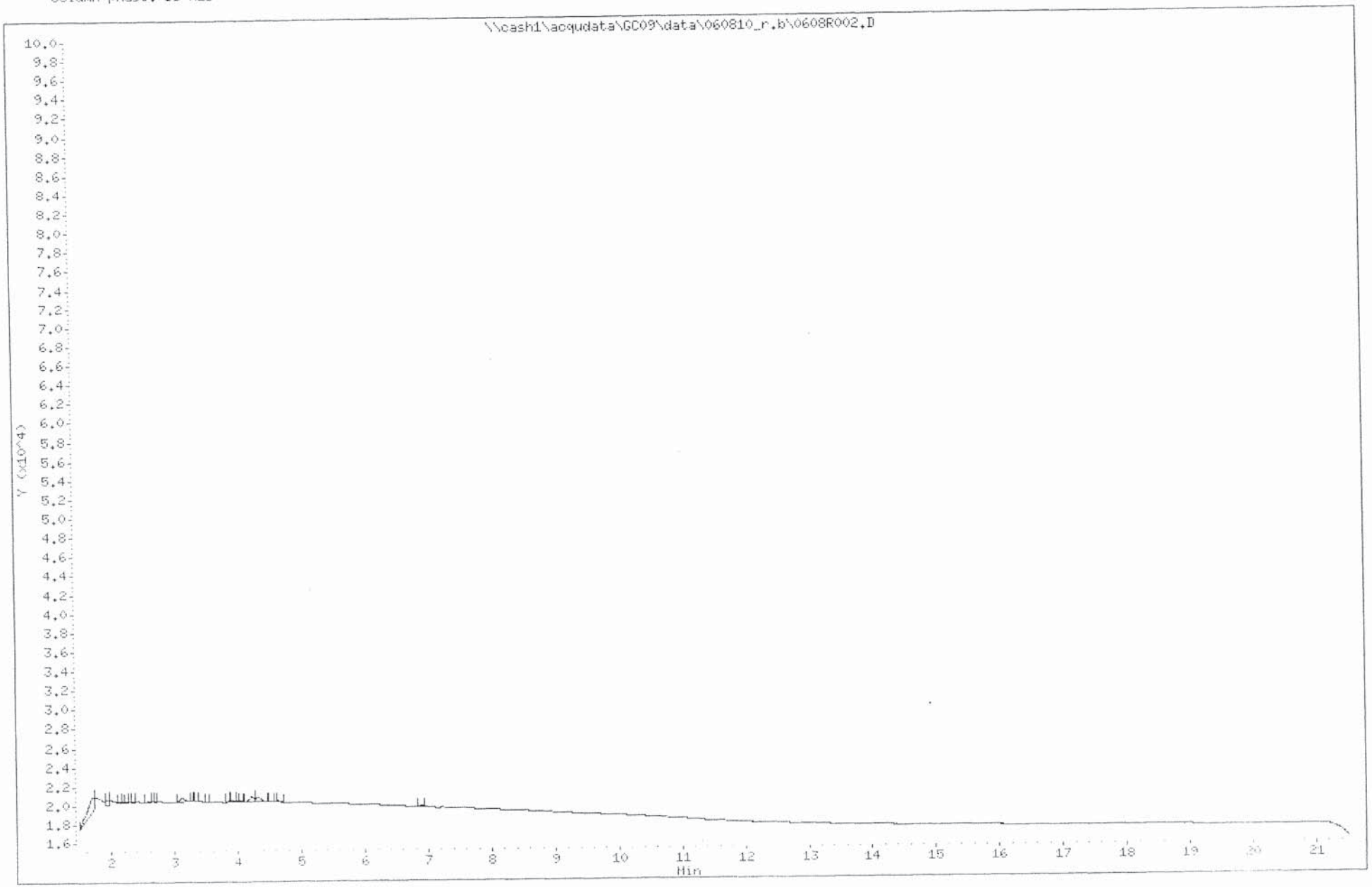


Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R002.D
Date : 08-JUN-2010 18:01
Client ID:
Sample Info: KWG1005547-1 | IB | 8082 PCB

Instrument: GC09.i

Operator: LHarris
Column diameter: 0.53

Column phase: DB-XLB



Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\0608F018.D	**Date Acquired:**	06/09/2010 00:59
Lab ID:	KWG1005547-3	**Date Quantitated:**	06/09/2010 10:29
RunType:	IB	**Batch ID:**	KWG1005547
Matrix:	NOT APPLICABLE	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: MMD

Secondary Review: 6/6/2010

Exception Report

Data File:	\\CASH1\ACQUDATA\GC09\DATA\060810_R.B\0608R018.D	**Date Acquired:**	06/09/2010 00:59
Lab ID:	KWG1005547-3	**Date Quantitated:**	06/09/2010 10:30
RunType:	IB	**Batch ID:**	KWG1005547
Matrix:	NOT APPLICABLE	**Analysis Method:**	8082
		MethodJoinID:	MJ579

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
ICAL Analyte Recovery	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	

Primary Review: [signature]

Secondary Review: [signature] 6/15/10

Quantitation Report

Bottle ID:		Tier:		Matrix:	NOT APPLICABL
Prod Code:	8082 PCB	Collect Date:		Receive Date:	06/08/2010

Analysis Lot:	KWG1005547	Prep Lot:	Report Group:
Analysis Method:	8082	Prep Method:	
Prep Ref:		Prep Date:	

Quant Method:	\\CASH1\ACQUDATA\GC09\DATA\060810.B\052710A_F.M	Calibration ID:	CAL9510
Title:			
MB Ref:		Method ID:	MJ579
		Quant based on Method	

Data File #1:	J:\GC09\DATA\060810.B\0608F018.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R018.D			Vial:	2
Acqu Date:	06/09/2010 00:59	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	IB			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005547-3				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Surrogate Compounds

Parameter Name	RT #1	RT #2	Resp #1	Respe #2	ng/mL #1	ng/mL #2			Rpt
Tetrachloro-m-xylene	0.00		0	0d		0.0000			NA
			%Recovery =		NA	NA	Limits =	21-114	
Decachlorobiphenyl	0.00		0	0		0.0000			NA
			%Recovery =		NA	NA	Limits =	36-113	

Target Compounds

Final Conc. Units: ug/L

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/L #1	ug/L #2	Rpt
Aroclor 1016			0	0	0.0000	0.0000			
Aroclor 1016 {1}			0	0	0.0000	0.0000			
Aroclor 1016 {2}			0	0	0.0000	0.0000			
Aroclor 1016 {3}			0	0	0.0000	0.0000			
Aroclor 1016 {4}			0	0	0.0000	0.0000			
Aroclor 1016 {5}			0	0	0.0000	0.0000			
Aroclor 1221			0	0	0.0000	0.0000			
Aroclor 1221 {1}			0	0	0.0000	0.0000			
Aroclor 1221 {2}			0	0	0.0000	0.0000			
Aroclor 1221 {3}			0	0	0.0000	0.0000			
Aroclor 1221 {4}			0	0	0.0000	0.0000			
Aroclor 1232			0	0	0.0000	0.0000			
Aroclor 1232 {1}			0	0	0.0000	0.0000			
Aroclor 1232 {2}			0	0	0.0000	0.0000			
Aroclor 1232 {3}			0	0	0.0000	0.0000			
Aroclor 1232 {4}			0	0	0.0000	0.0000			
Aroclor 1242			0	0	0.0000	0.0000			
Aroclor 1242 {1}			0	0	0.0000	0.0000			

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File #1:	J:\GC09\DATA\060810.B\0608F018.D			Instrument:	GC09.i
Data File #2:	\\cash1\acqudata\GC09\data\060810_r.b\0608R018.D			Vial:	2
Acqu Date:	06/09/2010 00:59	Quant Date:	06/09/2010 10:29	Dilution:	1.0
Run Type:	IB			Soln Conc. Units:	ng/mL
Lab ID:	KWG1005547-3				
Signal #1:	DB-35MS	Signal #2:	DB-XLB		

Target Compounds

Final Conc. Units: ug/L

Parameter Name	RT #1	RT #2	Resp #1	Resp #2	ng/mL #1	ng/mL #2	ug/L #1	ug/L #2	Rpt
Aroclor 1242 {2}			0	0	0.0000	0.0000			
Aroclor 1242 {3}			0	0	0.0000	0.0000			
Aroclor 1242 {4}			0	0	0.0000	0.0000			
Aroclor 1242 {5}			0	0	0.0000	0.0000			
Aroclor 1248			0	0	0.0000	0.0000			
Aroclor 1248 {1}			0d	0	0.0000	0.0000			
Aroclor 1248 {2}			0d	0	0.0000	0.0000			
Aroclor 1248 {3}			0d	0	0.0000	0.0000			
Aroclor 1248 {4}			0d	0	0.0000	0.0000			
Aroclor 1248 {5}			0d	0	0.0000	0.0000			
Aroclor 1254			0	0	0.0000	0.0000			
Aroclor 1254 {1}			0	0	0.0000	0.0000			
Aroclor 1254 {2}			0	0	0.0000	0.0000			
Aroclor 1254 {3}			0	0	0.0000	0.0000			
Aroclor 1254 {4}			0	0	0.0000	0.0000			
Aroclor 1254 {5}			0	0	0.0000	0.0000			
Aroclor 1260			0	0	0.0000	0.0000			
Aroclor 1260 {1}			0	0	0.0000	0.0000			
Aroclor 1260 {2}			0	0	0.0000	0.0000			
Aroclor 1260 {3}			0	0	0.0000	0.0000			
Aroclor 1260 {4}			0	0	0.0000	0.0000			
Aroclor 1260 {5}			0	0	0.0000	0.0000			
Aroclor 1262			0	0	0.0000	0.0000			
Aroclor 1262 {1}			0	0	0.0000	0.0000			
Aroclor 1262 {2}			0	0	0.0000	0.0000			
Aroclor 1262 {3}			0	0	0.0000	0.0000			
Aroclor 1262 {4}			0	0	0.0000	0.0000			
Aroclor 1262 {5}			0	0	0.0000	0.0000			
Aroclor 1268			0	0	0.0000	0.0000			
Aroclor 1268 {1}			0	0	0.0000	0.0000			
Aroclor 1268 {2}			0	0	0.0000	0.0000			
Aroclor 1268 {3}			0	0	0.0000	0.0000			
Aroclor 1268 {4}			0	0	0.0000	0.0000			
Aroclor 1268 {5}			0	0	0.0000	0.0000			

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File: \\cash1\acqudata\GC09\data\060810.b\0608F018.D
Report Date: 09-Jun-2010 10:29

Laboratory Name

Sample #1 : \\cash1\acqudata\GC09\data\060810.b\0608F018.D
Sample #2 : \\cash1\acqudata\GC09\data\060810_r.b\0608R018.D
Inj Date : 09-JUN-2010 00:59
Sample Info: KWG1005547-3 | IB | 8082 PCB
Misc Info : SEMIVOA GC\W1005547\3-IB.H
Cal Date : 09-JUN-2010 08:48
Operator : LHarris
Inst ID : GC09.i
Dil Factor : 1.000000

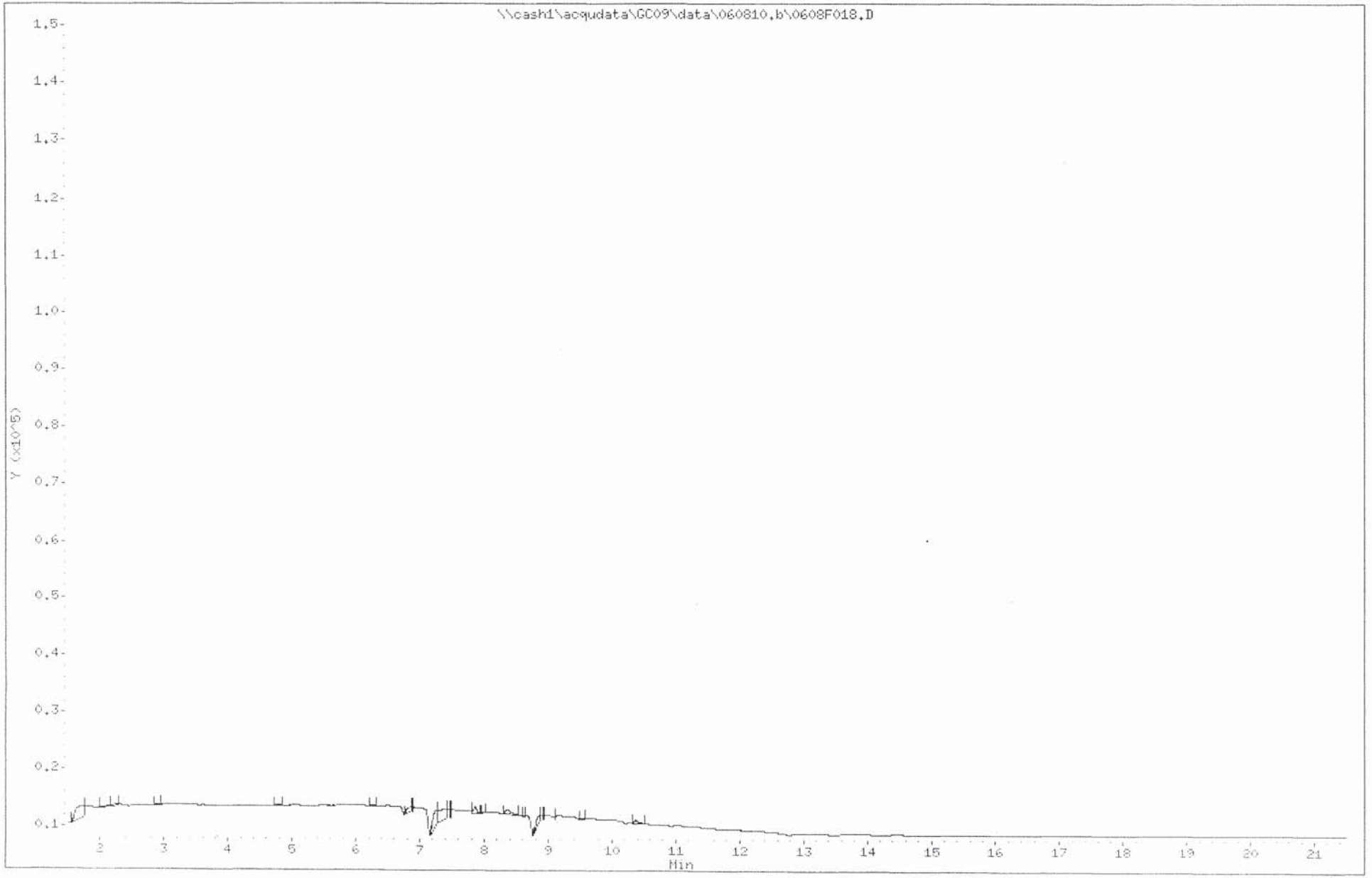
Method #1 : \\Cash1\Acqudata\GC09\data\060810.b\052710A_f.m
Method #2 : \\Cash1\Acqudata\GC09\data\060810_r.b\052710A_r.m
Sub List #1 : ALL.SUB
Sub List #2 : ALL.SUB
Col #1 Phase : DB-35MS
Col #2 Phase : DB-XLB

Compound	RT#1	RT#2	Resp#1	Resp#2	Conc#1	Conc#2	Target Range	Ratio

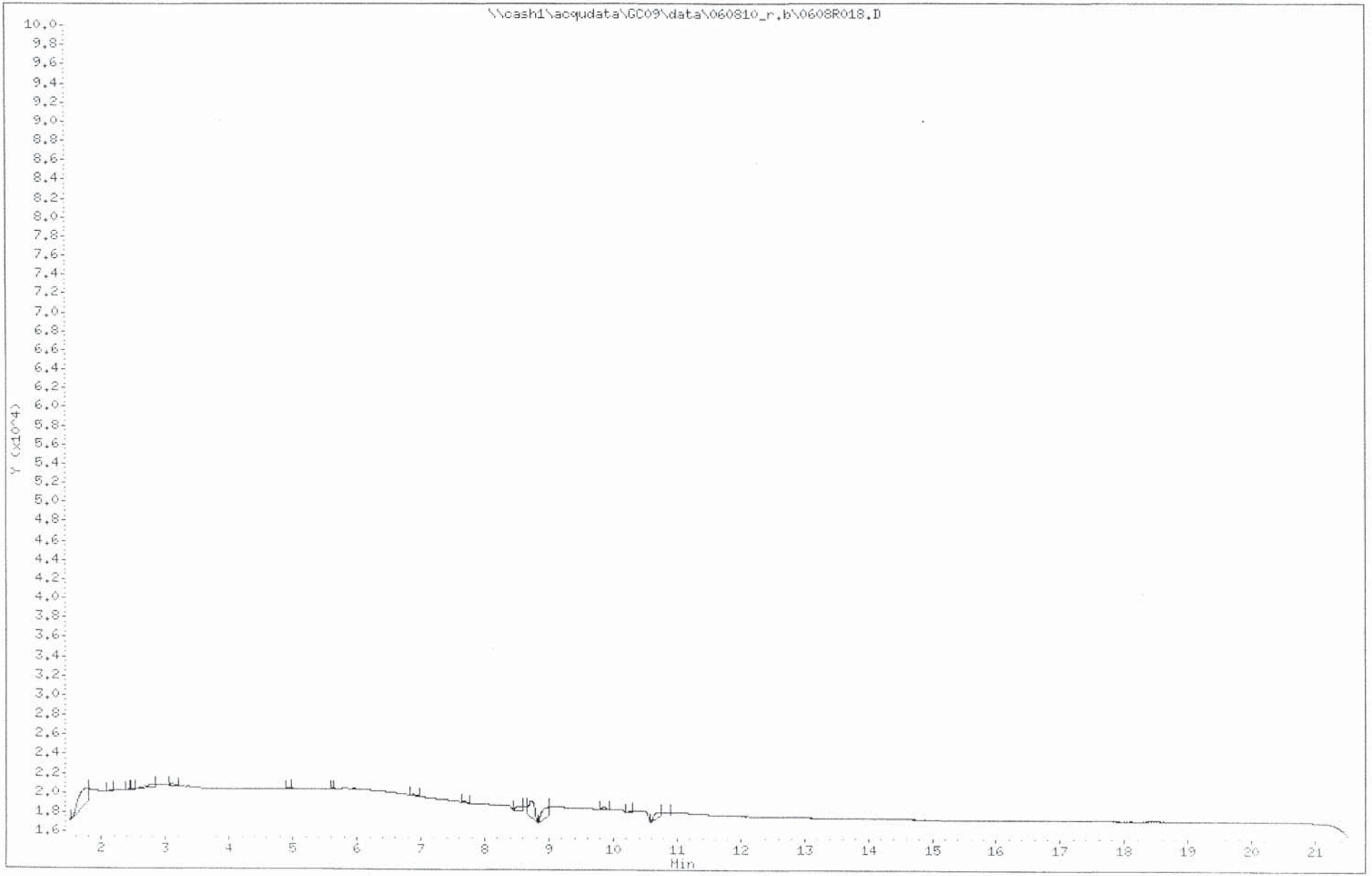
Data File: \\cash1\acqudata\GC09\data\060810.b\0608F018.D
Date : 09-JUN-2010 00:59
Client ID:
Sample Info: KWG1005547-3 | IB | 8082 PCB

Instrument: GC09.i

Operator: LHarris
Column phase: DB-35MS
Column diameter: 0.53



Data File: \\cash1\acqudata\GC09\data\060810_r.b\0608R018.D
Date : 09-JUN-2010 00:59
Client ID:
Instrument: GC09.i
Sample Info: KWG1005547-3 | IB | 8082 PCB
Operator: LHarris
Column phase: DB-XLB
Column diameter: 0.53



Volatile Organic Compounds

Organic Analysis:
Volatile Organic Compounds

Summary Package

Sample and QC Results

COLUMBIA ANALYTICAL SERVICES, INC.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244

Cover Page - Organic Analysis Data Package
Volatile Organic Compounds

Sample Name	Lab Code	Date Collected	Date Received
D-4-16	K1005244-003	05/19/2010	05/21/2010
D-4-16MS	KWG1004900-1	05/19/2010	05/21/2010
D-4-16DMS	KWG1004900-2	05/19/2010	05/21/2010

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed in the case narrative. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on floppy diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: K Krasnow Name: K Krasnow

Date: 6/11/10 Title: Scientist

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16
Lab Code: K1005244-003

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	ND	U	5.6	0.51	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	ND	U	5.6	0.35	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	ND	U	5.6	0.69	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	ND	U	5.6	0.35	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	ND	U	5.6	0.24	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	ND	U	5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
Acetone	**13**	J	23	4.3	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	ND	U	5.6	0.29	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**1.7**	J	12	0.48	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	ND	U	5.6	0.39	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	ND	U	5.6	0.23	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	ND	U	5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	ND	U	5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**4.0**	J	23	1.1	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	ND	U	5.6	0.16	1	05/25/10	05/25/10	KWG1004900	
Chloroform	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	ND	U	5.6	0.39	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	ND	U	5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	ND	U	5.6	0.30	1	05/25/10	05/25/10	KWG1004900	
Benzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	ND	U	5.6	0.17	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	ND	U	5.6	0.30	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	ND	U	5.6	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	ND	U	5.6	0.17	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	ND	U	23	0.82	1	05/25/10	05/25/10	KWG1004900	
Toluene	ND	U	5.6	0.20	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	ND	U	5.6	0.38	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	ND	U	5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	ND	U	5.6	0.29	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	ND	U	23	1.1	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	ND	U	5.6	0.23	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16
Lab Code: K1005244-003

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	ND	U	5.6	0.21	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	ND	U	23	0.30	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	ND	U	5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	ND	U	5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	ND	U	5.6	0.41	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	ND	U	5.6	0.15	1	05/25/10	05/25/10	KWG1004900	
Styrene	ND	U	5.6	0.15	1	05/25/10	05/25/10	KWG1004900	
Bromoform	ND	U	5.6	0.34	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	ND	U	23	0.15	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	ND	U	5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	ND	U	23	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	ND	U	5.6	0.42	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	ND	U	23	0.32	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	ND	U	23	0.40	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	ND	U	23	0.29	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	ND	U	23	0.37	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	ND	U	23	0.31	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	ND	U	23	0.35	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	ND	U	23	0.31	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	ND	U	5.6	0.27	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	ND	U	5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	ND	U	23	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	ND	U	23	0.73	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	ND	U	23	0.26	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	ND	U	23	0.24	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	ND	U	23	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	ND	U	23	0.24	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Collected:**	05/19/2010
Sample Matrix:	Misc. solid	**Date Received:**	05/21/2010

Volatile Organic Compounds

Sample Name:	D-4-16	**Units:**	ug/Kg
Lab Code:	K1005244-003	**Basis:**	Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	100	83-128	05/25/10	Acceptable
Toluene-d8	108	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	103	77-124	05/25/10	Acceptable

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Method Blank
Lab Code: KWG1004900-5

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	ND	U	5.0	0.46	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	ND	U	5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	ND	U	5.0	0.62	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	ND	U	5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	ND	U	5.0	0.21	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	ND	U	5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Acetone	**6.3**	J	20	3.9	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	ND	U	5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**1.5**	J	10	0.43	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	ND	U	5.0	0.35	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	ND	U	5.0	0.20	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	ND	U	5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	ND	U	5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**2.3**	J	20	0.99	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	ND	U	5.0	0.14	1	05/25/10	05/25/10	KWG1004900	
Chloroform	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	ND	U	5.0	0.35	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	ND	U	5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	ND	U	5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
Benzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	ND	U	5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	ND	U	5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	ND	U	5.0	0.28	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	ND	U	5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	ND	U	20	0.74	1	05/25/10	05/25/10	KWG1004900	
Toluene	ND	U	5.0	0.18	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	ND	U	5.0	0.34	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	ND	U	5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	ND	U	5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	ND	U	20	0.93	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	ND	U	5.0	0.20	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Method Blank
Lab Code: KWG1004900-5

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	ND	U	5.0	0.19	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	ND	U	20	0.27	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	ND	U	5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	ND	U	5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	ND	U	5.0	0.37	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	ND	U	5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Styrene	ND	U	5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Bromoform	ND	U	5.0	0.30	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	ND	U	20	0.13	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	ND	U	5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	ND	U	20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	ND	U	5.0	0.38	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	ND	U	20	0.29	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	ND	U	20	0.36	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	ND	U	20	0.26	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	ND	U	20	0.33	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	ND	U	20	0.28	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	ND	U	20	0.31	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	ND	U	20	0.28	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	ND	U	5.0	0.24	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	ND	U	5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	ND	U	20	0.32	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	ND	U	20	0.66	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	ND	U	20	0.23	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	ND	U	20	0.21	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	ND	U	20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	ND	U	20	0.21	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Collected:**	NA
Sample Matrix:	Misc. solid	**Date Received:**	NA

Volatile Organic Compounds

Sample Name:	Method Blank	**Units:**	ug/Kg
Lab Code:	KWG1004900-5	**Basis:**	Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	100	83-128	05/25/10	Acceptable
Toluene-d8	110	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	104	77-124	05/25/10	Acceptable

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244

Surrogate Recovery Summary
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: PERCENT
Level: Low

Sample Name	Lab Code	Sur1	Sur2	Sur3
D-4-16	K1005244-003	100	108	103
Method Blank	KWG1004900-5	100	110	104
D-4-16MS	KWG1004900-1	99	109	108
D-4-16DMS	KWG1004900-2	102	114	107
Lab Control Sample	KWG1004900-3	94	105	110
Duplicate Lab Control Sample	KWG1004900-4	100	115	109

Surrogate Recovery Control Limits (%)

Sur1 =	Dibromofluoromethane	83-128
Sur2 =	Toluene-d8	90-125
Sur3 =	4-Bromofluorobenzene	77-124

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 05/25/2010
Time Analyzed: 10:34

Internal Standard Area and RT Summary
Volatile Organic Compounds

File ID: J:\MS24\DATA\052510\0525F004.D
Instrument ID: MS24
Analysis Method: 8260B

Lab Code: KWG1004880-2
Analysis Lot: KWG1004880

		Fluorobenzene		Chlorobenzene-d5		1,4-Dichlorobenzene-d4	
		Area	**RT**	**Area**	**RT**	**Area**	**RT**
	Results ==>	367,346	4.02	157,615	6.65	137,921	8.98
	Upper Limit ==>	734,692	4.52	315,230	7.15	275,842	9.48
	Lower Limit ==>	183,673	3.52	78,808	6.15	68,961	8.48
	ICAL Result ==>	321,724	4.02	137,440	6.65	130,674	8.98
Associated Analyses							
Lab Control Sample	KWG1004900-3	418,826	4.02	167,585	6.65	148,678	8.98
Duplicate Lab Control Sample	KWG1004900-4	355,415	4.02	159,438	6.65	144,430	8.98
Method Blank	KWG1004900-5	348,737	4.02	149,494	6.65	128,420	8.98
D-4-16	K1005244-003	342,019	4.02	146,554	6.65	116,925	8.98
D-4-16MS	KWG1004900-1	420,529	4.02	177,982	6.65	145,052	8.98
D-4-16DMS	KWG1004900-2	367,092	4.02	163,841	6.65	139,592	8.98

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	05/25/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	05/25/2010

Matrix Spike/Duplicate Matrix Spike Summary
Volatile Organic Compounds

Sample Name:	D-4-16	**Units:**	ug/Kg
Lab Code:	K1005244-003	**Basis:**	Dry
Extraction Method:	EPA 5030A	**Level:**	Low
Analysis Method:	8260B	**Extraction Lot:**	KWG1004900

Analyte Name	Sample Result	D-4-16MS KWG1004900-1 Matrix Spike			D-4-16DMS KWG1004900-2 Duplicate Matrix Spike			%Rec Limits	RPD	RPD Limit
		Result	Expected	%Rec	Result	Expected	%Rec			
Vinyl Chloride	ND	49.8	54.8	91	55.8	55.1	101	12-160	11	40
1,1-Dichloroethene	ND	44.0	54.8	80	49.5	55.1	90	31-153	12	40
Chloroform	ND	46.5	54.8	85	49.8	55.1	90	30-138	7	40
Carbon Tetrachloride	ND	47.8	54.8	87	51.3	55.1	93	10-144	7	40
Benzene	ND	46.7	54.8	85	49.3	55.1	89	30-137	5	40
Trichloroethene (TCE)	ND	51.6	54.8	94	52.9	55.1	96	18-145	3	40
Bromodichloromethane	ND	46.1	54.8	84	52.1	55.1	95	14-146	12	40
Toluene	ND	45.9	54.8	84	50.7	55.1	92	15-135	10	40
1,1,2-Trichloroethane	ND	48.5	54.8	88	51.7	55.1	94	18-141	7	40
2-Hexanone	ND	527	274	192 *	532	276	193 *	20-128	1	40
Chlorobenzene	ND	44.0	54.8	80	45.7	55.1	83	15-124	4	40
Ethylbenzene	ND	46.4	54.8	85	48.5	55.1	88	13-128	4	40
1,2,3-Trichloropropane	ND	71.0	54.8	130	68.7	55.1	125	10-159	3	40
2-Chlorotoluene	ND	52.7	54.8	96	51.6	55.1	94	10-140	2	40
1,2-Dichlorobenzene	ND	45.3	54.8	83	45.7	55.1	83	10-124	1	40
Naphthalene	ND	44.4	54.8	81	42.8	55.1	78	10-127	4	40

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Extracted: 05/25/2010
Date Analyzed: 05/25/2010

Lab Control Spike/Duplicate Lab Control Spike Summary
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry
Level: Low
Extraction Lot: KWG1004900

	Lab Control Sample KWG1004900-3 Lab Control Spike				Duplicate Lab Control Sample KWG1004900-4 Duplicate Lab Control Spike					
Analyte Name	**Result**	**Expected**	**%Rec**		**Result**	**Expected**	**%Rec**	**%Rec Limits**	**RPD**	**RPD Limit**
Dichlorodifluoromethane	36.3	50.0	73		42.5	50.0	85	38-160	16	40
Chloromethane	39.1	50.0	78		44.8	50.0	90	37-146	14	40
Vinyl Chloride	40.0	50.0	80		45.8	50.0	92	54-127	14	40
Bromomethane	26.7	50.0	53		31.0	50.0	62	22-180	15	40
Chloroethane	37.0	50.0	74		43.1	50.0	86	39-147	15	40
Trichlorofluoromethane	30.1	50.0	60		34.7	50.0	69	51-140	14	40
1,1-Dichloroethene	36.0	50.0	72		41.4	50.0	83	64-152	14	40
Acetone	176	250	70		204	250	82	43-124	15	40
Carbon Disulfide	80.7	100	81		93.4	100	93	55-141	15	40
Methylene Chloride	37.1	50.0	74		44.0	50.0	88	56-131	17	40
trans-1,2-Dichloroethene	37.4	50.0	75	*	42.8	50.0	86	77-127	14	40
1,1-Dichloroethane	38.1	50.0	76		41.9	50.0	84	59-137	9	40
2,2-Dichloropropane	40.3	50.0	81		43.9	50.0	88	74-134	9	40
cis-1,2-Dichloroethene	39.5	50.0	79		43.0	50.0	86	62-138	8	40
2-Butanone (MEK)	208	250	83		231	250	92	54-116	10	40
Bromochloromethane	40.9	50.0	82		45.1	50.0	90	59-130	10	40
Chloroform	41.5	50.0	83		45.0	50.0	90	61-137	8	40
1,1,1-Trichloroethane (TCA)	40.9	50.0	82		44.7	50.0	89	59-146	9	40
Carbon Tetrachloride	40.5	50.0	81		44.2	50.0	88	60-152	9	40
1,1-Dichloropropene	39.0	50.0	78		42.3	50.0	85	52-142	8	40
Benzene	41.8	50.0	84		44.2	50.0	88	56-148	6	40
1,2-Dichloroethane (EDC)	40.0	50.0	80		44.2	50.0	88	57-139	10	40
Trichloroethene (TCE)	42.8	50.0	86		46.2	50.0	92	76-124	8	40
1,2-Dichloropropane	41.3	50.0	83		46.6	50.0	93	58-135	12	40
Dibromomethane	39.8	50.0	80		45.2	50.0	90	56-133	13	40
Bromodichloromethane	41.9	50.0	84		48.0	50.0	96	61-143	14	40
cis-1,3-Dichloropropene	37.5	50.0	75		42.8	50.0	86	58-138	13	40
4-Methyl-2-pentanone (MIBK)	216	250	86		253	250	101	44-140	16	40
Toluene	40.9	50.0	82		47.0	50.0	94	75-117	14	40
trans-1,3-Dichloropropene	52.4	50.0	105		53.6	50.0	107	63-121	2	40
1,1,2-Trichloroethane	42.9	50.0	86		44.5	50.0	89	56-133	4	40
Tetrachloroethene (PCE)	43.4	50.0	87		43.9	50.0	88	66-126	1	40
2-Hexanone	230	250	92		238	250	95	54-122	4	40
1,3-Dichloropropane	44.0	50.0	88		45.5	50.0	91	56-134	3	40
Dibromochloromethane	47.3	50.0	95		48.5	50.0	97	58-140	3	40
1,2-Dibromoethane (EDB)	44.2	50.0	88		45.4	50.0	91	55-134	3	40

Results flagged with an asterisk (*) indicate values outside control criteria.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Extracted: 05/25/2010
Date Analyzed: 05/25/2010

Lab Control Spike/Duplicate Lab Control Spike Summary
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry
Level: Low
Extraction Lot: KWG1004900

	Lab Control Sample KWG1004900-3 Lab Control Spike			Duplicate Lab Control Sample KWG1004900-4 Duplicate Lab Control Spike					
Analyte Name	**Result**	**Expected**	**%Rec**	**Result**	**Expected**	**%Rec**	**%Rec Limits**	**RPD**	**RPD Limit**
Chlorobenzene	43.4	50.0	87	44.5	50.0	89	70-116	3	40
Ethylbenzene	44.9	50.0	90	45.8	50.0	92	70-118	2	40
1,1,1,2-Tetrachloroethane	47.5	50.0	95	48.3	50.0	97	58-137	2	40
m,p-Xylenes	94.7	100	95	96.1	100	96	58-140	2	40
o-Xylene	46.5	50.0	93	47.7	50.0	95	57-139	3	40
Styrene	47.8	50.0	96	48.5	50.0	97	62-135	1	40
Bromoform	52.5	50.0	105	53.6	50.0	107	62-134	2	40
Isopropylbenzene	48.7	50.0	97	48.6	50.0	97	55-129	0	40
1,1,2,2-Tetrachloroethane	39.3	50.0	79	44.1	50.0	88	53-133	11	40
Bromobenzene	47.9	50.0	96	46.8	50.0	94	52-130	2	40
n-Propylbenzene	49.0	50.0	98	48.6	50.0	97	57-143	1	40
1,2,3-Trichloropropane	45.2	50.0	90	44.8	50.0	90	53-134	1	40
2-Chlorotoluene	49.9	50.0	100	49.0	50.0	98	66-129	2	40
1,3,5-Trimethylbenzene	49.5	50.0	99	49.3	50.0	99	53-141	0	40
4-Chlorotoluene	50.5	50.0	101	50.0	50.0	100	51-134	1	40
tert-Butylbenzene	47.9	50.0	96	48.6	50.0	97	67-131	2	40
1,2,4-Trimethylbenzene	47.4	50.0	95	48.5	50.0	97	55-140	2	40
sec-Butylbenzene	46.4	50.0	93	46.8	50.0	94	55-146	1	40
4-Isopropyltoluene	46.7	50.0	93	47.6	50.0	95	52-140	2	40
1,3-Dichlorobenzene	47.2	50.0	94	47.9	50.0	96	52-131	1	40
1,4-Dichlorobenzene	46.2	50.0	92	46.8	50.0	94	52-128	1	40
n-Butylbenzene	47.9	50.0	96	48.8	50.0	98	53-139	2	40
1,2-Dichlorobenzene	46.3	50.0	93	46.7	50.0	93	53-128	1	40
1,2-Dibromo-3-chloropropane	50.4	50.0	101	52.1	50.0	104	55-127	3	40
1,2,4-Trichlorobenzene	50.2	50.0	100	50.5	50.0	101	57-136	1	40
Hexachlorobutadiene	54.3	50.0	109	54.2	50.0	108	54-140	0	40
Naphthalene	50.4	50.0	101	50.9	50.0	102	54-134	1	40
1,2,3-Trichlorobenzene	52.6	50.0	105	52.5	50.0	105	52-138	0	40

Results flagged with an asterisk (*) indicate values outside control criteria.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	05/25/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	05/25/2010
		Time Analyzed:	12:39

Method Blank Summary
Volatile Organic Compounds

Sample Name:	Method Blank	**File ID:**	J:\MS24\DATA\052510\0525F009.D
Lab Code:	KWG1004900-5	**Instrument ID:**	MS24
Extraction Method:	EPA 5030A	**Level:**	Low
Analysis Method:	8260B	**Extraction Lot:**	KWG1004900

This Method Blank applies to the following analyses:

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed
Lab Control Sample	KWG1004900-3	J:\MS24\DATA\052510\0525F005.D	05/25/10	11:14
Duplicate Lab Control Sample	KWG1004900-4	J:\MS24\DATA\052510\0525F006.D	05/25/10	11:35
D-4-16	K1005244-003	J:\MS24\DATA\052510\0525F012.D	05/25/10	13:42
D-4-16MS	KWG1004900-1	J:\MS24\DATA\052510\0525F013.D	05/25/10	14:40
D-4-16DMS	KWG1004900-2	J:\MS24\DATA\052510\0525F014.D	05/25/10	15:01

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	05/25/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	05/25/2010
		Time Analyzed:	11:14

Lab Control Sample Summary
Volatile Organic Compounds

Sample Name:	Lab Control Sample	**File ID:**	J:\MS24\DATA\052510\0525F005.D
Lab Code:	KWG1004900-3	**Instrument ID:**	MS24
Extraction Method:	EPA 5030A	**Level:**	Low
Analysis Method:	8260B	**Extraction Lot:**	KWG1004900

This Lab Control Sample applies to the following analyses:

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed
Method Blank	KWG1004900-5	J:\MS24\DATA\052510\0525F009.D	05/25/10	12:39
D-4-16	K1005244-003	J:\MS24\DATA\052510\0525F012.D	05/25/10	13:42
D-4-16MS	KWG1004900-1	J:\MS24\DATA\052510\0525F013.D	05/25/10	14:40
D-4-16DMS	KWG1004900-2	J:\MS24\DATA\052510\0525F014.D	05/25/10	15:01

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 05/25/2010
Time Analyzed: 08:42

Tune Summary
Volatile Organic Compounds

File ID: J:\MS24\DATA\052510\0525F002.D
Instrument ID: MS24
Column:

Analysis Method: 8260B
Analysis Lot: KWG1004880

Target Mass	Relative to Mass	Lower Limit%	Upper Limit%	Relative Abundance %	Raw Abundance	Result Pass/Fail
50	95	15	40	24.2	3284	PASS
75	95	30	60	55.9	7570	PASS
95	95	100	100	100.0	13552	PASS
96	95	5	9	8.7	1182	PASS
173	174	0	2	1.2	123	PASS
174	95	50	120	77.4	10491	PASS
175	174	5	9	8.4	877	PASS
176	174	95	101	97.2	10193	PASS
177	176	5	9	6.1	618	PASS

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed	Q
Continuing Calibration Verification	KWG1004880-2	J:\MS24\DATA\052510\0525F004.D	05/25/2010	10:34	
Lab Control Sample	KWG1004900-3	J:\MS24\DATA\052510\0525F005.D	05/25/2010	11:14	
Duplicate Lab Control Sample	KWG1004900-4	J:\MS24\DATA\052510\0525F006.D	05/25/2010	11:35	
Method Blank	KWG1004900-5	J:\MS24\DATA\052510\0525F009.D	05/25/2010	12:39	
D-4-16	K1005244-003	J:\MS24\DATA\052510\0525F012.D	05/25/2010	13:42	
D-4-16MS	KWG1004900-1	J:\MS24\DATA\052510\0525F013.D	05/25/2010	14:40	
D-4-16DMS	KWG1004900-2	J:\MS24\DATA\052510\0525F014.D	05/25/2010	15:01	

Results flagged with an asterisk (*) indicate the analysis performed outside specified tune window

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Level ID	File ID	Level ID	File ID
A	J:\MS24\DATA\042110\0421F005.D	F	J:\MS24\DATA\042110\0421F010.D
B	J:\MS24\DATA\042110\0421F006.D	G	J:\MS24\DATA\042110\0421F011.D
C	J:\MS24\DATA\042110\0421F007.D	H	J:\MS24\DATA\042110\0421F012.D
D	J:\MS24\DATA\042110\0421F008.D	I	J:\MS24\DATA\042110\0421F013.D
E	J:\MS24\DATA\042110\0421F009.D		

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
Dichlorodifluoromethane				B	2.0	0.330	C	5.0	0.312	D	10	0.299	E	20	0.321
	F	50	0.282	G	100	0.308	H	200	0.331	I	300	0.286			
† Chloromethane	A	1.0	0.506	B	2.0	0.397	C	5.0	0.378	D	10	0.369	E	20	0.379
†	F	50	0.350	G	100	0.385	H	200	0.404	I	300	0.354			
‡ Vinyl Chloride				B	2.0	0.364	C	5.0	0.349	D	10	0.351	E	20	0.360
‡	F	50	0.319	G	100	0.350	H	200	0.366	I	300	0.318			
Bromomethane							C	5.0	0.324	D	10	0.293	E	20	0.297
	F	50	0.233	G	100	0.281	H	200	0.304						
Chloroethane							C	5.0	0.228	D	10	0.221	E	20	0.239
	F	50	0.209	G	100	0.231	H	200	0.242	I	300	0.214			
Trichlorofluoromethane				B	2.0	0.487	C	5.0	0.472	D	10	0.476	E	20	0.506
	F	50	0.431	G	100	0.481	H	200	0.518	I	300	0.458			
‡ 1,1-Dichloroethene				B	2.0	0.244	C	5.0	0.199	D	10	0.203	E	20	0.219
‡	F	50	0.191	G	100	0.205	H	200	0.225	I	300	0.193			
Acetone										D	20	0.150	E	40	0.118
	F	100	0.108	G	200	0.107	H	400	0.115	I	600	0.101			
Carbon Disulfide	A	1.0	0.905	B	2.0	0.781	C	5.0	0.761	D	10	0.786	E	20	0.855
	F	50	0.742	G	100	0.847	H	200	0.923	I	300	0.826			
Methylene Chloride							C	5.0	0.322	D	10	0.295	E	20	0.319
	F	50	0.283	G	100	0.299	H	200	0.310	I	300	0.279			
trans-1,2-Dichloroethene	A	1.0	0.310	B	2.0	0.263	C	5.0	0.250	D	10	0.245	E	20	0.261
	F	50	0.236	G	100	0.261	H	200	0.282	I	300	0.254			
† 1,1-Dichloroethane	A	1.0	0.555	B	2.0	0.546	C	5.0	0.477	D	10	0.496	E	20	0.523
†	F	50	0.475	G	100	0.529	H	200	0.566	I	300	0.496			
2,2-Dichloropropane							C	5.0	0.339	D	10	0.338	E	20	0.355
	F	50	0.333	G	100	0.370	H	200	0.404	I	300	0.347			
cis-1,2-Dichloroethene	A	1.0	0.307	B	2.0	0.290	C	5.0	0.271	D	10	0.275	E	20	0.297
	F	50	0.285	G	100	0.301	H	200	0.318	I	300	0.278			
2-Butanone (MEK)							C	10	0.0436	D	20	0.0374	E	40	0.0364
	F	100	0.0352	G	200	0.0360	H	400	0.0396	I	600	0.0339			

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
Bromochloromethane							C	5.0	0.135	D	10	0.136	E	20	0.144
	F	50	0.138	G	100	0.149	H	200	0.158	I	300	0.140			
‡ Chloroform				B	2.0	0.483	C	5.0	0.452	D	10	0.465	E	20	0.498
‡	F	50	0.458	G	100	0.512	H	200	0.550	I	300	0.493			
1,1,1-Trichloroethane (TCA)				B	2.0	0.361	C	5.0	0.360	D	10	0.359	E	20	0.390
	F	50	0.343	G	100	0.387	H	200	0.414	I	300	0.373			
Carbon Tetrachloride							C	5.0	0.302	D	10	0.308	E	20	0.337
	F	50	0.287	G	100	0.327	H	200	0.362	I	300	0.324			
1,1-Dichloropropene	A	1.0	0.359	B	2.0	0.334	C	5.0	0.302	D	10	0.327	E	20	0.354
	F	50	0.299	G	100	0.339	H	200	0.371	I	300	0.334			
Benzene	A	1.0	1.10	B	2.0	1.02	C	5.0	0.979	D	10	1.01	E	20	1.13
	F	50	0.998	G	100	1.09	H	200	1.19	I	300	1.07			
1,2-Dichloroethane (EDC)	A	1.0	0.435	B	2.0	0.416	C	5.0	0.412	D	10	0.415	E	20	0.451
	F	50	0.415	G	100	0.442	H	200	0.472	I	300	0.431			
Trichloroethene (TCE)	A	1.0	0.304	B	2.0	0.298	C	5.0	0.265	D	10	0.257	E	20	0.300
	F	50	0.260	G	100	0.275	H	200	0.297	I	300	0.277			
‡ 1,2-Dichloropropane	A	1.0	0.315	B	2.0	0.283	C	5.0	0.254	D	10	0.252	E	20	0.295
‡	F	50	0.270	G	100	0.294	H	200	0.304	I	300	0.287			
Dibromomethane							C	5.0	0.173	D	10	0.165	E	20	0.190
	F	50	0.174	G	100	0.186	H	200	0.189	I	300	0.179			
Bromodichloromethane							C	5.0	0.312	D	10	0.310	E	20	0.368
	F	50	0.345	G	100	0.384	H	200	0.400	I	300	0.385			
cis-1,3-Dichloropropene							C	5.0	0.325	D	10	0.328	E	20	0.408
	F	50	0.384	G	100	0.445	H	200	0.475	I	300	0.431			
4-Methyl-2-pentanone (MIBK)							C	10	0.103	D	20	0.0913	E	40	0.108
	F	100	0.108	G	200	0.118	H	400	0.123	I	600	0.110			
‡ Toluene				B	2.0	0.664	C	5.0	0.579	D	10	0.599	E	20	0.693
‡	F	50	0.601	G	100	0.665	H	200	0.694	I	300	0.624			
trans-1,3-Dichloropropene							C	5.0	0.661	D	10	0.684	E	20	0.877
	F	50	0.809	G	100	0.868	H	200	0.964	I	300	0.932			
1,1,2-Trichloroethane	A	1.0	0.536	B	2.0	0.603	C	5.0	0.510	D	10	0.473	E	20	0.558
	F	50	0.482	G	100	0.491	H	200	0.522	I	300	0.485			
Tetrachloroethene (PCE)	A	1.0	0.506	B	2.0	0.541	C	5.0	0.473	D	10	0.481	E	20	0.537
	F	50	0.448	G	100	0.459	H	200	0.506	I	300	0.463			
2-Hexanone				B	4.0	0.0963	C	10	0.0808	D	20	0.0774	E	40	0.0846
	F	100	0.0845	G	200	0.0884	H	400	0.0973	I	600	0.0901			

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound** ‡ **CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
1,3-Dichloropropane	A	1.0	0.986	B	2.0	1.19	C	5.0	0.953	D	10	0.965	E	20	1.13
	F	50	0.984	G	100	1.02	H	200	1.06	I	300	0.993			
Dibromochloromethane							C	5.0	0.499	D	10	0.485	E	20	0.579
	F	50	0.554	G	100	0.620	H	200	0.674	I	300	0.644			
1,2-Dibromoethane (EDB)				B	2.0	0.622	C	5.0	0.506	D	10	0.527	E	20	0.591
	F	50	0.544	G	100	0.562	H	200	0.613	I	300	0.561			
† Chlorobenzene	A	1.0	1.72	B	2.0	1.94	C	5.0	1.61	D	10	1.62	E	20	1.81
†	F	50	1.63	G	100	1.73	H	200	1.82	I	300	1.65			
‡ Ethylbenzene				B	2.0	0.821	C	5.0	0.768	D	10	0.802	E	20	0.903
‡	F	50	0.795	G	100	0.845	H	200	0.936	I	300	0.855			
1,1,1,2-Tetrachloroethane							C	5.0	0.508	D	10	0.502	E	20	0.583
	F	50	0.545	G	100	0.582	H	200	0.663	I	300	0.603			
m,p-Xylenes	A	2.0	0.859	B	4.0	0.949	C	10	0.876	D	20	0.925	E	40	1.07
	F	100	0.985	G	200	1.06	H	400	1.15	I	600	1.04			
o-Xylene							C	5.0	0.883	D	10	0.912	E	20	1.03
	F	50	0.948	G	100	1.06	H	200	1.10	I	300	1.02			
Styrene							C	5.0	0.662	D	10	0.652	E	20	0.804
	F	50	0.776	G	100	0.842	H	200	0.887	I	300	0.837			
† Bromoform							C	5.0	0.286	D	10	0.283	E	20	0.339
†	F	50	0.354	G	100	0.381	H	200	0.417	I	300	0.403			
Isopropylbenzene							C	5.0	2.05	D	10	2.22	E	20	2.52
	F	50	2.46	G	100	2.48	H	200	2.68	I	300	2.40			
† 1,1,2,2-Tetrachloroethane				B	2.0	0.939	C	5.0	0.706	D	10	0.625	E	20	0.735
†	F	50	0.684	G	100	0.692	H	200	0.754	I	300	0.688			
Bromobenzene	A	1.0	0.742	B	2.0	0.959	C	5.0	0.715	D	10	0.664	E	20	0.825
	F	50	0.763	G	100	0.780	H	200	0.839	I	300	0.779			
n-Propylbenzene							C	5.0	2.69	D	10	2.78	E	20	3.44
	F	50	3.22	G	100	3.31	H	200	3.67	I	300	3.32			
1,2,3-Trichloropropane	A	1.0	0.278	B	2.0	0.354	C	5.0	0.305	D	10	0.249	E	20	0.277
	F	50	0.257	G	100	0.250	H	200	0.271	I	300	0.249			
2-Chlorotoluene				B	2.0	2.18	C	5.0	1.78	D	10	1.77	E	20	2.16
	F	50	1.98	G	100	2.01	H	200	2.18	I	300	2.01			
1,3,5-Trimethylbenzene							C	5.0	1.94	D	10	1.89	E	20	2.42
	F	50	2.27	G	100	2.34	H	200	2.55	I	300	2.34			
4-Chlorotoluene				B	2.0	2.23	C	5.0	1.81	D	10	1.76	E	20	2.13
	F	50	2.02	G	100	2.08	H	200	2.21	I	300	1.99			

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound**

‡ **CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
tert-Butylbenzene							C	5.0	1.75	D	10	1.68	E	20	2.09
	F	50	1.96	G	100	2.06	H	200	2.20	I	300	2.06			
1,2,4-Trimethylbenzene							C	5.0	1.95	D	10	1.86	E	20	2.47
	F	50	2.31	G	100	2.45	H	200	2.60	I	300	2.47			
sec-Butylbenzene							C	5.0	2.47	D	10	2.58	E	20	3.24
	F	50	2.87	G	100	3.08	H	200	3.26	I	300	3.06			
4-Isopropyltoluene							C	5.0	2.07	D	10	2.16	E	20	2.69
	F	50	2.40	G	100	2.52	H	200	2.74	I	300	2.57			
1,3-Dichlorobenzene	A	1.0	1.28	B	2.0	1.69	C	5.0	1.38	D	10	1.31	E	20	1.55
	F	50	1.41	G	100	1.43	H	200	1.53	I	300	1.45			
1,4-Dichlorobenzene	A	1.0	1.49	B	2.0	1.70	C	5.0	1.45	D	10	1.37	E	20	1.62
	F	50	1.41	G	100	1.47	H	200	1.53	I	300	1.45			
n-Butylbenzene							C	5.0	1.79	D	10	1.92	E	20	2.36
	F	50	2.09	G	100	2.27	H	200	2.43	I	300	2.29			
1,2-Dichlorobenzene	A	1.0	1.30	B	2.0	1.42	C	5.0	1.38	D	10	1.30	E	20	1.64
	F	50	1.39	G	100	1.44	H	200	1.51	I	300	1.42			
1,2-Dibromo-3-chloropropane				B	2.0	0.0926	C	5.0	0.115	D	10	0.0877	E	20	0.119
	F	50	0.0940	G	100	0.0982	H	200	0.131	I	300	0.133			
1,2,4-Trichlorobenzene							C	5.0	0.911	D	10	0.837	E	20	1.01
	F	50	0.687	G	100	0.884	H	200	1.07						
Hexachlorobutadiene	A	1.0	0.471	B	2.0	0.357	C	5.0	0.502	D	10	0.497	E	20	0.583
	F	50	0.356	G	100	0.380	H	200	0.592	I	300	0.319			
Naphthalene							C	5.0	2.01	D	10	1.89	E	20	2.31
	F	50	1.73	G	100	2.16	H	200	2.69	I	300	1.70			
1,2,3-Trichlorobenzene							C	5.0	0.732	D	10	0.848	E	20	1.05
	F	50	0.688	G	100	0.645	H	200	1.07						
Dibromofluoromethane				B	10	0.230	C	20	0.256	D	30	0.248	E	40	0.256
	F	50	0.240	G	70	0.257	H	80	0.265	I	100	0.248			
Toluene-d8				B	10	0.728	C	20	0.885	D	30	0.845	E	40	0.927
	F	50	0.844	G	70	0.939	H	80	0.958	I	100	0.887			
4-Bromofluorobenzene				B	10	0.690	C	20	0.767	D	30	0.720	E	40	0.795
	F	50	0.808	G	70	0.764	H	80	0.813	I	100	0.782			

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound**

‡ **CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Compound Type	Calibration Evaluation					RRF Evaluation		
		Fit Type	Eval.	Eval. Result	Q	Control Criteria	Average RRF	Q	Minimum RRF
Dichlorodifluoromethane	TRG	AverageRF	% RSD	6.1		≤ 15	0.309		0.01
† Chloromethane	TRG	AverageRF	% RSD	11.9		≤ 15	0.391		0.10
‡ Vinyl Chloride	MS	AverageRF	% RSD	5.4		≤ 15	0.347		0.01
Bromomethane	TRG	AverageRF	% RSD	10.6		≤ 15	0.289		0.01
Chloroethane	TRG	AverageRF	% RSD	5.5		≤ 15	0.226		0.01
Trichlorofluoromethane	TRG	AverageRF	% RSD	5.7		≤ 15	0.479		0.01
‡ 1,1-Dichloroethene	MS	AverageRF	% RSD	8.7		≤ 15	0.210		0.01
Acetone	TRG	AverageRF	% RSD	15.1	*	≤ 15	0.116		0.01
Carbon Disulfide	TRG	AverageRF	% RSD	7.6		≤ 15	0.825		0.01
Methylene Chloride	TRG	AverageRF	% RSD	5.7		≤ 15	0.301		0.01
trans-1,2-Dichloroethene	TRG	AverageRF	% RSD	8.4		≤ 15	0.263		0.01
† 1,1-Dichloroethane	TRG	AverageRF	% RSD	6.5		≤ 15	0.518		0.10
2,2-Dichloropropane	TRG	AverageRF	% RSD	7.0		≤ 15	0.355		0.01
cis-1,2-Dichloroethene	TRG	AverageRF	% RSD	5.4		≤ 15	0.291		0.01
2-Butanone (MEK)	TRG	AverageRF	% RSD	8.7		≤ 15	0.0374		0.01
Bromochloromethane	TRG	AverageRF	% RSD	5.9		≤ 15	0.143		0.01
‡ Chloroform	MS	AverageRF	% RSD	6.6		≤ 15	0.489		0.01
1,1,1-Trichloroethane (TCA)	TRG	AverageRF	% RSD	6.0		≤ 15	0.373		0.01
Carbon Tetrachloride	MS	AverageRF	% RSD	7.7		≤ 15	0.321		0.01
1,1-Dichloropropene	TRG	AverageRF	% RSD	7.2		≤ 15	0.335		0.01
Benzene	MS	AverageRF	% RSD	6.5		≤ 15	1.07		0.01
1,2-Dichloroethane (EDC)	TRG	AverageRF	% RSD	4.7		≤ 15	0.432		0.01
Trichloroethene (TCE)	MS	AverageRF	% RSD	6.6		≤ 15	0.282		0.01
‡ 1,2-Dichloropropane	TRG	AverageRF	% RSD	7.6		≤ 15	0.284		0.01
Dibromomethane	TRG	AverageRF	% RSD	5.2		≤ 15	0.179		0.01
Bromodichloromethane	MS	AverageRF	% RSD	10.1		≤ 15	0.358		0.01
cis-1,3-Dichloropropene	TRG	AverageRF	% RSD	14.4		≤ 15	0.399		0.01
4-Methyl-2-pentanone (MIBK)	TRG	AverageRF	% RSD	9.4		≤ 15	0.109		0.01
‡ Toluene	MS	AverageRF	% RSD	7.0		≤ 15	0.640		0.01
trans-1,3-Dichloropropene	TRG	AverageRF	% RSD	14.2		≤ 15	0.828		0.01
1,1,2-Trichloroethane	MS	AverageRF	% RSD	8.2		≤ 15	0.518		0.01
Tetrachloroethene (PCE)	TRG	AverageRF	% RSD	6.9		≤ 15	0.491		0.01
2-Hexanone	MS	AverageRF	% RSD	8.0		≤ 15	0.0874		0.01
1,3-Dichloropropane	TRG	AverageRF	% RSD	7.9		≤ 15	1.03		0.01
Dibromochloromethane	TRG	AverageRF	% RSD	12.4		≤ 15	0.579		0.01
1,2-Dibromoethane (EDB)	TRG	AverageRF	% RSD	7.2		≤ 15	0.566		0.01
† Chlorobenzene	MS	AverageRF	% RSD	6.6		≤ 15	1.73		0.30
‡ Ethylbenzene	MS	AverageRF	% RSD	6.8		≤ 15	0.841		0.01
1,1,1,2-Tetrachloroethane	TRG	AverageRF	% RSD	9.9		≤ 15	0.570		0.01
m,p-Xylenes	TRG	AverageRF	% RSD	9.8		≤ 15	0.991		0.01
o-Xylene	TRG	AverageRF	% RSD	8.1		≤ 15	0.995		0.01
Styrene	TRG	AverageRF	% RSD	11.6		≤ 15	0.780		0.01

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound** ‡ **CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Compound Type	Calibration Evaluation					RRF Evaluation		
		Fit Type	Eval.	Eval. Result	Q	Control Criteria	Average RRF	Q	Minimum RRF
† Bromoform	TRG	AverageRF	% RSD	15.1	*	≤ 15	0.352		0.10
Isopropylbenzene	TRG	AverageRF	% RSD	8.6		≤ 15	2.40		0.01
† 1,1,2,2-Tetrachloroethane	TRG	AverageRF	% RSD	12.9		≤ 15	0.728		0.30
Bromobenzene	TRG	AverageRF	% RSD	10.7		≤ 15	0.785		0.01
n-Propylbenzene	TRG	AverageRF	% RSD	10.9		≤ 15	3.20		0.01
1,2,3-Trichloropropane	MS	AverageRF	% RSD	12.4		≤ 15	0.277		0.01
2-Chlorotoluene	MS	AverageRF	% RSD	8.3		≤ 15	2.01		0.01
1,3,5-Trimethylbenzene	TRG	AverageRF	% RSD	10.8		≤ 15	2.25		0.01
4-Chlorotoluene	TRG	AverageRF	% RSD	8.5		≤ 15	2.03		0.01
tert-Butylbenzene	TRG	AverageRF	% RSD	9.6		≤ 15	1.97		0.01
1,2,4-Trimethylbenzene	TRG	AverageRF	% RSD	12.4		≤ 15	2.30		0.01
sec-Butylbenzene	TRG	AverageRF	% RSD	10.6		≤ 15	2.94		0.01
4-Isopropyltoluene	TRG	AverageRF	% RSD	10.4		≤ 15	2.45		0.01
1,3-Dichlorobenzene	TRG	AverageRF	% RSD	8.8		≤ 15	1.45		0.01
1,4-Dichlorobenzene	TRG	AverageRF	% RSD	6.9		≤ 15	1.50		0.01
n-Butylbenzene	TRG	AverageRF	% RSD	11.1		≤ 15	2.17		0.01
1,2-Dichlorobenzene	MS	AverageRF	% RSD	7.3		≤ 15	1.42		0.01
1,2-Dibromo-3-chloropropane	TRG	AverageRF	% RSD	16.5	*	≤ 15	0.109		0.01
1,2,4-Trichlorobenzene	TRG	AverageRF	% RSD	15.1	*	≤ 15	0.901		0.01
Hexachlorobutadiene	TRG	AverageRF	% RSD	22.5	*	≤ 15	0.451		0.01
Naphthalene	MS	AverageRF	% RSD	17.0	*	≤ 15	2.07		0.01
1,2,3-Trichlorobenzene	TRG	AverageRF	% RSD	22.0	*	≤ 15	0.839		0.01
Dibromofluoromethane	SURR	AverageRF	% RSD	4.4		≤ 15	0.250		0.01
Toluene-d8	SURR	AverageRF	% RSD	8.3		≤ 15	0.877		0.01
4-Bromofluorobenzene	SURR	AverageRF	% RSD	5.6		≤ 15	0.767		0.01

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010
Date Analyzed: 04/21/2010

Second Source Calibration Verification
Volatile Organic Compounds

Calibration Type: Internal Standard
Analysis Method: 8260B

Calibration ID: CAL9404
Units: PPB

File ID: J:\MS24\DATA\042110\0421F017.D

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Dichlorodifluoromethane	50	48	0.309	0.298	-4	NA	± 40 %	AverageRF
† Chloromethane	50	43	0.391	0.336	-14	NA	± 40 %	AverageRF
‡ Vinyl Chloride	50	46	0.347	0.319	-8	NA	± 20 %	AverageRF
Bromomethane	50	45	0.289	0.257	-11	NA	± 40 %	AverageRF
Chloroethane	50	41	0.226	0.188	-17	NA	± 40 %	AverageRF
Trichlorofluoromethane	50	42	0.479	0.404	-16	NA	± 30 %	AverageRF
‡ 1,1-Dichloroethene	50	51	0.210	0.214	2	NA	± 20 %	AverageRF
Acetone	250	210	0.116	0.0983	-16	NA	± 30 %	AverageRF
Carbon Disulfide	100	96	0.825	0.791	-4	NA	± 30 %	AverageRF
Methylene Chloride	50	46	0.301	0.275	-9	NA	± 30 %	AverageRF
trans-1,2-Dichloroethene	50	49	0.263	0.256	-3	NA	± 30 %	AverageRF
† 1,1-Dichloroethane	50	47	0.518	0.485	-6	NA	± 30 %	AverageRF
2,2-Dichloropropane	50	51	0.355	0.359	1	NA	± 30 %	AverageRF
cis-1,2-Dichloroethene	50	46	0.291	0.268	-8	NA	± 30 %	AverageRF
2-Butanone (MEK)	250	230	0.0374	0.0348	-7	NA	± 30 %	AverageRF
Bromochloromethane	50	48	0.143	0.137	-4	NA	± 30 %	AverageRF
‡ Chloroform	50	48	0.489	0.472	-3	NA	± 20 %	AverageRF
1,1,1-Trichloroethane (TCA)	50	51	0.373	0.380	2	NA	± 30 %	AverageRF
Carbon Tetrachloride	50	51	0.321	0.325	1	NA	± 30 %	AverageRF
1,1-Dichloropropene	50	48	0.335	0.321	-4	NA	± 30 %	AverageRF
Benzene	50	48	1.07	1.02	-4	NA	± 30 %	AverageRF
1,2-Dichloroethane (EDC)	50	47	0.432	0.407	-6	NA	± 30 %	AverageRF
Trichloroethene (TCE)	50	51	0.282	0.287	2	NA	± 30 %	AverageRF
‡ 1,2-Dichloropropane	50	50	0.284	0.282	-1	NA	± 20 %	AverageRF
Dibromomethane	50	50	0.179	0.178	0	NA	± 30 %	AverageRF
Bromodichloromethane	50	51	0.358	0.364	2	NA	± 30 %	AverageRF
cis-1,3-Dichloropropene	50	45	0.399	0.361	-10	NA	± 30 %	AverageRF
4-Methyl-2-pentanone (MIBK)	250	250	0.109	0.108	-1	NA	± 30 %	AverageRF
‡ Toluene	50	50	0.640	0.634	-1	NA	± 20 %	AverageRF
trans-1,3-Dichloropropene	50	58	0.828	0.953	15	NA	± 30 %	AverageRF
1,1,2-Trichloroethane	50	51	0.518	0.527	2	NA	± 30 %	AverageRF
Tetrachloroethene (PCE)	50	53	0.491	0.524	7	NA	± 30 %	AverageRF
2-Hexanone	250	260	0.0874	0.0904	3	NA	± 30 %	AverageRF
1,3-Dichloropropane	50	51	1.03	1.05	1	NA	± 30 %	AverageRF
Dibromochloromethane	50	51	0.579	0.591	2	NA	± 30 %	AverageRF
1,2-Dibromoethane (EDB)	50	52	0.566	0.591	5	NA	± 30 %	AverageRF
† Chlorobenzene	50	50	1.73	1.72	0	NA	± 30 %	AverageRF
‡ Ethylbenzene	50	53	0.841	0.897	7	NA	± 20 %	AverageRF
1,1,1,2-Tetrachloroethane	50	52	0.570	0.591	4	NA	± 30 %	AverageRF
m,p-Xylenes	100	110	0.991	1.11	12	NA	± 30 %	AverageRF

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010
Date Analyzed: 04/21/2010

Second Source Calibration Verification
Volatile Organic Compounds

Calibration Type: Internal Standard
Analysis Method: 8260B

Calibration ID: CAL9404
Units: PPB

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
o-Xylene	50	54	0.995	1.08	8	NA	± 30 %	AverageRF
Styrene	50	54	0.780	0.848	9	NA	± 30 %	AverageRF
† Bromoform	50	54	0.352	0.383	9	NA	± 30 %	AverageRF
Isopropylbenzene	50	58	2.40	2.76	15	NA	± 30 %	AverageRF
† 1,1,2,2-Tetrachloroethane	50	47	0.728	0.686	-6	NA	± 30 %	AverageRF
Bromobenzene	50	47	0.785	0.734	-7	NA	± 30 %	AverageRF
n-Propylbenzene	50	52	3.20	3.32	4	NA	± 30 %	AverageRF
1,2,3-Trichloropropane	50	46	0.277	0.256	-7	NA	± 30 %	AverageRF
2-Chlorotoluene	50	51	2.01	2.04	2	NA	± 30 %	AverageRF
1,3,5-Trimethylbenzene	50	53	2.25	2.37	5	NA	± 30 %	AverageRF
4-Chlorotoluene	50	50	2.03	2.01	-1	NA	± 30 %	AverageRF
tert-Butylbenzene	50	53	1.97	2.09	6	NA	± 30 %	AverageRF
1,2,4-Trimethylbenzene	50	53	2.30	2.44	6	NA	± 30 %	AverageRF
sec-Butylbenzene	50	54	2.94	3.18	8	NA	± 30 %	AverageRF
4-Isopropyltoluene	50	56	2.45	2.73	12	NA	± 30 %	AverageRF
1,3-Dichlorobenzene	50	52	1.45	1.52	5	NA	± 30 %	AverageRF
1,4-Dichlorobenzene	50	51	1.50	1.54	2	NA	± 30 %	AverageRF
n-Butylbenzene	50	56	2.17	2.41	12	NA	± 30 %	AverageRF
1,2-Dichlorobenzene	50	50	1.42	1.41	-1	NA	± 30 %	AverageRF
1,2-Dibromo-3-chloropropane	50	57	0.109	0.124	14	NA	± 30 %	AverageRF
1,2,4-Trichlorobenzene	50	53	0.901	0.961	7	NA	± 30 %	AverageRF
Hexachlorobutadiene	50	59	0.451	0.536	19	NA	± 30 %	AverageRF
Naphthalene	50	54	2.07	2.22	7	NA	± 30 %	AverageRF
1,2,3-Trichlorobenzene	50	60	0.839	0.999	19	NA	± 30 %	AverageRF

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 05/25/2010

Continuing Calibration Verification Summary
Volatile Organic Compounds

Calibration Type: Internal Standard
Analysis Method: 8260B

Calibration Date: 04/21/2010
Calibration ID: CAL9404
Analysis Lot: KWG1004880
Units: PPB

File ID: J:\MS24\DATA\052510\0525F004.D

Analyte Name	Expected	Result	Min RF	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
Dichlorodifluoromethane	50	36	0.01	0.309	0.221	-29	NA	± 40 %	AverageRF
† Chloromethane	50	43	0.10	0.391	0.340	-13	NA	± 40 %	AverageRF
‡ Vinyl Chloride	50	42	0.01	0.347	0.294	-15	NA	± 20 %	AverageRF
Bromomethane	50	34	0.01	0.289	0.199	-31	NA	± 40 %	AverageRF
Chloroethane	50	46	0.01	0.226	0.206	-9	NA	± 40 %	AverageRF
Trichlorofluoromethane	50	40	0.01	0.479	0.380	-21	NA	± 30 %	AverageRF
‡ 1,1-Dichloroethene	50	45	0.01	0.210	0.188	-10	NA	± 20 %	AverageRF
Acetone	100	92	0.01	0.116	0.107	-8	NA	± 30 %	AverageRF
Carbon Disulfide	50	48	0.01	0.825	0.794	-4	NA	± 30 %	AverageRF
Methylene Chloride	50	52	0.01	0.301	0.315	5	NA	± 30 %	AverageRF
trans-1,2-Dichloroethene	50	49	0.01	0.263	0.258	-2	NA	± 30 %	AverageRF
† 1,1-Dichloroethane	50	50	0.10	0.518	0.518	0	NA	± 30 %	AverageRF
2,2-Dichloropropane	50	49	0.01	0.355	0.347	-2	NA	± 30 %	AverageRF
cis-1,2-Dichloroethene	50	49	0.01	0.291	0.286	-2	NA	± 30 %	AverageRF
2-Butanone (MEK)	100	92	0.01	0.0374	0.0343	-8	NA	± 30 %	AverageRF
Bromochloromethane	50	49	0.01	0.143	0.140	-2	NA	± 30 %	AverageRF
‡ Chloroform	50	50	0.01	0.489	0.490	0	NA	± 20 %	AverageRF
1,1,1-Trichloroethane (TCA)	50	48	0.01	0.373	0.360	-3	NA	± 30 %	AverageRF
Carbon Tetrachloride	50	48	0.01	0.321	0.308	-4	NA	± 30 %	AverageRF
1,1-Dichloropropene	50	50	0.01	0.335	0.337	0	NA	± 30 %	AverageRF
Benzene	50	51	0.01	1.07	1.08	1	NA	± 30 %	AverageRF
1,2-Dichloroethane (EDC)	50	49	0.01	0.432	0.423	-2	NA	± 30 %	AverageRF
Trichloroethene (TCE)	50	52	0.01	0.282	0.294	4	NA	± 30 %	AverageRF
‡ 1,2-Dichloropropane	50	50	0.01	0.284	0.284	0	NA	± 20 %	AverageRF
Dibromomethane	50	47	0.01	0.179	0.170	-5	NA	± 30 %	AverageRF
Bromodichloromethane	50	52	0.01	0.358	0.371	4	NA	± 30 %	AverageRF
cis-1,3-Dichloropropene	50	54	0.01	0.399	0.434	9	NA	± 30 %	AverageRF
4-Methyl-2-pentanone (MIBK)	100	89	0.01	0.109	0.0973	-11	NA	± 30 %	AverageRF
‡ Toluene	50	52	0.01	0.640	0.663	4	NA	± 20 %	AverageRF
trans-1,3-Dichloropropene	50	54	0.01	0.828	0.894	8	NA	± 30 %	AverageRF
1,1,2-Trichloroethane	50	47	0.01	0.518	0.486	-6	NA	± 30 %	AverageRF
Tetrachloroethene (PCE)	50	51	0.01	0.491	0.498	2	NA	± 30 %	AverageRF
2-Hexanone	100	93	0.01	0.0874	0.0817	-7	NA	± 30 %	AverageRF
1,3-Dichloropropane	50	48	0.01	1.03	1.00	-3	NA	± 30 %	AverageRF
Dibromochloromethane	50	52	0.01	0.579	0.605	5	NA	± 30 %	AverageRF
1,2-Dibromoethane (EDB)	50	49	0.01	0.566	0.549	-3	NA	± 30 %	AverageRF
† Chlorobenzene	50	51	0.30	1.73	1.77	2	NA	± 30 %	AverageRF
‡ Ethylbenzene	50	52	0.01	0.841	0.880	5	NA	± 20 %	AverageRF
1,1,1,2-Tetrachloroethane	50	53	0.01	0.570	0.599	5	NA	± 30 %	AverageRF

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 05/25/2010

Continuing Calibration Verification Summary
Volatile Organic Compounds

Calibration Type: Internal Standard
Analysis Method: 8260B

Calibration Date: 04/21/2010
Calibration ID: CAL9404
Analysis Lot: KWG1004880
Units: PPB

Analyte Name	Expected	Result	Min RF	Average RF	CCV RF	%D	%Drift	Criteria	Curve Fit
m,p-Xylenes	100	110	0.01	0.991	1.09	10	NA	± 30 %	AverageRF
o-Xylene	50	52	0.01	0.995	1.04	5	NA	± 30 %	AverageRF
Styrene	50	53	0.01	0.780	0.832	7	NA	± 30 %	AverageRF
† Bromoform	50	53	0.10	0.352	0.375	7	NA	± 30 %	AverageRF
Isopropylbenzene	50	53	0.01	2.40	2.54	6	NA	± 30 %	AverageRF
† 1,1,2,2-Tetrachloroethane	50	41	0.30	0.728	0.590	-19	NA	± 30 %	AverageRF
Bromobenzene	50	55	0.01	0.785	0.856	9	NA	± 30 %	AverageRF
n-Propylbenzene	50	58	0.01	3.20	3.69	15	NA	± 30 %	AverageRF
1,2,3-Trichloropropane	50	45	0.01	0.277	0.251	-9	NA	± 30 %	AverageRF
2-Chlorotoluene	50	55	0.01	2.01	2.20	9	NA	± 30 %	AverageRF
1,3,5-Trimethylbenzene	50	56	0.01	2.25	2.50	11	NA	± 30 %	AverageRF
4-Chlorotoluene	50	57	0.01	2.03	2.32	14	NA	± 30 %	AverageRF
tert-Butylbenzene	50	53	0.01	1.97	2.09	6	NA	± 30 %	AverageRF
1,2,4-Trimethylbenzene	50	55	0.01	2.30	2.53	10	NA	± 30 %	AverageRF
sec-Butylbenzene	50	54	0.01	2.94	3.18	8	NA	± 30 %	AverageRF
4-Isopropyltoluene	50	54	0.01	2.45	2.63	7	NA	± 30 %	AverageRF
1,3-Dichlorobenzene	50	53	0.01	1.45	1.54	7	NA	± 30 %	AverageRF
1,4-Dichlorobenzene	50	52	0.01	1.50	1.55	3	NA	± 30 %	AverageRF
n-Butylbenzene	50	56	0.01	2.17	2.43	12	NA	± 30 %	AverageRF
1,2-Dichlorobenzene	50	51	0.01	1.42	1.45	2	NA	± 30 %	AverageRF
1,2-Dibromo-3-chloropropane	50	47	0.01	0.109	0.103	-5	NA	± 30 %	AverageRF
1,2,4-Trichlorobenzene	50	55	0.01	0.901	0.998	11	NA	± 30 %	AverageRF
Hexachlorobutadiene	50	62	0.01	0.451	0.555	23	NA	± 30 %	AverageRF
Naphthalene	50	49	0.01	2.07	2.05	-1	NA	± 30 %	AverageRF
1,2,3-Trichlorobenzene	50	55	0.01	0.839	0.931	11	NA	± 30 %	AverageRF
Dibromofluoromethane	50	48	0.01	0.250	0.240	-4	NA	± 30 %	AverageRF
Toluene-d8	50	55	0.01	0.877	0.961	10	NA	± 30 %	AverageRF
4-Bromofluorobenzene	50	56	0.01	0.767	0.854	11	NA	± 30 %	AverageRF

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound** ‡ **CCC Compound**

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244

Analysis Run Log
Volatile Organic Compounds

Analysis Method: 8260B

Analysis Lot: KWG1004880
Instrument ID: MS24

File ID	Sample Name	Lab Code	Date Analysis Started	Start Time	Q	Date Analysis Finished	Finish Time
0525F002.D	GC/MS Tuning - Bromofluorobenzene	KWG1004880-1	5/25/2010	08:42		5/25/2010	08:55
0525F004.D	Continuing Calibration Verification	KWG1004880-2	5/25/2010	10:34		5/25/2010	10:47
0525F005.D	Lab Control Sample	KWG1004900-3	5/25/2010	11:14		5/25/2010	11:27
0525F006.D	Duplicate Lab Control Sample	KWG1004900-4	5/25/2010	11:35		5/25/2010	11:48
0525F009.D	Method Blank	KWG1004900-5	5/25/2010	12:39		5/25/2010	12:52
0525F010.D	ZZZZZZ	ZZZZZZ	5/25/2010	13:00		5/25/2010	13:13
0525F011.D	ZZZZZZ	ZZZZZZ	5/25/2010	13:21		5/25/2010	13:34
0525F012.D	D-4-16	K1005244-003	5/25/2010	13:42		5/25/2010	13:55
0525F013.D	D-4-16MS	KWG1004900-1	5/25/2010	14:40		5/25/2010	14:53
0525F014.D	D-4-16DMS	KWG1004900-2	5/25/2010	15:01		5/25/2010	15:14
0525F016.D	ZZZZZZ	ZZZZZZ	5/25/2010	15:43		5/25/2010	15:56
0525F017.D	ZZZZZZ	ZZZZZZ	5/25/2010	16:04		5/25/2010	16:17

Results flagged with an asterisk (*) indicate the holding time was exceeded for the analysis

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Extracted: 05/25/2010

Extraction Prep Log
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Extraction Lot: KWG1004900
Level: Low

Sample Name	Lab Code	Date Collected	Date Received	Sample Amount	Final Volume	% Solids	Note
D-4-16	K1005244-003	05/19/10	05/21/10	5.00g	5.0ml	90.7	
Method Blank	KWG1004900-5	NA	NA	5.00g	5.0ml	NA	
D-4-16MS	KWG1004900-1	05/19/10	05/21/10	5.03g	5.0ml	90.7	
D-4-16DMS	KWG1004900-2	05/19/10	05/21/10	5.00g	5.0ml	90.7	
Lab Control Sample	KWG1004900-3	NA	NA	5.00g	5.0ml	NA	
Duplicate Lab Control Sample	KWG1004900-4	NA	NA	5.00g	5.0ml	NA	

Results flagged with an asterisk (*) indicate the holding time was exceeded for the analysis

Organic Analysis:
Volatile Organic Compounds

Validation Package

Organic Analysis:
Volatile Organic Compounds

Validation Package

QC Reports

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244

Surrogate Recovery Summary
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: PERCENT
Level: Low

Sample Name	Lab Code	Sur1	Sur2	Sur3
Duplicate Lab Control Sample	KWG1004900-4	100	115	109
D-4-16	K1005244-003	100	108	103
Method Blank	KWG1004900-5	100	110	104
D-4-16MS	KWG1004900-1	99	109	108
D-4-16DMS	KWG1004900-2	102	114	107
Lab Control Sample	KWG1004900-3	94	105	110

Surrogate Recovery Control Limits (%)

Sur1 = Dibromofluoromethane	83-128
Sur2 = Toluene-d8	90-125
Sur3 = 4-Bromofluorobenzene	77-124

Results flagged with an asterisk (*) indicate values outside control criteria.

Results flagged with a pound (#) indicate the control criteria is not applicable.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 05/25/2010
Time Analyzed: 10:34

Internal Standard Area and RT Summary
Volatile Organic Compounds

File ID: J:\MS24\DATA\052510\0525F004.D
Instrument ID: MS24
Analysis Method: 8260B

Lab Code: KWG1004880-2
Analysis Lot: KWG1004880

		Fluorobenzene		Chlorobenzene-d5		1,4-Dichlorobenzene-d4	
		Area	**RT**	**Area**	**RT**	**Area**	**RT**
	Results ==>	367,346	4.02	157,615	6.65	137,921	8.98
	Upper Limit ==>	734,692	4.52	315,230	7.15	275,842	9.48
	Lower Limit ==>	183,673	3.52	78,808	6.15	68,961	8.48
	ICAL Result ==>	321,724	4.02	137,440	6.65	130,674	8.98
Associated Analyses							
Lab Control Sample	KWG1004900-3	418,826	4.02	167,585	6.65	148,678	8.98
Duplicate Lab Control Sample	KWG1004900-4	355,415	4.02	159,438	6.65	144,430	8.98
Method Blank	KWG1004900-5	348,737	4.02	149,494	6.65	128,420	8.98
D-4-16	K1005244-003	342,019	4.02	146,554	6.65	116,925	8.98
D-4-16MS	KWG1004900-1	420,529	4.02	177,982	6.65	145,052	8.98
D-4-16DMS	KWG1004900-2	367,092	4.02	163,841	6.65	139,592	8.98

Results flagged with an asterisk (*) indicate values outside control criteria.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	05/25/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	05/25/2010

Matrix Spike/Duplicate Matrix Spike Summary
Volatile Organic Compounds

Sample Name:	D-4-16	**Units:**	ug/Kg
Lab Code:	K1005244-003	**Basis:**	Dry
Extraction Method:	EPA 5030A	**Level:**	Low
Analysis Method:	8260B	**Extraction Lot:**	KWG1004900

Analyte Name	Sample Result	D-4-16MS KWG1004900-1 Matrix Spike			D-4-16DMS KWG1004900-2 Duplicate Matrix Spike			%Rec Limits	RPD	RPD Limit
		Result	Expected	%Rec	Result	Expected	%Rec			
Vinyl Chloride	ND	49.8	54.8	91	55.8	55.1	101	12-160	11	40
1,1-Dichloroethene	ND	44.0	54.8	80	49.5	55.1	90	31-153	12	40
Chloroform	ND	46.5	54.8	85	49.8	55.1	90	30-138	7	40
Carbon Tetrachloride	ND	47.8	54.8	87	51.3	55.1	93	10-144	7	40
Benzene	ND	46.7	54.8	85	49.3	55.1	89	30-137	5	40
Trichloroethene (TCE)	ND	51.6	54.8	94	52.9	55.1	96	18-145	3	40
Bromodichloromethane	ND	46.1	54.8	84	52.1	55.1	95	14-146	12	40
Toluene	ND	45.9	54.8	84	50.7	55.1	92	15-135	10	40
1,1,2-Trichloroethane	ND	48.5	54.8	88	51.7	55.1	94	18-141	7	40
2-Hexanone	ND	527	274	192 *	532	276	193 *	20-128	1	40
Chlorobenzene	ND	44.0	54.8	80	45.7	55.1	83	15-124	4	40
Ethylbenzene	ND	46.4	54.8	85	48.5	55.1	88	13-128	4	40
1,2,3-Trichloropropane	ND	71.0	54.8	130	68.7	55.1	125	10-159	3	40
2-Chlorotoluene	ND	52.7	54.8	96	51.6	55.1	94	10-140	2	40
1,2-Dichlorobenzene	ND	45.3	54.8	83	45.7	55.1	83	10-124	1	40
Naphthalene	ND	44.4	54.8	81	42.8	55.1	78	10-127	4	40

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Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Extracted: 05/25/2010
Date Analyzed: 05/25/2010

Lab Control Spike/Duplicate Lab Control Spike Summary
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry
Level: Low
Extraction Lot: KWG1004900

	Lab Control Sample KWG1004900-3 Lab Control Spike				Duplicate Lab Control Sample KWG1004900-4 Duplicate Lab Control Spike					
Analyte Name	**Result**	**Expected**	**%Rec**		**Result**	**Expected**	**%Rec**	**%Rec Limits**	**RPD**	**RPD Limit**
Dichlorodifluoromethane	36.3	50.0	73		42.5	50.0	85	38-160	16	40
Chloromethane	39.1	50.0	78		44.8	50.0	90	37-146	14	40
Vinyl Chloride	40.0	50.0	80		45.8	50.0	92	54-127	14	40
Bromomethane	26.7	50.0	53		31.0	50.0	62	22-180	15	40
Chloroethane	37.0	50.0	74		43.1	50.0	86	39-147	15	40
Trichlorofluoromethane	30.1	50.0	60		34.7	50.0	69	51-140	14	40
1,1-Dichloroethene	36.0	50.0	72		41.4	50.0	83	64-152	14	40
Acetone	176	250	70		204	250	82	43-124	15	40
Carbon Disulfide	80.7	100	81		93.4	100	93	55-141	15	40
Methylene Chloride	37.1	50.0	74		44.0	50.0	88	56-131	17	40
trans-1,2-Dichloroethene	37.4	50.0	75	*	42.8	50.0	86	77-127	14	40
1,1-Dichloroethane	38.1	50.0	76		41.9	50.0	84	59-137	9	40
2,2-Dichloropropane	40.3	50.0	81		43.9	50.0	88	74-134	9	40
cis-1,2-Dichloroethene	39.5	50.0	79		43.0	50.0	86	62-138	8	40
2-Butanone (MEK)	208	250	83		231	250	92	54-116	10	40
Bromochloromethane	40.9	50.0	82		45.1	50.0	90	59-130	10	40
Chloroform	41.5	50.0	83		45.0	50.0	90	61-137	8	40
1,1,1-Trichloroethane (TCA)	40.9	50.0	82		44.7	50.0	89	59-146	9	40
Carbon Tetrachloride	40.5	50.0	81		44.2	50.0	88	60-152	9	40
1,1-Dichloropropene	39.0	50.0	78		42.3	50.0	85	52-142	8	40
Benzene	41.8	50.0	84		44.2	50.0	88	56-148	6	40
1,2-Dichloroethane (EDC)	40.0	50.0	80		44.2	50.0	88	57-139	10	40
Trichloroethene (TCE)	42.8	50.0	86		46.2	50.0	92	76-124	8	40
1,2-Dichloropropane	41.3	50.0	83		46.6	50.0	93	58-135	12	40
Dibromomethane	39.8	50.0	80		45.2	50.0	90	56-133	13	40
Bromodichloromethane	41.9	50.0	84		48.0	50.0	96	61-143	14	40
cis-1,3-Dichloropropene	37.5	50.0	75		42.8	50.0	86	58-138	13	40
4-Methyl-2-pentanone (MIBK)	216	250	86		253	250	101	44-140	16	40
Toluene	40.9	50.0	82		47.0	50.0	94	75-117	14	40
trans-1,3-Dichloropropene	52.4	50.0	105		53.6	50.0	107	63-121	2	40
1,1,2-Trichloroethane	42.9	50.0	86		44.5	50.0	89	56-133	4	40
Tetrachloroethene (PCE)	43.4	50.0	87		43.9	50.0	88	66-126	1	40
2-Hexanone	230	250	92		238	250	95	54-122	4	40
1,3-Dichloropropane	44.0	50.0	88		45.5	50.0	91	56-134	3	40
Dibromochloromethane	47.3	50.0	95		48.5	50.0	97	58-140	3	40
1,2-Dibromoethane (EDB)	44.2	50.0	88		45.4	50.0	91	55-134	3	40

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Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Extracted: 05/25/2010
Date Analyzed: 05/25/2010

Lab Control Spike/Duplicate Lab Control Spike Summary
Volatile Organic Compounds

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry
Level: Low
Extraction Lot: KWG1004900

	Lab Control Sample KWG1004900-3 Lab Control Spike			Duplicate Lab Control Sample KWG1004900-4 Duplicate Lab Control Spike					
Analyte Name	**Result**	**Expected**	**%Rec**	**Result**	**Expected**	**%Rec**	**%Rec Limits**	**RPD**	**RPD Limit**
Chlorobenzene	43.4	50.0	87	44.5	50.0	89	70-116	3	40
Ethylbenzene	44.9	50.0	90	45.8	50.0	92	70-118	2	40
1,1,1,2-Tetrachloroethane	47.5	50.0	95	48.3	50.0	97	58-137	2	40
m,p-Xylenes	94.7	100	95	96.1	100	96	58-140	2	40
o-Xylene	46.5	50.0	93	47.7	50.0	95	57-139	3	40
Styrene	47.8	50.0	96	48.5	50.0	97	62-135	1	40
Bromoform	52.5	50.0	105	53.6	50.0	107	62-134	2	40
Isopropylbenzene	48.7	50.0	97	48.6	50.0	97	55-129	0	40
1,1,2,2-Tetrachloroethane	39.3	50.0	79	44.1	50.0	88	53-133	11	40
Bromobenzene	47.9	50.0	96	46.8	50.0	94	52-130	2	40
n-Propylbenzene	49.0	50.0	98	48.6	50.0	97	57-143	1	40
1,2,3-Trichloropropane	45.2	50.0	90	44.8	50.0	90	53-134	1	40
2-Chlorotoluene	49.9	50.0	100	49.0	50.0	98	66-129	2	40
1,3,5-Trimethylbenzene	49.5	50.0	99	49.3	50.0	99	53-141	0	40
4-Chlorotoluene	50.5	50.0	101	50.0	50.0	100	51-134	1	40
tert-Butylbenzene	47.9	50.0	96	48.6	50.0	97	67-131	2	40
1,2,4-Trimethylbenzene	47.4	50.0	95	48.5	50.0	97	55-140	2	40
sec-Butylbenzene	46.4	50.0	93	46.8	50.0	94	55-146	1	40
4-Isopropyltoluene	46.7	50.0	93	47.6	50.0	95	52-140	2	40
1,3-Dichlorobenzene	47.2	50.0	94	47.9	50.0	96	52-131	1	40
1,4-Dichlorobenzene	46.2	50.0	92	46.8	50.0	94	52-128	1	40
n-Butylbenzene	47.9	50.0	96	48.8	50.0	98	53-139	2	40
1,2-Dichlorobenzene	46.3	50.0	93	46.7	50.0	93	53-128	1	40
1,2-Dibromo-3-chloropropane	50.4	50.0	101	52.1	50.0	104	55-127	3	40
1,2,4-Trichlorobenzene	50.2	50.0	100	50.5	50.0	101	57-136	1	40
Hexachlorobutadiene	54.3	50.0	109	54.2	50.0	108	54-140	0	40
Naphthalene	50.4	50.0	101	50.9	50.0	102	54-134	1	40
1,2,3-Trichlorobenzene	52.6	50.0	105	52.5	50.0	105	52-138	0	40

Results flagged with an asterisk (*) indicate values outside control criteria.

Percent recoveries and relative percent differences (RPD) are determined by the software using values in the calculation which have not been rounded.

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	05/25/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	05/25/2010
		Time Analyzed:	12:39

Method Blank Summary
Volatile Organic Compounds

Sample Name:	Method Blank	**File ID:**	J:\MS24\DATA\052510\0525F009.D
Lab Code:	KWG1004900-5	**Instrument ID:**	MS24
Extraction Method:	EPA 5030A	**Level:**	Low
Analysis Method:	8260B	**Extraction Lot:**	KWG1004900

This Method Blank applies to the following analyses:

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed
Lab Control Sample	KWG1004900-3	J:\MS24\DATA\052510\0525F005.D	05/25/10	11:14
Duplicate Lab Control Sample	KWG1004900-4	J:\MS24\DATA\052510\0525F006.D	05/25/10	11:35
D-4-16	K1005244-003	J:\MS24\DATA\052510\0525F012.D	05/25/10	13:42
D-4-16MS	KWG1004900-1	J:\MS24\DATA\052510\0525F013.D	05/25/10	14:40
D-4-16DMS	KWG1004900-2	J:\MS24\DATA\052510\0525F014.D	05/25/10	15:01

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Report

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Extracted:**	05/25/2010
Sample Matrix:	Misc. solid	**Date Analyzed:**	05/25/2010
		Time Analyzed:	11:14

Lab Control Sample Summary
Volatile Organic Compounds

Sample Name:	Lab Control Sample	**File ID:**	J:\MS24\DATA\052510\0525F005.D
Lab Code:	KWG1004900-3	**Instrument ID:**	MS24
Extraction Method:	EPA 5030A	**Level:**	Low
Analysis Method:	8260B	**Extraction Lot:**	KWG1004900

This Lab Control Sample applies to the following analyses:

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed
Method Blank	KWG1004900-5	J:\MS24\DATA\052510\0525F009.D	05/25/10	12:39
D-4-16	K1005244-003	J:\MS24\DATA\052510\0525F012.D	05/25/10	13:42
D-4-16MS	KWG1004900-1	J:\MS24\DATA\052510\0525F013.D	05/25/10	14:40
D-4-16DMS	KWG1004900-2	J:\MS24\DATA\052510\0525F014.D	05/25/10	15:01

Organic Analysis:
Volatile Organic Compounds

Validation Package

Raw Data

COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16
Lab Code: K1005244-003

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	ND	U	5.6	0.51	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	ND	U	5.6	0.35	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	ND	U	5.6	0.69	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	ND	U	5.6	0.35	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	ND	U	5.6	0.24	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	ND	U	5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
Acetone	**13**	J	23	4.3	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	ND	U	5.6	0.29	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**1.7**	J	12	0.48	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	ND	U	5.6	0.39	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	ND	U	5.6	0.23	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	ND	U	5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	ND	U	5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**4.0**	J	23	1.1	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	ND	U	5.6	0.16	1	05/25/10	05/25/10	KWG1004900	
Chloroform	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	ND	U	5.6	0.39	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	ND	U	5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	ND	U	5.6	0.30	1	05/25/10	05/25/10	KWG1004900	
Benzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	ND	U	5.6	0.17	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	ND	U	5.6	0.30	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	ND	U	5.6	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	ND	U	5.6	0.17	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	ND	U	23	0.82	1	05/25/10	05/25/10	KWG1004900	
Toluene	ND	U	5.6	0.20	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	ND	U	5.6	0.38	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	ND	U	5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	ND	U	5.6	0.29	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	ND	U	23	1.1	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	ND	U	5.6	0.23	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16
Lab Code: K1005244-003

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 5030A
Analysis Method: 8260B

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	ND	U	5.6	0.21	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	ND	U	23	0.30	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	ND	U	5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	ND	U	5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	ND	U	5.6	0.41	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	ND	U	5.6	0.15	1	05/25/10	05/25/10	KWG1004900	
Styrene	ND	U	5.6	0.15	1	05/25/10	05/25/10	KWG1004900	
Bromoform	ND	U	5.6	0.34	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	ND	U	23	0.15	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	ND	U	5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	ND	U	23	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	ND	U	5.6	0.42	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	ND	U	23	0.32	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	ND	U	23	0.40	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	ND	U	23	0.29	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	ND	U	23	0.37	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	ND	U	23	0.31	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	ND	U	23	0.35	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	ND	U	23	0.31	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	ND	U	5.6	0.27	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	ND	U	5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	ND	U	23	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	ND	U	5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	ND	U	23	0.73	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	ND	U	23	0.26	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	ND	U	23	0.24	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	ND	U	23	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	ND	U	23	0.24	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16
Lab Code: K1005244-003

Units: ug/Kg
Basis: Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	100	83-128	05/25/10	Acceptable
Toluene-d8	108	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	103	77-124	05/25/10	Acceptable

Comments:

Exception Report

Data File: J:\MS24\DATA\052510\0525F012.D
Lab ID: K1005244-003
RunType: SMPL
Matrix: MISC. SOLID

Date Acquired: 05/25/2010 13:42
Date Quantitated: 05/25/2010 15:05
Batch ID: KWG1004880
Analysis Method: 8260B
ListJoinID: LJ1601

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Window	NA	NA	NA	x	
Analytical Holding Time	NA	NA	NA	x	
Preparation Holding Time	NA	NA	NA	x	
Pre-Preparation Holding Time	NA	NA	NA	x	
ICAL Pass/Fail	NA	NA	NA	x	
ICAL Average RSD	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Initial Calibration Minimum RF	NA	NA	NA	x	
Initial Calibration SPCC/CCC	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Continuing Calibration Minimum RF	NA	NA	NA	x	
Continuing Calibration SPCC/CCC	NA	NA	NA	x	
Method Blank	NA	NA	NA	x	
MB Surrogate Recovery	NA	NA	NA	x	
Lab Control Spike	NA	NA	NA		x
Duplicate Lab Control Spike	NA	NA	NA	x	
Internal Standards	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Relative Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Analyte Exceptions

Exception Categories	Analyte Name	Result	Low Limit	High Limit	Corrective Action
Lab Control Spike	trans-1,2-Dichloroethene	75	77	127	adv cmpd

Primary Review: KL 5/25/10

Secondary Review: Cam 5/25/10

Printed: 05/25/2010 15:25:31
u:\Stealth\Crystal.rpt\except2.rpt

Quantitation Report

Bottle ID:		**Tier:**	V	**Matrix:**	MISC. SOLID
Prod Code:	8260B VOC_FP	**Collect Date:**	05/19/2010	**Receive Date:**	05/21/2010

Analysis Lot:	KWG1004880	**Prep Lot:**	KWG1004900	**Report Group:**	K1005244
Analysis Method:	8260B	**Prep Method:**	EPA 5030A		
Prep Ref:	911424	**Prep Date:**	05/25/2010		

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:	Volatile Organic Compounds	**Report List ID:**	LJ1601
Tune Ref:	J:\MS24\DATA\052510\0525F002.D	**Method ID:**	MJ120
MB Ref:	J:\MS24\DATA\052510\0525F009.D	**Quant based on Report List**	

Data File:	J:\MS24\DATA\052510\0525F012.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 13:42	**Quant Date:**	05/25/2010 15:05	**Vial:**	11
Run Type:	SMPL			**Dilution:**	1.0
Lab ID:	K1005244-003			**Soln Conc. Units:**	PPB

Internal Standard Compounds

IS Ref	Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Area Criteria
1	Fluorobenzene	4.02	0.00	96	342019	50.00	OK
2	Chlorobenzene-d5	6.65	0.00	82	146554	50.00	OK
3	1,4-Dichlorobenzene-d4	8.98	0.00	152	116925	50.00	OK

Surrogate Compounds

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	%Rec	%Rec Limits		Rpt?
1	Dibromofluoromethane	3.53	-0.01	0.00	113	85826	50.21	100	83-128	OK	
1	Toluene-d8	5.31	0.00	0.00	98	323161	53.90	108	90-125	OK	
2	4-Bromofluorobenzene	7.83	0.00	0.00	95	115686	51.43	103	77-124	OK	

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Dichlorodifluoromethane				85	0		0.25	U	
1	Chloromethane	0.99		0.00	50	944	0.3500	0.51	U	
1	Vinyl Chloride				62	0		0.35	U	
1	Bromomethane				96	0		0.69	U	
1	Chloroethane				64	0		0.35	U	
1	Trichlorofluoromethane				101	0		0.24	U	
1	1,1-Dichloroethene				96	0		0.26	U	
1	Acetone	1.98		0.00	43	9363	11.76	13	J	
1	Carbon Disulfide	2.01		0.00	76	775	0.1400	0.29	U	
1	Methylene Chloride	2.27		0.00	84	3164	1.54	1.7	J	
1	trans-1,2-Dichloroethene				96	0		0.39	U	
1	1,1-Dichloroethane				63	0		0.23	U	
1	2,2-Dichloropropane				77	0d		0.40	U	
1	cis-1,2-Dichloroethene				96	0		0.26	U	
1	2-Butanone (MEK)	3.21		0.00	72	921	3.60	4.0	J	

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E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F012.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 13:42	**Quant Date:**	05/25/2010 15:05	**Vial:**	11
Run Type:	SMPL			**Dilution:**	1.0
Lab ID:	K1005244-003			**Soln Conc. Units:**	PPB

Target Compounds — **Final Conc. Units:** ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Bromochloromethane				128	0		0.16	U	
1	Chloroform				83	0		0.25	U	
1	1,1,1-Trichloroethane (TCA)				97	0		0.39	U	
1	Carbon Tetrachloride				117	0		0.40	U	
1	1,1-Dichloropropene				75	0		0.30	U	
1	Benzene				78	0		0.25	U	
1	1,2-Dichloroethane (EDC)				62	0d		0.17	U	
1	Trichloroethene (TCE)				95	0		0.30	U	
1	1,2-Dichloropropane				63	0		0.25	U	
1	Dibromomethane				93	0		0.31	U	
1	Bromodichloromethane				83	0		0.17	U	
1	cis-1,3-Dichloropropene				75	0		0.25	U	
1	4-Methyl-2-pentanone (MIBK)				58	0d		0.82	U	
1	Toluene	5.37		0.00	92	649	0.1500	0.20	U	
2	trans-1,3-Dichloropropene				75	0		0.38	U	
2	1,1,2-Trichloroethane				83	0		0.26	U	
2	Tetrachloroethene (PCE)				164	0		0.29	U	
2	2-Hexanone				57	0d		1.1	U	
2	1,3-Dichloropropane				76	0		0.23	U	
2	Dibromochloromethane				129	0		0.21	U	
2	1,2-Dibromoethane (EDB)				107	0		0.30	U	
2	Chlorobenzene				112	0		0.28	U	
2	Ethylbenzene				106	0		0.25	U	
2	1,1,1,2-Tetrachloroethane				131	0		0.28	U	
2	m,p-Xylenes				106	0		0.41	U	
2	o-Xylene				106	0		0.15	U	
2	Styrene				103	0		0.15	U	
2	Bromoform				173	0		0.34	U	
2	Isopropylbenzene				105	0		0.15	U	
3	1,1,2,2-Tetrachloroethane				83	0		0.40	U	
3	Bromobenzene				156	0		0.25	U	
3	n-Propylbenzene				91	0		0.40	U	
3	1,2,3-Trichloropropane				110	0		0.42	U	
3	2-Chlorotoluene				91	0		0.32	U	
3	1,3,5-Trimethylbenzene				105	0		0.40	U	
3	4-Chlorotoluene				91	0		0.29	U	
3	tert-Butylbenzene				119	0		0.37	U	
3	1,2,4-Trimethylbenzene				105	0		0.31	U	
3	sec-Butylbenzene				105	0		0.35	U	
3	4-Isopropyltoluene				119	0		0.31	U	
3	1,3-Dichlorobenzene				146	0		0.27	U	
3	1,4-Dichlorobenzene				146	0		0.28	U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
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NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F012.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 13:42	**Quant Date:**	05/25/2010 15:05	**Vial:**	11
Run Type:	SMPL			**Dilution:**	1.0
Lab ID:	K1005244-003			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
3	n-Butylbenzene				91	0		0.36	U	
3	1,2-Dichlorobenzene				146	0		0.25	U	
3	1,2-Dibromo-3-chloropropane				155	0		0.73	U	
3	1,2,4-Trichlorobenzene				180	0		0.26	U	
3	Hexachlorobutadiene				225	0		0.24	U	
3	Naphthalene				128	0		0.40	U	
3	1,2,3-Trichlorobenzene				180	0		0.24	U	

Prep Amount:	5.00 g	**Dilution:**	1.0
Prep Final Vol:	5.0 ml	**Unit Factor:**	1
Solids:	90.7 %		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

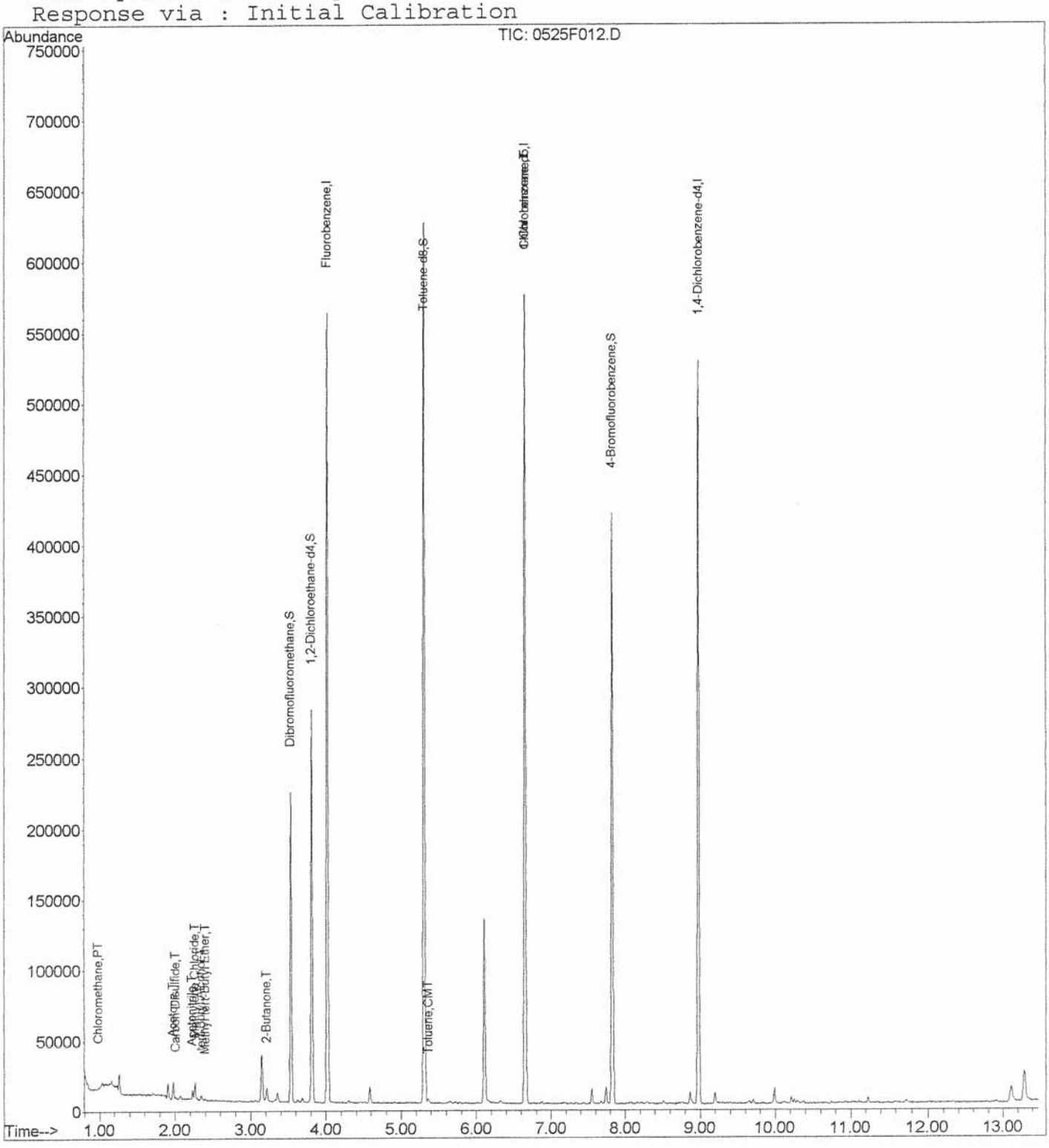
Quantitation Report (QT Reviewed)

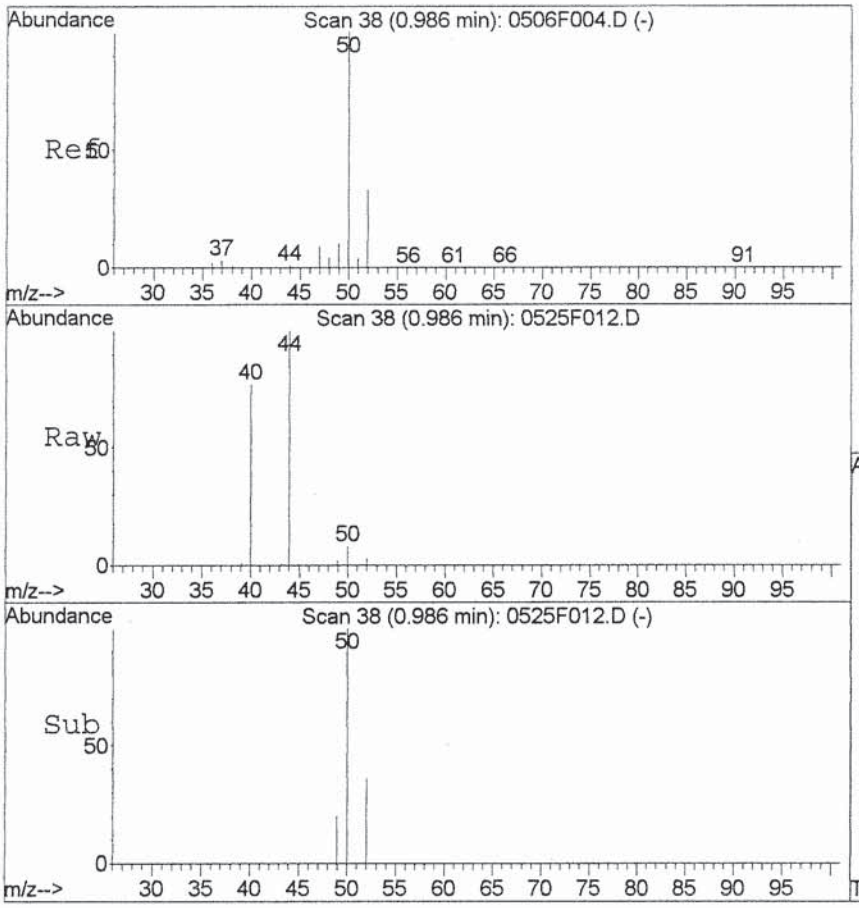
Data File : J:\MS24\DATA\052510\0525F012.D
Acq On : 25 May 2010 1:42 pm
Sample : K5244-003
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 15:05 2010

Vial: 11
Operator: KR
Inst : MS24
Multiplr: 1.00

Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Initial Calibration

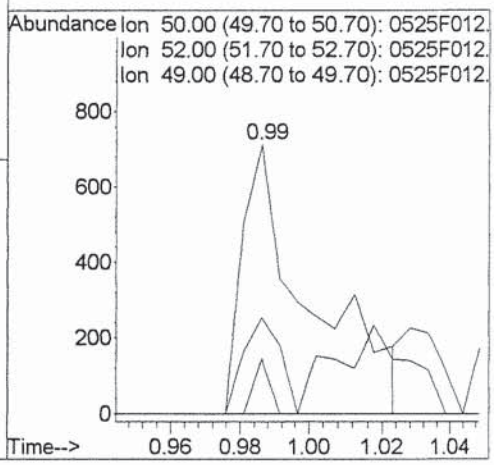


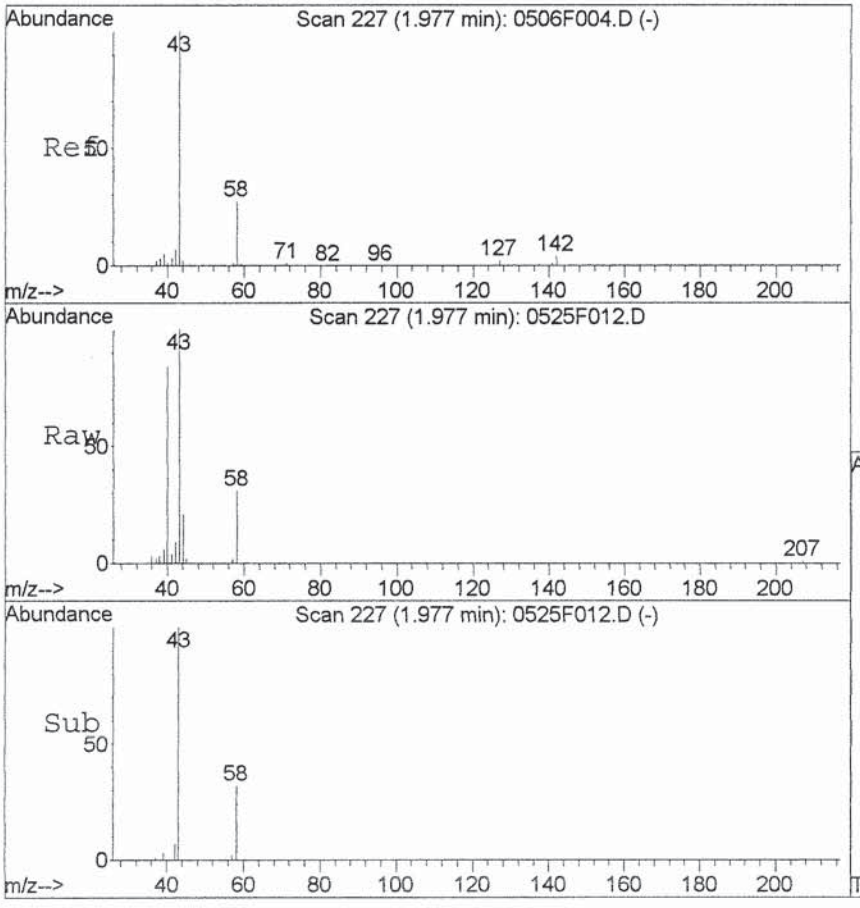


#3
Chloromethane
Concen: 0.35 PPB
RT: 0.99 min Scan# 38
Delta R.T. 0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 50 Resp: 944

Ion	Ratio	Lower	Upper
50	100		
52	35.5	2.6	62.6
49	20.4	0.0	40.5

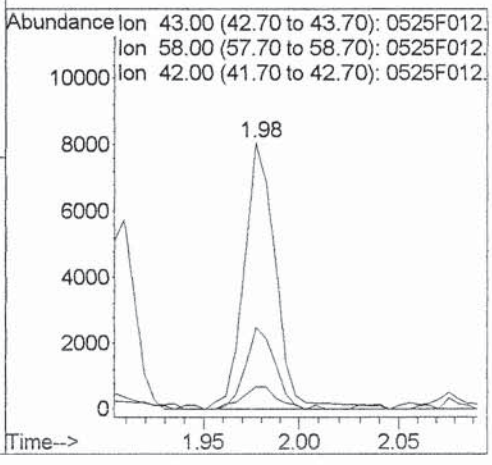


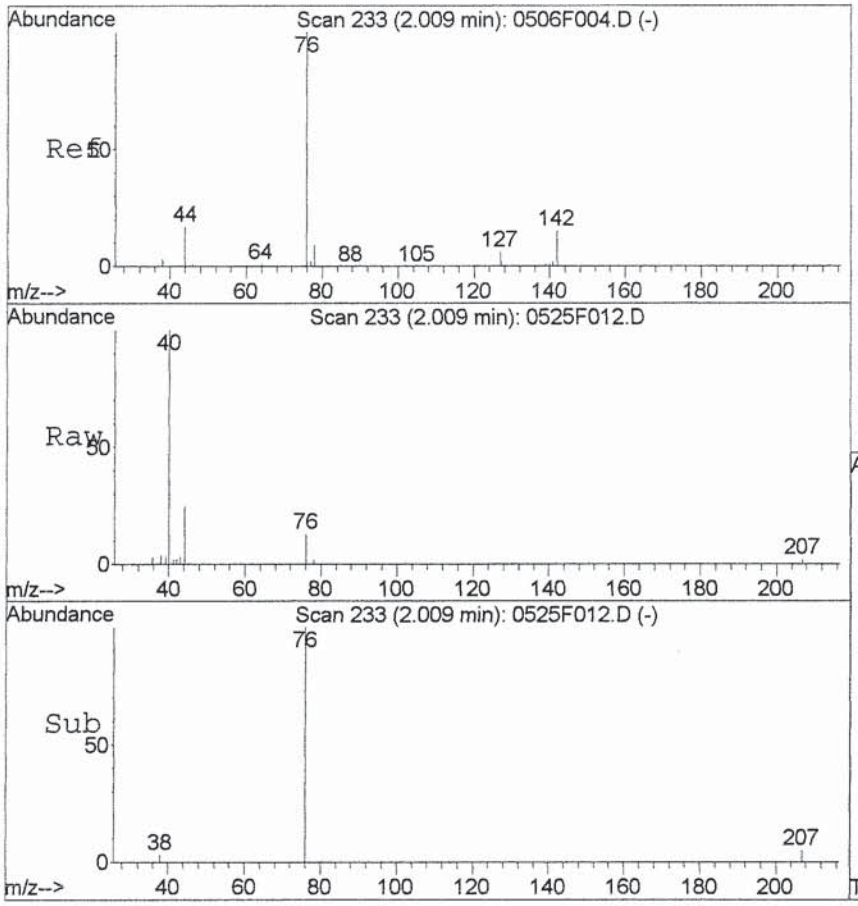


#13
Acetone
Concen: 11.76 PPB
RT: 1.98 min Scan# 227
Delta R.T. 0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 43 Resp: 9363

Ion	Ratio	Lower	Upper
43	100		
58	30.9	0.0	57.0
42	8.6	0.0	37.4

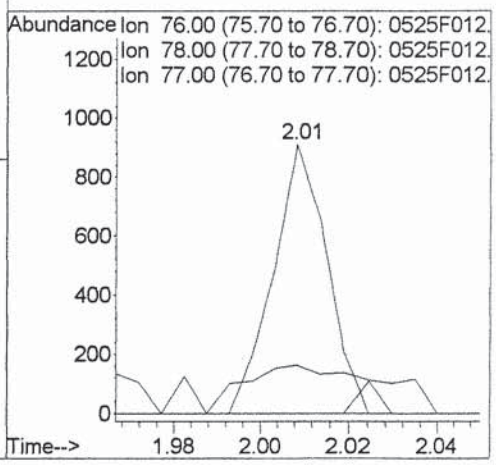


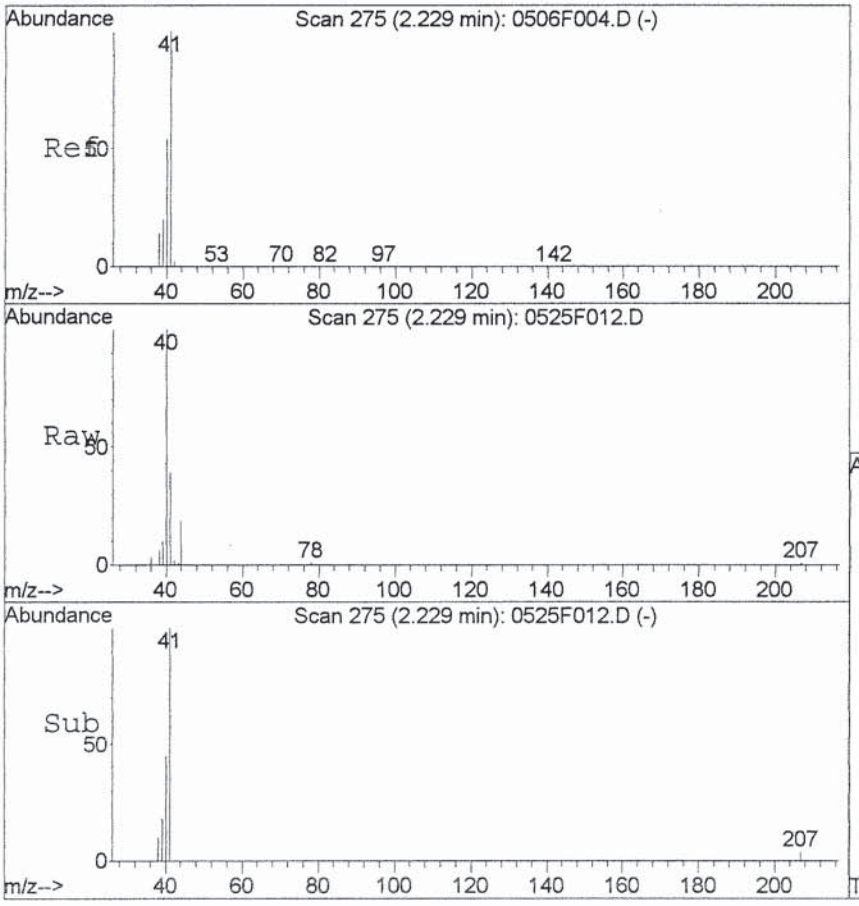


#15
Carbon Disulfide
Concen: 0.14 PPB
RT: 2.01 min Scan# 233
Delta R.T. 0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 76 Resp: 775

Ion	Ratio	Lower	Upper
76	100		
78	17.9	0.0	38.7
77	0.0	0.0	32.5

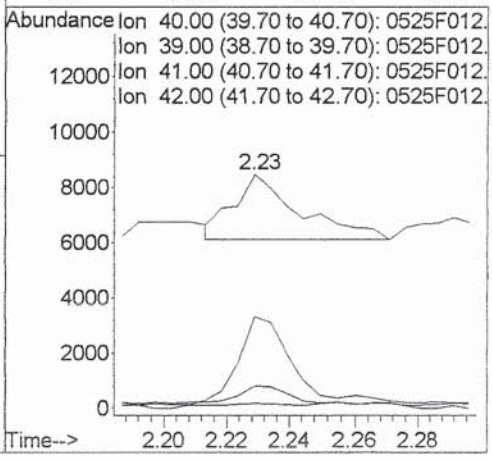


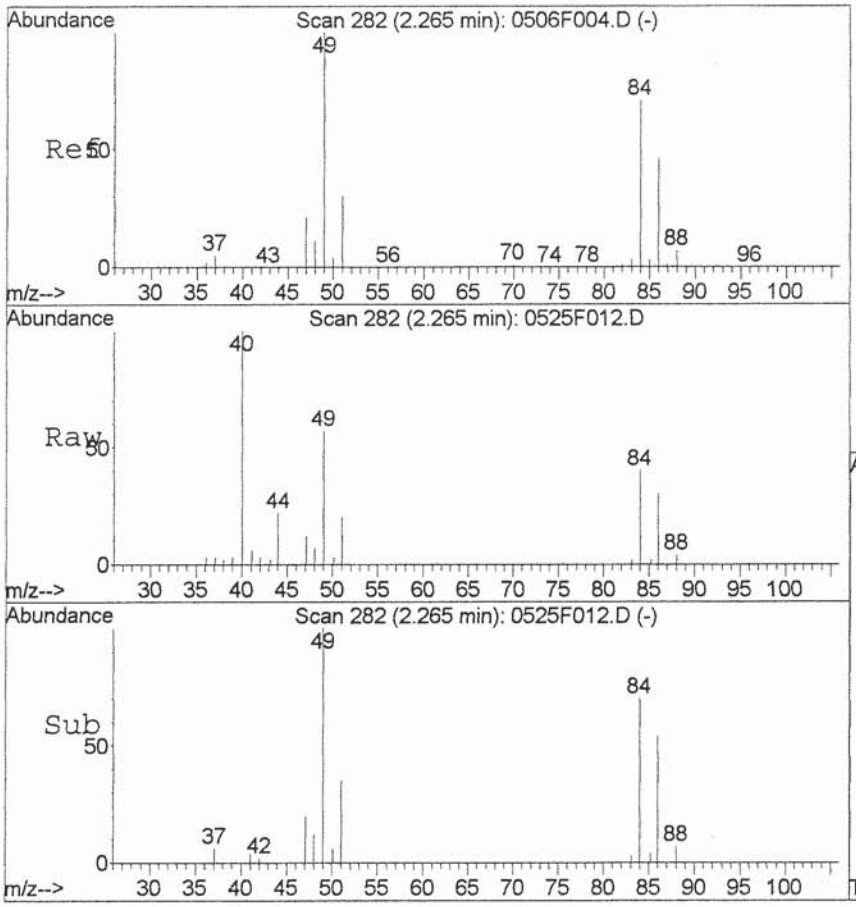


#17
Acetonitrile
Concen: 21.95 PPB
RT: 2.23 min Scan# 275
Delta R.T. 0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 40 Resp: 3433

Ion	Ratio	Lower	Upper
40	100		
39	28.8	6.4	66.4
41	115.2	159.7	219.7#
42	0.0	0.0	34.7

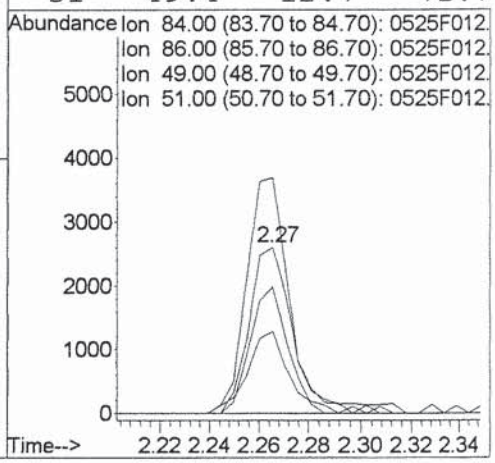


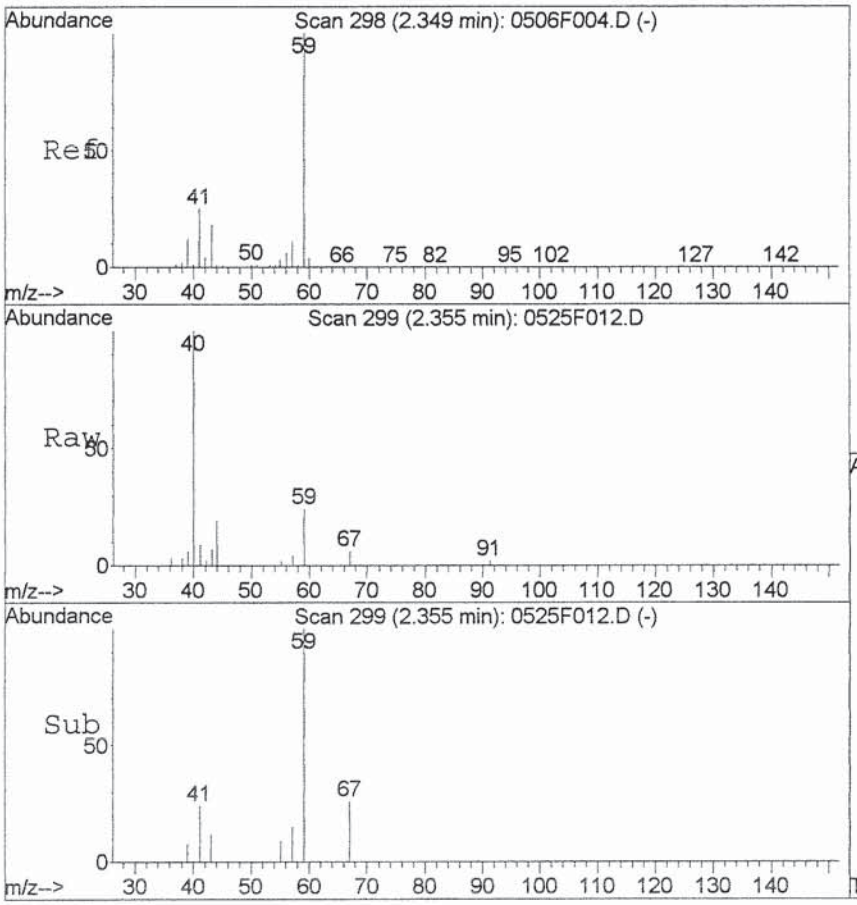


#18
Methylene Chloride
Concen: 1.54 PPB
RT: 2.27 min Scan# 282
Delta R.T. 0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 84 Resp: 3164

Ion	Ratio	Lower	Upper
84	100		
86	76.3	34.4	94.4
49	142.0	109.9	169.9
51	49.4	12.7	72.7

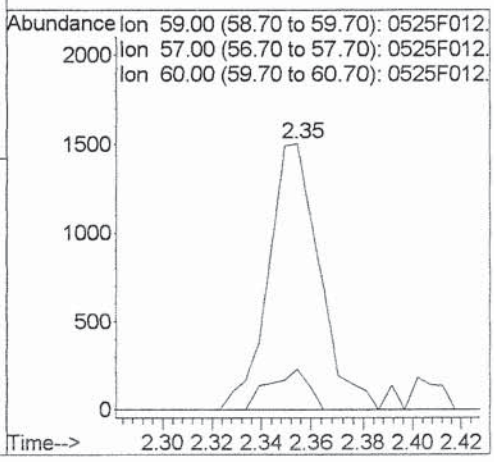


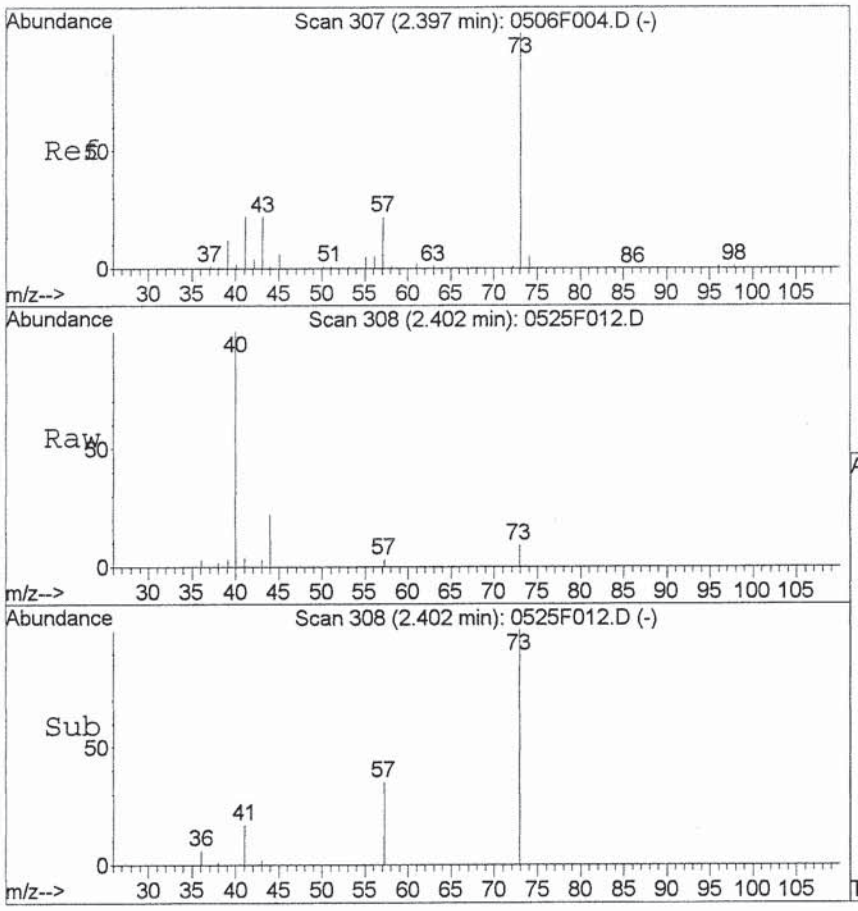


#19
tert-Butyl Alcohol
Concen: 7.47 PPB
RT: 2.35 min Scan# 299
Delta R.T. 0.01 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 59 Resp: 2133

Ion	Ratio	Lower	Upper
59	100		
57	15.4	0.0	41.1
60	0.0	0.0	33.6

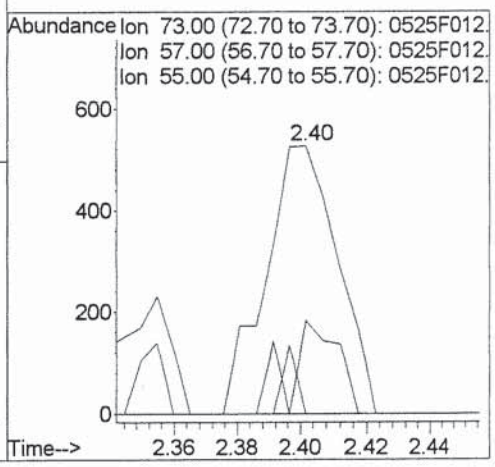


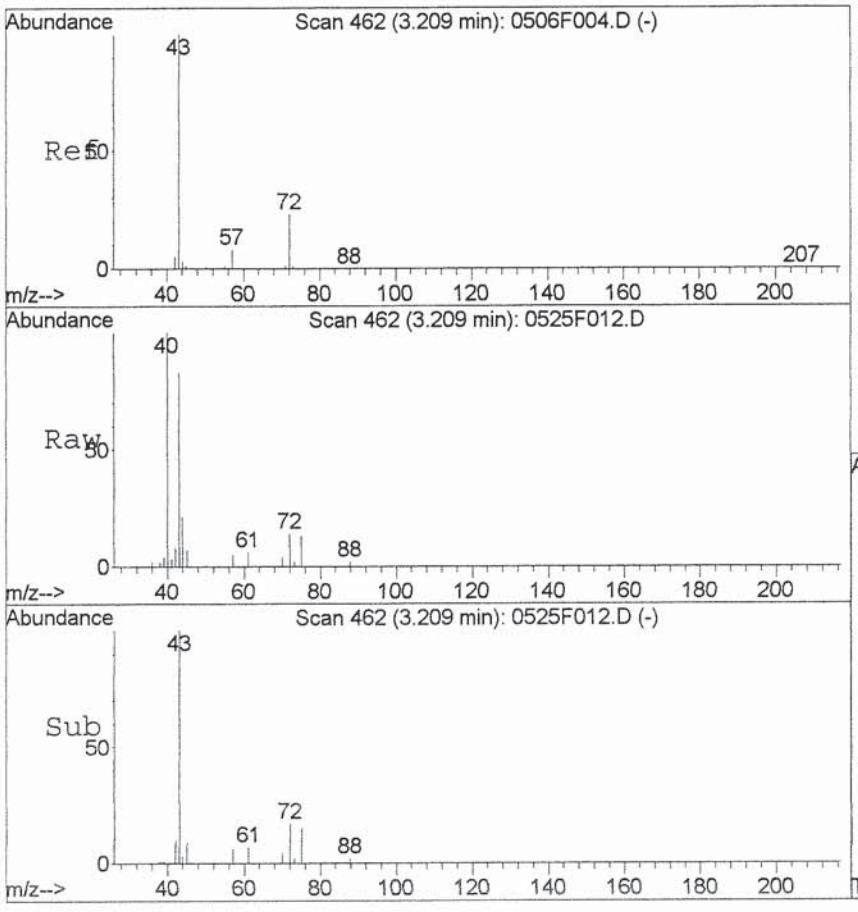


#21
Methyl tert-Butyl Ether
Concen: 0.14 PPB
RT: 2.40 min Scan# 308
Delta R.T. 0.01 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 73 Resp: 819

Ion	Ratio	Lower	Upper
73	100		
57	34.7	0.0	53.3
55	0.0	0.0	35.5

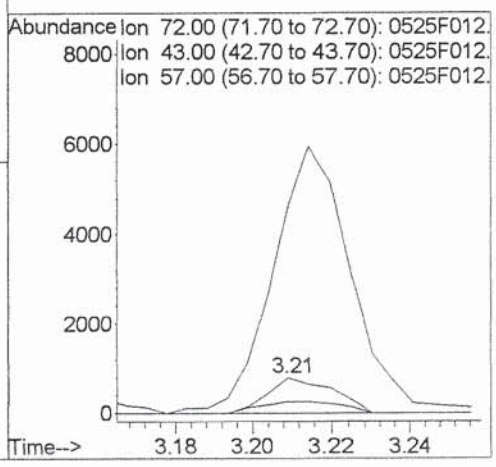


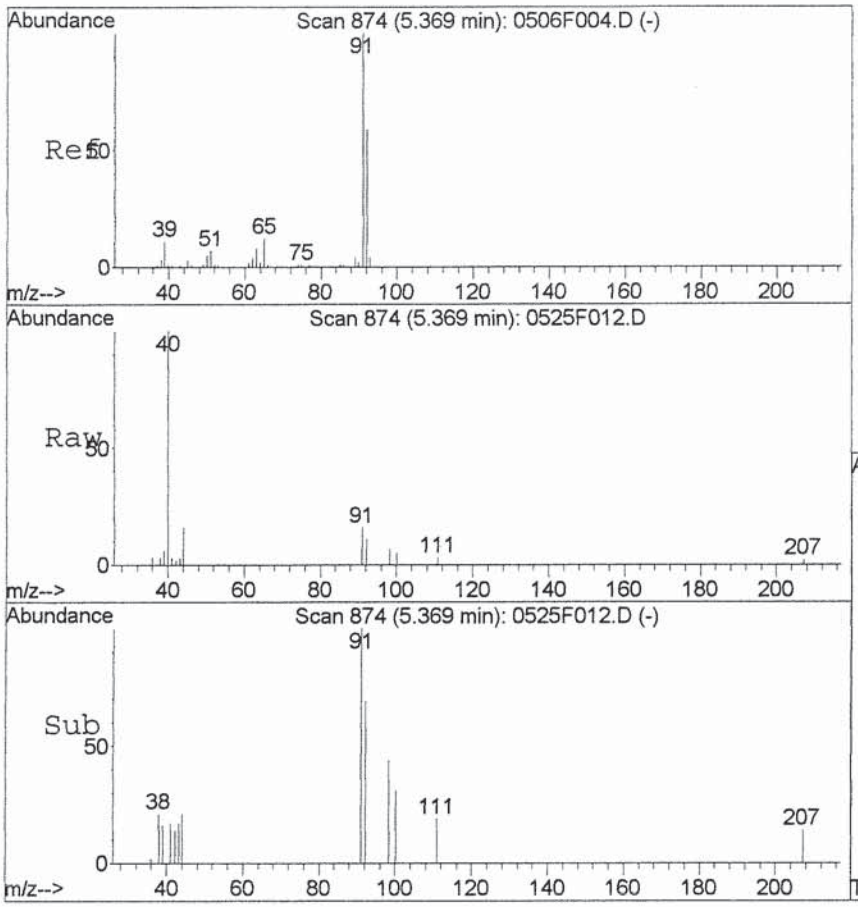


#31
2-Butanone
Concen: 3.60 PPB
RT: 3.21 min Scan# 462
Delta R.T. -0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 72 Resp: 921

Ion	Ratio	Lower	Upper
72	100		
43	566.6	403.8	463.8#
57	33.7	3.3	63.3

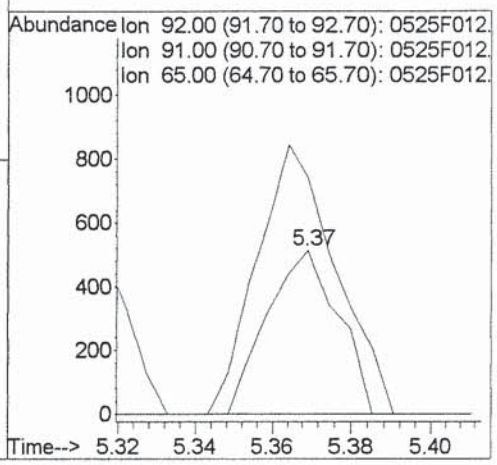


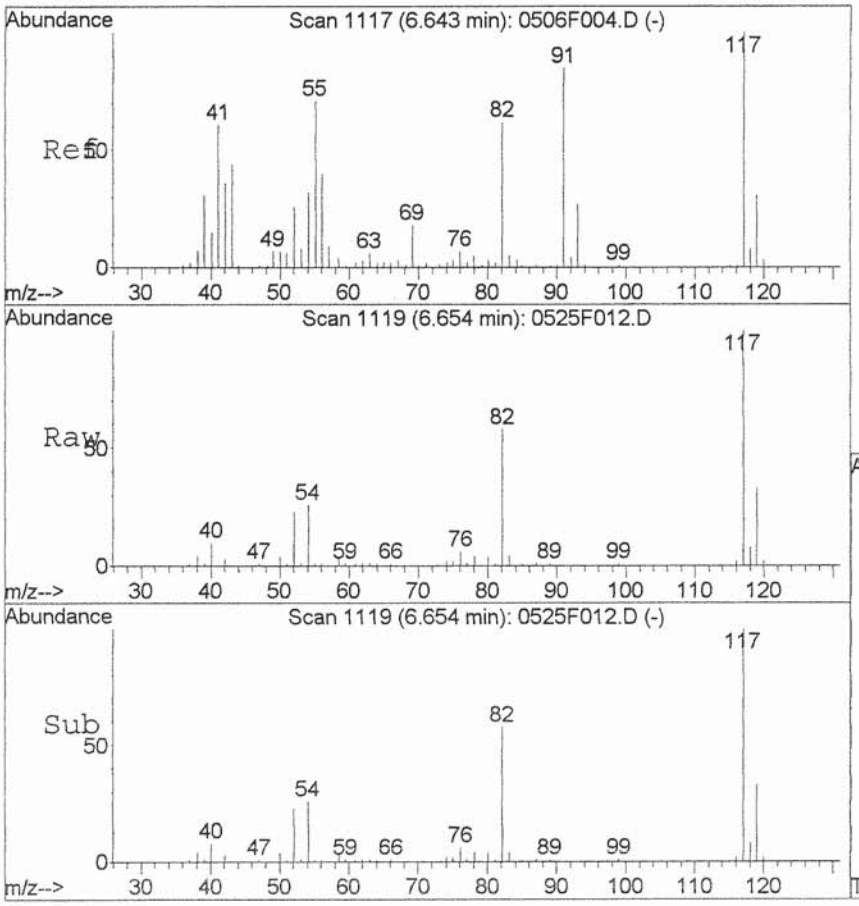


#58
Toluene
Concen: 0.15 PPB
RT: 5.37 min Scan# 874
Delta R.T. 0.00 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 92 Resp: 649

Ion	Ratio	Lower	Upper
92	100		
91	144.6	138.9	198.9
65	0.0	0.0	50.5

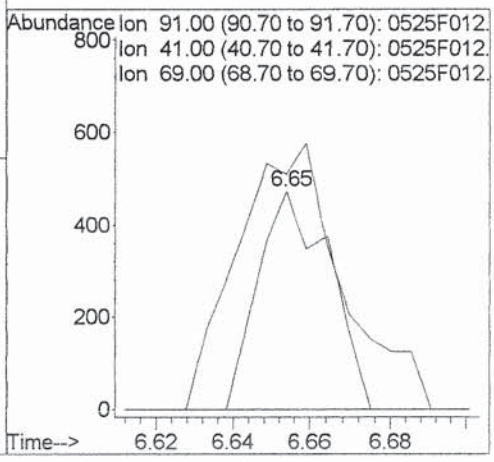




#69
1-Chlorohexane
Concen: 0.29 PPB
RT: 6.65 min Scan# 1119
Delta R.T. 0.01 min
Lab File: 0525F012.D
Acq: 25 May 2010 1:42 pm

Tgt Ion: 91 Resp: 600

Ion	Ratio	Lower	Upper
91	100		
41	81.4	41.3	101.3
69	0.0	0.0	51.4



COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Method Blank
Lab Code: KWG1004900-5

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	ND	U	5.0	0.46	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	ND	U	5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	ND	U	5.0	0.62	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	ND	U	5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	ND	U	5.0	0.21	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	ND	U	5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Acetone	**6.3**	J	20	3.9	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	ND	U	5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**1.5**	J	10	0.43	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	ND	U	5.0	0.35	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	ND	U	5.0	0.20	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	ND	U	5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	ND	U	5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**2.3**	J	20	0.99	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	ND	U	5.0	0.14	1	05/25/10	05/25/10	KWG1004900	
Chloroform	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	ND	U	5.0	0.35	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	ND	U	5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	ND	U	5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
Benzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	ND	U	5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	ND	U	5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	ND	U	5.0	0.28	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	ND	U	5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	ND	U	20	0.74	1	05/25/10	05/25/10	KWG1004900	
Toluene	ND	U	5.0	0.18	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	ND	U	5.0	0.34	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	ND	U	5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	ND	U	5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	ND	U	20	0.93	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	ND	U	5.0	0.20	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Method Blank
Lab Code: KWG1004900-5

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	ND	U	5.0	0.19	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	ND	U	20	0.27	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	ND	U	5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	ND	U	5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	ND	U	5.0	0.37	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	ND	U	5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Styrene	ND	U	5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Bromoform	ND	U	5.0	0.30	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	ND	U	20	0.13	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	ND	U	5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	ND	U	20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	ND	U	5.0	0.38	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	ND	U	20	0.29	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	ND	U	20	0.36	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	ND	U	20	0.26	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	ND	U	20	0.33	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	ND	U	20	0.28	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	ND	U	20	0.31	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	ND	U	20	0.28	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	ND	U	5.0	0.24	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	ND	U	5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	ND	U	20	0.32	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	ND	U	5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	ND	U	20	0.66	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	ND	U	20	0.23	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	ND	U	20	0.21	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	ND	U	20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	ND	U	20	0.21	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Method Blank
Lab Code: KWG1004900-5

Units: ug/Kg
Basis: Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	100	83-128	05/25/10	Acceptable
Toluene-d8	110	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	104	77-124	05/25/10	Acceptable

Comments:

Exception Report

Data File: J:\MS24\DATA\052510\0525F009.D
Lab ID: KWG1004900-5
RunType: MB
Matrix: SEDIMENT

Date Acquired: 05/25/2010 12:39
Date Quantitated: 05/25/2010 14:57
Batch ID: KWG1004880
Analysis Method: 8260B
MethodJoinID: MJ120

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Window	NA	NA	NA	x	
Analytical Holding Time	NA	NA	NA	x	
ICAL Pass/Fail	NA	NA	NA	x	
ICAL Average RSD	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Initial Calibration Minimum RF	NA	NA	NA		x
Initial Calibration SPCC/CCC	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Continuing Calibration Minimum RF	NA	NA	NA		x
Continuing Calibration SPCC/CCC	NA	NA	NA	x	
Internal Standards	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Relative Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Analyte Exceptions

Exception Categories	Analyte Name	Result	Low Limit	High Limit	Corrective Action
Initial Calibration Minimum RF	1,4-Dioxane	0.0037	0.01	NA	NT
Continuing Calibration Minimum RF	1,4-Dioxane	0.0034	0.01	NA	J

Primary Review: 5/25/10

Secondary Review: 5/25/10

Printed: 05/25/2010 15:15:38
u:\Stealth\Crystal.rpt\except2.rpt

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	SEDIMENT
Prod Code:	8260B VOC_FP	**Collect Date:**		**Receive Date:**	05/25/2010

Analysis Lot:	KWG1004880	**Prep Lot:**	KWG1004900	**Report Group:**	
Analysis Method:	8260B	**Prep Method:**	EPA 5030A		
Prep Ref:	911432	**Prep Date:**	05/25/2010		

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:			
Tune Ref:	J:\MS24\DATA\052510\0525F002.D	**Method ID:**	MJ120
MB Ref:		**Quant based on Method**	

Data File:	J:\MS24\DATA\052510\0525F009.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 12:39	**Quant Date:**	05/25/2010 14:57	**Vial:**	8
Run Type:	MB			**Dilution:**	1.0
Lab ID:	KWG1004900-5			**Soln Conc. Units:**	PPB

Internal Standard Compounds

IS Ref	Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Area Criteria
1	Fluorobenzene	4.02	0.00	96	348737	50.00	OK
2	Chlorobenzene-d5	6.65	0.00	82	149494	50.00	OK
3	1,4-Dichlorobenzene-d4	8.98	0.00	152	128420	50.00	OK

Surrogate Compounds

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	%Rec	%Rec Limits		Rpt?
1	Dibromofluoromethane	3.53	-0.01	0.00	113	86905	49.86	100	43-153	OK	
1	1,2-Dichloroethane-d4	3.81	0.00	0.00	65	116863	46.45	93	64-142	OK	
1	Toluene-d8	5.31	0.00	0.00	98	336543	55.05	110	49-151	OK	
2	4-Bromofluorobenzene	7.83	0.00	0.00	95	118745	51.75	104	33-150	OK	

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Dichlorodifluoromethane				85	0		0.22	U	
1	Chloromethane				50	0		0.46	U	
1	Vinyl Chloride				62	0		0.31	U	
1	Bromomethane				96	0		0.62	U	
1	Chloroethane				64	0		0.31	U	
1	Dichlorofluoromethane (CFC 21				67	0		0.35	U	
1	Trichlorofluoromethane				101	0		0.21	U	
1	Diethyl Ether				59	0		5.0	U	
1	Acrolein				56	0		2.6	U	
1	Trichlorotrifluoroethane				151	0		0.56	U	
1	1,1-Dichloroethene				96	0		0.23	U	
1	Acetone	1.98		0.00	43	5072	6.25	6.25	J	
1	Iodomethane (Methyl Iodide)				142	0		0.57	U	
1	Carbon Disulfide				76	0		0.26	U	

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F009.D		**Instrument:** MS24
Acqu Date:	05/25/2010 12:39	**Quant Date:** 05/25/2010 14:57	**Vial:** 8
Run Type:	MB		**Dilution:** 1.0
Lab ID:	KWG1004900-5		**Soln Conc. Units:** PPB

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	3-Chloro-1-propene				76	0		0.32	U	
1	Acetonitrile				40	0d		23	U	
1	Methylene Chloride	2.27		0.00	84	3082	1.47	1.47	J	
1	tert-Butyl Alcohol	2.36	0.01	0.00	59	1048	3.60	3.60	J	
1	Acrylonitrile				53	0		1.2	U	
1	Methyl tert-Butyl Ether	2.40		0.00	73	914	0.1500	0.18	U	
1	trans-1,2-Dichloroethene				96	0		0.35	U	
1	n-Hexane				57	0		0.22	U	
1	Diisopropyl Ether				45	0		0.16	U	
1	1,1-Dichloroethane				63	0		0.20	U	
1	Vinyl Acetate				86	0		0.92	U	
1	Chloroprene				53	0		0.76	U	
1	tert-Butyl Ethyl Ether				59	0		0.14	U	
1	2,2-Dichloropropane				77	0		0.36	U	
1	cis-1,2-Dichloroethene				96	0		0.23	U	
1	2-Butanone (MEK)	3.21		0.00	72	609	2.33	2.33	J	
1	Propionitrile				54	0		1.2	U	
1	Methacrylonitrile				67	0d		0.62	U	
1	Bromochloromethane				128	0		0.14	U	
1	Chloroform				83	0		0.22	U	
1	1,1,1-Trichloroethane (TCA)				97	0		0.35	U	
1	Carbon Tetrachloride				117	0		0.36	U	
1	1,1-Dichloropropene				75	0		0.27	U	
1	Isobutanol				43	0d		12	U	
1	Benzene				78	0		0.22	U	
1	1,2-Dichloroethane (EDC)				62	0d		0.15	U	
1	tert-Amyl Methyl Ether				55	0		0.32	U	
1	Trichloroethene (TCE)				95	0		0.27	U	
1	1,2-Dichloropropane				63	0		0.22	U	
1	Dibromomethane				93	0		0.28	U	
1	Methyl Methacrylate				69	0		0.39	U	
1	1,4-Dioxane				88	0		14	U	
1	Bromodichloromethane				83	0		0.15	U	
1	2-Nitropropane				41	0d		1.1	U	
1	2-Chloroethyl Vinyl Ether				63	0		0.48	U	
1	cis-1,3-Dichloropropene				75	0		0.22	U	
1	4-Methyl-2-pentanone (MIBK)				58	0d		0.74	U	
1	Toluene	5.36	-0.01	0.00	92	644	0.1400	0.18	U	
2	trans-1,3-Dichloropropene				75	0		0.34	U	
2	Ethyl Methacrylate				69	0d		0.34	U	
2	1,1,2-Trichloroethane				83	0		0.23	U	
2	Tetrachloroethene (PCE)				164	0		0.26	U	

U: Undetected at or above MDL
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B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F009.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 12:39	**Quant Date:**	05/25/2010 14:57	**Vial:**	8
Run Type:	MB			**Dilution:**	1.0
Lab ID:	KWG1004900-5			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2	2-Hexanone				57	0d		0.93	U	
2	1,3-Dichloropropane				76	0		0.20	U	
2	Dibromochloromethane				129	0		0.19	U	
2	1,2-Dibromoethane (EDB)				107	0		0.27	U	
2	1-Chlorohexane	6.65	0.01	0.00	91	682	0.3200	0.320	J	
2	Chlorobenzene				112	0		0.25	U	
2	Ethylbenzene				106	0		0.22	U	
2	1,1,1,2-Tetrachloroethane				131	0		0.25	U	
2	m,p-Xylenes				106	0		0.37	U	
2	o-Xylene				106	0		0.13	U	
2	Styrene				103	0		0.13	U	
2	Bromoform				173	0		0.30	U	
2	Isopropylbenzene				105	0		0.13	U	
2	cis-1,4-Dichloro-2-butene				89	0		1.2	U	
3	1,1,2,2-Tetrachloroethane				83	0		0.36	U	
3	trans-1,4-Dichloro-2-butene				53	0		0.27	U	
3	Bromobenzene				156	0		0.22	U	
3	n-Propylbenzene				91	0		0.36	U	
3	1,2,3-Trichloropropane				110	0		0.38	U	
3	2-Chlorotoluene				91	0		0.29	U	
3	1,3,5-Trimethylbenzene				105	0		0.36	U	
3	4-Chlorotoluene				91	0		0.26	U	
3	tert-Butylbenzene				119	0		0.33	U	
3	1,2,4-Trimethylbenzene				105	0d		0.28	U	
3	sec-Butylbenzene				105	0d		0.31	U	
3	4-Isopropyltoluene				119	0		0.28	U	
3	1,3-Dichlorobenzene				146	0		0.24	U	
3	1,4-Dichlorobenzene				146	0		0.25	U	
3	n-Butylbenzene				91	0		0.32	U	
3	1,2-Dichlorobenzene				146	0		0.22	U	
3	1,2-Dibromo-3-chloropropane				155	0		0.66	U	
3	1,2,4-Trichlorobenzene				180	0		0.23	U	
3	Hexachlorobutadiene				225	0		0.21	U	
3	Naphthalene	11.08	-0.01	0.00	128	881	0.1700	0.36	U	
3	1,2,3-Trichlorobenzene				180	0		0.21	U	
	4-Vinyl-1-cyclohexene				0	0		5.0	U	NR
	Benzyl Chloride				0	0		5.0	U	NR
	Hexachloroethane				0	0		1.0	U	NR
	2,2,4-Trimethylpentane				0	0		5.0	U	NR
	1,3-Butadiene				0	0		2.0	U	NR
	1,1,2-Trifluoroethane				0	0		10	U	NR
	n-Butyl Alcohol				0	0		250	U	NR

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?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F009.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 12:39	**Quant Date:**	05/25/2010 14:57	**Vial:**	8
Run Type:	MB			**Dilution:**	1.0
Lab ID:	KWG1004900-5			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2-Propanol			0	0		200	U	NR
Ethyl Alcohol			0	0		200	U	NR
2,2-Dichloro-1,1,1-trifluoroetha			0	0		5.0	U	NR
Cyclohexanone			0	0		10	U	NR
2-Methylpentane			0	0		5.0	U	NR
n-Heptane			0	0		5.0	U	NR
3-Methylpentane			0	0		5.0	U	NR
Methylcyclopentane			0	0		5.0	U	NR
Ethyl Acetate			0	0		20	U	NR

Prep Amount:	5.00 g	**Dilution:**	1.0
Prep Final Vol:	5.0 ml	**Unit Factor:**	1
Solids:	%		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

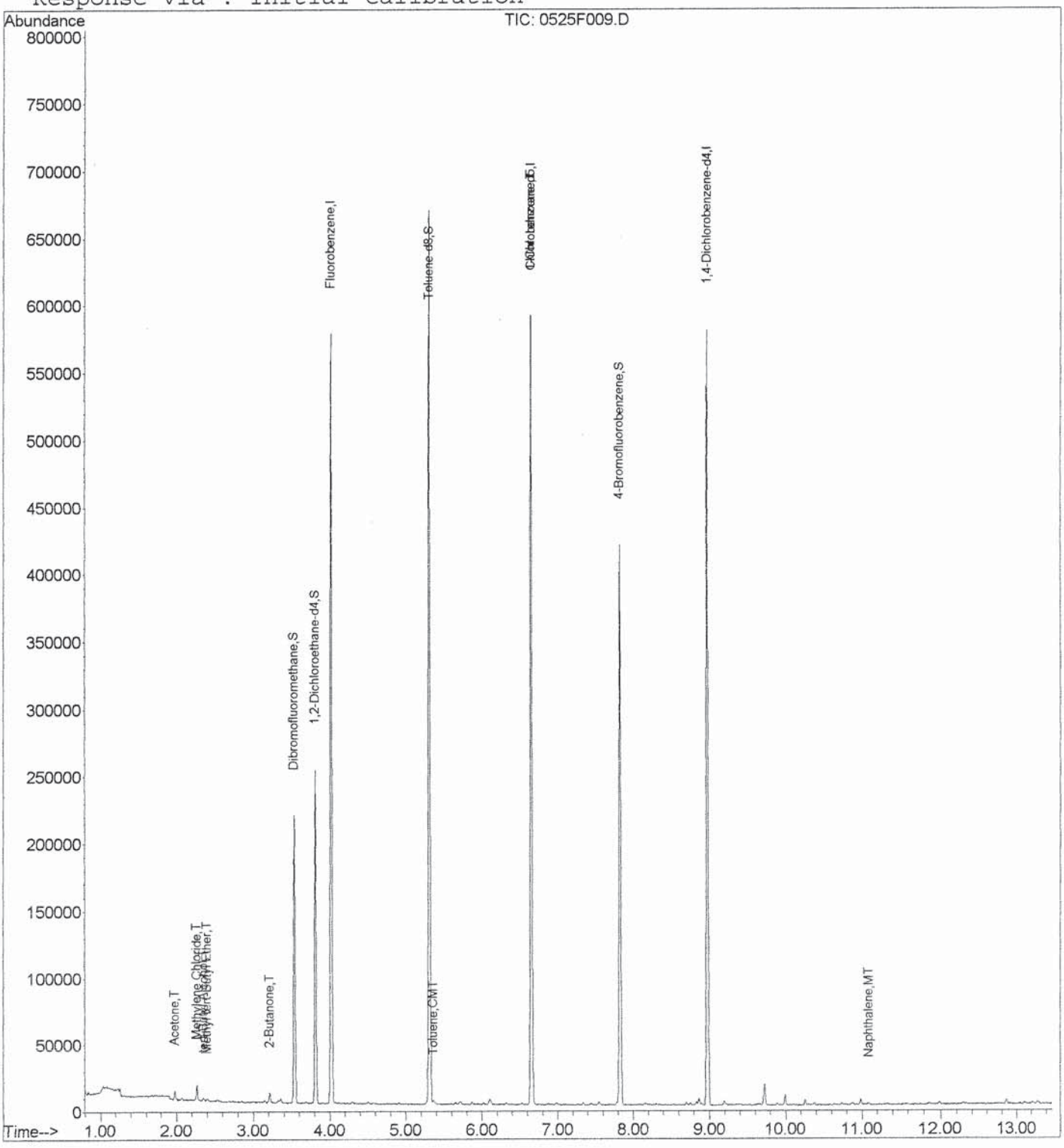
D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

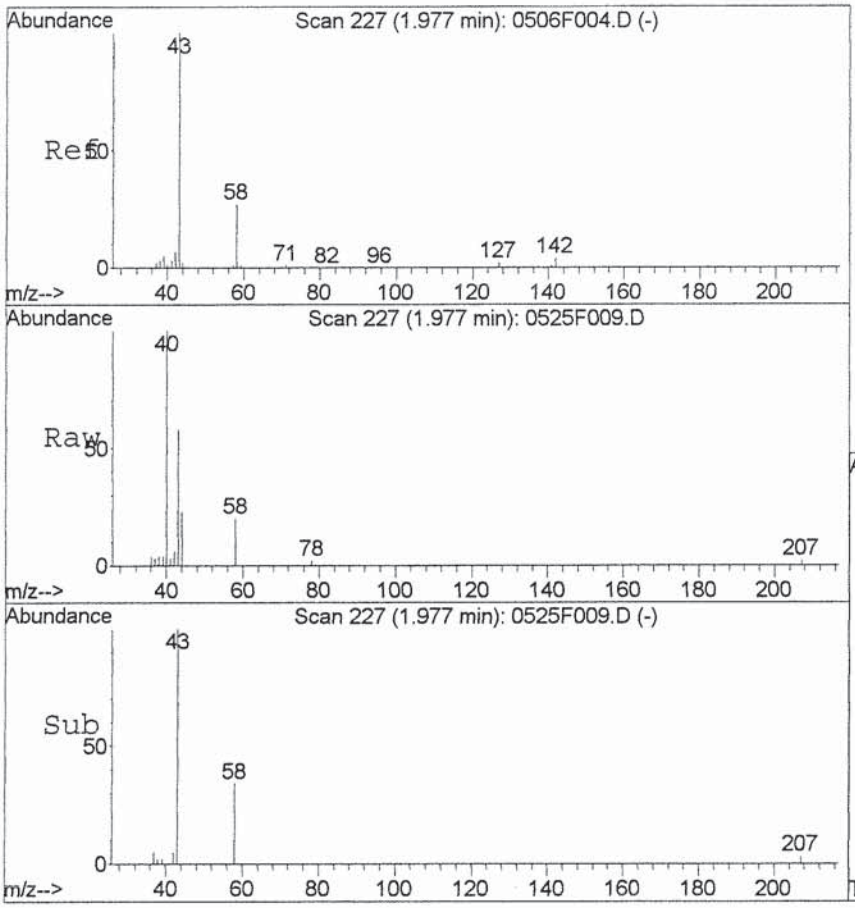
*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\052510\0525F009.D Vial: 8
Acq On : 25 May 2010 12:39 pm Operator: KR
Sample : MB Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 14:57 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Initial Calibration

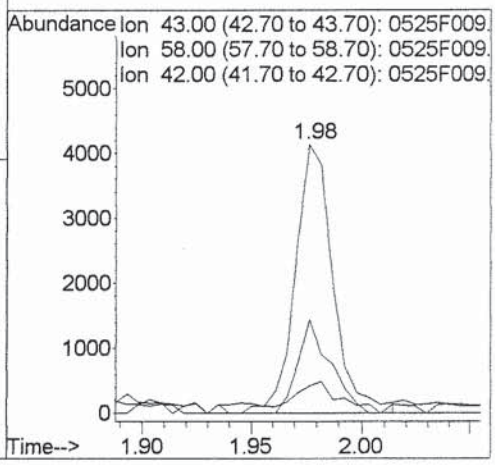


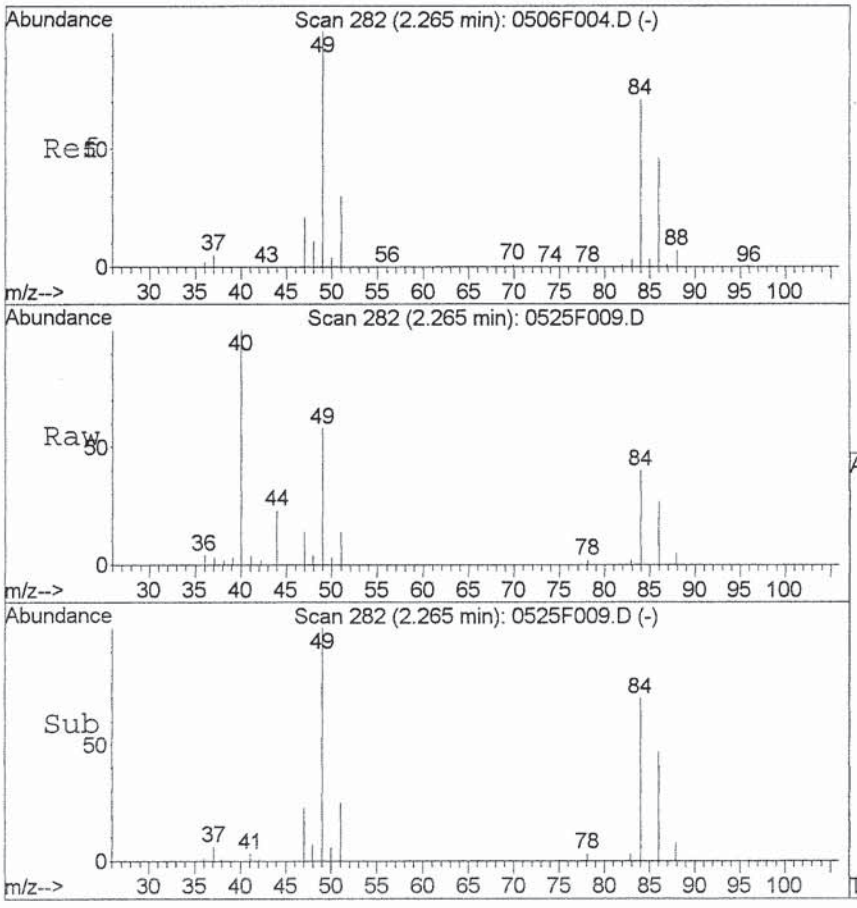


#13
Acetone
Concen: 6.25 PPB
RT: 1.98 min Scan# 227
Delta R.T. -0.00 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 43 Resp: 5072

Ion	Ratio	Lower	Upper
43	100		
58	34.7	0.0	57.0
42	10.1	0.0	37.4

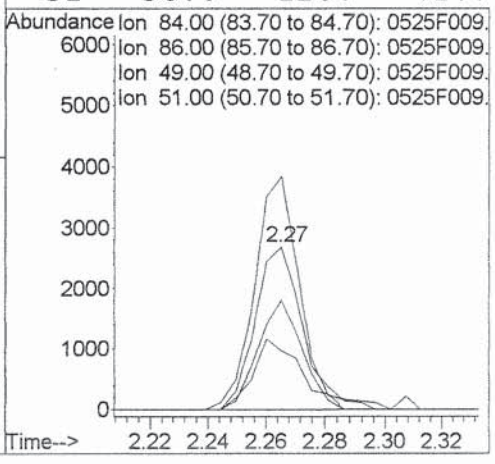


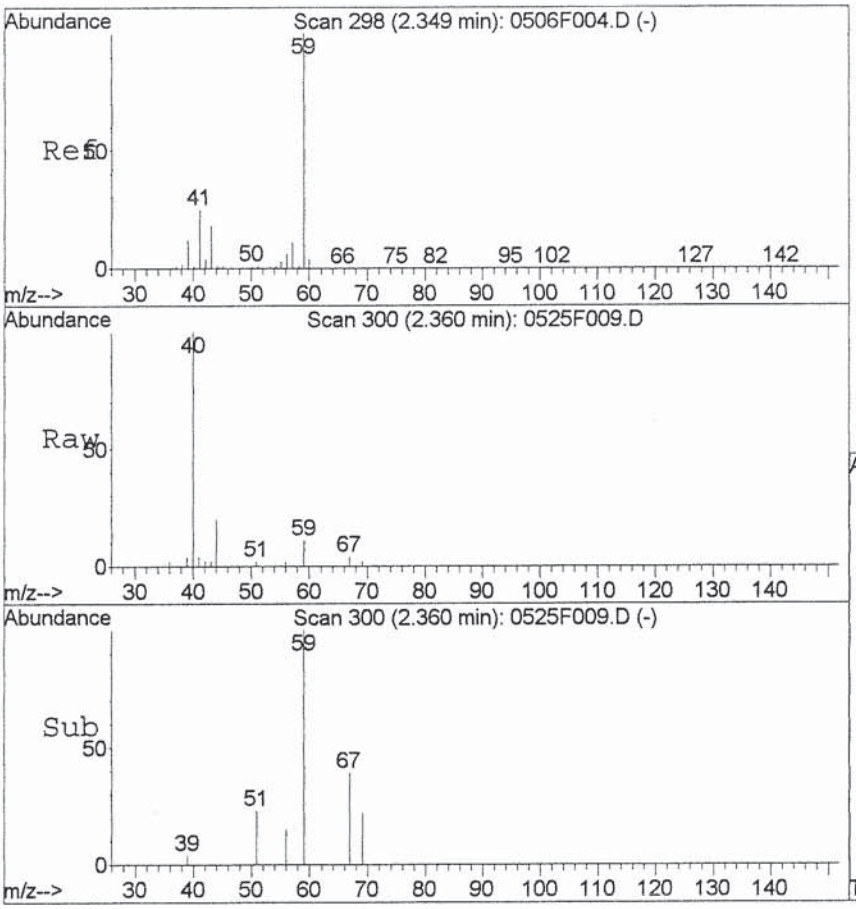


#18
Methylene Chloride
Concen: 1.47 PPB
RT: 2.27 min Scan# 282
Delta R.T. -0.00 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 84 Resp: 3082

Ion	Ratio	Lower	Upper
84	100		
86	67.3	34.4	94.4
49	143.2	109.9	169.9
51	36.0	12.7	72.7

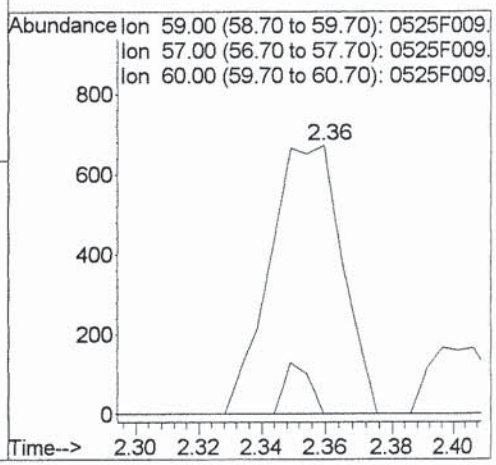


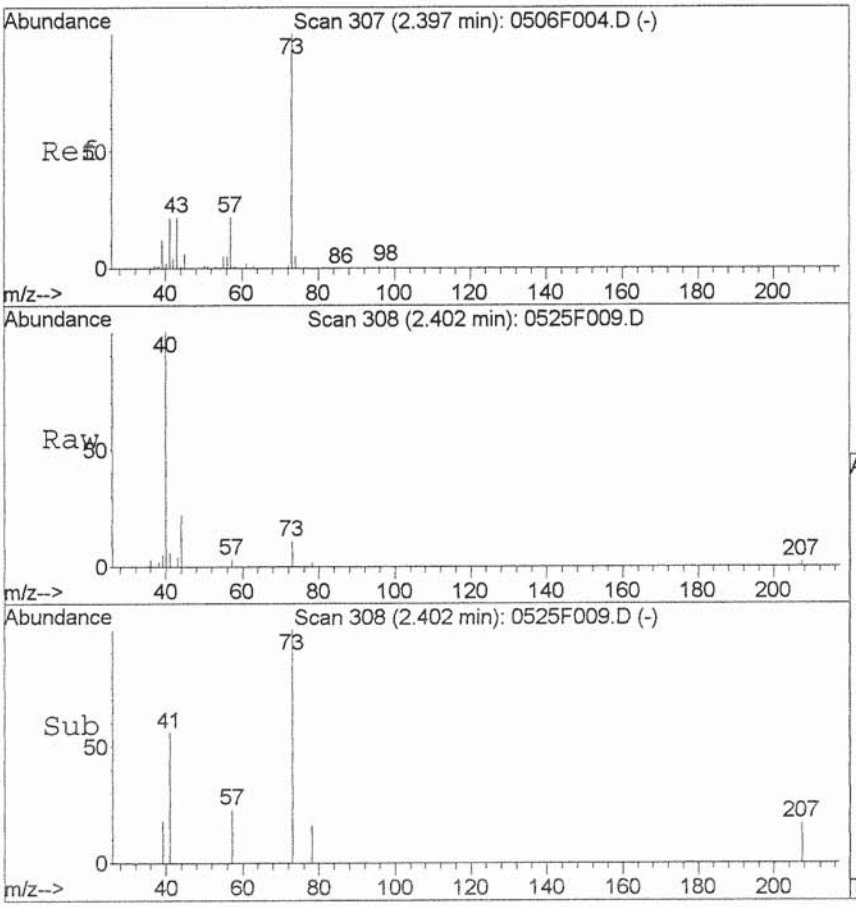


#19
tert-Butyl Alcohol
Concen: 3.60 PPB
RT: 2.36 min Scan# 300
Delta R.T. 0.01 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 59 Resp: 1048

Ion	Ratio	Lower	Upper
59	100		
57	0.0	0.0	41.1
60	0.0	0.0	33.6

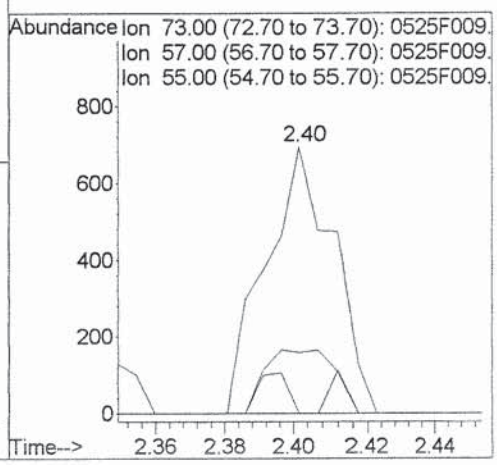


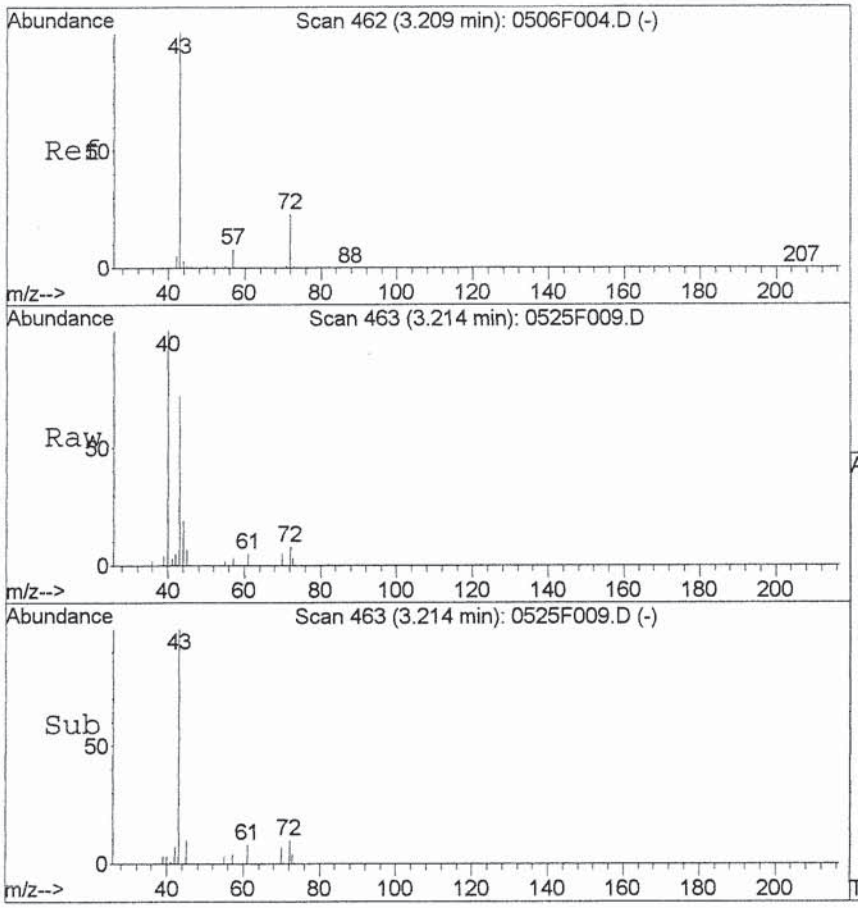


#21
Methyl tert-Butyl Ether
Concen: 0.15 PPB
RT: 2.40 min Scan# 308
Delta R.T. 0.01 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 73 Resp: 914

Ion	Ratio	Lower	Upper
73	100		
57	22.9	0.0	53.3
55	0.0	0.0	35.5

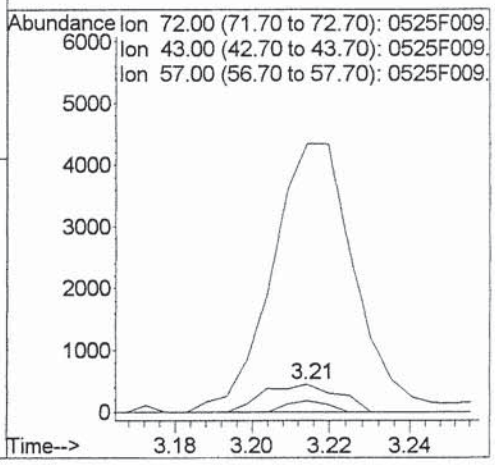


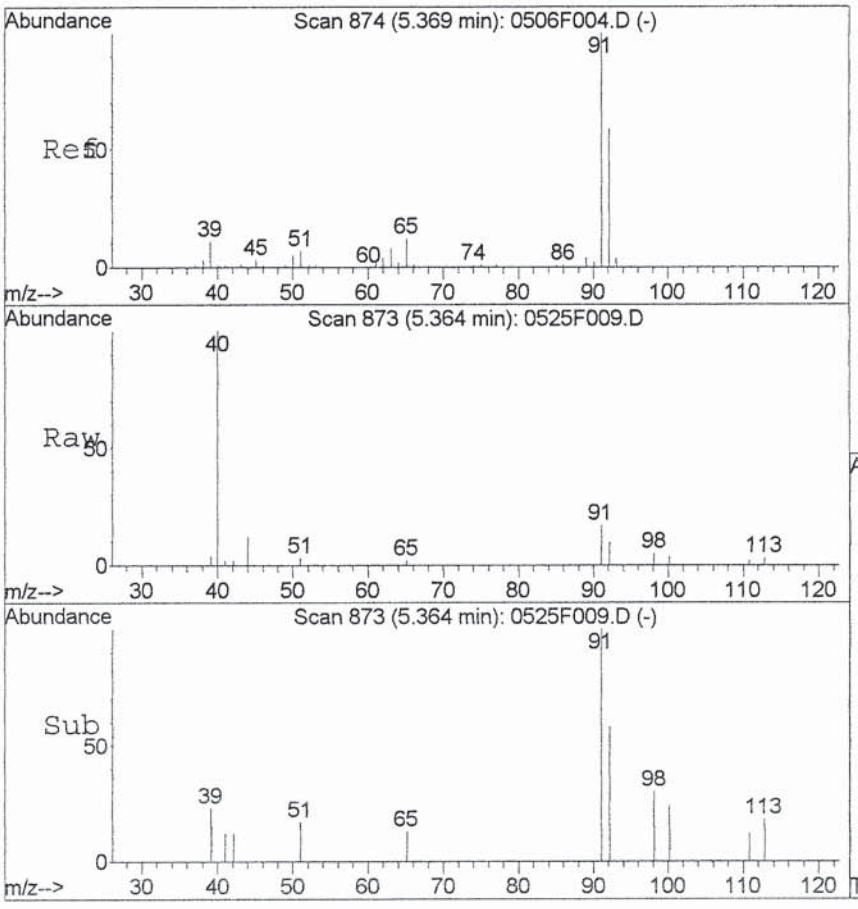


#31
2-Butanone
Concen: 2.33 PPB
RT: 3.21 min Scan# 463
Delta R.T. 0.01 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 72 Resp: 609

Ion	Ratio	Lower	Upper
72	100		
43	916.2	403.8	463.8#
57	41.2	3.3	63.3

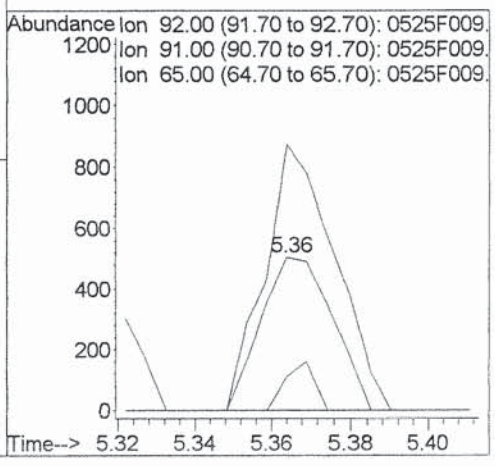


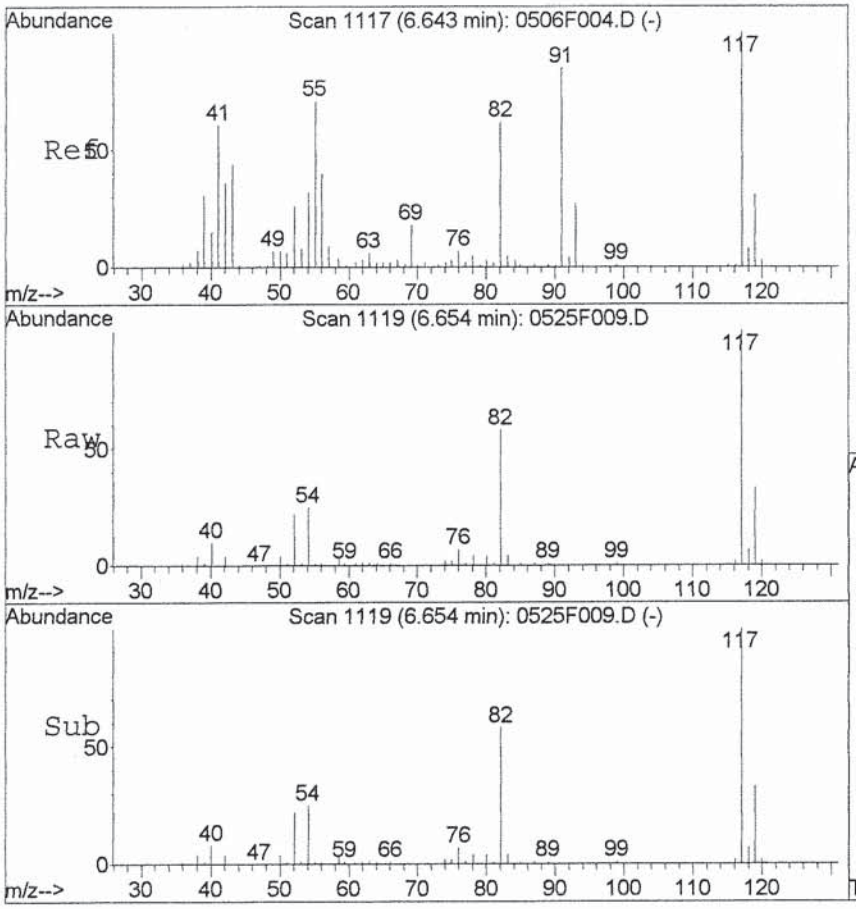


#58
Toluene
Concen: 0.14 PPB
RT: 5.36 min Scan# 873
Delta R.T. -0.01 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 92 Resp: 644

Ion	Ratio	Lower	Upper
92	100		
91	173.1	138.9	198.9
65	22.0	0.0	50.5

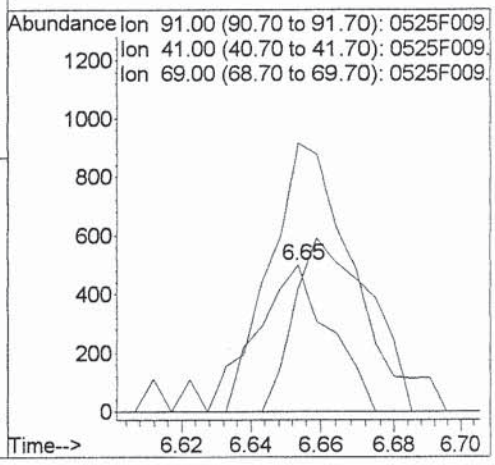


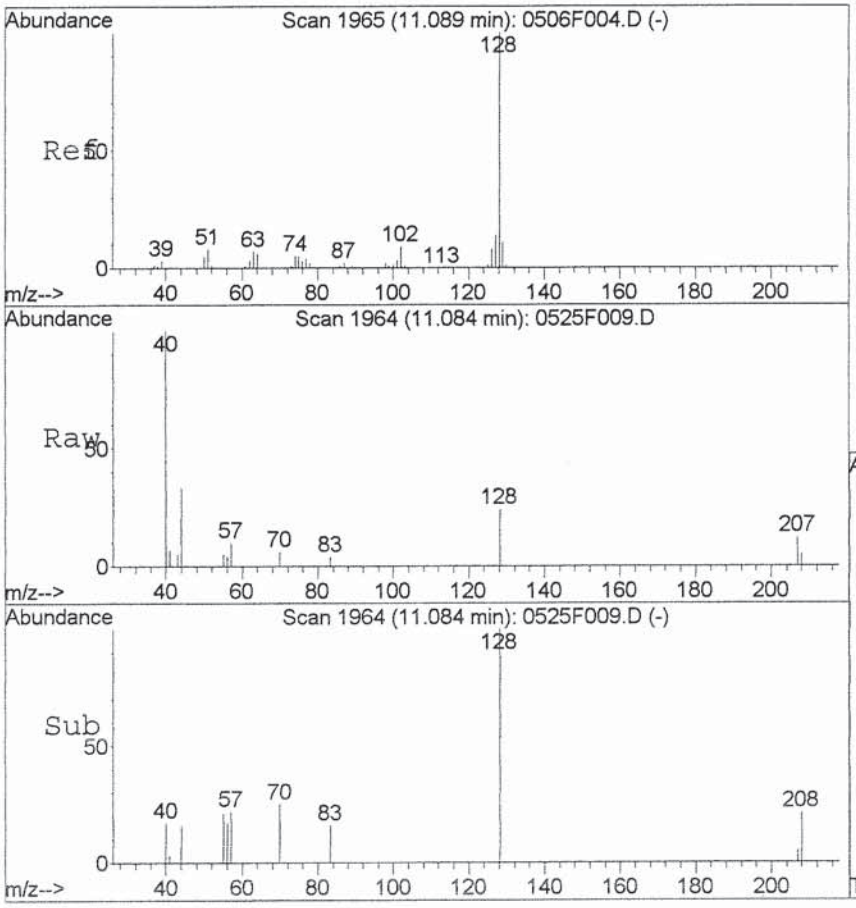


#69
1-Chlorohexane
Concen: 0.32 PPB
RT: 6.65 min Scan# 1119
Delta R.T. 0.01 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion: 91 Resp: 682

Ion	Ratio	Lower	Upper
91	100		
41	183.0	41.3	101.3#
69	84.0	0.0	51.4#

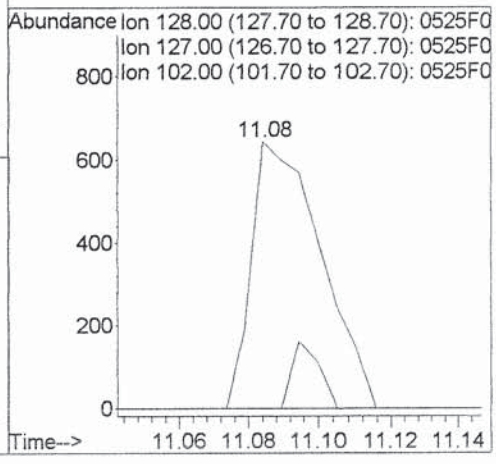




#101
Naphthalene
Concen: 0.17 PPB
RT: 11.08 min Scan# 1964
Delta R.T. -0.01 min
Lab File: 0525F009.D
Acq: 25 May 2010 12:39 pm

Tgt Ion:128 Resp: 881

Ion	Ratio	Lower	Upper
128	100		
127	0.0	0.0	43.7
102	0.0	0.0	38.7



COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16MS
Lab Code: KWG1004900-1

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 5030A
Analysis Method: 8260B

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	**48.0**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	**47.8**		5.5	0.51	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	**49.8**		5.5	0.34	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	**31.9**		5.5	0.68	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	**45.1**		5.5	0.34	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	**37.4**		5.5	0.24	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	**44.0**		5.5	0.26	1	05/25/10	05/25/10	KWG1004900	
Acetone	**297**		22	4.3	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	**98.5**		5.5	0.29	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**41.5**		11	0.48	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	**44.4**		5.5	0.39	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	**44.0**		5.5	0.22	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	**49.4**		5.5	0.40	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	**44.9**		5.5	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**412**		22	1.1	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	**46.4**		5.5	0.16	1	05/25/10	05/25/10	KWG1004900	
Chloroform	**46.5**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	**48.3**		5.5	0.39	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	**47.8**		5.5	0.40	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	**46.9**		5.5	0.30	1	05/25/10	05/25/10	KWG1004900	
Benzene	**46.7**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	**45.4**		5.5	0.17	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	**51.6**		5.5	0.30	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	**45.5**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	**45.7**		5.5	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	**46.1**		5.5	0.17	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	**41.8**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	**469**		22	0.82	1	05/25/10	05/25/10	KWG1004900	
Toluene	**45.9**		5.5	0.20	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	**57.1**		5.5	0.38	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	**48.5**		5.5	0.26	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	**47.7**		5.5	0.29	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	**527**		22	1.1	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	**49.2**		5.5	0.22	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16MS
Lab Code: KWG1004900-1

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 5030A
Analysis Method: 8260B

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	**50.7**		5.5	0.21	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	**52.2**		22	0.30	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	**44.0**		5.5	0.28	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	**46.4**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	**47.9**		5.5	0.28	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	**96.3**		5.5	0.41	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	**45.9**		5.5	0.15	1	05/25/10	05/25/10	KWG1004900	
Styrene	**45.8**		5.5	0.15	1	05/25/10	05/25/10	KWG1004900	
Bromoform	**59.9**		5.5	0.33	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	**49.1**		22	0.15	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	**54.6**		5.5	0.40	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	**50.4**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	**52.9**		22	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	**71.0**		5.5	0.42	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	**52.7**		22	0.32	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	**50.3**		22	0.40	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	**52.3**		22	0.29	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	**49.5**		22	0.37	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	**46.7**		22	0.31	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	**44.4**		22	0.34	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	**44.0**		22	0.31	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	**47.0**		5.5	0.27	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	**46.2**		5.5	0.28	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	**41.8**		22	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	**45.3**		5.5	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	**91.0**		22	0.73	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	**38.9**		22	0.26	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	**27.9**		22	0.24	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	**44.4**		22	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	**40.0**		22	0.24	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Collected:**	05/19/2010
Sample Matrix:	Misc. solid	**Date Received:**	05/21/2010

Volatile Organic Compounds

Sample Name:	D-4-16MS	**Units:**	ug/Kg
Lab Code:	KWG1004900-1	**Basis:**	Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	99	83-128	05/25/10	Acceptable
Toluene-d8	109	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	108	77-124	05/25/10	Acceptable

Comments:

Exception Report

Data File: J:\MS24\DATA\052510\0525F013.D
Lab ID: KWG1004900-1 -- K1005244-003MS
RunType: MS
Matrix: MISC. SOLID

Date Acquired: 05/25/2010 14:40
Date Quantitated: 05/25/2010 15:26
Batch ID: KWG1004880
Analysis Method: 8260B
MethodJoinID: MJ120

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Window	NA	NA	NA	x	
Analytical Holding Time	NA	NA	NA	x	
ICAL Pass/Fail	NA	NA	NA	x	
ICAL Average RSD	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Initial Calibration Minimum RF	NA	NA	NA		x
Initial Calibration SPCC/CCC	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Continuing Calibration Minimum RF	NA	NA	NA		x
Continuing Calibration SPCC/CCC	NA	NA	NA	x	
Internal Standards	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Relative Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Analyte Exceptions

Exception Categories	Analyte Name	Result	Low Limit	High Limit	Corrective Action
Initial Calibration Minimum RF	1,4-Dioxane	0.0037	0.01	NA	NC
Continuing Calibration Minimum RF	1,4-Dioxane	0.0034	0.01	NA	J

Primary Review: KL 5/25/10

Secondary Review: Comm 5/25/10

Printed: 05/25/2010 15:34:17
u:\Stealth\Crystal.rpt\except2.rpt

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	MISC. SOLID
Prod Code:	8260B VOC_FP	**Collect Date:**		**Receive Date:**	05/25/2010

Analysis Lot:	KWG1004880	**Prep Lot:**	KWG1004900	**Report Group:**	
Analysis Method:	8260B	**Prep Method:**	EPA 5030A		
Prep Ref:	911428	**Prep Date:**	05/25/2010		

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:			
Tune Ref:	J:\MS24\DATA\052510\0525F002.D	**Method ID:**	MJ120
MB Ref:	J:\MS24\DATA\052510\0525F009.D	**Quant based on Method**	

Data File:	J:\MS24\DATA\052510\0525F013.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 14:40	**Quant Date:**	05/25/2010 15:26	**Vial:**	12
Run Type:	MS			**Dilution:**	1.0
Lab ID:	KWG1004900-1 -- K1005244-003MS			**Soln Conc. Units:**	PPB

Internal Standard Compounds

IS Ref	Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Area Criteria
1	Fluorobenzene	4.02	0.00	96	420529	50.00	OK
2	Chlorobenzene-d5	6.65	0.00	82	177982	50.00	OK
3	1,4-Dichlorobenzene-d4	8.98	0.00	152	145052	50.00	OK

Surrogate Compounds

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	%Rec	%Rec Limits		Rpt?
1	Dibromofluoromethane	3.53	-0.01	0.00	113	103877	49.42	99	83-128	OK	
1	1,2-Dichloroethane-d4	3.81	0.00	0.00	65	150773	49.69	99	64-142	OK	
1	Toluene-d8	5.31	0.00	0.00	98	400179	54.28	109	90-125	OK	
2	4-Bromofluorobenzene	7.83	0.00	0.00	95	147981	54.17	108	77-124	OK	

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Dichlorodifluoromethane	0.85	-0.01	0.00	85	113599	43.76	48.0		
1	Chloromethane	0.98	-0.01	0.00	50	143441	43.58	47.8		
1	Vinyl Chloride	1.05	-0.01	0.00	62	132654	45.43	49.8		
1	Bromomethane	1.28	-0.02	0.00	96	70612**m**	29.07	31.9		
1	Chloroethane	1.35	-0.02	0.00	64	78331	41.16	45.1		
1	Dichlorofluoromethane (CFC 21	1.49	-0.03	-0.01	67	501	0.1000	0.110	J	
1	Trichlorofluoromethane	1.50	-0.01	0.00	101	137442	34.14	37.4		
1	Diethyl Ether	1.71	-0.02	0.00	59	95014	39.45	43.2		
1	Acrolein	1.86	-0.01	0.00	56	33362	71.74	78.6	J	
1	Trichlorotrifluoroethane	1.85	-0.01	0.00	151	64690	42.03	46.1		
1	1,1-Dichloroethene	1.87	-0.01	0.00	96	70955	40.18	44.0		
1	Acetone	1.97	-0.01	0.00	43	265777	271.44	297		
1	Iodomethane (Methyl Iodide)	1.99	-0.01	0.00	142	168877	70.02	76.7		
1	Carbon Disulfide	2.00	-0.01	0.00	76	623432	89.85	98.5		

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F013.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 14:40	**Quant Date:**	05/25/2010 15:26	**Vial:**	12
Run Type:	MS			**Dilution:**	1.0
Lab ID:	KWG1004900-1 -- K1005244-003MS			**Soln Conc. Units:**	PPB

Target Compounds **Final Conc. Units:** ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	3-Chloro-1-propene	2.15	-0.01	0.00	76	85567	67.75	74.3		
1	Acetonitrile	2.23		0.00	40	153139	796.45	873		
1	Methylene Chloride	2.26	-0.01	0.00	84	95864	37.87	41.5		
1	tert-Butyl Alcohol	2.35		0.00	59	357066	1,017	1110		
1	Acrylonitrile	2.50		0.00	53	140851	143.93	158		
1	Methyl tert-Butyl Ether	2.39	-0.01	0.00	73	340655	46.03	50.4		
1	trans-1,2-Dichloroethene	2.41	-0.01	0.00	96	89547	40.54	44.4		
1	n-Hexane	2.53	-0.01	0.00	57	187046	61.79	67.7		
1	Diisopropyl Ether	2.73	-0.01	0.00	45	673974	81.00	88.8		
1	1,1-Dichloroethane	2.74	-0.01	0.00	63	175003	40.17	44.0		
1	Vinyl Acetate	2.78		0.00	86	95632	235.87	259		
1	Chloroprene	2.77	-0.01	0.00	53	252214	85.10	93.3		
1	tert-Butyl Ethyl Ether	2.99		0.00	59	653179	86.47	94.8		
1	2,2-Dichloropropane	3.14	-0.01	0.00	77	134518	45.04	49.4		
1	cis-1,2-Dichloroethene	3.17	-0.01	0.00	96	100350	40.95	44.9		
1	2-Butanone (MEK)	3.20	-0.01	0.00	72	118259	375.48	412		
1	Propionitrile	3.32		0.00	54	35726	91.60	100		
1	Methacrylonitrile	3.40		0.00	67	106414	102.74	113		
1	Bromochloromethane	3.36		0.00	128	50816	42.30	46.4		
1	Chloroform	3.40	-0.01	0.00	83	174260	42.39	46.5		
1	1,1,1-Trichloroethane (TCA)	3.50	-0.01	0.00	97	138429	44.09	48.3		
1	Carbon Tetrachloride	3.59		0.00	117	117874	43.66	47.8		
1	1,1-Dichloropropene	3.62	-0.01	0.00	75	120710	42.79	46.9		
1	Isobutanol	3.78		0.00	43	218735	1,701	1860		
1	Benzene	3.78	-0.01	0.00	78	381793	42.61	46.7		
1	1,2-Dichloroethane (EDC)	3.87		0.00	62	150521	41.41	45.4		
1	tert-Amyl Methyl Ether	3.84		0.00	55	173409	89.33	97.9		
1	Trichloroethene (TCE)	4.29		0.00	95	111416	47.04	51.6		
1	1,2-Dichloropropane	4.52	-0.01	0.00	63	99053	41.51	45.5		
1	Dibromomethane	4.62		0.00	93	62942	41.74	45.7		
1	Methyl Methacrylate	4.60		0.00	69	167608	100.35	110		
1	1,4-Dioxane	4.61	-0.01	0.00	88	37265	1,208	1320		
1	Bromodichloromethane	4.75		0.00	83	126654	42.10	46.1		
1	2-Nitropropane	5.02		0.00	41	65143	104.20	114		
1	2-Chloroethyl Vinyl Ether	5.02		0.00	63	87931	59.61	65.3		
1	cis-1,3-Dichloropropene	5.13		0.00	75	128164	38.16	41.8		
1	4-Methyl-2-pentanone (MIBK)	5.26	-0.01	0.00	58	392298	428.08	469		
1	Toluene	5.36	-0.01	0.00	92	225224	41.85	45.9		
2	trans-1,3-Dichloropropene	5.65		0.00	75	153524	52.11	57.1		
2	Ethyl Methacrylate	5.67		0.00	69	264051	94.24	103		
2	1,1,2-Trichloroethane	5.81		0.00	83	81501	44.21	48.5		
2	Tetrachloroethene (PCE)	5.82		0.00	164	76033	43.55	47.7		

U: Undetected at or above MDL
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m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F013.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 14:40	**Quant Date:**	05/25/2010 15:26	**Vial:**	12
Run Type:	MS			**Dilution:**	1.0
Lab ID:	KWG1004900-1 -- K1005244-003MS			**Soln Conc. Units:**	PPB

Target Compounds — **Final Conc. Units:** ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2	2-Hexanone	6.02		0.00	57	149611	480.75	527		
2	1,3-Dichloropropane	5.97		0.00	76	164795	44.88	49.2		
2	Dibromochloromethane	6.13		0.00	129	95451	46.30	50.7		
2	1,2-Dibromoethane (EDB)	6.25		0.00	107	95897	47.62	52.2		
2	1-Chlorohexane	6.64		0.00	91	117661	46.49	51.0		
2	Chlorobenzene	6.68		0.00	112	246909	40.18	44.0		
2	Ethylbenzene	6.76		0.00	106	126782	42.37	46.4		
2	1,1,1,2-Tetrachloroethane	6.77		0.00	131	88617	43.71	47.9		
2	m,p-Xylenes	6.87		0.00	106	309984	87.91	96.3		
2	o-Xylene	7.27		0.00	106	148170	41.85	45.9		
2	Styrene	7.30		0.00	103	116134**m**	41.83	45.8		
2	Bromoform	7.51		0.00	173	68377	54.61	59.9		
2	Isopropylbenzene	7.62		0.00	105	382936	44.78	49.1		
2	cis-1,4-Dichloro-2-butene	7.80		0.00	89	35903	148.92	163		
3	1,1,2,2-Tetrachloroethane	8.03		0.00	83	105216	49.82	54.6		
3	trans-1,4-Dichloro-2-butene	8.10		0.00	53	90172	180.36	198		
3	Bromobenzene	7.96		0.00	156	104730	45.98	50.4		
3	n-Propylbenzene	8.04		0.00	91	448455	48.26	52.9		
3	1,2,3-Trichloropropane	8.07		0.00	110	51982	64.78	71.0		
3	2-Chlorotoluene	8.15		0.00	91	279905	48.05	52.7		
3	1,3,5-Trimethylbenzene	8.23		0.00	105	299447	45.86	50.3		
3	4-Chlorotoluene	8.27		0.00	91	280512	47.69	52.3		
3	tert-Butylbenzene	8.54		0.00	119	258293	45.16	49.5		
3	1,2,4-Trimethylbenzene	8.60		0.00	105	284264	42.57	46.7		
3	sec-Butylbenzene	8.76		0.00	105	345304	40.52	44.4		
3	4-Isopropyltoluene	8.91		0.00	119	285646	40.17	44.0		
3	1,3-Dichlorobenzene	8.90		0.00	146	180030	42.87	47.0		
3	1,4-Dichlorobenzene	9.00		0.00	146	183336	42.15	46.2		
3	n-Butylbenzene	9.31		0.00	91	239372	38.11	41.8		
3	1,2-Dichlorobenzene	9.36		0.00	146	170709	41.35	45.3		
3	1,2-Dibromo-3-chloropropane	10.14		0.00	155	26206	83.06	91.0		
3	1,2,4-Trichlorobenzene	10.86		0.00	180	92691	35.47	38.9		
3	Hexachlorobutadiene	10.96		0.00	225	33262	25.44	27.9		
3	Naphthalene	11.09		0.00	128	243693	40.55	44.4		
3	1,2,3-Trichlorobenzene	11.32		0.00	180	88832	36.50	40.0		
	4-Vinyl-1-cyclohexene				0	0		5.5	UJ	NR
	Benzyl Chloride				0	0		5.5	UJ	NR
	Hexachloroethane				0	0		1.1	UJ	NR
	2,2,4-Trimethylpentane				0	0		5.5	UJ	NR
	1,3-Butadiene				0	0		2.2	UJ	NR
	1,1,2-Trifluoroethane				0	0		11	UJ	NR
	n-Butyl Alcohol				0	0		270	U	NR

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F013.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 14:40	**Quant Date:**	05/25/2010 15:26	**Vial:**	12
Run Type:	MS			**Dilution:**	1.0
Lab ID:	KWG1004900-1 -- K1005244-003MS			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2-Propanol			0	0		220	U	NR
Ethyl Alcohol			0	0		220	U	NR
2,2-Dichloro-1,1,1-trifluoroetha			0	0		5.5	UJ	NR
Cyclohexanone			0	0		11	UJ	NR
Methylcyclopentane			0	0		5.5	UJ	NR
n-Heptane			0	0		5.5	UJ	NR
3-Methylpentane			0	0		5.5	UJ	NR
2-Methylpentane			0	0		5.5	UJ	NR
Ethyl Acetate			0	0		22	UJ	NR

Prep Amount:	5.03 g	**Dilution:**	1.0
Prep Final Vol:	5.0 ml	**Unit Factor:**	1
Solids:	90.7 %		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

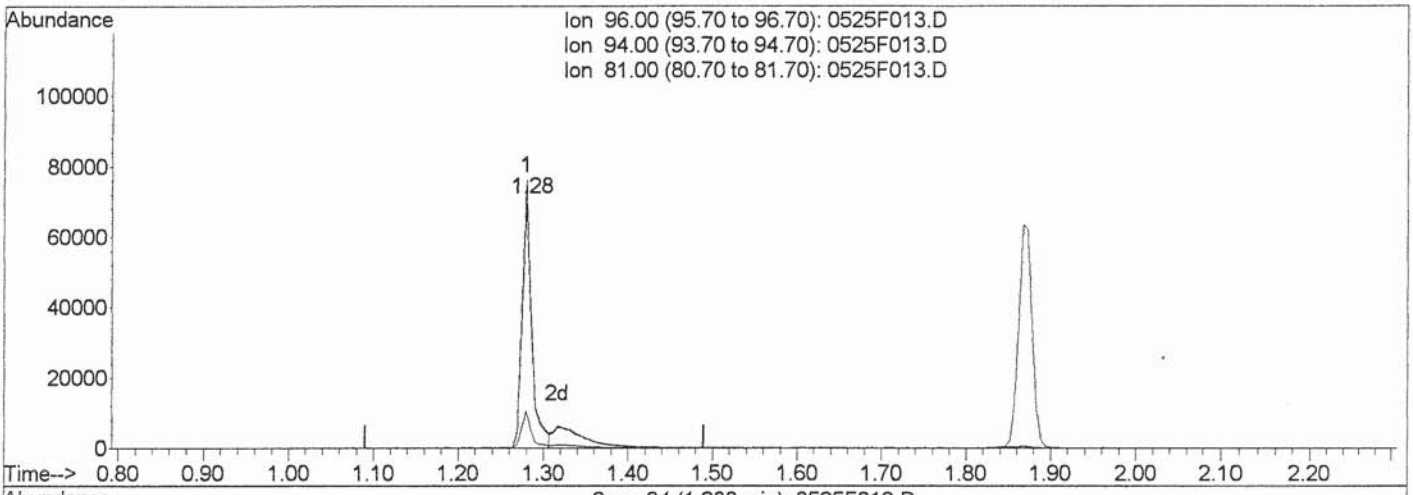
D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

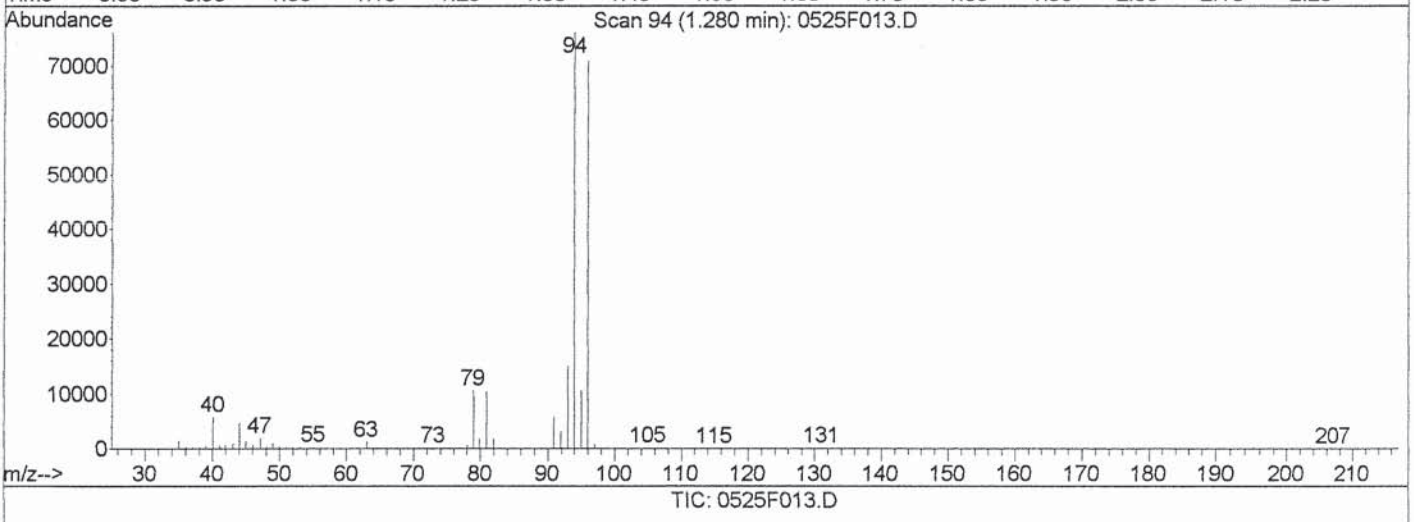
*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F013.D Vial: 12
Acq On : 25 May 2010 2:40 pm Operator: KR
Sample : K5244-003MS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:26 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F013.D

(5) Bromomethane (T)

1.28min 22.59PPB

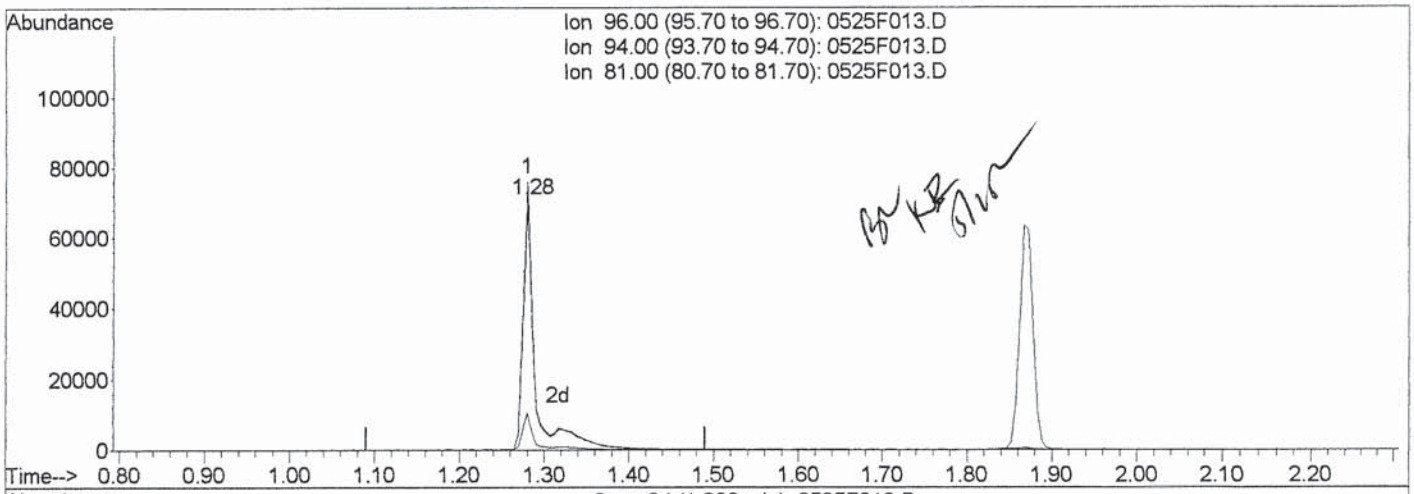
response 54864

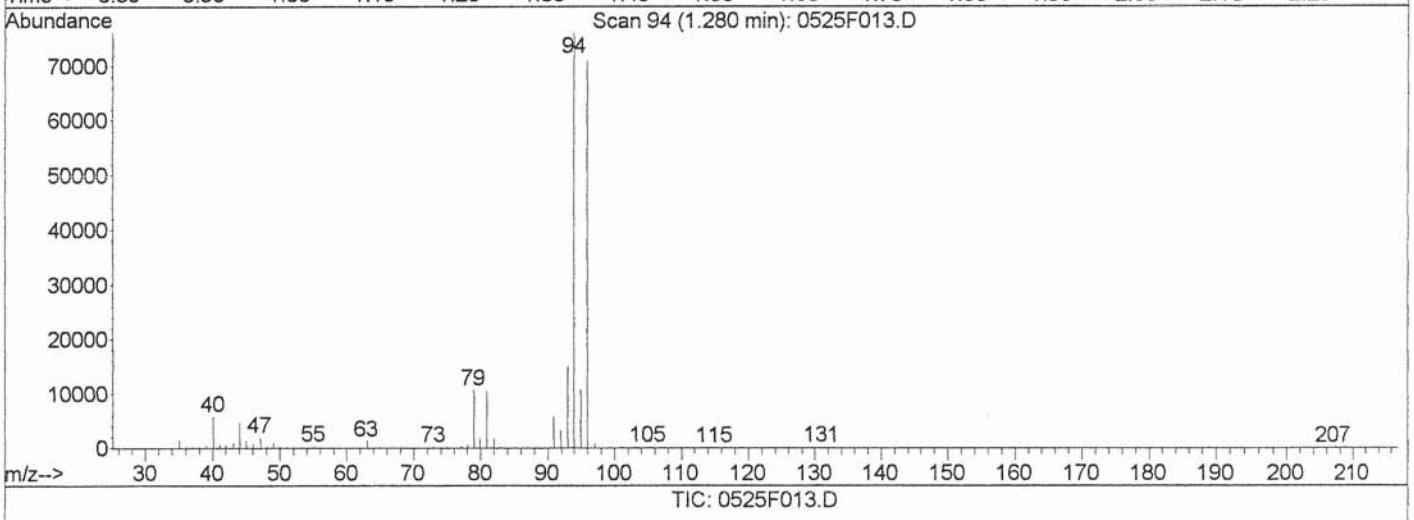
Ion	Exp%	Act%
96.00	100	100
94.00	106.40	107.24
81.00	18.10	14.55
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F013.D Vial: 12
Acq On : 25 May 2010 2:40 pm Operator: KR
Sample : K5244-003MS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:26 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F013.D

(5) Bromomethane (T)

1.28min 29.07PPB m

response 70612

Ion	Exp%	Act%
96.00	100	100
94.00	106.40	107.24
81.00	18.10	14.78
0.00	0.00	0.00

5/25/10

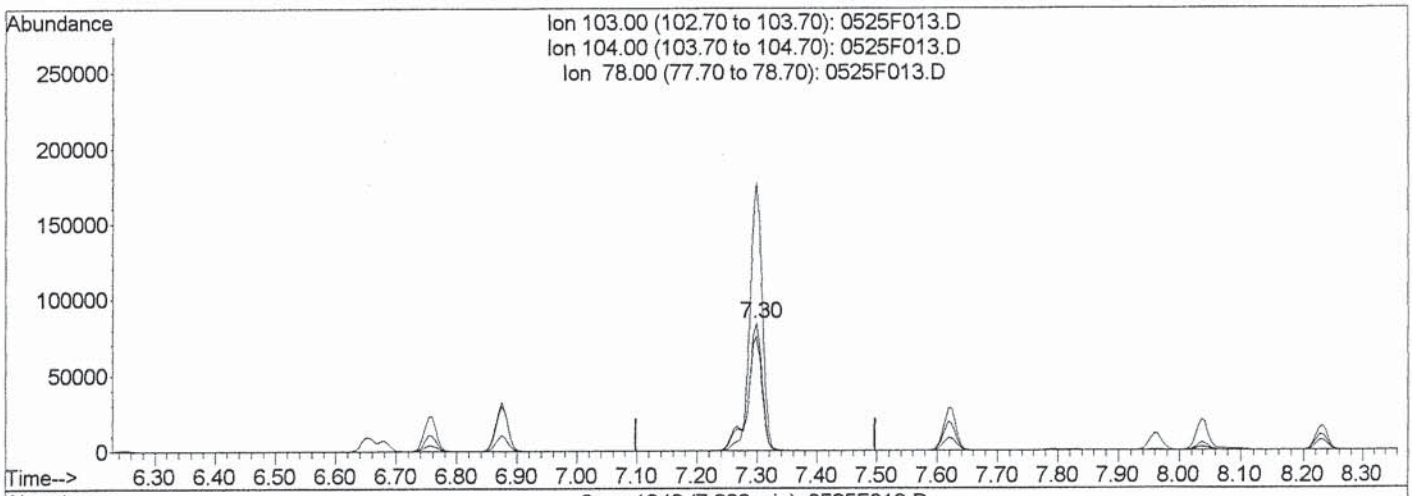
0525F013.D 042110MS24SOIL.M Tue May 25 15:26:40 2010

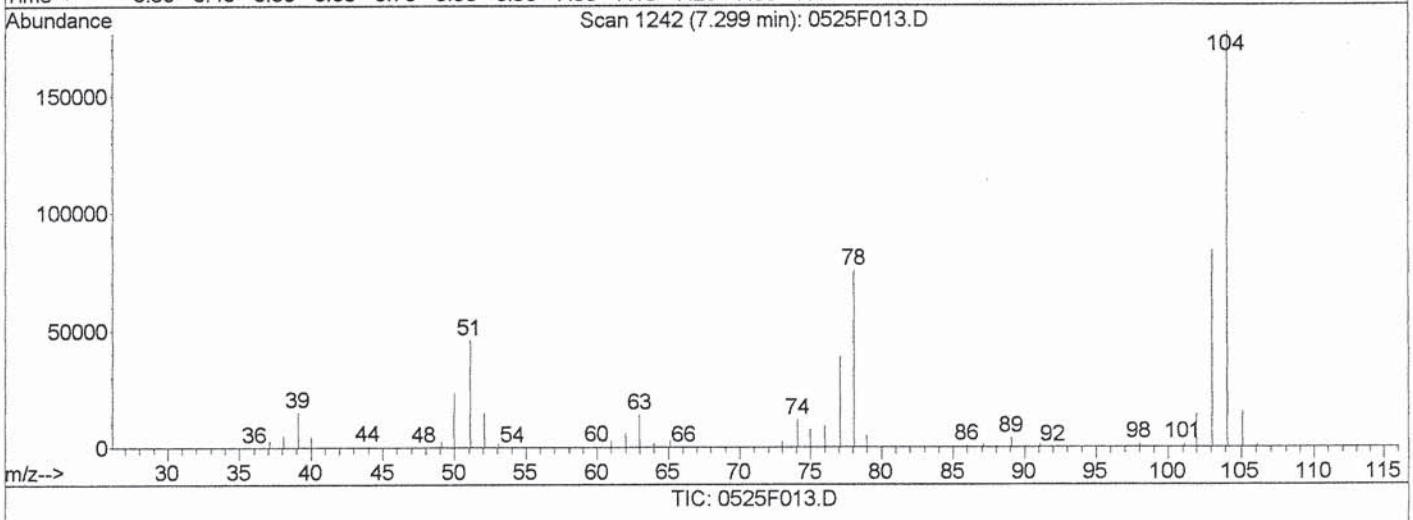
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F013.D
Acq On : 25 May 2010 2:40 pm
Sample : K5244-003MS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 15:26 2010

Vial: 12
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F013.D

(75) Styrene (T)

7.30min 49.48PPB

response 137379

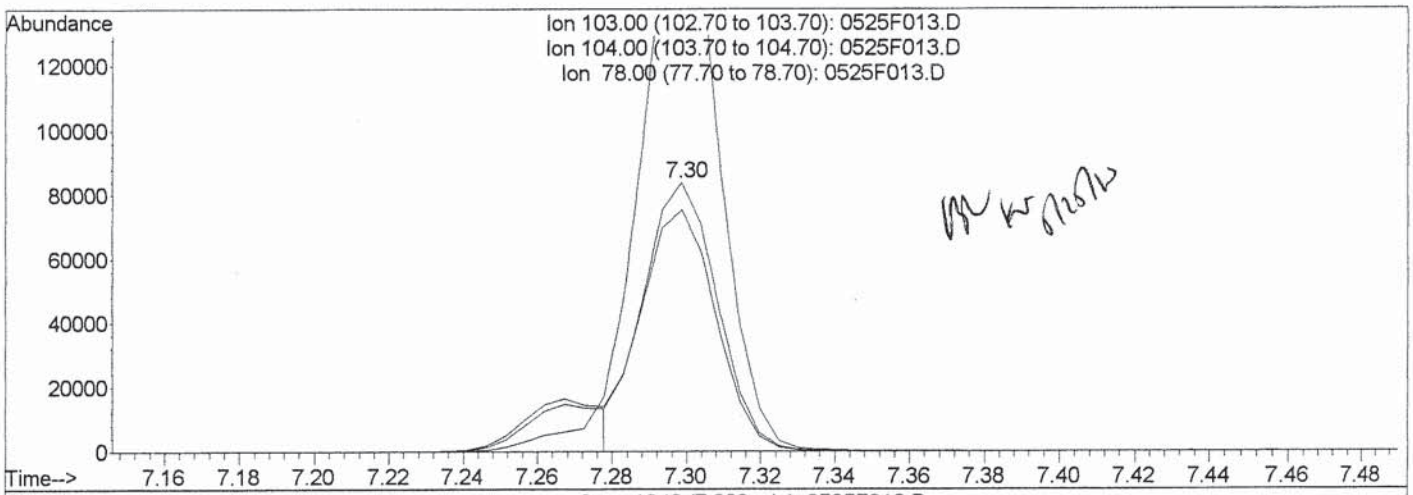
Ion	Exp%	Act%
103.00	100	100
104.00	208.80	211.36
78.00	88.30	89.86
0.00	0.00	0.00

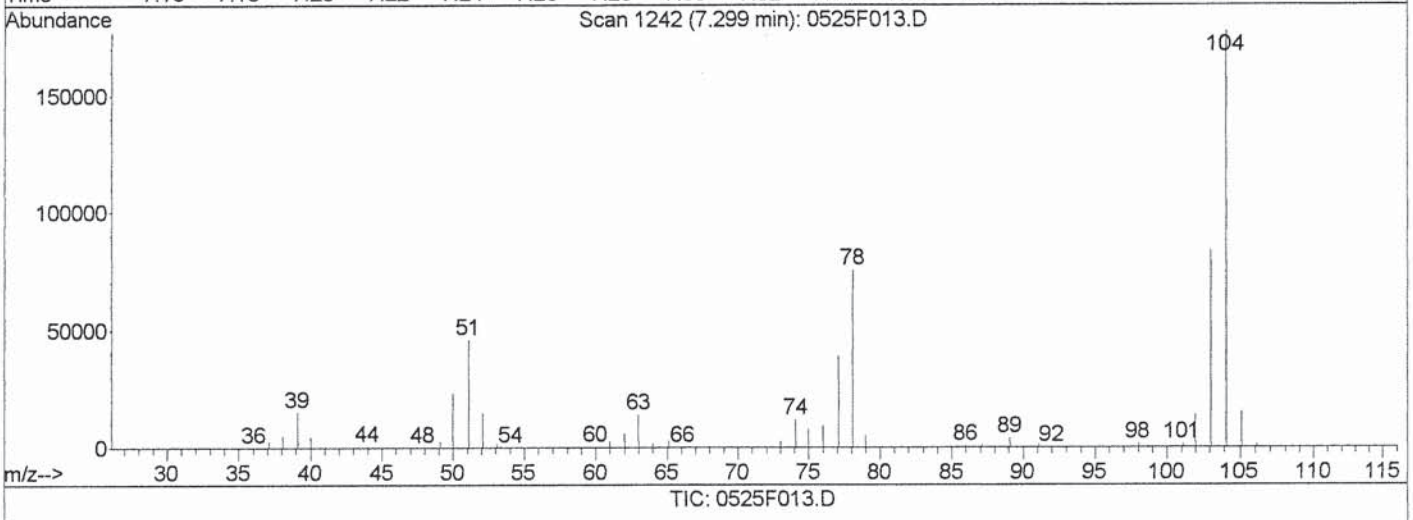
0525F013.D 042110MS24SOIL.M Tue May 25 15:26:48 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F013.D Vial: 12
Acq On : 25 May 2010 2:40 pm Operator: KR
Sample : K5244-003MS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:26 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F013.D

(75) Styrene (T)

7.30min 41.83PPB m

response 116134

Ion	Exp%	Act%
103.00	100	100
104.00	208.80	211.36
78.00	88.30	89.86
0.00	0.00	0.00

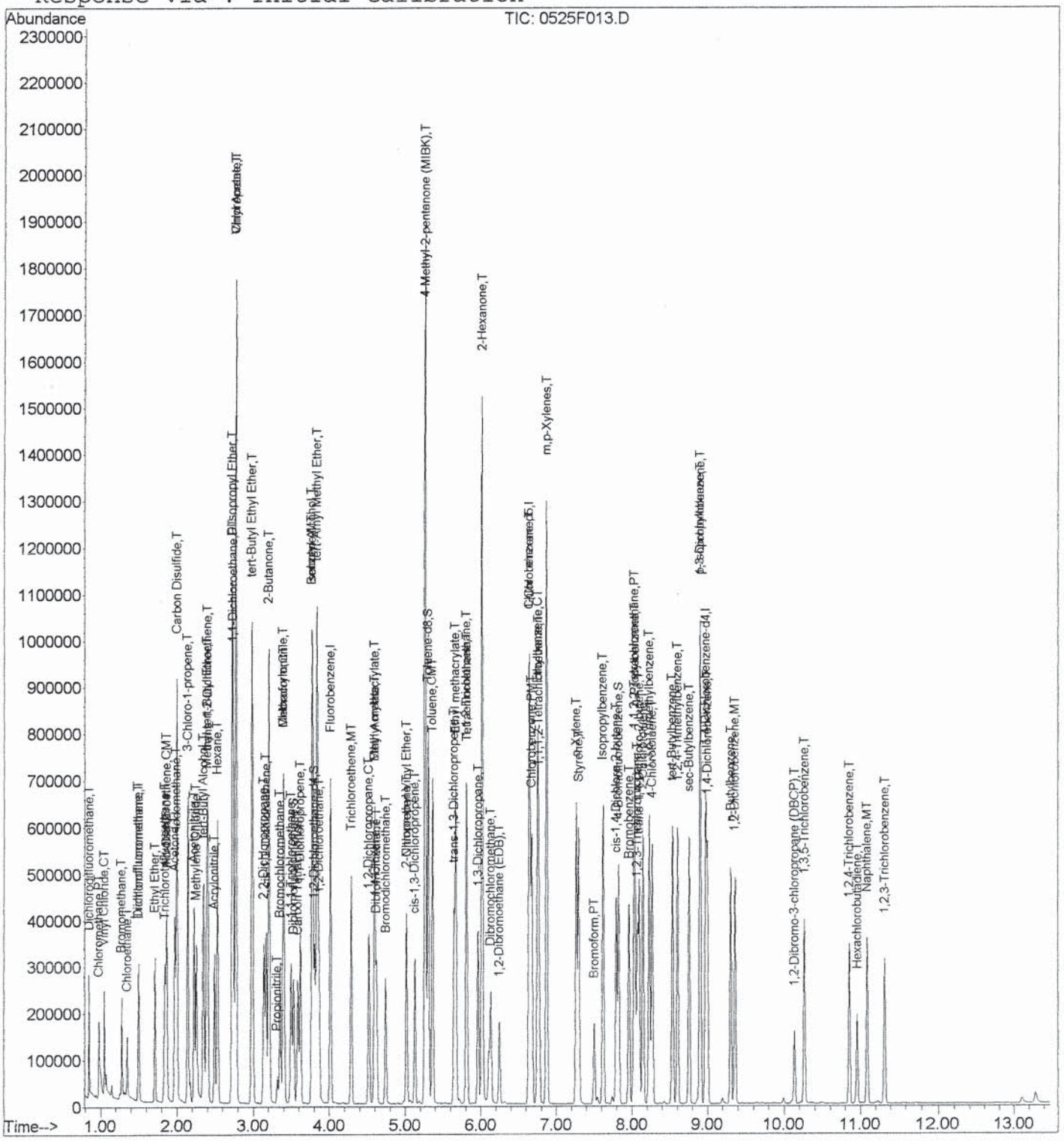
Lcma 5/25/10

0525F013.D 042110MS24SOIL.M Tue May 25 15:26:56 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\052510\0525F013.D Vial: 12
Acq On : 25 May 2010 2:40 pm Operator: KR
Sample : K5244-003MS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:26 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Initial Calibration



Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16DMS
Lab Code: KWG1004900-2

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 5030A
Analysis Method: 8260B

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	**51.5**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	**52.7**		5.6	0.51	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	**55.8**		5.6	0.35	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	**34.2**		5.6	0.69	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	**48.2**		5.6	0.35	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	**41.5**		5.6	0.24	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	**49.5**		5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
Acetone	**331**		23	4.3	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	**108**		5.6	0.29	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**47.9**		12	0.48	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	**49.6**		5.6	0.39	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	**50.0**		5.6	0.23	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	**53.0**		5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	**47.2**		5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**428**		23	1.1	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	**49.7**		5.6	0.16	1	05/25/10	05/25/10	KWG1004900	
Chloroform	**49.8**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	**52.4**		5.6	0.39	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	**51.3**		5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	**50.0**		5.6	0.30	1	05/25/10	05/25/10	KWG1004900	
Benzene	**49.3**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	**50.1**		5.6	0.17	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	**52.9**		5.6	0.30	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	**51.3**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	**51.7**		5.6	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	**52.1**		5.6	0.17	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	**46.8**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	**507**		23	0.82	1	05/25/10	05/25/10	KWG1004900	
Toluene	**50.7**		5.6	0.20	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	**60.8**		5.6	0.38	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	**51.7**		5.6	0.26	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	**48.9**		5.6	0.29	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	**532**		23	1.1	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	**52.4**		5.6	0.23	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16DMS
Lab Code: KWG1004900-2

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 5030A
Analysis Method: 8260B

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	**54.2**		5.6	0.21	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	**55.1**		23	0.30	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	**45.7**		5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	**48.5**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	**50.4**		5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	**99.5**		5.6	0.41	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	**46.8**		5.6	0.15	1	05/25/10	05/25/10	KWG1004900	
Styrene	**46.8**		5.6	0.15	1	05/25/10	05/25/10	KWG1004900	
Bromoform	**62.3**		5.6	0.34	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	**49.8**		23	0.15	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	**62.4**		5.6	0.40	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	**49.5**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	**51.6**		23	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	**68.7**		5.6	0.42	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	**51.6**		23	0.32	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	**50.1**		23	0.40	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	**50.5**		23	0.29	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	**49.2**		23	0.37	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	**47.2**		23	0.31	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	**46.1**		23	0.35	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	**45.8**		23	0.31	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	**47.4**		5.6	0.27	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	**46.3**		5.6	0.28	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	**44.5**		23	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	**45.7**		5.6	0.25	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	**90.3**		23	0.73	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	**40.2**		23	0.26	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	**34.5**		23	0.24	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	**42.8**		23	0.40	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	**40.3**		23	0.24	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: 05/19/2010
Date Received: 05/21/2010

Volatile Organic Compounds

Sample Name: D-4-16DMS
Lab Code: KWG1004900-2

Units: ug/Kg
Basis: Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	102	83-128	05/25/10	Acceptable
Toluene-d8	114	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	107	77-124	05/25/10	Acceptable

Comments:

Exception Report

Data File: J:\MS24\DATA\052510\0525F014.D
Lab ID: KWG1004900-2 -- K1005244-003DMS
RunType: DMS
Matrix: MISC. SOLID

Date Acquired: 05/25/2010 15:01
Date Quantitated: 05/25/2010 15:27
Batch ID: KWG1004880
Analysis Method: 8260B
MethodJoinID: MJ120

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Window	NA	NA	NA	x	
Analytical Holding Time	NA	NA	NA	x	
ICAL Pass/Fail	NA	NA	NA	x	
ICAL Average RSD	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Initial Calibration Minimum RF	NA	NA	NA		x
Initial Calibration SPCC/CCC	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Continuing Calibration Minimum RF	NA	NA	NA		x
Continuing Calibration SPCC/CCC	NA	NA	NA	x	
Internal Standards	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Relative Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Analyte Exceptions

Exception Categories	Analyte Name	Result	Low Limit	High Limit	Corrective Action
Initial Calibration Minimum RF	1,4-Dioxane	0.0037	0.01	NA	H
Continuing Calibration Minimum RF	1,4-Dioxane	0.0034	0.01	NA	I

Primary Review: [signature] 5/25/10

Secondary Review: [signature] 5/25/10

Printed: 05/25/2010 15:34:20
u:\Stealth\Crystal.rpt\except2.rpt

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	MISC. SOLID
Prod Code:	8260B VOC_FP	**Collect Date:**		**Receive Date:**	06/10/2010

Analysis Lot:	KWG1004880	**Prep Lot:**	KWG1004900	**Report Group:**	
Analysis Method:	8260B	**Prep Method:**	EPA 5030A		
Prep Ref:	911429	**Prep Date:**	05/25/2010		

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:			
Tune Ref:	J:\MS24\DATA\052510\0525F002.D	**Method ID:**	MJ120
MB Ref:	J:\MS24\DATA\052510\0525F009.D	**Quant based on Method**	

Data File:	J:\MS24\DATA\052510\0525F014.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 15:01	**Quant Date:**	05/25/2010 15:27	**Vial:**	13
Run Type:	DMS			**Dilution:**	1.0
Lab ID:	KWG1004900-2 -- K1005244-003DMS			**Soln Conc. Units:**	PPB

Internal Standard Compounds

IS Ref	Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Area Criteria
1	Fluorobenzene	4.02	0.00	96	367092	50.00	OK
2	Chlorobenzene-d5	6.65	0.00	82	163841	50.00	OK
3	1,4-Dichlorobenzene-d4	8.98	0.00	152	139592	50.00	OK

Surrogate Compounds

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	%Rec	%Rec Limits		Rpt?
1	Dibromofluoromethane	3.54	0.00	0.00	113	93825	51.14	102	83-128	OK	
1	1,2-Dichloroethane-d4	3.81	0.00	0.00	65	135333	51.10	102	64-142	OK	
1	Toluene-d8	5.31	0.00	0.00	98	367458	57.10	114	90-125	OK	
2	4-Bromofluorobenzene	7.83	0.00	0.00	95	134383	53.44	107	77-124	OK	

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Dichlorodifluoromethane	0.86		0.00	85	105865	46.71	51.5		
1	Chloromethane	0.99		0.00	50	137235	47.77	52.7		
1	Vinyl Chloride	1.06		0.00	62	128936	50.59	55.8		
1	Bromomethane	1.29	-0.01	0.00	96	65717**m**	31.00	34.2		
1	Chloroethane	1.36	-0.01	0.00	64	72618**m**	43.71	48.2		
1	Dichlorofluoromethane (CFC 21	1.52		0.00	67	549	0.1300	0.143	J	
1	Trichlorofluoromethane	1.51		0.00	101	132449	37.68	41.5		
1	Diethyl Ether	1.73		0.00	59	96212	45.77	50.5		
1	Acrolein	1.87		0.00	56	28197	69.46	76.6	J	
1	Trichlorotrifluoroethane	1.86		0.00	151	61620	45.87	50.6		
1	1,1-Dichloroethene	1.88		0.00	96	69162	44.87	49.5		
1	Acetone	1.98		0.00	43	256495	300.09	331		
1	Iodomethane (Methyl Iodide)	2.00		0.00	142	175694	83.45	92.0		
1	Carbon Disulfide	2.01		0.00	76	594146	98.09	108		

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F014.D		**Instrument:** MS24
Acqu Date:	05/25/2010 15:01	**Quant Date:** 05/25/2010 15:27	**Vial:** 13
Run Type:	DMS		**Dilution:** 1.0
Lab ID:	KWG1004900-2 -- K1005244-003DMS		**Soln Conc. Units:** PPB

Target Compounds — **Final Conc. Units:** ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	3-Chloro-1-propene	2.16		0.00	76	82277	74.62	82.3		
1	Acetonitrile	2.23		0.00	40	152807	910.41	1000		
1	Methylene Chloride	2.27		0.00	84	96110	43.49	47.9		
1	tert-Butyl Alcohol	2.35		0.00	59	326546	1,065	1170		
1	Acrylonitrile	2.50		0.00	53	137008	160.38	177		
1	Methyl tert-Butyl Ether	2.40		0.00	73	340660	52.73	58.1		
1	trans-1,2-Dichloroethene	2.42		0.00	96	86691	44.96	49.6		
1	n-Hexane	2.54		0.00	57	189997	71.90	79.3		
1	Diisopropyl Ether	2.74		0.00	45	669491	92.18	102		
1	1,1-Dichloroethane	2.75		0.00	63	172414	45.33	50.0		
1	Vinyl Acetate	2.78		0.00	86	90374	255.35	282		
1	Chloroprene	2.78		0.00	53	241038	93.17	103		
1	tert-Butyl Ethyl Ether	2.99		0.00	59	650569	98.66	109		
1	2,2-Dichloropropane	3.15		0.00	77	125283	48.06	53.0		
1	cis-1,2-Dichloroethene	3.18		0.00	96	91625	42.83	47.2		
1	2-Butanone (MEK)	3.21		0.00	72	106632	387.85	428		
1	Propionitrile	3.32		0.00	54	33468	98.30	108		
1	Methacrylonitrile	3.40		0.00	67	95221	105.32	116		
1	Bromochloromethane	3.36		0.00	128	47317	45.12	49.7		
1	Chloroform	3.41		0.00	83	162046	45.15	49.8		
1	1,1,1-Trichloroethane (TCA)	3.51		0.00	97	130329	47.56	52.4		
1	Carbon Tetrachloride	3.59		0.00	117	109580	46.50	51.3		
1	1,1-Dichloropropene	3.63		0.00	75	111676	45.35	50.0		
1	Isobutanol	3.78		0.00	43	187426	1,669	1840		
1	Benzene	3.79		0.00	78	349433	44.68	49.3		
1	1,2-Dichloroethane (EDC)	3.87		0.00	62	144077	45.41	50.1		
1	tert-Amyl Methyl Ether	3.85	0.01	0.00	55	164336	96.98	107		
1	Trichloroethene (TCE)	4.29		0.00	95	99187	47.98	52.9		
1	1,2-Dichloropropane	4.53		0.00	63	96892	46.51	51.3		
1	Dibromomethane	4.62		0.00	93	61781	46.93	51.7		
1	Methyl Methacrylate	4.60		0.00	69	160701	110.22	122		
1	1,4-Dioxane	4.62		0.00	88	35089	1,303	1440		
1	Bromodichloromethane	4.75		0.00	83	124125	47.26	52.1		
1	2-Nitropropane	5.02		0.00	41	61458	112.62	124		
1	2-Chloroethyl Vinyl Ether	5.02		0.00	63	85096	66.08	72.9		
1	cis-1,3-Dichloropropene	5.13		0.00	75	124553	42.49	46.8		
1	4-Methyl-2-pentanone (MIBK)	5.27		0.00	58	368189	460.26	507		
1	Toluene	5.37		0.00	92	215866	45.95	50.7		
2	trans-1,3-Dichloropropene	5.65		0.00	75	149518	55.14	60.8		
2	Ethyl Methacrylate	5.67		0.00	69	258051	100.04	110		
2	1,1,2-Trichloroethane	5.81		0.00	83	79601	46.90	51.7		
2	Tetrachloroethene (PCE)	5.82		0.00	164	71278	44.35	48.9		

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N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

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#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F014.D		**Instrument:** MS24
Acqu Date:	05/25/2010 15:01	**Quant Date:** 05/25/2010 15:27	**Vial:** 13
Run Type:	DMS		**Dilution:** 1.0
Lab ID:	KWG1004900-2 -- K1005244-003DMS		**Soln Conc. Units:** PPB

Target Compounds — **Final Conc. Units:** ug/Kg Dry Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2	2-Hexanone	6.02		0.00	57	138222	482.48	532		
2	1,3-Dichloropropane	5.97		0.00	76	160506	47.49	52.4		
2	Dibromochloromethane	6.13		0.00	129	93346	49.19	54.2		
2	1,2-Dibromoethane (EDB)	6.25		0.00	107	92687	50.00	55.1		
2	1-Chlorohexane	6.64		0.00	91	109319	46.93	51.7		
2	Chlorobenzene	6.68		0.00	112	234345	41.43	45.7		
2	Ethylbenzene	6.76		0.00	106	121240	44.01	48.5		
2	1,1,1,2-Tetrachloroethane	6.77		0.00	131	85278	45.69	50.4		
2	m,p-Xylenes	6.87		0.00	106	293018	90.27	99.5		
2	o-Xylene	7.27		0.00	106	138371	42.46	46.8		
2	Styrene	7.30		0.00	103	108579**m**	42.49	46.8		
2	Bromoform	7.51		0.00	173	65132	56.50	62.3		
2	Isopropylbenzene	7.62		0.00	105	355229	45.13	49.8		
2	cis-1,4-Dichloro-2-butene	7.80		0.00	89	31298	141.03	155		
3	1,1,2,2-Tetrachloroethane	8.03		0.00	83	115054	56.61	62.4		
3	trans-1,4-Dichloro-2-butene	8.10		0.00	53	83261	173.05	191		
3	Bromobenzene	7.96		0.00	156	98504	44.94	49.5		
3	n-Propylbenzene	8.04		0.00	91	418498	46.80	51.6		
3	1,2,3-Trichloropropane	8.07		0.00	110	48145	62.34	68.7		
3	2-Chlorotoluene	8.15		0.00	91	262507	46.82	51.6		
3	1,3,5-Trimethylbenzene	8.23		0.00	105	285321	45.41	50.1		
3	4-Chlorotoluene	8.27		0.00	91	259197	45.79	50.5		
3	tert-Butylbenzene	8.54		0.00	119	245772	44.65	49.2		
3	1,2,4-Trimethylbenzene	8.60		0.00	105	274937	42.79	47.2		
3	sec-Butylbenzene	8.76		0.00	105	343178	41.84	46.1		
3	4-Isopropyltoluene	8.91		0.00	119	284224	41.53	45.8		
3	1,3-Dichlorobenzene	8.90		0.00	146	173595	42.95	47.4		
3	1,4-Dichlorobenzene	9.00		0.00	146	175633	41.96	46.3		
3	n-Butylbenzene	9.31		0.00	91	244049	40.38	44.5		
3	1,2-Dichlorobenzene	9.36		0.00	146	164707	41.46	45.7		
3	1,2-Dibromo-3-chloropropane	10.14		0.00	155	24863	81.88	90.3		
3	1,2,4-Trichlorobenzene	10.85	-0.01	0.00	180	91633	36.44	40.2		
3	Hexachlorobutadiene	10.96		0.00	225	39384	31.30	34.5		
3	Naphthalene	11.09		0.00	128	224567	38.83	42.8		
3	1,2,3-Trichlorobenzene	11.32		0.00	180	85690	36.59	40.3		
	4-Vinyl-1-cyclohexene				0	0		5.5	UJ	NR
	Benzyl Chloride				0	0		5.5	UJ	NR
	Hexachloroethane				0	0		1.1	UJ	NR
	2,2,4-Trimethylpentane				0	0		5.5	UJ	NR
	1,3-Butadiene				0	0		2.2	UJ	NR
	1,1,2-Trifluoroethane				0	0		11	UJ	NR
	n-Butyl Alcohol				0	0		280	U	NR

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#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F014.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 15:01	**Quant Date:**	05/25/2010 15:27	**Vial:**	13
Run Type:	DMS			**Dilution:**	1.0
Lab ID:	KWG1004900-2 -- K1005244-003DMS			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Dry Weight

Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2-Propanol			0	0		220	U	NR
Ethyl Alcohol			0	0		220	U	NR
2,2-Dichloro-1,1,1-trifluoroetha			0	0		5.5	UJ	NR
Cyclohexanone			0	0		11	UJ	NR
Methylcyclopentane			0	0		5.5	UJ	NR
n-Heptane			0	0		5.5	UJ	NR
3-Methylpentane			0	0		5.5	UJ	NR
2-Methylpentane			0	0		5.5	UJ	NR
Ethyl Acetate			0	0		22	UJ	NR

Prep Amount:	5.00 g	**Dilution:**	1.0
Prep Final Vol:	5.0 ml	**Unit Factor:**	1
Solids:	90.7 %		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
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E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

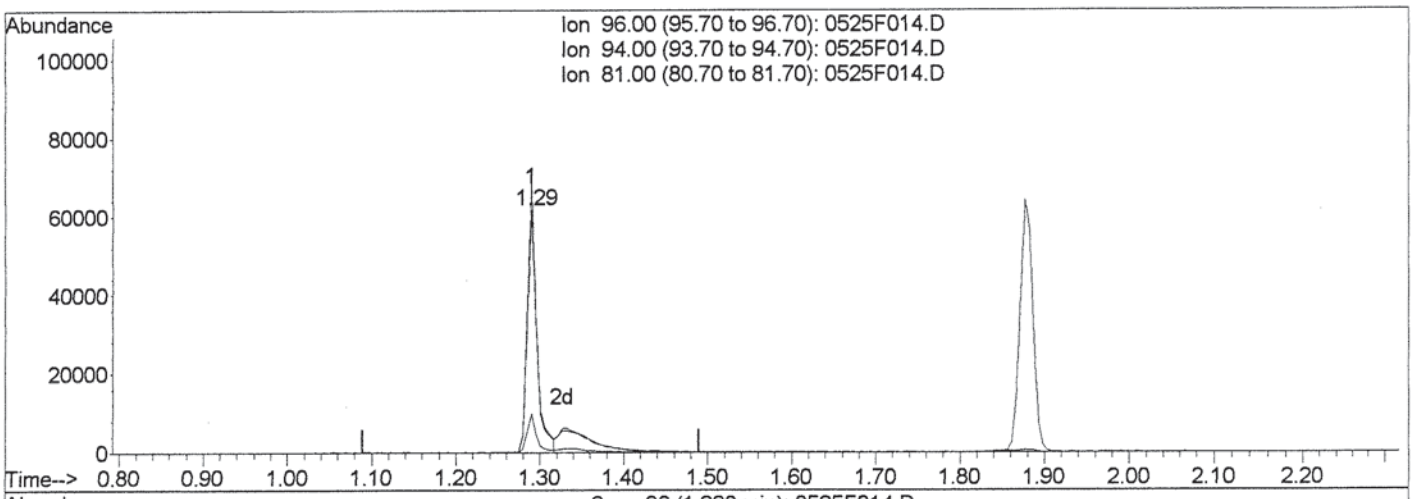
D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

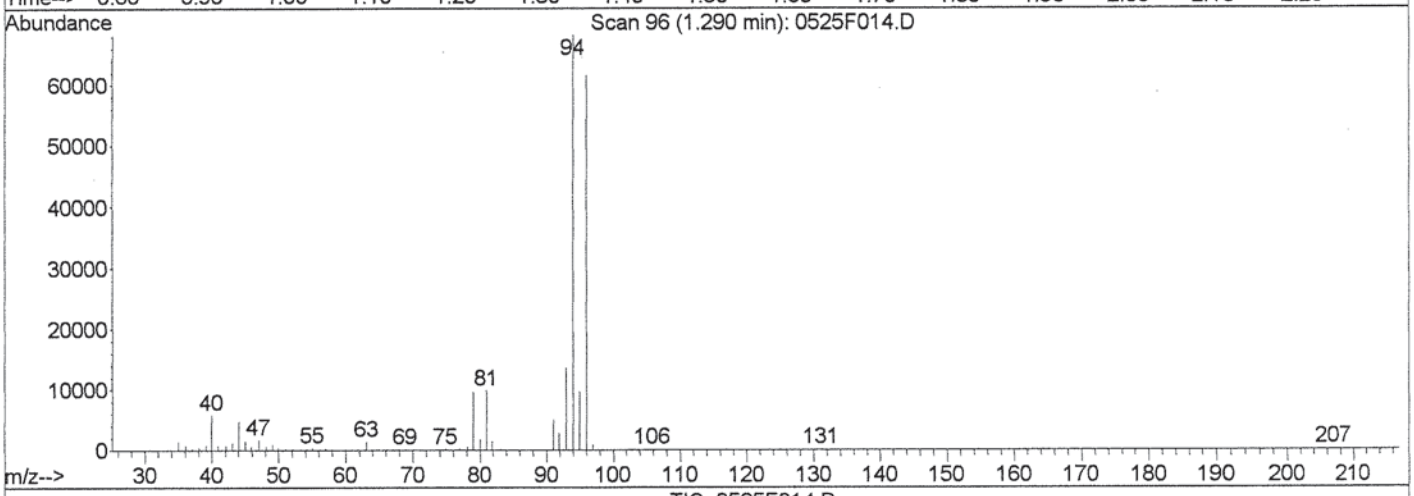
*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F014.D Vial: 13
Acq On : 25 May 2010 3:01 pm Operator: KR
Sample : K5244-003DMS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F014.D

(5) Bromomethane (T)

1.29min 22.86PPB

response 48469

Ion	Exp%	Act%
96.00	100	100
94.00	106.40	110.52
81.00	18.10	15.76
0.00	0.00	0.00

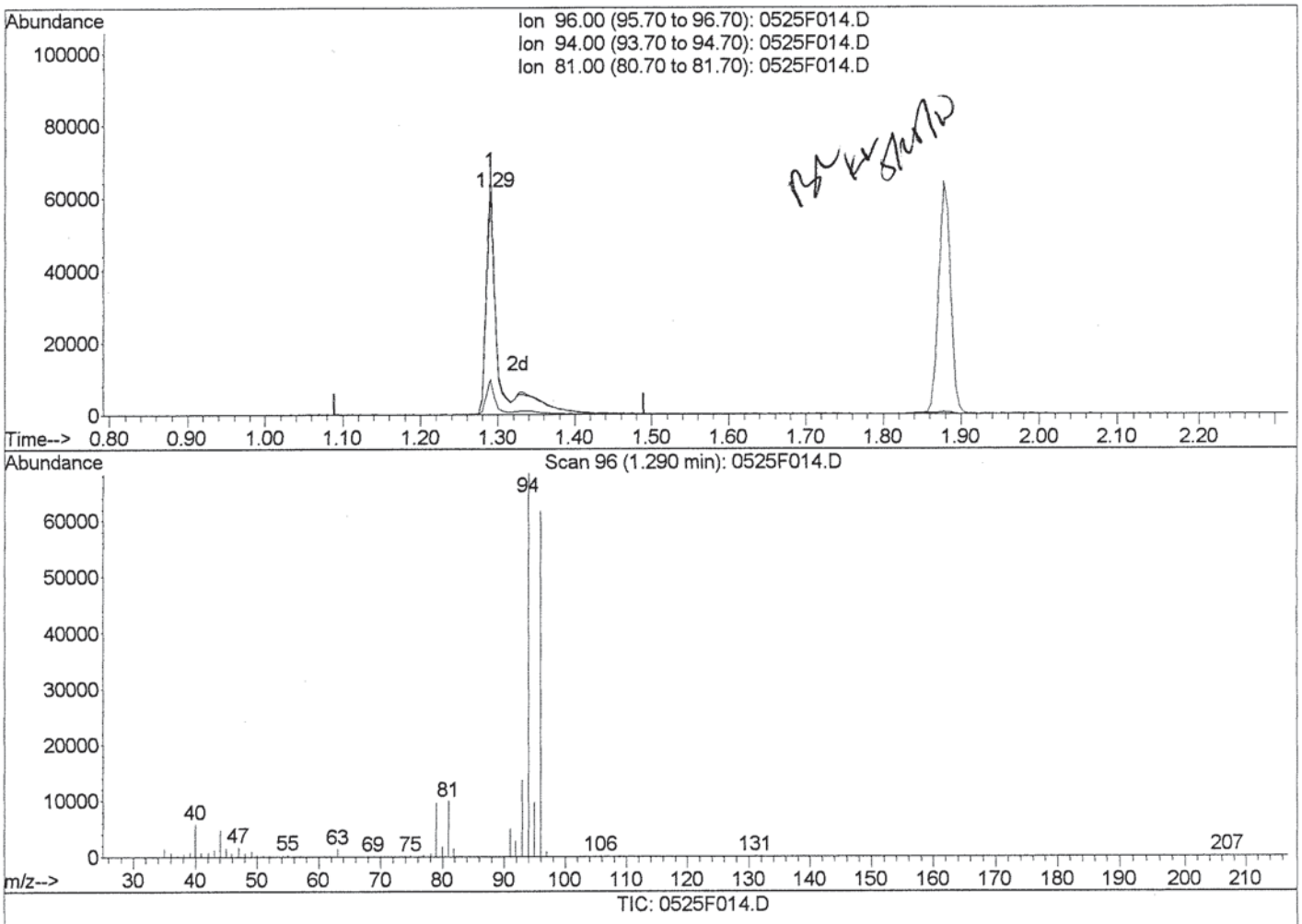
0525F014.D 042110MS24SOIL.M Tue May 25 15:27:25 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F014.D
Acq On : 25 May 2010 3:01 pm
Sample : K5244-003DMS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010

Vial: 13
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration



TIC: 0525F014.D

(5) Bromomethane (T)

1.29min 31.00PPB m

response 65717

Ion	Exp%	Act%
96.00	100	100
94.00	106.40	110.78
81.00	18.10	15.95
0.00	0.00	0.00

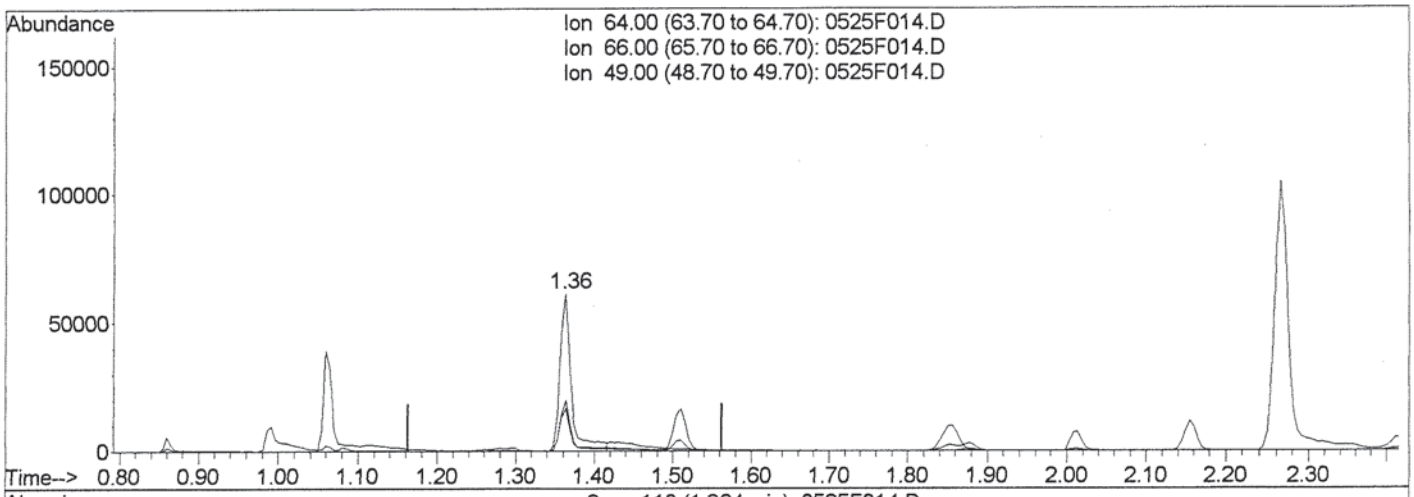
Lmm 5/25/10

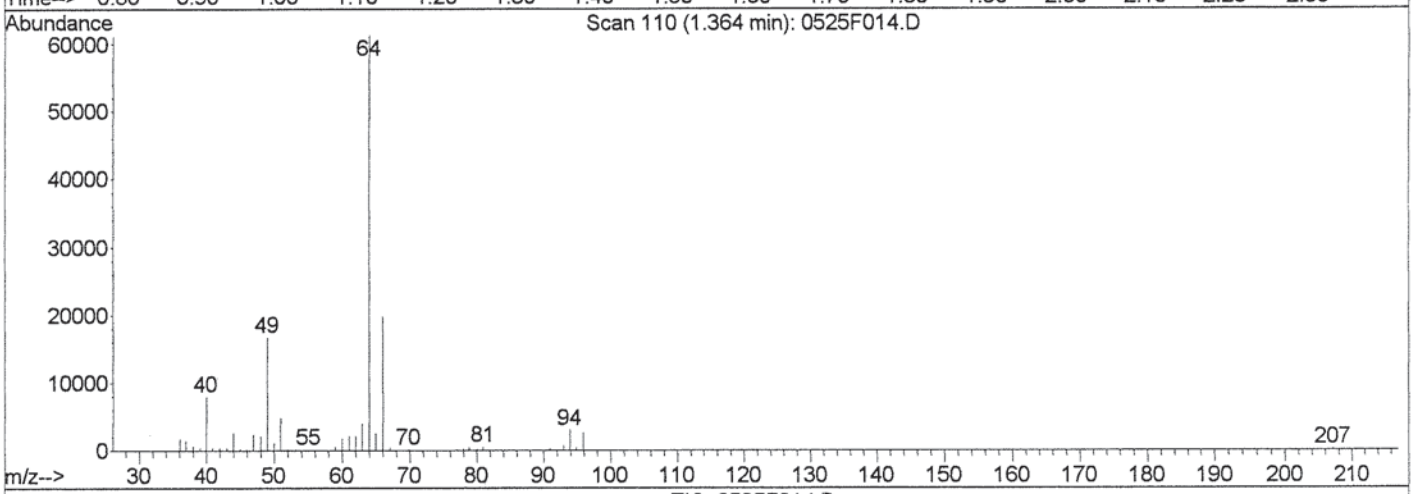
0525F014.D 042110MS24SOIL.M Tue May 25 15:27:29 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F014.D Vial: 13
Acq On : 25 May 2010 3:01 pm Operator: KR
Sample : K5244-003DMS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F014.D

(6) Chloroethane (T)

1.36min 37.90PPB

response 62961

Ion	Exp%	Act%
64.00	100	100
66.00	32.10	32.03
49.00	27.50	27.23
0.00	0.00	0.00

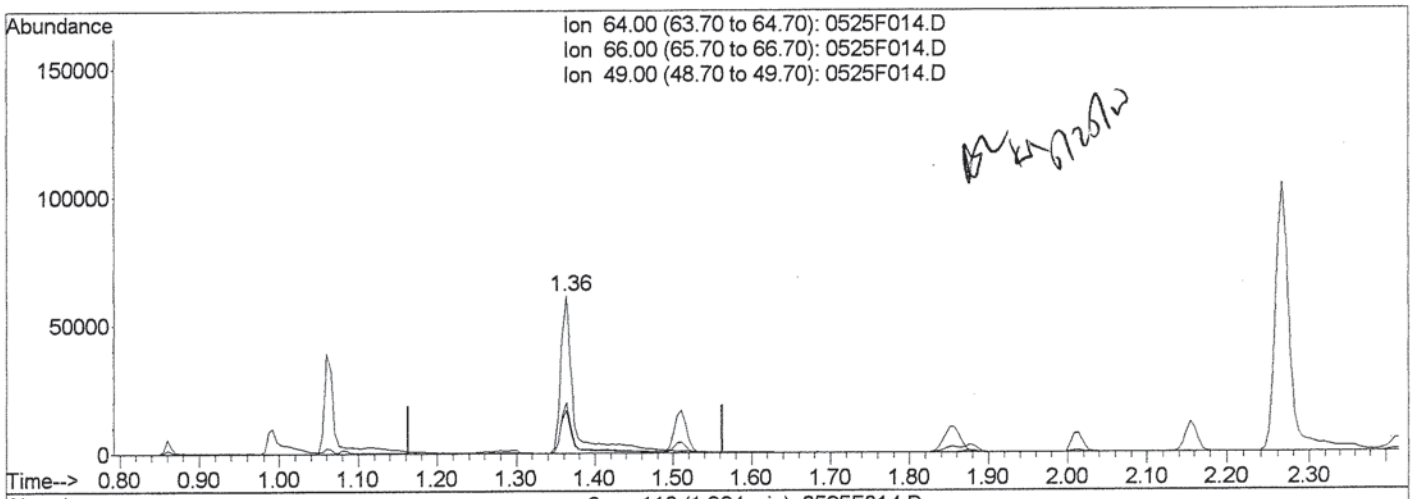
0525F014.D 042110MS24SOIL.M Tue May 25 15:27:34 2010

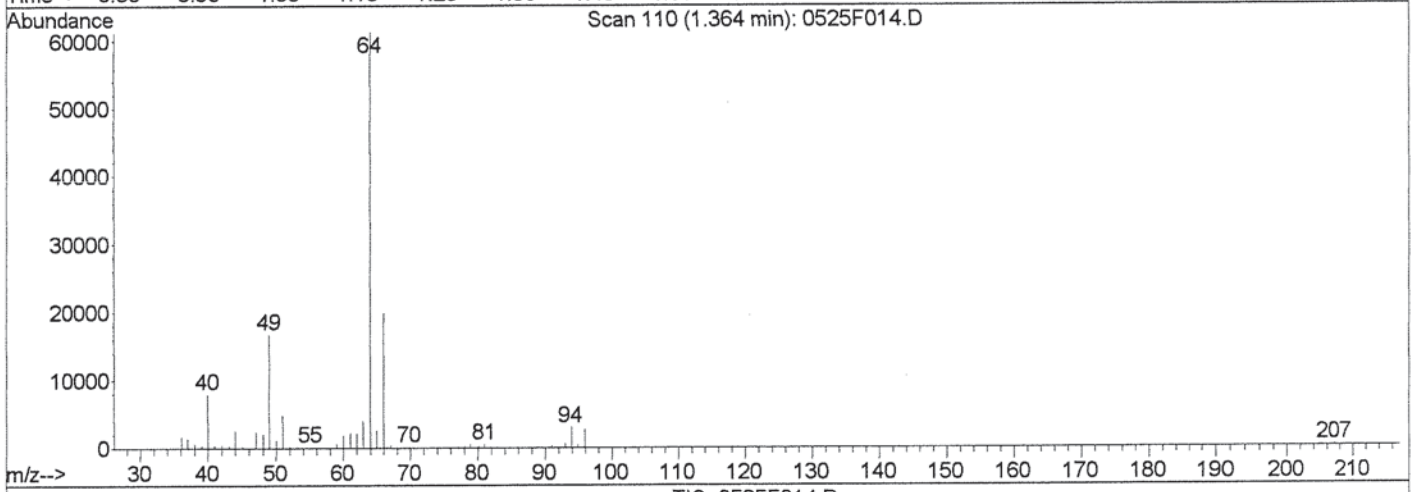
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F014.D
Acq On : 25 May 2010 3:01 pm
Sample : K5244-003DMS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010

Vial: 13
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F014.D

(6) Chloroethane (T)

1.36min 43.71PPB m

response 72618

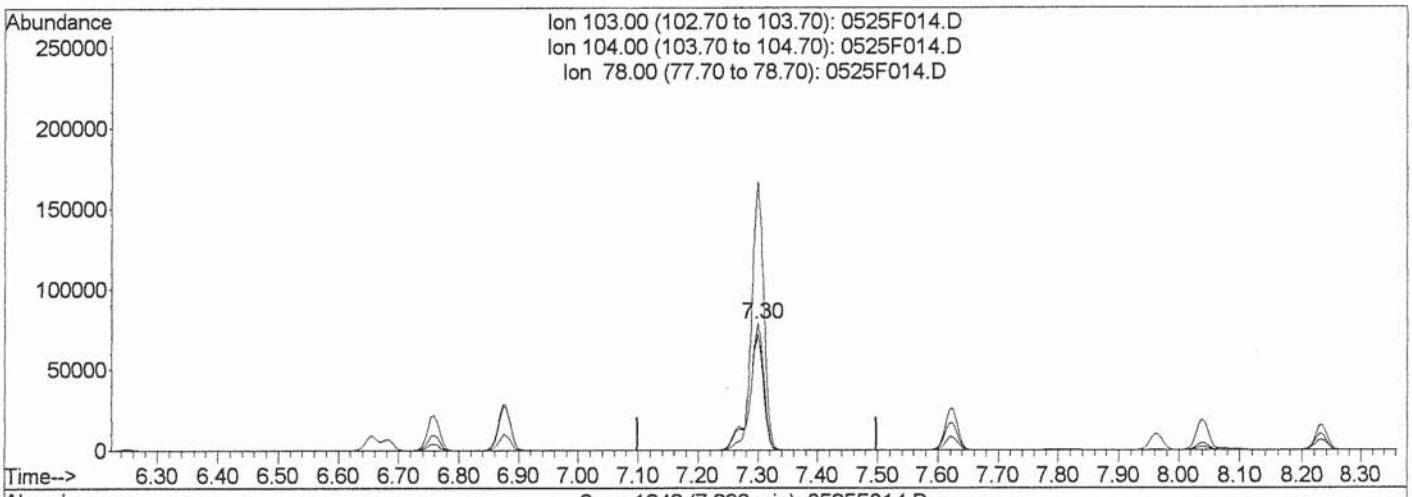
Ion	Exp%	Act%
64.00	100	100
66.00	32.10	32.28
49.00	27.50	27.23
0.00	0.00	0.00

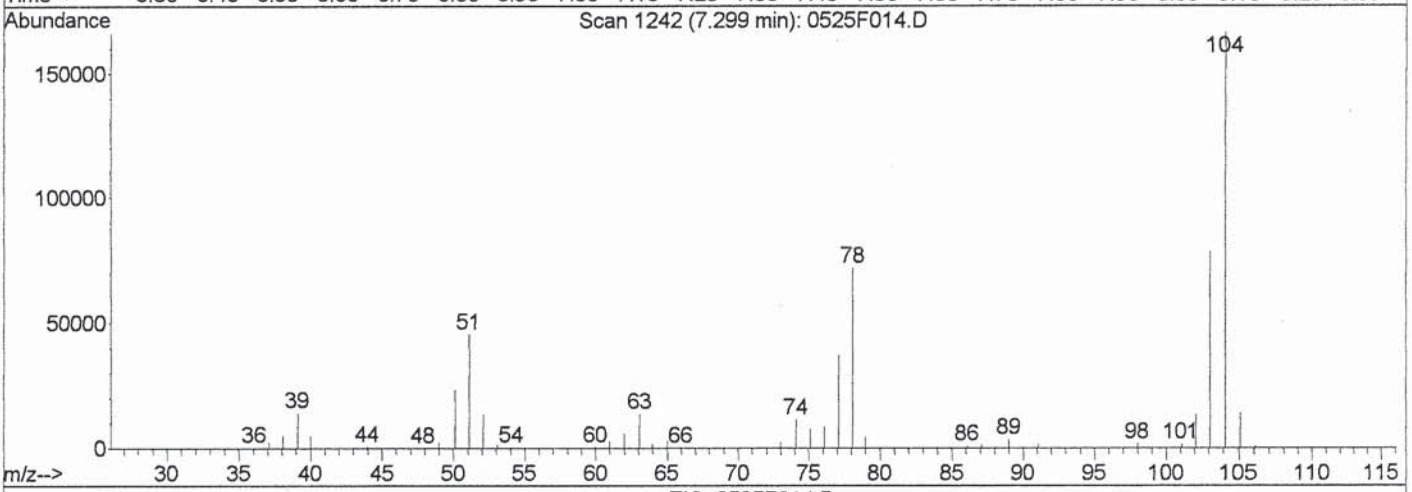
0525F014.D 042110MS24SOIL.M Tue May 25 15:27:39 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F014.D Vial: 13
Acq On : 25 May 2010 3:01 pm Operator: KR
Sample : K5244-003DMS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F014.D

(75) Styrene (T)

7.30min 50.12PPB

response 128079

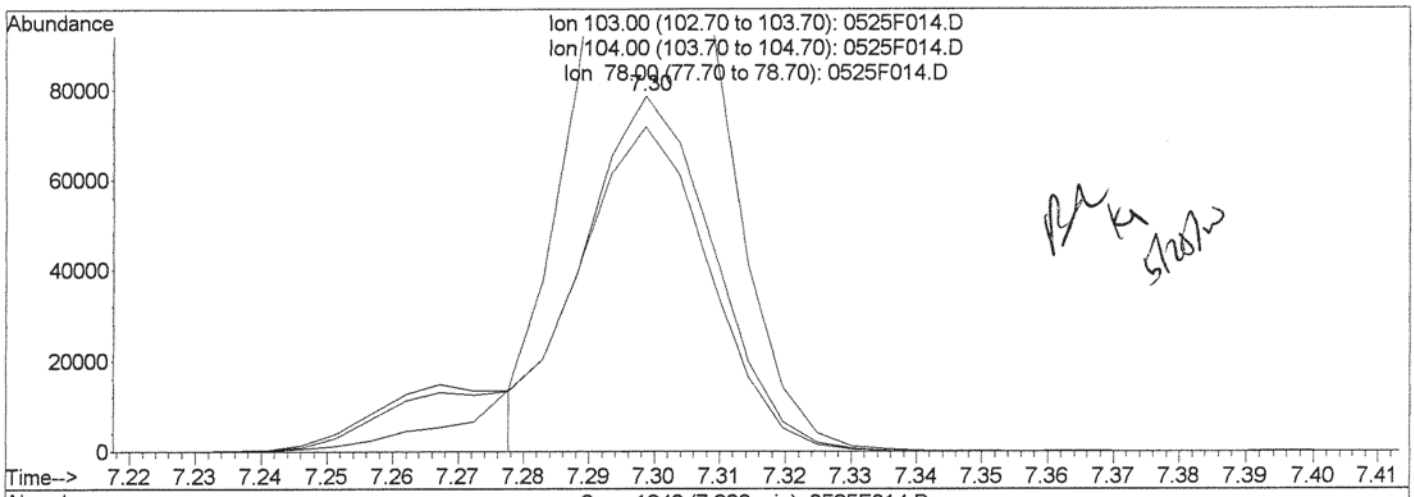
Ion	Exp%	Act%
103.00	100	100
104.00	208.80	211.90
78.00	88.30	91.31
0.00	0.00	0.00

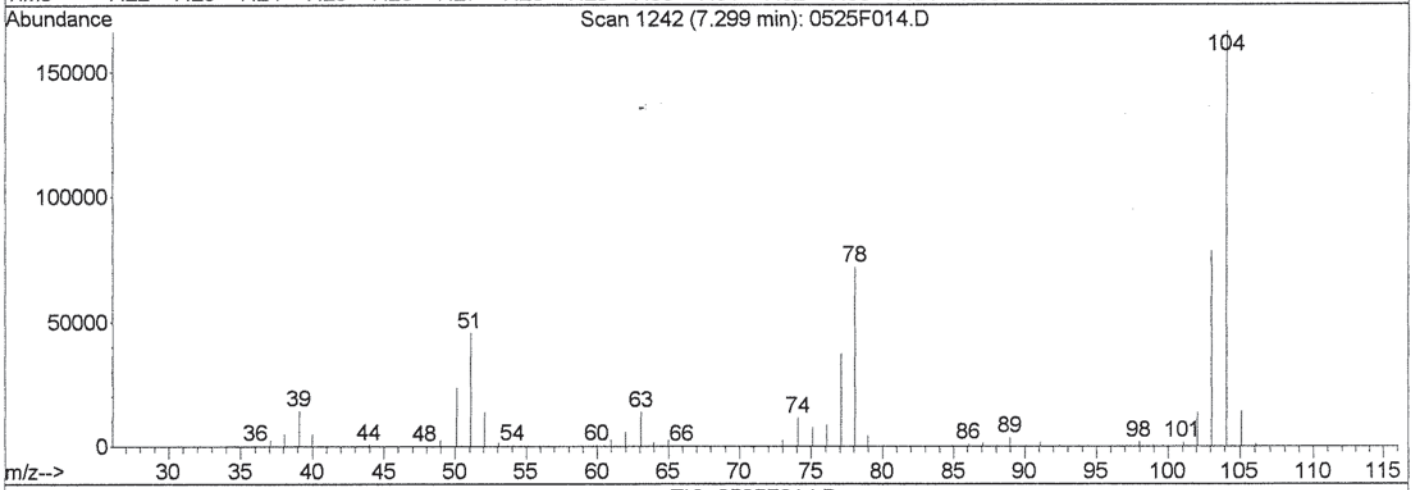
0525F014.D 042110MS24SOIL.M Tue May 25 15:27:46 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F014.D Vial: 13
Acq On : 25 May 2010 3:01 pm Operator: KR
Sample : K5244-003DMS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F014.D

(75) Styrene (T)

7.30min 42.49PPB m

response 108579

Ion	Exp%	Act%
103.00	100	100
104.00	208.80	211.90
78.00	88.30	91.31
0.00	0.00	0.00

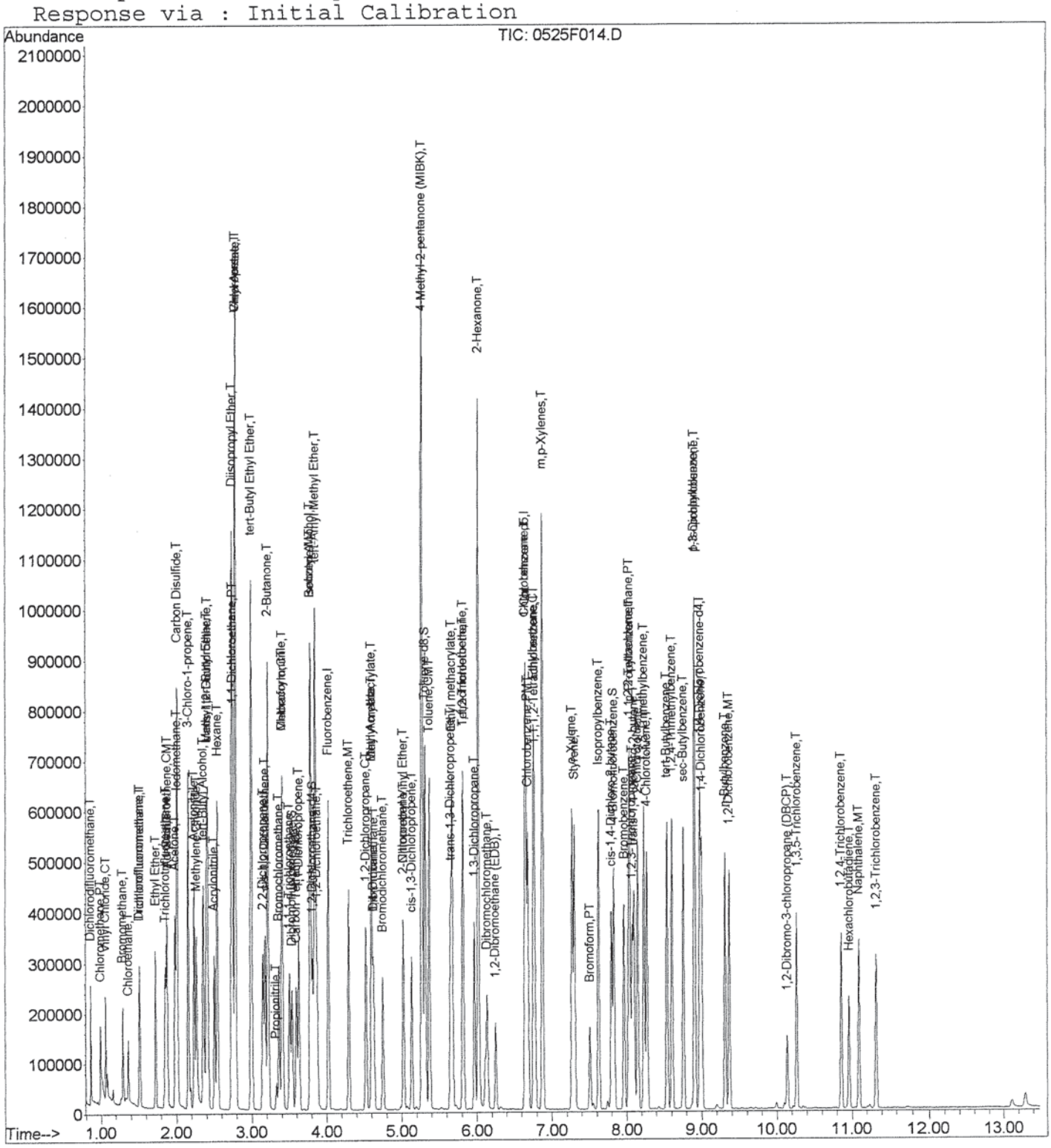
Lmm 5/25/10

0525F014.D 042110MS24SOIL.M Tue May 25 15:27:53 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\052510\0525F014.D Vial: 13
Acq On : 25 May 2010 3:01 pm Operator: KR
Sample : K5244-003DMS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 15:27 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Initial Calibration



COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Lab Control Sample
Lab Code: KWG1004900-3

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	**36.3**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	**39.1**		5.0	0.46	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	**40.0**		5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	**26.7**		5.0	0.62	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	**37.0**		5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	**30.1**		5.0	0.21	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	**36.0**		5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Acetone	**176**		20	3.9	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	**80.7**		5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**37.1**		10	0.43	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	**37.4**		5.0	0.35	1	05/25/10	05/25/10	KWG1004900	*
1,1-Dichloroethane	**38.1**		5.0	0.20	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	**40.3**		5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	**39.5**		5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**208**		20	0.99	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	**40.9**		5.0	0.14	1	05/25/10	05/25/10	KWG1004900	
Chloroform	**41.5**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	**40.9**		5.0	0.35	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	**40.5**		5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	**39.0**		5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
Benzene	**41.8**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	**40.0**		5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	**42.8**		5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	**41.3**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	**39.8**		5.0	0.28	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	**41.9**		5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	**37.5**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	**216**		20	0.74	1	05/25/10	05/25/10	KWG1004900	
Toluene	**40.9**		5.0	0.18	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	**52.4**		5.0	0.34	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	**42.9**		5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	**43.4**		5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	**230**		20	0.93	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	**44.0**		5.0	0.20	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Lab Control Sample
Lab Code: KWG1004900-3

Units: ug/Kg
Basis: Dry

Extraction Method: EPA 5030A
Analysis Method: 8260B

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	**47.3**		5.0	0.19	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	**44.2**		20	0.27	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	**43.4**		5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	**44.9**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	**47.5**		5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	**94.7**		5.0	0.37	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	**46.5**		5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Styrene	**47.8**		5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Bromoform	**52.5**		5.0	0.30	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	**48.7**		20	0.13	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	**39.3**		5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	**47.9**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	**49.0**		20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	**45.2**		5.0	0.38	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	**49.9**		20	0.29	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	**49.5**		20	0.36	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	**50.5**		20	0.26	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	**47.9**		20	0.33	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	**47.4**		20	0.28	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	**46.4**		20	0.31	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	**46.7**		20	0.28	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	**47.2**		5.0	0.24	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	**46.2**		5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	**47.9**		20	0.32	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	**46.3**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	**50.4**		20	0.66	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	**50.2**		20	0.23	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	**54.3**		20	0.21	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	**50.4**		20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	**52.6**		20	0.21	1	05/25/10	05/25/10	KWG1004900	

* See Case Narrative

Comments:

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Collected:**	NA
Sample Matrix:	Misc. solid	**Date Received:**	NA

Volatile Organic Compounds

Sample Name:	Lab Control Sample	**Units:**	ug/Kg
Lab Code:	KWG1004900-3	**Basis:**	Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	94	83-128	05/25/10	Acceptable
Toluene-d8	105	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	110	77-124	05/25/10	Acceptable

Comments:

Exception Report

Data File: J:\MS24\DATA\052510\0525F005.D
Lab ID: KWG1004900-3
RunType: LCS
Matrix: SEDIMENT

Date Acquired: 05/25/2010 11:14
Date Quantitated: 05/25/2010 14:52
Batch ID: KWG1004880
Analysis Method: 8260B
MethodJoinID: MJ120

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Window	NA	NA	NA	x	
Analytical Holding Time	NA	NA	NA	x	
ICAL Pass/Fail	NA	NA	NA	x	
ICAL Average RSD	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Initial Calibration Minimum RF	NA	NA	NA		x
Initial Calibration SPCC/CCC	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Continuing Calibration Minimum RF	NA	NA	NA		x
Continuing Calibration SPCC/CCC	NA	NA	NA	x	
Internal Standards	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Relative Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Analyte Exceptions

Exception Categories	Analyte Name	Result	Low Limit	High Limit	Corrective Action
Initial Calibration Minimum RF	1,4-Dioxane	0.0037	0.01	NA	KJ
Continuing Calibration Minimum RF	1,4-Dioxane	0.0034	0.01	NA	J

Primary Review: KLS 5/26/10

Secondary Review: CMM 5/25/10

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	SEDIMENT
Prod Code:	8260B VOC_FP	**Collect Date:**		**Receive Date:**	05/25/2010

Analysis Lot:	KWG1004880	**Prep Lot:**	KWG1004900	**Report Group:**	
Analysis Method:	8260B	**Prep Method:**	EPA 5030A		
Prep Ref:	911430	**Prep Date:**	05/25/2010		

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:			
Tune Ref:	J:\MS24\DATA\052510\0525F002.D	**Method ID:**	MJ120
MB Ref:	J:\MS24\DATA\052510\0525F009.D	**Quant based on Method**	

Data File:	J:\MS24\DATA\052510\0525F005.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:14	**Quant Date:**	05/25/2010 14:52	**Vial:**	4
Run Type:	LCS			**Dilution:**	1.0
Lab ID:	KWG1004900-3			**Soln Conc. Units:**	PPB

Internal Standard Compounds

IS Ref	Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Area Criteria
1	Fluorobenzene	4.02	0.00	96	418826	50.00	OK
2	Chlorobenzene-d5	6.65	0.00	82	167585	50.00	OK
3	1,4-Dichlorobenzene-d4	8.98	0.00	152	148678	50.00	OK

Surrogate Compounds

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	%Rec	%Rec Limits		Rpt?
1	Dibromofluoromethane	3.53	-0.01	0.00	113	97882	46.76	94	43-153	OK	
1	1,2-Dichloroethane-d4	3.81	0.00	0.00	65	130622	43.23	86	64-142	OK	
1	Toluene-d8	5.31	0.00	0.00	98	386523	52.64	105	49-151	OK	
2	4-Bromofluorobenzene	7.83	0.00	0.00	95	141811	55.13	110	33-150	OK	

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Dichlorodifluoromethane	0.85	-0.01	0.00	85	93940	36.33	36.3		
1	Chloromethane	0.99		0.00	50	128054	39.07	39.1		
1	Vinyl Chloride	1.06		0.00	62	116250	39.98	40.0		
1	Bromomethane	1.28	-0.02	0.00	96	64549m	26.68	26.7		
1	Chloroethane	1.36	-0.01	0.00	64	70171	37.02	37.0		
1	Dichlorofluoromethane (CFC 21	1.86	0.34	0.08	67	733	0.1500	0.150	J	
1	Trichlorofluoromethane	1.50	-0.01	0.00	101	120539	30.06	30.1		
1	Diethyl Ether	1.72	-0.01	0.00	59	89591	37.35	37.4		
1	Acrolein	1.87		0.00	56	30793	66.49	66.5	J	
1	Trichlorotrifluoroethane	1.85	-0.01	0.00	151	57423	37.46	37.5		
1	1,1-Dichloroethene	1.88		0.00	96	63244	35.96	36.0		
1	Acetone	1.98		0.00	43	171424	175.79	176		
1	Iodomethane (Methyl Iodide)	1.99	-0.01	0.00	142	170741	71.08	71.1		
1	Carbon Disulfide	2.01		0.00	76	557733	80.70	80.7		

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F005.D		
Acqu Date:	05/25/2010 11:14	**Quant Date:**	05/25/2010 14:52
Run Type:	LCS		
Lab ID:	KWG1004900-3		

Instrument:	MS24
Vial:	4
Dilution:	1.0
Soln Conc. Units:	PPB

Target Compounds **Final Conc. Units:** ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	3-Chloro-1-propene	2.15	-0.01	0.00	76	81477	64.77	64.8		
1	Acetonitrile	2.23		0.00	40	135744	708.85	709		
1	Methylene Chloride	2.26	-0.01	0.00	84	93446	37.06	37.1		
1	tert-Butyl Alcohol	2.35		0.00	59	126740	362.41	362		
1	Acrylonitrile	2.50		0.00	53	105585	108.33	108		
1	Methyl tert-Butyl Ether	2.40		0.00	73	302945	41.10	41.1		
1	trans-1,2-Dichloroethene	2.41	-0.01	0.00	96	82158	37.35	37.4		
1	n-Hexane	2.54		0.00	57	182056	60.39	60.4		
1	Diisopropyl Ether	2.73	-0.01	0.00	45	633926	76.50	76.5		
1	1,1-Dichloroethane	2.75		0.00	63	165226	38.08	38.1		
1	Vinyl Acetate	2.78		0.00	86	75221	186.28	186		
1	Chloroprene	2.78		0.00	53	232884	78.90	78.9		
1	tert-Butyl Ethyl Ether	2.99		0.00	59	607207	80.71	80.7		
1	2,2-Dichloropropane	3.15		0.00	77	119819	40.28	40.3		
1	cis-1,2-Dichloroethene	3.17	-0.01	0.00	96	96307	39.46	39.5		
1	2-Butanone (MEK)	3.21		0.00	72	65231	207.95	208		
1	Propionitrile	3.32		0.00	54	25937	66.77	66.8		
1	Methacrylonitrile	3.40		0.00	67	72797	70.57	70.6		
1	Bromochloromethane	3.36		0.00	128	48989	40.94	40.9		
1	Chloroform	3.41		0.00	83	169934	41.50	41.5		
1	1,1,1-Trichloroethane (TCA)	3.50	-0.01	0.00	97	127810	40.88	40.9		
1	Carbon Tetrachloride	3.59		0.00	117	108965	40.53	40.5		
1	1,1-Dichloropropene	3.62	-0.01	0.00	75	109683	39.04	39.0		
1	Isobutanol	3.78		0.00	43	86901	678.35	678		
1	Benzene	3.79		0.00	78	372961	41.80	41.8		
1	1,2-Dichloroethane (EDC)	3.87		0.00	62	144784	39.99	40.0		
1	tert-Amyl Methyl Ether	3.84		0.00	55	163950	84.80	84.8		
1	Trichloroethene (TCE)	4.29		0.00	95	100882	42.77	42.8		
1	1,2-Dichloropropane	4.52	-0.01	0.00	63	98099	41.27	41.3		
1	Dibromomethane	4.62		0.00	93	59837	39.84	39.8		
1	Methyl Methacrylate	4.60		0.00	69	114382	68.76	68.8		
1	1,4-Dioxane	4.62		0.00	88	24461	795.98	796		
1	Bromodichloromethane	4.75		0.00	83	125589	41.91	41.9		
1	2-Nitropropane	5.02		0.00	41	43041	69.13	69.1		
1	2-Chloroethyl Vinyl Ether	5.02		0.00	63	65320	44.46	44.5		
1	cis-1,3-Dichloropropene	5.13		0.00	75	125501	37.52	37.5		
1	4-Methyl-2-pentanone (MIBK)	5.26	-0.01	0.00	58	197121	215.97	216		
1	Toluene	5.36	-0.01	0.00	92	219274	40.91	40.9		
2	trans-1,3-Dichloropropene	5.65		0.00	75	145282	52.38	52.4		
2	Ethyl Methacrylate	5.67		0.00	69	210382	79.74	79.7		
2	1,1,2-Trichloroethane	5.81		0.00	83	74501	42.92	42.9		
2	Tetrachloroethene (PCE)	5.82		0.00	164	71345	43.40	43.4		

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
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N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
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*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F005.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:14	**Quant Date:**	05/25/2010 14:52	**Vial:**	4
Run Type:	LCS			**Dilution:**	1.0
Lab ID:	KWG1004900-3			**Soln Conc. Units:**	PPB

Target Compounds — **Final Conc. Units:** ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2	2-Hexanone	6.02		0.00	57	67292	229.65	230		
2	1,3-Dichloropropane	5.97		0.00	76	152238	44.03	44.0		
2	Dibromochloromethane	6.13		0.00	129	91877	47.33	47.3		
2	1,2-Dibromoethane (EDB)	6.25		0.00	107	83813	44.20	44.2		
2	1-Chlorohexane	6.64		0.00	91	113281	47.54	47.5		
2	Chlorobenzene	6.68		0.00	112	250826	43.35	43.4		
2	Ethylbenzene	6.76		0.00	106	126408	44.86	44.9		
2	1,1,1,2-Tetrachloroethane	6.77		0.00	131	90599	47.46	47.5		
2	m,p-Xylenes	6.87		0.00	106	314370	94.68	94.7		
2	o-Xylene	7.27		0.00	106	154993	46.50	46.5		
2	Styrene	7.30		0.00	103	124929m	47.79	47.8		
2	Bromoform	7.51		0.00	173	61934	52.53	52.5		
2	Isopropylbenzene	7.62		0.00	105	392329	48.73	48.7		
2	cis-1,4-Dichloro-2-butene	7.80		0.00	89	25161	110.84	111		
3	1,1,2,2-Tetrachloroethane	8.03		0.00	83	85096	39.31	39.3		
3	trans-1,4-Dichloro-2-butene	8.10		0.00	53	60236	117.54	118		
3	Bromobenzene	7.96		0.00	156	111713	47.85	47.9		
3	n-Propylbenzene	8.04		0.00	91	466564	48.99	49.0		
3	1,2,3-Trichloropropane	8.07		0.00	110	37204	45.23	45.2		
3	2-Chlorotoluene	8.15		0.00	91	298225	49.94	49.9		
3	1,3,5-Trimethylbenzene	8.23		0.00	105	331375	49.52	49.5		
3	4-Chlorotoluene	8.27		0.00	91	304369	50.48	50.5		
3	tert-Butylbenzene	8.54		0.00	119	280554	47.86	47.9		
3	1,2,4-Trimethylbenzene	8.60		0.00	105	324183	47.37	47.4		
3	sec-Butylbenzene	8.76		0.00	105	405317	46.40	46.4		
3	4-Isopropyltoluene	8.91		0.00	119	340079	46.66	46.7		
3	1,3-Dichlorobenzene	8.90		0.00	146	203378	47.24	47.2		
3	1,4-Dichlorobenzene	9.00		0.00	146	206016	46.21	46.2		
3	n-Butylbenzene	9.31		0.00	91	308038	47.85	47.9		
3	1,2-Dichlorobenzene	9.36		0.00	146	195804	46.27	46.3		
3	1,2-Dibromo-3-chloropropane	10.14		0.00	155	16290	50.37	50.4		
3	1,2,4-Trichlorobenzene	10.86		0.00	180	134358	50.16	50.2		
3	Hexachlorobutadiene	10.96		0.00	225	72715	54.25	54.3		
3	Naphthalene	11.09		0.00	128	310313	50.38	50.4		
3	1,2,3-Trichlorobenzene	11.32		0.00	180	131110	52.56	52.6		
	4-Vinyl-1-cyclohexene				0	0		5.0	UJ	NR
	Benzyl Chloride				0	0		5.0	UJ	NR
	Hexachloroethane				0	0		1.0	UJ	NR
	2,2,4-Trimethylpentane				0	0		5.0	UJ	NR
	1,3-Butadiene				0	0		2.0	UJ	NR
	1,1,2-Trifluoroethane				0	0		10	UJ	NR
	n-Butyl Alcohol				0	0		250	U	NR

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F005.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:14	**Quant Date:**	05/25/2010 14:52	**Vial:**	4
Run Type:	LCS			**Dilution:**	1.0
Lab ID:	KWG1004900-3			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2-Propanol			0	0		200	U	NR
Ethyl Alcohol			0	0		200	U	NR
2,2-Dichloro-1,1,1-trifluoroetha			0	0		5.0	UJ	NR
Cyclohexanone			0	0		10	UJ	NR
2-Methylpentane			0	0		5.0	UJ	NR
n-Heptane			0	0		5.0	UJ	NR
3-Methylpentane			0	0		5.0	UJ	NR
Methylcyclopentane			0	0		5.0	UJ	NR
Ethyl Acetate			0	0		20	UJ	NR

Prep Amount:	5.00 g	**Dilution:**	1.0
Prep Final Vol:	5.0 ml	**Unit Factor:**	1
Solids:	%		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

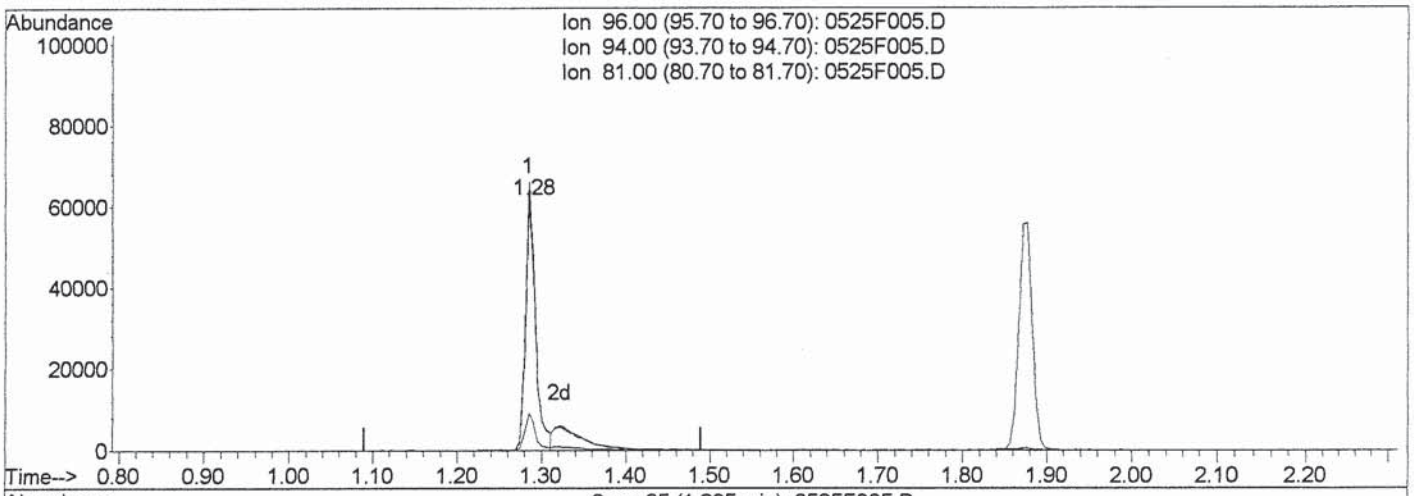
D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

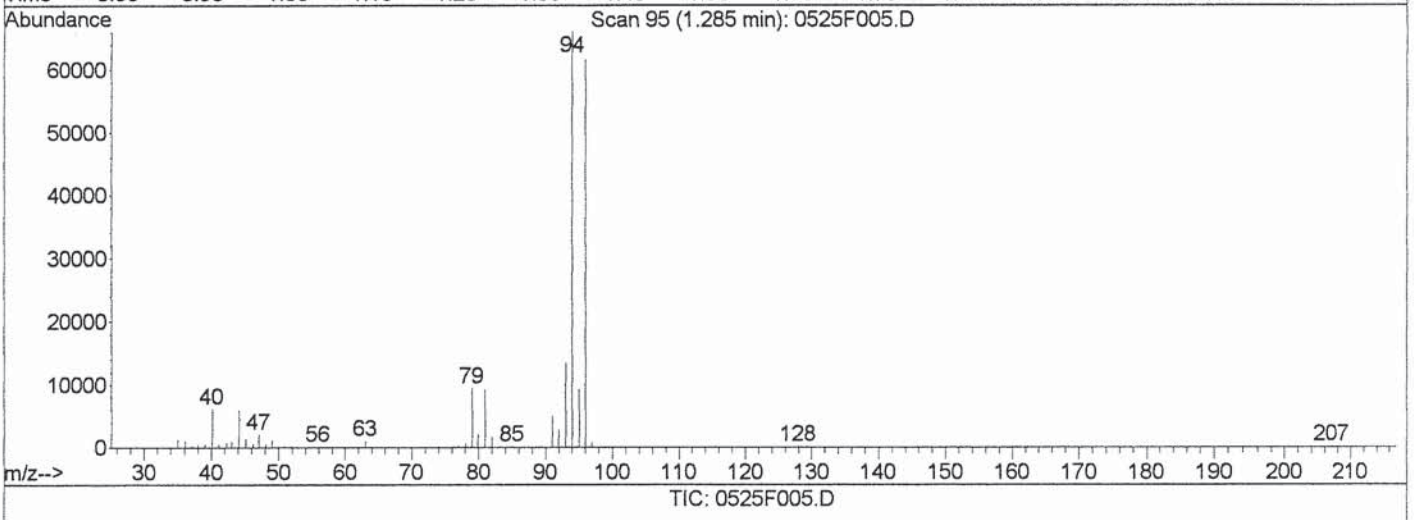
*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F005.D Vial: 4
Acq On : 25 May 2010 11:14 am Operator: KR
Sample : LCS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 14:52 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F005.D

(5) Bromomethane (T)

1.28min 20.81PPB

response 50347

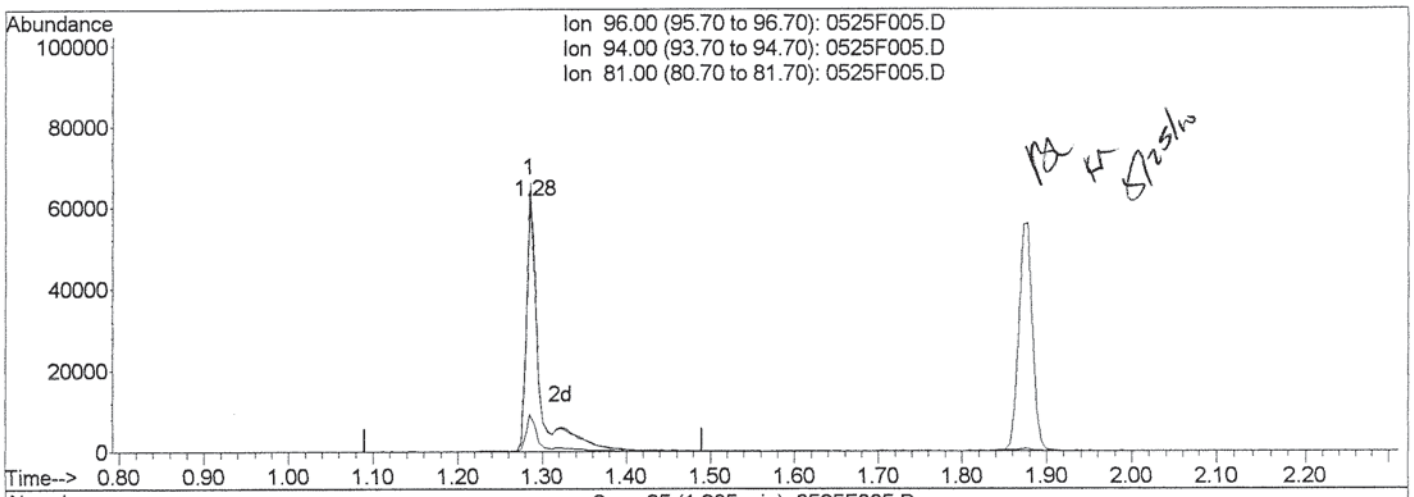
Ion	Exp%	Act%
96.00	100	100
94.00	106.40	106.81
81.00	18.10	14.66
0.00	0.00	0.00

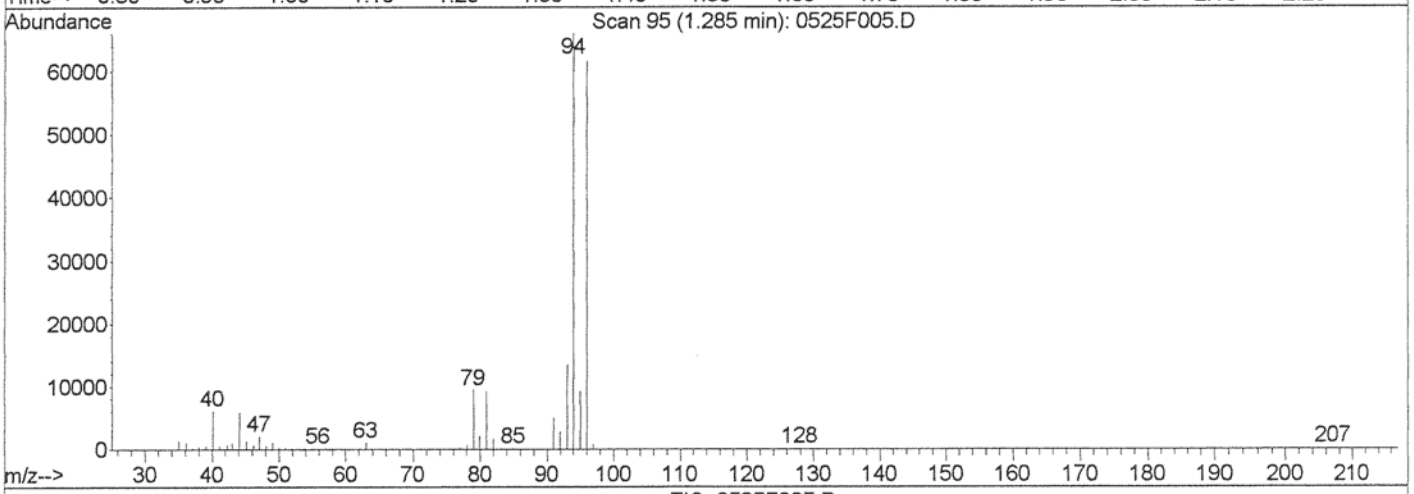
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F005.D
Acq On : 25 May 2010 11:14 am
Sample : LCS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 14:52 2010

Vial: 4
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F005.D

(5) Bromomethane (T)

1.28min 26.68PPB m

response 64549

Ion	Exp%	Act%
96.00	100	100
94.00	106.40	107.10
81.00	18.10	14.86
0.00	0.00	0.00

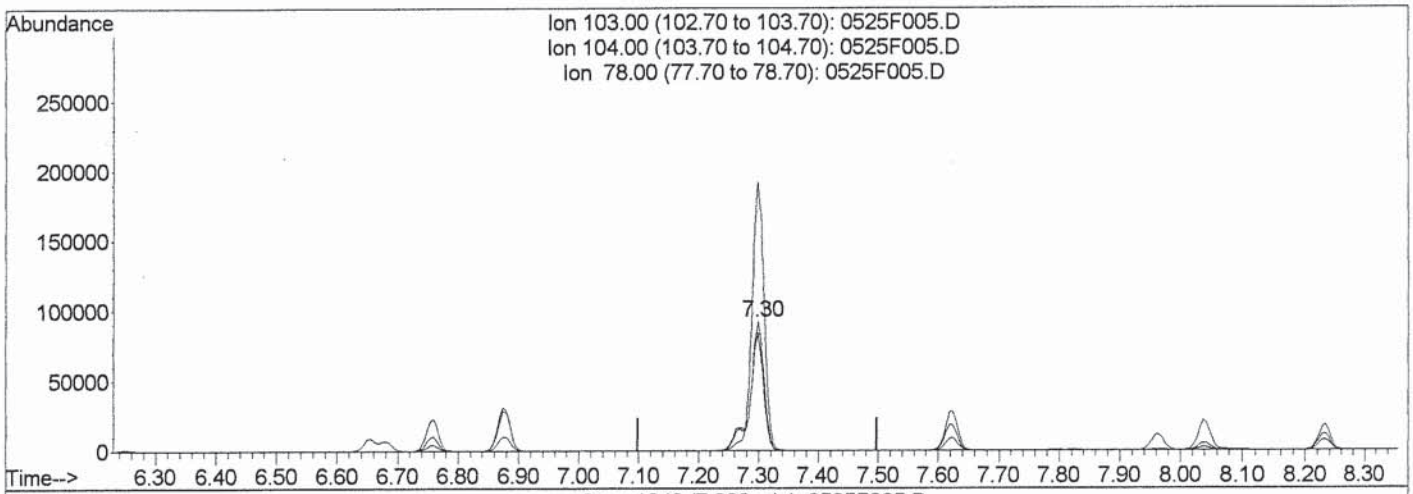
0525F005.D 042110MS24SOIL.M Tue May 25 14:52:37 2010

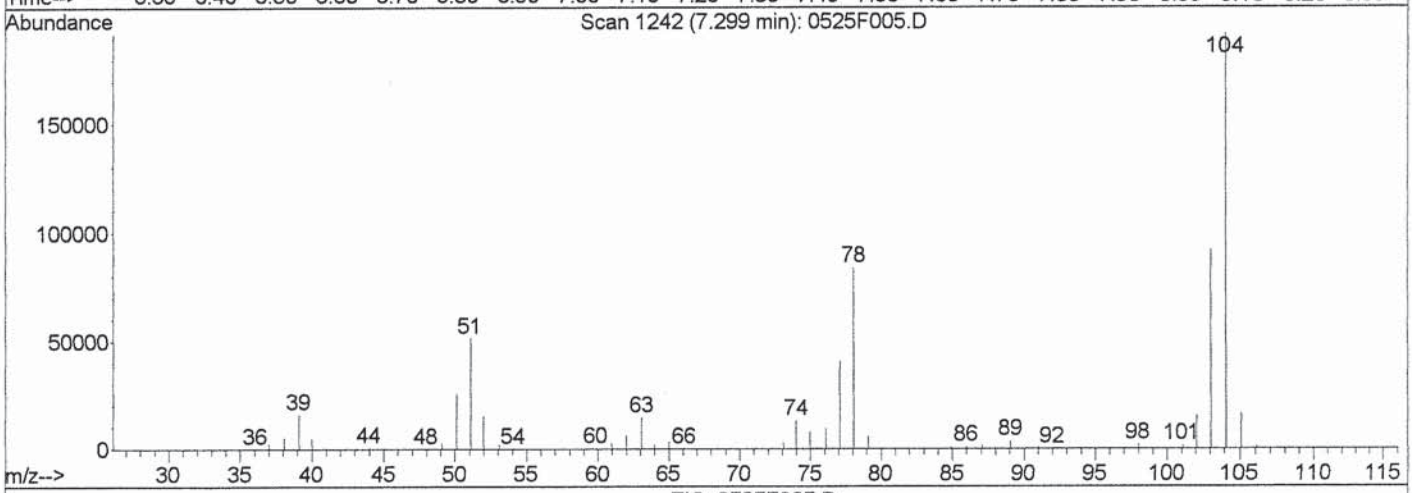
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F005.D
Acq On : 25 May 2010 11:14 am
Sample : LCS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 14:52 2010

Vial: 4
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F005.D

(75) Styrene (T)

7.30min 56.09PPB

response 146633

Ion	Exp%	Act%
103.00	100	100
104.00	208.80	208.82
78.00	88.30	91.09
0.00	0.00	0.00

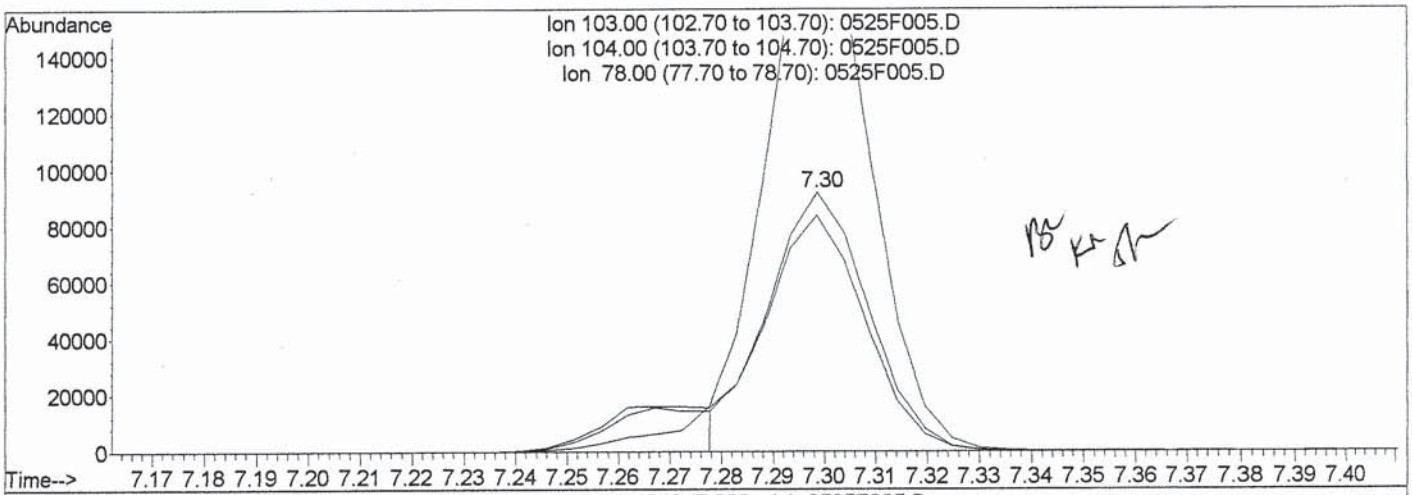
0525F005.D 042110MS24SOIL.M Tue May 25 14:52:46 2010

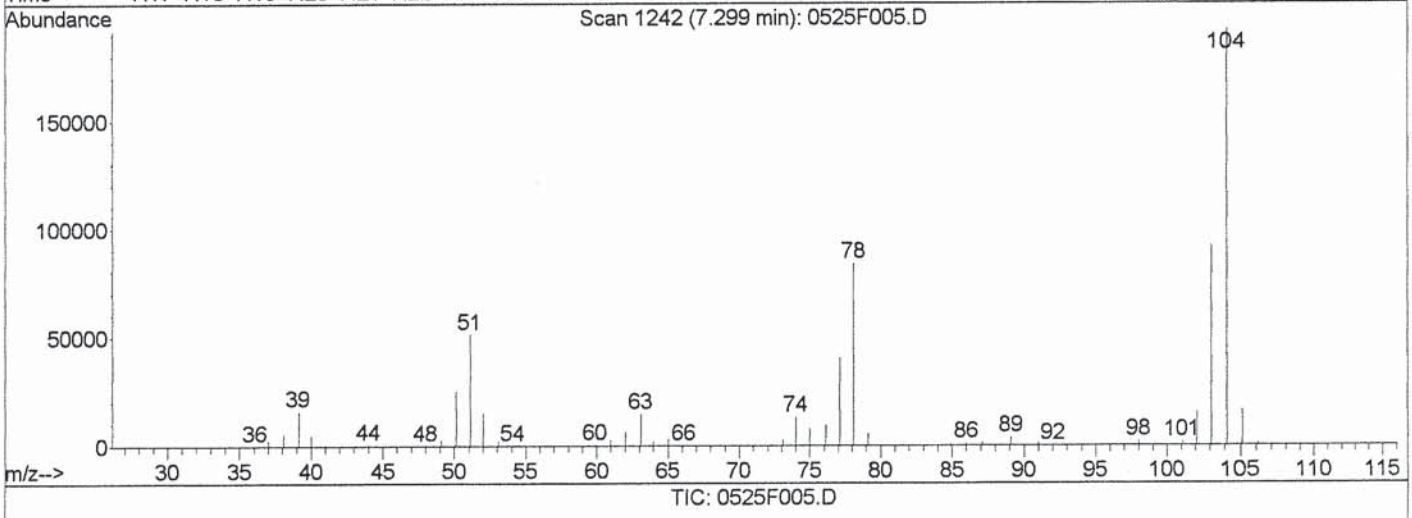
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F005.D
Acq On : 25 May 2010 11:14 am
Sample : LCS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 14:52 2010

Vial: 4
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F005.D

(75) Styrene (T)

7.30min 47.79PPB m

response 124929

Ion	Exp%	Act%
103.00	100	100
104.00	208.80	208.82
78.00	88.30	91.09
0.00	0.00	0.00

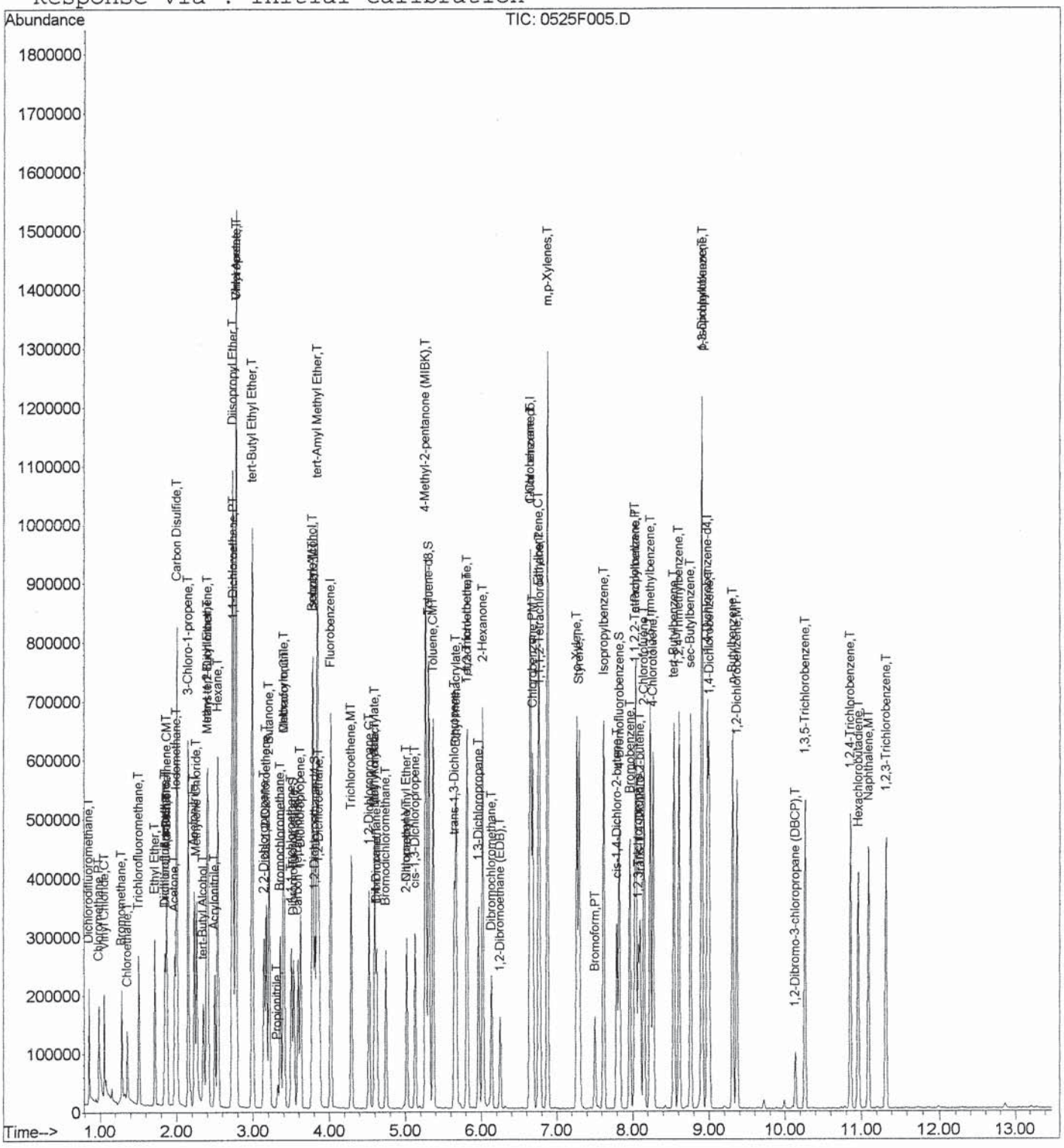
5/25/10

0525F005.D 042110MS24SOIL.M Tue May 25 14:52:55 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\052510\0525F005.D Vial: 4
Acq On : 25 May 2010 11:14 am Operator: KR
Sample : LCS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 14:52 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Initial Calibration



COLUMBIA ANALYTICAL SERVICES, INC.

Analytical Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Duplicate Lab Control Sample
Lab Code: KWG1004900-4

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dichlorodifluoromethane	**42.5**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Chloromethane	**44.8**		5.0	0.46	1	05/25/10	05/25/10	KWG1004900	
Vinyl Chloride	**45.8**		5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Bromomethane	**31.0**		5.0	0.62	1	05/25/10	05/25/10	KWG1004900	
Chloroethane	**43.1**		5.0	0.31	1	05/25/10	05/25/10	KWG1004900	
Trichlorofluoromethane	**34.7**		5.0	0.21	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethene	**41.4**		5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Acetone	**204**		20	3.9	1	05/25/10	05/25/10	KWG1004900	
Carbon Disulfide	**93.4**		5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
Methylene Chloride	**44.0**		10	0.43	1	05/25/10	05/25/10	KWG1004900	
trans-1,2-Dichloroethene	**42.8**		5.0	0.35	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloroethane	**41.9**		5.0	0.20	1	05/25/10	05/25/10	KWG1004900	
2,2-Dichloropropane	**43.9**		5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
cis-1,2-Dichloroethene	**43.0**		5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
2-Butanone (MEK)	**231**		20	0.99	1	05/25/10	05/25/10	KWG1004900	
Bromochloromethane	**45.1**		5.0	0.14	1	05/25/10	05/25/10	KWG1004900	
Chloroform	**45.0**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1-Trichloroethane (TCA)	**44.7**		5.0	0.35	1	05/25/10	05/25/10	KWG1004900	
Carbon Tetrachloride	**44.2**		5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
1,1-Dichloropropene	**42.3**		5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
Benzene	**44.2**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloroethane (EDC)	**44.2**		5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
Trichloroethene (TCE)	**46.2**		5.0	0.27	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichloropropane	**46.6**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
Dibromomethane	**45.2**		5.0	0.28	1	05/25/10	05/25/10	KWG1004900	
Bromodichloromethane	**48.0**		5.0	0.15	1	05/25/10	05/25/10	KWG1004900	
cis-1,3-Dichloropropene	**42.8**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
4-Methyl-2-pentanone (MIBK)	**253**		20	0.74	1	05/25/10	05/25/10	KWG1004900	
Toluene	**47.0**		5.0	0.18	1	05/25/10	05/25/10	KWG1004900	
trans-1,3-Dichloropropene	**53.6**		5.0	0.34	1	05/25/10	05/25/10	KWG1004900	
1,1,2-Trichloroethane	**44.5**		5.0	0.23	1	05/25/10	05/25/10	KWG1004900	
Tetrachloroethene (PCE)	**43.9**		5.0	0.26	1	05/25/10	05/25/10	KWG1004900	
2-Hexanone	**238**		20	0.93	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichloropropane	**45.5**		5.0	0.20	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601
Sample Matrix: Misc. solid

Service Request: K1005244
Date Collected: NA
Date Received: NA

Volatile Organic Compounds

Sample Name: Duplicate Lab Control Sample
Lab Code: KWG1004900-4

Extraction Method: EPA 5030A
Analysis Method: 8260B

Units: ug/Kg
Basis: Dry

Level: Low

Analyte Name	Result	Q	MRL	MDL	Dilution Factor	Date Extracted	Date Analyzed	Extraction Lot	Note
Dibromochloromethane	**48.5**		5.0	0.19	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromoethane (EDB)	**45.4**		20	0.27	1	05/25/10	05/25/10	KWG1004900	
Chlorobenzene	**44.5**		5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
Ethylbenzene	**45.8**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,1,1,2-Tetrachloroethane	**48.3**		5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
m,p-Xylenes	**96.1**		5.0	0.37	1	05/25/10	05/25/10	KWG1004900	
o-Xylene	**47.7**		5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Styrene	**48.5**		5.0	0.13	1	05/25/10	05/25/10	KWG1004900	
Bromoform	**53.6**		5.0	0.30	1	05/25/10	05/25/10	KWG1004900	
Isopropylbenzene	**48.6**		20	0.13	1	05/25/10	05/25/10	KWG1004900	
1,1,2,2-Tetrachloroethane	**44.1**		5.0	0.36	1	05/25/10	05/25/10	KWG1004900	
Bromobenzene	**46.8**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
n-Propylbenzene	**48.6**		20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichloropropane	**44.8**		5.0	0.38	1	05/25/10	05/25/10	KWG1004900	
2-Chlorotoluene	**49.0**		20	0.29	1	05/25/10	05/25/10	KWG1004900	
1,3,5-Trimethylbenzene	**49.3**		20	0.36	1	05/25/10	05/25/10	KWG1004900	
4-Chlorotoluene	**50.0**		20	0.26	1	05/25/10	05/25/10	KWG1004900	
tert-Butylbenzene	**48.6**		20	0.33	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trimethylbenzene	**48.5**		20	0.28	1	05/25/10	05/25/10	KWG1004900	
sec-Butylbenzene	**46.8**		20	0.31	1	05/25/10	05/25/10	KWG1004900	
4-Isopropyltoluene	**47.6**		20	0.28	1	05/25/10	05/25/10	KWG1004900	
1,3-Dichlorobenzene	**47.9**		5.0	0.24	1	05/25/10	05/25/10	KWG1004900	
1,4-Dichlorobenzene	**46.8**		5.0	0.25	1	05/25/10	05/25/10	KWG1004900	
n-Butylbenzene	**48.8**		20	0.32	1	05/25/10	05/25/10	KWG1004900	
1,2-Dichlorobenzene	**46.7**		5.0	0.22	1	05/25/10	05/25/10	KWG1004900	
1,2-Dibromo-3-chloropropane	**52.1**		20	0.66	1	05/25/10	05/25/10	KWG1004900	
1,2,4-Trichlorobenzene	**50.5**		20	0.23	1	05/25/10	05/25/10	KWG1004900	
Hexachlorobutadiene	**54.2**		20	0.21	1	05/25/10	05/25/10	KWG1004900	
Naphthalene	**50.9**		20	0.36	1	05/25/10	05/25/10	KWG1004900	
1,2,3-Trichlorobenzene	**52.5**		20	0.21	1	05/25/10	05/25/10	KWG1004900	

Comments:

Client:	Exponent	**Service Request:**	K1005244
Project:	Heglar - Kronquist/0907194.000.0601	**Date Collected:**	NA
Sample Matrix:	Misc. solid	**Date Received:**	NA

Volatile Organic Compounds

Sample Name:	Duplicate Lab Control Sample	**Units:**	ug/Kg
Lab Code:	KWG1004900-4	**Basis:**	Dry

Surrogate Name	%Rec	Control Limits	Date Analyzed	Note
Dibromofluoromethane	100	83-128	05/25/10	Acceptable
Toluene-d8	115	90-125	05/25/10	Acceptable
4-Bromofluorobenzene	109	77-124	05/25/10	Acceptable

Comments:

Exception Report

Data File: J:\MS24\DATA\052510\0525F006.D
Lab ID: KWG1004900-4
RunType: DLCS
Matrix: SEDIMENT

Date Acquired: 05/25/2010 11:35
Date Quantitated: 05/25/2010 14:53
Batch ID: KWG1004880
Analysis Method: 8260B
MethodJoinID: MJ120

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Window	NA	NA	NA	x	
Analytical Holding Time	NA	NA	NA	x	
ICAL Pass/Fail	NA	NA	NA	x	
ICAL Average RSD	NA	NA	NA	x	
ICAL Analyte Recovery	NA	NA	NA	x	
Initial Calibration Minimum RF	NA	NA	NA		x
Initial Calibration SPCC/CCC	NA	NA	NA	x	
Second Source ICAL Verification	NA	NA	NA	x	
Calibration Verification Pass/Fail	NA	NA	NA	x	
Continuing Calibration Recovery	NA	NA	NA	x	
Continuing Calibration Minimum RF	NA	NA	NA		x
Continuing Calibration SPCC/CCC	NA	NA	NA	x	
Internal Standards	NA	NA	NA	x	
Surrogates	NA	NA	NA	x	
Analyte Co-elution	NA	NA	NA	x	
Retention Time	NA	NA	NA	x	
Relative Retention Time	NA	NA	NA	x	
Below Lowest ICAL Level	NA	NA	NA	x	
Std MRL Unsupported by ICAL	NA	NA	NA	x	
Above Highest ICAL Level	NA	NA	NA	x	
Enviroquant/Stealth Calibration Check	NA	NA	NA	x	
Overdiluted Analysis	NA	NA	NA	x	

Analyte Exceptions

Exception Categories	Analyte Name	Result	Low Limit	High Limit	Corrective Action
Initial Calibration Minimum RF	1,4-Dioxane	0.0037	0.01	NA	
Continuing Calibration Minimum RF	1,4-Dioxane	0.0034	0.01	NA	

Primary Review: [signature]

Secondary Review: [signature] 5/25/10

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	SEDIMENT
Prod Code:	8260B VOC_FP	**Collect Date:**		**Receive Date:**	05/25/2010

Analysis Lot:	KWG1004880	**Prep Lot:**	KWG1004900	**Report Group:**	
Analysis Method:	8260B	**Prep Method:**	EPA 5030A		
Prep Ref:	911431	**Prep Date:**	05/25/2010		

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:			
Tune Ref:	J:\MS24\DATA\052510\0525F002.D	**Method ID:**	MJ120
MB Ref:	J:\MS24\DATA\052510\0525F009.D	**Quant based on Method**	

Data File:	J:\MS24\DATA\052510\0525F006.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:35	**Quant Date:**	05/25/2010 14:53	**Vial:**	5
Run Type:	DLCS			**Dilution:**	1.0
Lab ID:	KWG1004900-4			**Soln Conc. Units:**	PPB

Internal Standard Compounds

IS Ref	Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Area Criteria
1	Fluorobenzene	4.02	0.00	96	355415	50.00	OK
2	Chlorobenzene-d5	6.65	0.00	82	159438	50.00	OK
3	1,4-Dichlorobenzene-d4	8.98	0.00	152	144430	50.00	OK

Surrogate Compounds

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	%Rec	%Rec Limits		Rpt?
1	Dibromofluoromethane	3.53	-0.01	0.00	113	88941	50.07	100	43-153	OK	
1	1,2-Dichloroethane-d4	3.81	0.00	0.00	65	118957	46.39	93	64-142	OK	
1	Toluene-d8	5.31	0.00	0.00	98	357374	57.36	115	49-151	OK	
2	4-Bromofluorobenzene	7.83	0.00	0.00	95	132928	54.32	109	33-150	OK	

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	Dichlorodifluoromethane	0.85	-0.01	0.00	85	93270	42.51	42.5		
1	Chloromethane	0.99		0.00	50	124589	44.79	44.8		
1	Vinyl Chloride	1.05	-0.01	0.00	62	113073	45.82	45.8		
1	Bromomethane	1.28	-0.02	0.00	96	63578m	30.97	31.0		
1	Chloroethane	1.36	-0.01	0.00	64	69293	43.08	43.1		
1	Dichlorofluoromethane (CFC 21	1.36	-0.16	-0.04	67	501	0.1200	0.120	J	
1	Trichlorofluoromethane	1.51		0.00	101	118021	34.68	34.7		
1	Diethyl Ether	1.72	-0.01	0.00	59	89151	43.80	43.8		
1	Acrolein	1.87		0.00	56	30969	78.80	78.8	J	
1	Trichlorotrifluoroethane	1.85	-0.01	0.00	151	55378	42.58	42.6		
1	1,1-Dichloroethene	1.87	-0.01	0.00	96	61839	41.43	41.4		
1	Acetone	1.98		0.00	43	168911	204.12	204		
1	Iodomethane (Methyl Iodide)	1.99	-0.01	0.00	142	173493	85.11	85.1		
1	Carbon Disulfide	2.01		0.00	76	547516	93.36	93.4		

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D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

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#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F006.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:35	**Quant Date:**	05/25/2010 14:53	**Vial:**	5
Run Type:	DLCS			**Dilution:**	1.0
Lab ID:	KWG1004900-4			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
1	3-Chloro-1-propene	2.15	-0.01	0.00	76	79396	74.38	74.4		
1	Acetonitrile	2.23		0.00	40	133076	818.90	819		
1	Methylene Chloride	2.26	-0.01	0.00	84	94029	43.95	44.0		
1	tert-Butyl Alcohol	2.35		0.00	59	123186	415.09	415		
1	Acrylonitrile	2.50		0.00	53	106159	128.35	128		
1	Methyl tert-Butyl Ether	2.40		0.00	73	302646	48.38	48.4		
1	trans-1,2-Dichloroethene	2.41	-0.01	0.00	96	79888	42.80	42.8		
1	n-Hexane	2.54		0.00	57	177690	69.45	69.5		
1	Diisopropyl Ether	2.73	-0.01	0.00	45	596830	84.87	84.9		
1	1,1-Dichloroethane	2.75		0.00	63	154157	41.86	41.9		
1	Vinyl Acetate	2.78		0.00	86	71020	207.26	207		
1	Chloroprene	2.78		0.00	53	216092	86.27	86.3		
1	tert-Butyl Ethyl Ether	2.99		0.00	59	571111	89.46	89.5		
1	2,2-Dichloropropane	3.15		0.00	77	110919	43.94	43.9		
1	cis-1,2-Dichloroethene	3.18		0.00	96	88963	42.95	43.0		
1	2-Butanone (MEK)	3.21		0.00	72	61376	230.57	231		
1	Propionitrile	3.32		0.00	54	24202	73.42	73.4		
1	Methacrylonitrile	3.40		0.00	67	67590	77.21	77.2		
1	Bromochloromethane	3.36		0.00	128	45769	45.08	45.1		
1	Chloroform	3.41		0.00	83	156269	44.97	45.0		
1	1,1,1-Trichloroethane (TCA)	3.50	-0.01	0.00	97	118480	44.65	44.7		
1	Carbon Tetrachloride	3.59		0.00	117	100789	44.17	44.2		
1	1,1-Dichloropropene	3.62	-0.01	0.00	75	100924	42.33	42.3		
1	Isobutanol	3.78		0.00	43	83188	765.23	765		
1	Benzene	3.79		0.00	78	334954	44.24	44.2		
1	1,2-Dichloroethane (EDC)	3.87		0.00	62	135823	44.21	44.2		
1	tert-Amyl Methyl Ether	3.84		0.00	55	148350	90.42	90.4		
1	Trichloroethene (TCE)	4.29		0.00	95	92543	46.23	46.2		
1	1,2-Dichloropropane	4.53		0.00	63	93960	46.59	46.6		
1	Dibromomethane	4.62		0.00	93	57668	45.24	45.2		
1	Methyl Methacrylate	4.60		0.00	69	111290	78.84	78.8		
1	1,4-Dioxane	4.62		0.00	88	20712	794.23	794		
1	Bromodichloromethane	4.75		0.00	83	121996	47.98	48.0		
1	2-Nitropropane	5.02		0.00	41	42024	79.54	79.5		
1	2-Chloroethyl Vinyl Ether	5.02		0.00	63	64550	51.77	51.8		
1	cis-1,3-Dichloropropene	5.13		0.00	75	121428	42.78	42.8		
1	4-Methyl-2-pentanone (MIBK)	5.27		0.00	58	196132	253.23	253		
1	Toluene	5.37		0.00	92	213725	46.99	47.0		
2	trans-1,3-Dichloropropene	5.65		0.00	75	141549	53.64	53.6		
2	Ethyl Methacrylate	5.67		0.00	69	207312	82.59	82.6		
2	1,1,2-Trichloroethane	5.81		0.00	83	73440	44.47	44.5		
2	Tetrachloroethene (PCE)	5.82		0.00	164	68729	43.94	43.9		

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d: Compound manually deleted
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#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F006.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:35	**Quant Date:**	05/25/2010 14:53	**Vial:**	5
Run Type:	DLCS			**Dilution:**	1.0
Lab ID:	KWG1004900-4			**Soln Conc. Units:**	PPB

Target Compounds **Final Conc. Units:** ug/Kg Wet Weight

IS Ref	Parameter Name	RT	RT Dev	RRT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2	2-Hexanone	6.02		0.00	57	66464	238.41	238		
2	1,3-Dichloropropane	5.97		0.00	76	149682	45.51	45.5		
2	Dibromochloromethane	6.13		0.00	129	89651	48.54	48.5		
2	1,2-Dibromoethane (EDB)	6.25		0.00	107	81903	45.40	45.4		
2	1-Chlorohexane	6.64		0.00	91	109032	48.09	48.1		
2	Chlorobenzene	6.68		0.00	112	244714	44.46	44.5		
2	Ethylbenzene	6.76		0.00	106	122742	45.79	45.8		
2	1,1,1,2-Tetrachloroethane	6.77		0.00	131	87642	48.26	48.3		
2	m,p-Xylenes	6.87		0.00	106	303682	96.13	96.1		
2	o-Xylene	7.27		0.00	106	151400	47.74	47.7		
2	Styrene	7.30		0.00	103	120521**m**	48.46	48.5		
2	Bromoform	7.51		0.00	173	60095	53.57	53.6		
2	Isopropylbenzene	7.62		0.00	105	372430	48.62	48.6		
2	cis-1,4-Dichloro-2-butene	7.80		0.00	89	23418	108.43	108		
3	1,1,2,2-Tetrachloroethane	8.03		0.00	83	92728	44.09	44.1		
3	trans-1,4-Dichloro-2-butene	8.10		0.00	53	57395	115.30	115		
3	Bromobenzene	7.96		0.00	156	106106	46.79	46.8		
3	n-Propylbenzene	8.04		0.00	91	449946	48.63	48.6		
3	1,2,3-Trichloropropane	8.07		0.00	110	35782	44.78	44.8		
3	2-Chlorotoluene	8.15		0.00	91	284295	49.01	49.0		
3	1,3,5-Trimethylbenzene	8.23		0.00	105	320553	49.31	49.3		
3	4-Chlorotoluene	8.27		0.00	91	292877	50.00	50.0		
3	tert-Butylbenzene	8.54		0.00	119	276985	48.64	48.6		
3	1,2,4-Trimethylbenzene	8.60		0.00	105	322331	48.48	48.5		
3	sec-Butylbenzene	8.76		0.00	105	396758	46.75	46.8		
3	4-Isopropyltoluene	8.91		0.00	119	336922	47.59	47.6		
3	1,3-Dichlorobenzene	8.90		0.00	146	200190	47.87	47.9		
3	1,4-Dichlorobenzene	9.00		0.00	146	202634	46.78	46.8		
3	n-Butylbenzene	9.31		0.00	91	305408	48.83	48.8		
3	1,2-Dichlorobenzene	9.36		0.00	146	191871	46.68	46.7		
3	1,2-Dibromo-3-chloropropane	10.14		0.00	155	16371	52.11	52.1		
3	1,2,4-Trichlorobenzene	10.86		0.00	180	131439	50.51	50.5		
3	Hexachlorobutadiene	10.96		0.00	225	70504	54.15	54.2		
3	Naphthalene	11.09		0.00	128	304749	50.93	50.9		
3	1,2,3-Trichlorobenzene	11.32		0.00	180	127235	52.51	52.5		
	4-Vinyl-1-cyclohexene				0	0		5.0	UJ	NR
	Benzyl Chloride				0	0		5.0	UJ	NR
	Hexachloroethane				0	0		1.0	UJ	NR
	2,2,4-Trimethylpentane				0	0		5.0	UJ	NR
	1,3-Butadiene				0	0		2.0	UJ	NR
	1,1,2-Trifluoroethane				0	0		10	UJ	NR
	n-Butyl Alcohol				0	0		250	U	NR

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d: Compound manually deleted
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?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File:	J:\MS24\DATA\052510\0525F006.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 11:35	**Quant Date:**	05/25/2010 14:53	**Vial:**	5
Run Type:	DLCS			**Dilution:**	1.0
Lab ID:	KWG1004900-4			**Soln Conc. Units:**	PPB

Target Compounds

Final Conc. Units: ug/Kg Wet Weight

Parameter Name	RT	RT Dev	Quant Mass	Response	Solution Conc	Final Conc	Q	Rpt?
2-Propanol			0	0		200	U	NR
Ethyl Alcohol			0	0		200	U	NR
2,2-Dichloro-1,1,1-trifluoroetha			0	0		5.0	UJ	NR
Cyclohexanone			0	0		10	UJ	NR
2-Methylpentane			0	0		5.0	UJ	NR
n-Heptane			0	0		5.0	UJ	NR
3-Methylpentane			0	0		5.0	UJ	NR
Methylcyclopentane			0	0		5.0	UJ	NR
Ethyl Acetate			0	0		20	UJ	NR

Prep Amount:	5.00 g	**Dilution:**	1.0
Prep Final Vol:	5.0 ml	**Unit Factor:**	1
Solids:	%		

Final Concentration = ((Soln Conc x Prep Final Vol x Dilution) / (Prep Amount x Solids)) x Unit Factor

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E: Analyte concentration above high point of ICAL
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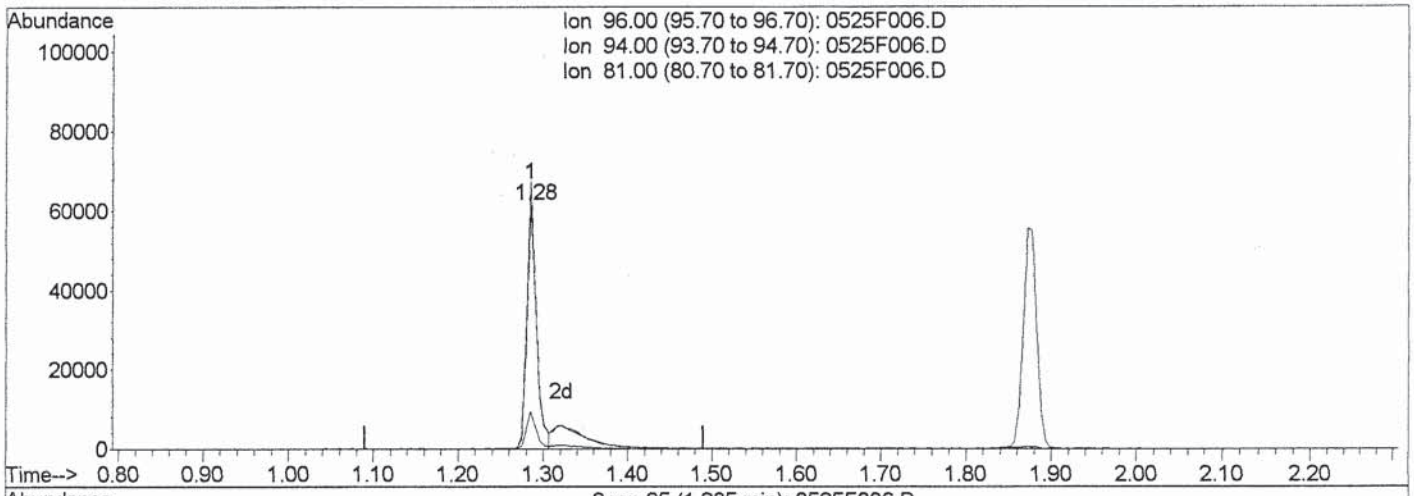
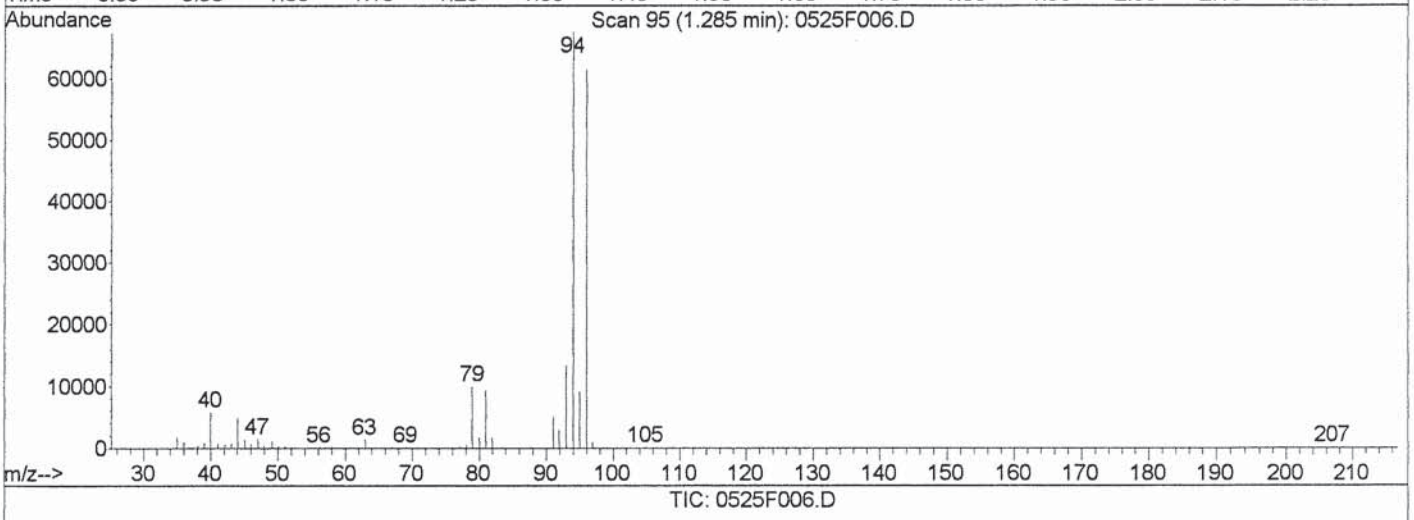
*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F006.D Vial: 5
Acq On : 25 May 2010 11:35 am Operator: KR
Sample : DLCS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 14:53 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F006.D

(5) Bromomethane (T)

1.28min 23.33PPB

response 47899

Ion	Exp%	Act%
96.00	100	100
94.00	106.40	109.46
81.00	18.10	14.79
0.00	0.00	0.00

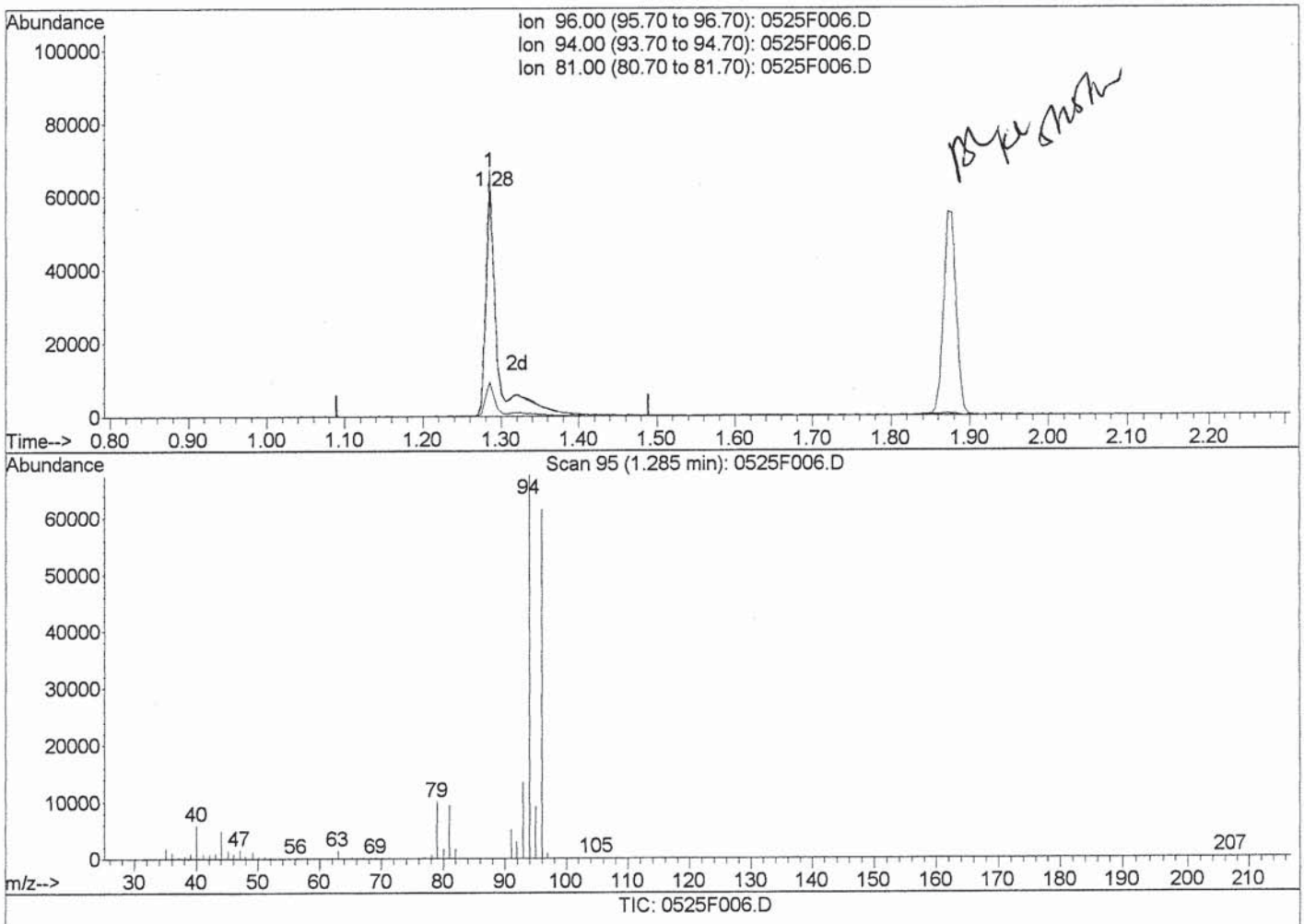
0525F006.D 042110MS24SOIL.M Tue May 25 14:53:29 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F006.D
Acq On : 25 May 2010 11:35 am
Sample : DLCS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 14:53 2010

Vial: 5
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration



TIC: 0525F006.D

(5) Bromomethane (T)

1.28min 30.97PPB m

response 63578

Ion	Exp%	Act%
96.00	100	100
94.00	106.40	109.74
81.00	18.10	15.27
0.00	0.00	0.00

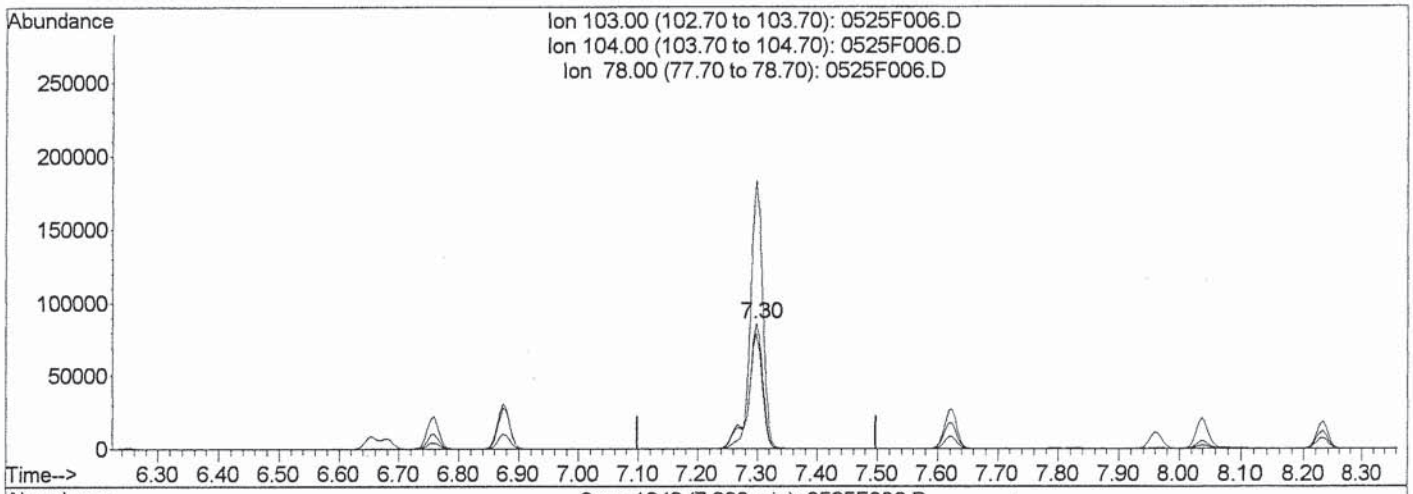
Anne 5/25/10

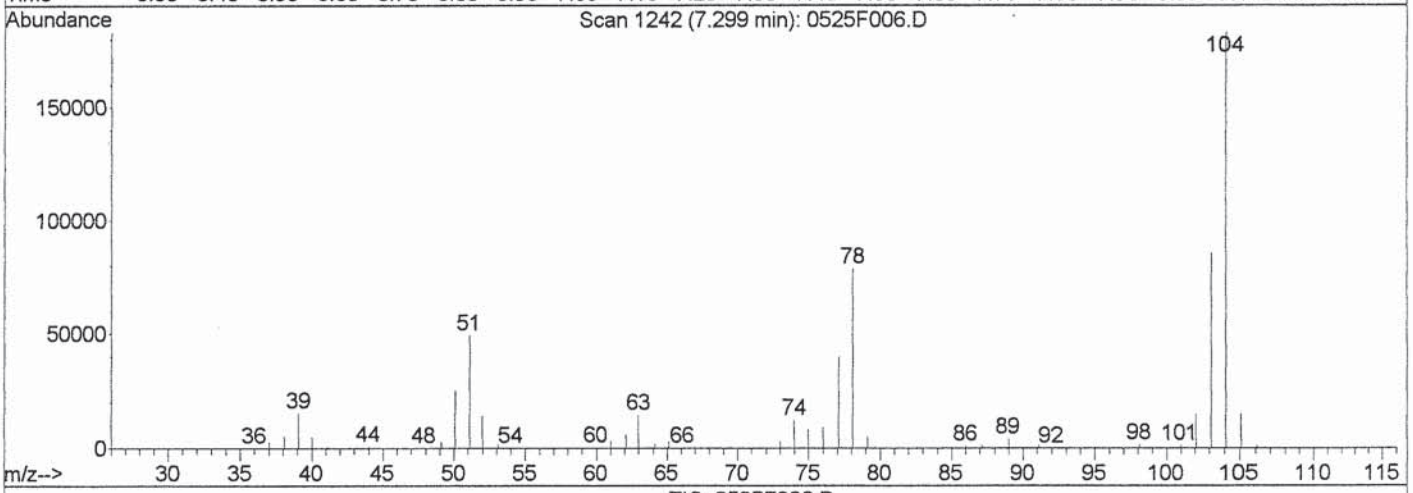
0525F006.D 042110MS24SOIL.M Tue May 25 14:53:35 2010

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F006.D Vial: 5
Acq On : 25 May 2010 11:35 am Operator: KR
Sample : DLCS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 14:53 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F006.D

(75) Styrene (T)

7.30min 56.95PPB

response 141646

Ion	Exp%	Act%
103.00	100	100
104.00	208.80	213.79
78.00	88.30	92.06
0.00	0.00	0.00

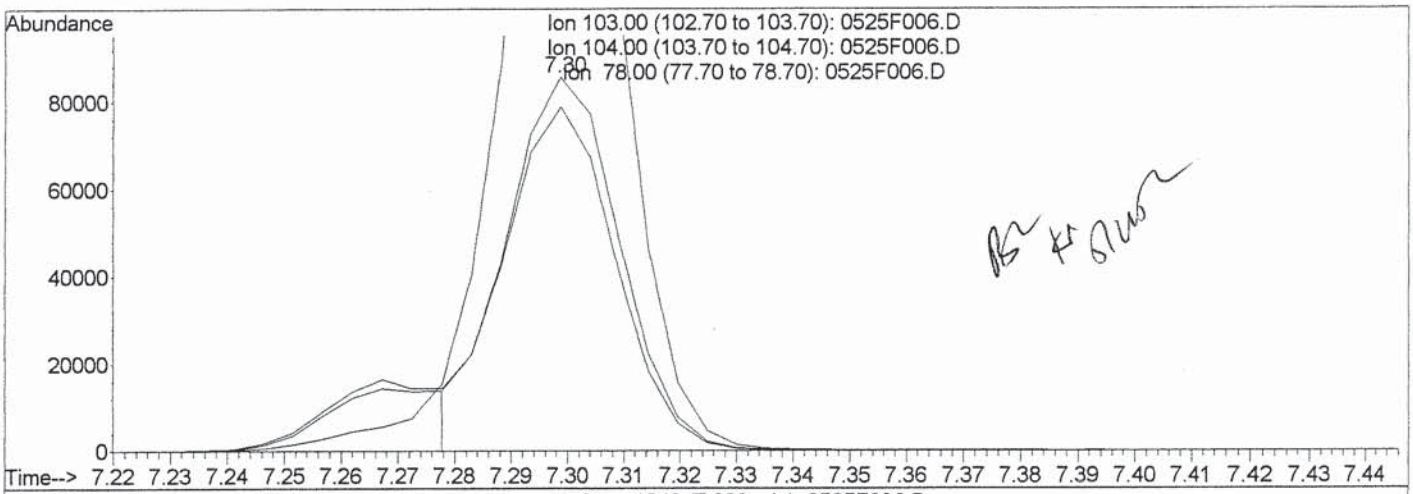
0525F006.D 042110MS24SOIL.M Tue May 25 14:53:45 2010

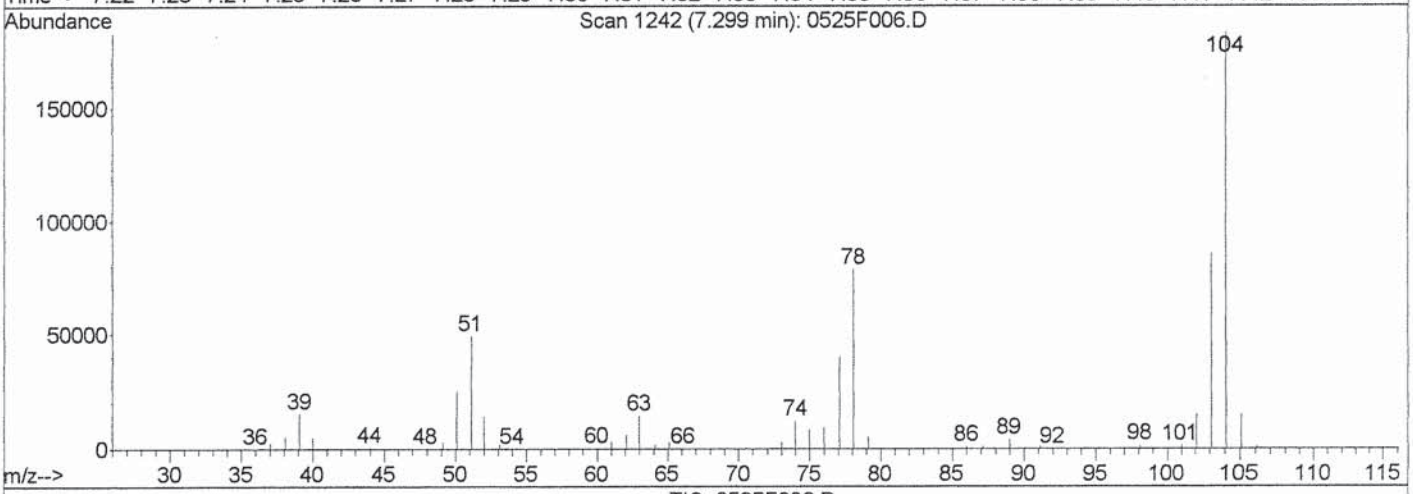
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\052510\0525F006.D
Acq On : 25 May 2010 11:35 am
Sample : DLCS
Misc :
MS Integration Params: rteint.p
Quant Time: May 25 14:53 2010

Vial: 5
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Single Level Calibration





TIC: 0525F006.D

(75) Styrene (T)

7.30min 48.46PPB m

response 120521

Ion	Exp%	Act%
103.00	100	100
104.00	208.80	213.79
78.00	88.30	92.06
0.00	0.00	0.00

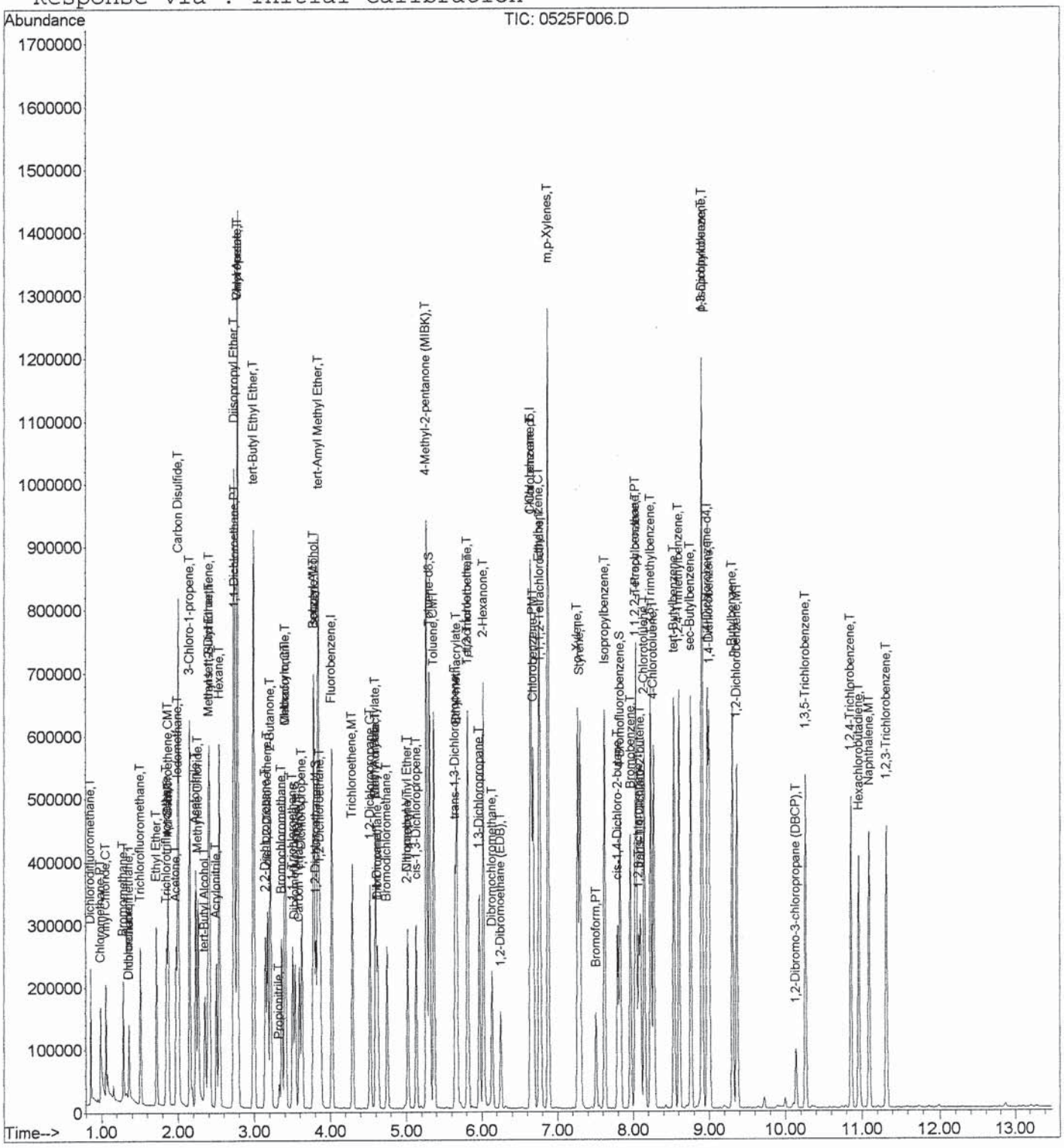
Conc 5/25/10

0525F006.D 042110MS24SOIL.M Tue May 25 14:53:51 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\052510\0525F006.D Vial: 5
Acq On : 25 May 2010 11:35 am Operator: KR
Sample : DLCS Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 25 14:53 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Mon May 10 10:34:02 2010
Response via : Initial Calibration



Organic Analysis:
Volatile Organic Compounds

Validation Package

Standards Data

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Date Analyzed: 05/25/2010
Time Analyzed: 08:42

Tune Summary
Volatile Organic Compounds

File ID: J:\MS24\DATA\052510\0525F002.D
Instrument ID: MS24
Column:

Analysis Method: 8260B
Analysis Lot: KWG1004880

Target Mass	Relative to Mass	Lower Limit%	Upper Limit%	Relative Abundance %	Raw Abundance	Result Pass/Fail
50	95	15	40	24.2	3284	PASS
75	95	30	60	55.9	7570	PASS
95	95	100	100	100.0	13552	PASS
96	95	5	9	8.7	1182	PASS
173	174	0	2	1.2	123	PASS
174	95	50	120	77.4	10491	PASS
175	174	5	9	8.4	877	PASS
176	174	95	101	97.2	10193	PASS
177	176	5	9	6.1	618	PASS

Sample Name	Lab Code	File ID	Date Analyzed	Time Analyzed	Q
Continuing Calibration Verification	KWG1004880-2	J:\MS24\DATA\052510\0525F004.D	05/25/2010	10:34	
Lab Control Sample	KWG1004900-3	J:\MS24\DATA\052510\0525F005.D	05/25/2010	11:14	
Duplicate Lab Control Sample	KWG1004900-4	J:\MS24\DATA\052510\0525F006.D	05/25/2010	11:35	
Method Blank	KWG1004900-5	J:\MS24\DATA\052510\0525F009.D	05/25/2010	12:39	
D-4-16	K1005244-003	J:\MS24\DATA\052510\0525F012.D	05/25/2010	13:42	
D-4-16MS	KWG1004900-1	J:\MS24\DATA\052510\0525F013.D	05/25/2010	14:40	
D-4-16DMS	KWG1004900-2	J:\MS24\DATA\052510\0525F014.D	05/25/2010	15:01	

Results flagged with an asterisk (*) indicate the analysis performed outside specified tune window

Exception Report

Data File: J:\MS24\DATA\052510\0525F002.D
Lab ID: KWG1004880-1
RunType: BFB
Matrix: SOIL

Date Acquired: 05/25/2010 08:42
Date Quantitated:
Batch ID: KWG1004880
Analysis Method: BFB
ListJoinID: LJ774

Sample Exceptions

Exception Categories	Result	Low Limit	High Limit	Pass	Fail
Tune Ion Ratio	NA	NA	NA	x	

Primary Review: KL 5/28/10

Secondary Review: Cmm 5/25/10

Printed: 05/25/2010 10:10:19
u:\Stealth\Crystal.rpt\except2.rpt

Quantitation Report

Bottle ID:		**Tier:**		**Matrix:**	SOIL
Prod Code:	8260B	**Collect Date:**		**Receive Date:**	05/25/2010

Analysis Lot:	KWG1004880	**Prep Lot:**		**Report Group:**	
Analysis Method:	BFB	**Prep Method:**			
Prep Ref:		**Prep Date:**			

Quant Method:	J:\MS24\METHODS\042110MS24SO	**Calibration ID:**	CAL9404
Title:	GC/MS Tuning Evaluation	**Report List ID:**	LJ774
Tune Ref:		**Method ID:**	MJ159
MB Ref:		**Quant based on Report List**	

Data File:	J:\MS24\DATA\052510\0525F002.D			**Instrument:**	MS24
Acqu Date:	05/25/2010 08:42	**Quant Date:**		**Vial:**	2
Run Type:	BFB			**Dilution:**	1.0
Lab ID:	KWG1004880-1			**Soln Conc. Units:**	

Tune Results

Target Mass	Relative to Mass	Lower Limit%	Upper Limit%	Relative Abundance %	Raw Abundance	Result Pass/Fail
50	95	15	40	24.2	3284	Pass
75	95	30	60	55.9	7570	Pass
95	95	100	100	100.0	13552	Pass
96	95	5	9	8.7	1182	Pass
173	174	0	2	1.2	123	Pass
174	95	50	120	77.4	10491	Pass
175	174	5	9	8.4	877	Pass
176	174	95	101	97.2	10193	Pass
177	176	5	9	6.1	618	Pass

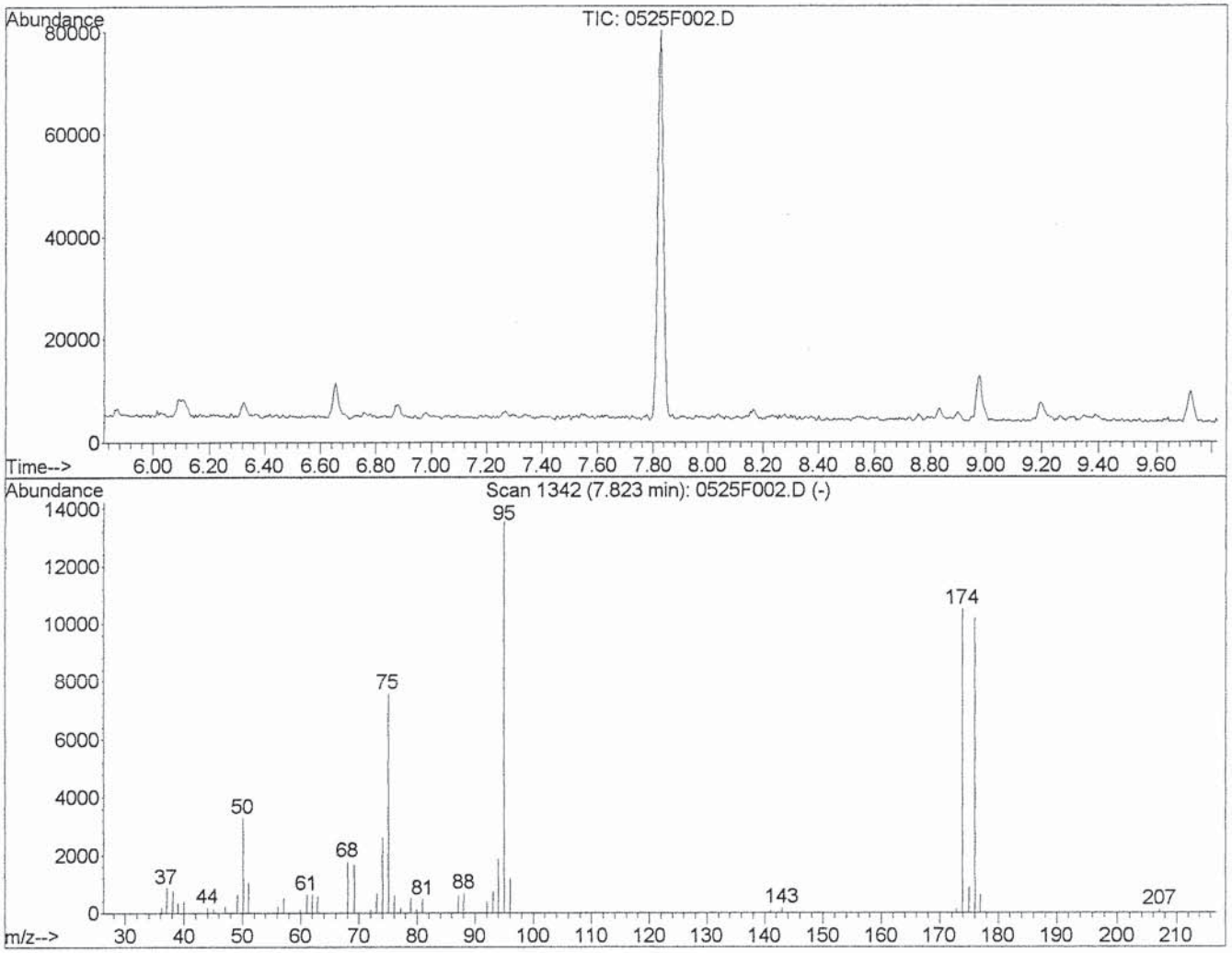
U: Undetected at or above MDL
J: Analyte detected above MDL, but below MRL
B: Hit above MRL also found in Method Blank
E: Analyte concentration above high point of ICAL
N: Presumptive evidence of compound

D: Result from dilution
m: Manual integration performed
d: Compound manually deleted
NR: Analyte not reported from this analysis

*: Result fails acceptance criteria
#: Acceptance criteria not applicable
?: Insufficient information to determine acceptance
e: Result >= MRL, but MRL less than low point of ICAL
c: check for co-elution

Data File : J:\MS24\DATA\052510\0525F002.D
Acq On : 25 May 2010 8:42 am
Sample : BFB NEW TUNE
Misc :
MS Integration Params: rteint.p
Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B

Vial: 2
Operator: KR
Inst : MS24
Multiplr: 1.00



Spectrum Information: Scan 1342 Head of apex - 1335 Scan KR 5/26/10

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	24.2	3284	PASS
75	95	30	60	55.9	7570	PASS
95	95	100	100	100.0	13552	PASS
96	95	5	9	8.7	1182	PASS
173	174	0.00	2	1.2	123	PASS
174	95	50	120	77.4	10491	PASS
175	174	5	9	8.4	877	PASS
176	174	95	101	97.2	10193	PASS
177	176	5	9	6.1	618	PASS

0525F002.D 042110MS24SOIL.M Tue May 25 09:43:47 2010

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Level ID	File ID
A	J:\MS24\DATA\042110\0421F005.D
B	J:\MS24\DATA\042110\0421F006.D
C	J:\MS24\DATA\042110\0421F007.D
D	J:\MS24\DATA\042110\0421F008.D
E	J:\MS24\DATA\042110\0421F009.D
F	J:\MS24\DATA\042110\0421F010.D
G	J:\MS24\DATA\042110\0421F011.D
H	J:\MS24\DATA\042110\0421F012.D
I	J:\MS24\DATA\042110\0421F013.D

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
Dichlorodifluoromethane				B	2.0	0.330	C	5.0	0.312	D	10	0.299	E	20	0.321
	F	50	0.282	G	100	0.308	H	200	0.331	I	300	0.286			
† Chloromethane	A	1.0	0.506	B	2.0	0.397	C	5.0	0.378	D	10	0.369	E	20	0.379
†	F	50	0.350	G	100	0.385	H	200	0.404	I	300	0.354			
‡ Vinyl Chloride				B	2.0	0.364	C	5.0	0.349	D	10	0.351	E	20	0.360
‡	F	50	0.319	G	100	0.350	H	200	0.366	I	300	0.318			
Bromomethane							C	5.0	0.324	D	10	0.293	E	20	0.297
	F	50	0.233	G	100	0.281	H	200	0.304						
Chloroethane							C	5.0	0.228	D	10	0.221	E	20	0.239
	F	50	0.209	G	100	0.231	H	200	0.242	I	300	0.214			
Trichlorofluoromethane				B	2.0	0.487	C	5.0	0.472	D	10	0.476	E	20	0.506
	F	50	0.431	G	100	0.481	H	200	0.518	I	300	0.458			
‡ 1,1-Dichloroethene				B	2.0	0.244	C	5.0	0.199	D	10	0.203	E	20	0.219
‡	F	50	0.191	G	100	0.205	H	200	0.225	I	300	0.193			
Acetone										D	20	0.150	E	40	0.118
	F	100	0.108	G	200	0.107	H	400	0.115	I	600	0.101			
Carbon Disulfide	A	1.0	0.905	B	2.0	0.781	C	5.0	0.761	D	10	0.786	E	20	0.855
	F	50	0.742	G	100	0.847	H	200	0.923	I	300	0.826			
Methylene Chloride							C	5.0	0.322	D	10	0.295	E	20	0.319
	F	50	0.283	G	100	0.299	H	200	0.310	I	300	0.279			
trans-1,2-Dichloroethene	A	1.0	0.310	B	2.0	0.263	C	5.0	0.250	D	10	0.245	E	20	0.261
	F	50	0.236	G	100	0.261	H	200	0.282	I	300	0.254			
† 1,1-Dichloroethane	A	1.0	0.555	B	2.0	0.546	C	5.0	0.477	D	10	0.496	E	20	0.523
†	F	50	0.475	G	100	0.529	H	200	0.566	I	300	0.496			
2,2-Dichloropropane							C	5.0	0.339	D	10	0.338	E	20	0.355
	F	50	0.333	G	100	0.370	H	200	0.404	I	300	0.347			
cis-1,2-Dichloroethene	A	1.0	0.307	B	2.0	0.290	C	5.0	0.271	D	10	0.275	E	20	0.297
	F	50	0.285	G	100	0.301	H	200	0.318	I	300	0.278			
2-Butanone (MEK)							C	10	0.0436	D	20	0.0374	E	40	0.0364
	F	100	0.0352	G	200	0.0360	H	400	0.0396	I	600	0.0339			

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
Bromochloromethane							C	5.0	0.135	D	10	0.136	E	20	0.144
	F	50	0.138	G	100	0.149	H	200	0.158	I	300	0.140			
‡ Chloroform				B	2.0	0.483	C	5.0	0.452	D	10	0.465	E	20	0.498
‡	F	50	0.458	G	100	0.512	H	200	0.550	I	300	0.493			
1,1,1-Trichloroethane (TCA)				B	2.0	0.361	C	5.0	0.360	D	10	0.359	E	20	0.390
	F	50	0.343	G	100	0.387	H	200	0.414	I	300	0.373			
Carbon Tetrachloride							C	5.0	0.302	D	10	0.308	E	20	0.337
	F	50	0.287	G	100	0.327	H	200	0.362	I	300	0.324			
1,1-Dichloropropene	A	1.0	0.359	B	2.0	0.334	C	5.0	0.302	D	10	0.327	E	20	0.354
	F	50	0.299	G	100	0.339	H	200	0.371	I	300	0.334			
Benzene	A	1.0	1.10	B	2.0	1.02	C	5.0	0.979	D	10	1.01	E	20	1.13
	F	50	0.998	G	100	1.09	H	200	1.19	I	300	1.07			
1,2-Dichloroethane (EDC)	A	1.0	0.435	B	2.0	0.416	C	5.0	0.412	D	10	0.415	E	20	0.451
	F	50	0.415	G	100	0.442	H	200	0.472	I	300	0.431			
Trichloroethene (TCE)	A	1.0	0.304	B	2.0	0.298	C	5.0	0.265	D	10	0.257	E	20	0.300
	F	50	0.260	G	100	0.275	H	200	0.297	I	300	0.277			
‡ 1,2-Dichloropropane	A	1.0	0.315	B	2.0	0.283	C	5.0	0.254	D	10	0.252	E	20	0.295
‡	F	50	0.270	G	100	0.294	H	200	0.304	I	300	0.287			
Dibromomethane							C	5.0	0.173	D	10	0.165	E	20	0.190
	F	50	0.174	G	100	0.186	H	200	0.189	I	300	0.179			
Bromodichloromethane							C	5.0	0.312	D	10	0.310	E	20	0.368
	F	50	0.345	G	100	0.384	H	200	0.400	I	300	0.385			
cis-1,3-Dichloropropene							C	5.0	0.325	D	10	0.328	E	20	0.408
	F	50	0.384	G	100	0.445	H	200	0.475	I	300	0.431			
4-Methyl-2-pentanone (MIBK)							C	10	0.103	D	20	0.0913	E	40	0.108
	F	100	0.108	G	200	0.118	H	400	0.123	I	600	0.110			
‡ Toluene				B	2.0	0.664	C	5.0	0.579	D	10	0.599	E	20	0.693
‡	F	50	0.601	G	100	0.665	H	200	0.694	I	300	0.624			
trans-1,3-Dichloropropene							C	5.0	0.661	D	10	0.684	E	20	0.877
	F	50	0.809	G	100	0.868	H	200	0.964	I	300	0.932			
1,1,2-Trichloroethane	A	1.0	0.536	B	2.0	0.603	C	5.0	0.510	D	10	0.473	E	20	0.558
	F	50	0.482	G	100	0.491	H	200	0.522	I	300	0.485			
Tetrachloroethene (PCE)	A	1.0	0.506	B	2.0	0.541	C	5.0	0.473	D	10	0.481	E	20	0.537
	F	50	0.448	G	100	0.459	H	200	0.506	I	300	0.463			
2-Hexanone				B	4.0	0.0963	C	10	0.0808	D	20	0.0774	E	40	0.0846
	F	100	0.0845	G	200	0.0884	H	400	0.0973	I	600	0.0901			

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
1,3-Dichloropropane	A	1.0	0.986	B	2.0	1.19	C	5.0	0.953	D	10	0.965	E	20	1.13
	F	50	0.984	G	100	1.02	H	200	1.06	I	300	0.993			
Dibromochloromethane							C	5.0	0.499	D	10	0.485	E	20	0.579
	F	50	0.554	G	100	0.620	H	200	0.674	I	300	0.644			
1,2-Dibromoethane (EDB)				B	2.0	0.622	C	5.0	0.506	D	10	0.527	E	20	0.591
	F	50	0.544	G	100	0.562	H	200	0.613	I	300	0.561			
† Chlorobenzene	A	1.0	1.72	B	2.0	1.94	C	5.0	1.61	D	10	1.62	E	20	1.81
†	F	50	1.63	G	100	1.73	H	200	1.82	I	300	1.65			
‡ Ethylbenzene				B	2.0	0.821	C	5.0	0.768	D	10	0.802	E	20	0.903
‡	F	50	0.795	G	100	0.845	H	200	0.936	I	300	0.855			
1,1,1,2-Tetrachloroethane							C	5.0	0.508	D	10	0.502	E	20	0.583
	F	50	0.545	G	100	0.582	H	200	0.663	I	300	0.603			
m,p-Xylenes	A	2.0	0.859	B	4.0	0.949	C	10	0.876	D	20	0.925	E	40	1.07
	F	100	0.985	G	200	1.06	H	400	1.15	I	600	1.04			
o-Xylene							C	5.0	0.883	D	10	0.912	E	20	1.03
	F	50	0.948	G	100	1.06	H	200	1.10	I	300	1.02			
Styrene							C	5.0	0.662	D	10	0.652	E	20	0.804
	F	50	0.776	G	100	0.842	H	200	0.887	I	300	0.837			
† Bromoform							C	5.0	0.286	D	10	0.283	E	20	0.339
†	F	50	0.354	G	100	0.381	H	200	0.417	I	300	0.403			
Isopropylbenzene							C	5.0	2.05	D	10	2.22	E	20	2.52
	F	50	2.46	G	100	2.48	H	200	2.68	I	300	2.40			
† 1,1,2,2-Tetrachloroethane				B	2.0	0.939	C	5.0	0.706	D	10	0.625	E	20	0.735
†	F	50	0.684	G	100	0.692	H	200	0.754	I	300	0.688			
Bromobenzene	A	1.0	0.742	B	2.0	0.959	C	5.0	0.715	D	10	0.664	E	20	0.825
	F	50	0.763	G	100	0.780	H	200	0.839	I	300	0.779			
n-Propylbenzene							C	5.0	2.69	D	10	2.78	E	20	3.44
	F	50	3.22	G	100	3.31	H	200	3.67	I	300	3.32			
1,2,3-Trichloropropane	A	1.0	0.278	B	2.0	0.354	C	5.0	0.305	D	10	0.249	E	20	0.277
	F	50	0.257	G	100	0.250	H	200	0.271	I	300	0.249			
2-Chlorotoluene				B	2.0	2.18	C	5.0	1.78	D	10	1.77	E	20	2.16
	F	50	1.98	G	100	2.01	H	200	2.18	I	300	2.01			
1,3,5-Trimethylbenzene							C	5.0	1.94	D	10	1.89	E	20	2.42
	F	50	2.27	G	100	2.34	H	200	2.55	I	300	2.34			
4-Chlorotoluene				B	2.0	2.23	C	5.0	1.81	D	10	1.76	E	20	2.13
	F	50	2.02	G	100	2.08	H	200	2.21	I	300	1.99			

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF	Level ID	Amt	RRF
tert-Butylbenzene							C	5.0	1.75	D	10	1.68	E	20	2.09
	F	50	1.96	G	100	2.06	H	200	2.20	I	300	2.06			
1,2,4-Trimethylbenzene							C	5.0	1.95	D	10	1.86	E	20	2.47
	F	50	2.31	G	100	2.45	H	200	2.60	I	300	2.47			
sec-Butylbenzene							C	5.0	2.47	D	10	2.58	E	20	3.24
	F	50	2.87	G	100	3.08	H	200	3.26	I	300	3.06			
4-Isopropyltoluene							C	5.0	2.07	D	10	2.16	E	20	2.69
	F	50	2.40	G	100	2.52	H	200	2.74	I	300	2.57			
1,3-Dichlorobenzene	A	1.0	1.28	B	2.0	1.69	C	5.0	1.38	D	10	1.31	E	20	1.55
	F	50	1.41	G	100	1.43	H	200	1.53	I	300	1.45			
1,4-Dichlorobenzene	A	1.0	1.49	B	2.0	1.70	C	5.0	1.45	D	10	1.37	E	20	1.62
	F	50	1.41	G	100	1.47	H	200	1.53	I	300	1.45			
n-Butylbenzene							C	5.0	1.79	D	10	1.92	E	20	2.36
	F	50	2.09	G	100	2.27	H	200	2.43	I	300	2.29			
1,2-Dichlorobenzene	A	1.0	1.30	B	2.0	1.42	C	5.0	1.38	D	10	1.30	E	20	1.64
	F	50	1.39	G	100	1.44	H	200	1.51	I	300	1.42			
1,2-Dibromo-3-chloropropane				B	2.0	0.0926	C	5.0	0.115	D	10	0.0877	E	20	0.119
	F	50	0.0940	G	100	0.0982	H	200	0.131	I	300	0.133			
1,2,4-Trichlorobenzene							C	5.0	0.911	D	10	0.837	E	20	1.01
	F	50	0.687	G	100	0.884	H	200	1.07						
Hexachlorobutadiene	A	1.0	0.471	B	2.0	0.357	C	5.0	0.502	D	10	0.497	E	20	0.583
	F	50	0.356	G	100	0.380	H	200	0.592	I	300	0.319			
Naphthalene							C	5.0	2.01	D	10	1.89	E	20	2.31
	F	50	1.73	G	100	2.16	H	200	2.69	I	300	1.70			
1,2,3-Trichlorobenzene							C	5.0	0.732	D	10	0.848	E	20	1.05
	F	50	0.688	G	100	0.645	H	200	1.07						
Dibromofluoromethane				B	10	0.230	C	20	0.256	D	30	0.248	E	40	0.256
	F	50	0.240	G	70	0.257	H	80	0.265	I	100	0.248			
Toluene-d8				B	10	0.728	C	20	0.885	D	30	0.845	E	40	0.927
	F	50	0.844	G	70	0.939	H	80	0.958	I	100	0.887			
4-Bromofluorobenzene				B	10	0.690	C	20	0.767	D	30	0.720	E	40	0.795
	F	50	0.808	G	70	0.764	H	80	0.813	I	100	0.782			

Results flagged with an asterisk (*) indicate values outside control criteria.

† **SPCC Compound** ‡ **CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Compound Type	Calibration Evaluation					RRF Evaluation		
		Fit Type	Eval.	Eval. Result	Q	Control Criteria	Average RRF	Q	Minimum RRF
Dichlorodifluoromethane	TRG	AverageRF	% RSD	6.1		≤ 15	0.309		0.01
† Chloromethane	TRG	AverageRF	% RSD	11.9		≤ 15	0.391		0.10
‡ Vinyl Chloride	MS	AverageRF	% RSD	5.4		≤ 15	0.347		0.01
Bromomethane	TRG	AverageRF	% RSD	10.6		≤ 15	0.289		0.01
Chloroethane	TRG	AverageRF	% RSD	5.5		≤ 15	0.226		0.01
Trichlorofluoromethane	TRG	AverageRF	% RSD	5.7		≤ 15	0.479		0.01
‡ 1,1-Dichloroethene	MS	AverageRF	% RSD	8.7		≤ 15	0.210		0.01
Acetone	TRG	AverageRF	% RSD	15.1	*	≤ 15	0.116		0.01
Carbon Disulfide	TRG	AverageRF	% RSD	7.6		≤ 15	0.825		0.01
Methylene Chloride	TRG	AverageRF	% RSD	5.7		≤ 15	0.301		0.01
trans-1,2-Dichloroethene	TRG	AverageRF	% RSD	8.4		≤ 15	0.263		0.01
† 1,1-Dichloroethane	TRG	AverageRF	% RSD	6.5		≤ 15	0.518		0.10
2,2-Dichloropropane	TRG	AverageRF	% RSD	7.0		≤ 15	0.355		0.01
cis-1,2-Dichloroethene	TRG	AverageRF	% RSD	5.4		≤ 15	0.291		0.01
2-Butanone (MEK)	TRG	AverageRF	% RSD	8.7		≤ 15	0.0374		0.01
Bromochloromethane	TRG	AverageRF	% RSD	5.9		≤ 15	0.143		0.01
‡ Chloroform	MS	AverageRF	% RSD	6.6		≤ 15	0.489		0.01
1,1,1-Trichloroethane (TCA)	TRG	AverageRF	% RSD	6.0		≤ 15	0.373		0.01
Carbon Tetrachloride	MS	AverageRF	% RSD	7.7		≤ 15	0.321		0.01
1,1-Dichloropropene	TRG	AverageRF	% RSD	7.2		≤ 15	0.335		0.01
Benzene	MS	AverageRF	% RSD	6.5		≤ 15	1.07		0.01
1,2-Dichloroethane (EDC)	TRG	AverageRF	% RSD	4.7		≤ 15	0.432		0.01
Trichloroethene (TCE)	MS	AverageRF	% RSD	6.6		≤ 15	0.282		0.01
‡ 1,2-Dichloropropane	TRG	AverageRF	% RSD	7.6		≤ 15	0.284		0.01
Dibromomethane	TRG	AverageRF	% RSD	5.2		≤ 15	0.179		0.01
Bromodichloromethane	MS	AverageRF	% RSD	10.1		≤ 15	0.358		0.01
cis-1,3-Dichloropropene	TRG	AverageRF	% RSD	14.4		≤ 15	0.399		0.01
4-Methyl-2-pentanone (MIBK)	TRG	AverageRF	% RSD	9.4		≤ 15	0.109		0.01
‡ Toluene	MS	AverageRF	% RSD	7.0		≤ 15	0.640		0.01
trans-1,3-Dichloropropene	TRG	AverageRF	% RSD	14.2		≤ 15	0.828		0.01
1,1,2-Trichloroethane	MS	AverageRF	% RSD	8.2		≤ 15	0.518		0.01
Tetrachloroethene (PCE)	TRG	AverageRF	% RSD	6.9		≤ 15	0.491		0.01
2-Hexanone	MS	AverageRF	% RSD	8.0		≤ 15	0.0874		0.01
1,3-Dichloropropane	TRG	AverageRF	% RSD	7.9		≤ 15	1.03		0.01
Dibromochloromethane	TRG	AverageRF	% RSD	12.4		≤ 15	0.579		0.01
1,2-Dibromoethane (EDB)	TRG	AverageRF	% RSD	7.2		≤ 15	0.566		0.01
† Chlorobenzene	MS	AverageRF	% RSD	6.6		≤ 15	1.73		0.30
‡ Ethylbenzene	MS	AverageRF	% RSD	6.8		≤ 15	0.841		0.01
1,1,1,2-Tetrachloroethane	TRG	AverageRF	% RSD	9.9		≤ 15	0.570		0.01
m,p-Xylenes	TRG	AverageRF	% RSD	9.8		≤ 15	0.991		0.01
o-Xylene	TRG	AverageRF	% RSD	8.1		≤ 15	0.995		0.01
Styrene	TRG	AverageRF	% RSD	11.6		≤ 15	0.780		0.01

Results flagged with an asterisk (*) indicate values outside control criteria.

† ***SPCC Compound*** ‡ **CCC Compound**

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010

Initial Calibration Summary
Volatile Organic Compounds

Calibration ID: CAL9404
Instrument ID: MS24

Column: MS

Analyte Name	Compound Type	Calibration Evaluation					RRF Evaluation		
		Fit Type	Eval.	Eval. Result	Q	Control Criteria	Average RRF	Q	Minimum RRF
† Bromoform	TRG	AverageRF	% RSD	15.1	*	≤ 15	0.352		0.10
Isopropylbenzene	TRG	AverageRF	% RSD	8.6		≤ 15	2.40		0.01
† 1,1,2,2-Tetrachloroethane	TRG	AverageRF	% RSD	12.9		≤ 15	0.728		0.30
Bromobenzene	TRG	AverageRF	% RSD	10.7		≤ 15	0.785		0.01
n-Propylbenzene	TRG	AverageRF	% RSD	10.9		≤ 15	3.20		0.01
1,2,3-Trichloropropane	MS	AverageRF	% RSD	12.4		≤ 15	0.277		0.01
2-Chlorotoluene	MS	AverageRF	% RSD	8.3		≤ 15	2.01		0.01
1,3,5-Trimethylbenzene	TRG	AverageRF	% RSD	10.8		≤ 15	2.25		0.01
4-Chlorotoluene	TRG	AverageRF	% RSD	8.5		≤ 15	2.03		0.01
tert-Butylbenzene	TRG	AverageRF	% RSD	9.6		≤ 15	1.97		0.01
1,2,4-Trimethylbenzene	TRG	AverageRF	% RSD	12.4		≤ 15	2.30		0.01
sec-Butylbenzene	TRG	AverageRF	% RSD	10.6		≤ 15	2.94		0.01
4-Isopropyltoluene	TRG	AverageRF	% RSD	10.4		≤ 15	2.45		0.01
1,3-Dichlorobenzene	TRG	AverageRF	% RSD	8.8		≤ 15	1.45		0.01
1,4-Dichlorobenzene	TRG	AverageRF	% RSD	6.9		≤ 15	1.50		0.01
n-Butylbenzene	TRG	AverageRF	% RSD	11.1		≤ 15	2.17		0.01
1,2-Dichlorobenzene	MS	AverageRF	% RSD	7.3		≤ 15	1.42		0.01
1,2-Dibromo-3-chloropropane	TRG	AverageRF	% RSD	16.5	*	≤ 15	0.109		0.01
1,2,4-Trichlorobenzene	TRG	AverageRF	% RSD	15.1	*	≤ 15	0.901		0.01
Hexachlorobutadiene	TRG	AverageRF	% RSD	22.5	*	≤ 15	0.451		0.01
Naphthalene	MS	AverageRF	% RSD	17.0	*	≤ 15	2.07		0.01
1,2,3-Trichlorobenzene	TRG	AverageRF	% RSD	22.0	*	≤ 15	0.839		0.01
Dibromofluoromethane	SURR	AverageRF	% RSD	4.4		≤ 15	0.250		0.01
Toluene-d8	SURR	AverageRF	% RSD	8.3		≤ 15	0.877		0.01
4-Bromofluorobenzene	SURR	AverageRF	% RSD	5.6		≤ 15	0.767		0.01

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

COLUMBIA ANALYTICAL SERVICES, INC.

QA/QC Results

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010
Date Analyzed: 04/21/2010

Second Source Calibration Verification
Volatile Organic Compounds

Calibration Type: Internal Standard
Analysis Method: 8260B

Calibration ID: CAL9404
Units: PPB

File ID: J:\MS24\DATA\042110\0421F017.D

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
Dichlorodifluoromethane	50	48	0.309	0.298	-4	NA	± 40 %	AverageRF
† Chloromethane	50	43	0.391	0.336	-14	NA	± 40 %	AverageRF
‡ Vinyl Chloride	50	46	0.347	0.319	-8	NA	± 20 %	AverageRF
Bromomethane	50	45	0.289	0.257	-11	NA	± 40 %	AverageRF
Chloroethane	50	41	0.226	0.188	-17	NA	± 40 %	AverageRF
Trichlorofluoromethane	50	42	0.479	0.404	-16	NA	± 30 %	AverageRF
‡ 1,1-Dichloroethene	50	51	0.210	0.214	2	NA	± 20 %	AverageRF
Acetone	250	210	0.116	0.0983	-16	NA	± 30 %	AverageRF
Carbon Disulfide	100	96	0.825	0.791	-4	NA	± 30 %	AverageRF
Methylene Chloride	50	46	0.301	0.275	-9	NA	± 30 %	AverageRF
trans-1,2-Dichloroethene	50	49	0.263	0.256	-3	NA	± 30 %	AverageRF
† 1,1-Dichloroethane	50	47	0.518	0.485	-6	NA	± 30 %	AverageRF
2,2-Dichloropropane	50	51	0.355	0.359	1	NA	± 30 %	AverageRF
cis-1,2-Dichloroethene	50	46	0.291	0.268	-8	NA	± 30 %	AverageRF
2-Butanone (MEK)	250	230	0.0374	0.0348	-7	NA	± 30 %	AverageRF
Bromochloromethane	50	48	0.143	0.137	-4	NA	± 30 %	AverageRF
‡ Chloroform	50	48	0.489	0.472	-3	NA	± 20 %	AverageRF
1,1,1-Trichloroethane (TCA)	50	51	0.373	0.380	2	NA	± 30 %	AverageRF
Carbon Tetrachloride	50	51	0.321	0.325	1	NA	± 30 %	AverageRF
1,1-Dichloropropene	50	48	0.335	0.321	-4	NA	± 30 %	AverageRF
Benzene	50	48	1.07	1.02	-4	NA	± 30 %	AverageRF
1,2-Dichloroethane (EDC)	50	47	0.432	0.407	-6	NA	± 30 %	AverageRF
Trichloroethene (TCE)	50	51	0.282	0.287	2	NA	± 30 %	AverageRF
‡ 1,2-Dichloropropane	50	50	0.284	0.282	-1	NA	± 20 %	AverageRF
Dibromomethane	50	50	0.179	0.178	0	NA	± 30 %	AverageRF
Bromodichloromethane	50	51	0.358	0.364	2	NA	± 30 %	AverageRF
cis-1,3-Dichloropropene	50	45	0.399	0.361	-10	NA	± 30 %	AverageRF
4-Methyl-2-pentanone (MIBK)	250	250	0.109	0.108	-1	NA	± 30 %	AverageRF
‡ Toluene	50	50	0.640	0.634	-1	NA	± 20 %	AverageRF
trans-1,3-Dichloropropene	50	58	0.828	0.953	15	NA	± 30 %	AverageRF
1,1,2-Trichloroethane	50	51	0.518	0.527	2	NA	± 30 %	AverageRF
Tetrachloroethene (PCE)	50	53	0.491	0.524	7	NA	± 30 %	AverageRF
2-Hexanone	250	260	0.0874	0.0904	3	NA	± 30 %	AverageRF
1,3-Dichloropropane	50	51	1.03	1.05	1	NA	± 30 %	AverageRF
Dibromochloromethane	50	51	0.579	0.591	2	NA	± 30 %	AverageRF
1,2-Dibromoethane (EDB)	50	52	0.566	0.591	5	NA	± 30 %	AverageRF
† Chlorobenzene	50	50	1.73	1.72	0	NA	± 30 %	AverageRF
‡ Ethylbenzene	50	53	0.841	0.897	7	NA	± 20 %	AverageRF
1,1,1,2-Tetrachloroethane	50	52	0.570	0.591	4	NA	± 30 %	AverageRF
m,p-Xylenes	100	110	0.991	1.11	12	NA	± 30 %	AverageRF

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Client: Exponent
Project: Heglar - Kronquist/0907194.000.0601

Service Request: K1005244
Calibration Date: 04/21/2010
Date Analyzed: 04/21/2010

Second Source Calibration Verification
Volatile Organic Compounds

Calibration Type: Internal Standard
Analysis Method: 8260B

Calibration ID: CAL9404
Units: PPB

Analyte Name	Expected	Result	Average RF	SSV RF	%D	%Drift	Criteria	Curve Fit
o-Xylene	50	54	0.995	1.08	8	NA	± 30 %	AverageRF
Styrene	50	54	0.780	0.848	9	NA	± 30 %	AverageRF
† Bromoform	50	54	0.352	0.383	9	NA	± 30 %	AverageRF
Isopropylbenzene	50	58	2.40	2.76	15	NA	± 30 %	AverageRF
† 1,1,2,2-Tetrachloroethane	50	47	0.728	0.686	-6	NA	± 30 %	AverageRF
Bromobenzene	50	47	0.785	0.734	-7	NA	± 30 %	AverageRF
n-Propylbenzene	50	52	3.20	3.32	4	NA	± 30 %	AverageRF
1,2,3-Trichloropropane	50	46	0.277	0.256	-7	NA	± 30 %	AverageRF
2-Chlorotoluene	50	51	2.01	2.04	2	NA	± 30 %	AverageRF
1,3,5-Trimethylbenzene	50	53	2.25	2.37	5	NA	± 30 %	AverageRF
4-Chlorotoluene	50	50	2.03	2.01	-1	NA	± 30 %	AverageRF
tert-Butylbenzene	50	53	1.97	2.09	6	NA	± 30 %	AverageRF
1,2,4-Trimethylbenzene	50	53	2.30	2.44	6	NA	± 30 %	AverageRF
sec-Butylbenzene	50	54	2.94	3.18	8	NA	± 30 %	AverageRF
4-Isopropyltoluene	50	56	2.45	2.73	12	NA	± 30 %	AverageRF
1,3-Dichlorobenzene	50	52	1.45	1.52	5	NA	± 30 %	AverageRF
1,4-Dichlorobenzene	50	51	1.50	1.54	2	NA	± 30 %	AverageRF
n-Butylbenzene	50	56	2.17	2.41	12	NA	± 30 %	AverageRF
1,2-Dichlorobenzene	50	50	1.42	1.41	-1	NA	± 30 %	AverageRF
1,2-Dibromo-3-chloropropane	50	57	0.109	0.124	14	NA	± 30 %	AverageRF
1,2,4-Trichlorobenzene	50	53	0.901	0.961	7	NA	± 30 %	AverageRF
Hexachlorobutadiene	50	59	0.451	0.536	19	NA	± 30 %	AverageRF
Naphthalene	50	54	2.07	2.22	7	NA	± 30 %	AverageRF
1,2,3-Trichlorobenzene	50	60	0.839	0.999	19	NA	± 30 %	AverageRF

Results flagged with an asterisk (*) indicate values outside control criteria.

† SPCC Compound **‡ CCC Compound**

Date: 4/24/10

Columbia Analytical Services, Inc.

Injection Log

MS24 - Agilent 5975

Tune File: ~~tune.sun~~

By: KL

New Tune: NO

IS/SS Std. ID: 5604A-30C

4/24/10 Cal ~~8260~~ 9404 8260

CCV Std ID: See prepsheet

ICAL Date: 4/24/10 Cal 9402 DOD

MS/DMS/LCS/ICV Std ID: 1

Second RV: 4/26/10

BFB Std. ID: 5604A-18F

LIMS ID:

	Sample Name	File Name	Method	Dilution	pH	R	Comments
1	IB	0424F001	8260.M				
2	BFB	2		1ml 5ml			
3	IB	3					
4	IB	4					
5	Soil Ical 1	5					
6	2	6					
7	5	7					
8	10	8					
9	20	9					
10	50	10					
11	100	11					
12	200	12					
13	300	13					
14	ID	14					
15	IB	15					
16	ICV	16					
17	ICV	17					
18							
19							
20							
21							
22							
23							
24							
25							
26							
27							

INTIAL CALIBRATION CURVE

Date 4/21/10 Analysis 8260 soil Appendix

Prepared By KL Instrument MS24

Stock Solution #1 SLUA 20F Analytes Surrogates Init. Concentr: 100 ppm

Stock Solution #2 J J3A Analytes 8260 soil CCV mix Init. Concentr: 100/200/500

~~Stock Solution #3~~ ~~Analytes Ethyl Acrylate~~ ~~Init. Concentration 1000~~

Stock Solution #4 Analytes

#	Aliquot of Stock Solution #1	Final Conc. of #1 (µg/L)	Aliquot of Stock Solution #2	Final Conc. of #2 (µg/L)	Aliquot of Stock Solution #3	Final Conc. of #3 (µg/L)	Aliquot of Stock Solution #4 (µL)	Final Conc. of #4 (µg/L)
1			0.5	1				
2	5	10	1	2				
3	10	20	2.5	5	0.5	10		
4	15	30	5	10	1	20		
5	20	40	10	20	2.5	50		
6	25	50	25	50	5	100		
7	35	70	50	100	10	200		
8	40	80	100	200	20	400		
9	50	100	150	300				

4/21/10

ICV std#: Restek ICV 50uL 31A

Oxygenates 25uL 28B

M

Freon 21 25uL 22E

Appx 20uL 220

Acrolein 50uL 29E

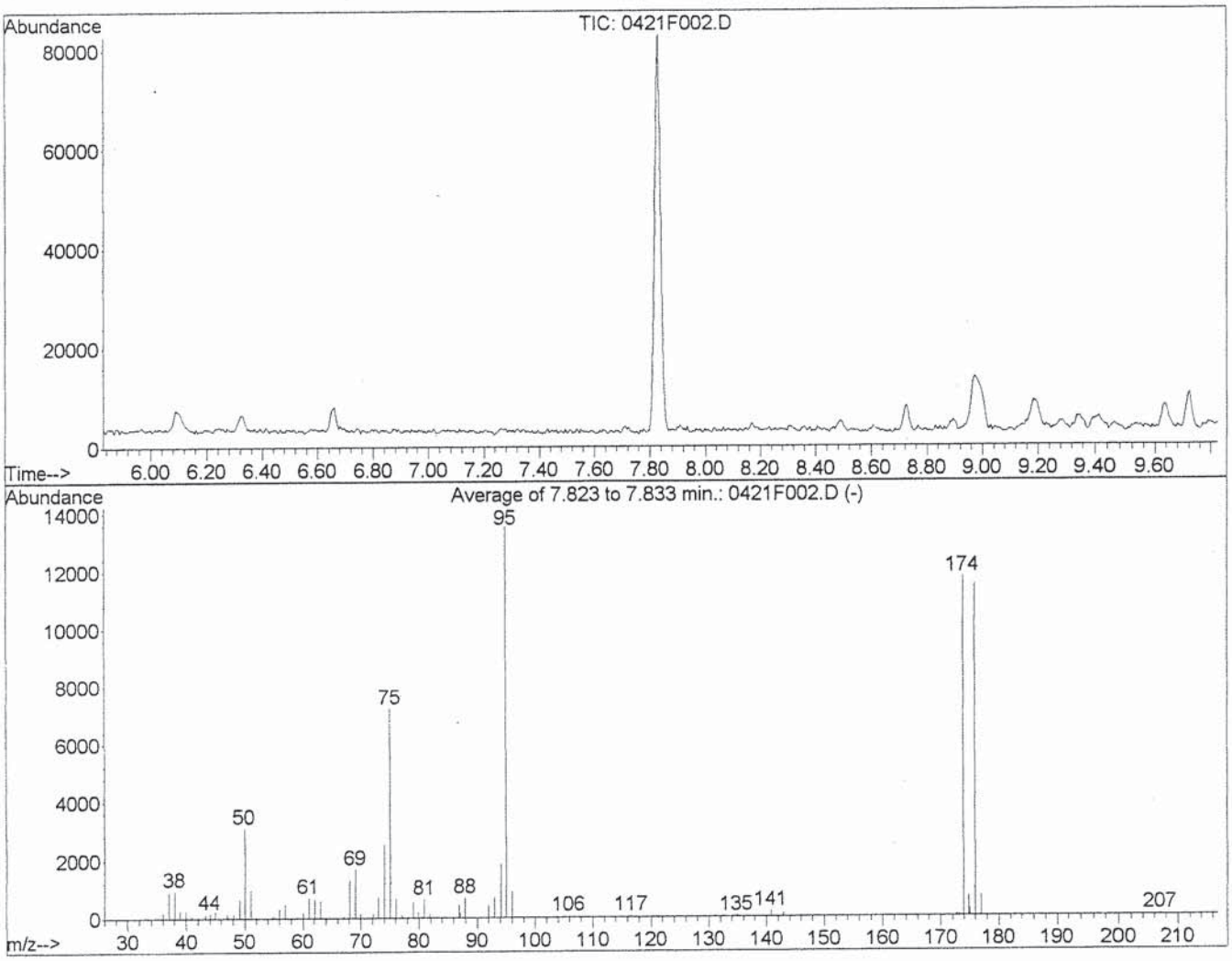
~~Ethyl Acrylate 2.5uL~~ KL 4/21/10

~~Eth/Prop. Oxide 5uL~~

1099

Data File : J:\MS24\DATA\042110\0421F002.D
Acq On : 21 Apr 2010 9:48 am
Sample : BFB
Misc :
MS Integration Params: rteint.p
Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B

Vial: 3
Operator: KR
Inst : MS24
Multiplr: 1.00



AutoFind: Scans 1342, 1343, 1344; Background Corrected with Scan 1335

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	22.8	3095	PASS
75	95	30	60	53.3	7226	PASS
95	95	100	100	100.0	13552	PASS
96	95	5	9	6.7	906	PASS
173	174	0.00	2	0.6	70	PASS
174	95	50	120	86.7	11746	PASS
175	174	5	9	5.9	698	PASS
176	174	95	101	97.8	11493	PASS
177	176	5	9	6.0	693	PASS

0421F002.D 042110MS24SOIL.M Wed Apr 21 13:43:25 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F004.D
Acq On : 21 Apr 2010 10:36 am
Sample : IB
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:43:32 2010

Vial: 6
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Thu Apr 01 15:44:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	4.02	96	300532	50.00	PPB	0.00
59) Chlorobenzene-d5	6.65	82	122711	50.00	PPB	0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	105253	50.00	PPB	0.00
System Monitoring Compounds						
38) Dibromofluoromethane	0.00	113	0	0.00	PPB	
Spiked Amount 50.000			Recovery =	0.00%		
42) 1,2-Dichloroethane-d4	3.81	65	784	0.42	PPB	-0.01
Spiked Amount 50.000			Recovery =	0.84%		
57) Toluene-d8	5.31	98	1786	0.38	PPB	0.00
Spiked Amount 50.000			Recovery =	0.76%		
79) 4-Bromofluorobenzene	7.83	95	831	0.46	PPB	0.00
Spiked Amount 50.000			Recovery =	0.92%		
Target Compounds						Qvalue
13) Acetone	1.98	43	7387	9.48	PPB	95
18) Methylene Chloride	2.27	84	1568	0.98	PPB	80
19) tert-Butyl Alcohol	2.35	59	1686	6.58	PPB	95
58) Toluene	5.36	92	1474	0.40	PPB #	67

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4/21/10

(#) = qualifier out of range (m) = manual integration
0421F004.D 042110MS24SOIL.M Wed Apr 21 13:44:43 2010 Page 1

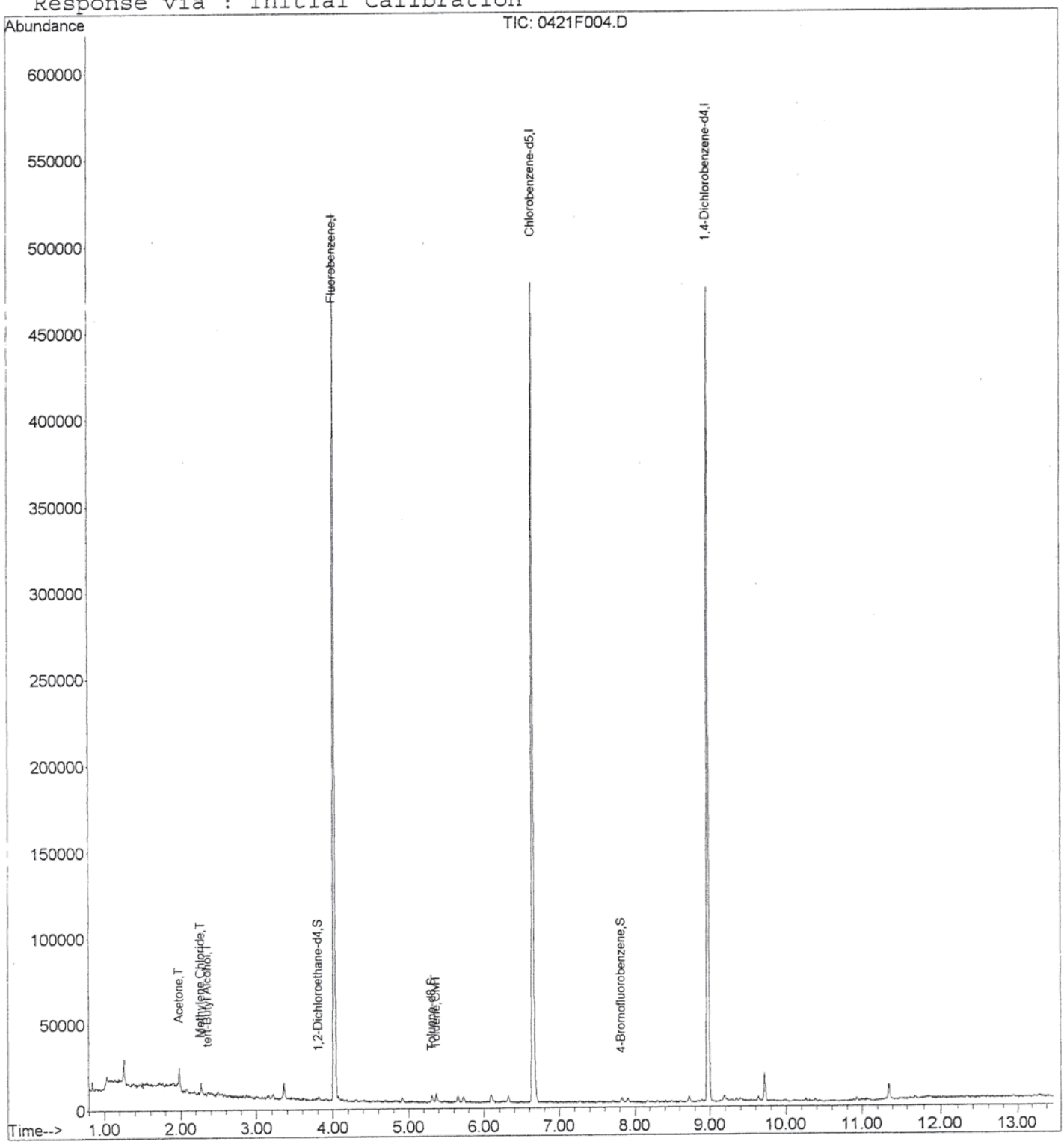
Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F004.D
Acq On : 21 Apr 2010 10:36 am
Sample : IB
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:44 2010

Vial: 6
Operator: KR
Inst : MS24
Multiplr: 1.00

Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Thu Apr 01 15:44:07 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F005.D
Acq On : 21 Apr 2010 12:29 pm
Sample : SOIL ICAL 1
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:53:05 2010

Vial: 7
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

KR 4/21/10

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene	4.02	96	323147	50.00	PPB		0.00
59) Chlorobenzene-d5	6.65	82	128828	50.00	PPB		0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	114981	50.00	PPB		0.00
System Monitoring Compounds							
38) Dibromofluoromethane	0.00	113	0	0.00	PPB		
Spiked Amount 50.000			Recovery =	0.00%			
42) 1,2-Dichloroethane-d4	0.00	65	0d	0.00	PPB		
Spiked Amount 50.000			Recovery =	0.00%			
57) Toluene-d8	0.00	98	0d	0.00	PPB		
Spiked Amount 50.000			Recovery =	0.00%			
79) 4-Bromofluorobenzene	0.00	95	0d	0.00	PPB		
Spiked Amount 50.000			Recovery =	0.00%			
Target Compounds							Qvalue
2) Dichlorodifluoromethane	0.86	85	2254	1.62	PPB		91
3) Chloromethane	0.99	50	3271	1.72	PPB		95
4) Vinyl Chloride	1.06	62	2660	1.43	PPB		93
5) Bromomethane	1.29	96	3186	2.61	PPB		84
6) Chloroethane	1.36	64	1719	1.48	PPB		75
7) Dichlorofluoromethane	1.52	67	3847	1.24	PPB		95
8) Trichlorofluoromethane	1.51	101	3200	1.16	PPB		92
9) Ethyl Ether	1.73	59	1901	1.26	PPB		96
10) Acrolein	1.87	56	2029	6.76	PPB		80
11) Trichlorotrifluoroethane	1.86	151	1215	1.07	PPB		88
12) 1,1-Dichloroethene	1.88	96	1932	1.61	PPB	#	75
13) Acetone	1.98	43	8541	10.20	PPB		100
14) Iodomethane	2.00	142	2202	16.21	PPB		98
15) Carbon Disulfide	2.01	76	5847	1.28	PPB		95
16) 3-Chloro-1-propene	2.15	76	1174	1.42	PPB		91
17) Acetonitrile	2.24	40	3399	24.31	PPB	#	86
18) Methylene Chloride	2.26	84	3837	2.23	PPB		88
19) tert-Butyl Alcohol	2.36	59	2997	10.88	PPB		84
20) Acrylonitrile	2.50	53	1742	2.64	PPB		85
21) Methyl tert-Butyl Ether	2.40	73	12089	2.44	PPB		94
22) trans-1,2-Dichloroethene	2.41	96	2005	1.36	PPB	#	76
23) Hexane	2.54	57	2720	1.33	PPB		88
24) Diisopropyl Ether	2.73	45	7163	1.36	PPB		91
25) 1,1-Dichloroethane	2.75	63	3584	1.27	PPB		95
27) Chloroprene	2.78	53	4786	2.22	PPB		91
28) tert-Butyl Ethyl Ether	2.99	59	5946	1.19	PPB		96
29) 2,2-Dichloropropane	3.15	77	2653	1.25	PPB		98

(#) = qualifier out of range (m) = manual integration

4/26/10

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F005.D Vial: 7
Acq On : 21 Apr 2010 12:29 pm Operator: KR
Sample : SOIL ICAL 1 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 13:53:05 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
30)	cis-1,2-Dichloroethene	3.18	96	1984	1.20	PPB		84
31)	2-Butanone	3.22	72	982	4.22	PPB	#	53
33)	Methacrylonitrile	3.40	67	1662	2.21	PPB		84
34)	Bromochloromethane	3.36	128	1103	1.38	PPB		87
36)	Chloroform	3.41	83	3354	1.18	PPB		98
37)	1,1,1-Trichloroethane	3.51	97	2725	1.20	PPB		90
39)	Carbon Tetrachloride	3.59	117	2181	1.11	PPB		96
40)	1,1-Dichloropropene	3.62	75	2318	1.19	PPB		98
41)	Isobutyl Alcohol	3.78	43	1822	19.79	PPB		95
43)	Benzene	3.79	78	7110	1.19	PPB		98
44)	1,2-Dichloroethane	3.87	62	2813	1.20	PPB		94
45)	tert-Amyl Methyl Ether	3.84	55	1790	1.46	PPB		86
46)	Trichloroethene	4.29	95	1963	1.25	PPB		88
48)	1,2-Dichloropropane	4.52	63	2039	1.27	PPB		81
49)	Dibromomethane	4.62	93	1224	1.21	PPB		91
50)	Methyl methacrylate	4.60	69	1281	1.03	PPB		96
52)	Bromodichloromethane	4.75	83	2436	1.13	PPB		98
53)	2-Nitropropane	5.03	41	1786	4.19	PPB	#	61
54)	2-Chloroethyl Vinyl Ether	5.02	63	1052	0.95	PPB		90
55)	cis-1,3-Dichloropropene	5.13	75	2250	0.91	PPB		92
56)	4-Methyl-2-pentanone (MIBK	5.27	58	1282	1.89	PPB		89
58)	Toluene	5.37	92	4747	1.19	PPB		99
61)	trans-1,3-Dichloropropene	5.65	75	1831	0.84	PPB		82
62)	Ethyl methacrylate	5.67	69	1624	0.81	PPB		93
63)	1,1,2-Trichloroethane	5.80	83	1382	1.19	PPB		92
64)	Tetrachloroethene	5.82	164	1304	1.08	PPB		86
66)	1,3-Dichloropropane	5.97	76	2540	1.08	PPB		87
67)	Dibromochloromethane	6.13	129	1340	0.89	PPB		89
68)	1,2-Dibromoethane (EDB)	6.25	107	1317	0.96	PPB		86
69)	1-Chlorohexane	6.65	91	1954	1.04	PPB		76
70)	Chlorobenzene	6.68	112	4444	1.09	PPB		83
71)	Ethylbenzene	6.76	106	1803	0.84	PPB		89
72)	1,1,1,2-Tetrachloroethane	6.77	131	1499	1.02	PPB		87
73)	m,p-Xylenes	6.87	106	4427	1.76	PPB		95
74)	o-Xylene	7.27	106	2489	0.98	PPB	#	81
75)	Styrene	7.30	103	1744	0.86	PPB		92
76)	Bromoform	7.50	173	770	0.81	PPB		80
77)	Isopropylbenzene	7.62	105	5114	0.82	PPB		92
81)	1,1,2,2-Tetrachloroethane	8.02	83	1511	0.97	PPB		68
83)	Bromobenzene	7.96	156	1707	1.05	PPB		88
84)	n-Propylbenzene	8.04	91	6555	0.92	PPB		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F005.D
Acq On : 21 Apr 2010 12:29 pm
Sample : SOIL ICAL 1
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:53:05 2010

Vial: 7
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

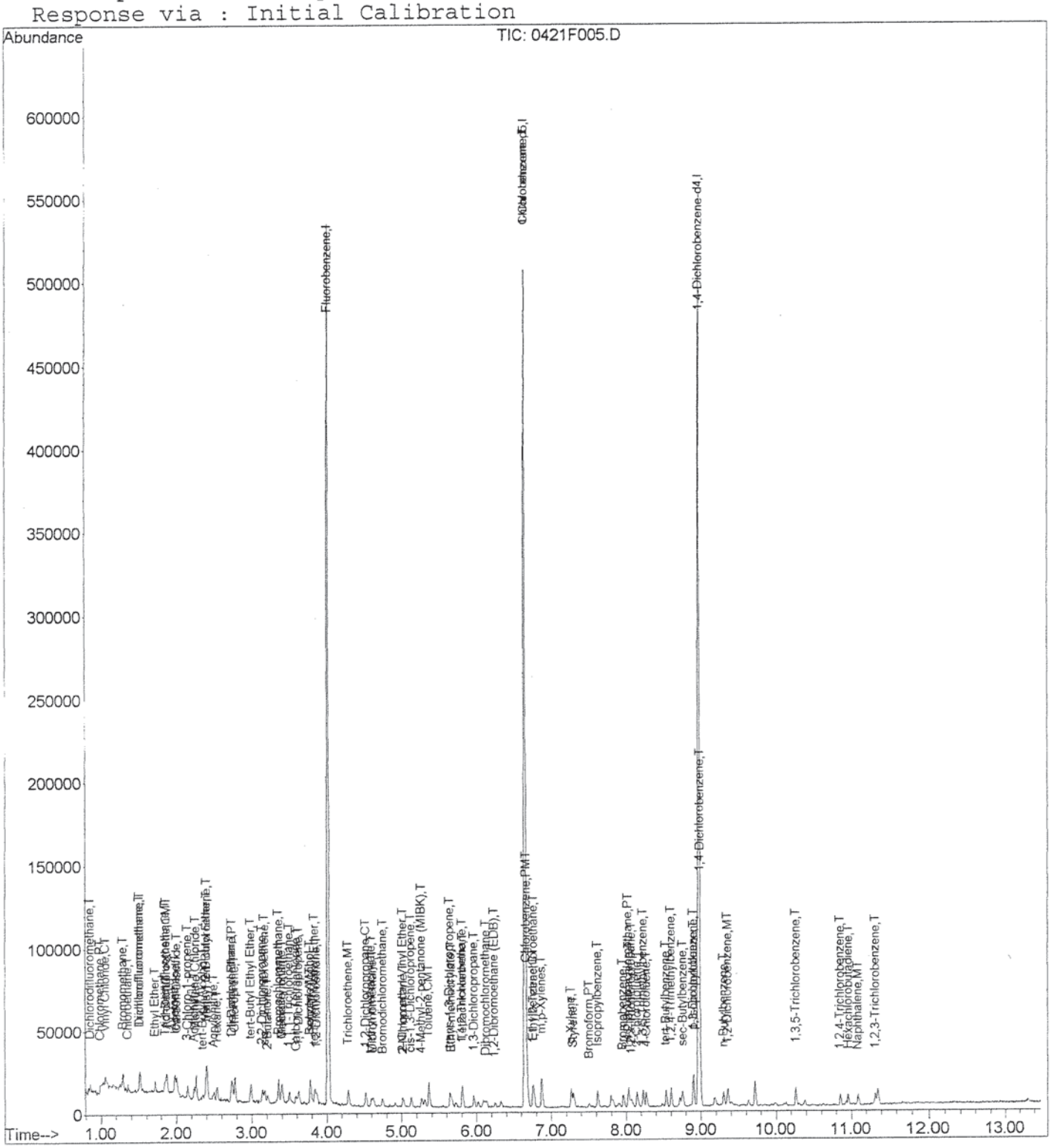
	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
85)	1,2,3-Trichloropropane	8.07	110	639	1.20	PPB		87
86)	2-Chlorotoluene	8.15	91	4420	1.05	PPB		91
87)	1,3,5-Trimethylbenzene	8.23	105	4698	0.94	PPB		94
88)	4-Chlorotoluene	8.27	91	4108	0.94	PPB		92
89)	tert-Butylbenzene	8.54	119	3977	0.90	PPB		92
90)	1,2,4-Trimethylbenzene	8.60	105	4455	0.89	PPB		96
91)	sec-Butylbenzene	8.76	105	5561	0.87	PPB		91
92)	p-Isopropyltoluene	8.91	119	5066	0.95	PPB		97
93)	1,3-Dichlorobenzene	8.90	146	2933	1.01	PPB		89
94)	1,4-Dichlorobenzene	9.00	146	3426	1.13	PPB		97
95)	n-Butylbenzene	9.31	91	4041	0.88	PPB		85
96)	1,2-Dichlorobenzene	9.36	146	3000	1.06	PPB		84
98)	1,3,5-Trichlorobenzene	10.27	180	2269	1.03	PPB		79
99)	1,2,4-Trichlorobenzene	10.85	180	1942	0.97	PPB		95
100)	Hexachlorobutadiene	10.96	225	1083	0.94	PPB		88
101)	Naphthalene	11.09	128	3928	0.77	PPB		96
102)	1,2,3-Trichlorobenzene	11.32	180	1620	0.88	PPB	#	74

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F005.D Vial: 7
Acq On : 21 Apr 2010 12:29 pm Operator: KR
Sample : SOIL ICAL 1 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 13:56 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F006.D
Acq On : 21 Apr 2010 12:50 pm
Sample : SOIL ICAL 2
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:57:32 2010

Vial: 8
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

KR 4/21/10

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene	4.02	96	338381	50.00	PPB		0.00
59) Chlorobenzene-d5	6.65	82	132058	50.00	PPB		0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	117168	50.00	PPB		0.00
System Monitoring Compounds							
38) Dibromofluoromethane	3.54	113	15547	10.26	PPB		0.00
Spiked Amount 50.000			Recovery =	20.52%			
42) 1,2-Dichloroethane-d4	3.81	65	22688	10.82	PPB		0.00
Spiked Amount 50.000			Recovery =	21.64%			
57) Toluene-d8	5.31	98	49280	9.21	PPB		0.00
Spiked Amount 50.000			Recovery =	18.42%			
79) 4-Bromofluorobenzene	7.83	95	18220	9.28	PPB		0.00
Spiked Amount 50.000			Recovery =	18.56%			
Target Compounds							Qvalue
2) Dichlorodifluoromethane	0.86	85	4473	3.07	PPB		98
3) Chloromethane	0.99	50	5380	2.70	PPB		98
4) Vinyl Chloride	1.06	62	4926	2.53	PPB		99
5) Bromomethane	1.29	96	5054	3.95	PPB		97
6) Chloroethane	1.36	64	3462	2.85	PPB		94
7) Dichlorofluoromethane	1.52	67	7161	2.21	PPB		98
8) Trichlorofluoromethane	1.51	101	6585	2.29	PPB		98
9) Ethyl Ether	1.72	59	3686	2.34	PPB		93
10) Acrolein	1.87	56	3928	12.50	PPB		78
11) Trichlorotrifluoroethane	1.86	151	2545	2.14	PPB		97
12) 1,1-Dichloroethene	1.88	96	3308	2.63	PPB		99
13) Acetone	1.98	43	7645	8.72	PPB		95
14) Iodomethane	2.00	142	4065	17.26	PPB		86
15) Carbon Disulfide	2.01	76	10567	2.22	PPB		99
16) 3-Chloro-1-propene	2.16	76	2203	2.55	PPB		91
17) Acetonitrile	2.23	40	6438	43.98	PPB	#	96
18) Methylene Chloride	2.27	84	5104	2.83	PPB		92
19) tert-Butyl Alcohol	2.35	59	4047	14.03	PPB		96
20) Acrylonitrile	2.50	53	3346	4.85	PPB		89
21) Methyl tert-Butyl Ether	2.40	73	22676	4.37	PPB		97
22) trans-1,2-Dichloroethene	2.42	96	3563	2.31	PPB		92
23) Hexane	2.54	57	5227	2.45	PPB		98
24) Diisopropyl Ether	2.74	45	13145	2.38	PPB		95
25) 1,1-Dichloroethane	2.75	63	7386	2.49	PPB		92
26) Vinyl Acetate	2.78	86	648	2.00	PPB	#	90
27) Chloroprene	2.78	53	8948	3.96	PPB		96
28) tert-Butyl Ethyl Ether	2.99	59	11340	2.16	PPB		92

(#) = qualifier out of range (m) = manual integration
0421F006.D 042110MS24SOIL.M Wed Apr 21 13:59:50 2010 4/26/10 Page 1

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F006.D
Acq On : 21 Apr 2010 12:50 pm
Sample : SOIL ICAL 2
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:57:32 2010

Vial: 8
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	2,2-Dichloropropane	3.15	77	4678	2.10	PPB		94
30)	cis-1,2-Dichloroethene	3.17	96	3922	2.26	PPB		92
31)	2-Butanone	3.21	72	1413	5.80	PPB	#	68
32)	Propionitrile	3.32	54	1578	5.39	PPB		87
33)	Methacrylonitrile	3.40	67	3335	4.24	PPB		88
34)	Bromochloromethane	3.36	128	1894	2.27	PPB		90
36)	Chloroform	3.41	83	6542	2.21	PPB		95
37)	1,1,1-Trichloroethane	3.51	97	4883	2.05	PPB		98
39)	Carbon Tetrachloride	3.60	117	4026	1.96	PPB		88
40)	1,1-Dichloropropene	3.63	75	4522	2.22	PPB		91
41)	Isobutyl Alcohol	3.78	43	4237	43.95	PPB		94
43)	Benzene	3.79	78	13820	2.21	PPB		92
44)	1,2-Dichloroethane	3.87	62	5633	2.29	PPB		96
45)	tert-Amyl Methyl Ether	3.84	55	3288	2.56	PPB	#	78
46)	Trichloroethene	4.29	95	4036	2.46	PPB		93
48)	1,2-Dichloropropane	4.53	63	3827	2.28	PPB		90
49)	Dibromomethane	4.62	93	2314	2.19	PPB		96
50)	Methyl methacrylate	4.60	69	2707	2.08	PPB		95
51)	1,4-Dioxane	4.62	88	1169	43.31	PPB		90
52)	Bromodichloromethane	4.75	83	4204	1.86	PPB		91
53)	2-Nitropropane	5.02	41	2926	6.56	PPB		89
54)	2-Chloroethyl Vinyl Ether	5.02	63	2309	1.98	PPB		92
55)	cis-1,3-Dichloropropene	5.13	75	4438	1.71	PPB		94
56)	4-Methyl-2-pentanone (MIBK	5.27	58	3195	4.50	PPB	#	73
58)	Toluene	5.36	92	8985	2.15	PPB		98
61)	trans-1,3-Dichloropropene	5.64	75	4320	1.94	PPB		95
62)	Ethyl methacrylate	5.67	69	3992	1.94	PPB		91
63)	1,1,2-Trichloroethane	5.81	83	3187	2.67	PPB		93
64)	Tetrachloroethene	5.81	164	2858	2.30	PPB		85
65)	2-Hexanone	6.02	57	1017	4.42	PPB	#	61
66)	1,3-Dichloropropane	5.97	76	6308	2.62	PPB		93
67)	Dibromochloromethane	6.13	129	3292	2.13	PPB		97
68)	1,2-Dibromoethane (EDB)	6.25	107	3286	2.34	PPB		91
69)	1-Chlorohexane	6.64	91	4006	2.07	PPB		82
70)	Chlorobenzene	6.69	112	10264	2.46	PPB		92
71)	Ethylbenzene	6.76	106	4337	1.97	PPB	#	73
72)	1,1,1,2-Tetrachloroethane	6.77	131	3398	2.27	PPB		91
73)	m,p-Xylenes	6.88	106	10030	3.88	PPB		89
74)	o-Xylene	7.27	106	5287	2.02	PPB		93
75)	Styrene	7.30	103	3780m	1.82	PPB		
76)	Bromoform	7.51	173	1922	1.96	PPB		88

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F006.D Vial: 8
Acq On : 21 Apr 2010 12:50 pm Operator: KR
Sample : SOIL ICAL 2 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 13:57:32 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
77)	Isopropylbenzene	7.63	105	11760	1.85	PPB		94
78)	cis-1,4-Dichloro-2-butene	7.79	89	630	3.43	PPB		90
81)	1,1,2,2-Tetrachloroethane	8.03	83	4403	2.77	PPB		98
82)	trans-1,4-Dichloro-2-buten	8.10	53	1026	2.63	PPB		75
83)	Bromobenzene	7.96	156	4495	2.70	PPB		89
84)	n-Propylbenzene	8.04	91	14655	2.02	PPB		97
85)	1,2,3-Trichloropropane	8.07	110	1658	3.06	PPB	#	40
86)	2-Chlorotoluene	8.15	91	10214	2.38	PPB		94
87)	1,3,5-Trimethylbenzene	8.23	105	10604	2.09	PPB		99
88)	4-Chlorotoluene	8.27	91	10443	2.35	PPB		92
89)	tert-Butylbenzene	8.54	119	10169	2.26	PPB		97
90)	1,2,4-Trimethylbenzene	8.61	105	10047	1.96	PPB		90
91)	sec-Butylbenzene	8.76	105	13324	2.05	PPB		99
92)	p-Isopropyltoluene	8.91	119	10662	1.96	PPB		97
93)	1,3-Dichlorobenzene	8.90	146	7905	2.67	PPB		97
94)	1,4-Dichlorobenzene	9.00	146	7975	2.59	PPB		95
95)	n-Butylbenzene	9.31	91	8129	1.73	PPB		97
96)	1,2-Dichlorobenzene	9.36	146	6672	2.31	PPB		87
97)	1,2-Dibromo-3-chloropropan	10.14	155	434m	1.46	PPB		
98)	1,3,5-Trichlorobenzene	10.26	180	4524	2.02	PPB		86
99)	1,2,4-Trichlorobenzene	10.86	180	3374	1.65	PPB		84
100)	Hexachlorobutadiene	10.96	225	1675	1.43	PPB		99
101)	Naphthalene	11.09	128	7922	1.53	PPB		97
102)	1,2,3-Trichlorobenzene	11.32	180	3809	2.04	PPB		84

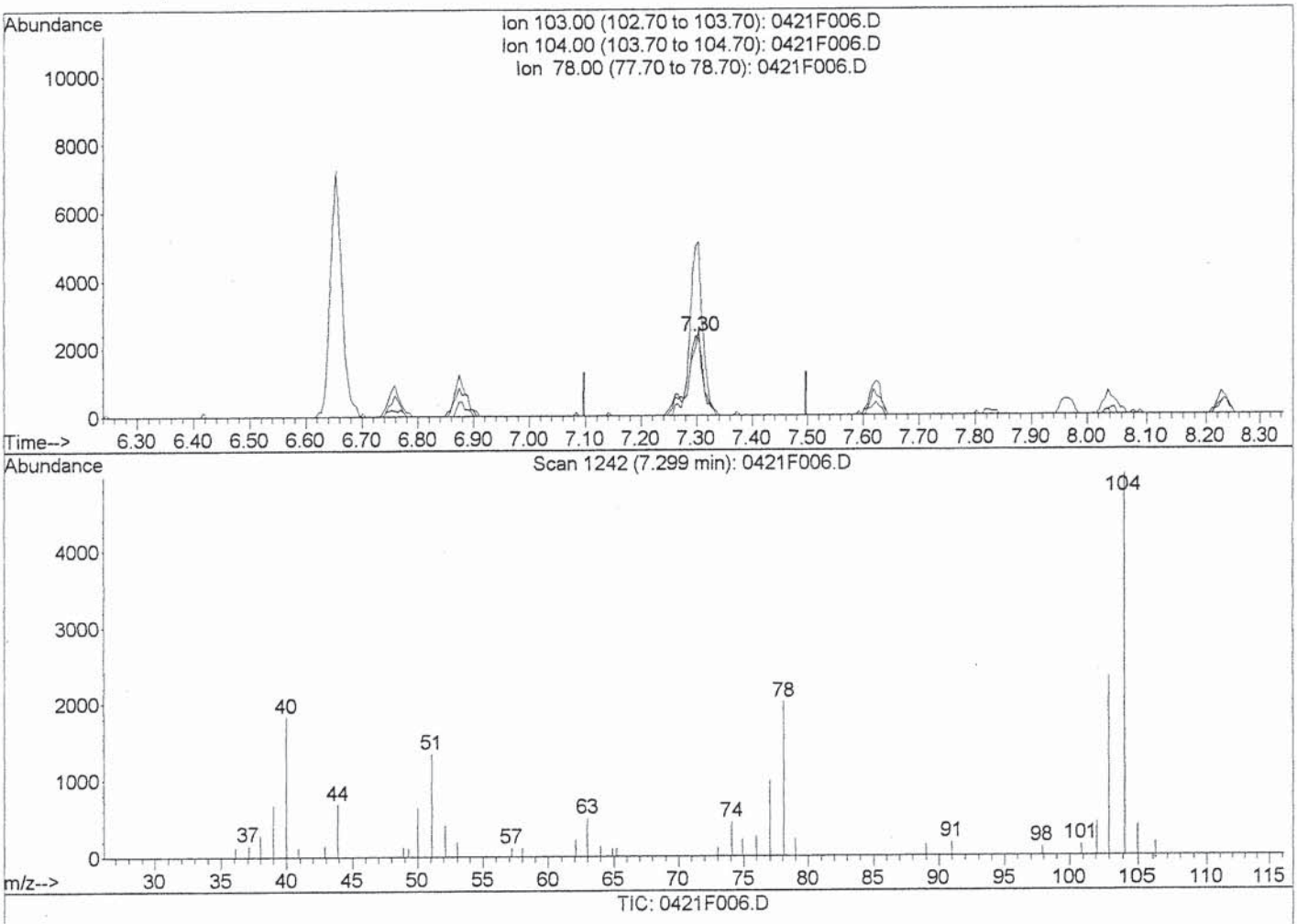
(#) = qualifier out of range (m) = manual integration

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F006.D
Acq On : 21 Apr 2010 12:50 pm
Sample : SOIL ICAL 2
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:58 2010

Vial: 8
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration



TIC: 0421F006.D

(75) Styrene (T)

7.30min 2.18PPB

response 4525

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	213.40
78.00	90.20	85.96
0.00	0.00	0.00

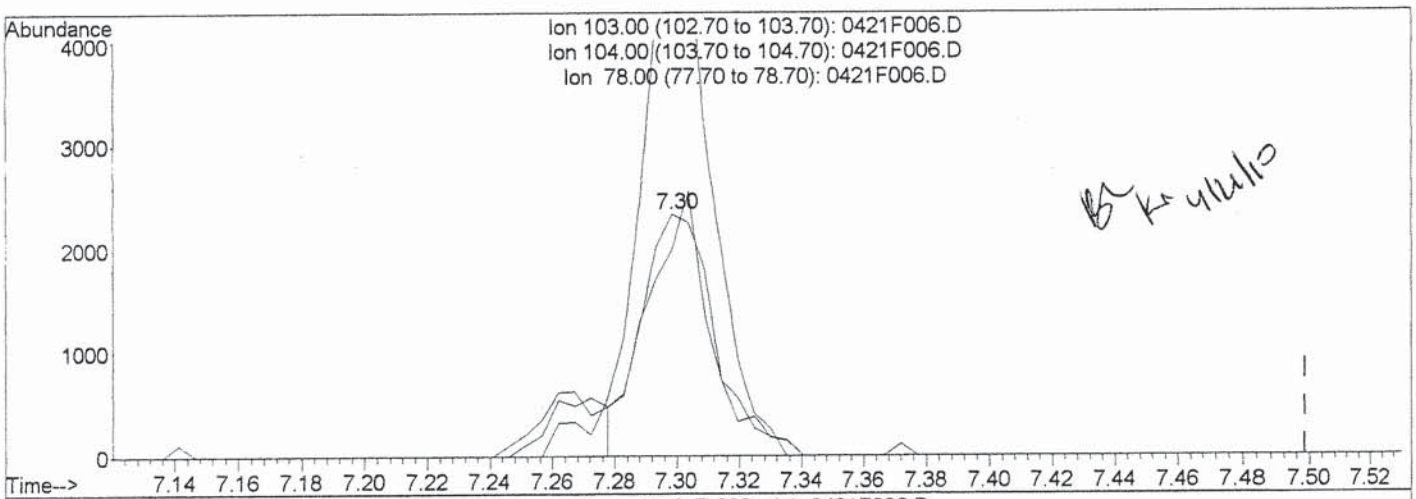
0421F006.D 042110MS24SOIL.M Wed Apr 21 13:58:43 2010

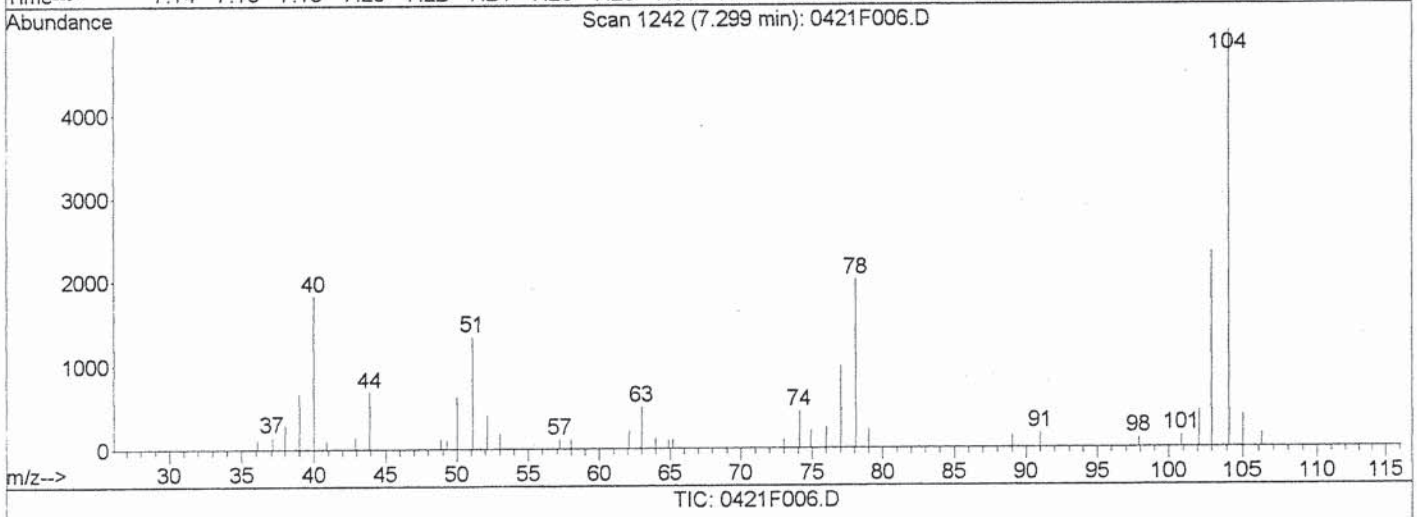
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F006.D
Acq On : 21 Apr 2010 12:50 pm
Sample : SOIL ICAL 2
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:58 2010

Vial: 8
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F006.D

(75) Styrene (T)

7.30min 1.82PPB m

response 3780

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	213.40
78.00	90.20	85.96
0.00	0.00	0.00

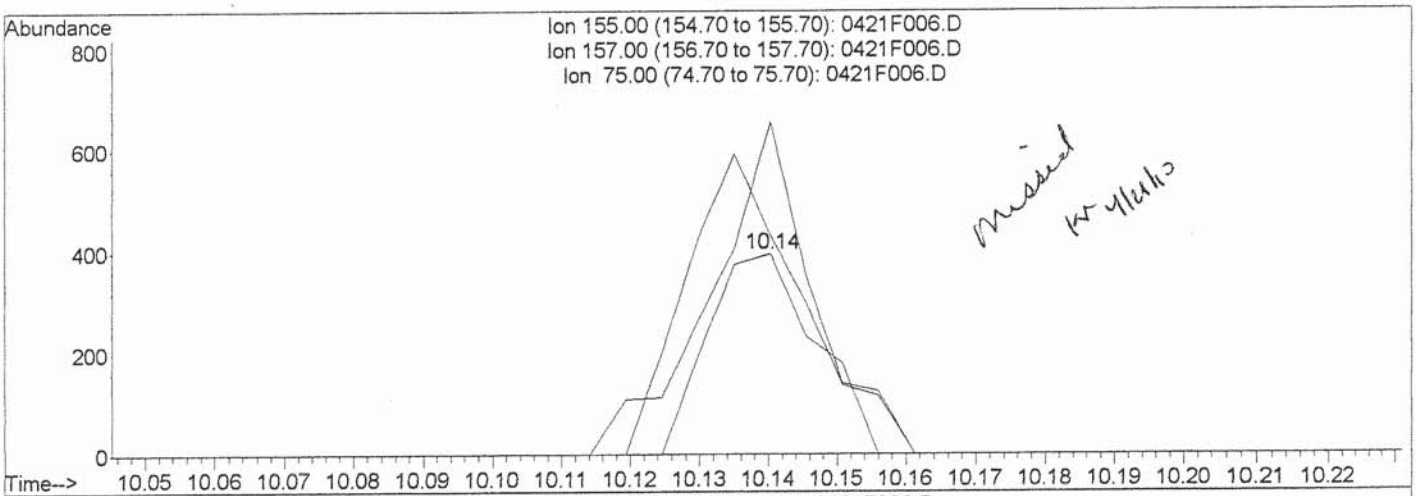
0421F006.D 042110MS24SOIL.M Wed Apr 21 13:58:50 2010

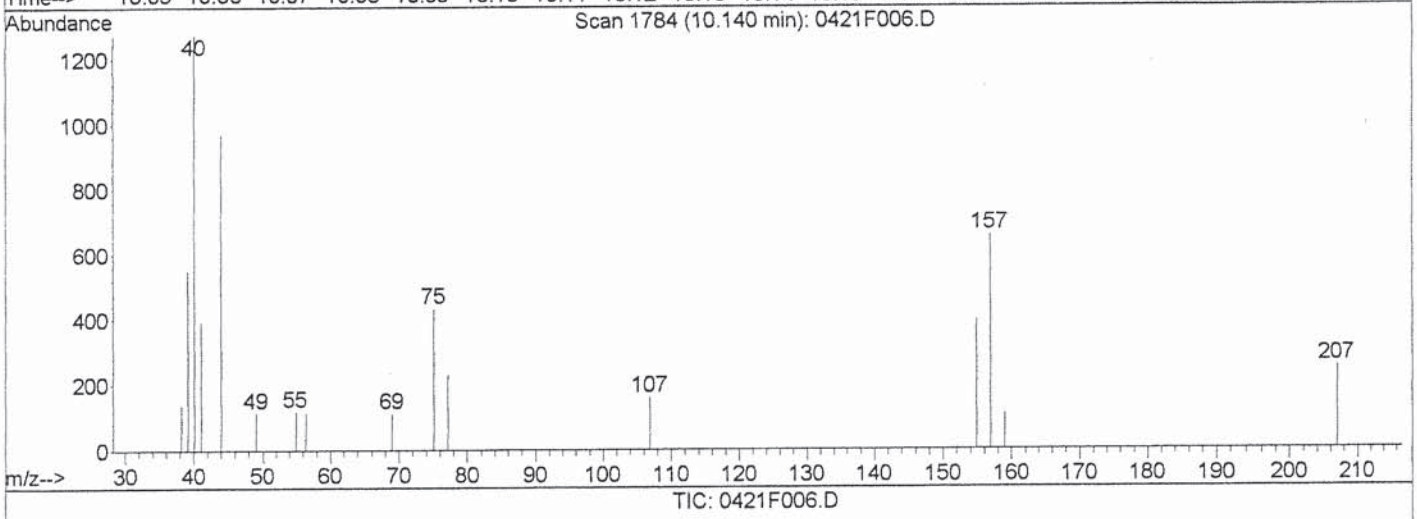
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F006.D
Acq On : 21 Apr 2010 12:50 pm
Sample : SOIL ICAL 2
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 13:59 2010

Vial: 8
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F006.D

(97) 1,2-Dibromo-3-chloropropane (DBCP) (T)

10.14min 1.46PPB m

response 434

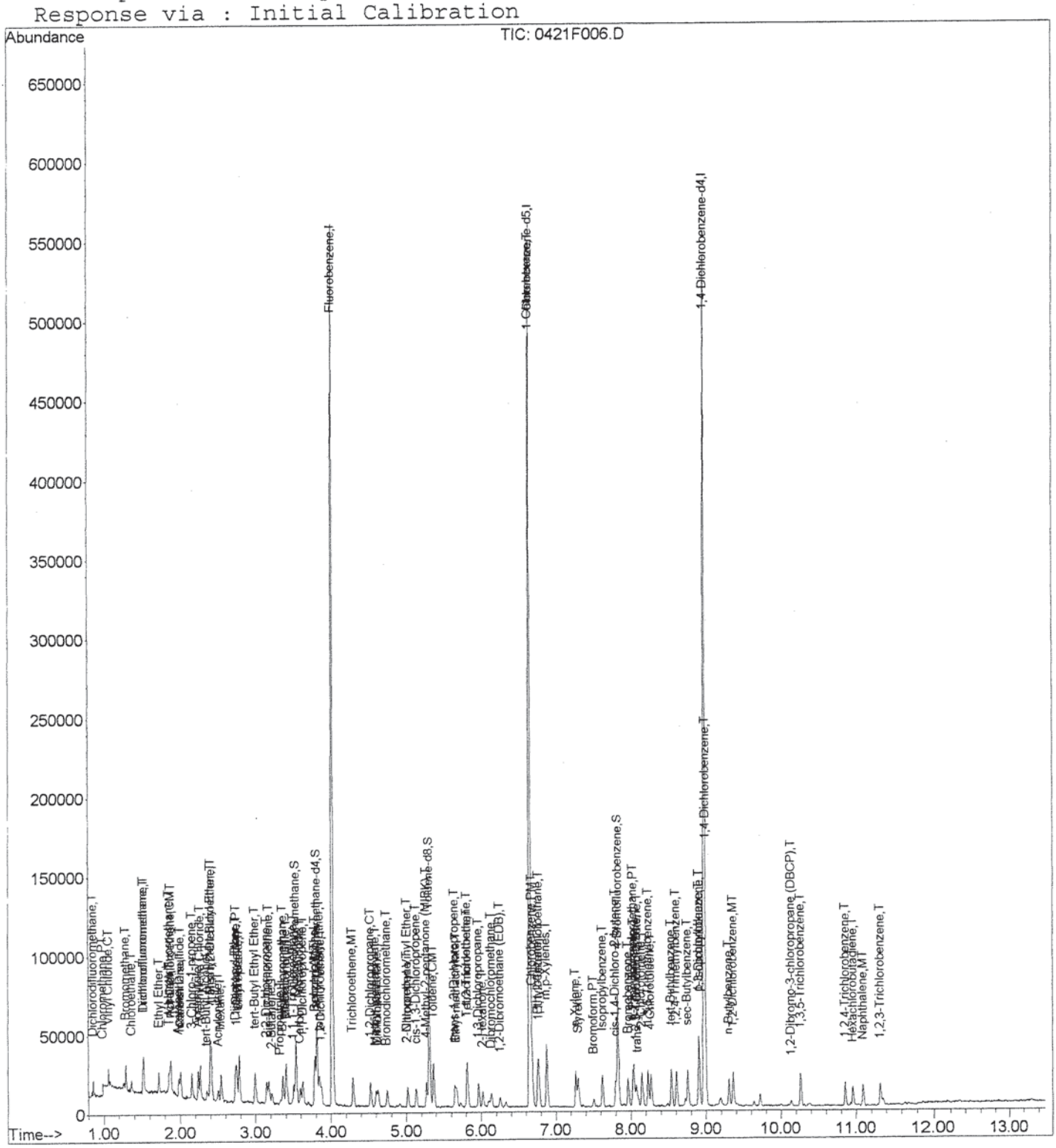
Ion	Exp%	Act%
155.00	100	100
157.00	109.90	165.91#
75.00	94.00	109.09
0.00	0.00	0.00

0421F006.D 042110MS24SOIL.M Wed Apr 21 13:59:09 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F006.D Vial: 8
Acq On : 21 Apr 2010 12:50 pm Operator: KR
Sample : SOIL ICAL 2 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 13:59 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F007.D
Acq On : 21 Apr 2010 1:11 pm
Sample : SOIL ICAL 5
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:00:00 2010

Vial: 9
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

KA 4/21/10

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	4.02	96	313503	50.00	PPB	0.00
59) Chlorobenzene-d5	6.65	82	130398	50.00	PPB	0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	118928	50.00	PPB	0.00
System Monitoring Compounds						
38) Dibromofluoromethane	3.53	113	32060	22.83	PPB	0.00
Spiked Amount 50.000			Recovery =	45.66%		
42) 1,2-Dichloroethane-d4	3.81	65	47416	24.41	PPB	0.00
Spiked Amount 50.000			Recovery =	48.82%		
57) Toluene-d8	5.31	98	110919	22.37	PPB	0.00
Spiked Amount 50.000			Recovery =	44.74%		
79) 4-Bromofluorobenzene	7.83	95	40005	20.64	PPB	0.00
Spiked Amount 50.000			Recovery =	41.28%		
Target Compounds						Qvalue
2) Dichlorodifluoromethane	0.86	85	9768	7.23	PPB	100
3) Chloromethane	0.99	50	11862	6.43	PPB	100
4) Vinyl Chloride	1.06	62	10952	6.07	PPB	100
5) Bromomethane	1.29	96	10155	8.57	PPB	100
6) Chloroethane	1.36	64	7161	6.36	PPB	100
7) Dichlorofluoromethane	1.52	67	17703	5.90	PPB	100
8) Trichlorofluoromethane	1.51	101	14806	5.55	PPB	100
9) Ethyl Ether	1.73	59	9108	6.24	PPB	100
10) Acrolein	1.87	56	8726	29.98	PPB	100
11) Trichlorotrifluoroethane	1.85	151	5642	5.13	PPB	100
12) 1,1-Dichloroethene	1.88	96	6250	5.36	PPB	100
13) Acetone	1.98	43	13412	16.50	PPB	100
14) Iodomethane	2.00	142	11565	22.29	PPB	100
15) Carbon Disulfide	2.01	76	23870	5.40	PPB	100
16) 3-Chloro-1-propene	2.16	76	4656	5.81	PPB	100
17) Acetonitrile	2.23	40	14160	104.41	PPB	100
18) Methylene Chloride	2.27	84	10099	6.04	PPB	100
19) tert-Butyl Alcohol	2.35	59	7871	29.46	PPB	100
20) Acrylonitrile	2.50	53	7036	11.00	PPB	100
21) Methyl tert-Butyl Ether	2.40	73	51350	10.69	PPB	100
22) trans-1,2-Dichloroethene	2.42	96	7850	5.50	PPB	100
23) Hexane	2.54	57	10972	5.55	PPB	100
24) Diisopropyl Ether	2.73	45	27973	5.48	PPB	100
25) 1,1-Dichloroethane	2.75	63	14956	5.45	PPB	100
26) Vinyl Acetate	2.78	86	1610	5.36	PPB	100
27) Chloroprene	2.78	53	20626	9.85	PPB	100
28) tert-Butyl Ethyl Ether	2.99	59	26133	5.38	PPB	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F007.D Vial: 9
Acq On : 21 Apr 2010 1:11 pm Operator: KR
Sample : SOIL ICAL 5 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:00:00 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
29)	2,2-Dichloropropane	3.15	77	10630	5.16	PPB	100
30)	cis-1,2-Dichloroethene	3.18	96	8500	5.29	PPB	100
31)	2-Butanone	3.21	72	2735	12.12	PPB	100
32)	Propionitrile	3.33	54	3098	11.42	PPB	100
33)	Methacrylonitrile	3.40	67	7222	9.90	PPB	100
34)	Bromochloromethane	3.36	128	4232	5.47	PPB	100
36)	Chloroform	3.41	83	14178	5.16	PPB	100
37)	1,1,1-Trichloroethane	3.50	97	11289	5.12	PPB	100
39)	Carbon Tetrachloride	3.59	117	9463	4.98	PPB	100
40)	1,1-Dichloropropene	3.63	75	9468	5.01	PPB	100
41)	Isobutyl Alcohol	3.78	43	9799	109.70	PPB	100
43)	Benzene	3.79	78	30684	5.30	PPB	100
44)	1,2-Dichloroethane	3.87	62	12920	5.66	PPB	100
45)	tert-Amyl Methyl Ether	3.84	55	6547	5.50	PPB	100
46)	Trichloroethene	4.29	95	8322	5.48	PPB	100
48)	1,2-Dichloropropane	4.53	63	7951	5.12	PPB	100
49)	Dibromomethane	4.62	93	5427	5.55	PPB	100
50)	Methyl methacrylate	4.60	69	5291	4.38	PPB	100
51)	1,4-Dioxane	4.62	88	2390	95.58	PPB	100
52)	Bromodichloromethane	4.75	83	9773	4.66	PPB	100
53)	2-Nitropropane	5.02	41	5725	13.85	PPB	100
54)	2-Chloroethyl Vinyl Ether	5.02	63	4851	4.50	PPB	100
55)	cis-1,3-Dichloropropene	5.13	75	10175	4.22	PPB	100
56)	4-Methyl-2-pentanone (MIBK	5.27	58	6486	9.85	PPB	100
58)	Toluene	5.37	92	18140	4.69	PPB	100
61)	trans-1,3-Dichloropropene	5.65	75	8613	3.91	PPB	100
62)	Ethyl methacrylate	5.67	69	8507	4.18	PPB	100
63)	1,1,2-Trichloroethane	5.81	83	6646	5.65	PPB	100
64)	Tetrachloroethene	5.82	164	6162	5.03	PPB	100
65)	2-Hexanone	6.02	57	2108	9.28	PPB	100
66)	1,3-Dichloropropane	5.97	76	12429	5.22	PPB	100
67)	Dibromochloromethane	6.13	129	6504	4.25	PPB	100
68)	1,2-Dibromoethane (EDB)	6.25	107	6597	4.76	PPB	100
69)	1-Chlorohexane	6.64	91	8271	4.34	PPB	100
70)	Chlorobenzene	6.68	112	20964	5.08	PPB	100
71)	Ethylbenzene	6.75	106	10014	4.60	PPB	100
72)	1,1,1,2-Tetrachloroethane	6.77	131	6623	4.47	PPB	100
73)	m,p-Xylenes	6.88	106	22854	8.96	PPB	100
74)	o-Xylene	7.27	106	11517	4.46	PPB	100
75)	Styrene	7.30	103	8634m	4.21	PPB	
76)	Bromoform	7.51	173	3723	3.85	PPB	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F007.D Vial: 9
Acq On : 21 Apr 2010 1:11 pm Operator: KR
Sample : SOIL ICAL 5 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:00:00 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
77)	Isopropylbenzene	7.62	105	26752	4.26	PPB	100
78)	cis-1,4-Dichloro-2-butene	7.80	89	1486	8.20	PPB	100
81)	1,1,2,2-Tetrachloroethane	8.02	83	8402	5.22	PPB	100
82)	trans-1,4-Dichloro-2-buten	8.10	53	1918	4.85	PPB	100
83)	Bromobenzene	7.96	156	8500	5.04	PPB	100
84)	n-Propylbenzene	8.04	91	31949	4.34	PPB	100
85)	1,2,3-Trichloropropane	8.08	110	3627	6.59	PPB	100
86)	2-Chlorotoluene	8.15	91	21114	4.84	PPB	100
87)	1,3,5-Trimethylbenzene	8.23	105	23126	4.49	PPB	100
88)	4-Chlorotoluene	8.27	91	21506	4.77	PPB	100
89)	tert-Butylbenzene	8.54	119	20786	4.55	PPB	100
90)	1,2,4-Trimethylbenzene	8.60	105	23148	4.46	PPB	100
91)	sec-Butylbenzene	8.76	105	29359	4.45	PPB	100
92)	p-Isopropyltoluene	8.91	119	24653	4.47	PPB	100
93)	1,3-Dichlorobenzene	8.90	146	16425	5.46	PPB	100
94)	1,4-Dichlorobenzene	9.00	146	17226	5.52	PPB	100
95)	n-Butylbenzene	9.31	91	21264	4.47	PPB	100
96)	1,2-Dichlorobenzene	9.36	146	16439	5.62	PPB	100
97)	1,2-Dibromo-3-chloropropan	10.14	155	1364	4.51	PPB	100
98)	1,3,5-Trichlorobenzene	10.26	180	12141	5.34	PPB	100
99)	1,2,4-Trichlorobenzene	10.86	180	10839	5.23	PPB	100
100)	Hexachlorobutadiene	10.96	225	5965	5.02	PPB	100
101)	Naphthalene	11.09	128	23953	4.56	PPB	100
102)	1,2,3-Trichlorobenzene	11.32	180	8700	4.58	PPB	100

(#) = qualifier out of range (m) = manual integration

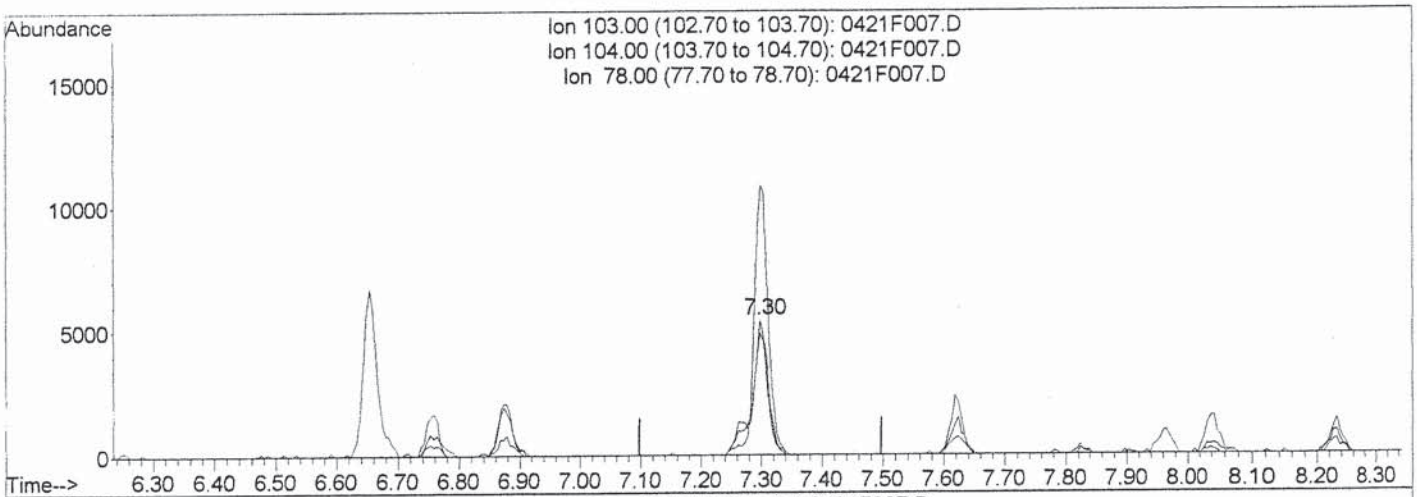
Quantitation Report (Qedit)

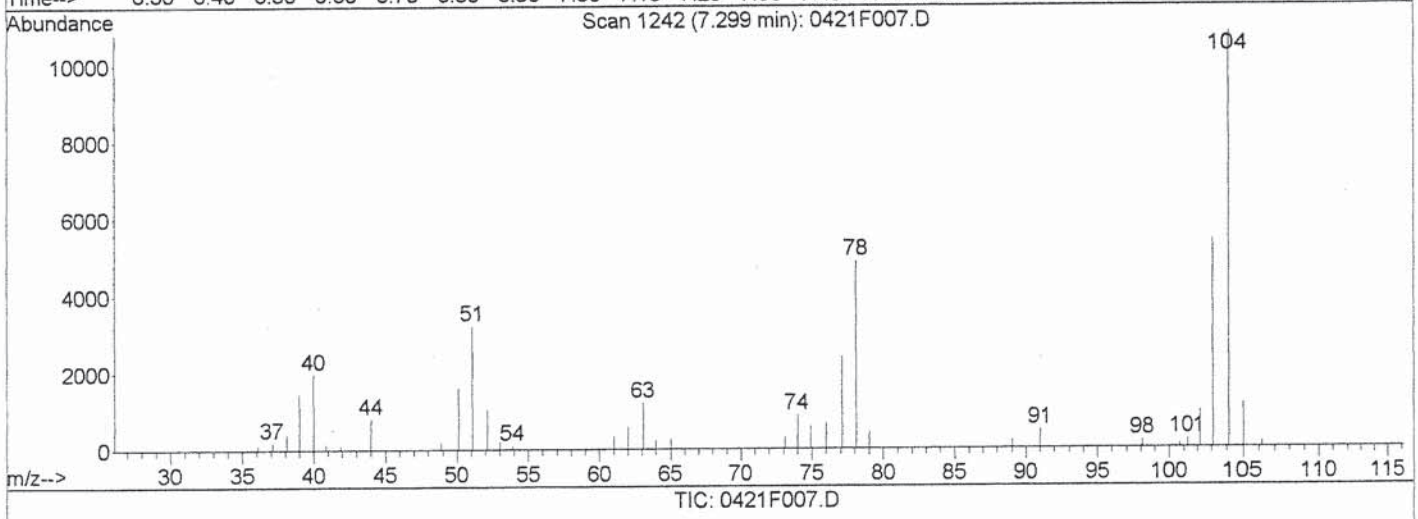
Data File : J:\MS24\DATA\042110\0421F007.D
Acq On : 21 Apr 2010 1:11 pm
Sample : SOIL ICAL 5
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:00 2010

Vial: 9
Operator: KR
Inst : MS24
Multiplr: 1.00

Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F007.D

(75) Styrene (T)

7.30min 4.83PPB

response 9894

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	200.48
78.00	90.20	90.18
0.00	0.00	0.00

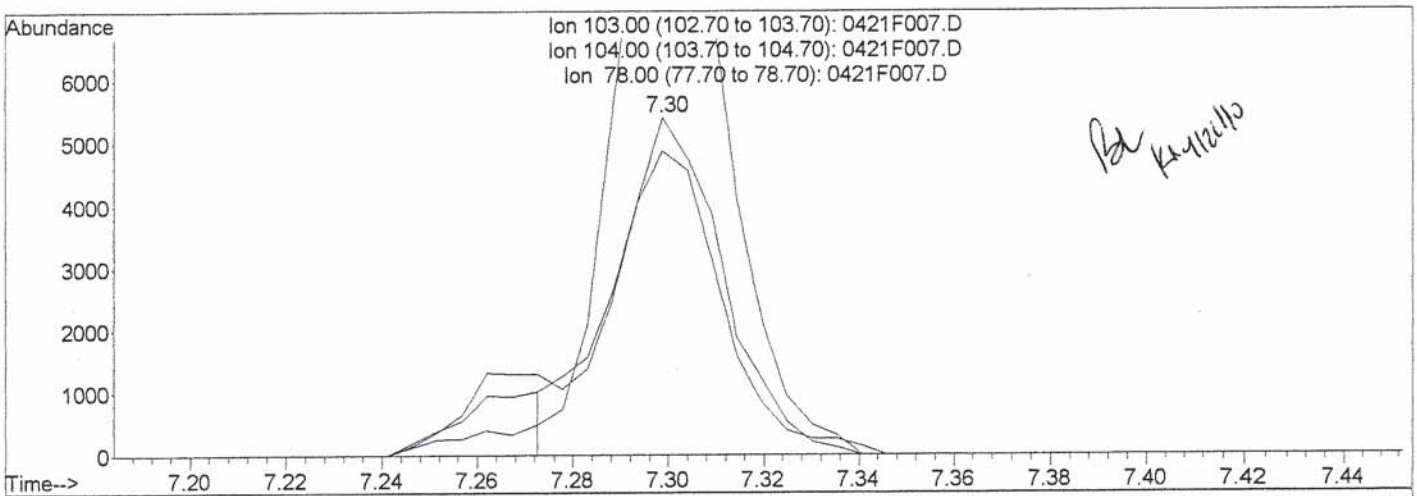
0421F007.D 042110MS24SOIL.M Wed Apr 21 14:01:04 2010

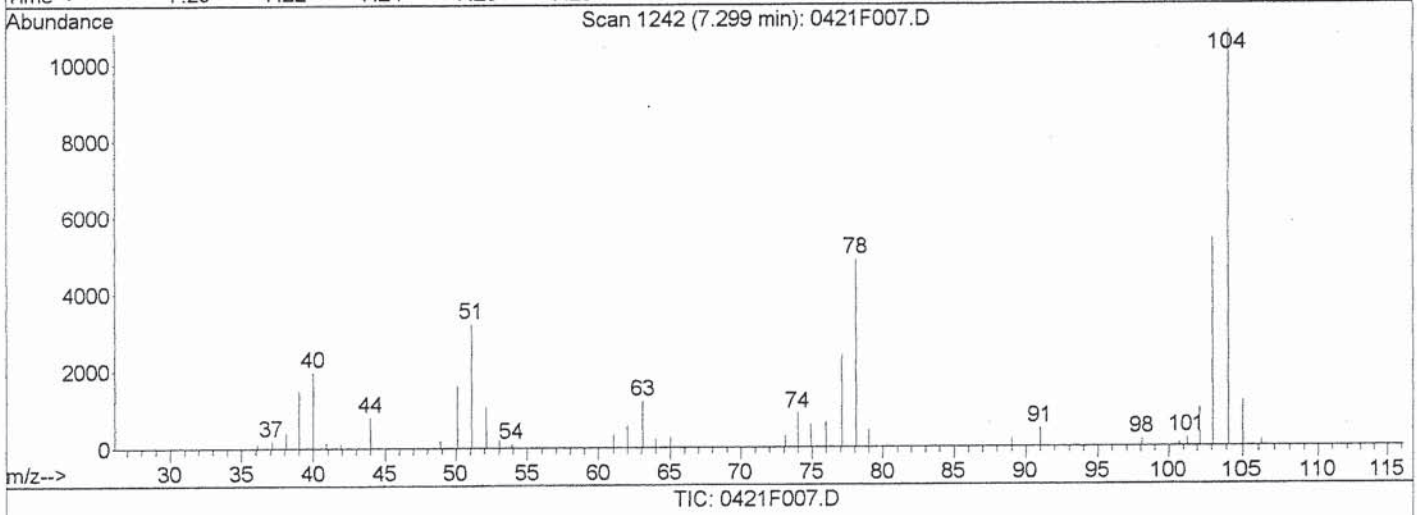
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F007.D
Acq On : 21 Apr 2010 1:11 pm
Sample : SOIL ICAL 5
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:01 2010

Vial: 9
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F007.D

(75) Styrene (T)

7.30min 4.21PPB m

response 8634

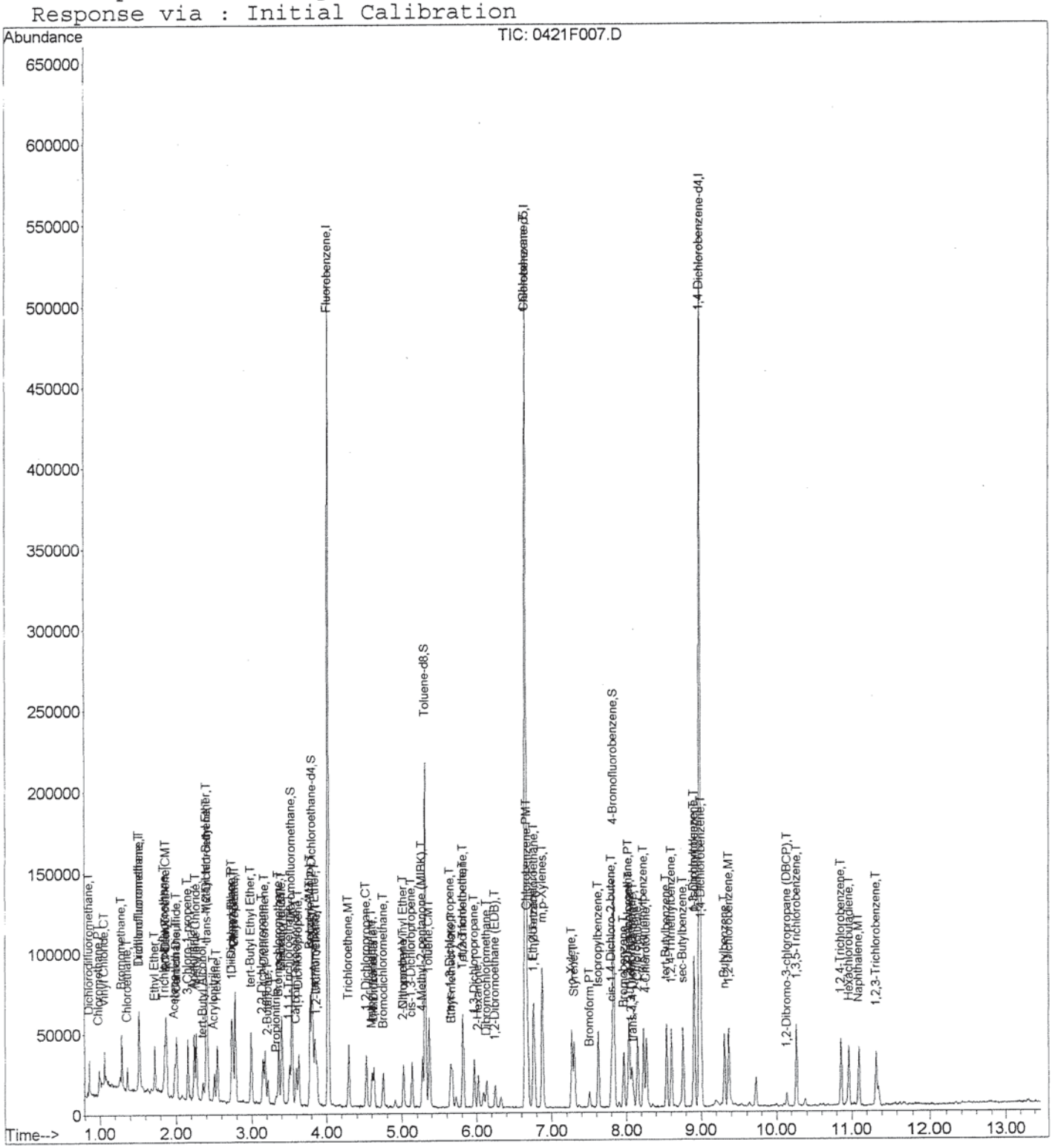
Ion	Exp%	Act%
103.00	100	100
104.00	200.50	200.48
78.00	90.20	90.18
0.00	0.00	0.00

0421F007.D 042110MS24SOIL.M Wed Apr 21 14:01:12 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F007.D Vial: 9
Acq On : 21 Apr 2010 1:11 pm Operator: KR
Sample : SOIL ICAL 5 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:01 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F008.D
Acq On : 21 Apr 2010 1:32 pm
Sample : SOIL ICAL 10
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:03:35 2010

Vial: 10
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene	4.01	96	323083	50.00	PPB		-0.01
59) Chlorobenzene-d5	6.65	82	130801	50.00	PPB		0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	119901	50.00	PPB		0.00
System Monitoring Compounds							
38) Dibromofluoromethane	3.52	113	48118	33.25	PPB		-0.01
Spiked Amount 50.000			Recovery =	66.50%			
42) 1,2-Dichloroethane-d4	3.80	65	69679	34.81	PPB		-0.01
Spiked Amount 50.000			Recovery =	69.62%			
57) Toluene-d8	5.30	98	163753	32.04	PPB		-0.01
Spiked Amount 50.000			Recovery =	64.08%			
79) 4-Bromofluorobenzene	7.82	95	56484	29.06	PPB		0.00
Spiked Amount 50.000			Recovery =	58.12%			
Target Compounds							Qvalue
2) Dichlorodifluoromethane	0.83	85	19345	13.90	PPB		98
3) Chloromethane	0.95	50	23839	12.53	PPB		98
4) Vinyl Chloride	1.03	62	22675	12.20	PPB		99
5) Bromomethane	1.26	96	18959	15.52	PPB		98
6) Chloroethane	1.33	64	14284	12.31	PPB		98
7) Dichlorofluoromethane	1.49	67	36517	11.81	PPB		98
8) Trichlorofluoromethane	1.48	101	30775	11.20	PPB		94
9) Ethyl Ether	1.70	59	17787	11.82	PPB		94
10) Acrolein	1.85	56	16308	54.36	PPB		87
11) Trichlorotrifluoroethane	1.83	151	11822	10.42	PPB		89
12) 1,1-Dichloroethene	1.85	96	13116	10.91	PPB		91
13) Acetone	1.96	43	19401	23.16	PPB		93
14) Iodomethane	1.97	142	27325	31.90	PPB		98
15) Carbon Disulfide	1.99	76	50785	11.15	PPB		99
16) 3-Chloro-1-propene	2.13	76	9413	11.41	PPB	#	84
17) Acetonitrile	2.21	40	29413	210.44	PPB		90
18) Methylene Chloride	2.24	84	19045	11.05	PPB		92
19) tert-Butyl Alcohol	2.34	59	12465	45.27	PPB		93
20) Acrylonitrile	2.48	53	13666	20.74	PPB		94
21) Methyl tert-Butyl Ether	2.38	73	102464	20.69	PPB		98
22) trans-1,2-Dichloroethene	2.40	96	15814	10.75	PPB		94
23) Hexane	2.52	57	22363	10.98	PPB		91
24) Diisopropyl Ether	2.72	45	58294	11.07	PPB		97
25) 1,1-Dichloroethane	2.73	63	32045	11.33	PPB		94
26) Vinyl Acetate	2.77	86	2695	8.71	PPB	#	65
27) Chloroprene	2.76	53	43022	19.94	PPB		94
28) tert-Butyl Ethyl Ether	2.98	59	55994	11.19	PPB		93

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F008.D
Acq On : 21 Apr 2010 1:32 pm
Sample : SOIL ICAL 10
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:03:35 2010

Vial: 10
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	2,2-Dichloropropane	3.13	77	21839	10.28	PPB		94
30)	cis-1,2-Dichloroethene	3.16	96	17790	10.74	PPB		93
31)	2-Butanone	3.20	72	4833	20.79	PPB		94
32)	Propionitrile	3.31	54	5734	20.51	PPB		93
33)	Methacrylonitrile	3.39	67	15090	20.08	PPB		93
34)	Bromochloromethane	3.35	128	8778	11.00	PPB	#	77
36)	Chloroform	3.40	83	30022	10.61	PPB		93
37)	1,1,1-Trichloroethane	3.49	97	23177	10.19	PPB		92
39)	Carbon Tetrachloride	3.58	117	19885	10.16	PPB		95
40)	1,1-Dichloropropene	3.61	75	21145	10.86	PPB		92
41)	Isobutyl Alcohol	3.77	43	17641	191.64	PPB		93
43)	Benzene	3.78	78	65104	10.91	PPB		98
44)	1,2-Dichloroethane	3.86	62	26823	11.41	PPB		98
45)	tert-Amyl Methyl Ether	3.83	55	13579	11.08	PPB		93
46)	Trichloroethene	4.28	95	16622	10.62	PPB		94
48)	1,2-Dichloropropane	4.52	63	16277	10.17	PPB		98
49)	Dibromomethane	4.62	93	10654	10.57	PPB		93
50)	Methyl methacrylate	4.59	69	11003	8.84	PPB		90
51)	1,4-Dioxane	4.62	88	4220	163.76	PPB		93
52)	Bromodichloromethane	4.74	83	20045	9.27	PPB		96
53)	2-Nitropropane	5.02	41	10891	25.57	PPB		95
54)	2-Chloroethyl Vinyl Ether	5.01	63	9818	8.84	PPB		95
55)	cis-1,3-Dichloropropene	5.13	75	21169	8.52	PPB		96
56)	4-Methyl-2-pentanone (MIBK	5.26	58	11805	17.40	PPB		98
58)	Toluene	5.36	92	38684	9.70	PPB		89
61)	trans-1,3-Dichloropropene	5.64	75	17890	8.10	PPB		87
62)	Ethyl methacrylate	5.67	69	17405	8.53	PPB		95
63)	1,1,2-Trichloroethane	5.80	83	12376	10.49	PPB		95
64)	Tetrachloroethene	5.81	164	12582	10.23	PPB		88
65)	2-Hexanone	6.02	57	4051	17.78	PPB		96
66)	1,3-Dichloropropane	5.96	76	25248	10.58	PPB		82
67)	Dibromochloromethane	6.13	129	12677	8.26	PPB		92
68)	1,2-Dibromoethane (EDB)	6.25	107	13787	9.91	PPB		94
69)	1-Chlorohexane	6.64	91	17246	9.01	PPB		95
70)	Chlorobenzene	6.68	112	42471	10.26	PPB		99
71)	Ethylbenzene	6.75	106	20968	9.60	PPB	#	83
72)	1,1,1,2-Tetrachloroethane	6.77	131	13144	8.85	PPB		95
73)	m,p-Xylenes	6.87	106	48376	18.90	PPB		94
74)	o-Xylene	7.27	106	23869	9.22	PPB		88
75)	Styrene	7.30	103	17049m	8.30	PPB		
76)	Bromoform	7.50	173	7400	7.62	PPB		85

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F008.D Vial: 10
Acq On : 21 Apr 2010 1:32 pm Operator: KR
Sample : SOIL ICAL 10 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:03:35 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
77)	Isopropylbenzene	7.62	105	58080	9.22	PPB		94
78)	cis-1,4-Dichloro-2-butene	7.80	89	2275	12.51	PPB	#	70
81)	1,1,2,2-Tetrachloroethane	8.02	83	14993	9.23	PPB		89
82)	trans-1,4-Dichloro-2-buten	8.10	53	3199	8.03	PPB		79
83)	Bromobenzene	7.96	156	15929	9.36	PPB		96
84)	n-Propylbenzene	8.04	91	66776	9.00	PPB		97
85)	1,2,3-Trichloropropane	8.07	110	5970	10.76	PPB	#	59
86)	2-Chlorotoluene	8.15	91	42547	9.68	PPB		97
87)	1,3,5-Trimethylbenzene	8.23	105	45414	8.74	PPB		95
88)	4-Chlorotoluene	8.27	91	42142	9.26	PPB		92
89)	tert-Butylbenzene	8.54	119	40406	8.77	PPB		96
90)	1,2,4-Trimethylbenzene	8.60	105	44697	8.54	PPB		87
91)	sec-Butylbenzene	8.76	105	61961	9.32	PPB		98
92)	p-Isopropyltoluene	8.91	119	51821	9.33	PPB		98
93)	1,3-Dichlorobenzene	8.90	146	31467	10.37	PPB		93
94)	1,4-Dichlorobenzene	9.00	146	32874	10.44	PPB		96
95)	n-Butylbenzene	9.31	91	45994	9.58	PPB		93
96)	1,2-Dichlorobenzene	9.36	146	31243	10.59	PPB		96
97)	1,2-Dibromo-3-chloropropan	10.14	155	2102	6.90	PPB		93
98)	1,3,5-Trichlorobenzene	10.26	180	23957	10.46	PPB		89
99)	1,2,4-Trichlorobenzene	10.86	180	20067	9.60	PPB		97
100)	Hexachlorobutadiene	10.96	225	11917	9.95	PPB		97
101)	Naphthalene	11.09	128	45398	8.57	PPB		95
102)	1,2,3-Trichlorobenzene	11.32	180	20338	10.62	PPB		89

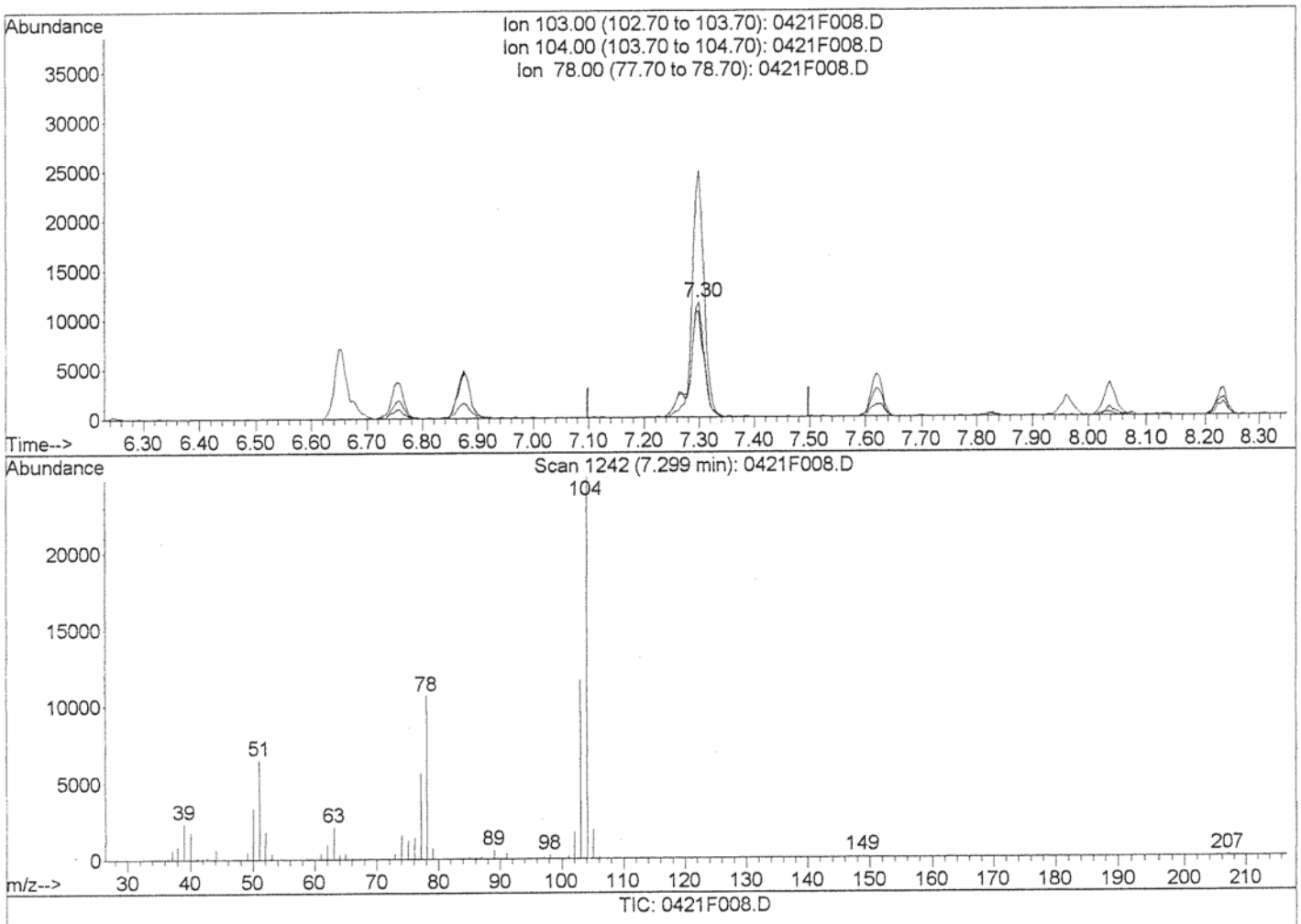
(#) = qualifier out of range (m) = manual integration

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F008.D
Acq On : 21 Apr 2010 1:32 pm
Sample : SOIL ICAL 10
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:05 2010

Vial: 10
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration



TIC: 0421F008.D

(75) Styrene (T)

7.30min 10.16PPB

response 20883

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	214.38
78.00	90.20	91.45
0.00	0.00	0.00

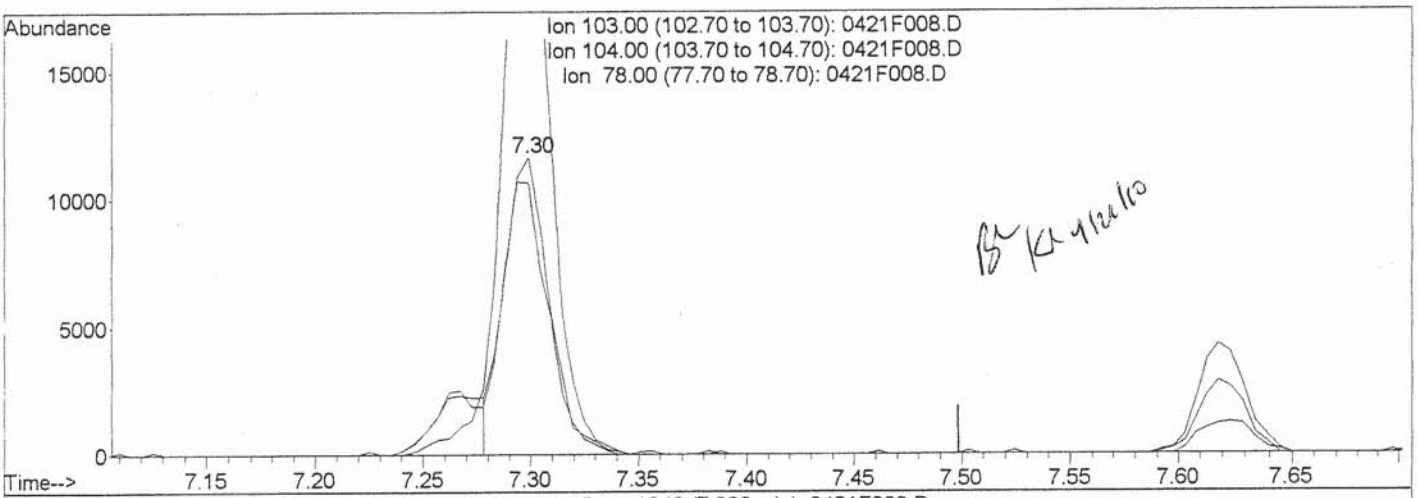
0421F008.D 042110MS24SOIL.M Wed Apr 21 14:05:42 2010

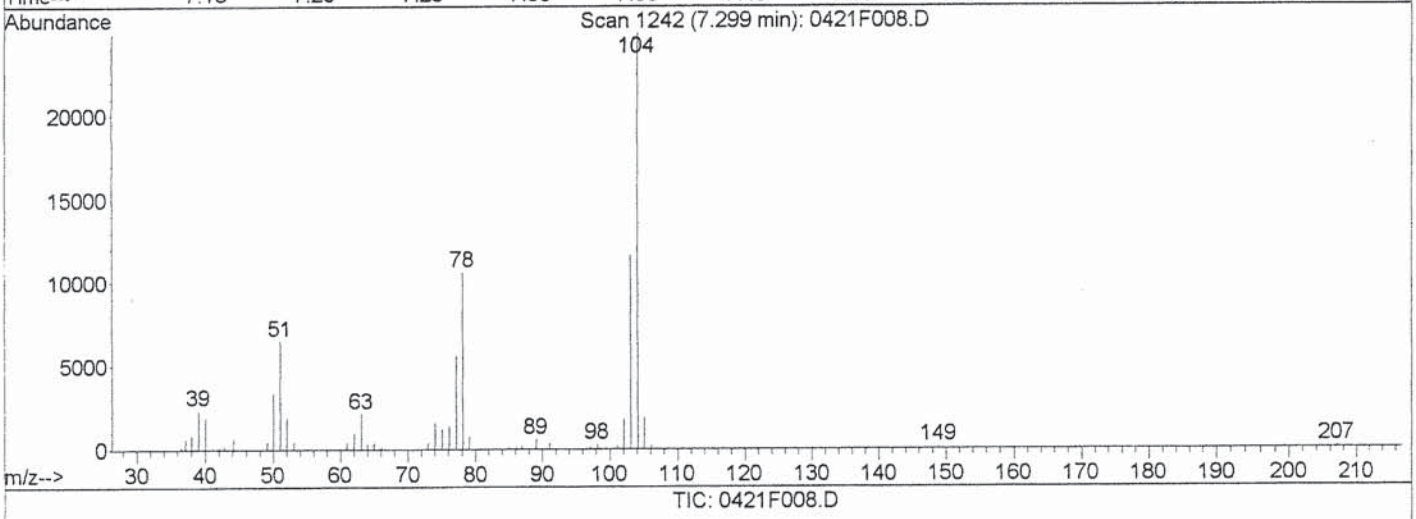
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F008.D
Acq On : 21 Apr 2010 1:32 pm
Sample : SOIL ICAL 10
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:05 2010

Vial: 10
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F008.D

(75) Styrene (T)

7.30min 8.30PPB m

response 17049

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	214.38
78.00	90.20	91.45
0.00	0.00	0.00

0421F008.D 042110MS24SOIL.M Wed Apr 21 14:05:55 2010

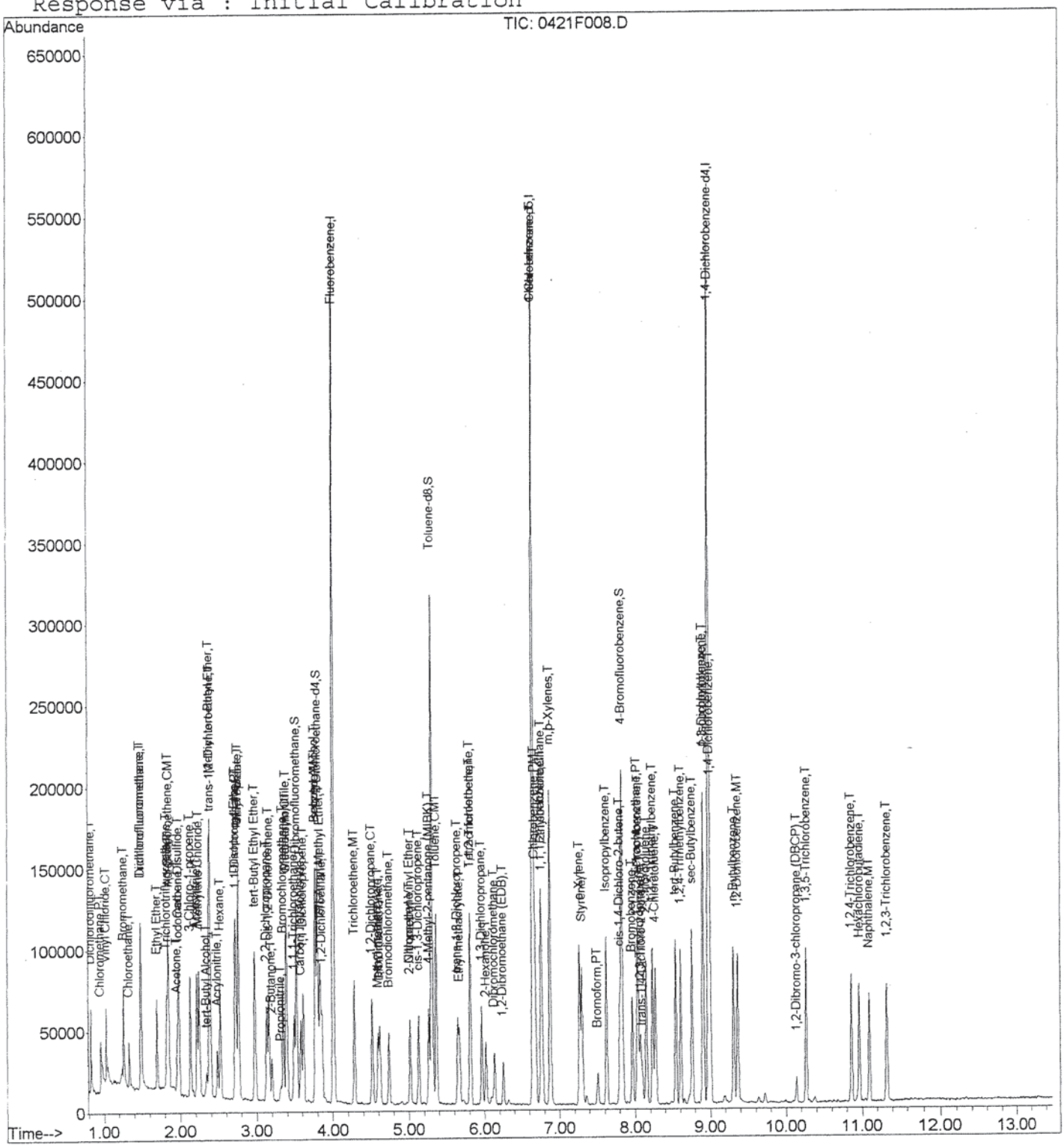
Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F008.D
Acq On : 21 Apr 2010 1:32 pm
Sample : SOIL ICAL 10
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:05 2010

Vial: 10
Operator: KR
Inst : MS24
Multiplr: 1.00

Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F009.D
Acq On : 21 Apr 2010 1:53 pm
Sample : SOIL ICAL 20
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:07:08 2010

Vial: 11
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene	4.02	96	327246	50.00	PPB		0.00
59) Chlorobenzene-d5	6.65	82	133223	50.00	PPB		0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	115766	50.00	PPB		0.00
System Monitoring Compounds							
38) Dibromofluoromethane	3.53	113	67092	45.76	PPB		0.00
Spiked Amount 50.000			Recovery =	91.52%			
42) 1,2-Dichloroethane-d4	3.81	65	97235	47.96	PPB		0.00
Spiked Amount 50.000			Recovery =	95.92%			
57) Toluene-d8	5.31	98	242612	46.86	PPB		0.00
Spiked Amount 50.000			Recovery =	93.72%			
79) 4-Bromofluorobenzene	7.83	95	84748	42.80	PPB		0.00
Spiked Amount 50.000			Recovery =	85.60%			
Target Compounds							Qvalue
2) Dichlorodifluoromethane	0.86	85	42049	29.83	PPB		97
3) Chloromethane	0.99	50	49571	25.73	PPB		94
4) Vinyl Chloride	1.06	62	47132	25.04	PPB		99
5) Bromomethane	1.29	96	38903	31.44	PPB		98
6) Chloroethane	1.36	64	31345	26.66	PPB		98
7) Dichlorofluoromethane	1.52	67	77999	24.91	PPB		99
8) Trichlorofluoromethane	1.51	101	66297	23.81	PPB		96
9) Ethyl Ether	1.73	59	39484	25.91	PPB		96
10) Acrolein	1.87	56	37052	121.94	PPB		91
11) Trichlorotrifluoroethane	1.86	151	25452	22.15	PPB		93
12) 1,1-Dichloroethene	1.88	96	28722	23.60	PPB		95
13) Acetone	1.98	43	30935	36.47	PPB		95
14) Iodomethane	2.00	142	75294	61.05	PPB		98
15) Carbon Disulfide	2.01	76	111882	24.26	PPB		97
16) 3-Chloro-1-propene	2.16	76	20331	24.32	PPB		99
17) Acetonitrile	2.24	40	61025	431.06	PPB		96
18) Methylene Chloride	2.27	84	41796	23.95	PPB		98
19) tert-Butyl Alcohol	2.36	59	24835	89.06	PPB		97
20) Acrylonitrile	2.50	53	30359	45.49	PPB		97
21) Methyl tert-Butyl Ether	2.40	73	236732	47.20	PPB		98
22) trans-1,2-Dichloroethene	2.42	96	34154	22.93	PPB		89
23) Hexane	2.54	57	47877	23.20	PPB		97
24) Diisopropyl Ether	2.74	45	130995	24.57	PPB		97
25) 1,1-Dichloroethane	2.75	63	68481	23.91	PPB		95
26) Vinyl Acetate	2.78	86	6461	20.63	PPB	#	75
27) Chloroprene	2.78	53	91164	41.71	PPB		97
28) tert-Butyl Ethyl Ether	2.99	59	120577	23.78	PPB		96

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F009.D
Acq On : 21 Apr 2010 1:53 pm
Sample : SOIL ICAL 20
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:07:08 2010

Vial: 11
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	2,2-Dichloropropane	3.15	77	46459	21.59	PPB		95
30)	cis-1,2-Dichloroethene	3.18	96	38918	23.19	PPB		99
31)	2-Butanone	3.21	72	9531	40.47	PPB		89
32)	Propionitrile	3.32	54	11659	41.18	PPB		93
33)	Methacrylonitrile	3.40	67	31272	41.07	PPB		93
34)	Bromochloromethane	3.36	128	18867	23.35	PPB		97
36)	Chloroform	3.41	83	65143	22.72	PPB		99
37)	1,1,1-Trichloroethane	3.50	97	51008	22.15	PPB		92
39)	Carbon Tetrachloride	3.59	117	44174	22.28	PPB		95
40)	1,1-Dichloropropene	3.62	75	46361	23.50	PPB		90
41)	Isobutyl Alcohol	3.78	43	36913	395.89	PPB		96
43)	Benzene	3.79	78	147543	24.41	PPB		98
44)	1,2-Dichloroethane	3.87	62	59077	24.81	PPB		98
45)	tert-Amyl Methyl Ether	3.84	55	30741	24.76	PPB		91
46)	Trichloroethene	4.29	95	39299	24.78	PPB		97
47)	Ethyl Acrylate	4.37	55	521	0.24	PPB	#	76
48)	1,2-Dichloropropane	4.53	63	38558	23.78	PPB		98
49)	Dibromomethane	4.62	93	24825	24.32	PPB		94
50)	Methyl methacrylate	4.60	69	26321	20.88	PPB		91
51)	1,4-Dioxane	4.62	88	9778	374.62	PPB		97
52)	Bromodichloromethane	4.75	83	48148	21.99	PPB		95
53)	2-Nitropropane	5.02	41	25875	59.99	PPB		91
54)	2-Chloroethyl Vinyl Ether	5.02	63	23080	20.51	PPB		93
55)	cis-1,3-Dichloropropene	5.13	75	53351	21.20	PPB		90
56)	4-Methyl-2-pentanone (MIBK	5.27	58	28334	41.23	PPB		99
58)	Toluene	5.36	92	90753	22.47	PPB		99
61)	trans-1,3-Dichloropropene	5.65	75	46711	20.77	PPB		84
62)	Ethyl methacrylate	5.67	69	44466	21.38	PPB		99
63)	1,1,2-Trichloroethane	5.81	83	29747	24.75	PPB		95
64)	Tetrachloroethene	5.82	164	28624	22.85	PPB		86
65)	2-Hexanone	6.02	57	9019	38.87	PPB	#	69
66)	1,3-Dichloropropane	5.97	76	59989	24.68	PPB		89
67)	Dibromochloromethane	6.13	129	30849	19.74	PPB		95
68)	1,2-Dibromoethane (EDB)	6.25	107	31497	22.23	PPB		95
69)	1-Chlorohexane	6.64	91	38965	20.00	PPB		90
70)	Chlorobenzene	6.68	112	96405	22.87	PPB		100
71)	Ethylbenzene	6.76	106	48138	21.64	PPB		86
72)	1,1,1,2-Tetrachloroethane	6.77	131	31064	20.54	PPB		92
73)	m,p-Xylenes	6.87	106	114110	43.77	PPB		95
74)	o-Xylene	7.27	106	55056	20.89	PPB		87
75)	Styrene	7.30	103	42839m	20.47	PPB		

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F009.D
Acq On : 21 Apr 2010 1:53 pm
Sample : SOIL ICAL 20
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:07:08 2010

Vial: 11
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
76)	Bromoform	7.51	173	18073	18.28	PPB		88
77)	Isopropylbenzene	7.62	105	134343	20.94	PPB		94
78)	cis-1,4-Dichloro-2-butene	7.79	89	6331	34.18	PPB	#	81
81)	1,1,2,2-Tetrachloroethane	8.02	83	34017	21.69	PPB		92
82)	trans-1,4-Dichloro-2-buten	8.10	53	8182	21.26	PPB		95
83)	Bromobenzene	7.96	156	38189	23.24	PPB		94
84)	n-Propylbenzene	8.04	91	159146	22.22	PPB		97
85)	1,2,3-Trichloropropane	8.07	110	12805	23.91	PPB	#	69
86)	2-Chlorotoluene	8.15	91	100004	23.57	PPB		94
87)	1,3,5-Trimethylbenzene	8.23	105	111868	22.29	PPB		98
88)	4-Chlorotoluene	8.27	91	98821	22.50	PPB		99
89)	tert-Butylbenzene	8.54	119	96911	21.77	PPB		98
90)	1,2,4-Trimethylbenzene	8.60	105	114160	22.59	PPB		88
91)	sec-Butylbenzene	8.76	105	149839	23.33	PPB		98
92)	p-Isopropyltoluene	8.91	119	124522	23.22	PPB		97
93)	1,3-Dichlorobenzene	8.90	146	72004	24.58	PPB		99
94)	1,4-Dichlorobenzene	9.00	146	75085	24.70	PPB		91
95)	n-Butylbenzene	9.31	91	109471	23.62	PPB		95
96)	1,2-Dichlorobenzene	9.36	146	75958	26.67	PPB		97
97)	1,2-Dibromo-3-chloropropan	10.14	155	5504	18.70	PPB		92
98)	1,3,5-Trichlorobenzene	10.26	180	55421	25.06	PPB		91
99)	1,2,4-Trichlorobenzene	10.85	180	46894	23.24	PPB		90
100)	Hexachlorobutadiene	10.96	225	26977	23.32	PPB		99
101)	Naphthalene	11.09	128	106995	20.92	PPB		98
102)	1,2,3-Trichlorobenzene	11.32	180	48715	26.35	PPB		87

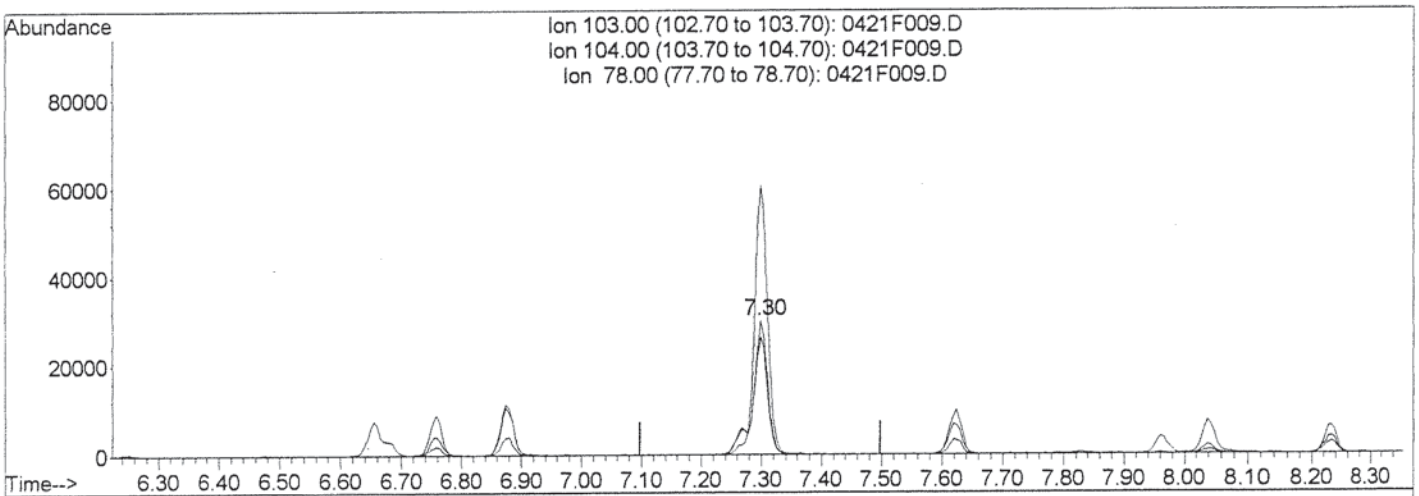
(#) = qualifier out of range (m) = manual integration

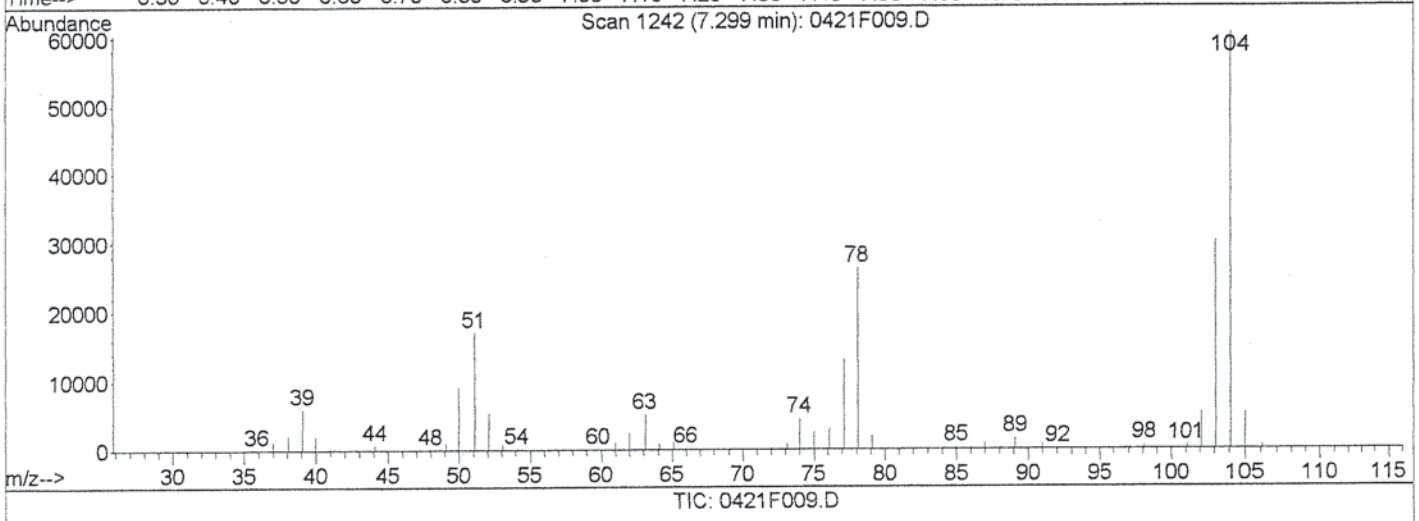
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F009.D
Acq On : 21 Apr 2010 1:53 pm
Sample : SOIL ICAL 20
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:07 2010

Vial: 11
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F009.D

(75) Styrene (T)

7.30min 24.23PPB

response 50720

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	201.01
78.00	90.20	87.08
0.00	0.00	0.00

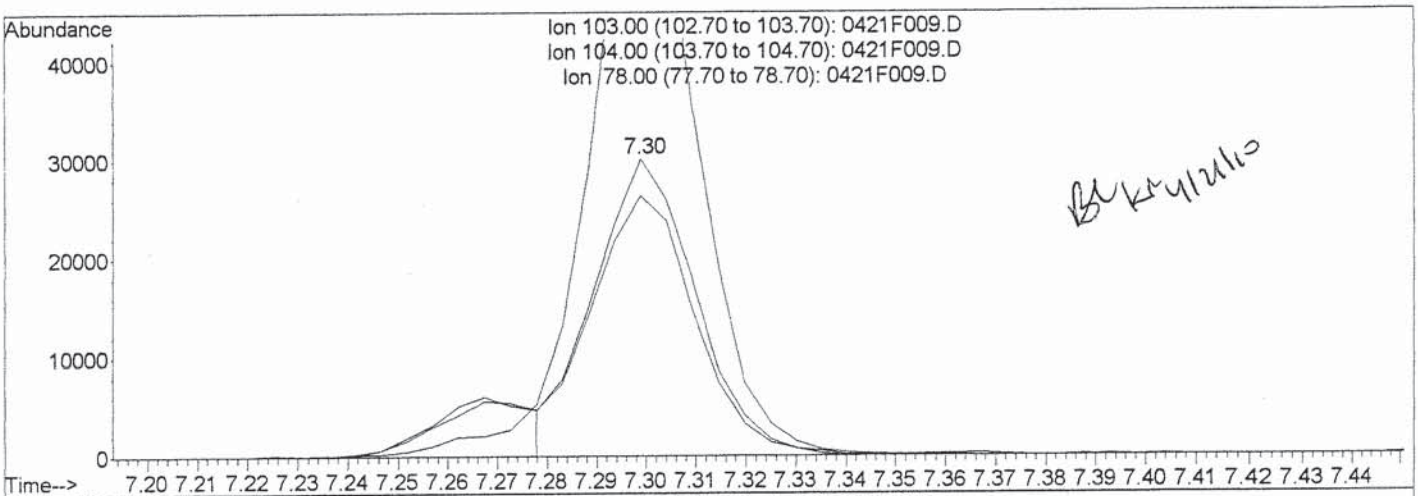
0421F009.D 042110MS24SOIL.M Wed Apr 21 14:08:02 2010

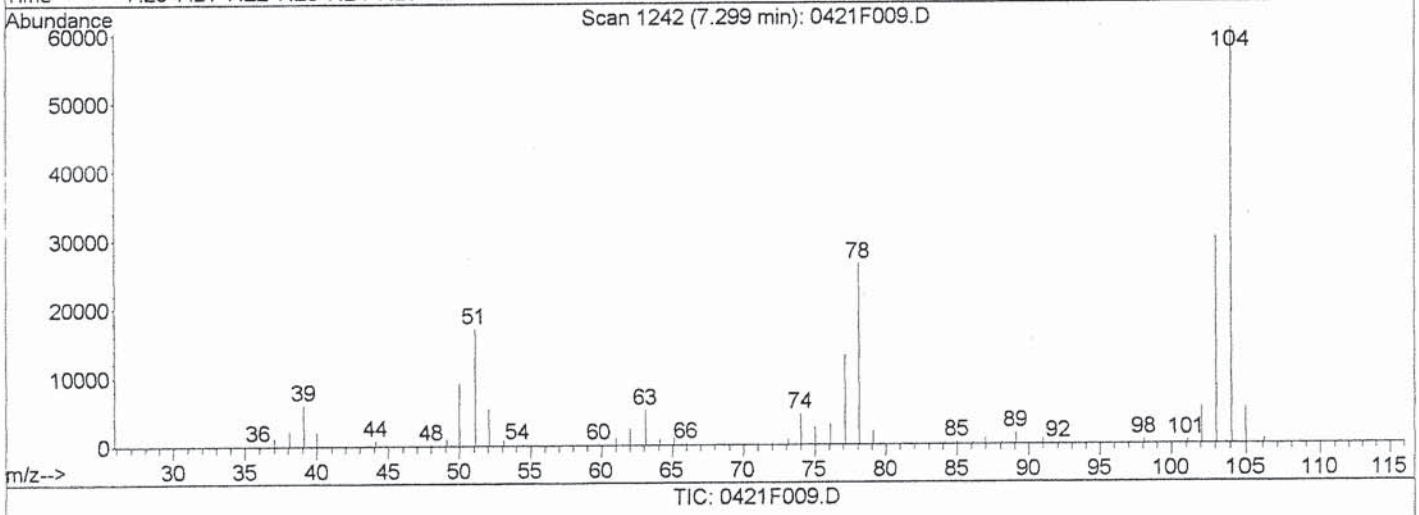
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F009.D
Acq On : 21 Apr 2010 1:53 pm
Sample : SOIL ICAL 20
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:08 2010

Vial: 11
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Single Level Calibration





TIC: 0421F009.D

(75) Styrene (T)

7.30min 20.47PPB m

response 42839

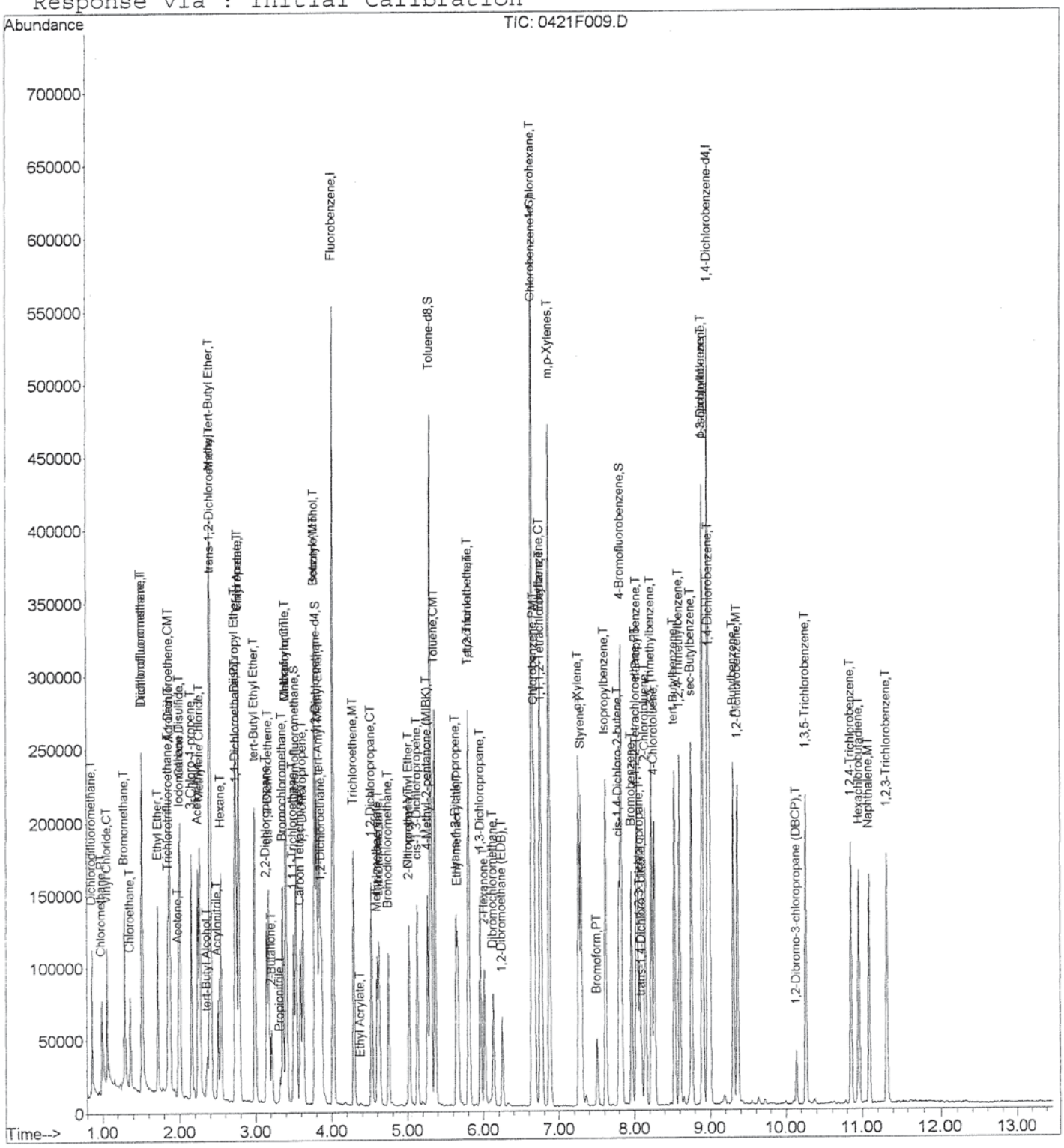
Ion	Exp%	Act%
103.00	100	100
104.00	200.50	201.01
78.00	90.20	87.49
0.00	0.00	0.00

0421F009.D 042110MS24SOIL.M Wed Apr 21 14:08:10 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F009.D Vial: 11
Acq On : 21 Apr 2010 1:53 pm Operator: KR
Sample : SOIL ICAL 20 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:08 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 13:52:07 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F010.D
Acq On : 21 Apr 2010 2:15 pm
Sample : SOIL ICAL 50
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:31:03 2010

Vial: 12
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:10:01 2010
Response via : Initial Calibration
DataAcq Meth : 8260

KR 4/21/10

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene	4.02	96	321724	50.00	PPB		0.00
59) Chlorobenzene-d5	6.65	82	137440	50.00	PPB		0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	130674	50.00	PPB		0.00
System Monitoring Compounds							
38) Dibromofluoromethane	3.53	113	77159	48.46	PPB		0.00
Spiked Amount 50.000			Recovery =	96.92%			
42) 1,2-Dichloroethane-d4	3.81	65	111265	47.89	PPB		0.00
Spiked Amount 50.000			Recovery =	95.78%			
57) Toluene-d8	5.31	98	271397	49.85	PPB		0.00
Spiked Amount 50.000			Recovery =	99.70%			
79) 4-Bromofluorobenzene	7.83	95	111120	54.41	PPB		0.00
Spiked Amount 50.000			Recovery =	108.82%			
Target Compounds							Qvalue
2) Dichlorodifluoromethane	0.85	85	90628	44.62	PPB		96
3) Chloromethane	0.98	50	112745	43.17	PPB		94
4) Vinyl Chloride	1.05	62	102580	44.77	PPB		96
5) Bromomethane	1.28	96	74995	38.23	PPB		98
6) Chloroethane	1.36	64	67234	45.50	PPB		95
7) Dichlorofluoromethane	1.51	67	171027	46.63	PPB		98
8) Trichlorofluoromethane	1.51	101	138536	44.36	PPB		96
9) Ethyl Ether	1.72	59	87788	47.57	PPB		96
10) Acrolein	1.87	56	83933	229.99	PPB		94
11) Trichlorotrifluoroethane	1.85	151	53517	44.63	PPB		98
12) 1,1-Dichloroethene	1.87	96	61320	40.90	PPB		98
13) Acetone	1.98	43	69245	66.95	PPB		98
14) Iodomethane	1.99	142	184618	125.94	PPB		97
15) Carbon Disulfide	2.00	76	238783	45.39	PPB		97
16) 3-Chloro-1-propene	2.15	76	44001	45.64	PPB		91
17) Acetonitrile	2.23	40	140525	945.06	PPB		97
18) Methylene Chloride	2.26	84	90903	45.27	PPB		95
19) tert-Butyl Alcohol	2.35	59	56566	188.51	PPB		96
20) Acrylonitrile	2.50	53	69152	90.72	PPB		96
21) Methyl tert-Butyl Ether	2.40	73	532941	96.56	PPB		98
22) trans-1,2-Dichloroethene	2.41	96	76042	44.44	PPB		88
23) Hexane	2.54	57	102835	42.76	PPB		97
24) Diisopropyl Ether	2.73	45	298570	47.59	PPB		96
25) 1,1-Dichloroethane	2.75	63	152770	45.72	PPB		97
26) Vinyl Acetate	2.78	86	14390	47.11	PPB	#	62
27) Chloroprene	2.78	53	210165	95.45	PPB		96
28) tert-Butyl Ethyl Ether	2.99	59	278977	49.50	PPB		94

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F010.D
Acq On : 21 Apr 2010 2:15 pm
Sample : SOIL ICAL 50
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:31:03 2010

Vial: 12
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:10:01 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	2,2-Dichloropropane	3.15	77	107169	48.42	PPB		99
30)	cis-1,2-Dichloroethene	3.17	96	91630	49.43	PPB		95
31)	2-Butanone	3.21	72	22659	89.97	PPB		91
32)	Propionitrile	3.32	54	27337	86.44	PPB		90
33)	Methacrylonitrile	3.40	67	75111	98.38	PPB		89
34)	Bromochloromethane	3.36	128	44320	49.80	PPB		87
36)	Chloroform	3.41	83	147468	48.30	PPB		98
37)	1,1,1-Trichloroethane	3.50	97	110512	46.76	PPB		95
39)	Carbon Tetrachloride	3.59	117	92304	45.44	PPB		93
40)	1,1-Dichloropropene	3.62	75	96165	44.58	PPB		96
41)	Isobutyl Alcohol	3.78	43	89901	946.75	PPB		91
43)	Benzene	3.78	78	321162	47.68	PPB		97
44)	1,2-Dichloroethane	3.87	62	133395	48.67	PPB		97
45)	tert-Amyl Methyl Ether	3.84	55	69334	45.90	PPB	#	77
46)	Trichloroethene	4.29	95	83749	45.67	PPB		87
48)	1,2-Dichloropropane	4.53	63	86870	48.27	PPB		98
49)	Dibromomethane	4.62	93	55942	49.43	PPB		92
50)	Methyl methacrylate	4.60	69	60392	52.13	PPB		94
51)	1,4-Dioxane	4.62	88	22968	990.42	PPB		98
52)	Bromodichloromethane	4.75	83	110857	52.22	PPB		99
53)	2-Nitropropane	5.02	41	63266	141.63	PPB		97
54)	2-Chloroethyl Vinyl Ether	5.02	63	53630	51.77	PPB		91
55)	cis-1,3-Dichloropropene	5.13	75	123457	54.32	PPB		93
56)	4-Methyl-2-pentanone (MIBK	5.27	58	69584	107.06	PPB		96
58)	Toluene	5.36	92	193408	47.44	PPB		97
61)	trans-1,3-Dichloropropene	5.65	75	111122	54.61	PPB		89
62)	Ethyl methacrylate	5.67	69	105513	53.51	PPB		97
63)	1,1,2-Trichloroethane	5.81	83	66254	44.96	PPB		93
64)	Tetrachloroethene	5.81	164	61640	44.18	PPB		91
65)	2-Hexanone	6.02	57	23232	99.68	PPB	#	64
66)	1,3-Dichloropropane	5.97	76	135279	47.10	PPB		90
67)	Dibromochloromethane	6.13	129	76188	53.22	PPB		95
68)	1,2-Dibromoethane (EDB)	6.25	107	74710	48.40	PPB		95
69)	1-Chlorohexane	6.64	91	87440	44.91	PPB		91
70)	Chlorobenzene	6.68	112	223708	46.73	PPB		99
71)	Ethylbenzene	6.76	106	109279	48.28	PPB	#	80
72)	1,1,1,2-Tetrachloroethane	6.77	131	74906	51.31	PPB		93
73)	m,p-Xylenes	6.87	106	270760	105.23	PPB		95
74)	o-Xylene	7.27	106	130278	50.26	PPB		88
75)	Styrene	7.30	103	106647m	54.96	PPB		
76)	Bromoform	7.51	173	48675	58.54	PPB		89

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F010.D Vial: 12
Acq On : 21 Apr 2010 2:15 pm Operator: KR
Sample : SOIL ICAL 50 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:31:03 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:10:01 2010
Response via : Initial Calibration
DataAcq Meth : 8260

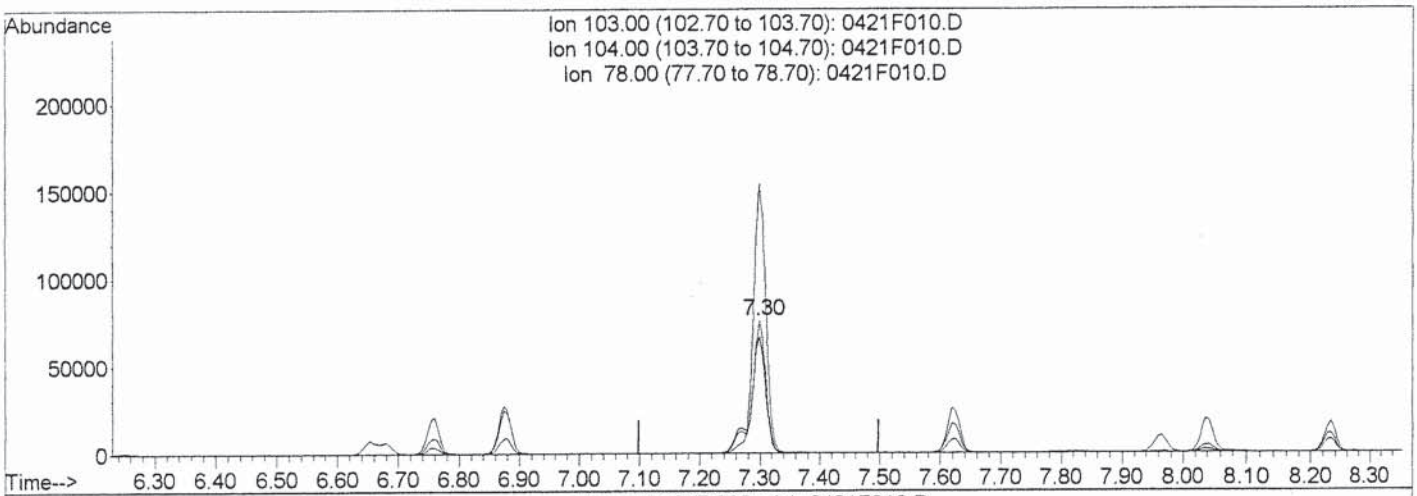
	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
77)	Isopropylbenzene	7.62	105	338628	54.41	PPB		96
78)	cis-1,4-Dichloro-2-butene	7.80	89	19697	130.58	PPB	#	79
81)	1,1,2,2-Tetrachloroethane	8.03	83	89412	45.53	PPB		90
82)	trans-1,4-Dichloro-2-buten	8.10	53	22104	53.83	PPB		84
83)	Bromobenzene	7.96	156	99691	48.84	PPB		93
84)	n-Propylbenzene	8.04	91	420480	54.18	PPB		97
85)	1,2,3-Trichloropropane	8.07	110	33552	43.90	PPB	#	57
86)	2-Chlorotoluene	8.15	91	258451	50.14	PPB		95
87)	1,3,5-Trimethylbenzene	8.23	105	297211	54.55	PPB		99
88)	4-Chlorotoluene	8.27	91	263827	50.93	PPB		96
89)	tert-Butylbenzene	8.54	119	256444	53.27	PPB		99
90)	1,2,4-Trimethylbenzene	8.60	105	301887	55.22	PPB		89
91)	sec-Butylbenzene	8.76	105	375122	51.95	PPB		99
92)	p-Isopropyltoluene	8.91	119	313474	51.98	PPB		96
93)	1,3-Dichlorobenzene	8.90	146	184199	48.87	PPB		99
94)	1,4-Dichlorobenzene	9.00	146	184788	46.32	PPB		90
95)	n-Butylbenzene	9.31	91	273563	51.73	PPB		95
96)	1,2-Dichlorobenzene	9.36	146	181618	49.26	PPB		98
97)	1,2-Dibromo-3-chloropropan	10.14	155	12286	45.53	PPB		80
98)	1,3,5-Trichlorobenzene	10.26	180	105104	38.90	PPB		97
99)	1,2,4-Trichlorobenzene	10.86	180	89767	39.47	PPB		94
100)	Hexachlorobutadiene	10.96	225	46548	36.96	PPB		94
101)	Naphthalene	11.09	128	226410	41.22	PPB		96
102)	1,2,3-Trichlorobenzene	11.32	180	89863	41.44	PPB		93

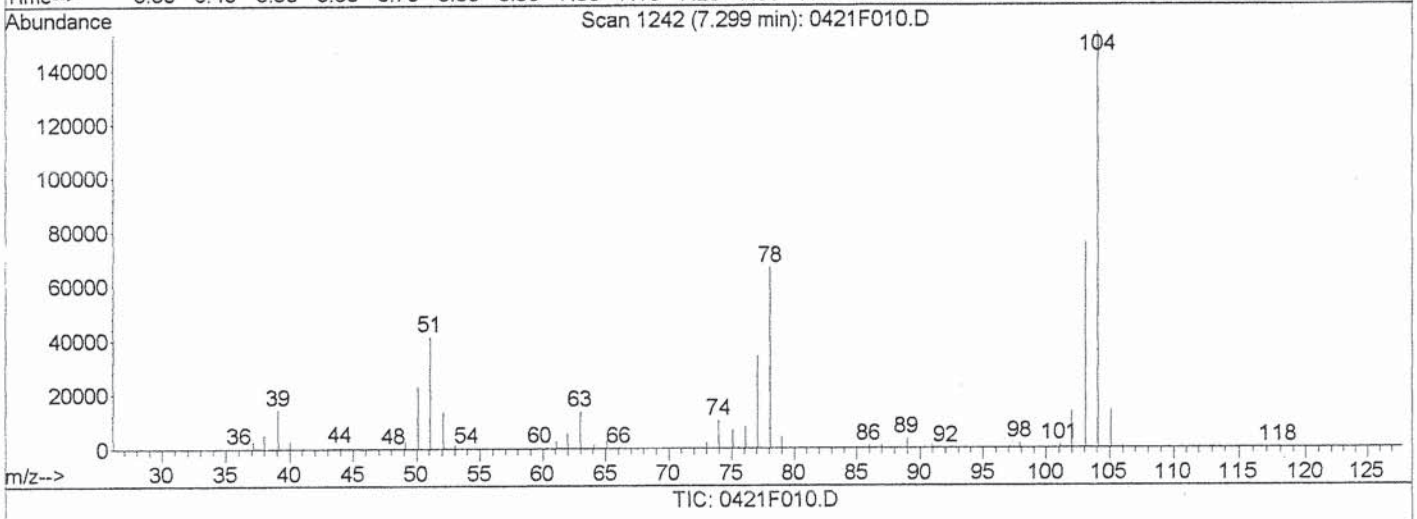
(#) = qualifier out of range (m) = manual integration

Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F010.D Vial: 12
Acq On : 21 Apr 2010 2:15 pm Operator: KR
Sample : SOIL ICAL 50 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:31 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:10:01 2010
Response via : Single Level Calibration





TIC: 0421F010.D

(75) Styrene (T)

7.30min 64.27PPB

response 124713

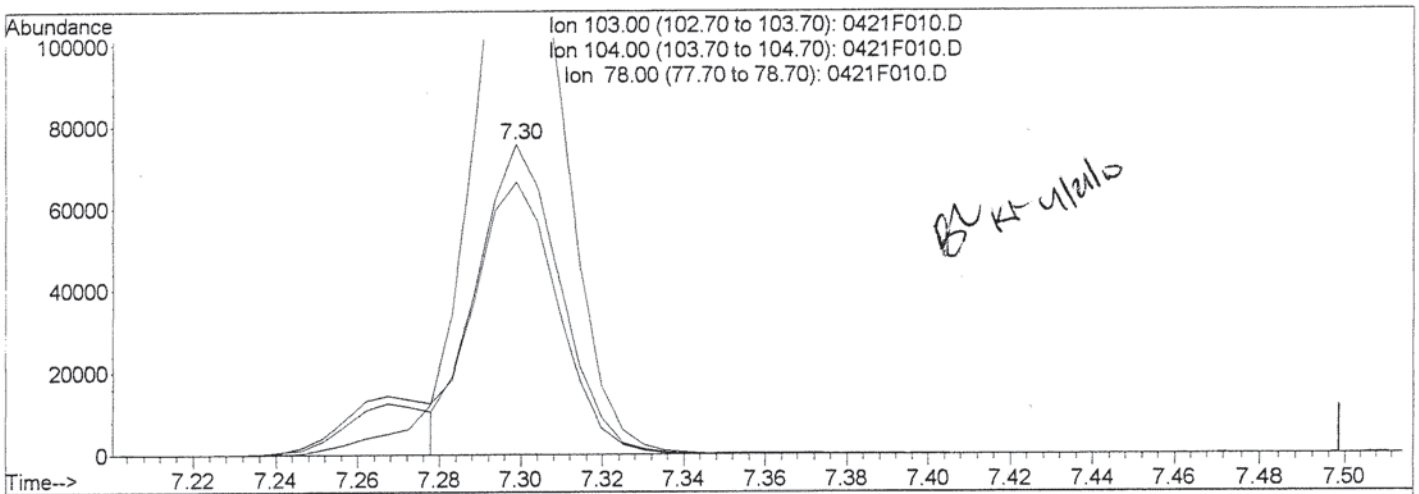
Ion	Exp%	Act%
103.00	100	100
104.00	200.50	203.21
78.00	90.20	87.77
0.00	0.00	0.00

0421F010.D 042110MS24SOIL.M Wed Apr 21 14:31:58 2010

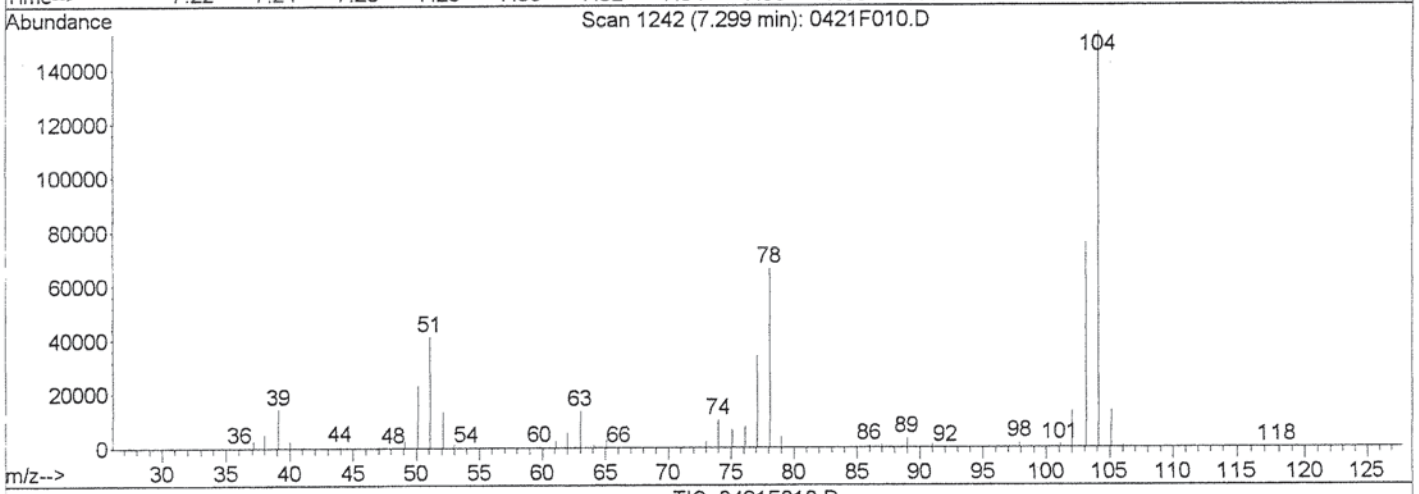
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F010.D Vial: 12
Acq On : 21 Apr 2010 2:15 pm Operator: KR
Sample : SOIL ICAL 50 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:32 2010 Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:10:01 2010
Response via : Single Level Calibration



BL KR 4/21/10



TIC: 0421F010.D

(75) Styrene (T)

7.30min 54.96PPB m

response 106647

Ion	Exp%	Act%
103.00	100	100
104.00	200.50	203.21
78.00	90.20	87.77
0.00	0.00	0.00

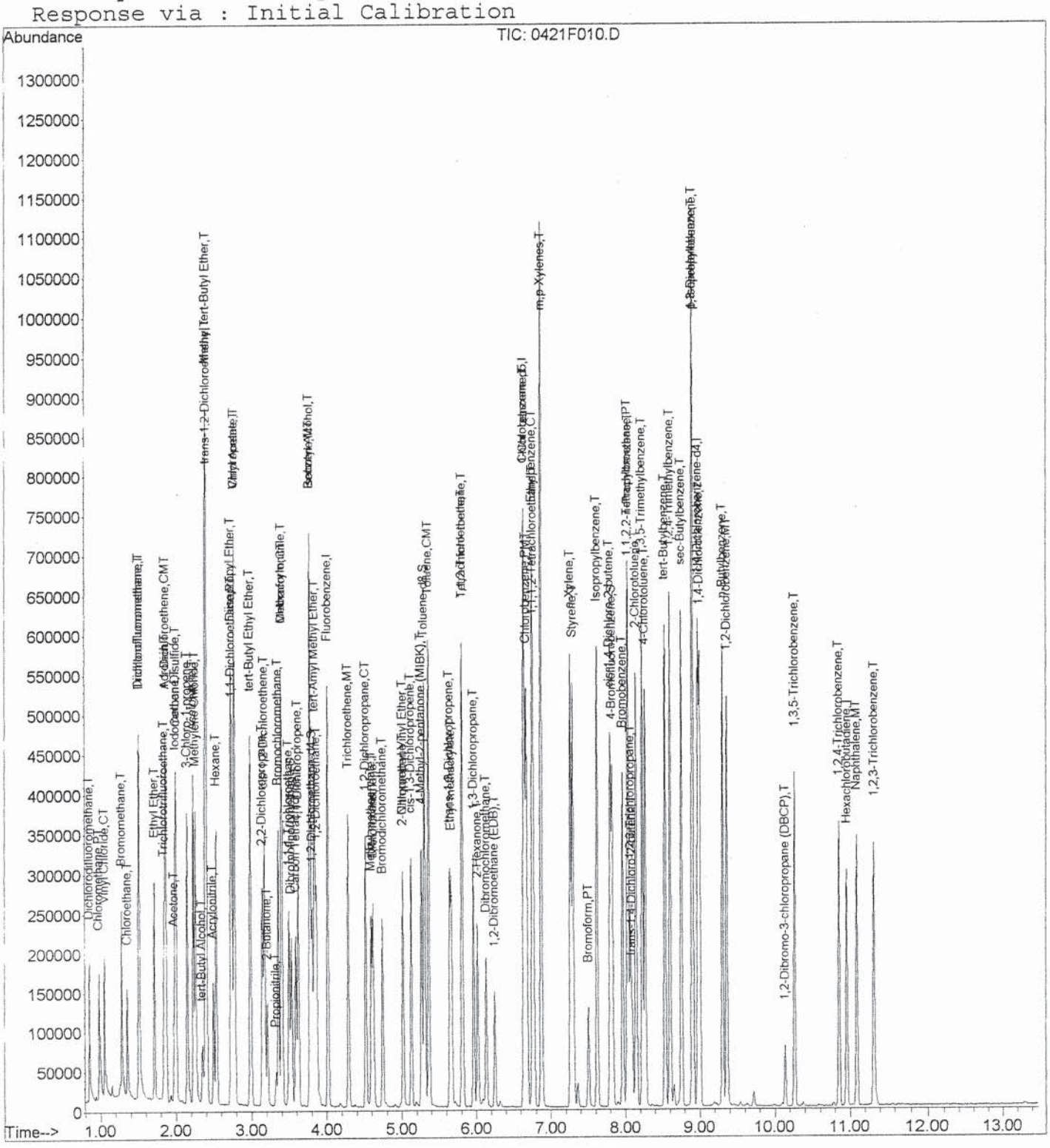
4/26/10

0421F010.D 042110MS24SOIL.M Wed Apr 21 14:32:04 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F010.D Vial: 12
Acq On : 21 Apr 2010 2:15 pm Operator: KR
Sample : SOIL ICAL 50 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:32 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:10:01 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F011.D
Acq On : 21 Apr 2010 2:36 pm
Sample : SOIL ICAL 100
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:52:53 2010

Vial: 13
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:32:32 2010
Response via : Initial Calibration
DataAcq Meth : 8260

KR 7/22/10

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene	4.02	96	326252	50.00	PPB		0.00
59) Chlorobenzene-d5	6.65	82	143154	50.00	PPB		0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	131664	50.00	PPB		0.00
System Monitoring Compounds							
38) Dibromofluoromethane	3.53	113	117239	73.06	PPB		0.00
Spiked Amount 50.000			Recovery = 146.12%				
42) 1,2-Dichloroethane-d4	3.81	65	163394	69.94	PPB		0.00
Spiked Amount 50.000			Recovery = 139.88%				
57) Toluene-d8	5.31	98	429109	77.78	PPB		0.00
Spiked Amount 50.000			Recovery = 155.56%				
79) 4-Bromofluorobenzene	7.83	95	153146	70.75	PPB		0.00
Spiked Amount 50.000			Recovery = 141.50%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	0.85	85	200916	99.69	PPB		97
3) Chloromethane	0.98	50	250911	96.94	PPB		95
4) Vinyl Chloride	1.05	62	228235	100.33	PPB		95
5) Bromomethane	1.28	96	183599	98.07	PPB		98
6) Chloroethane	1.35	64	150419	102.69	PPB		95
7) Dichlorofluoromethane	1.51	67	381235	103.67	PPB		96
8) Trichlorofluoromethane	1.50	101	313851	101.38	PPB		97
9) Ethyl Ether	1.71	59	189895	102.31	PPB		96
10) Acrolein	1.87	56	170835	467.86	PPB		92
11) Trichlorotrifluoroethane	1.85	151	116908	98.26	PPB		96
12) 1,1-Dichloroethene	1.87	96	133621	90.63	PPB		95
13) Acetone	1.98	43	139877	145.38	PPB		97
14) Iodomethane	1.99	142	445634	281.51	PPB		98
15) Carbon Disulfide	2.00	76	552601	105.21	PPB		98
16) 3-Chloro-1-propene	2.15	76	98970	103.49	PPB		99
17) Acetonitrile	2.23	40	293557	1968.47	PPB		96
18) Methylene Chloride	2.26	84	195378	98.28	PPB		95
19) tert-Butyl Alcohol	2.35	59	117463	405.99	PPB		95
20) Acrylonitrile	2.50	53	144912	190.42	PPB		93
21) Methyl tert-Butyl Ether	2.40	73	1166099	209.54	PPB		98
22) trans-1,2-Dichloroethene	2.41	96	170416	100.07	PPB		82
23) Hexane	2.54	57	223004	93.70	PPB		96
24) Diisopropyl Ether	2.73	45	667381	105.76	PPB		96
25) 1,1-Dichloroethane	2.75	63	344910	103.27	PPB		95
26) Vinyl Acetate	2.78	86	32095	105.13	PPB	#	79
27) Chloroprene	2.78	53	478463	215.93	PPB		94
28) tert-Butyl Ethyl Ether	2.99	59	611931	107.26	PPB		95

(#) = qualifier out of range (m) = manual integration
0421F011.D 042110MS24SOIL.M Wed Apr 21 14:53:46 2010 Page 1

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F011.D Vial: 13
Acq On : 21 Apr 2010 2:36 pm Operator: KR
Sample : SOIL ICAL 100 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:52:53 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:32:32 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	2,2-Dichloropropane	3.15	77	241199	108.32	PPB		97
30)	cis-1,2-Dichloroethene	3.17	96	196276	104.61	PPB		93
31)	2-Butanone	3.21	72	46997	188.75	PPB	#	87
32)	Propionitrile	3.32	54	57489	184.25	PPB		91
33)	Methacrylonitrile	3.40	67	167296	216.79	PPB		88
34)	Bromochloromethane	3.36	128	97528	108.17	PPB		92
36)	Chloroform	3.41	83	333797	108.56	PPB		96
37)	1,1,1-Trichloroethane	3.50	97	252256	106.63	PPB		95
39)	Carbon Tetrachloride	3.59	117	213316	105.97	PPB		94
40)	1,1-Dichloropropene	3.62	75	220874	102.83	PPB		94
41)	Isobutyl Alcohol	3.78	43	208799	2191.69	PPB		95
43)	Benzene	3.78	78	714040	105.34	PPB		97
44)	1,2-Dichloroethane	3.87	62	288349	104.20	PPB		98
45)	tert-Amyl Methyl Ether	3.84	55	147062	97.34	PPB	#	72
46)	Trichloroethene	4.29	95	179630	98.02	PPB		92
48)	1,2-Dichloropropane	4.53	63	192068	105.86	PPB		99
49)	Dibromomethane	4.62	93	121056	105.78	PPB		92
50)	Methyl methacrylate	4.60	69	141801	119.43	PPB		93
51)	1,4-Dioxane	4.62	88	49878	2126.07	PPB		99
52)	Bromodichloromethane	4.75	83	250846	115.24	PPB		97
53)	2-Nitropropane	5.02	41	158017	352.10	PPB		97
54)	2-Chloroethyl Vinyl Ether	5.02	63	124403	117.38	PPB		93
55)	cis-1,3-Dichloropropene	5.13	75	290398	123.33	PPB		93
56)	4-Methyl-2-pentanone (MIBK	5.26	58	154002	229.61	PPB		97
58)	Toluene	5.36	92	434142	106.10	PPB		96
61)	trans-1,3-Dichloropropene	5.65	75	248497	114.60	PPB		88
62)	Ethyl methacrylate	5.67	69	235477	112.67	PPB		99
63)	1,1,2-Trichloroethane	5.81	83	140620	93.18	PPB		95
64)	Tetrachloroethene	5.82	164	131482	92.27	PPB		87
65)	2-Hexanone	6.02	57	50618	208.65	PPB	#	66
66)	1,3-Dichloropropane	5.97	76	292573	98.76	PPB		88
67)	Dibromochloromethane	6.13	129	177378	117.08	PPB		96
68)	1,2-Dibromoethane (EDB)	6.25	107	160961	100.76	PPB		99
69)	1-Chlorohexane	6.64	91	204314	102.49	PPB		89
70)	Chlorobenzene	6.68	112	494364	100.24	PPB		98
71)	Ethylbenzene	6.76	106	242049	103.38	PPB	#	78
72)	1,1,1,2-Tetrachloroethane	6.77	131	166758	108.95	PPB		97
73)	m,p-Xylenes	6.87	106	608283	225.02	PPB		99
74)	o-Xylene	7.27	106	303878	112.41	PPB		94
75)	Styrene	7.30	103	240994m	116.35	PPB		
76)	Bromoform	7.51	173	108994	120.69	PPB		90

(#) = qualifier out of range (m) = manual integration
0421F011.D 042110MS24SOIL.M Wed Apr 21 14:53:46 2010 Page 2

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F011.D Vial: 13
Acq On : 21 Apr 2010 2:36 pm Operator: KR
Sample : SOIL ICAL 100 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:52:53 2010 Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:32:32 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
77)	Isopropylbenzene	7.62	105	710509	107.24	PPB		96
78)	cis-1,4-Dichloro-2-butene	7.80	89	42736	256.34	PPB	#	79
81)	1,1,2,2-Tetrachloroethane	8.03	83	182259	93.79	PPB		89
82)	trans-1,4-Dichloro-2-buten	8.10	53	46380	109.99	PPB		78
83)	Bromobenzene	7.96	156	205283	100.20	PPB		97
84)	n-Propylbenzene	8.04	91	872184	109.26	PPB		97
85)	1,2,3-Trichloropropane	8.07	110	65952	87.43	PPB	#	56
86)	2-Chlorotoluene	8.15	91	528945	101.79	PPB		98
87)	1,3,5-Trimethylbenzene	8.23	105	614976	109.53	PPB		98
88)	4-Chlorotoluene	8.27	91	547060	104.43	PPB		96
89)	tert-Butylbenzene	8.54	119	541258	109.80	PPB		98
90)	1,2,4-Trimethylbenzene	8.60	105	646423	114.37	PPB		90
91)	sec-Butylbenzene	8.76	105	811949	110.53	PPB		99
92)	p-Isopropyltoluene	8.91	119	663557	108.13	PPB		97
93)	1,3-Dichlorobenzene	8.90	146	376672	99.57	PPB		100
94)	1,4-Dichlorobenzene	9.00	146	387228	97.53	PPB		94
95)	n-Butylbenzene	9.31	91	597579	111.19	PPB		95
96)	1,2-Dichlorobenzene	9.36	146	378148	102.05	PPB		98
97)	1,2-Dibromo-3-chloropropan	10.14	155	25853	98.00	PPB		79
98)	1,3,5-Trichlorobenzene	10.26	180	271091	103.41	PPB		91
99)	1,2,4-Trichlorobenzene	10.86	180	232818	106.07	PPB		94
100)	Hexachlorobutadiene	10.96	225	100057	82.43	PPB		97
101)	Naphthalene	11.09	128	568190	109.04	PPB		97
102)	1,2,3-Trichlorobenzene	11.32	180	169842	80.01	PPB		92

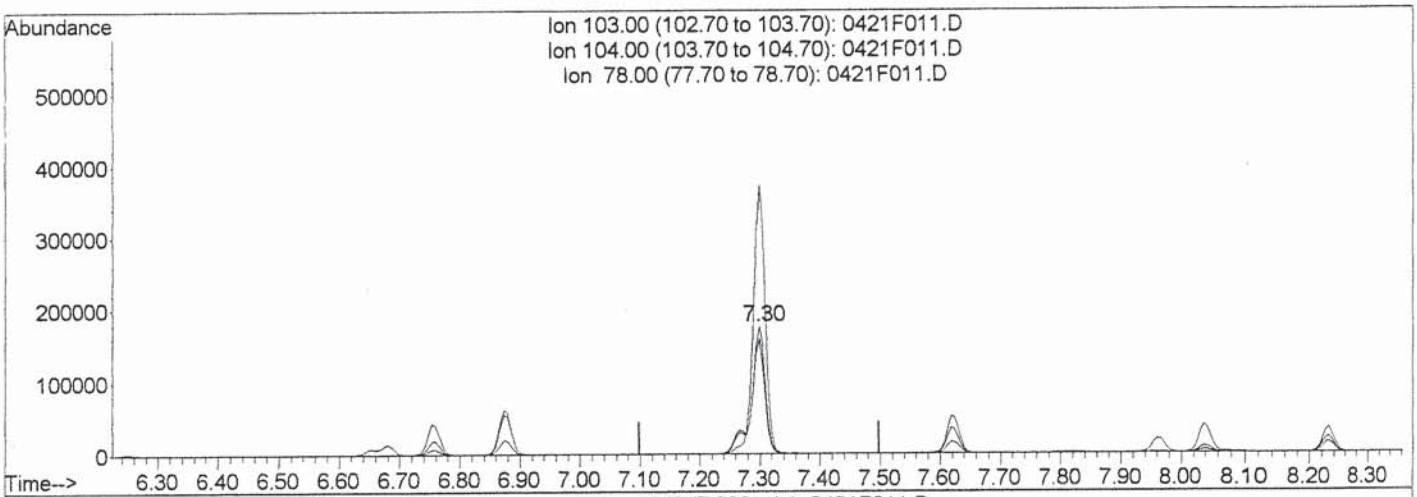
(#) = qualifier out of range (m) = manual integration

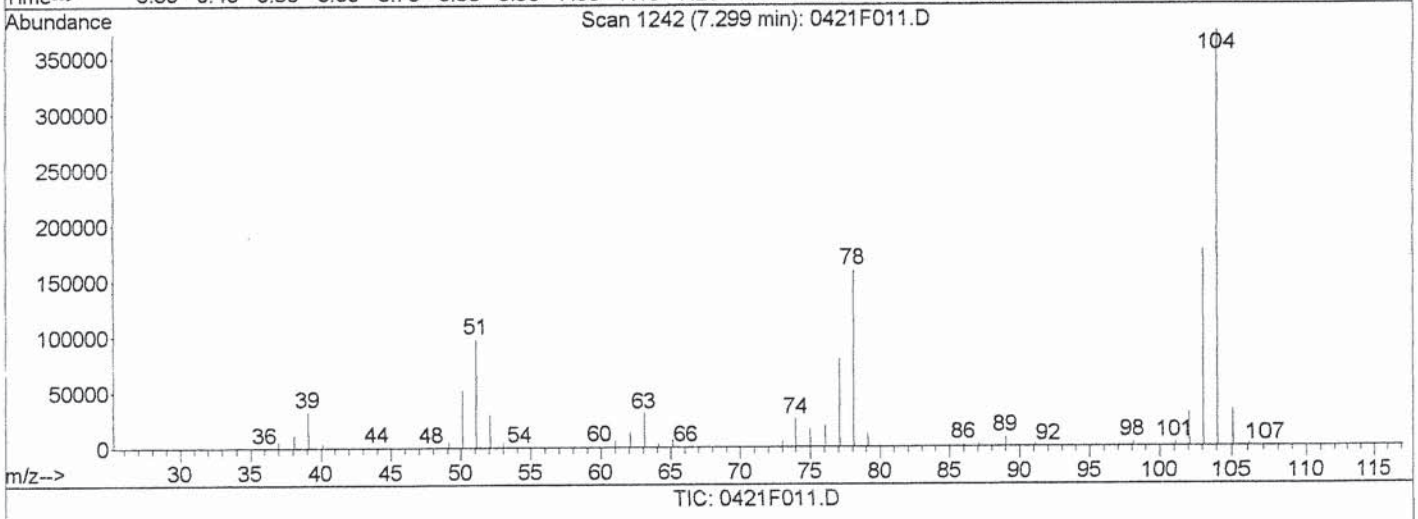
Quantitation Report (Qedit)

Data File : J:\MS24\DATA\042110\0421F011.D
Acq On : 21 Apr 2010 2:36 pm
Sample : SOIL ICAL 100
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:53 2010

Vial: 13
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:32:32 2010
Response via : Single Level Calibration





TIC: 0421F011.D

(75) Styrene (T)

7.30min 136.57PPB

response 282860

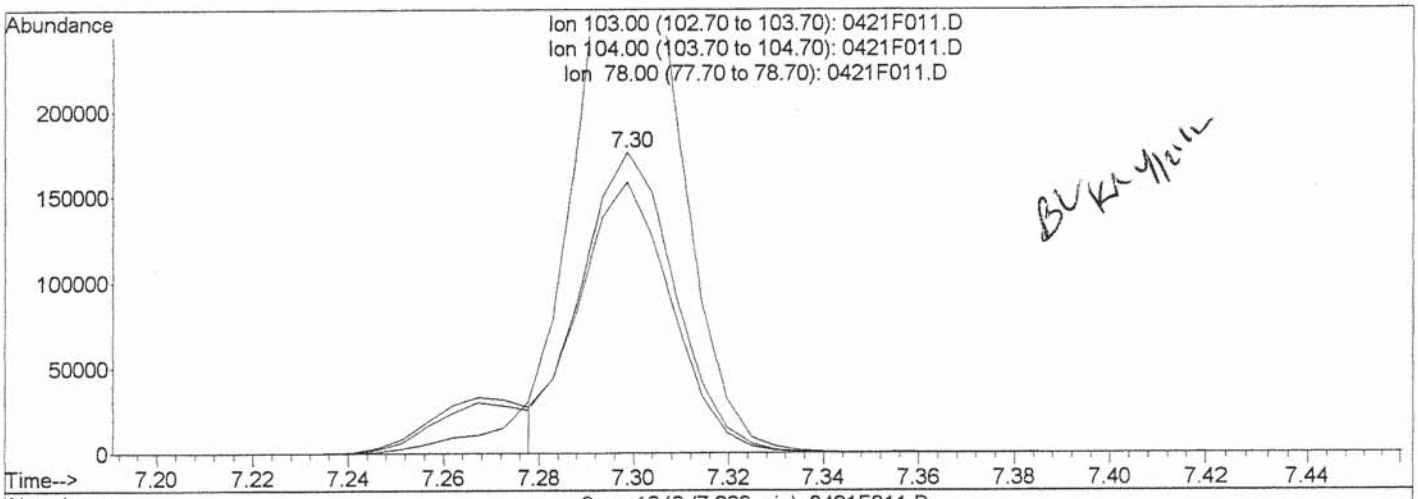
Ion	Exp%	Act%
103.00	100	100
104.00	200.50	212.00
78.00	90.20	90.08
0.00	0.00	0.00

Quantitation Report (Qedit)

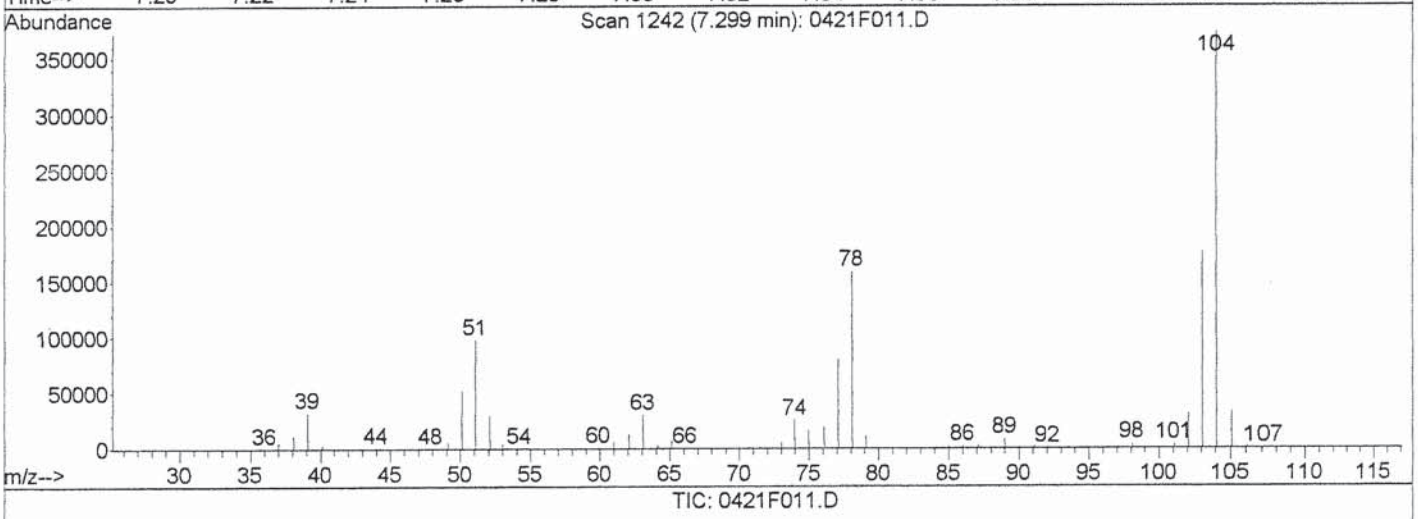
Data File : J:\MS24\DATA\042110\0421F011.D
Acq On : 21 Apr 2010 2:36 pm
Sample : SOIL ICAL 100
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 14:53 2010

Vial: 13
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: temp.res

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:32:32 2010
Response via : Single Level Calibration



BU KR 4/21/10



TIC: 0421F011.D

(75) Styrene (T)

7.30min 116.35PPB m

response 240994

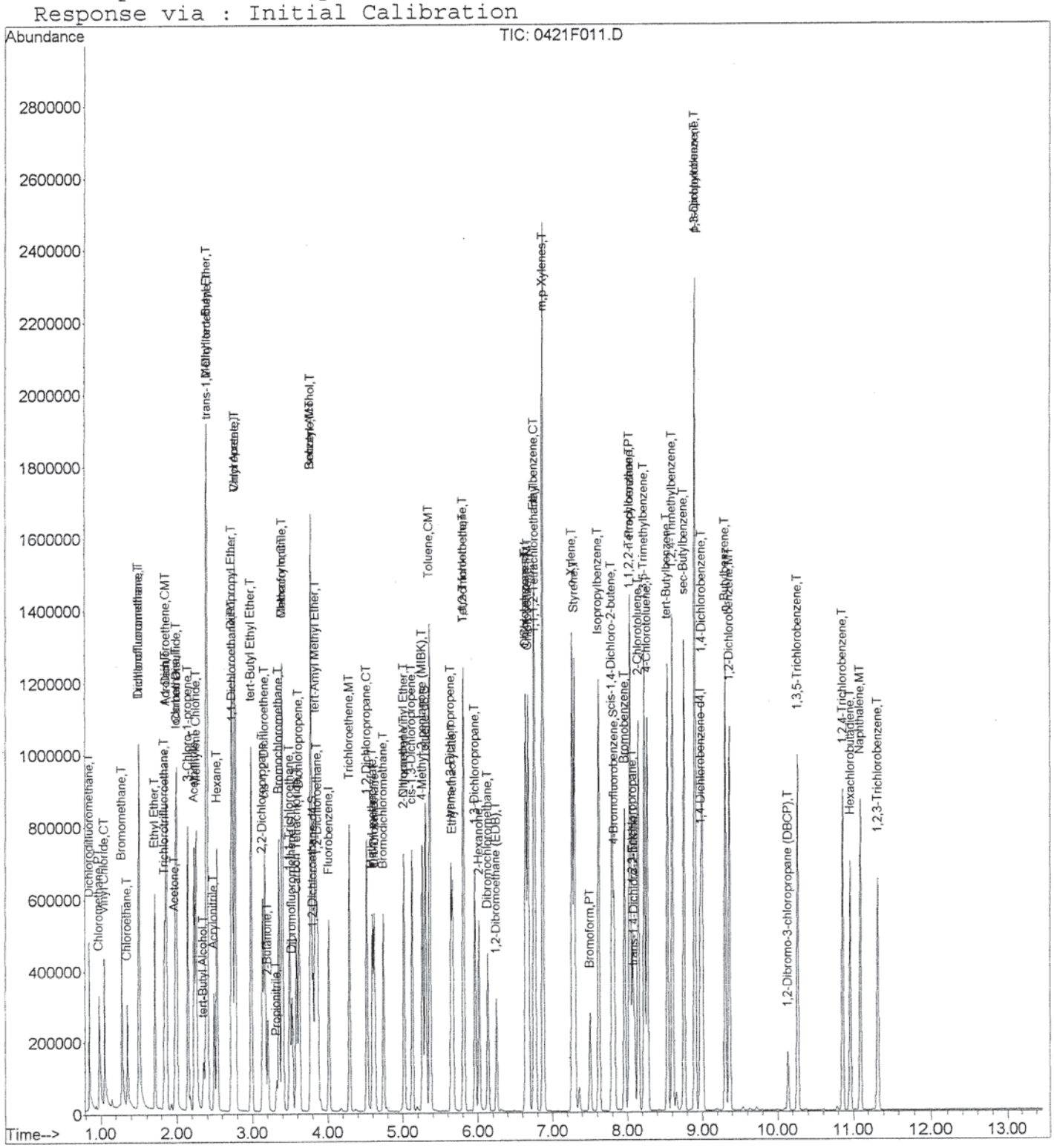
Ion	Exp%	Act%
103.00	100	100
104.00	200.50	212.00
78.00	90.20	90.08
0.00	0.00	0.00

0421F011.D 042110MS24SOIL.M Wed Apr 21 14:53:37 2010

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F011.D Vial: 13
Acq On : 21 Apr 2010 2:36 pm Operator: KR
Sample : SOIL ICAL 100 Inst : MS24
Misc : Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: Apr 21 14:53 2010 Quant Results File: 042110MS24SO

Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:32:32 2010
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F012.D
Acq On : 21 Apr 2010 2:57 pm
Sample : SOIL ICAL 200
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 15:32:48 2010

Vial: 14
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:54:43 2010
Response via : Initial Calibration
DataAcq Meth : 8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	4.01	96	330206	50.00	PPB	-0.01
59) Chlorobenzene-d5	6.65	82	146708	50.00	PPB	0.00
80) 1,4-Dichlorobenzene-d4	8.98	152	130673	50.00	PPB	0.00
System Monitoring Compounds						
38) Dibromofluoromethane	3.53	113	139985	85.56	PPB	0.00
Spiked Amount 50.000			Recovery =	171.12%		
42) 1,2-Dichloroethane-d4	3.81	65	202165	85.52	PPB	0.00
Spiked Amount 50.000			Recovery =	171.04%		
57) Toluene-d8	5.31	98	506060	88.98	PPB	0.00
Spiked Amount 50.000			Recovery =	177.96%		
79) 4-Bromofluorobenzene	7.83	95	190868	85.89	PPB	0.00
Spiked Amount 50.000			Recovery =	171.78%		
Target Compounds						Qvalue
2) Dichlorodifluoromethane	0.84	85	436976	214.33	PPB	97
3) Chloromethane	0.97	50	533074	204.38	PPB	95
4) Vinyl Chloride	1.04	62	483865	210.05	PPB	98
5) Bromomethane	1.27	96	401111	212.51	PPB	99
6) Chloroethane	1.34	64	319333	214.25	PPB	97
7) Dichlorofluoromethane	1.50	67	798428	213.39	PPB	95
8) Trichlorofluoromethane	1.49	101	684572	217.99	PPB	96
9) Ethyl Ether	1.70	59	397518	210.91	PPB	96
10) Acrolein	1.85	56	389260	1063.04	PPB	91
11) Trichlorotrifluoroethane	1.84	151	262892	218.94	PPB	94
12) 1,1-Dichloroethene	1.86	96	297437	214.20	PPB	95
13) Acetone	1.97	43	302643	379.45	PPB	98
14) Iodomethane	1.98	142	966238	557.62	PPB	99
15) Carbon Disulfide	1.99	76	1219159	227.65	PPB	97
16) 3-Chloro-1-propene	2.14	76	218476	224.15	PPB	97
17) Acetonitrile	2.22	40	648542	4308.09	PPB	97
18) Methylene Chloride	2.25	84	409896	204.42	PPB	96
19) tert-Butyl Alcohol	2.34	59	268026	944.90	PPB	96
20) Acrylonitrile	2.49	53	328562	429.52	PPB	95
21) Methyl tert-Butyl Ether	2.39	73	2618970	461.83	PPB	99
22) trans-1,2-Dichloroethene	2.40	96	373116	216.46	PPB	86
23) Hexane	2.53	57	497927	208.58	PPB	96
24) Diisopropyl Ether	2.72	45	1458779	226.54	PPB	96
25) 1,1-Dichloroethane	2.74	63	747986	220.24	PPB	96
26) Vinyl Acetate	2.77	86	70994	227.43	PPB #	66
27) Chloroprene	2.77	53	1087766	479.57	PPB	97
28) tert-Butyl Ethyl Ether	2.98	59	1327752	227.58	PPB	95

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : J:\MS24\DATA\042110\0421F012.D
Acq On : 21 Apr 2010 2:57 pm
Sample : SOIL ICAL 200
Misc :
MS Integration Params: rteint.p
Quant Time: Apr 21 15:32:48 2010

Vial: 14
Operator: KR
Inst : MS24
Multiplr: 1.00
Quant Results File: 042110MS24SOIL.

Quant Method : J:\MS24\METHODS\042110MS24SOIL.M (RTE Integrator)
Title : VOA MS24 EPA Method 8260B
Last Update : Wed Apr 21 14:54:43 2010
Response via : Initial Calibration
DataAcq Meth : 8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	2,2-Dichloropropane	3.14	77	533126	232.68	PPB		98
30)	cis-1,2-Dichloroethene	3.17	96	420285	219.87	PPB		96
31)	2-Butanone	3.20	72	104661	420.03	PPB	#	84
32)	Propionitrile	3.31	54	125804	403.67	PPB		92
33)	Methacrylonitrile	3.39	67	372164	469.92	PPB		89
34)	Bromochloromethane	3.35	128	208721	225.04	PPB		91
36)	Chloroform	3.40	83	726206	230.06	PPB		98
37)	1,1,1-Trichloroethane	3.50	97	546937	225.94	PPB		95
39)	Carbon Tetrachloride	3.59	117	478266	231.98	PPB		95
40)	1,1-Dichloropropene	3.62	75	489781	224.39	PPB		95
41)	Isobutyl Alcohol	3.77	43	465528	4752.07	PPB		95
43)	Benzene	3.78	78	1569828	227.09	PPB		98
44)	1,2-Dichloroethane	3.86	62	623023	221.12	PPB		98
45)	tert-Amyl Methyl Ether	3.84	55	321088	210.79	PPB	#	72
46)	Trichloroethene	4.28	95	392738	212.34	PPB		90
48)	1,2-Dichloropropane	4.52	63	401287	216.71	PPB		99
49)	Dibromomethane	4.62	93	249595	213.03	PPB		94
50)	Methyl methacrylate	4.60	69	298658	239.23	PPB		91
51)	1,4-Dioxane	4.61	88	102318	4255.47	PPB		97
52)	Bromodichloromethane	4.75	83	528554	232.82	PPB		96
53)	2-Nitropropane	5.02	41	363289	780.45	PPB		92
54)	2-Chloroethyl Vinyl Ether	5.02	63	269945	243.20	PPB		92
55)	cis-1,3-Dichloropropene	5.13	75	627929	251.73	PPB		94
56)	4-Methyl-2-pentanone (MIBK	5.26	58	325101	465.13	PPB		94
58)	Toluene	5.36	92	916260	219.01	PPB		95
61)	trans-1,3-Dichloropropene	5.65	75	565538	247.27	PPB		90
62)	Ethyl methacrylate	5.67	69	528893	240.83	PPB		99
63)	1,1,2-Trichloroethane	5.81	83	306558	200.16	PPB		93
64)	Tetrachloroethene	5.81	164	296977	205.63	PPB		90
65)	2-Hexanone	6.02	57	114173	455.94	PPB	#	69
66)	1,3-Dichloropropane	5.97	76	622357	205.36	PPB		88
67)	Dibromochloromethane	6.13	129	395747	246.47	PPB		97
68)	1,2-Dibromoethane (EDB)	6.25	107	359567	219.36	PPB		99
69)	1-Chlorohexane	6.64	91	467237	227.89	PPB		88
70)	Chlorobenzene	6.68	112	1067934	211.21	PPB		99
71)	Ethylbenzene	6.76	106	549447	227.70	PPB	#	77
72)	1,1,1,2-Tetrachloroethane	6.77	131	389221	243.78	PPB		96
73)	m,p-Xylenes	6.87	106	1351009	479.10	PPB		94
74)	o-Xylene	7.27	106	645306	227.29	PPB		92
75)	Styrene	7.30	103	520670m	237.52	PPB		
76)	Bromoform	7.51	173	244615	253.80	PPB		90

(#) = qualifier out of range (m) = manual integration