

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\
 Data File : BM019319.D
 Acq On : 22 Mar 2019 19:08
 Operator : JU/SJ
 Sample : K1968-14
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09A9

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	4.969	328	337	352	rBV	34584912	54018590	100.00%	27.672%
2	6.799	642	648	666	rVB	131746	238657	0.44%	0.122%
3	6.963	670	676	692	rBV	477090	825472	1.53%	0.423%
4	7.152	701	708	720	rBV	557463	913857	1.69%	0.468%
5	7.504	760	768	781	rBV	3513445	5508386	10.20%	2.822%
6	7.657	781	794	812	rVV	17349631	26755935	49.53%	13.706%
7	7.969	838	847	871	rBV	17200901	27303540	50.54%	13.987%
8	8.328	902	908	923	rBV	357660	658326	1.22%	0.337%
9	8.787	978	986	992	rBV	578440	1025158	1.90%	0.525%
10	9.104	1034	1040	1047	rBV	156921	252563	0.47%	0.129%
11	9.328	1073	1078	1082	rBV	57262	94311	0.17%	0.048%
12	9.375	1082	1086	1089	rVV2	57439	108874	0.20%	0.056%
13	9.416	1089	1093	1101	rVB	84450	145636	0.27%	0.075%
14	9.498	1101	1107	1121	rBV	472214	869155	1.61%	0.445%
15	10.028	1191	1197	1220	rBV	677984	1251345	2.32%	0.641%
16	10.275	1229	1239	1245	rBV	19902245	33486812	61.99%	17.154%
17	10.334	1245	1249	1254	rVV	69688	162560	0.30%	0.083%
18	10.404	1255	1261	1267	rVB	661324	1084365	2.01%	0.555%
19	10.781	1317	1325	1346	rBV	4880844	7800585	14.44%	3.996%
20	12.410	1596	1602	1603	rBV	135468	192291	0.36%	0.099%
21	12.439	1603	1607	1617	rVB	469287	766015	1.42%	0.392%
22	12.710	1646	1653	1663	rBV3	30148	62117	0.11%	0.032%
23	13.057	1705	1712	1723	rBV	2028796	3127806	5.79%	1.602%
24	13.322	1752	1757	1771	rVB	52155	109208	0.20%	0.056%
25	13.692	1813	1820	1827	rBV	1552272	2146854	3.97%	1.100%
26	13.751	1827	1830	1844	rVB3	56482	101163	0.19%	0.052%
27	13.969	1860	1867	1882	rBV	1670748	2537051	4.70%	1.300%
28	14.275	1912	1919	1927	rBV2	929329	1416199	2.62%	0.725%
29	14.539	1958	1964	1968	rBV3	35620	80807	0.15%	0.041%
30	14.586	1968	1972	1980	rVB	159958	249947	0.46%	0.128%
31	15.275	2082	2089	2104	rBV	2080163	2977742	5.51%	1.525%
32	15.410	2107	2112	2127	rBV	614978	916211	1.70%	0.469%
33	16.010	2209	2214	2219	rBV	46415	67888	0.13%	0.035%
34	17.033	2382	2388	2397	rBV	1171074	1631670	3.02%	0.836%

LSC Area Percent Report

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35	17.133	2397	2405	2420	rBV	2197164	3234484	5.99%	1.657%
36	19.439	2791	2797	2804	rBV	2887816	3803614	7.04%	1.948%
37	21.239	3098	3103	3112	rBV	1552810	2036666	3.77%	1.043%
38	21.803	3197	3199	3204	rVB3	50519	54330	0.10%	0.028%
39	22.262	3274	3277	3284	rVB2	95115	116206	0.22%	0.060%
40	22.756	3359	3361	3366	rVB	126671	155634	0.29%	0.080%
41	23.315	3450	3456	3469	rBV2	2558017	4551896	8.43%	2.332%
42	23.456	3474	3480	3491	rVV	1152453	2144658	3.97%	1.099%
43	23.950	3560	3564	3569	rVB3	127686	227095	0.42%	0.116%

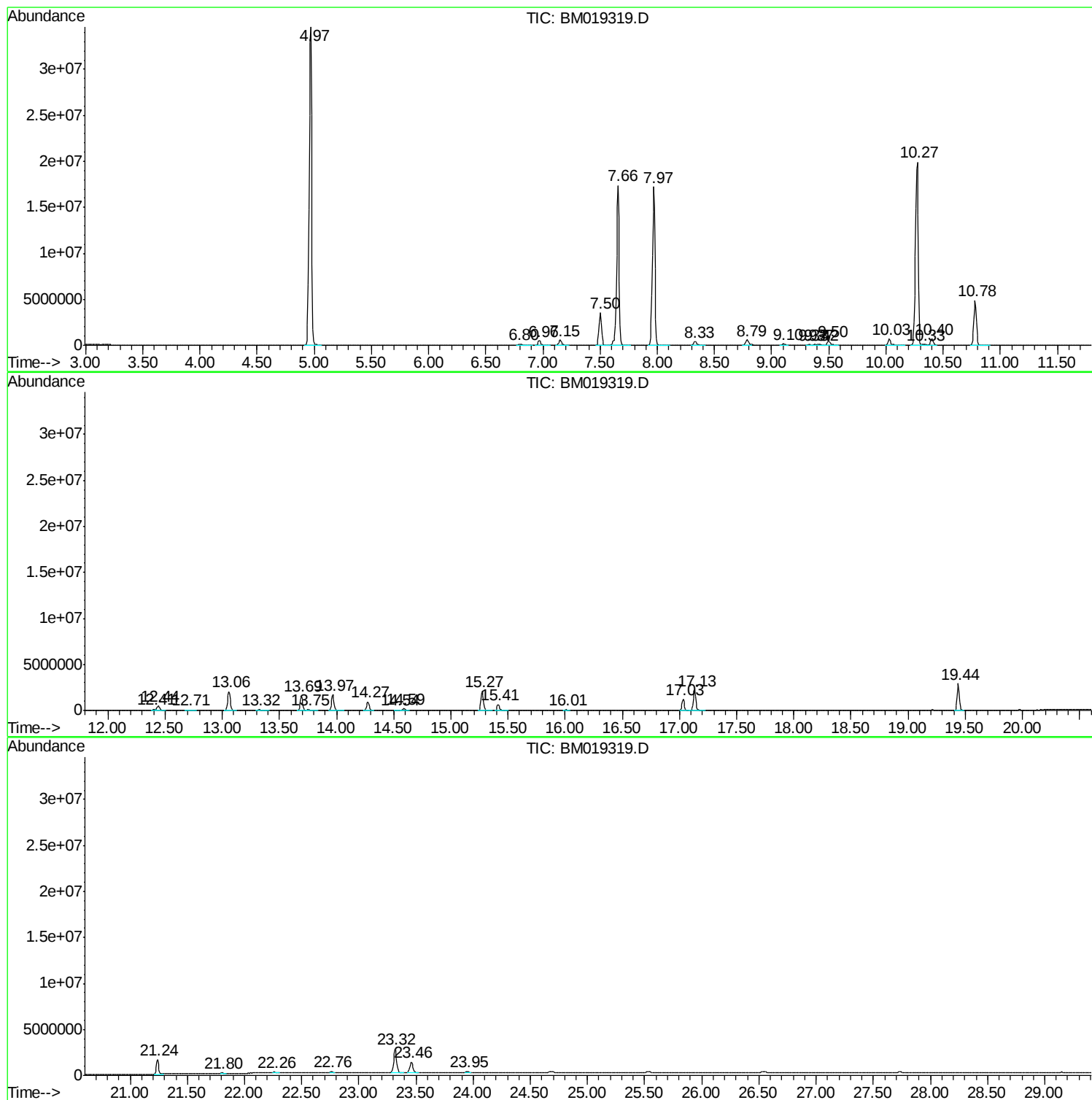
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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



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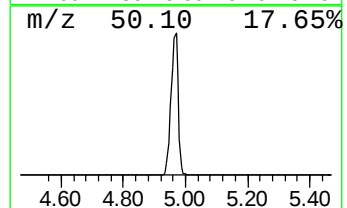
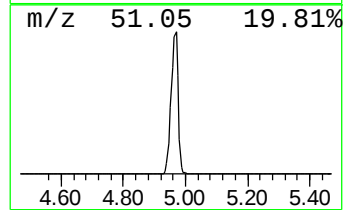
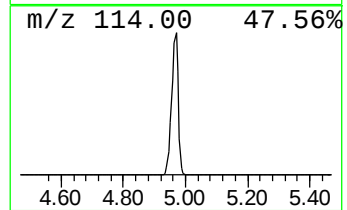
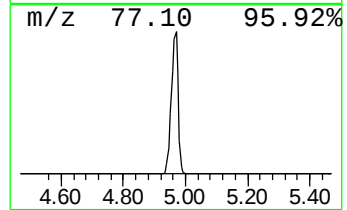
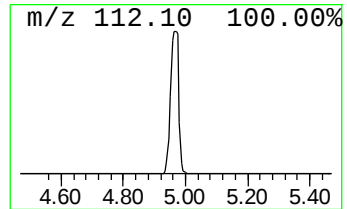
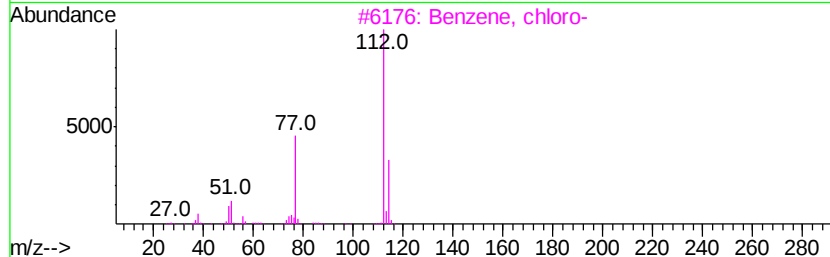
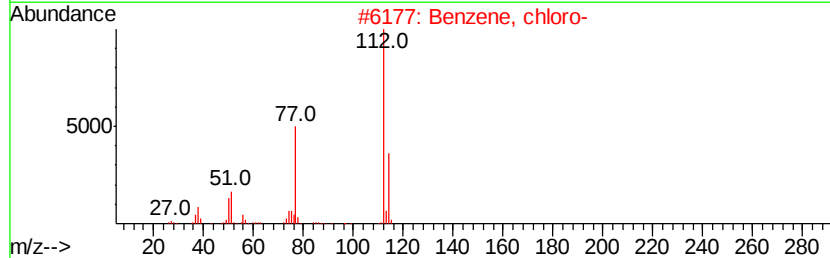
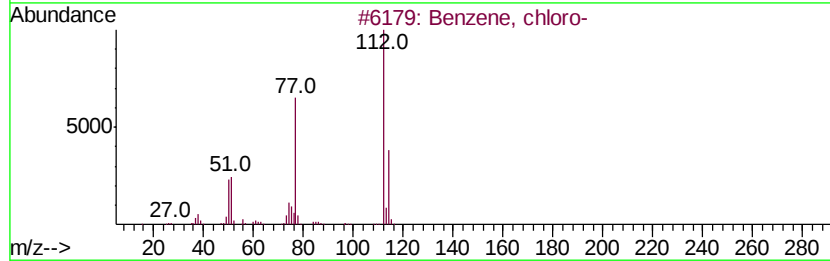
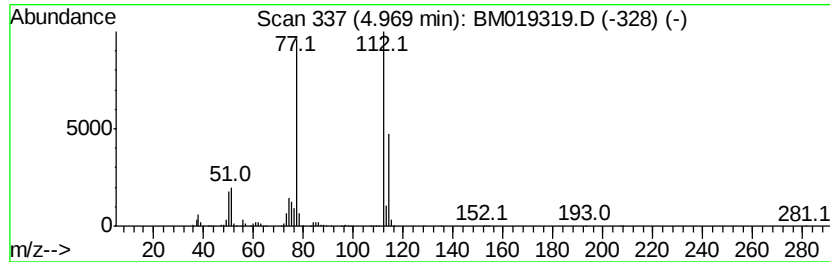
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	40.38 ng/ul	54018600	1,4-Dichlorobenzene-d4	7.62

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	93
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	90
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	90
5		Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	22



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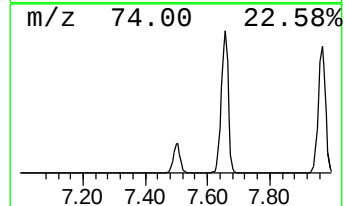
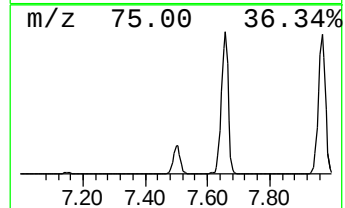
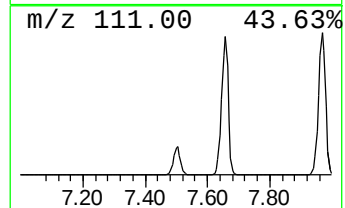
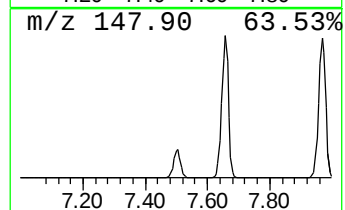
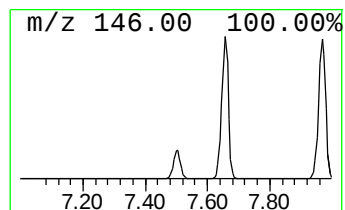
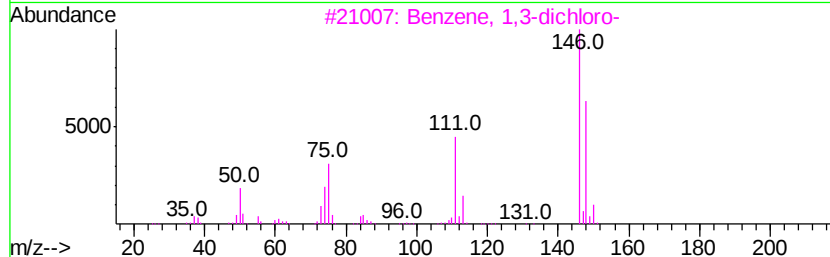
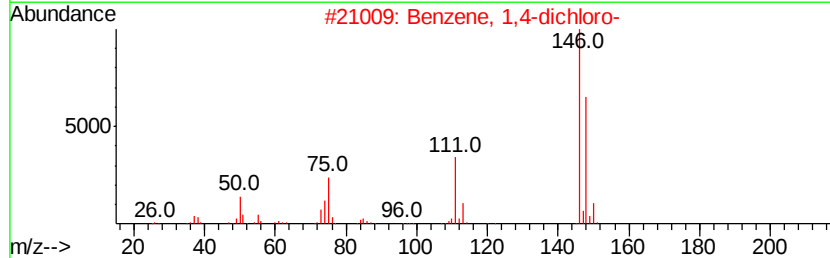
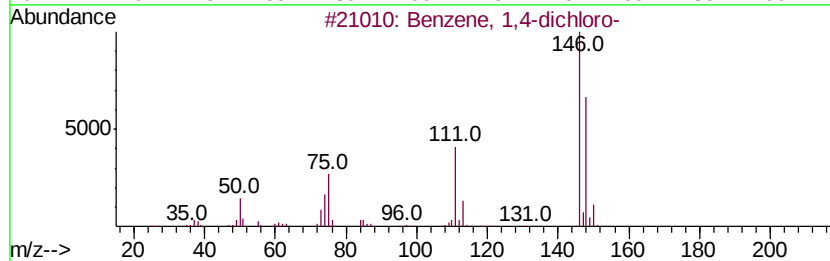
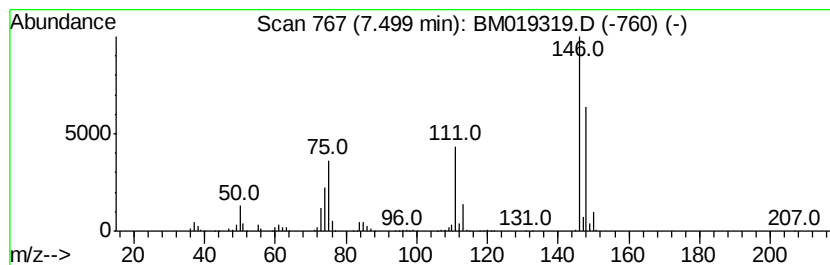
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Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	4.12 ng/ul	5508390	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95



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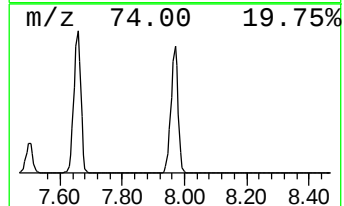
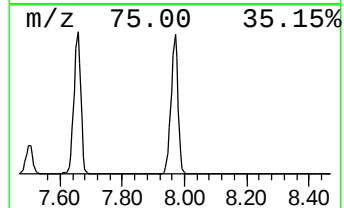
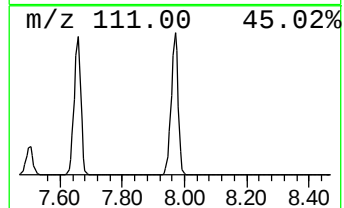
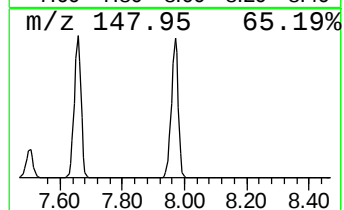
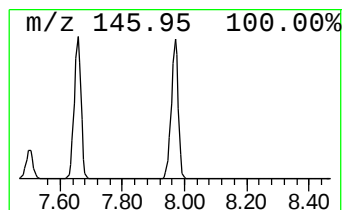
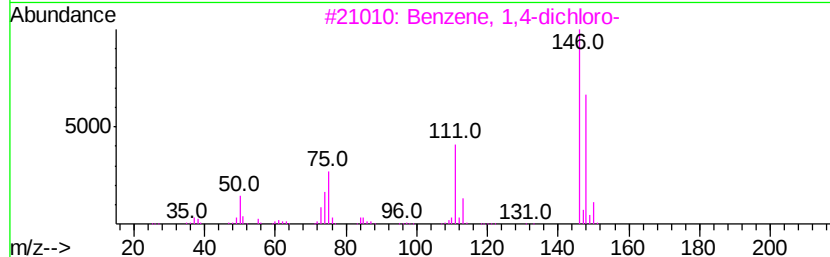
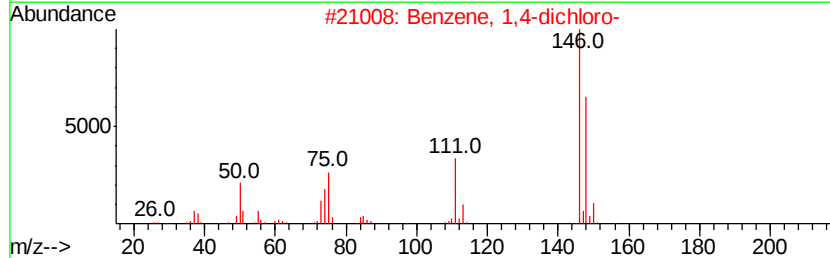
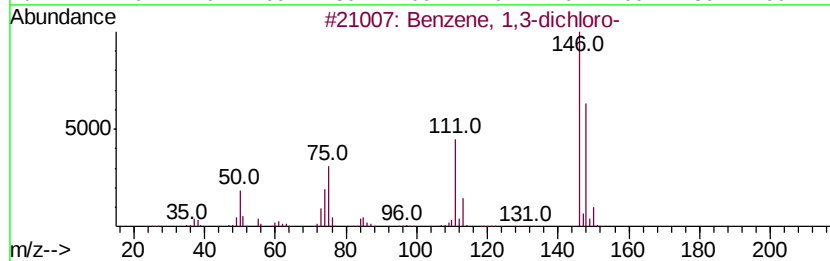
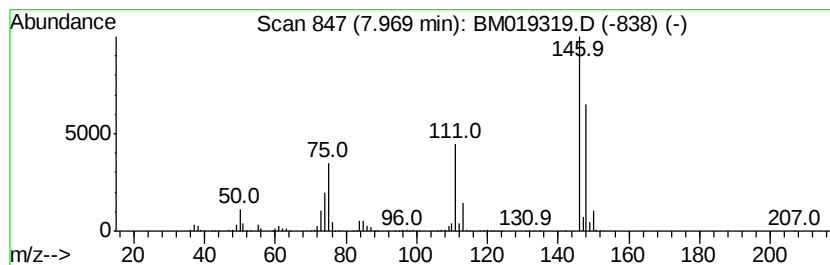
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	20.41 ng/ul	27303500	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
5		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95



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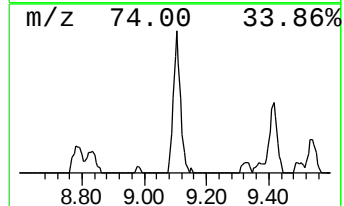
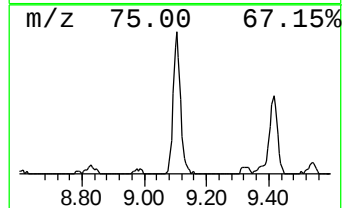
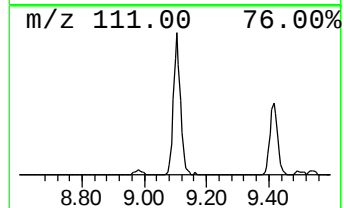
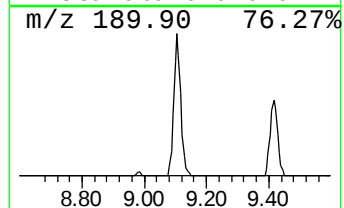
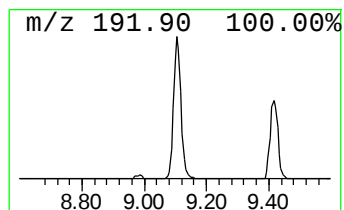
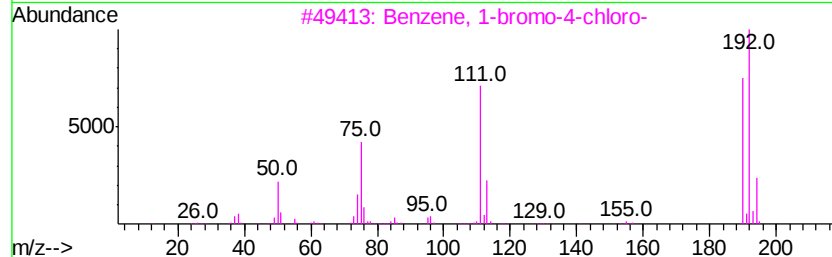
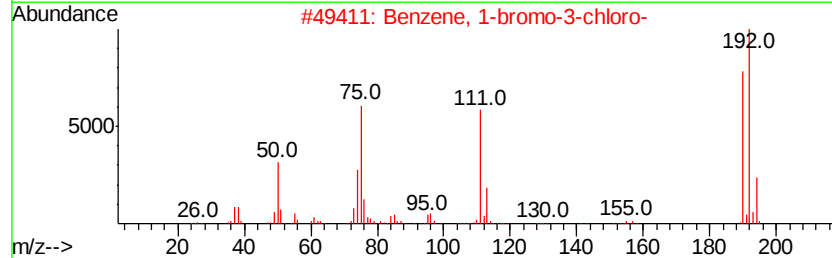
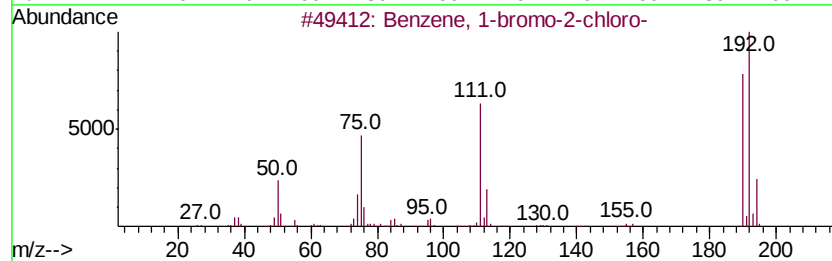
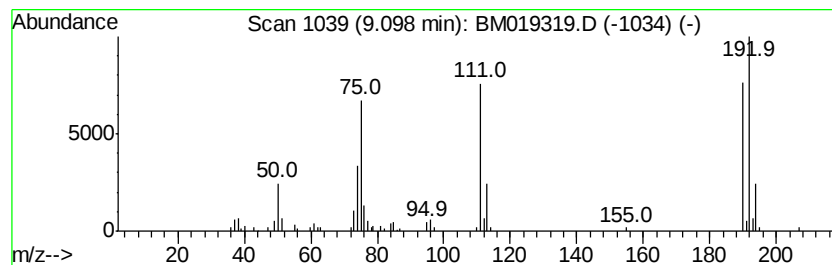
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Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	4.66 ng/ul	252563	Napthalene-d8	10.40

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	97
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	94
5		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	94



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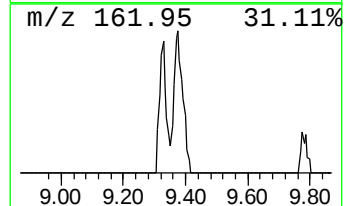
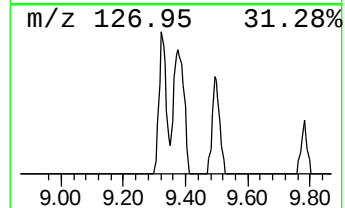
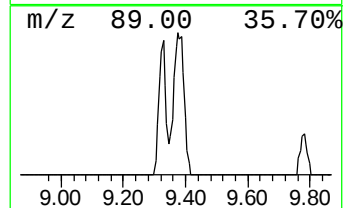
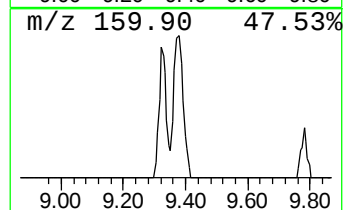
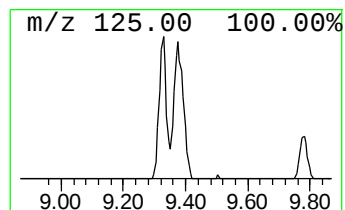
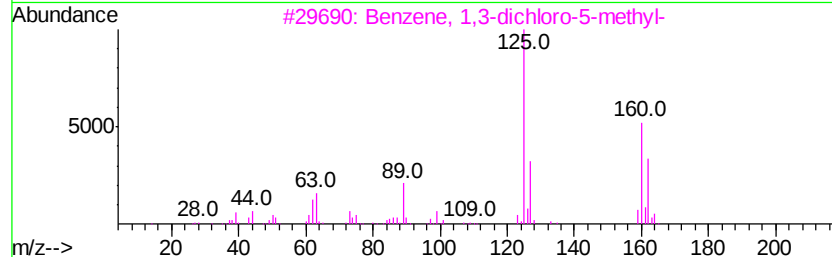
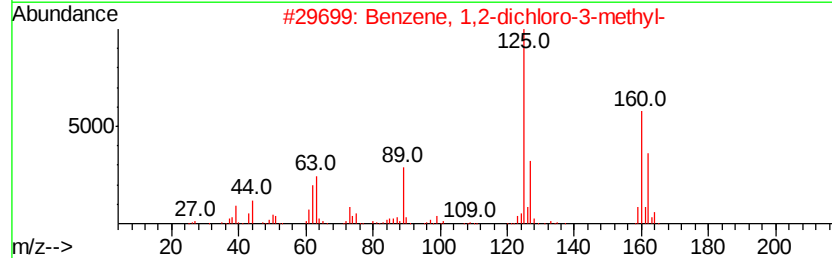
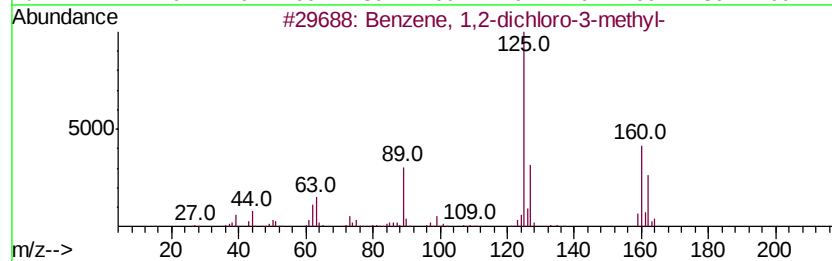
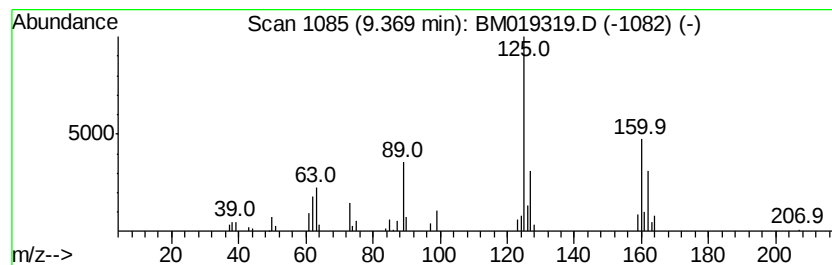
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2-dichloro-3-met... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	2.01 ng/ul	108874	Napthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
2		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
3		Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	97
4		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
5		Benzene, 2,4-dichloro-1-methyl-	160	C7H6Cl2	000095-73-8	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\
Data File : BM019319.D
Acq On : 22 Mar 2019 19:08
Operator : JU/SJ
Sample : K1968-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09A9

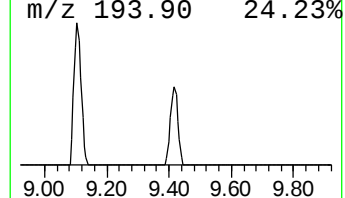
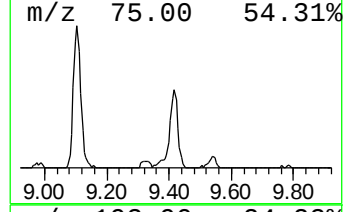
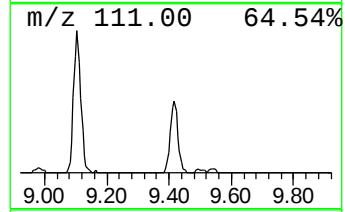
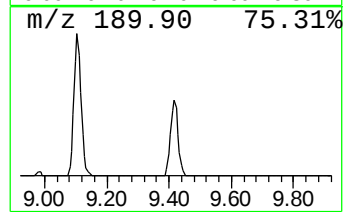
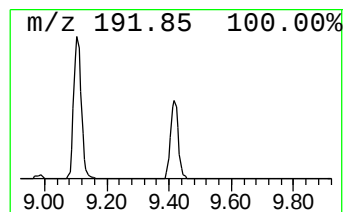
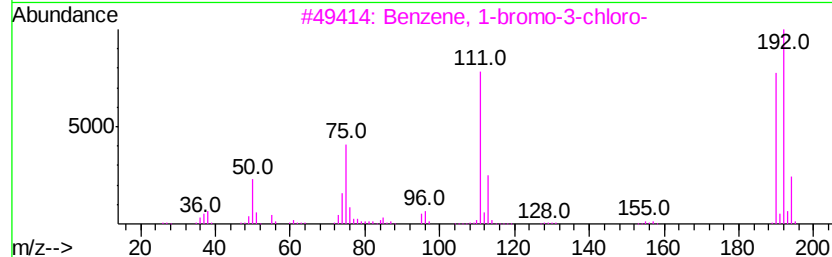
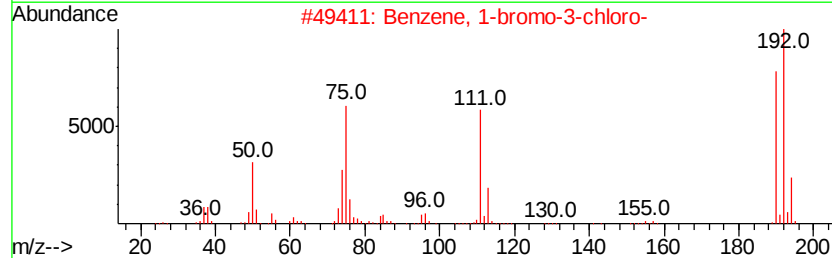
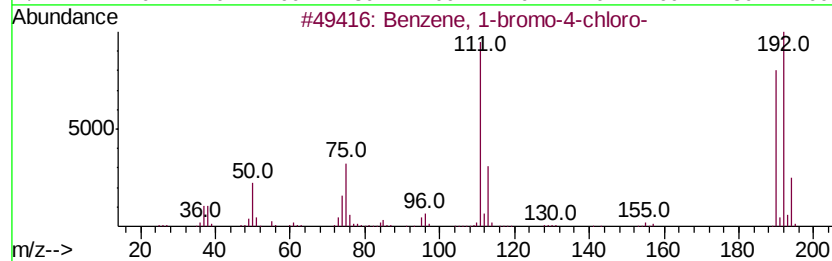
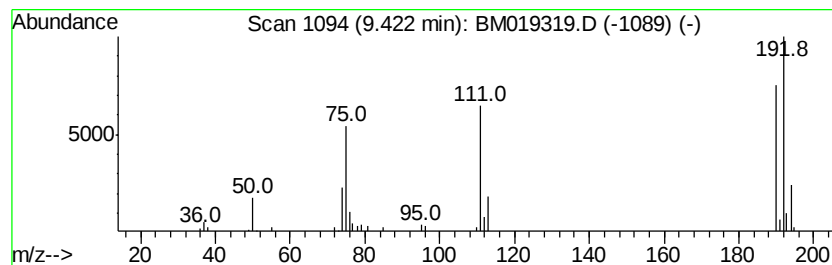
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-bromo-4-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.42	2.69 ng/ul	145636	Napthalene-d8	10.40

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\
Data File : BM019319.D
Acq On : 22 Mar 2019 19:08
Operator : JU/SJ
Sample : K1968-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

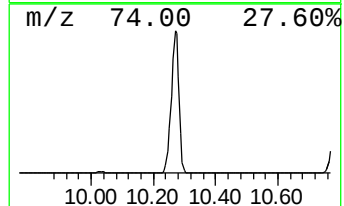
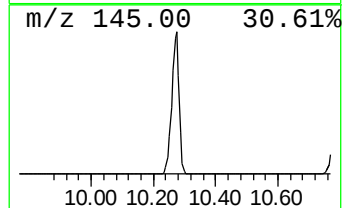
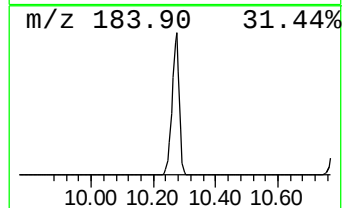
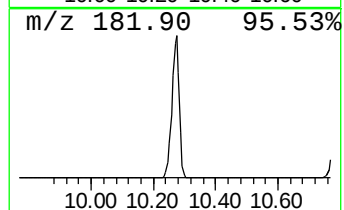
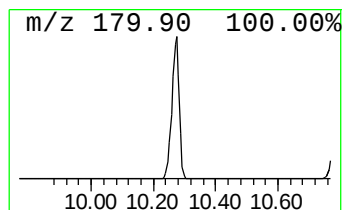
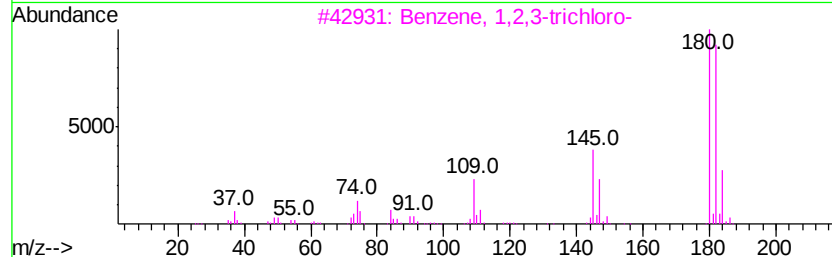
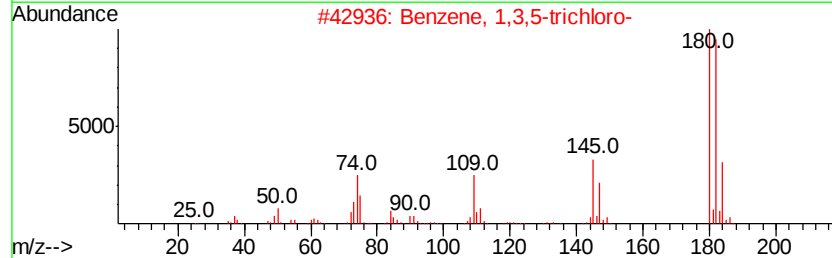
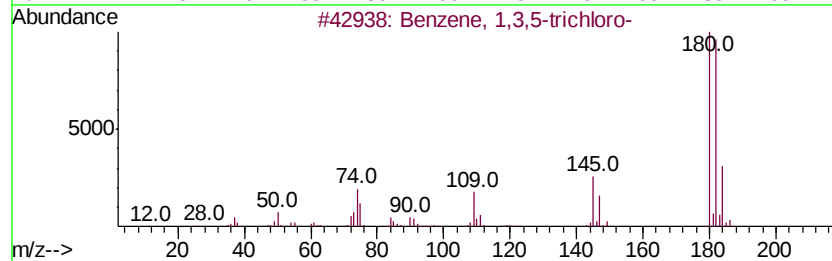
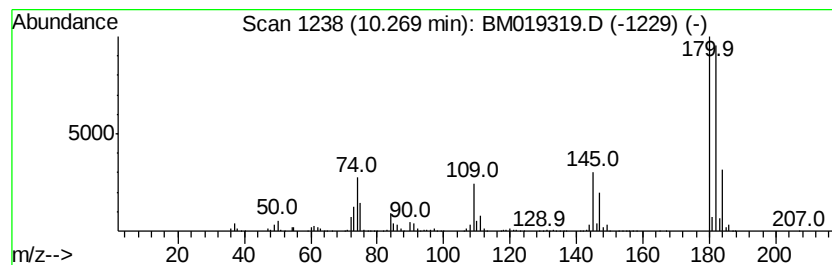
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	617.63 ng/ul	33486800	Naphthalene-d8	10.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 98
2		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 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, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 97



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\
Data File : BM019319.D
Acq On : 22 Mar 2019 19:08
Operator : JU/SJ
Sample : K1968-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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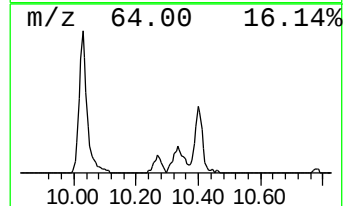
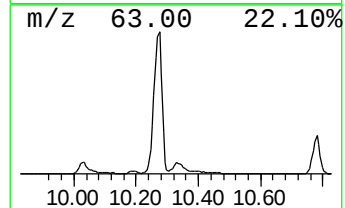
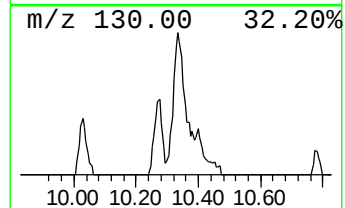
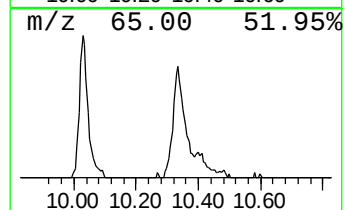
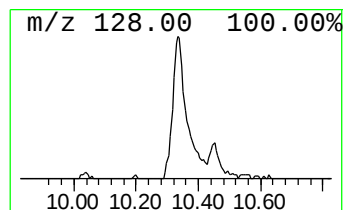
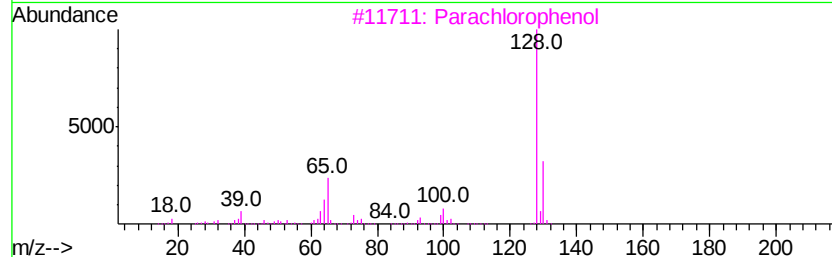
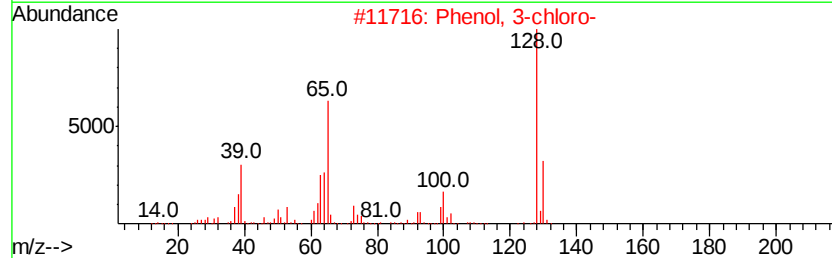
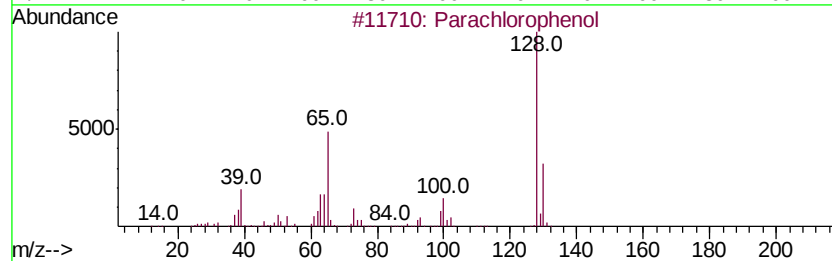
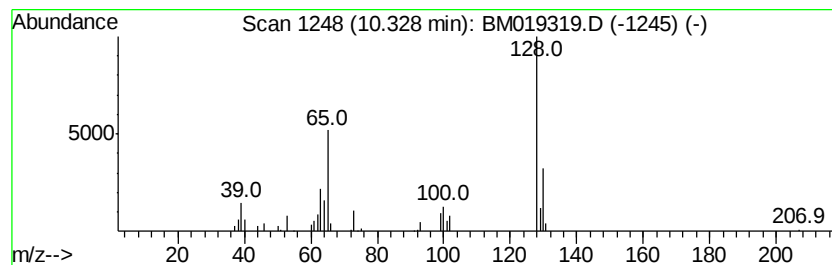
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Parachlorophenol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	3.00 ng/ul	162560	Naphthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Parachlorophenol	128	C6H5ClO	000106-48-9	94
2		Phenol, 3-chloro-	128	C6H5ClO	000108-43-0	94
3		Parachlorophenol	128	C6H5ClO	000106-48-9	94
4		Phenol, 3-chloro-	128	C6H5ClO	000108-43-0	87
5		p-Chlorophenyl N-(cyanomethyl)ca...	210	C9H7ClN2O2	1000240-68-8	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032219\
Data File : BM019319.D
Acq On : 22 Mar 2019 19:08
Operator : JU/SJ
Sample : K1968-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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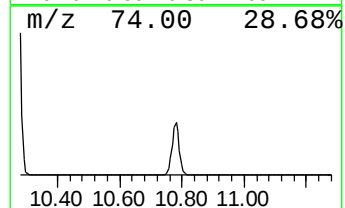
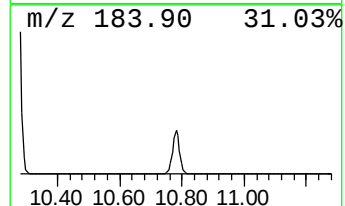
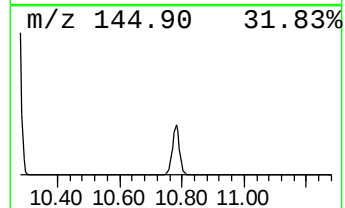
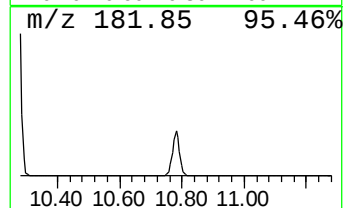
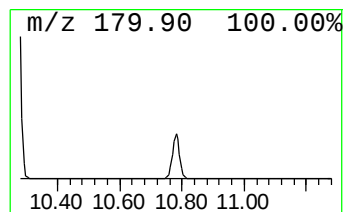
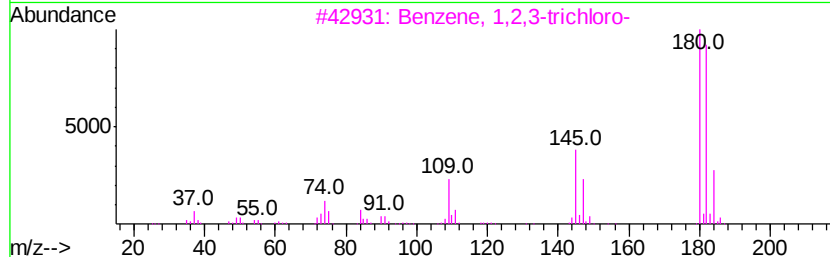
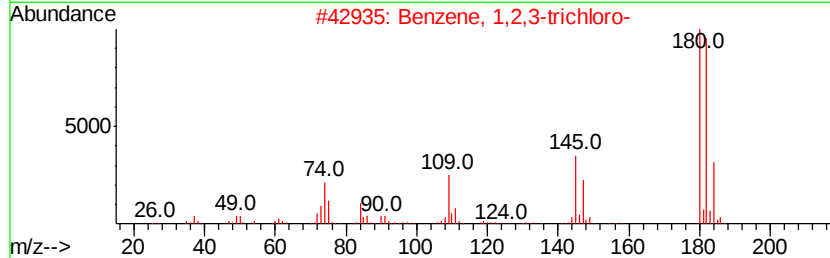
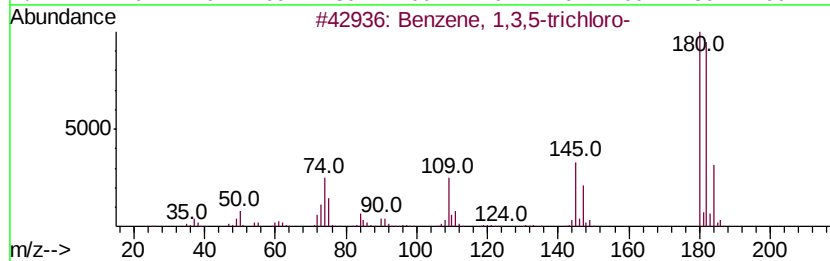
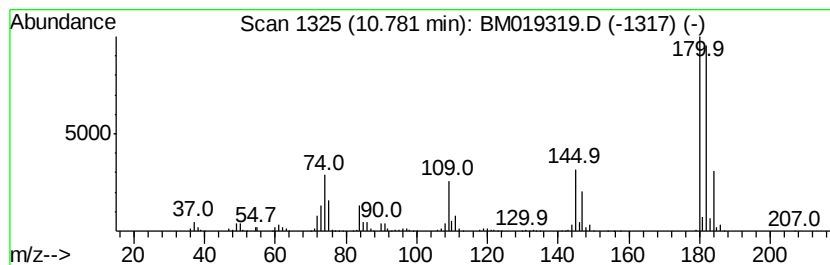
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	143.87 ng/ul	7800590	Naphthalene-d8	10.40

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, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	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 M\DATA\BM032219\
Data File : BM019319.D
Acq On : 22 Mar 2019 19:08
Operator : JU/SJ
Sample : K1968-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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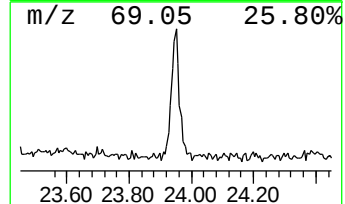
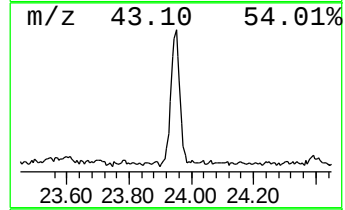
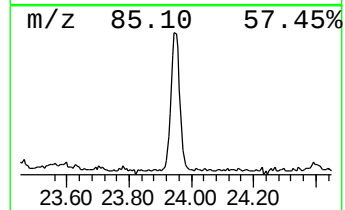
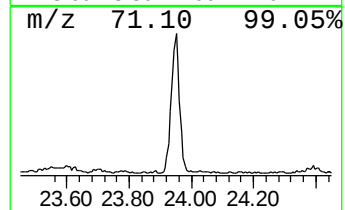
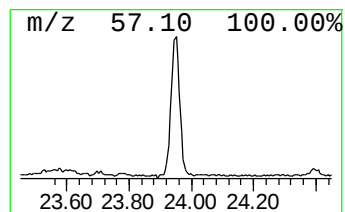
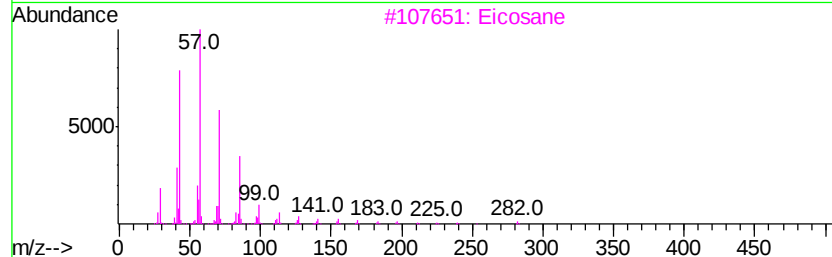
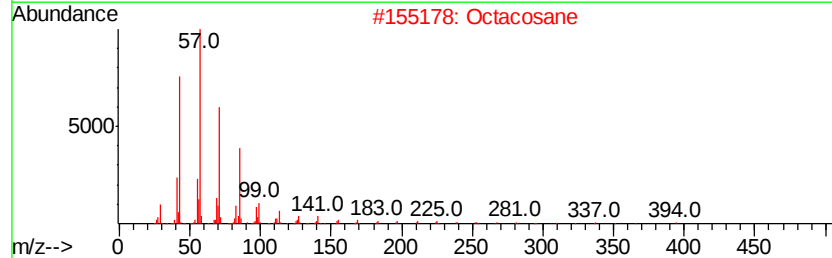
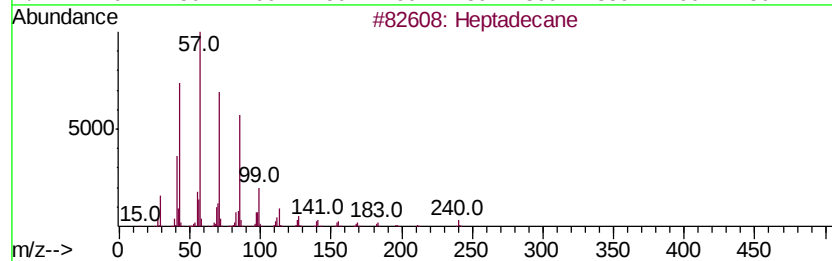
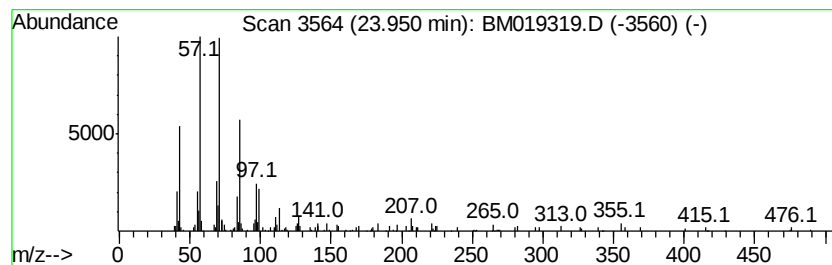
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.95	2.12 ng/ul	227095	Perylene-d12	23.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	72
2		Octacosane	394	C28H58	000630-02-4	72
3		Eicosane	282	C20H42	000112-95-8	70
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	64
5		Heptacosane	380	C27H56	000593-49-7	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032219\
Data File : BM019319.D
Acq On : 22 Mar 2019 19:08
Operator : JU/SJ
Sample : K1968-14
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09A9

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-	4.97	40.4	ng/ul	54018600	1	7.62	26755900	20.0
Benzene, 1,4-dich...	7.50	4.1	ng/ul	5508390	1	7.62	26755900	20.0
Benzene, 1,3-dich...	7.97	20.4	ng/ul	27303500	1	7.62	26755900	20.0
Benzene, 1-bromo-...	9.10	4.7	ng/ul	252563	2	10.40	1084370	20.0
Benzene, 1,2-dich...	9.37	2.0	ng/ul	108874	2	10.40	1084370	20.0
Benzene, 1-bromo-...	9.42	2.7	ng/ul	145636	2	10.40	1084370	20.0
Benzene, 1,3,5-tr...	10.27	617.6	ng/ul	33486800	2	10.40	1084370	20.0
Parachlorophenol	10.33	3.0	ng/ul	162560	2	10.40	1084370	20.0
Benzene, 1,2,3-tr...	10.78	143.9	ng/ul	7800590	2	10.40	1084370	20.0
(DEL) Alkane: Str...	23.95	2.1	ng/ul	227095	6	23.46	2144660	20.0