

Table 7  
Groundwater Analytical Results  
September 2021  
REXAM Beverage Can Company/Ball Corporation  
Kent, WA

Sample ID	Model Toxics Control Act - Groundwater  Method C Cleanup Levels	ETMW-2	ETMW-3	ETMW-7	MW-101	MW-102	MW-103	MW-104	MW-105R	Dup2	MW-106
Laboratory ID No.		109299-12	109299-02	109299-10	109299-15	109299-11	109299-03	109299-04	109299-13	109299-14	109299-16
Date Sampled		9/16/2021	9/15/2021	9/16/2021	9/16/2021	9/16/2021	9/15/2021	9/15/2021	9/16/2021	9/16/2021	9/16/2021
Well Location		Site	Site	Site	Site	Site	Site	Site	Site	Site	Site
<b>ANALYTICAL PARAMETERS</b>											
<b>Volatile Organic Compounds (µg/l)</b>											
Dichlorodifluoromethane	3,500	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Chloromethane	NS	<10	<10	<10	<10	<10	<10	<10	<10	<10	<10
Vinyl chloride	0.29	<0.02	<0.02	<b>2.7</b>	<b>1.3</b>	<b>0.081</b>	<b>0.26</b>	<0.02	<b>4.3</b>	<b>4.1</b>	<b>0.20</b>
Bromomethane	25	<5	<5	<5	<5	<5	<5	<5	<5	<5	<5
Chloroethane	830	<1	<1	<b>2,000 ve/2,200</b>	<b>6.9</b>	<b>94</b>	<1	<1	<b>320 ve/ 310</b>	<b>330 ve/ 300</b>	<b>85</b>
Trichlorofluoromethane	5,300	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Acetone	16,000	<50	<50	<50	<50	<50	<50	<50	<50	<50	<50
1,1-Dichloroethene	7.0	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Hexane	1,100	<5	<5	<5	<5	<5	<5	<5	<5	<5	<5
Methylene Chloride	5.0	<5	<5	<5	<5	<5	<5	<5	<5	<5	<5
Methyl tert-butyl ether (MTBE)	240	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl
trans-1,2-Dichloroethene	100	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,1-Dichloroethane	77	<1	<1	<b>2.1</b>	<1	<1	<1	<1	<b>100</b>	<b>100</b>	<1
2,2-Dichloropropane	NS	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl
cis-1,2-Dichloroethene	35	<1	<1	<1	<1	<1	<b>1.3</b>	<1	<1	<1	<1
Chloroform	14	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
2-Butanone (MEK)	11,000	<20	<20	<20	<20	<20	<20	<20	<20	<20	<20
1,2-Dichloroethane (EDC)	4.8	<0.2	<0.2	<b>6.9</b>	<0.2	<b>1.6</b>	<0.2	<0.2	<b>1.9</b>	<b>1.8</b>	<b>0.55</b>
1,1,1-Trichloroethane	200	<1	<1	<1	<1	<1	<1	<1	<b>1.6</b>	<b>1.6</b>	<1
1,1-Dichloropropene	NS	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Carbon tetrachloride	5.0	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
Benzene	5.0	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35
Trichloroethene	5.0	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
1,2-Dichloropropane	5.0	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Bromodichloromethane	7.1	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
Dibromomethane	180	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
4-Methyl-2-pentanone (MIBK)	1,400	<10	<10	<10	<10	<10	<10	<10	<10	<10	<10
cis-1,3-Dichloropropene	4.4 <sup>TOT</sup>	<0.4	<0.4	<0.4	<0.4	<0.4	<0.4	<0.4	<0.4	<0.4	<0.4
Toluene	1,000	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
trans-1,3-Dichloropropene	4.4 <sup>TOT</sup>	<0.4	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca
1,1,2-Trichloroethane	3.0	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
2-Hexanone (MBK)	88	<10	<10	<10	<10	<10	<10	<10	<10	<10	<10
1,3-Dichloropropane	350	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Tetrachloroethene	5.0	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Dibromochloromethane	5.2	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
1,2-Dibromoethane (EDB)	0.05	<0.01 j	<0.01 j	<0.01 j	<0.01 j	<0.01 j	<0.01 j	<0.01 j	<0.01 j	<0.01 j	<0.01 j
Chlorobenzene	100	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Ethylbenzene	700	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,1,1,2-Tetrachloroethane	17	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
m-Xylene & p-Xylene	3,500*	<2	<2	<2	<2	<2	<2	<2	<2	<2	<2
o-Xylene	3,500*	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Styrene	100	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1

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Sample ID	Model Toxics Control Act - Groundwater  Method C Cleanup Levels	ETMW-2	ETMW-3	ETMW-7	MW-101	MW-102	MW-103	MW-104	MW-105R	Dup2	MW-106
Laboratory ID No.		109299-12	109299-02	109299-10	109299-15	109299-11	109299-03	109299-04	109299-13	109299-14	109299-16
Date Sampled		9/16/2021	9/15/2021	9/16/2021	9/16/2021	9/16/2021	9/15/2021	9/15/2021	9/16/2021	9/16/2021	9/16/2021
Well Location		Site	Site	Site	Site	Site	Site	Site	Site	Site	Site
<b>ANALYTICAL PARAMETERS</b>											
Isopropylbenzene (Cumene)	1,800	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Bromoform	55	<5	<5	<5	<5	<5	<5	<5	<5	<5	<5
n-Propylbenzene	1,800	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Bromobenzene	140	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,3,5-Trimethylbenzene	180	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,1,2,2-Tetrachloroethane	2.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2
1,2,3-Trichloropropane	0.015	<0.072 j	<0.072 j	<0.072 j	<0.072 j	<0.072 j	<0.072 j	<0.072 j	<0.072 j	<0.072 j	<0.072 j
2-Chlorotoluene	350	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
4-Chlorotoluene	NS	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
tert-Butylbenzene	1,800	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,2,4-Trimethylbenzene	180	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
sec-Butylbenzene	1,800	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
p-Isopropyltoluene	NS	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,3-Dichlorobenzene	NS	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,4-Dichlorobenzene	75	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,2-Dichlorobenzene	600	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,2-Dibromo-3-Chloropropane	0.2	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j
1,2,4-Trichlorobenzene	15	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
Hexachlorobutadiene	5.6	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
Naphthalene	350	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,2,3-Trichlorobenzene	14	<1	<1	<1	<1	<1	<1	<1	<1	<1	<1
1,4-Dioxane	4.4	<b>0.72</b>	<0.4	<b>310</b>	<b>21</b>	<b>91</b>	<b>10</b>	<0.4	<b>60</b>	<b>51</b>	<b>150</b>
<b>Tentatively Identified Compounds (µg/l)</b>											
Benzene, 1-ethyl-2-methyl	NS										
Benzene, 1-ethyl-3-methyl	NS										<b>1.1</b>
Ethane, 1,2-dichloro-1,1,2-trifluoro	NS			<b>10</b>					<b>2.7</b>	<b>2.6</b>	
Butane, 1,3-dichloro-	NS			<b>3.4</b>		<b>3.4</b>			<b>3.0</b>	<b>3.0</b>	
Ethane, bromo	NS			<b>2.8</b>							
Ethane, (methylthio)	NS			<b>2.7</b>							
Ethane, 1,1'-thiobis	NS			<b>2.7</b>		<b>7.6</b>			<b>4.3</b>	<b>4.4</b>	
1,3-Dioxolane	NS			<b>1.3</b>							
Benzene, 4-ethyl-1,2-dimethyl-	NS	<b>1.2</b>									
1-Propene, 1,1,3-trichloro-2-methyl	NS								<b>1.6</b>		
Cyclopropane, 1,1-dimethyl	NS										
Benzene, 1,2,3-trimethyl	180										
Total Tentatively Identified Compounds	NS	<b>1.2</b>	0.0	<b>22.9</b>	0.0	<b>11.0</b>	0.0	0.0	<b>11.6</b>	<b>10.0</b>	<b>1.1</b>

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Laboratory ID No.		109299-12	109299-02	109299-10	109299-15	109299-11	109299-03	109299-04	109299-13	109299-14	109299-16
Date Sampled	Method C Cleanup Levels	9/16/2021	9/15/2021	9/16/2021	9/16/2021	9/16/2021	9/15/2021	9/15/2021	9/16/2021	9/16/2021	9/16/2021
Well Location		Site	Site	Site	Site	Site	Site	Site	Site	Site	Site
<b>ANALYTICAL PARAMETERS</b>											

Legend:

µg/l = micrograms per liter.

<0.00 = Not detected above Method Detection Limit.

NA = Not Analyzed

NS = No standard

Bolded values represent detections

Shaded cell denotes exceedance of Method C Cleanup Level

\* = Total xylenes

ve = The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

ca = The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

j = The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

jl = The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc = The presence of the analyte is likely due to laboratory contamination.

##<sup>TOT</sup> = Standard is for mixtures of cis- and trans- isomers.

NS = No default Method C established. Not a Site contaminant of concern.

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Kent, WA

Sample ID	Model Toxics Control Act - Groundwater  Method C Cleanup Levels	MW-107	MW-108	MW-113	Dup1	MW-201
Laboratory ID No.		109299-05	109299-18	109299-08	109299-06	109299-09
Date Sampled		9/15/2021	9/16/2021	9/15/2021	9/15/2021	9/16/2021
Well Location		Site	Site	Site	Site	Site
<b>ANALYTICAL PARAMETERS</b>						
<b>Volatile Organic Compounds (µg/l)</b>						
Dichlorodifluoromethane	3,500	<1	<1	<1	<1	<1
Chloromethane	NS	<10	<10	<10	<10	<10
Vinyl chloride	0.29	<0.02	<b>160 ve / 100</b>	<b>0.037</b>	<b>0.041</b>	<0.02
Bromomethane	25	<5	<5	<5	<5	<5
Chloroethane	830	<1	<b>2,400 ve / 2,300</b>	<1	<1	<1
Trichlorofluoromethane	5,300	<1	<1	<1	<1	<1
Acetone	16,000	<50	<50	<50	<50	<50
1,1-Dichloroethene	7.0	<1	<b>370 ve / 120</b>	<1	<1	<1
Hexane	1,100	<5	<5	<5	<5	<5
Methylene Chloride	5.0	<5	<5	<5	<5	<5
Methyl tert-butyl ether (MTBE)	240	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl
trans-1,2-Dichloroethene	100	<1	<b>1.3</b>	<1	<1	<1
1,1-Dichloroethane	77	<1	<b>1,200 ve / 1,200</b>	<1	<1	<1
2,2-Dichloropropane	NS	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl	<1 ca jl
cis-1,2-Dichloroethene	35	<1	<b>3.2</b>	<1	<1	<1
Chloroform	14	<1	<1	<1	<1	<1
2-Butanone (MEK)	11,000	<20	<20	<20	<20	<20
1,2-Dichloroethane (EDC)	4.8	<0.2	<b>5.8</b>	<0.2	<0.2	<0.2
1,1,1-Trichloroethane	200	<1	<b>1,500 ve / 1,700</b>	<1	<b>1.1</b>	<1
1,1-Dichloropropene	NS	<1	<1	<1	<1	<1
Carbon tetrachloride	5.0	<0.5	<0.5	<0.5	<0.5	<0.5
Benzene	5.0	<0.35	<b>0.46</b>	<0.35	<0.35	<0.35
Trichloroethene	5.0	<0.5	<b>1.1</b>	<0.5	<0.5	<0.5
1,2-Dichloropropane	5.0	<1	<1	<1	<1	<1
Bromodichloromethane	7.1	<0.5	<0.5	<0.5	<0.5	<0.5
Dibromomethane	180	<1	<1	<1	<1	<1
4-Methyl-2-pentanone (MIBK)	1,400	<10	<10	<10	<10	<10
cis-1,3-Dichloropropene	4.4 <sup>TOT</sup>	<0.4	<0.4	<0.4	<0.4	<0.4
Toluene	1,000	<1	<b>10</b>	<1	<1	<1
trans-1,3-Dichloropropene	4.4 <sup>TOT</sup>	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca	<0.4 ca
1,1,2-Trichloroethane	3.0	<0.5	<0.5	<0.5	<0.5	<0.5
2-Hexanone (MBK)	88	<10	<10	<10	<10	<10
1,3-Dichloropropane	350	<1	<1	<1	<1	<1
Tetrachloroethene	5.0	<1	<b>3.5</b>	<1	<1	<1
Dibromochloromethane	5.2	<0.5	<0.5	<0.5	<0.5	<0.5
1,2-Dibromoethane (EDB)	0.05	<0.01	<0.01 j	<0.01 j	<0.01 j	<0.01 j
Chlorobenzene	100	<1	<1	<1	<1	<1
Ethylbenzene	700	<1	<1	<1	<1	<1
1,1,1,2-Tetrachloroethane	17	<1	<1	<1	<1	<1
m-Xylene & p-Xylene	3,500*	<2	<2	<2	<2	<2
o-Xylene	3,500*	<1	<b>3.9</b>	<1	<1	<1
Styrene	100	<1	<1	<1	<1	<1

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Sample ID	Model Toxics Control Act - Groundwater	MW-107	MW-108	MW-113	Dup1	MW-201
Laboratory ID No.		109299-05	109299-18	109299-08	109299-06	109299-09
Date Sampled	Method C Cleanup Levels	9/15/2021	9/16/2021	9/15/2021	9/15/2021	9/16/2021
Well Location		Site	Site	Site	Site	Site
<b>ANALYTICAL PARAMETERS</b>						
Isopropylbenzene (Cumene)	1,800	<1	<1	<1	<1	<1
Bromoform	55	<5	<5	<5	<5	<5
n-Propylbenzene	1,800	<1	<1	<1	<1	<1
Bromobenzene	140	<1	<1	<1	<1	<1
1,3,5-Trimethylbenzene	180	<1	<b>1.6</b>	<1	<b>1.4</b>	<1
1,1,1,2-Tetrachloroethane	2.2	<0.2	<0.2	<0.2	<0.2	<0.2
1,2,3-Trichloropropane	0.015	<0.072	<0.072 j	<0.072 j	<0.072 j	<0.072 j
2-Chlorotoluene	350	<1	<1	<1	<1	<1
4-Chlorotoluene	NS	<1	<1	<1	<1	<1
tert-Butylbenzene	1,800	<1	<1	<1	<1	<1
1,2,4-Trimethylbenzene	180	<1	<b>4.7</b>	<b>3.0</b>	<b>4.8</b>	<1
sec-Butylbenzene	1,800	<1	<1	<1	<1	<1
p-Isopropyltoluene	NS	<1	<1	<1	<1	<1
1,3-Dichlorobenzene	NS	<1	<1	<1	<1	<1
1,4-Dichlorobenzene	75	<1	<1	<1	<1	<1
1,2-Dichlorobenzene	600	<1	<1	<1	<1	<1
1,2-Dibromo-3-Chloropropane	0.2	<0.14 j	<0.14 j	<0.14 j	<0.14 j	<0.14 j
1,2,4-Trichlorobenzene	15	<1	<1	<1	<1	<1
Hexachlorobutadiene	5.6	<0.5	<0.5	<0.5	<0.5	<0.5
Naphthalene	350	<1	<1	<1	<1	<1
1,2,3-Trichlorobenzene	14	<1	<1	<1	<1	<1
1,4-Dioxane	4.4	<0.4	<b>3,200</b>	<0.4	<0.4	<0.4
<b>Tentatively Identified Compounds (µg/l)</b>						
Benzene, 1-ethyl-2-methyl	NS		<b>4.5</b>		<b>3.4</b>	
Benzene, 1-ethyl-3-methyl	NS		<b>3.1</b>	<b>2.2</b>	<b>1.4</b>	
Ethane, 1,2-dichloro-1,1,2-trifluoro	NS		<b>6.9</b>			
Butane, 1,3-dichloro-	NS		<b>94</b>			
Ethane, bromo	NS					
Ethane, (methylthio)	NS		<b>10</b>			
Ethane, 1,1'-thiobis	NS					
1,3-Dioxolane	NS					
Benzene, 4-ethyl-1,2-dimethyl-	NS					
1-Propene, 1,1,3-trichloro-2-methyl	NS					
Cyclopropane, 1,1-dimethyl	NS		<b>42</b>			
Benzene, 1,2,3-trimethyl	180		<b>2.1</b>			
Total Tentatively Identified Compounds	NS	0.0	<b>162.6</b>	<b>2.2</b>	<b>4.8</b>	0.0

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Date Sampled	Method C Cleanup Levels	9/15/2021	9/16/2021	9/15/2021	9/15/2021	9/16/2021
Well Location		Site	Site	Site	Site	Site
<b>ANALYTICAL PARAMETERS</b>						

**Legend:**

µg/l = micrograms per liter.

<0.00 = Not detected above Method Detection Limit.

NA = Not Analyzed

NS = No standard

Bolded values represent detections

Shaded cell denotes exceedance of Method C Cleanup Level

\* = Total xylenes

ve = The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.

ca = The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.

j = The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.

jl = The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.

js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.

lc = The presence of the analyte is likely due to laboratory contamination.

##<sup>TOT</sup> = Standard is for mixtures of cis- and trans- isomers.

NS = No default Method C established. Not a Site contaminant of concern.

Table 7  
Groundwater Analytical Results  
September 2021  
REXAM Beverage Can Company/Ball Corporation  
Kent, WA

Sample ID	Model Toxics Control	MW4	MW5	Dup	FB	Trip Blank	FB-1
Laboratory ID No.	Act - Groundwater	109300-03	109300-01	109300-02	109299-07	109299-01	109299-17
Date Sampled	Method C Cleanup Levels	9/15/2021	9/15/2021	9/15/2021	9/15/2021	9/15/2021	9/16/2021
Well Location		Protective Coatings, Inc., 1215 2nd Ave. N.				Quality Control Samples	
<b>ANALYTICAL PARAMETERS</b>							
<b>Volatile Organic Compounds (µg/l)</b>							
Dichlorodifluoromethane	3,500	<1	<1	<1	<1	<1	<1
Chloromethane	NS	<10	<10	<10	<10	<10	<10
Vinyl chloride	0.29	<0.02	<b>0.51</b>	<b>0.38</b>	<0.02	<0.02	<0.02
Bromomethane	25	<5	<5	<5	<5	<5	<5
Chloroethane	830	<1	<1	<1	<1	<1	<1
Trichlorofluoromethane	5,300	<1	<1	<1	<1	<1	<1
Acetone	16,000	<50	<50	<50	<50	<50	<50
1,1-Dichloroethene	7.0	<1	<1	<1	<1	<1	<1
Hexane	1,100	<5	<5	<5	<5	<5	<5
Methylene Chloride	5.0	<5	<5	<5	<b>6.0 ca lc</b>	<5	<b>6.7 ca jl js lc</b>
Methyl tert-butyl ether (MTBE)	240	<1	<1	<1	<1 ca jl	<1 ca jl	<1 ca jl
trans-1,2-Dichloroethene	100	<1	<1	<1	<1	<1	<1
1,1-Dichloroethane	77	<1	<1	<1	<1	<1	<1
2,2-Dichloropropane	NS	<1	<1	<1	<1 ca jl	<1 ca jl	<1 ca jl
cis-1,2-Dichloroethene	35	<1	<1	<1	<1	<1	<1
Chloroform	14	<1	<1	<1	<1	<1	<1
2-Butanone (MEK)	11,000	<20	<20	<20	<20	<20	<20
1,2-Dichloroethane (EDC)	4.8	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2
1,1,1-Trichloroethane	200	<1	<1	<1	<1	<1	<1
1,1-Dichloropropene	NS	<1	<1	<1	<1	<1	<1
Carbon tetrachloride	5.0	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
Benzene	5.0	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35
Trichloroethene	5.0	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
1,2-Dichloropropane	5.0	<1	<1	<1	<1	<1	<1
Bromodichloromethane	7.1	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
Dibromomethane	180	<1	<1	<1	<1	<1	<1
4-Methyl-2-pentanone (MIBK)	1,400	<10	<10	<10	<10	<10	<10
cis-1,3-Dichloropropene	4.4 <sup>TOT</sup>	<0.4	<0.4	<0.4	<0.4	<0.4	<0.4
Toluene	1,000	<1	<1	<1	<1	<1	<1
trans-1,3-Dichloropropene	4.4 <sup>TOT</sup>	<0.4	<0.4	<0.4	<0.4 ca	<0.4 ca	<0.4 ca
1,1,2-Trichloroethane	3.0	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
2-Hexanone (MBK)	88	<10	<10	<10	<10	<10	<10
1,3-Dichloropropane	350	<1	<1	<1	<1	<1	<1
Tetrachloroethene	5.0	<1	<1	<1	<1	<1	<1
Dibromochloromethane	5.2	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
1,2-Dibromoethane (EDB)	0.05	<0.013 j	<0.013 j	<0.013 j	<0.01 j	<0.01 j	<0.01 j
Chlorobenzene	100	<1	<1	<1	<1	<1	<1
Ethylbenzene	700	<1	<1	<1	<1	<1	<1
1,1,1,2-Tetrachloroethane	17	<1	<1	<1	<1	<1	<1
m-Xylene & p-Xylene	3,500*	<2	<2	<2	<2	<2	<2
o-Xylene	3,500*	<1	<1	<1	<1	<1	<1
Styrene	100	<1	<1	<1	<1	<1	<1

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Sample ID	Model Toxics Control	MW4	MW5	Dup	FB	Trip Blank	FB-1
Laboratory ID No.	Act - Groundwater	109300-03	109300-01	109300-02	109299-07	109299-01	109299-17
Date Sampled	Method C Cleanup Levels	9/15/2021	9/15/2021	9/15/2021	9/15/2021	9/15/2021	9/16/2021
Well Location		Protective Coatings, Inc., 1215 2nd Ave. N.				Quality Control Samples	
<b>ANALYTICAL PARAMETERS</b>							
Isopropylbenzene (Cumene)	1,800	<1	<1	<1	<1	<1	<1
Bromoform	55	<5	<5	<5	<5	<5	<5
n-Propylbenzene	1,800	<1	<1	<1	<1	<1	<1
Bromobenzene	140	<1	<1	<1	<1	<1	<1
1,3,5-Trimethylbenzene	180	<1	<1	<1	<1	<1	<1
1,1,2,2-Tetrachloroethane	2.2	<0.2	<0.2	<0.2	<0.2	<0.2	<0.2
1,2,3-Trichloropropane	0.015	<0.065 j	<1	<0.065 j	<0.072 j	<0.072 j	<0.072 j
2-Chlorotoluene	350	<1	<1	<1	<1	<1	<1
4-Chlorotoluene	NS	<1	<1	<1	<1	<1	<1
tert-Butylbenzene	1,800	<1	<1	<1	<1	<1	<1
1,2,4-Trimethylbenzene	180	<1	<1	<1	<1	<1	<1
sec-Butylbenzene	1,800	<1	<1	<1	<1	<1	<1
p-Isopropyltoluene	NS	<1	<1	<1	<1	<1	<1
1,3-Dichlorobenzene	NS	<1	<1	<1	<1	<1	<1
1,4-Dichlorobenzene	75	<1	<1	<1	<1	<1	<1
1,2-Dichlorobenzene	600	<1	<1	<1	<1	<1	<1
1,2-Dibromo-3-Chloropropane	0.2	<1.3 j	<10	<1.3 j	<0.14 j	<0.14 j	<0.14 j
1,2,4-Trichlorobenzene	15	<1	<1	<1	<1	<1	<1
Hexachlorobutadiene	5.6	<0.5	<0.5	<0.5	<0.5	<0.5	<0.5
Naphthalene	350	<1	<1	<1	<1	<1	<1
1,2,3-Trichlorobenzene	14	<1	<1	<1	<1	<1	<1
1,4-Dioxane	4.4	<b>0.94</b>	<b>23</b>	<b>18</b>	<0.4	NA	<0.4
<b>Tentatively Identified Compounds (µg/l)</b>							
Benzene, 1-ethyl-2-methyl	NS						
Benzene, 1-ethyl-3-methyl	NS						
Ethane, 1,2-dichloro-1,1,2-trifluoro	NS						
Butane, 1,3-dichloro-	NS						
Ethane, bromo	NS						
Ethane, (methylthio)	NS						
Ethane, 1,1'-thiobis	NS						
1,3-Dioxolane	NS						
Benzene, 4-ethyl-1,2-dimethyl-	NS						
1-Propene, 1,1,3-trichloro-2-methyl	NS						
Cyclopropane, 1,1-dimethyl	NS						
Benzene, 1,2,3-trimethyl	180						
Total Tentatively Identified Compounds	NS	0.0	0.0	0.0	0.0	0.0	0.0

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Laboratory ID No.		109300-03	109300-01	109300-02	109299-07	109299-01	109299-17
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Well Location		Protective Coatings, Inc., 1215 2nd Ave. N.			Quality Control Samples		
<b>ANALYTICAL PARAMETERS</b>							
Legend:							

µg/l = micrograms per liter.  
 <0.00 = Not detected above Method Detection Limit.  
 NA = Not Analyzed  
 NS = No standard  
 Bolded values represent detections  
 Shaded cell denotes exceedance of Method C Cleanup Level  
 \* = Total xylenes  
 ve = The analyte response exceeded the valid instrument calibration range. The value reported is an estimate.  
 ca = The calibration results for the analyte were outside of acceptance criteria. The value reported is an estimate.  
 j = The analyte concentration is reported below the lowest calibration standard. The value reported is an estimate.  
 jl = The laboratory control sample(s) percent recovery and/or RPD were out of control limits. The reported concentration should be considered an estimate.  
  
 js - The surrogate associated with the analyte is out of control limits. The reported concentration should be considered an estimate.  
  
 lc = The presence of the analyte is likely due to laboratory contamination.  
 ##<sup>TOT</sup> = Standard is for mixtures of cis- and trans- isomers.  
 NS = No default Method C established. Not a Site contaminant of concern.