

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032819\
 Data File : BM019550.D
 Acq On : 28 Mar 2019 20:46
 Operator : JU/SJ
 Sample : K2069-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09C7

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.069	323	354	359	rBV7	54261941	497094338	100.00%	38.359%
2	6.281	554	560	574	rVB	466717	781216	0.16%	0.060%
3	6.969	668	677	691	rBV	275746	808694	0.16%	0.062%
4	7.128	698	704	718	rVB	388314	864569	0.17%	0.067%
5	7.510	754	769	780	rBV2	41682776	160919129	32.37%	12.418%
6	7.704	780	802	803	rBV5	50565635	244547254	49.20%	18.871%
7	8.022	836	856	857	rBV4	50617332	228430670	45.95%	17.627%
8	8.310	899	905	919	rVB	366925	642708	0.13%	0.050%
9	8.763	976	982	995	rBV	629202	1040743	0.21%	0.080%
10	9.081	1029	1036	1045	rBV	335598	530264	0.11%	0.041%
11	9.475	1096	1103	1122	rVB	527912	986361	0.20%	0.076%
12	10.004	1186	1193	1208	rBV	706302	1307959	0.26%	0.101%
13	10.269	1224	1238	1242	rBV2	45425134	107307279	21.59%	8.281%
14	10.381	1251	1257	1262	rBV	712723	1060154	0.21%	0.082%
15	10.757	1310	1321	1334	rBV	10492696	17353900	3.49%	1.339%
16	12.410	1591	1602	1614	rBV	1525757	2570696	0.52%	0.198%
17	13.027	1700	1707	1720	rVB	1219110	1917075	0.39%	0.148%
18	13.663	1809	1815	1822	rBV	1394937	1943605	0.39%	0.150%
19	13.933	1854	1861	1873	rBV	1756100	2455775	0.49%	0.190%
20	14.239	1907	1913	1922	rBV2	870667	1367841	0.28%	0.106%
21	15.239	2077	2083	2101	rBV	2074332	2961977	0.60%	0.229%
22	15.386	2103	2108	2121	rBV	568712	908571	0.18%	0.070%
23	16.998	2376	2382	2393	rBV2	1113494	1561900	0.31%	0.121%
24	17.098	2393	2399	2411	rBV2	2140375	3116497	0.63%	0.240%
25	19.409	2786	2792	2804	rBV	2731637	3710772	0.75%	0.286%
26	21.209	3093	3098	3106	rBV	1608995	2079632	0.42%	0.160%
27	23.274	3442	3449	3458	rBV2	2912087	5076636	1.02%	0.392%
28	23.415	3466	3473	3487	rVB	1364429	2556632	0.51%	0.197%

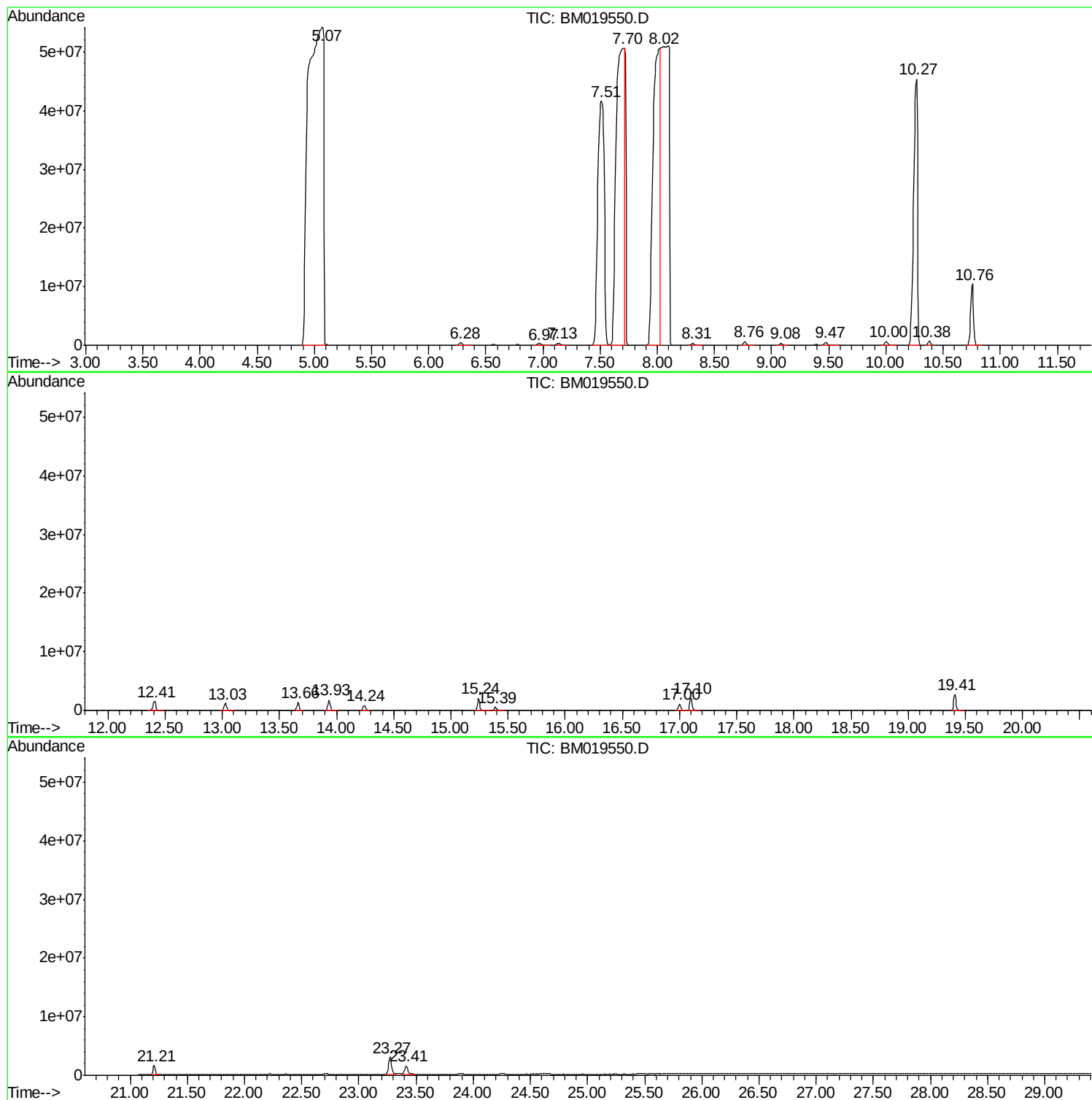
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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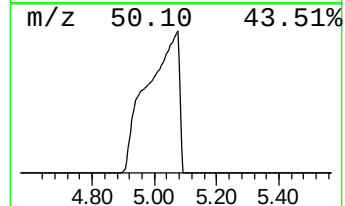
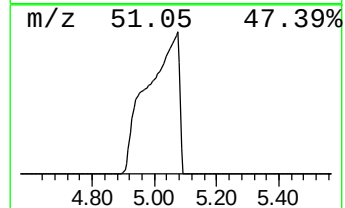
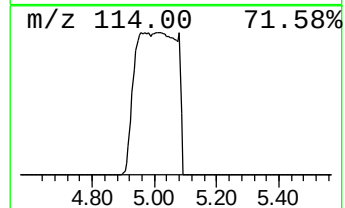
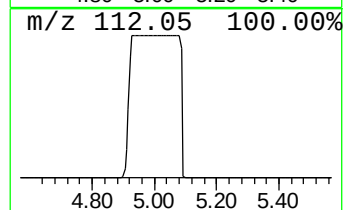
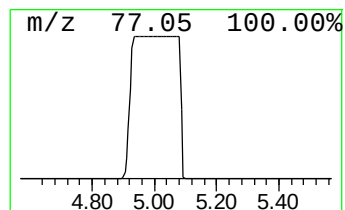
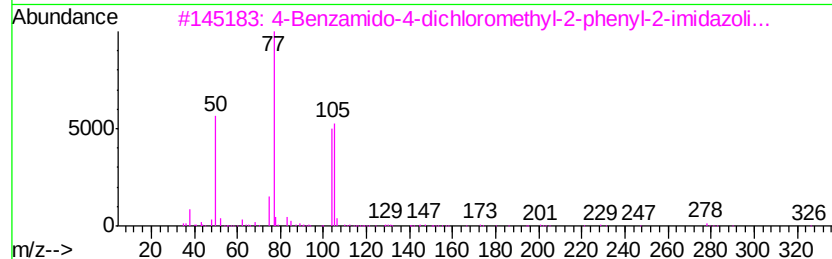
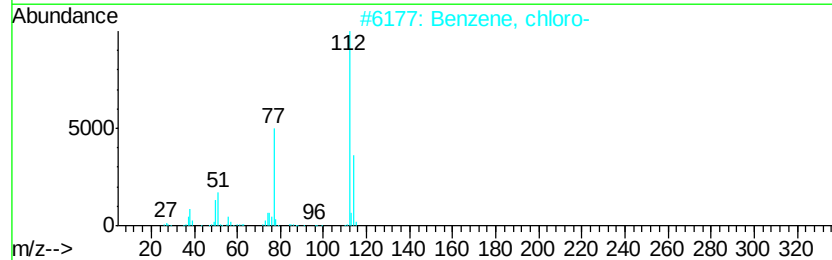
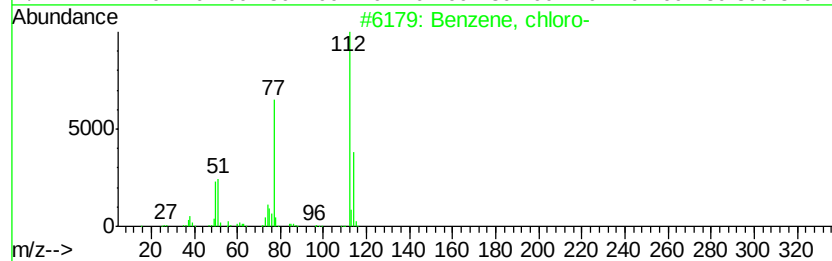
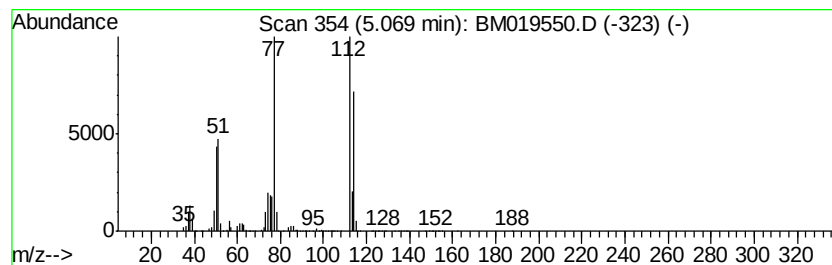
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	40.65 ng/ul	497094000	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	93
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	64
3		4-Benzamido-4-dichloromethyl-2-p...	361	C17H13Cl2N3O2	076543-29-8	16
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	16
5		Benzene, chloro-	112	C6H5Cl	000108-90-7	12



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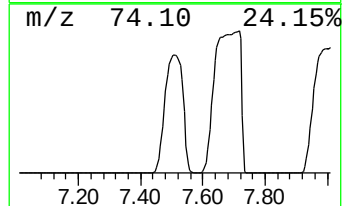
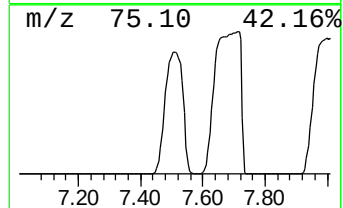
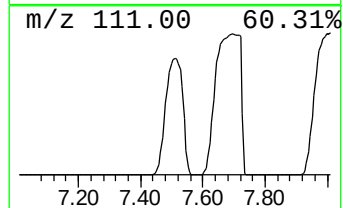
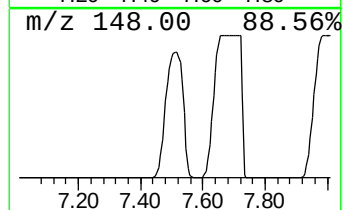
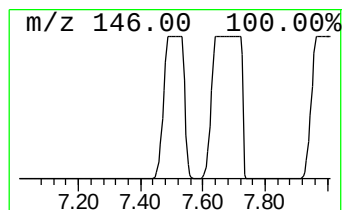
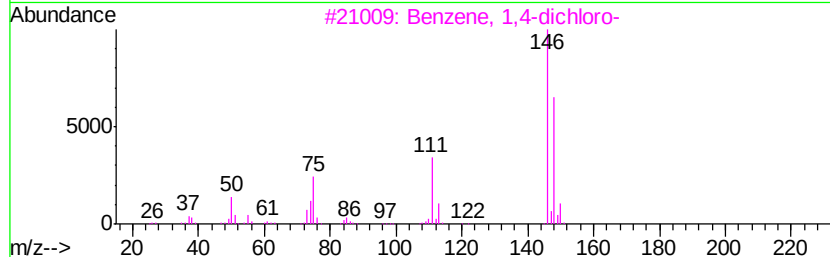
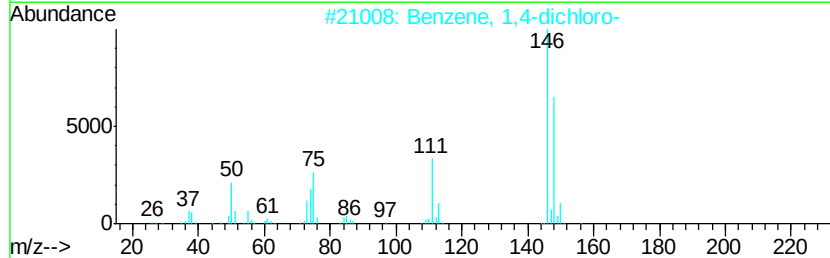
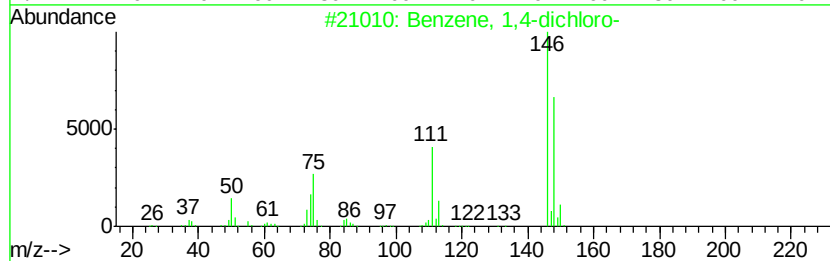
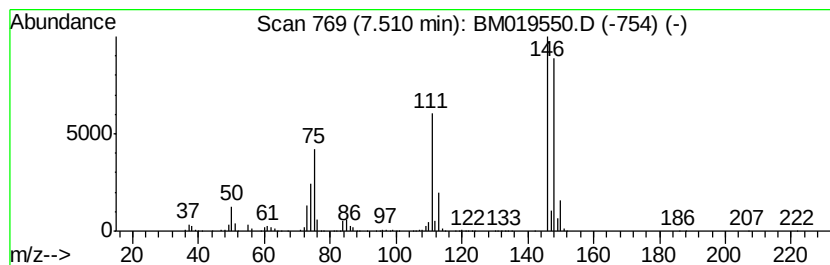
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.51	13.16 ng/ul	160919000	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
5		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93



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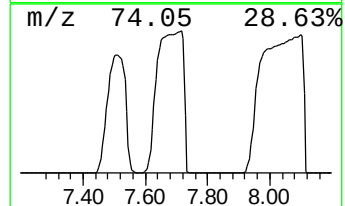
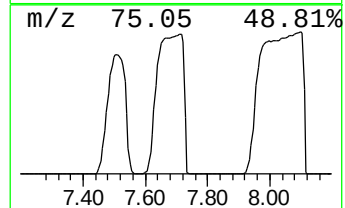
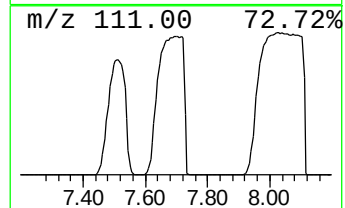
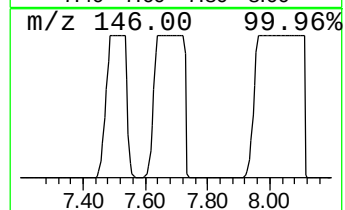
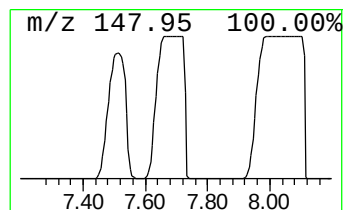
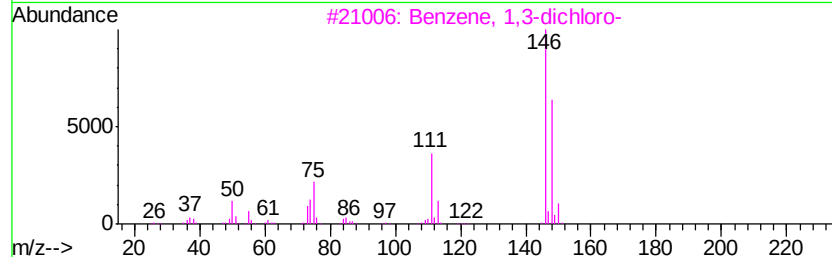
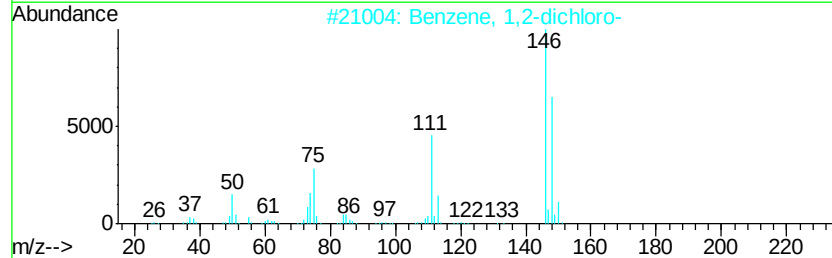
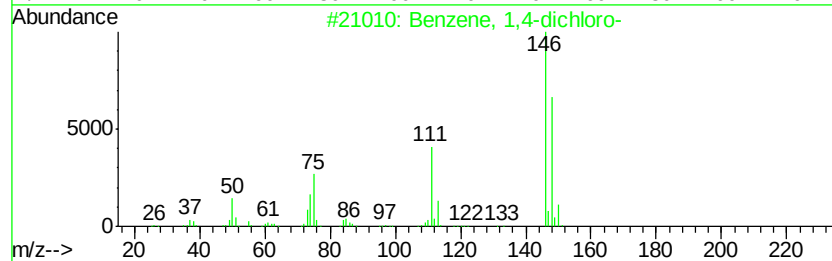
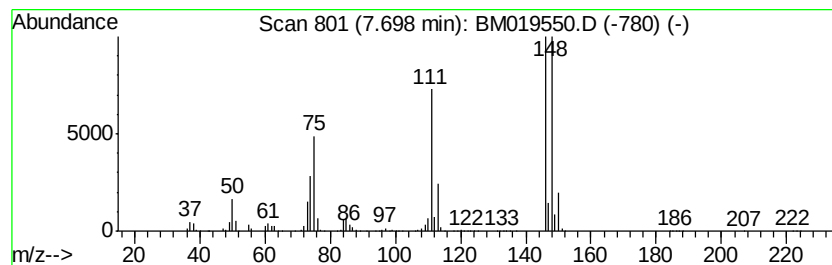
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	20.00 ng/ul	244547000	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
2		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	81
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	81
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81



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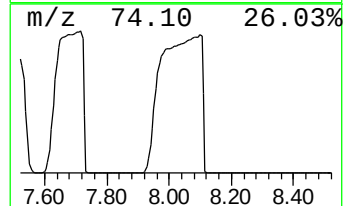
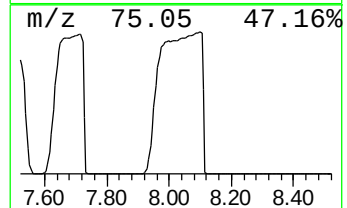
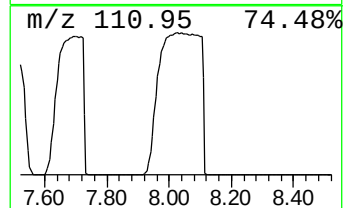
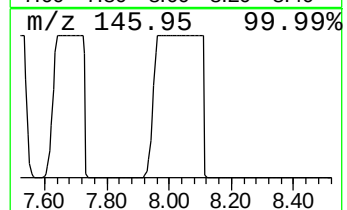
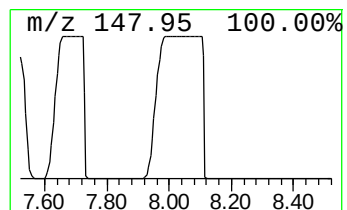
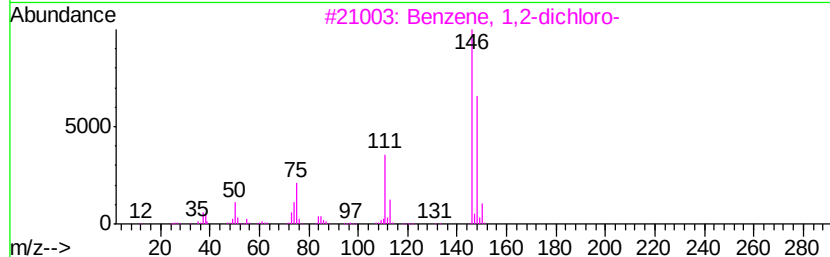
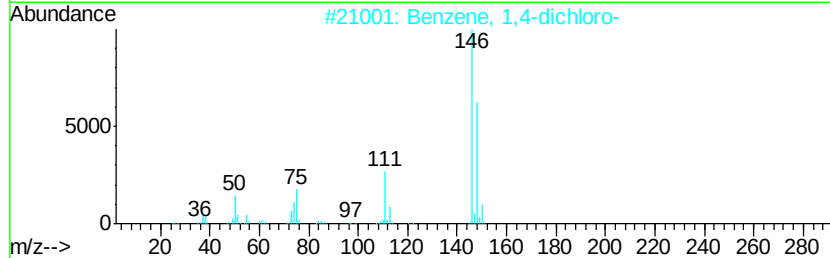
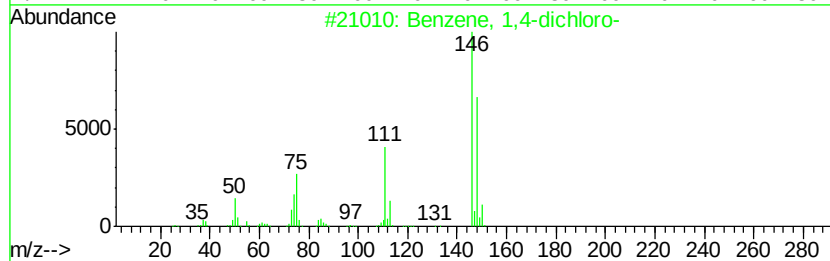
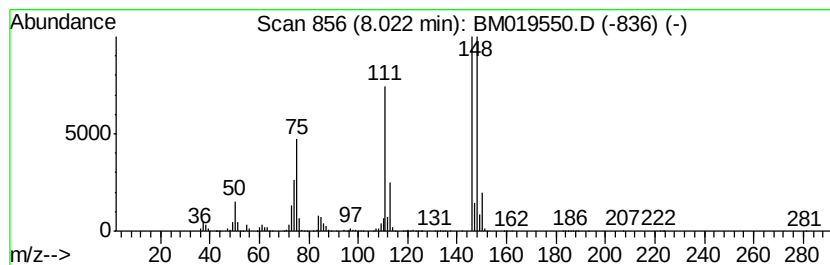
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Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	18.68 ng/ul	228431000	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	81
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	81



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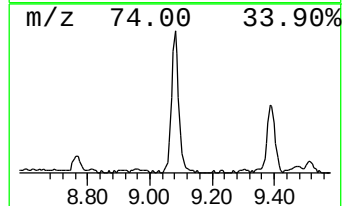
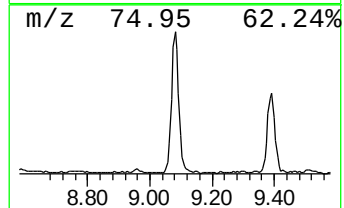
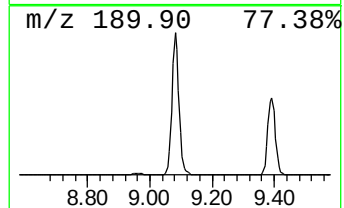
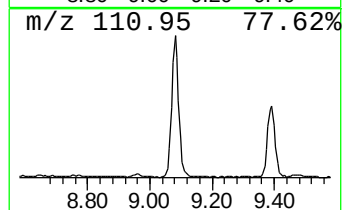
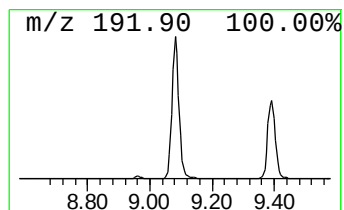
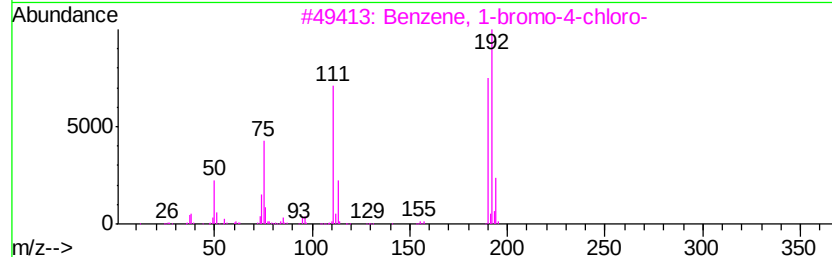
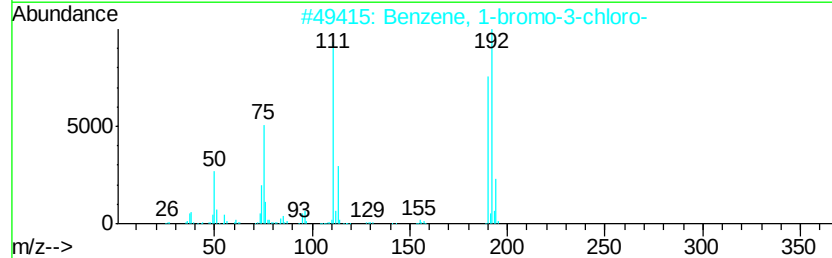
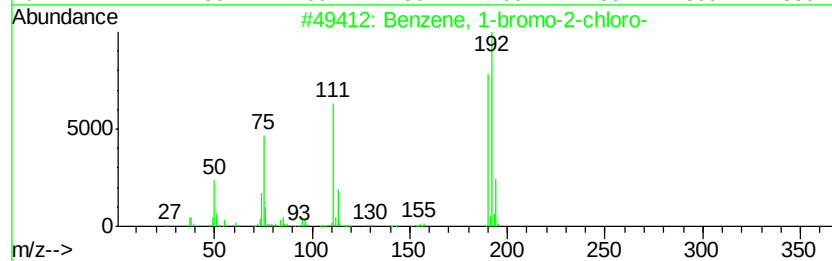
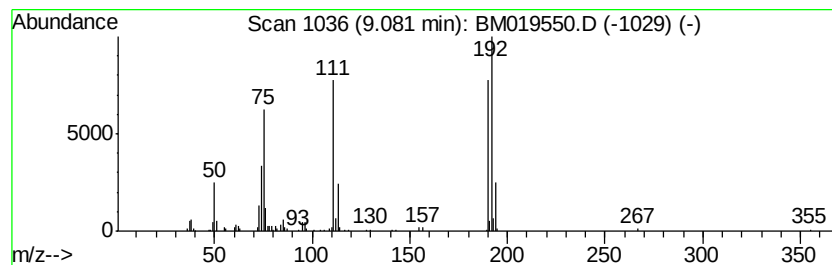
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-bromo-2-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.08	10.00 ng/ul	530264	Napthalene-d8	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96



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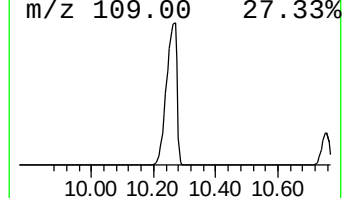
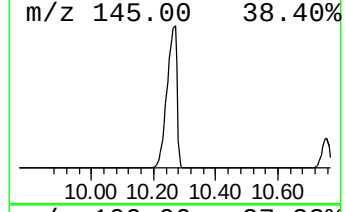
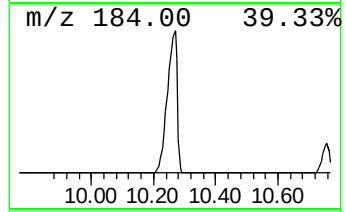
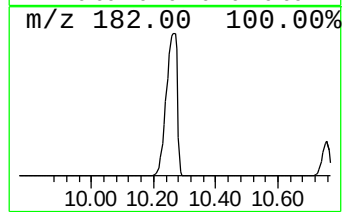
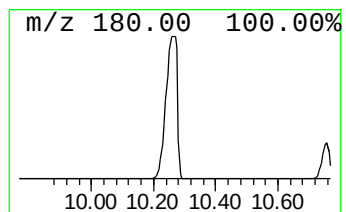
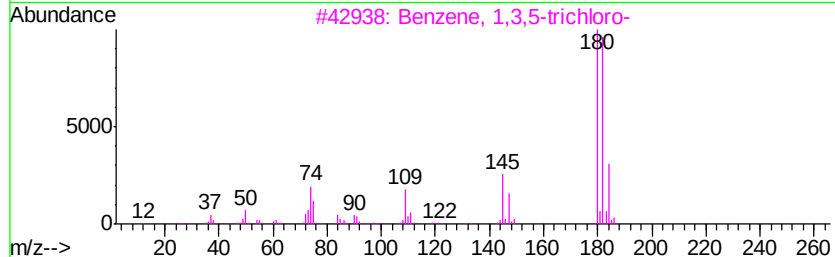
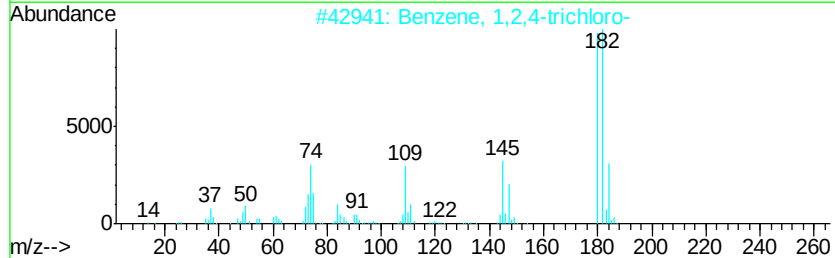
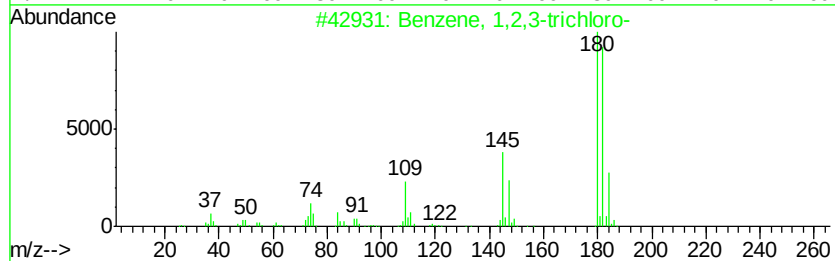
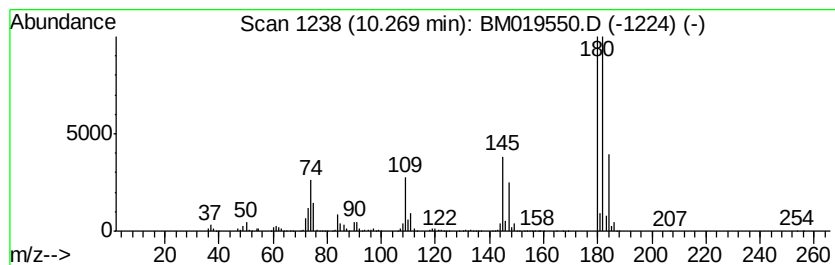
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,3-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	2024.37 ng/ul	107307000	Napthalene-d8	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
2		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	96
3		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95
4		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	95
5		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032819\
Data File : BM019550.D
Acq On : 28 Mar 2019 20:46
Operator : JU/SJ
Sample : K2069-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09C7

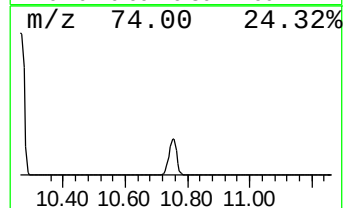
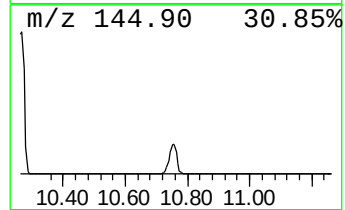
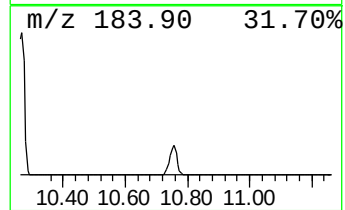
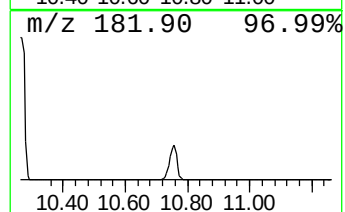
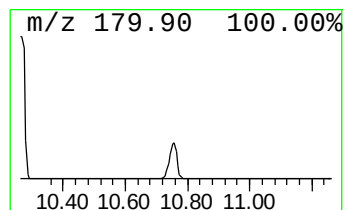
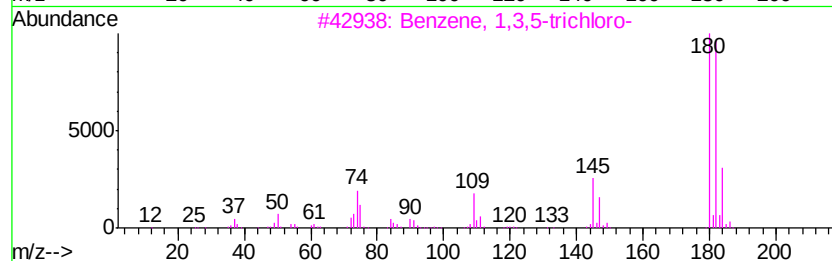
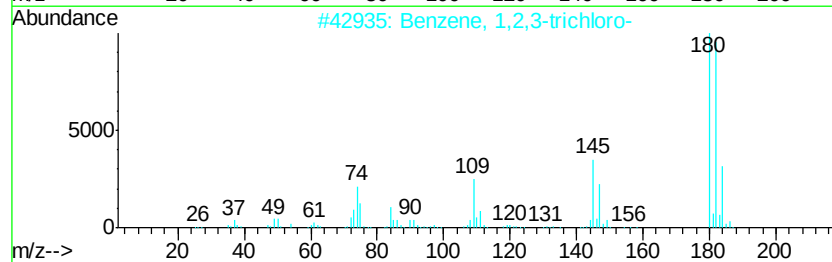
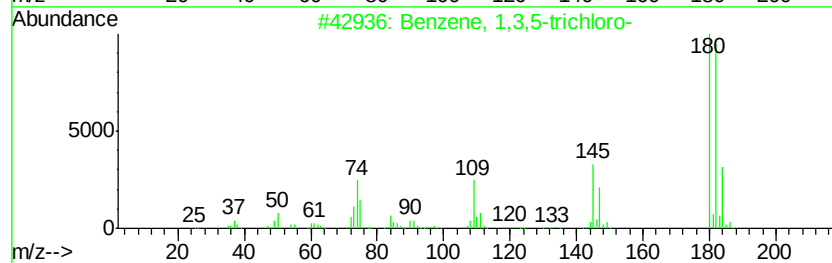
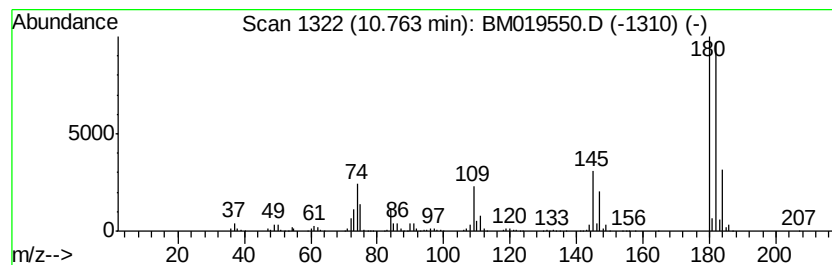
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,3,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	327.39 ng/ul	17353900	Naphthalene-d8	10.38

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032819\
Data File : BM019550.D
Acq On : 28 Mar 2019 20:46
Operator : JU/SJ
Sample : K2069-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09C7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.07	40.6	ng/ul	497094000	1	7.65	244547000	20.0
Benzene, 1,4-dich...	7.51	13.2	ng/ul	160919000	1	7.65	244547000	20.0
Benzene, 1,2-dich...	7.70	20.0	ng/ul	244547000	1	7.65	244547000	20.0
Benzene, 1,3-dich...	8.02	18.7	ng/ul	228431000	1	7.65	244547000	20.0
Benzene, 1-bromo-...	9.08	10.0	ng/ul	530264	2	10.38	1060150	20.0
Benzene, 1,2,3-tr...	10.27	2024.4	ng/ul	107307000	2	10.38	1060150	20.0
Benzene, 1,3,5-tr...	10.76	327.4	ng/ul	17353900	2	10.38	1060150	20.0