

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP012721\
 Data File : BP004647.D
 Acq On : 27 Jan 2021 12:46
 Operator : CG/JU
 Sample : M1173-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0FY6

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_P\METHODS\SOM-EPA-BP012721MA.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	5.169	319	337	342	rBV3	43733430	143985471	100.00%	29.784%
2	7.134	664	671	682	rBV	143614	242839	0.17%	0.050%
3	7.322	696	703	717	rBV	183850	327765	0.23%	0.068%
4	7.693	755	766	777	rBV	8138992	15540246	10.79%	3.215%
5	7.881	778	798	802	rBV2	39640987	111899515	77.72%	23.147%
6	8.216	835	855	859	rBV2	39746574	132195837	91.81%	27.345%
7	8.492	896	902	909	rBV	141053	218569	0.15%	0.045%
8	8.945	972	979	983	rBV	217090	389288	0.27%	0.081%
9	8.992	983	987	994	rVB	677771	1036256	0.72%	0.214%
10	9.281	1028	1036	1044	rBV	153960	257336	0.18%	0.053%
11	9.598	1085	1090	1095	rVB	86041	149920	0.10%	0.031%
12	9.669	1095	1102	1109	rBV	203282	345991	0.24%	0.072%
13	10.204	1185	1193	1202	rBV	295056	491442	0.34%	0.102%
14	10.481	1226	1240	1245	rBV	24642825	49257239	34.21%	10.189%
15	10.598	1253	1260	1265	rBV	194060	313560	0.22%	0.065%
16	10.981	1315	1325	1337	rVB	7046541	11618673	8.07%	2.403%
17	11.181	1352	1359	1367	rBV	448371	724604	0.50%	0.150%
18	11.398	1386	1396	1404	rBV	109889	195427	0.14%	0.040%
19	12.616	1598	1603	1611	rVB	538974	843830	0.59%	0.175%
20	13.228	1695	1707	1717	rBV	1719433	2563295	1.78%	0.530%
21	13.845	1799	1812	1817	rBV	579310	837014	0.58%	0.173%
22	14.128	1852	1860	1868	rVB	598220	870236	0.60%	0.180%
23	14.433	1906	1912	1922	rVB	331263	484861	0.34%	0.100%
24	15.433	2072	2082	2092	rBV	776974	1133844	0.79%	0.235%
25	15.539	2092	2100	2109	rBV	289006	386200	0.27%	0.080%
26	17.186	2373	2380	2387	rBV	442001	617586	0.43%	0.128%
27	17.286	2390	2397	2404	rBV	925426	1290883	0.90%	0.267%
28	18.539	2601	2610	2617	rBV	320096	426875	0.30%	0.088%
29	19.527	2772	2778	2784	rBV	1250657	1653309	1.15%	0.342%
30	21.268	3070	3074	3079	rBV	648123	765596	0.53%	0.158%
31	23.439	3436	3443	3451	rBV	856102	1613624	1.12%	0.334%
32	23.586	3462	3468	3478	rVB2	388235	762939	0.53%	0.158%

LSC Area Percent Report

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ClientSampleId :
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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\SOM-EPA-BP012721MA.M
Title : SVOA CALIBRATION

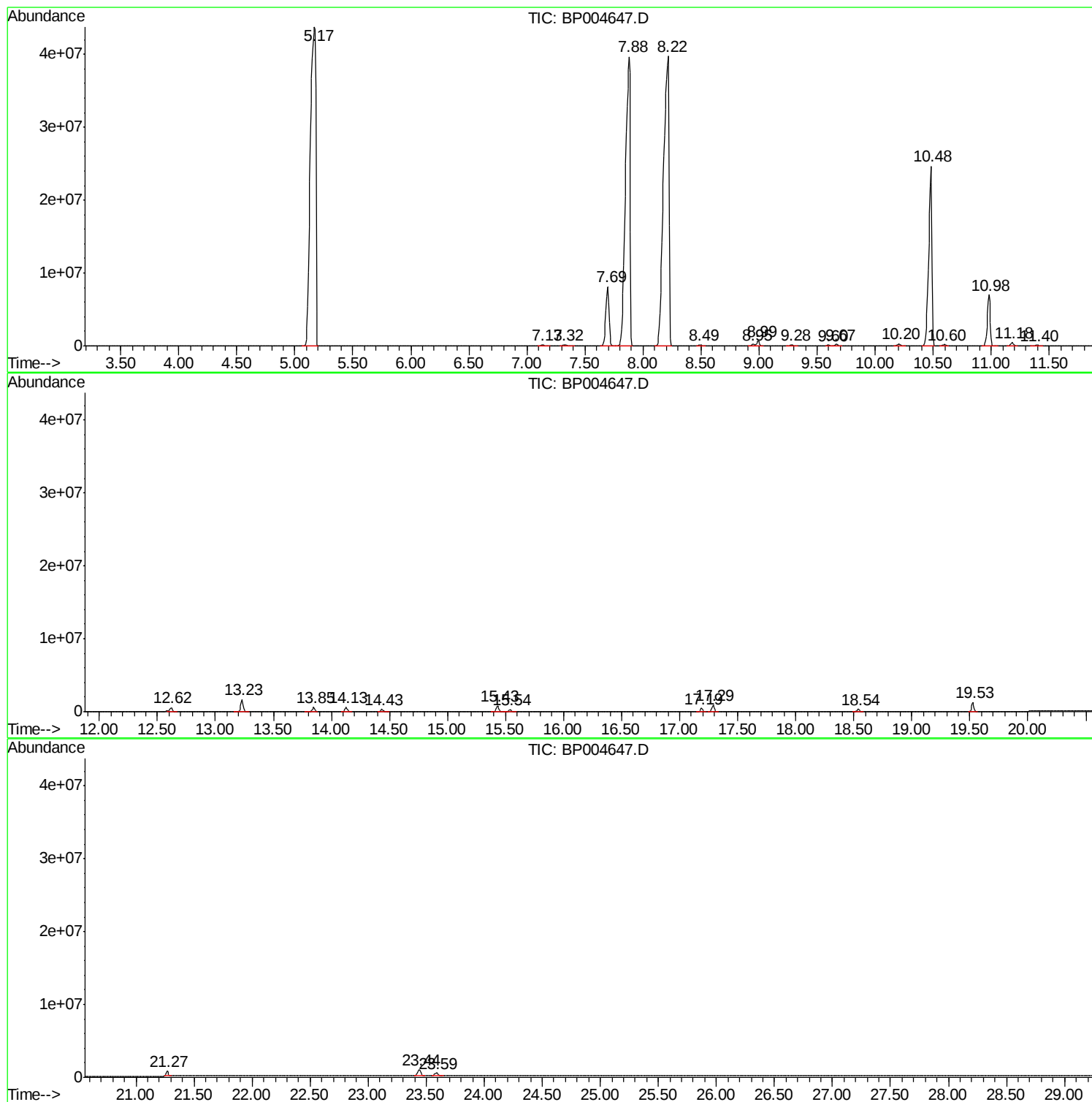
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP012721MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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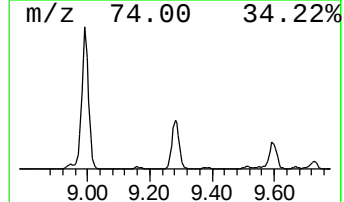
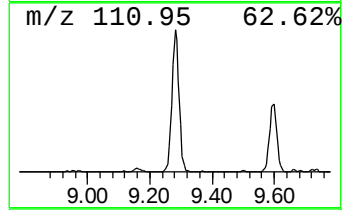
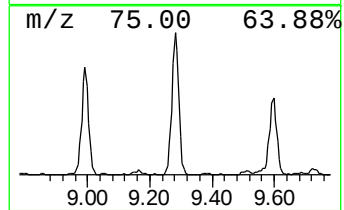
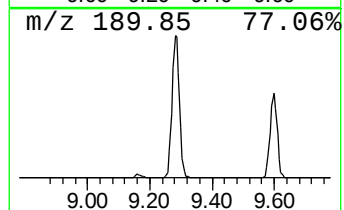
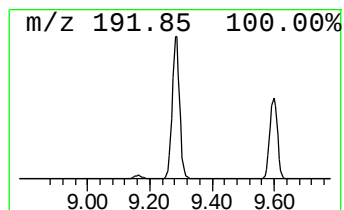
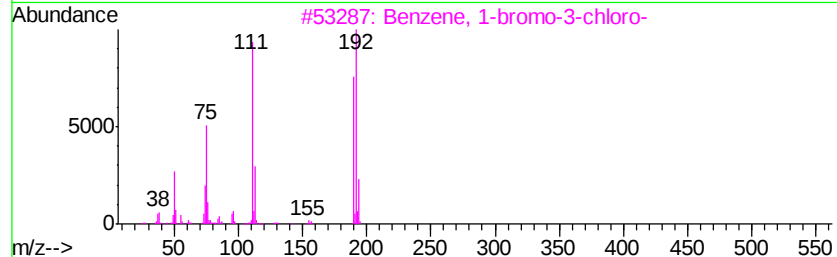
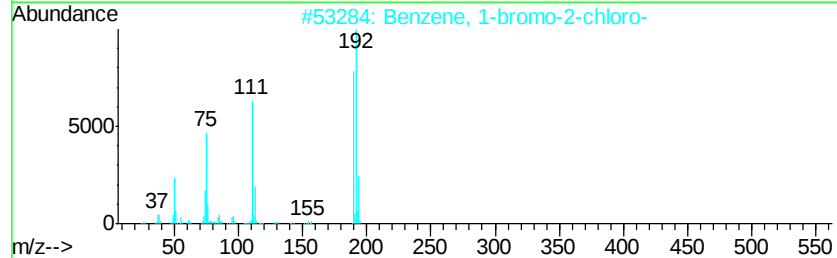
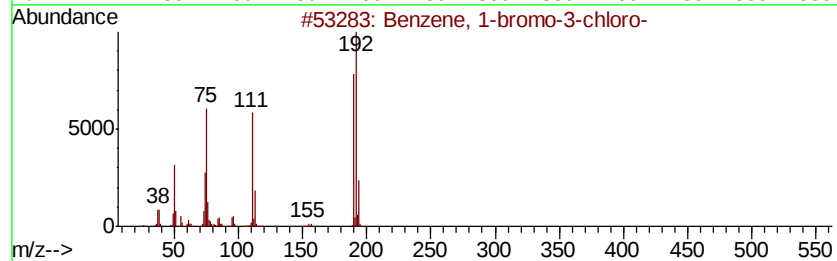
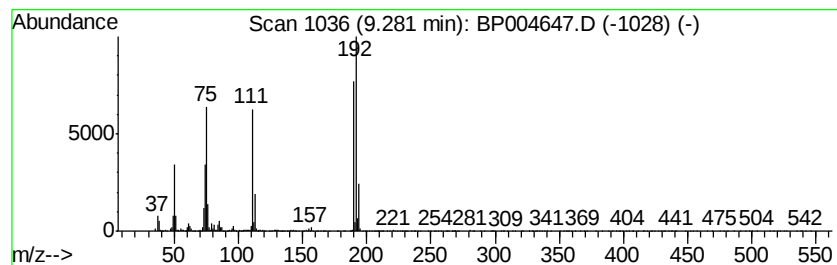
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-bromo-3-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	16.41 ng/ul	257336	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95
2		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	94
4		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94
5		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94



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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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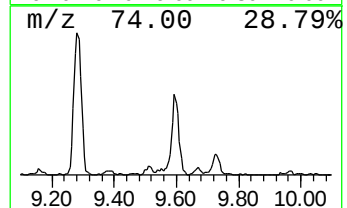
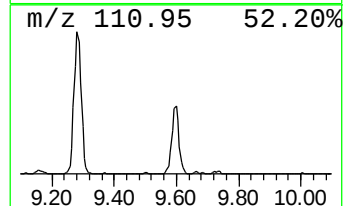
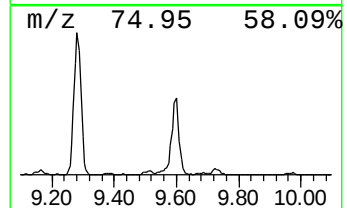
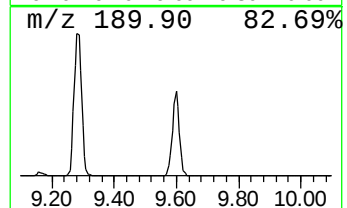
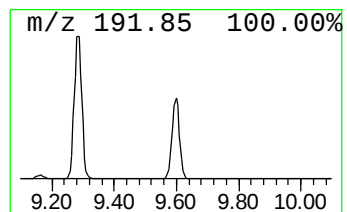
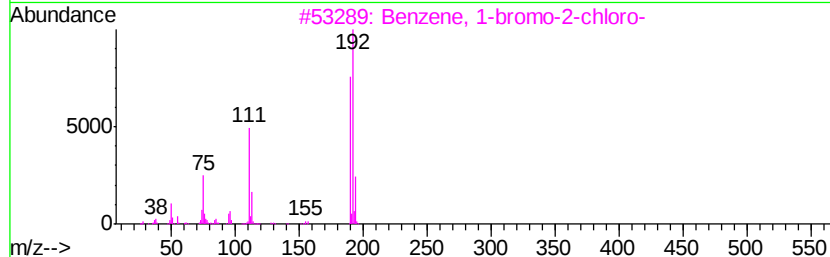
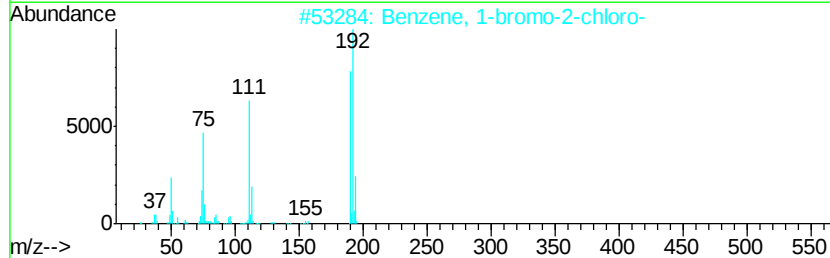
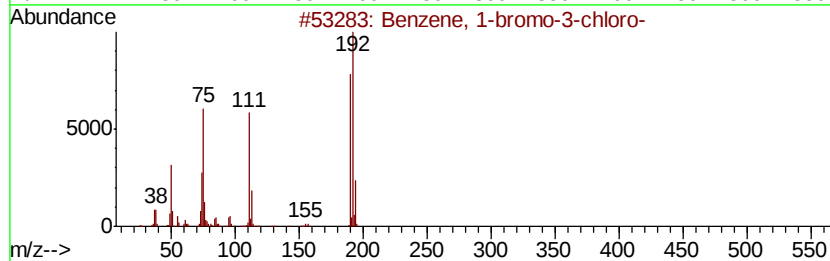
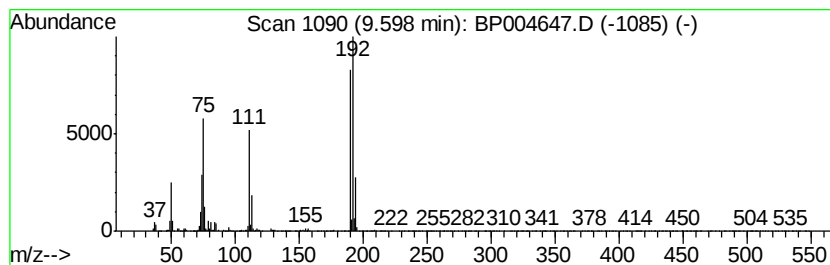
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	9.56 ng/ul	149920	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95
2		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	93
3		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	93
4		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	93
5		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	93



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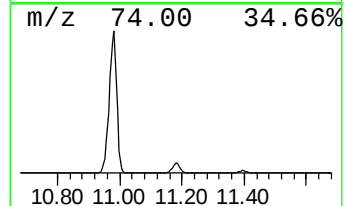
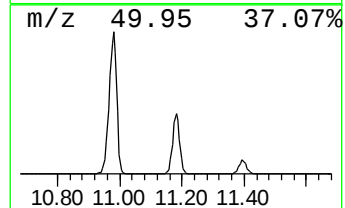
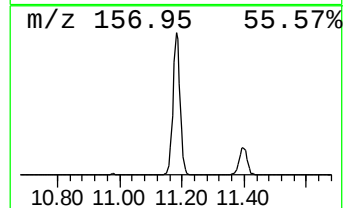
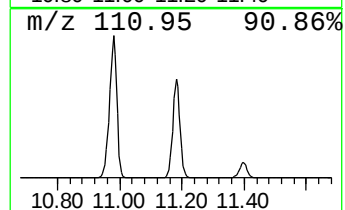
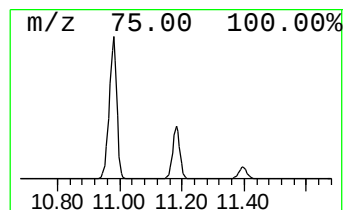
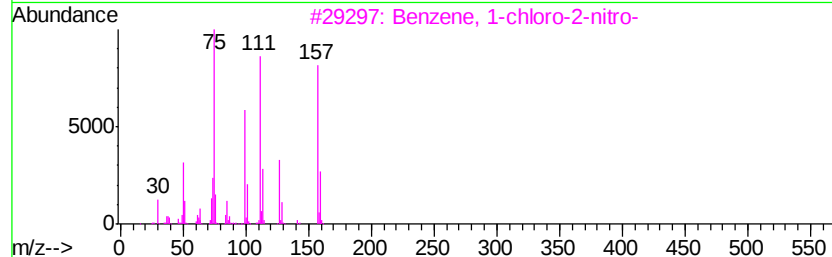
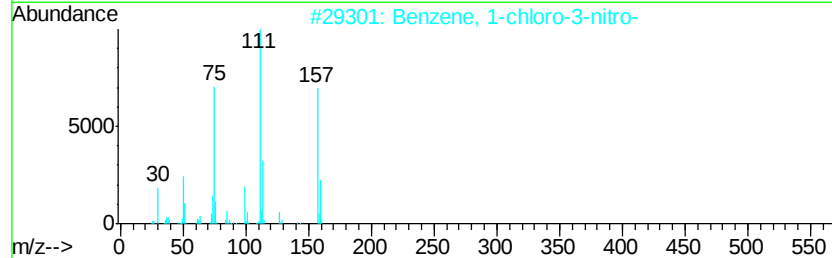
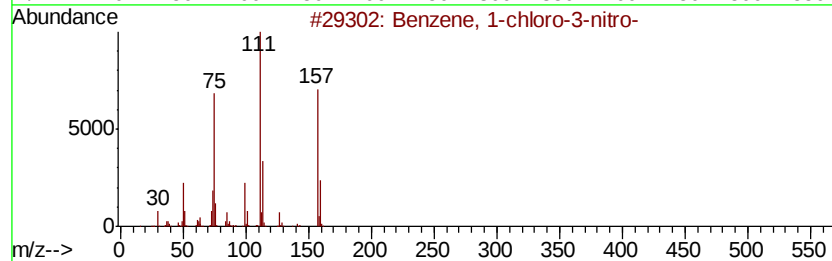
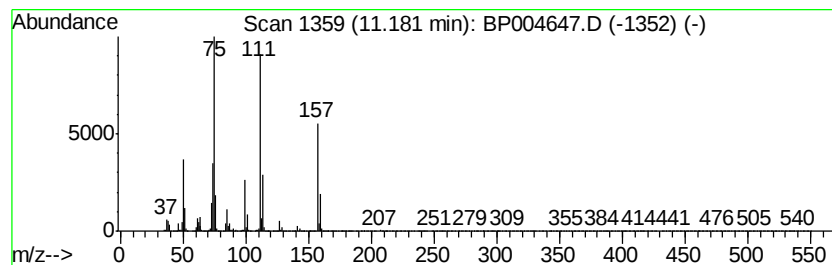
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.18	46.22 ng/ul	724604	Napthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	93
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	93
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91



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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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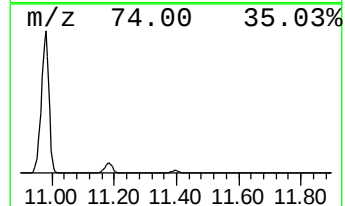
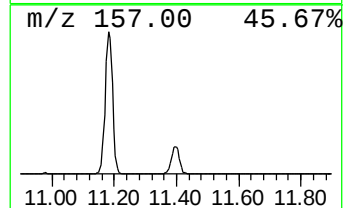
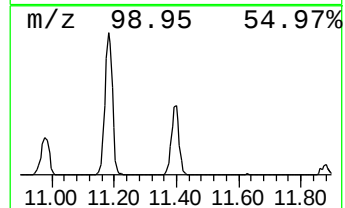
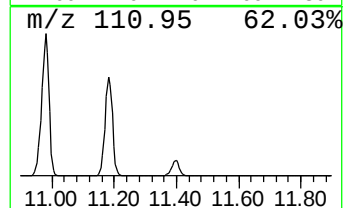
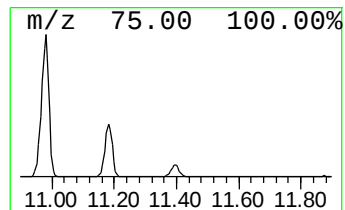
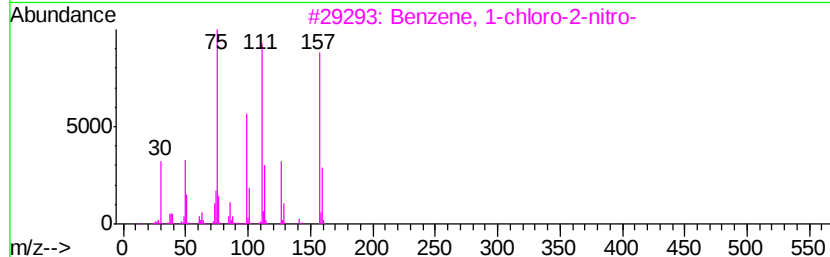
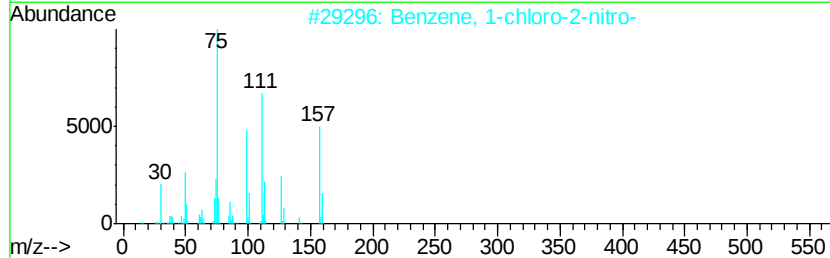
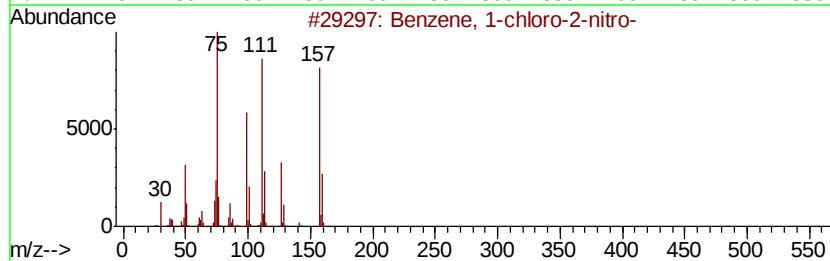
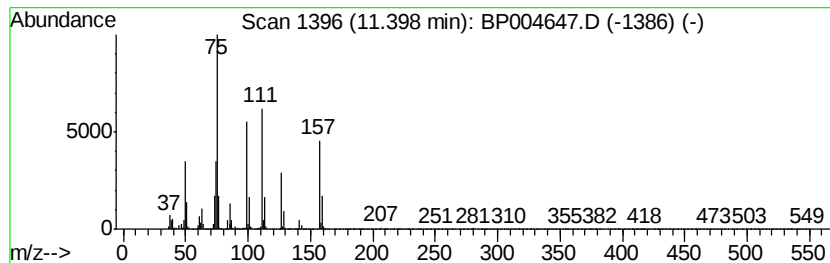
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP012721MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	12.47 ng/ul	195427	Naphthalene-d8	10.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	94
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	93
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	72
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	62
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	62



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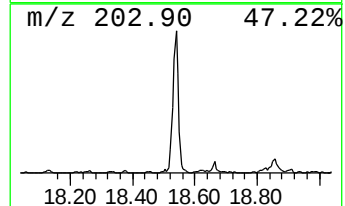
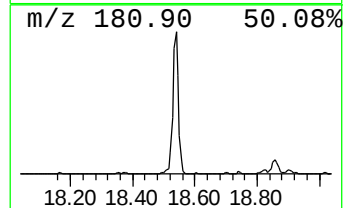
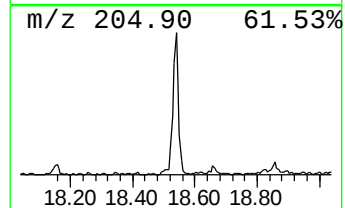
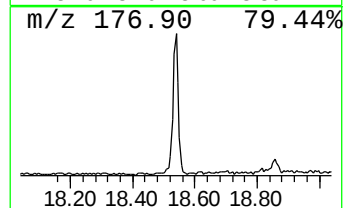
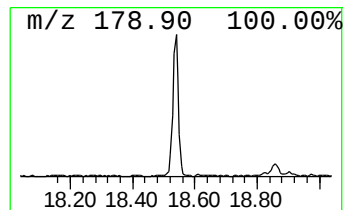
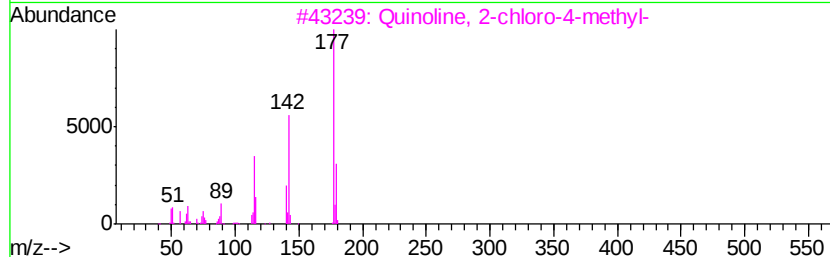
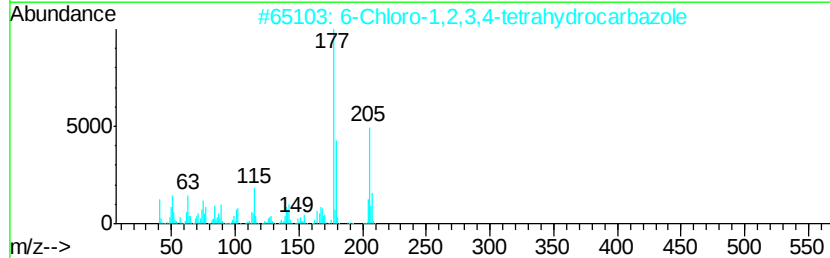
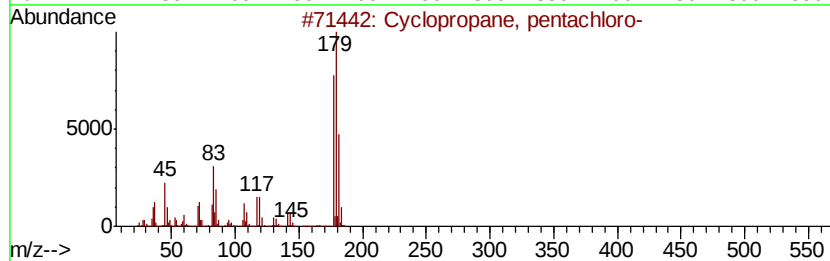
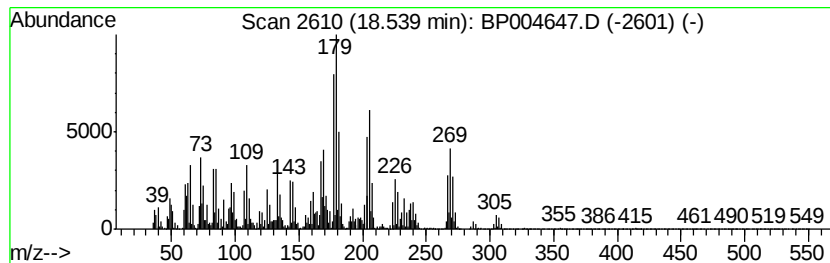
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.54	13.82 ng/ul	426875	Phenanthrene-d10	17.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	35
2		6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	20
3		Quinoline, 2-chloro-4-methyl-	177	C10H8ClN	000634-47-9	18
4		7-Chloro-4-hydroxyquinoline	179	C9H6ClNO	000086-99-7	11
5		Pyrido[3,2-d]pyrimidine-2,4(1H,3...	177	C8H7N3O2	002499-96-9	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP012721\
Data File : BP004647.D
Acq On : 27 Jan 2021 12:46
Operator : CG/JU
Sample : M1173-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0FY6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP012721MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-bromo-...	9.28	16.4	ng/ul	257336	2	10.60	313560	20.0
Benzene, 1-bromo-...	9.60	9.6	ng/ul	149920	2	10.60	313560	20.0
Benzene, 1-chloro...	11.18	46.2	ng/ul	724604	2	10.60	313560	20.0
Benzene, 1-chloro...	11.40	12.5	ng/ul	195427	2	10.60	313560	20.0
unknown-01	18.54	13.8	ng/ul	426875	4	17.19	617586	20.0