LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

Signal : TIC: BP010039.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.393	355	392	396	rBV3	53083863	439390784	100.00%	35.644%
2	6.869	633	643	649	rBV	520349	930191	0.21%	0.075%
3	7.134	677	688	703	rVB2	490938	1283151	0.29%	0.104%
4	7.293	703	715	727	rBV2	3491689	8742272	1.99%	0.709%
5	7.516	739	753	768	rVV2	790399	2751777	0.63%	0.223%
6	7.858	800	811	825	rBV	9321635	34236283	7.79%	2.777%
7	8.087	825	850	854	rVV4	48719354	297701678	67.75%	24.150%
8	8.434	882	909	912	rBV3	48325443	303683973	69.11%	24.635%
9	8.652	940	946	959	rVV	611342	1018517	0.23%	0.083%
10	9.110	1017	1024	1029	rBV	841696	1426623	0.32%	0.116%
11	9.775	1129	1137	1143	rBV	3385896	5926718	1.35%	0.481%
12	9.828	1143	1146	1157	rVB	717226	1146938	0.26%	0.093%
13	10.369	1231	1238	1244	rBV	987304	1753061	0.40%	0.142%
14	10.434	1244	1249	1257	rVB	715341	1236272	0.28%	0.100%
15	10.534	1257	1266	1273	rBV	1183491	2103989	0.48%	0.171%
16	10.634	1273	1283	1297	rVB	20179933	38883150	8.85%	3.154%
17	10.757	1297	1304	1313	rVB	899592	1450879	0.33%	0.118%
18	10.887	1313	1326	1331	rBV	16026016	37274834	8.48%	3.024%
19	11.146	1361	1370	1376	rVV	5190704	8716756	1.98%	0.707%
20	12.775	1640	1647	1657	rVB	739844	1208386	0.28%	0.098%
21	13.004	1680	1686	1693	rVV2	361407	639376	0.15%	0.052%
22	13.581	1776	1784	1791	rBV	966929	1504341	0.34%	0.122%
23	13.981	1845	1852	1866	rVB	2048622	2953680	0.67%	0.240%
24	14.263	1894	1900	1910	rVB	2234133	3229775	0.74%	0.262%
25	14.569	1946	1952	1963	rBV	1411111	2125878	0.48%	0.172%
26	14.757	1975	1984	1995	rBV	247467	449079	0.10%	0.036%
27	15.563	2115	2121	2132	rBV	2912597	4221263	0.96%	0.342%
28	15.675	2132	2140	2151	rBV	961822	1366510	0.31%	0.111%
29	17.322	2413	2420	2430	rBV	1862885	2696592	0.61%	0.219%
30	17.422	2430	2437	2446	rBV	3287667	4813879	1.10%	0.391%
31	18.663	2637	2648	2656	rBV	425929	585055	0.13%	0.047%
32	19.651	2809	2816	2828	rBV	4740269	6006885	1.37%	0.487%
33	21.110	3059	3064	3071	rBV	336381	477138	0.11%	0.039%
34	21.404	3108	3114	3131	rBV	2170165	3039609	0.69%	0.247%
35	23.698	3496	3504	3511	rBV2	2357499	4977585	1.13%	0.404%
36	23.857	3523	3531	3543	rVB2	1309940	2756097	0.63%	0.224%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

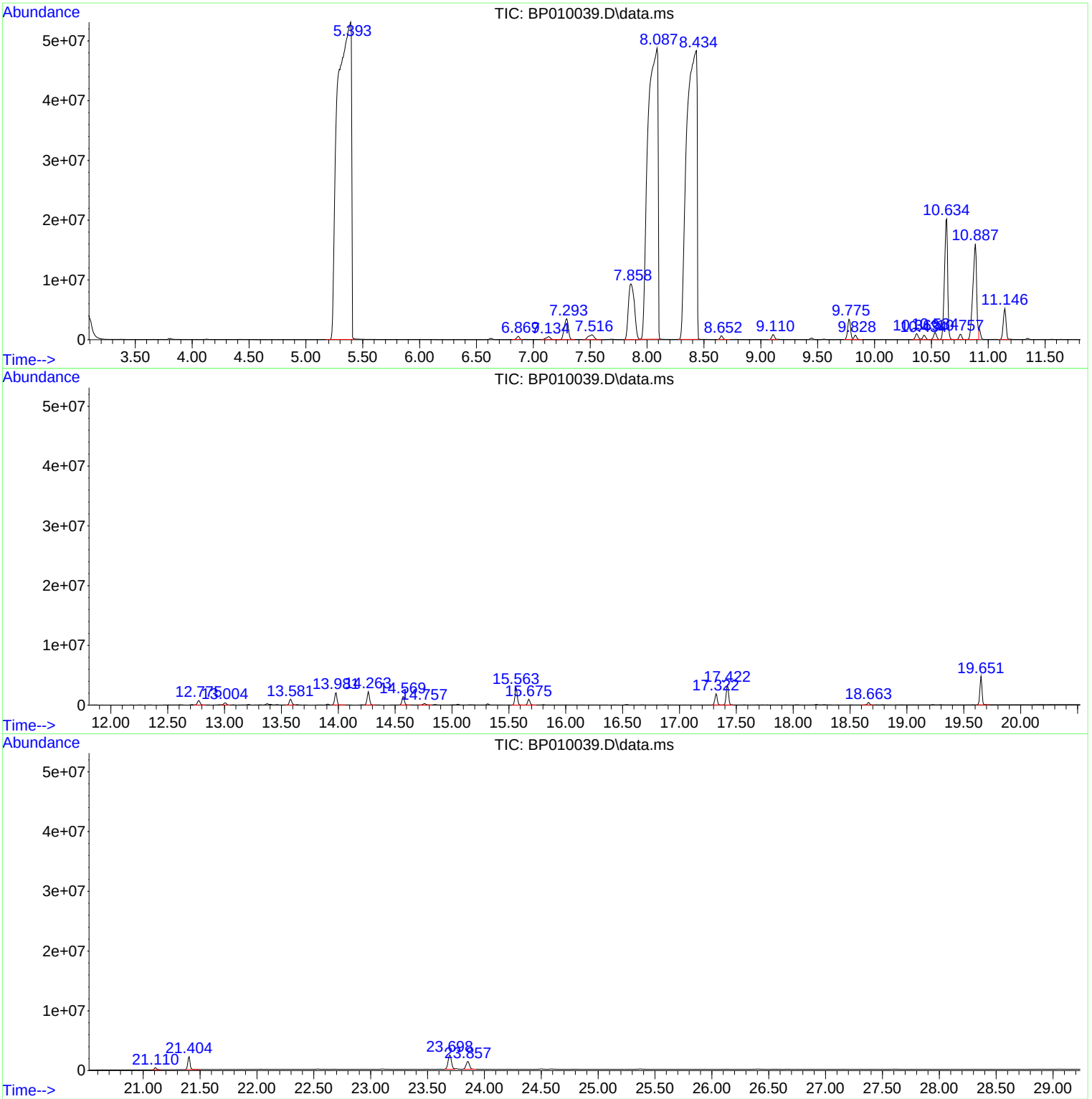
Sum of corrected areas: 1232708974

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
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Instrument :
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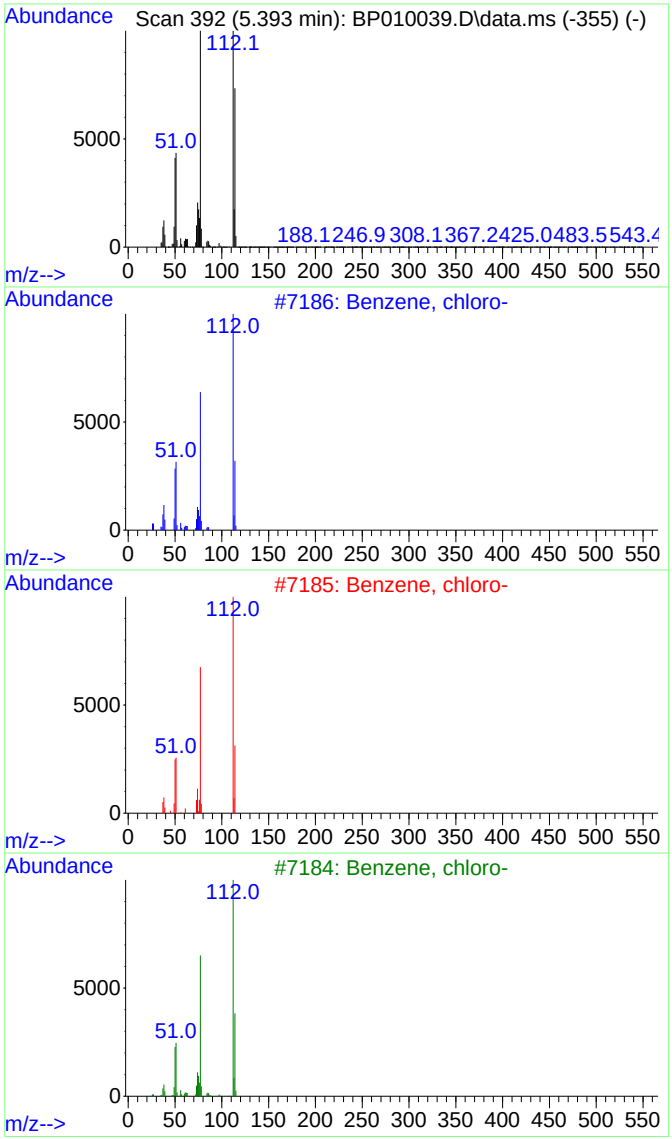
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

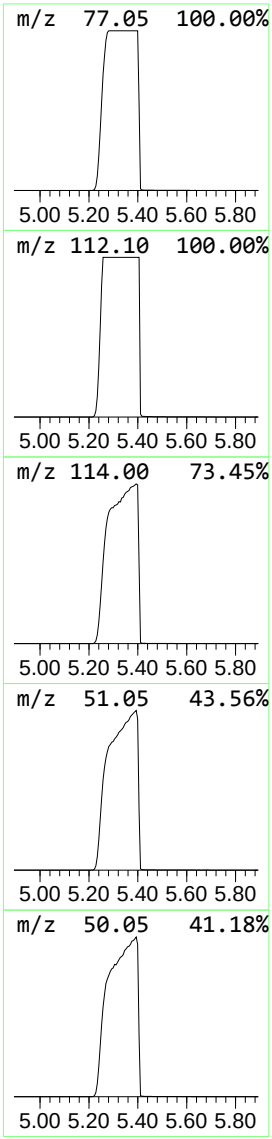
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.393	29.52 ng/ul	439391000	1,4-Dichlorobenzene-d4	7.999

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	72
5	Benzene, chloro-	112	C6H5Cl	000108-90-7	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

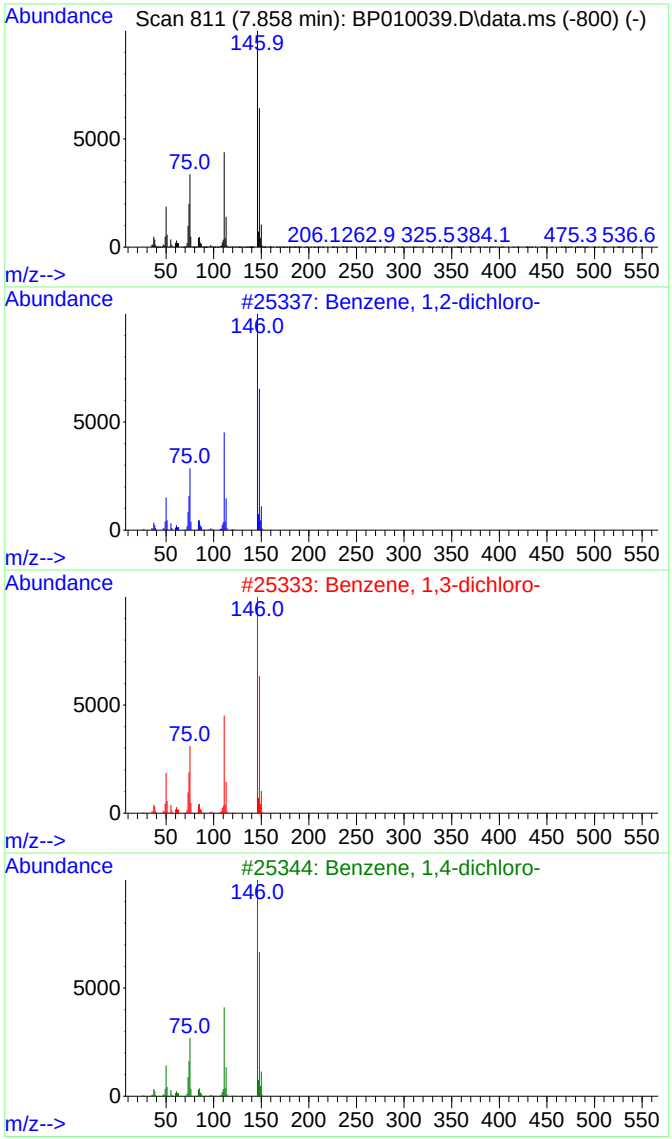
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

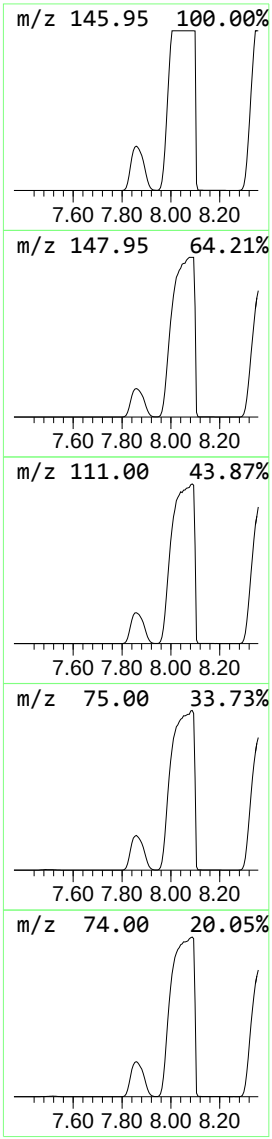
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.858	2.30 ng/ul	34236300	1,4-Dichlorobenzene-d4	7.999

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
4	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

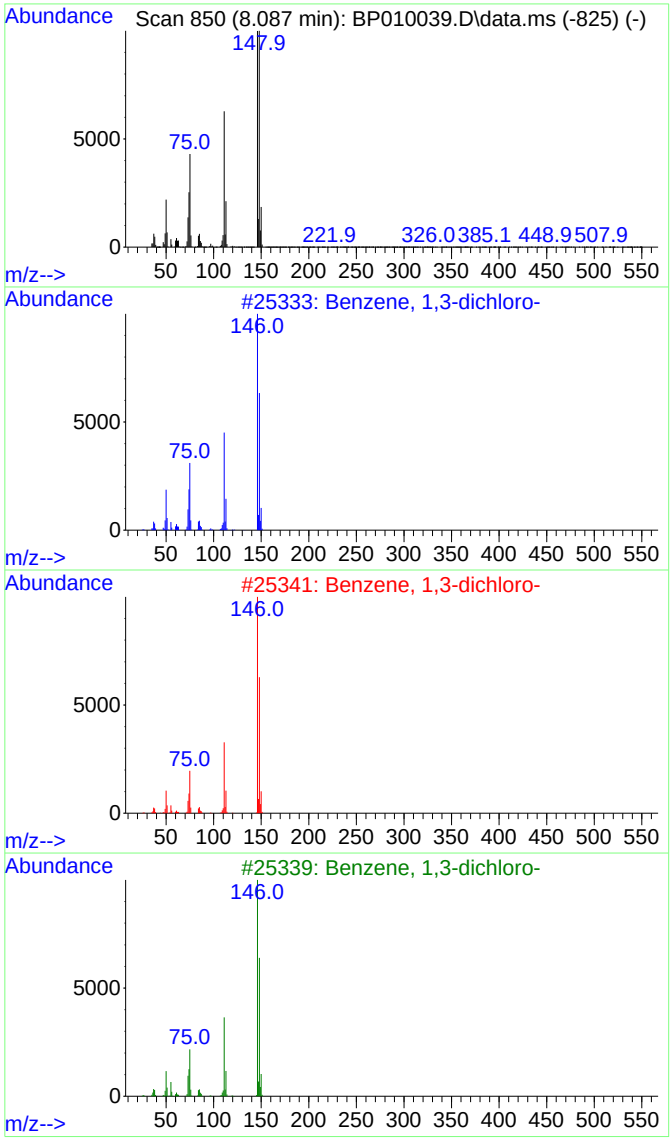
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

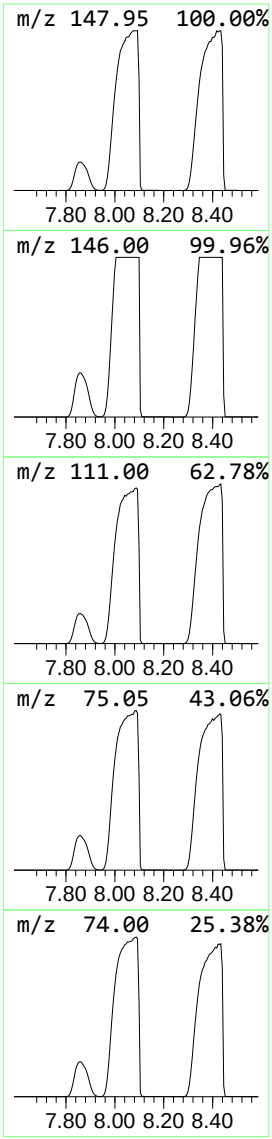
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.087	20.00 ng/ul	297702000	1,4-Dichlorobenzene-d4	7.999

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
2	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
3	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
4	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93
5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

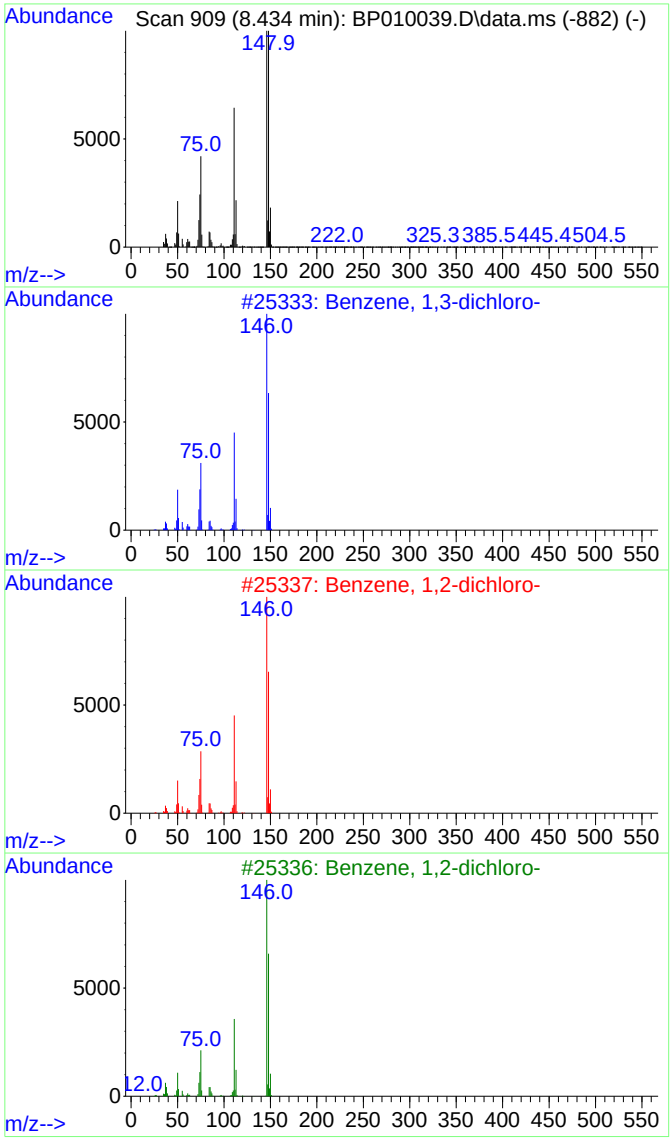
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

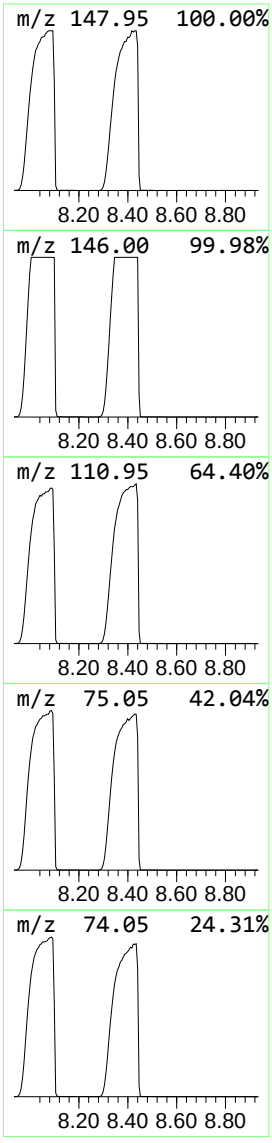
****

Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.434	20.40 ng/ul	303684000	1,4-Dichlorobenzene-d4	7.999

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
2	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
3	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
4	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
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Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

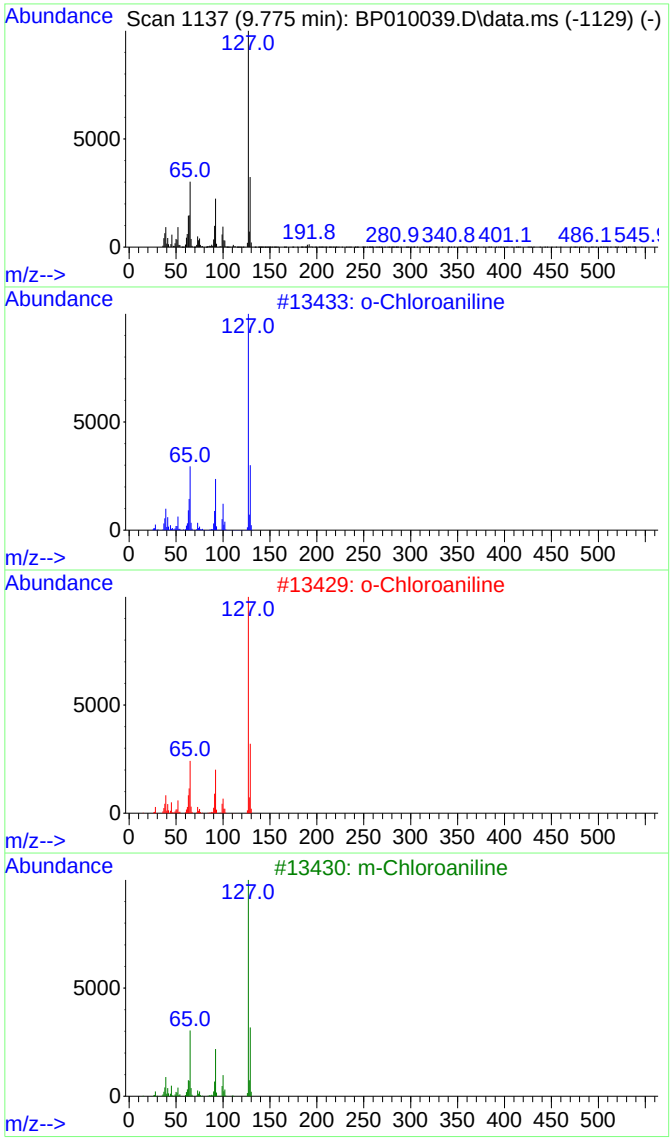
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

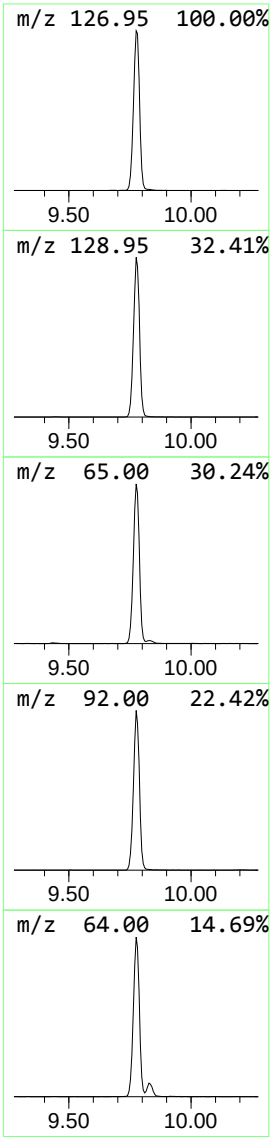
****

Peak Number 5 o-Chloroaniline Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.775	81.70 ng/ul	5926720	Naphthalene-d8	10.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Chloroaniline	127	C6H6ClN	000095-51-2	96
2	o-Chloroaniline	127	C6H6ClN	000095-51-2	96
3	m-Chloroaniline	127	C6H6ClN	000108-42-9	96
4	o-Chloroaniline	127	C6H6ClN	000095-51-2	95
5	m-Chloroaniline	127	C6H6ClN	000108-42-9	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

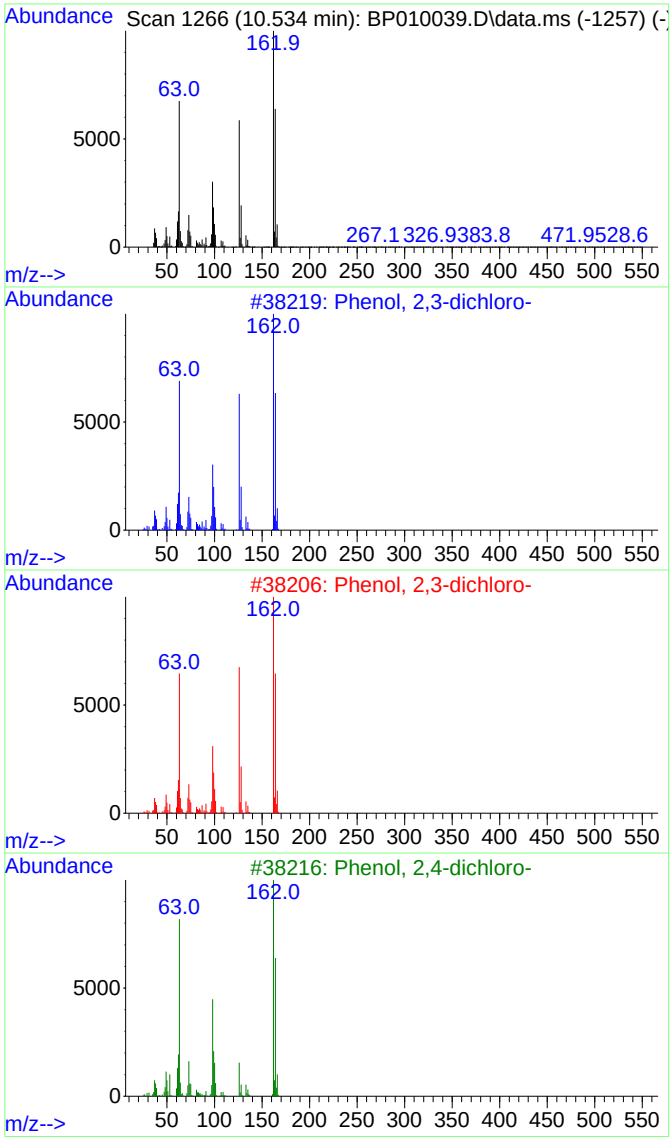
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

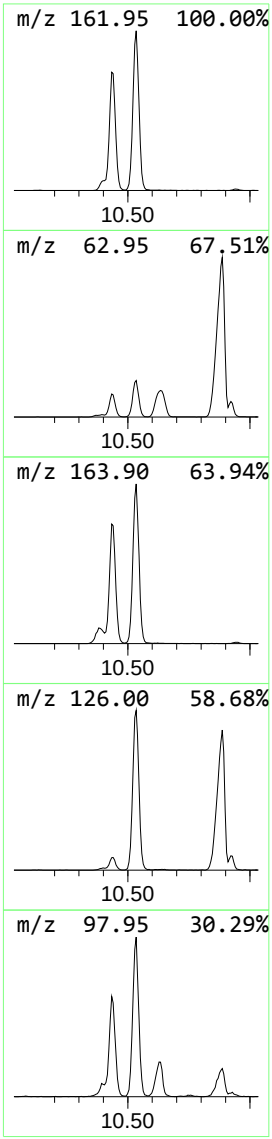
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Phenol, 2,3-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.534	29.00 ng/ul	2103990	Naphthalene-d8	10.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	99
2	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	98
3	Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	98
4	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	97
5	Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

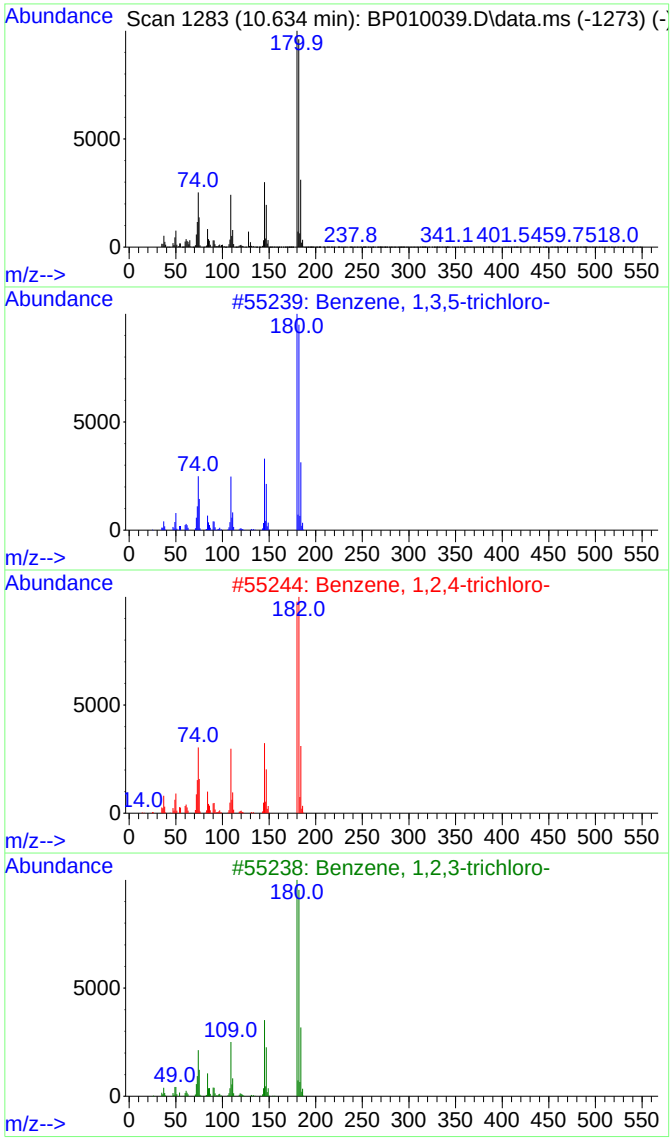
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

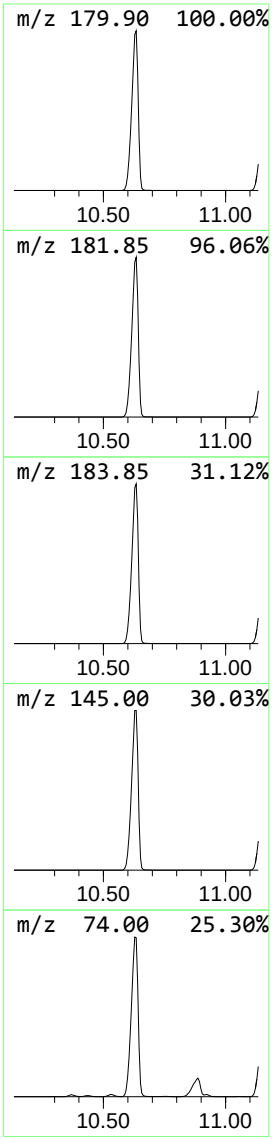
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.634	535.99 ng/ul	38883200	Naphthalene-d8	10.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
3	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97





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Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
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C0J22

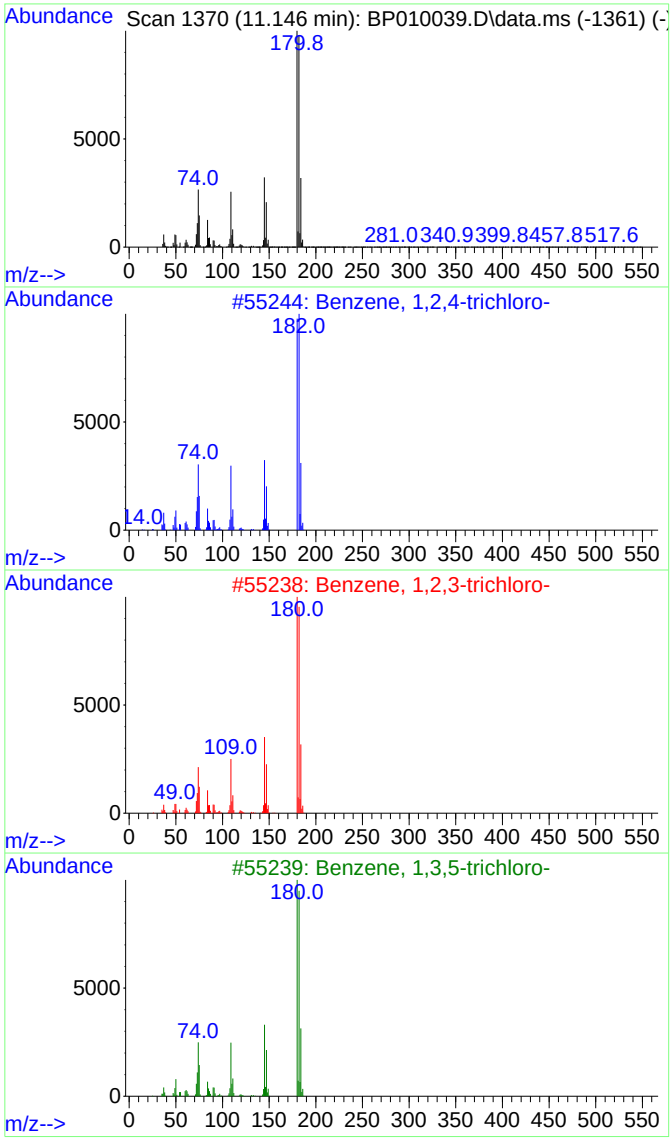
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

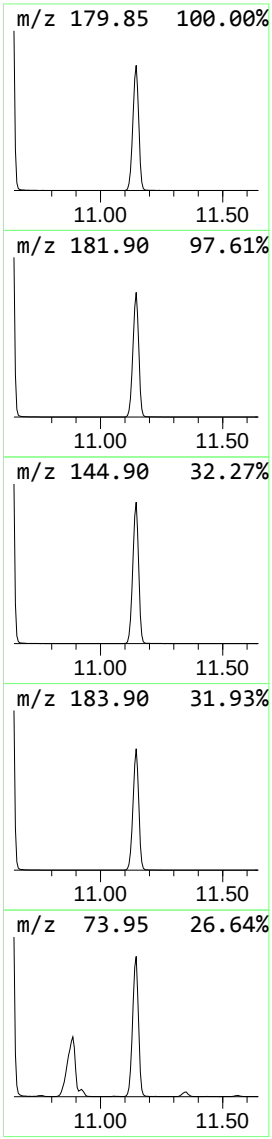
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,4-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.146	120.16 ng/ul	8716760	Naphthalene-d8	10.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
3	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
4	Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

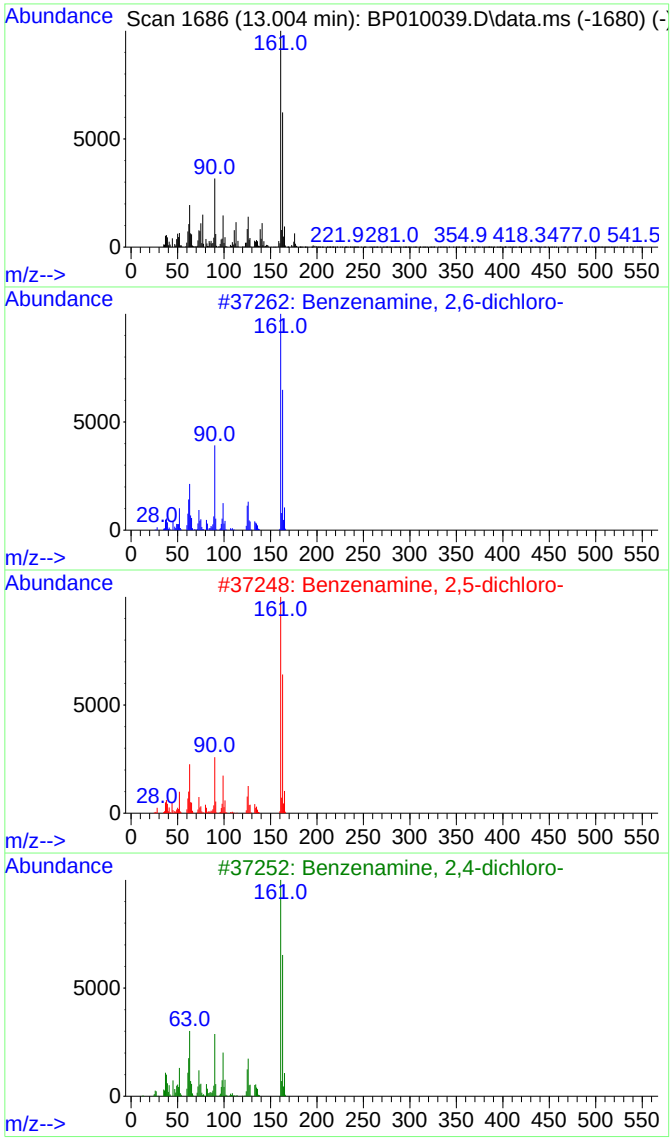
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

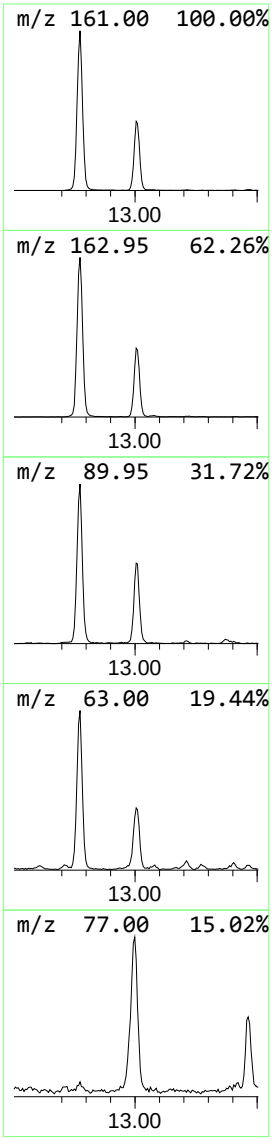
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzenamine, 2,6-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.004	6.02 ng/ul	639376	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	98
2	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	96
3	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	96
4	Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	95
5	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	95





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Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

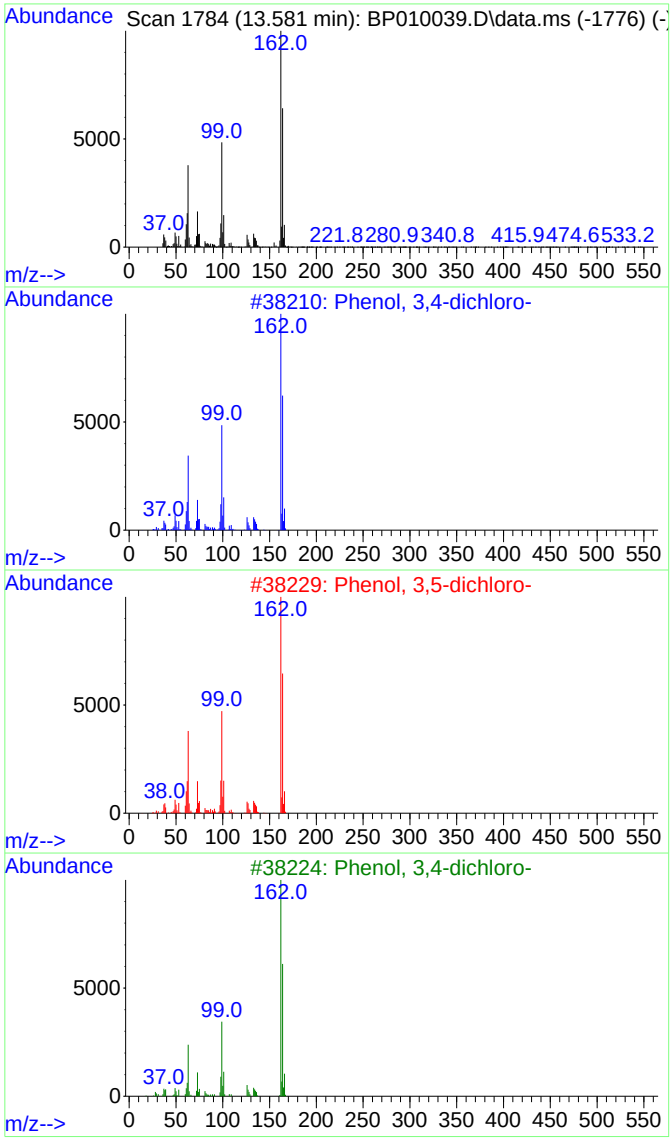
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

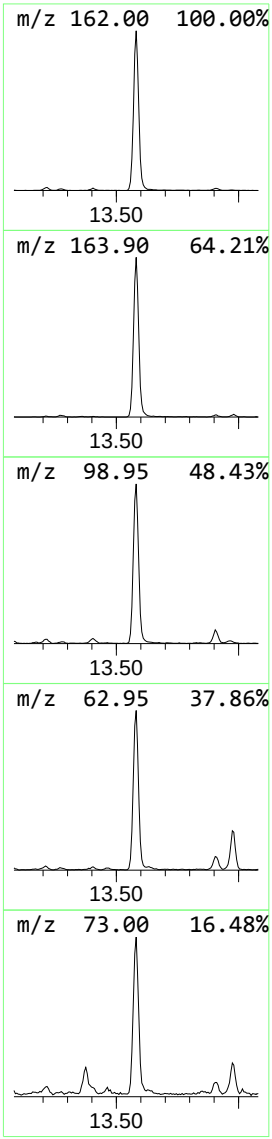
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Phenol, 3,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.581	14.15 ng/ul	1504340	Acenaphthene-d10	14.569

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
2	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
4	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	97
5	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

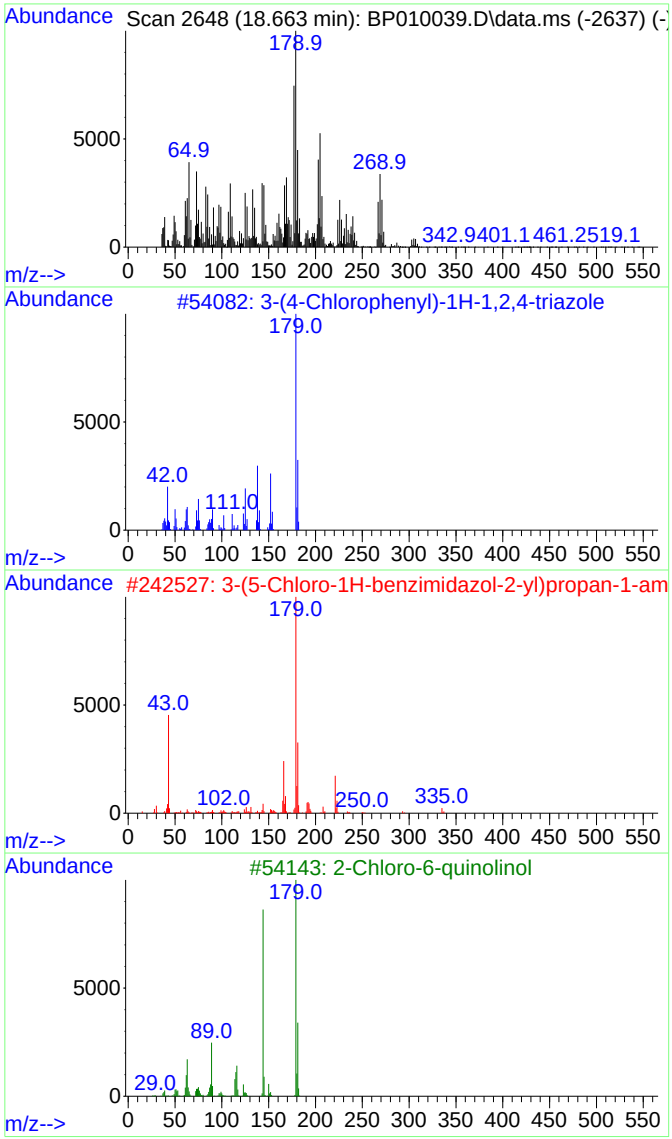
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

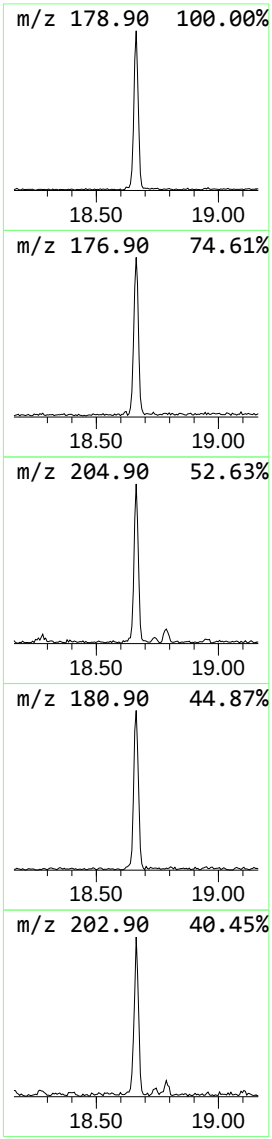
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.663	4.34 ng/ul	585055	Phenanthrene-d10	17.322

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(4-Chlorophenyl)-1H-1,2,4-tria...	179	C8H6ClN3	023195-59-7	18
2	3-(5-Chloro-1H-benzimidazol-2-yl...	335	C16H18ClN3O3	1000504-52-3	15
3	2-Chloro-6-quinolinol	179	C9H6ClNO	577967-89-6	14
4	trans-(2-Chlorovinyl)methyldieth...	194	C7H15ClO2Si	1000139-96-9	12
5	N-(4,6-Dichloro-1,3,5-triazin-2-...	220	C7H10Cl2N4	1000326-88-2	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

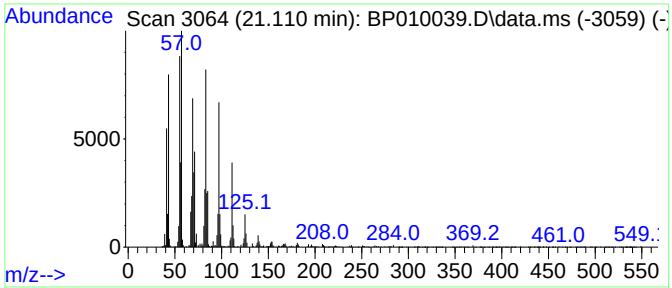
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

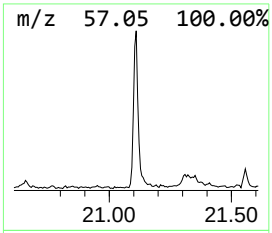
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

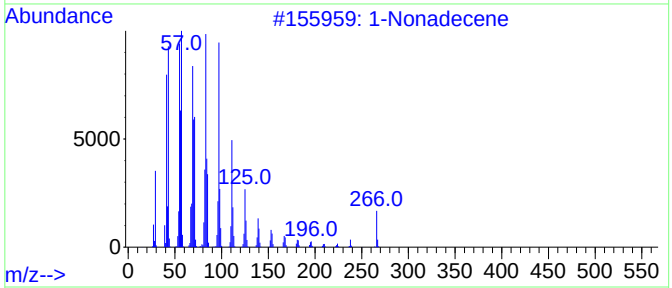
****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 11

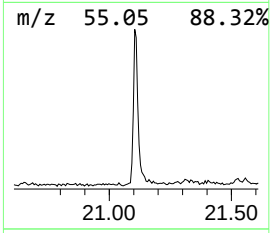
R.T.	EstConc	Area	Relative to ISTD	R.T.
21.110	3.14 ng/ul	477138	Chrysene-d12	21.404

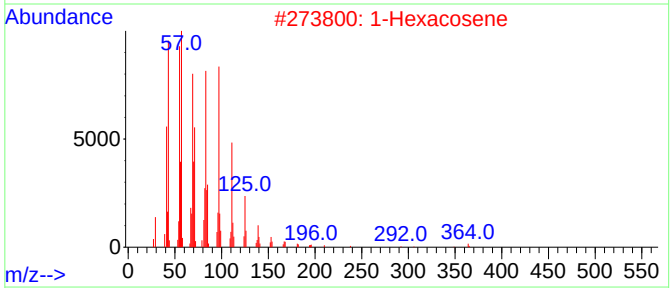
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	98
2	1-Hexacosene	364	C26H52	018835-33-1	94
3	1-Tetracosene	336	C24H48	010192-32-2	94
4	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94

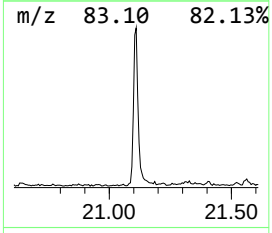


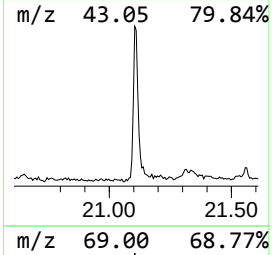


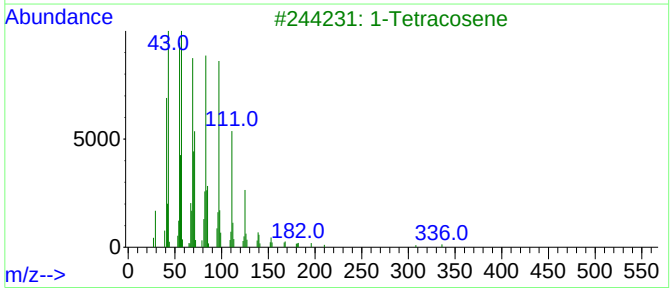


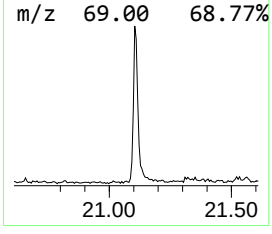












Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010039.D
Acq On : 22 Apr 2022 05:40
Operator : CG/JU
Sample : N2490-05
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J22

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, chloro-	5.393	29.5	ng/ul	439391000	1	7.999	297702000	20.0
Benzene, 1,2-di...	7.858	2.3	ng/ul	34236300	1	7.999	297702000	20.0
Benzene, 1,3-di...	8.087	20.0	ng/ul	297702000	1	7.999	297702000	20.0
unknown-01	8.434	20.4	ng/ul	303684000	1	7.999	297702000	20.0
o-Chloroaniline	9.775	81.7	ng/ul	5926720	2	10.757	1450880	20.0
Phenol, 2,3-dic...	10.534	29.0	ng/ul	2103990	2	10.757	1450880	20.0
Benzene, 1,3,5-...	10.634	536.0	ng/ul	38883200	2	10.757	1450880	20.0
Benzene, 1,2,4-...	11.146	120.2	ng/ul	8716760	2	10.757	1450880	20.0
Benzenamine, 2,...	13.004	6.0	ng/ul	639376	3	14.569	2125880	20.0
Phenol, 3,4-dic...	13.581	14.2	ng/ul	1504340	3	14.569	2125880	20.0
unknown-02	18.663	4.3	ng/ul	585055	4	17.322	2696590	20.0
(DEL) Alkane: S...	21.110	3.1	ng/ul	477138	5	21.404	3039610	20.0