

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
 Data File : BM019425.D
 Acq On : 25 Mar 2019 11:21
 Operator : JU/SJ
 Sample : K2027-10DL2 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0989DL2

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	4.951	327	334	352	rBV	12448723	17466926	29.50%	11.729%
2	6.963	671	676	691	rBV	33967	73531	0.12%	0.049%
3	7.145	702	707	717	rVB	45910	76905	0.13%	0.052%
4	7.492	759	766	779	rBV	1739640	2751328	4.65%	1.848%
5	7.651	779	793	820	rVV	22293912	36343658	61.38%	24.405%
6	7.969	834	847	853	rBV	31635495	59207200	100.00%	39.759%
7	8.775	979	984	986	rBV	52385	76491	0.13%	0.051%
8	8.816	986	991	1006	rVB	406898	778504	1.31%	0.523%
9	10.033	1191	1198	1211	rBV2	34748	85740	0.14%	0.058%
10	10.251	1227	1235	1252	rBV	9753337	15763103	26.62%	10.585%
11	10.386	1252	1258	1267	rVB	703960	1125357	1.90%	0.756%
12	10.769	1315	1323	1339	rBV	2194108	3718176	6.28%	2.497%
13	11.010	1355	1364	1384	rBV	295000	629278	1.06%	0.423%
14	11.227	1396	1401	1417	rBV	66575	159292	0.27%	0.107%
15	12.433	1596	1606	1617	rBV	211189	393639	0.66%	0.264%
16	13.045	1703	1710	1725	rBV	920771	1481218	2.50%	0.995%
17	13.686	1814	1819	1828	rBV	83547	152145	0.26%	0.102%
18	13.957	1857	1865	1873	rBV	130877	200202	0.34%	0.134%
19	14.263	1909	1917	1930	rBV2	951841	1437296	2.43%	0.965%
20	15.268	2081	2088	2101	rBV	162754	274885	0.46%	0.185%
21	17.021	2380	2386	2398	rBV2	1075915	1608328	2.72%	1.080%
22	17.127	2398	2404	2416	rVB	150256	269246	0.45%	0.181%
23	19.427	2790	2795	2805	rBV	215367	329031	0.56%	0.221%
24	21.221	3095	3100	3110	rBV	1334928	1901361	3.21%	1.277%
25	22.244	3271	3274	3279	rVB2	69281	75387	0.13%	0.051%
26	22.738	3355	3358	3362	rVB2	84524	104119	0.18%	0.070%
27	23.297	3448	3453	3460	rVB2	266273	481415	0.81%	0.323%
28	23.438	3471	3477	3491	rVB	1015898	1952783	3.30%	1.311%

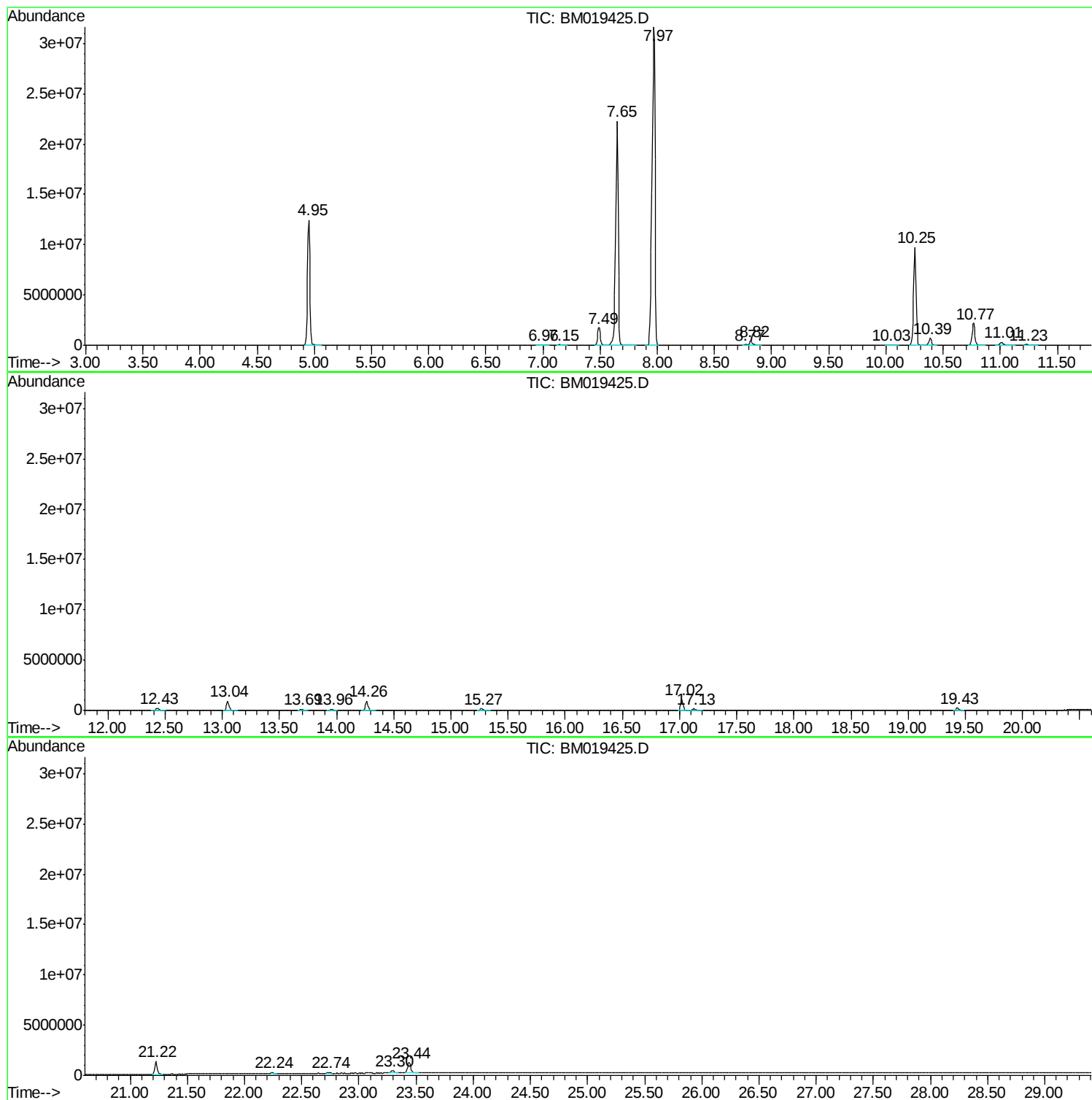
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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



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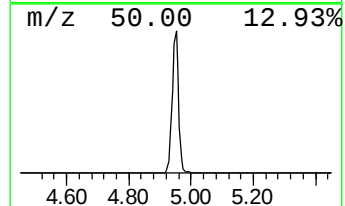
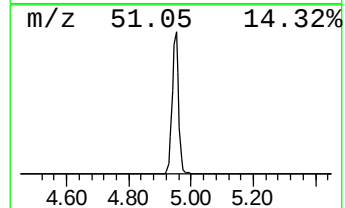
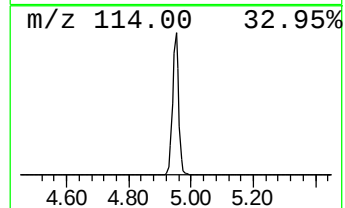
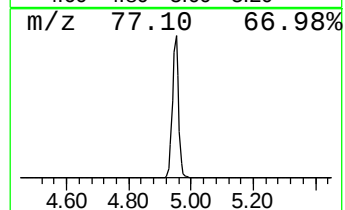
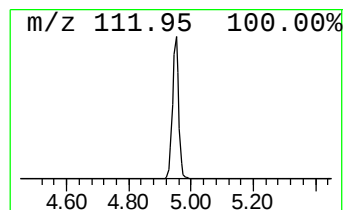
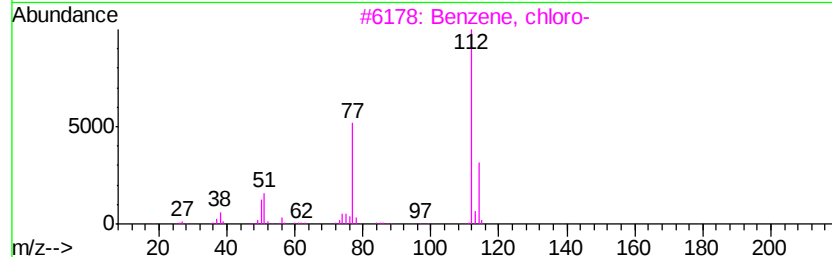
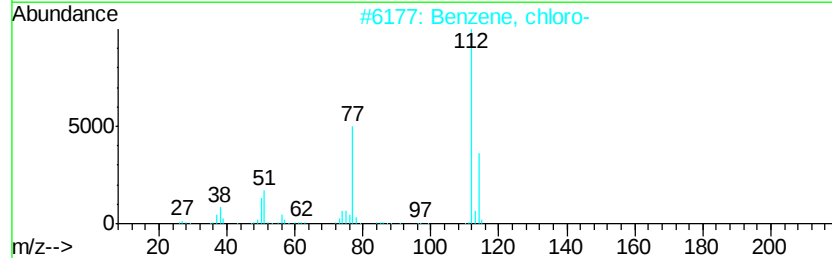
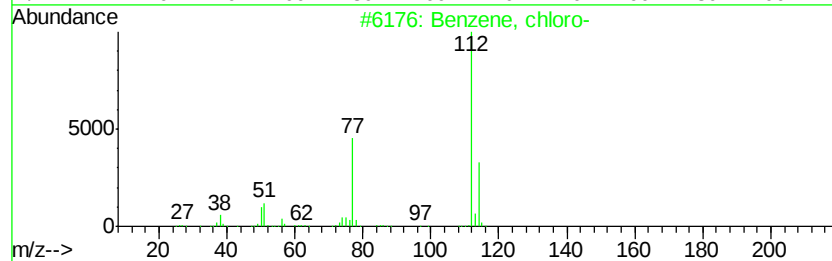
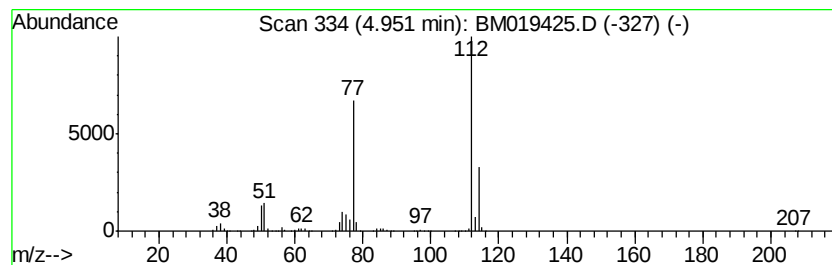
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	9.61 ng/ul	17466900	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	95
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	95
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5		Phenol, 4-fluoro-	112	C6H5FO	000371-41-5	38



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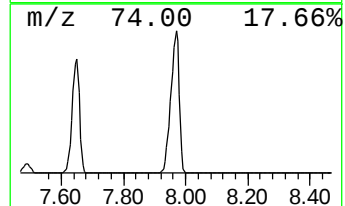
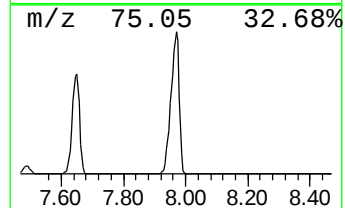
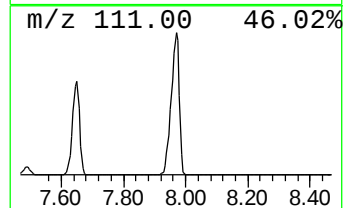
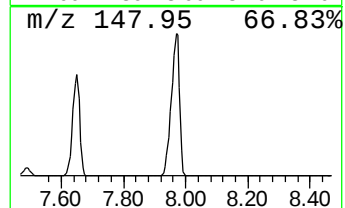
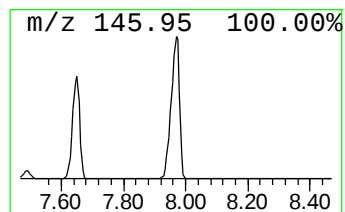
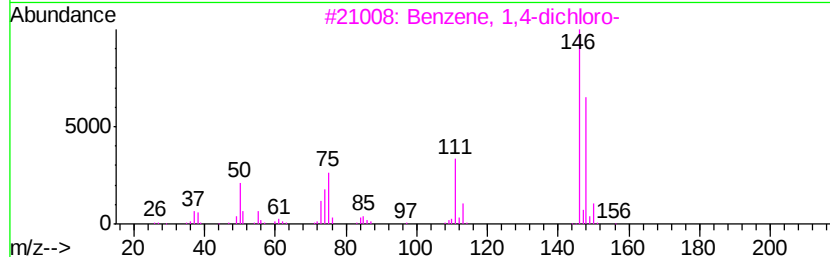
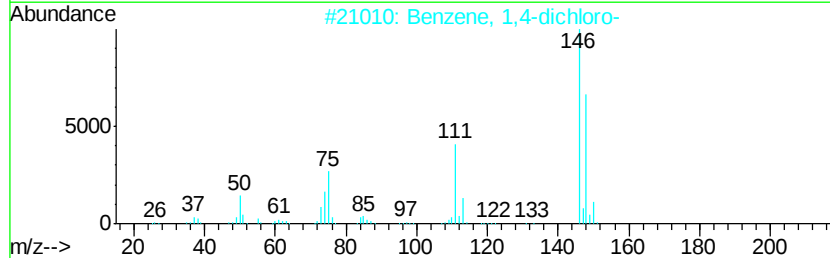
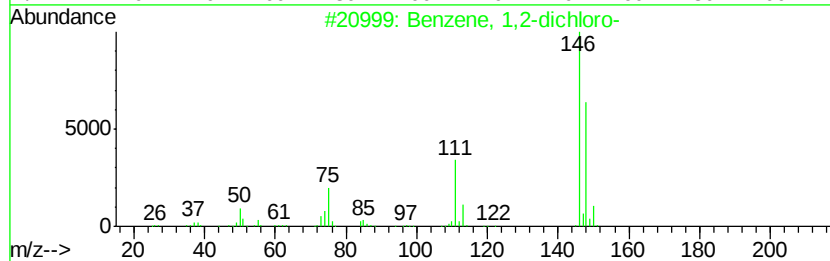
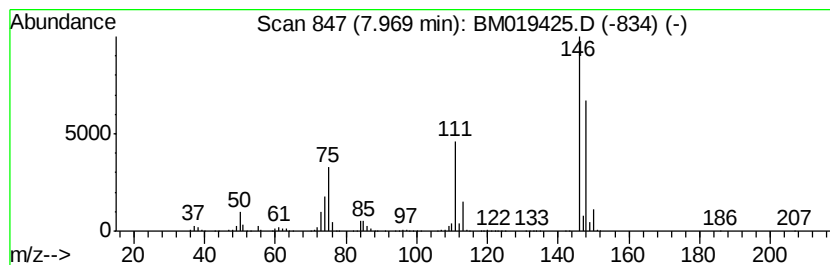
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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	32.58 ng/ul	59207200	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95



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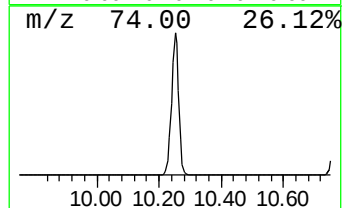
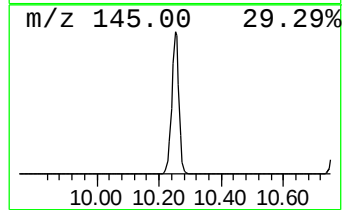
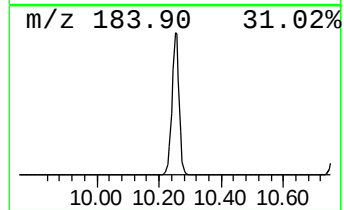
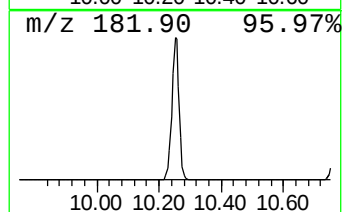
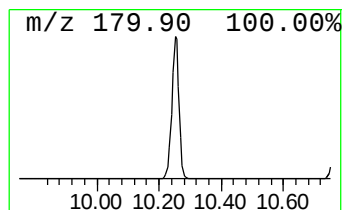
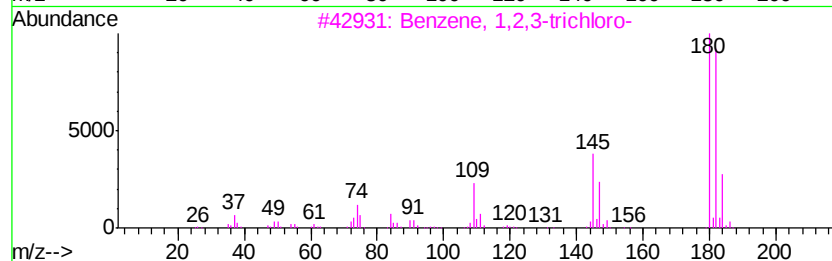
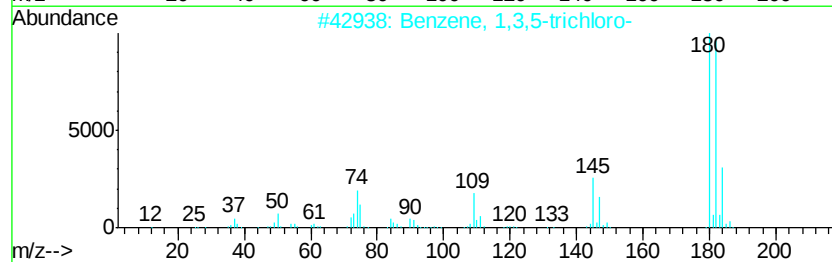
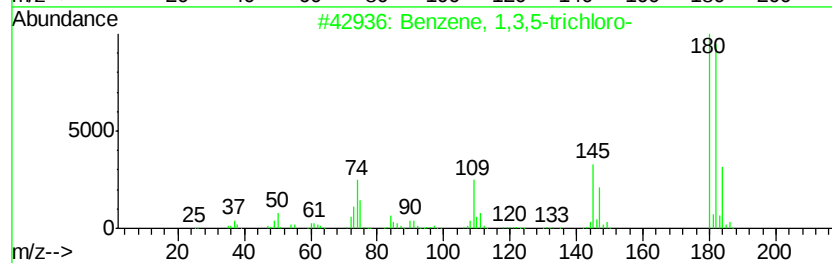
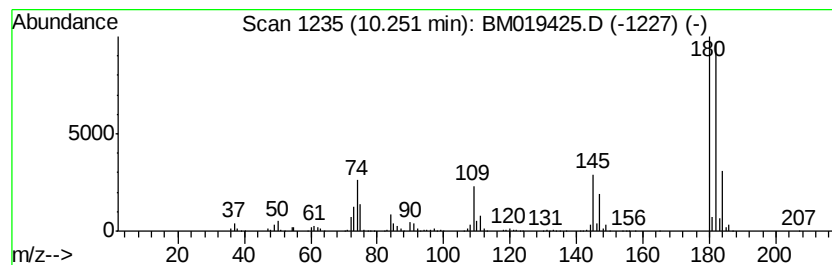
Instrument :
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	280.14 ng/ul	15763100	Napthalene-d8	10.39
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



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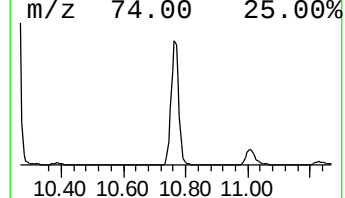
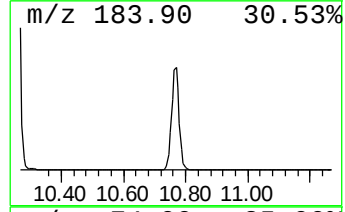
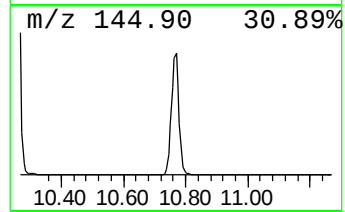
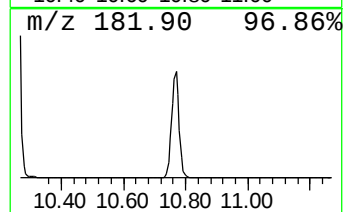
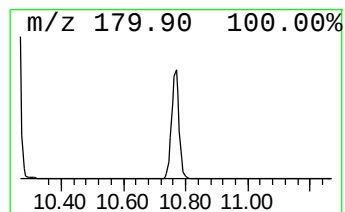
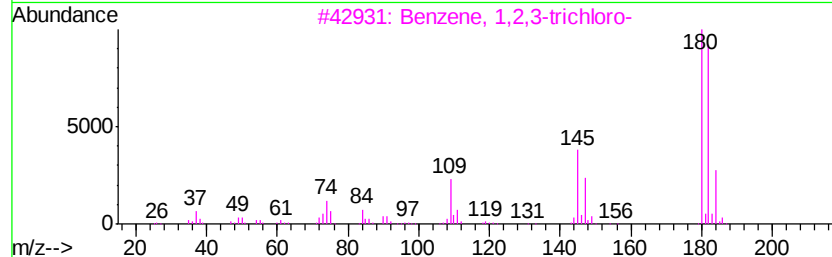
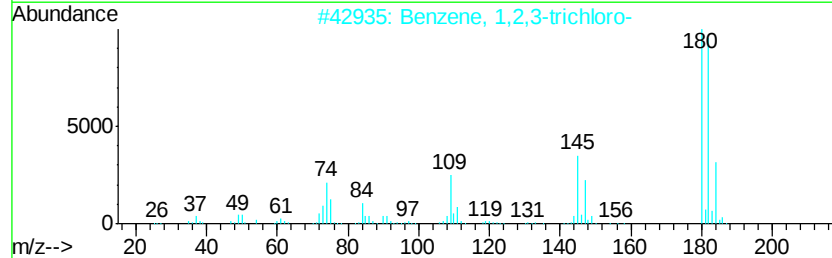
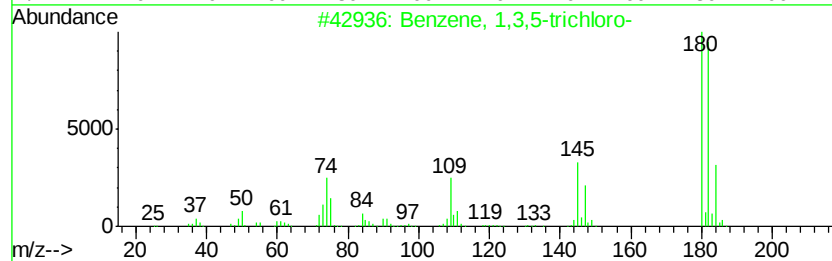
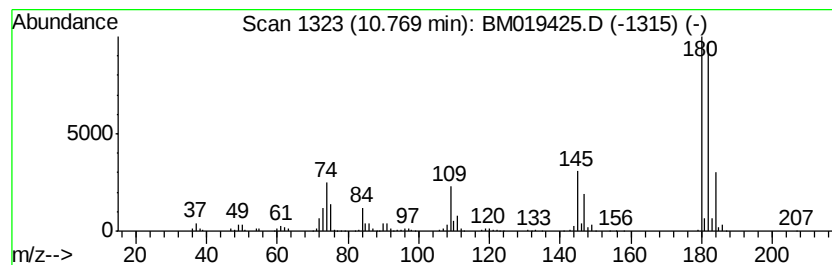
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Peak Number 4 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	66.08 ng/ul	3718180	Naphthalene-d8	10.39

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, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97



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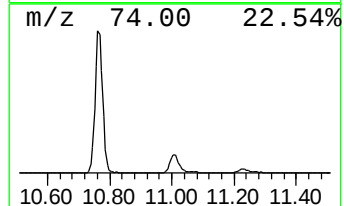
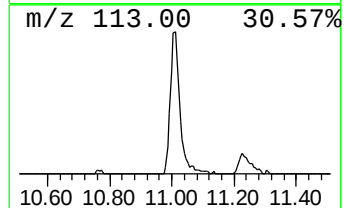
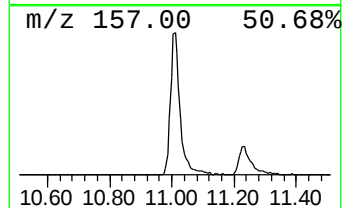
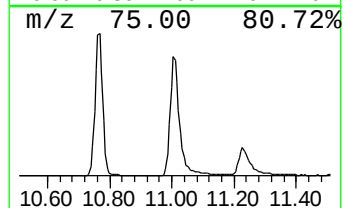
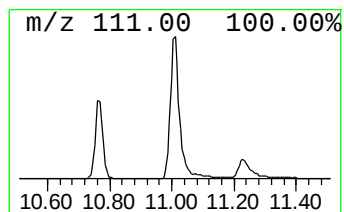
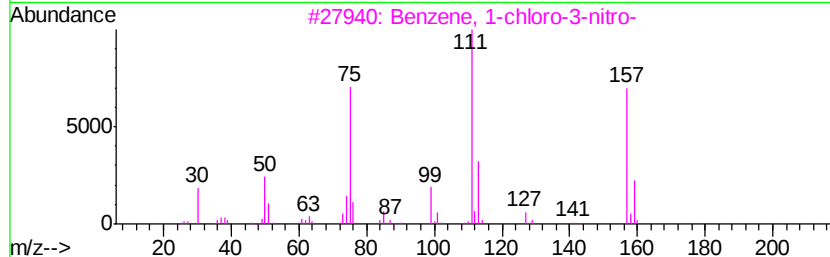
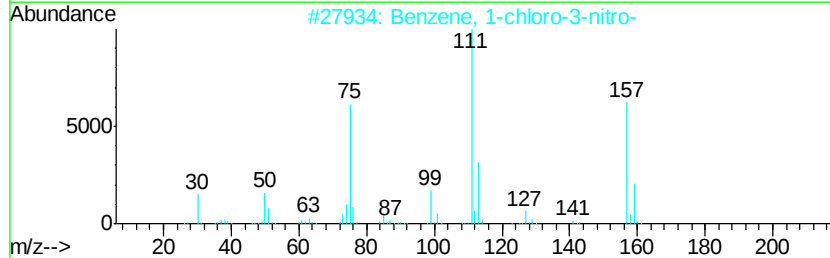
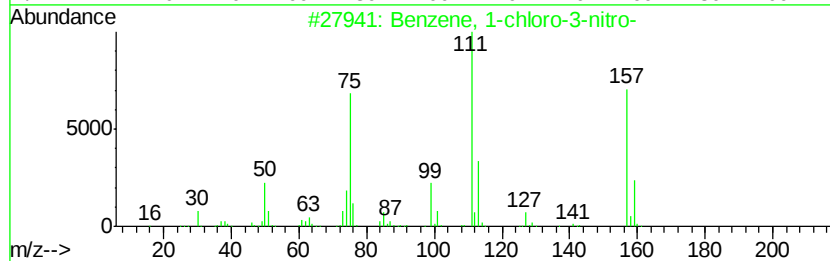
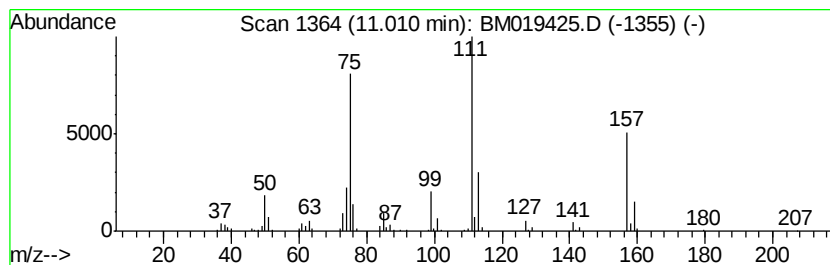
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Peak Number 5 Benzene, 1-chloro-3-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.01	11.18 ng/ul	629278	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	94
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	87



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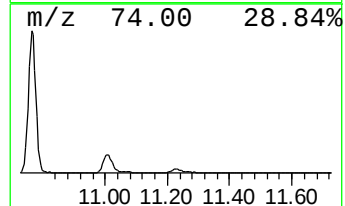
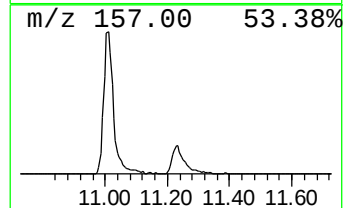
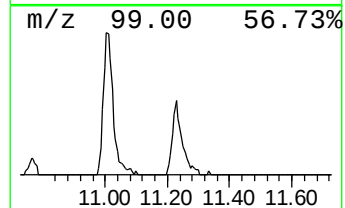
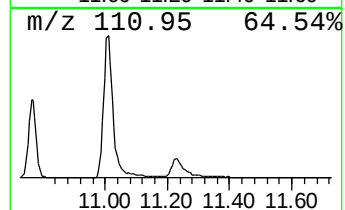
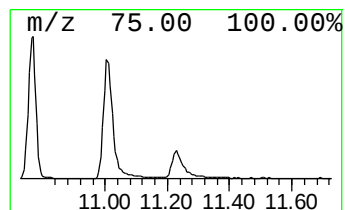
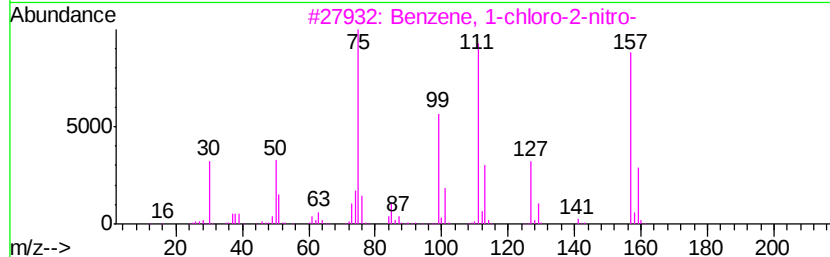
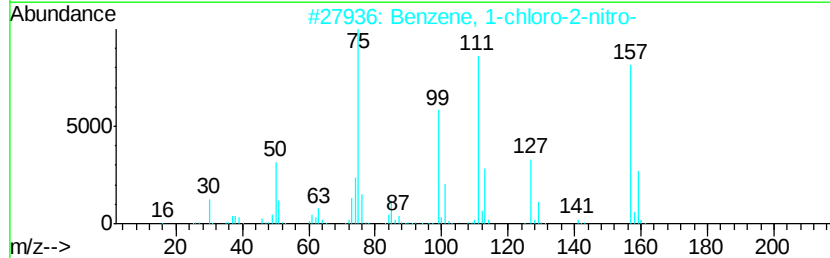
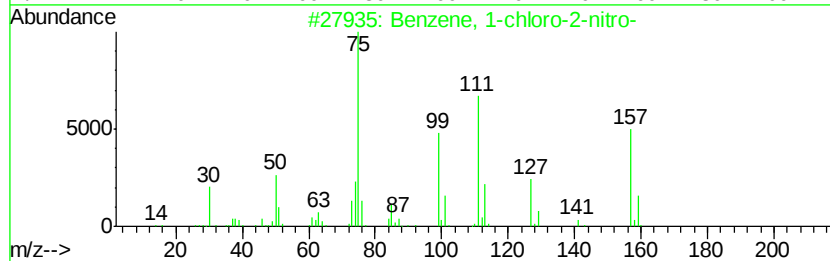
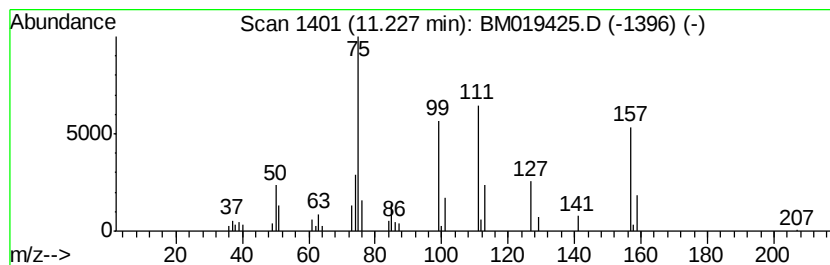
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-chloro-2-nitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	2.83 ng/ul	159292	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	93
5		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032519\
Data File : BM019425.D
Acq On : 25 Mar 2019 11:21
Operator : JU/SJ
Sample : K2027-10DL2 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0989DL2

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-	4.95	9.6	ng/ul	17466900	1	7.60	36343700	20.0
Benzene, 1,2-dich...	7.97	32.6	ng/ul	59207200	1	7.60	36343700	20.0
Benzene, 1,3,5-tr...	10.25	280.1	ng/ul	15763100	2	10.39	1125360	20.0
Benzene, 1,2,3-tr...	10.77	66.1	ng/ul	3718180	2	10.39	1125360	20.0
Benzene, 1-chloro...	11.01	11.2	ng/ul	629278	2	10.39	1125360	20.0
Benzene, 1-chloro...	11.23	2.8	ng/ul	159292	2	10.39	1125360	20.0