

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010053.D
 Acq On : 22 Apr 2022 13:30
 Operator : CG/JU
 Sample : N2473-09DL 5X
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0HZ1DL

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: BP010053.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.281	360	373	378	rBV2	42228354	106477889	100.00%	25.709%
2	6.617	595	600	607	rVB	67765	111480	0.10%	0.027%
3	7.269	705	711	723	rVB	154569	253116	0.24%	0.061%
4	7.475	738	746	755	rBV	147021	254075	0.24%	0.061%
5	7.840	797	808	820	rBV	9177805	16326007	15.33%	3.942%
6	8.011	820	837	842	rBV2	38206698	98910122	92.89%	23.882%
7	8.340	877	893	898	rBV	36173196	96225519	90.37%	23.234%
8	8.646	939	945	956	rVB	125281	209118	0.20%	0.050%
9	9.099	1014	1022	1024	rBV	201261	315099	0.30%	0.076%
10	9.140	1024	1029	1041	rVB	1275507	2181462	2.05%	0.527%
11	9.440	1072	1080	1091	rBV	248854	424816	0.40%	0.103%
12	9.675	1112	1120	1123	rBV	68838	115507	0.11%	0.028%
13	9.752	1129	1133	1140	rVV	143846	273460	0.26%	0.066%
14	9.822	1140	1145	1153	rVV	155462	305523	0.29%	0.074%
15	10.369	1230	1238	1247	rBV	221101	403289	0.38%	0.097%
16	10.634	1271	1283	1296	rBV	24665663	49728331	46.70%	12.007%
17	10.752	1296	1303	1311	rVB	1077480	1742405	1.64%	0.421%
18	10.875	1316	1324	1337	rBV	151596	278631	0.26%	0.067%
19	11.140	1360	1369	1386	rBV	9018379	15185364	14.26%	3.667%
20	11.340	1396	1403	1416	rBV	845740	1453043	1.36%	0.351%
21	11.557	1432	1440	1449	rBV	188170	332585	0.31%	0.080%
22	12.740	1630	1641	1642	rBV3	104779	214866	0.20%	0.052%
23	12.769	1642	1646	1654	rVB	480480	760051	0.71%	0.184%
24	13.375	1742	1749	1763	rBV	1788914	2669689	2.51%	0.645%
25	13.975	1842	1851	1857	rBV	558825	796584	0.75%	0.192%
26	14.263	1894	1900	1910	rVB	549201	799252	0.75%	0.193%
27	14.569	1940	1952	1966	rBV	1750781	2659567	2.50%	0.642%
28	14.893	2001	2007	2014	rVB	89415	128279	0.12%	0.031%
29	15.563	2114	2121	2134	rBV	718284	1083613	1.02%	0.262%
30	15.675	2135	2140	2150	rVB	144507	224190	0.21%	0.054%
31	16.540	2281	2287	2293	rBV3	92435	137472	0.13%	0.033%
32	17.322	2414	2420	2431	rBV	2243578	3203920	3.01%	0.774%
33	17.422	2431	2437	2449	rVB	782195	1150154	1.08%	0.278%
34	18.663	2639	2648	2655	rBV	663326	856080	0.80%	0.207%
35	19.651	2810	2816	2821	rBV	997568	1341600	1.26%	0.324%
36	21.404	3109	3114	3129	rBV	2396862	3150768	2.96%	0.761%

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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\SFAM-EPA-BP042122.M
Title : SVOA CALIBRATION

37	23.698	3498	3504	3510	rBV2	423353	858666	0.81%	0.207%
38	23.857	3524	3531	3543	rVB2	1180314	2620368	2.46%	0.633%

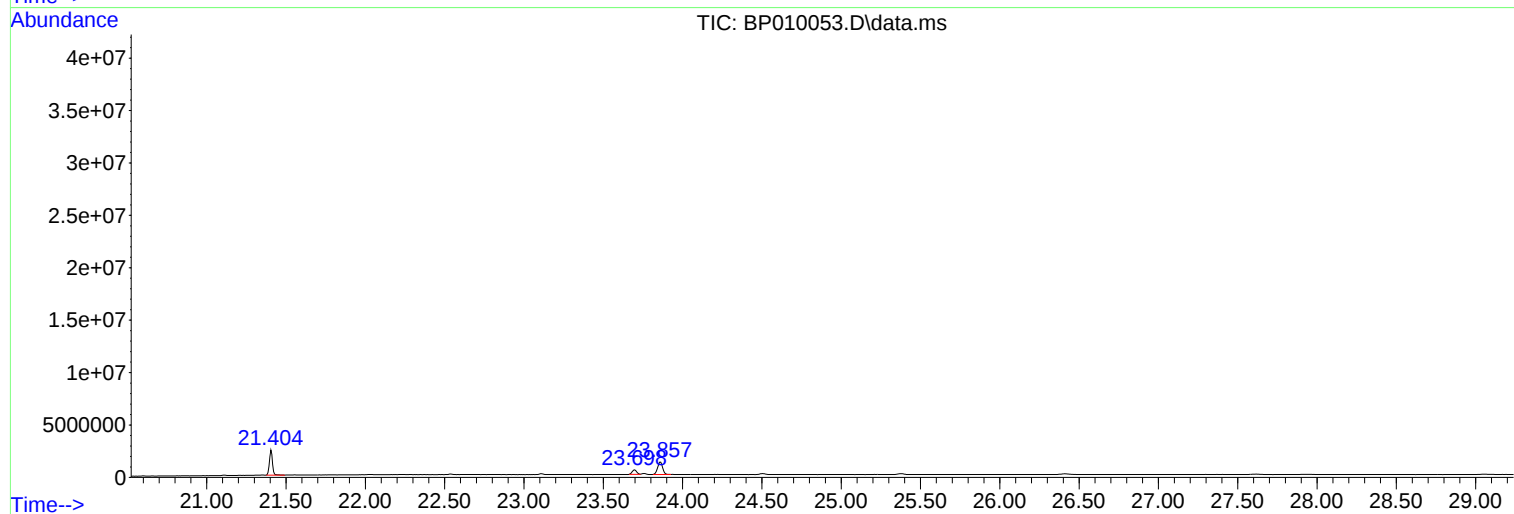
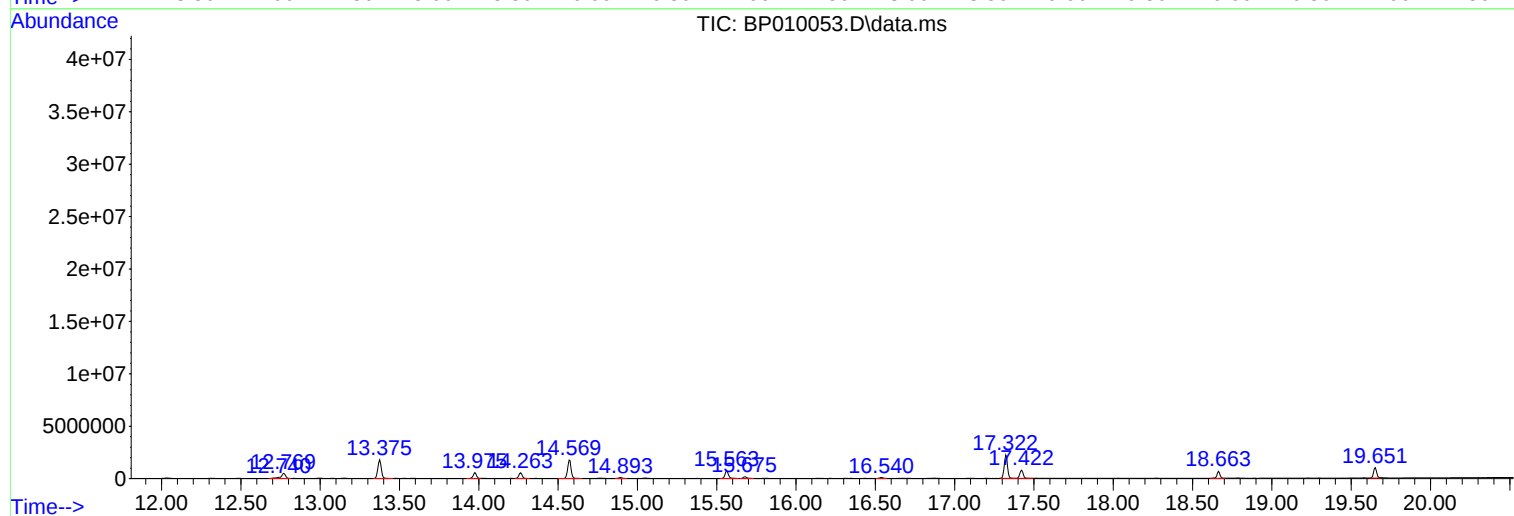
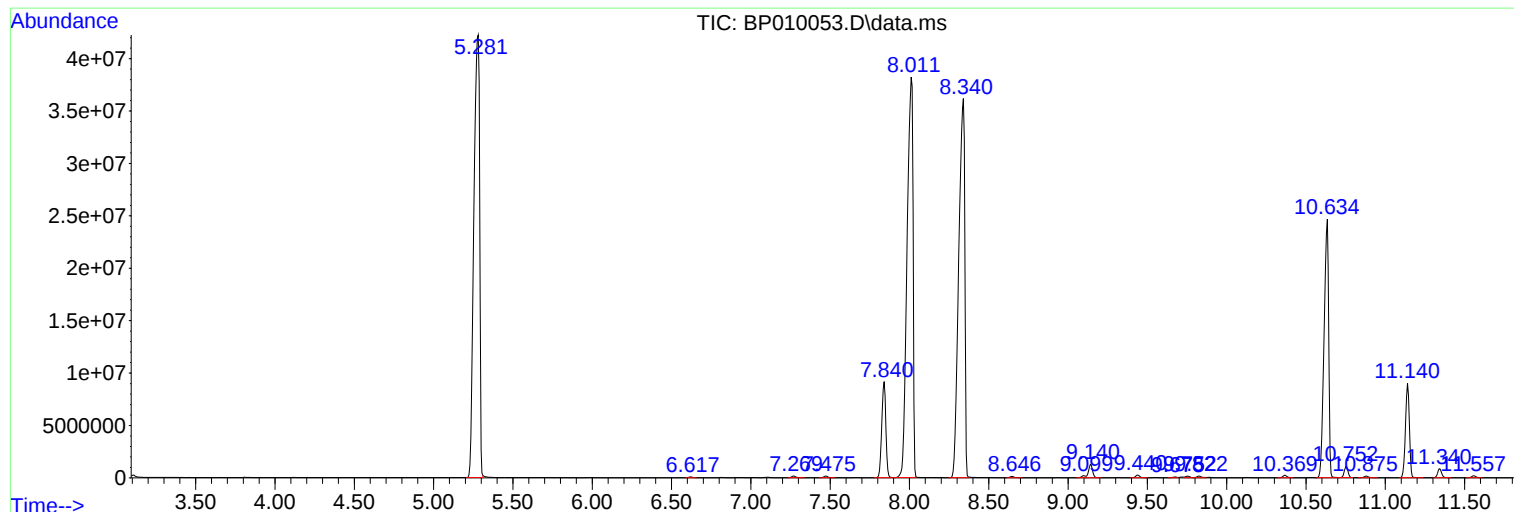
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ClientSampleId :
C0HZ1DL

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



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Misc :
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Instrument :
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ClientSampleId :
C0HZ1DL

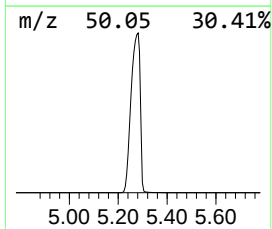
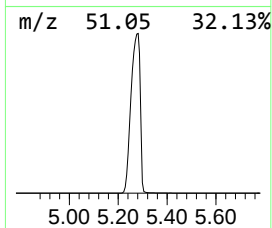
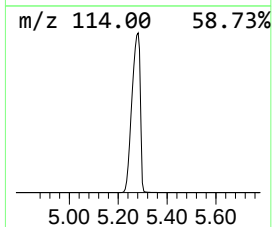
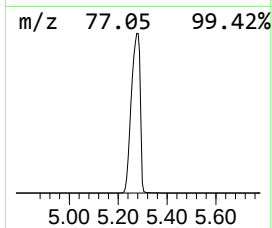
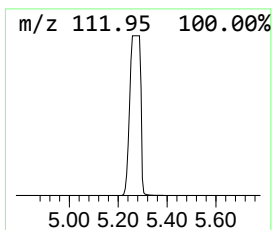
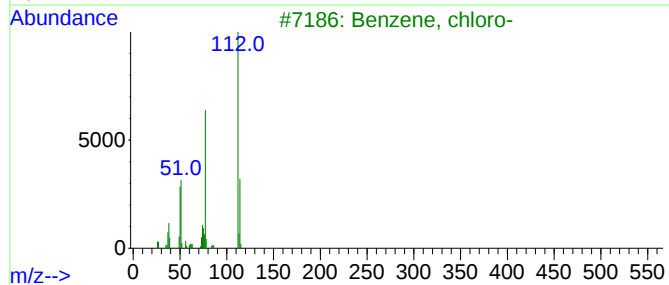
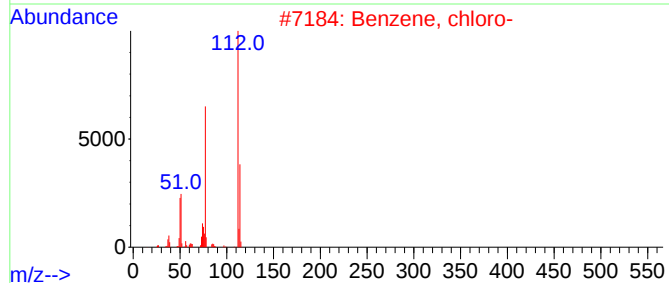
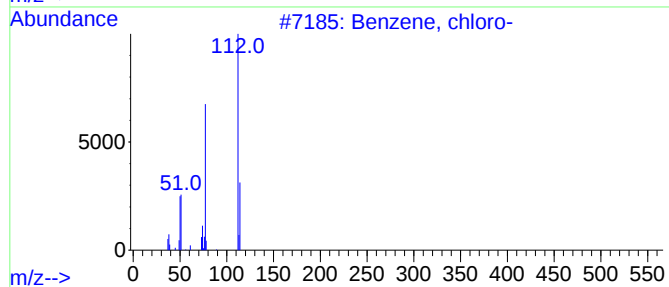
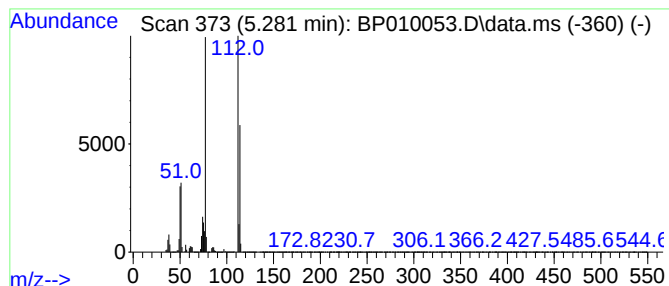
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.281	21.53 ng/ul	106478000	1,4-Dichlorobenzene-d4	7.958

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	93
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	90



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Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HZ1DL

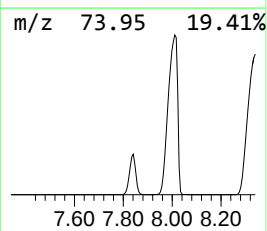
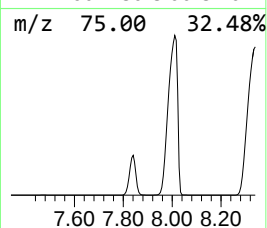
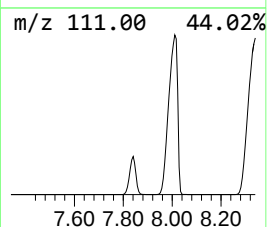
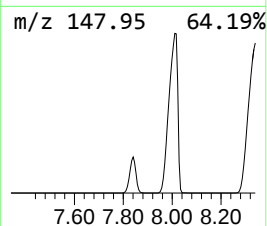
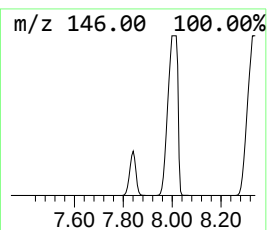
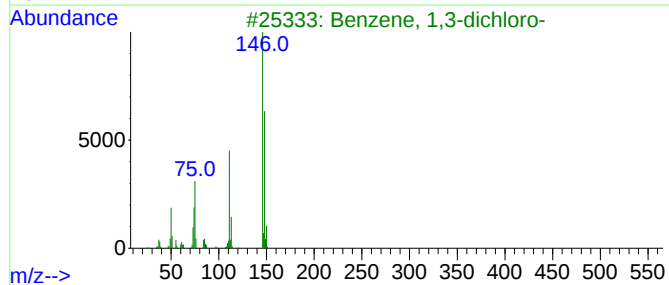
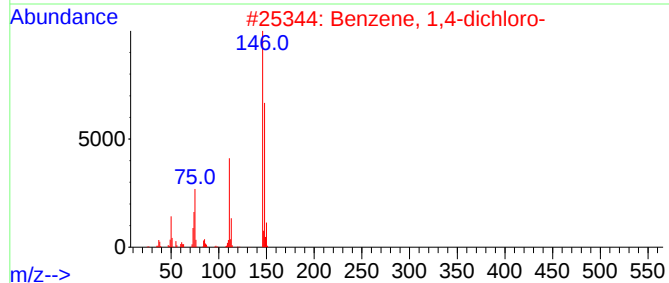
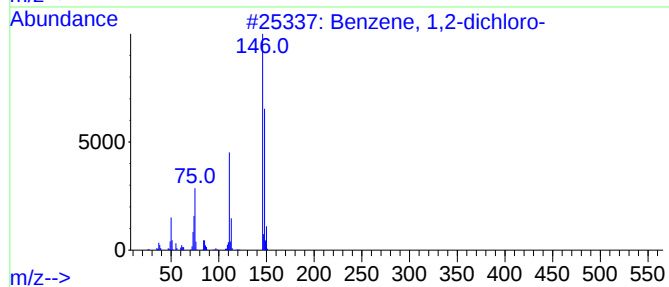
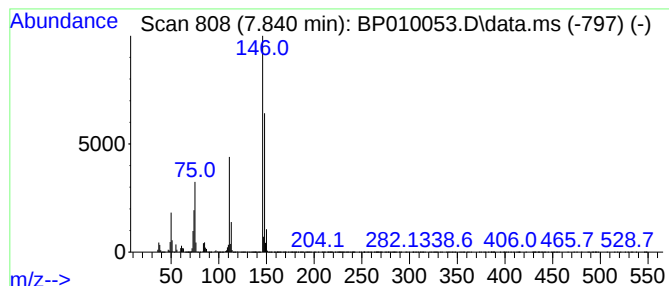
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.840	3.30 ng/ul	16326000	1,4-Dichlorobenzene-d4	7.958

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



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Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0HZ1DL

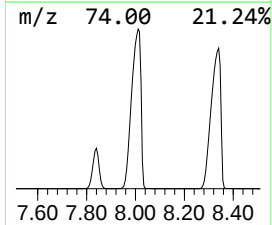
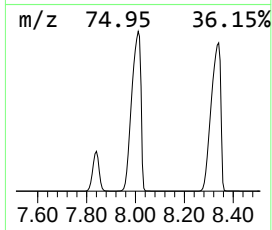
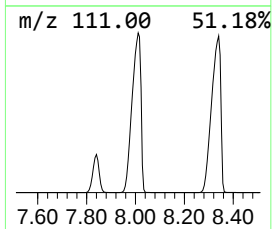
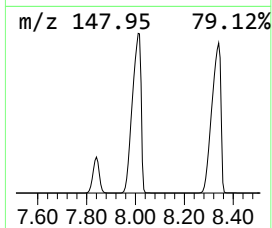
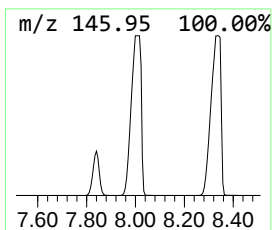
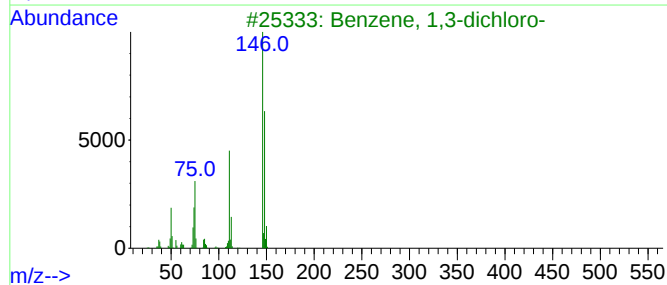
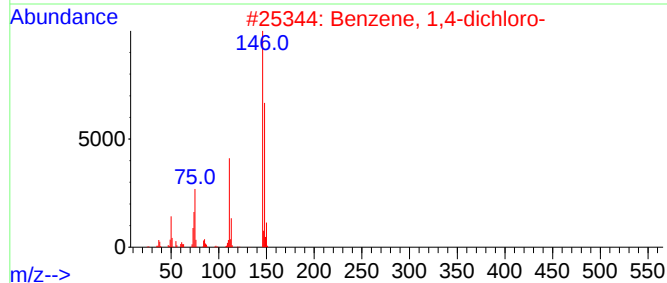
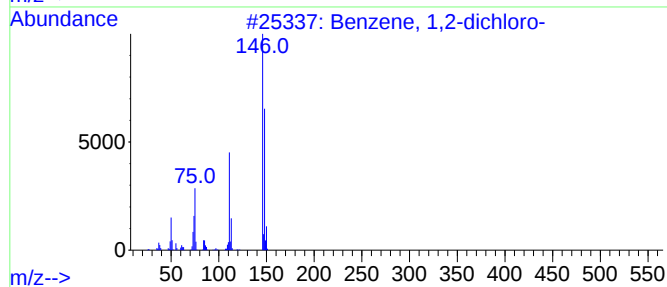
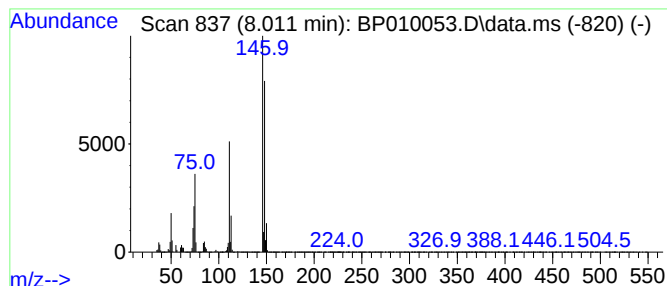
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.011	20.00 ng/ul	98910100	1,4-Dichlorobenzene-d4	7.958

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



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ALS Vial : 49 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C0HZ1DL

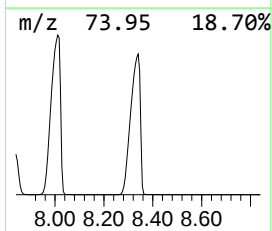
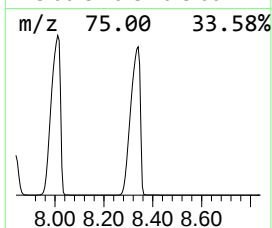
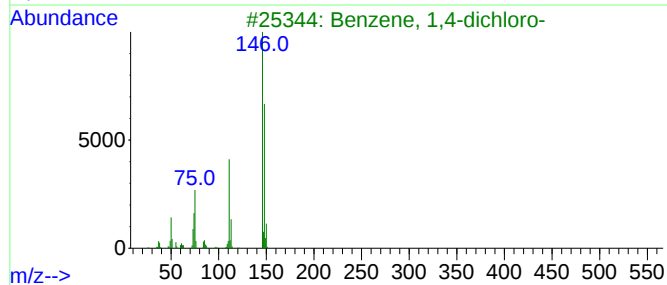
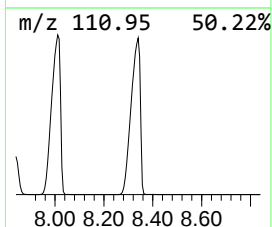
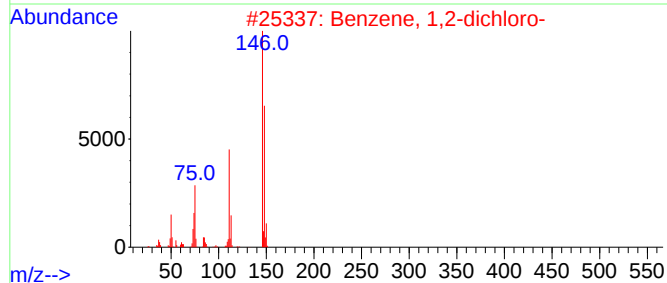
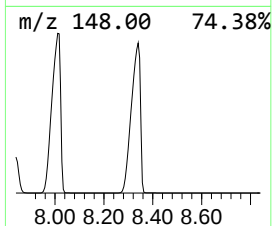
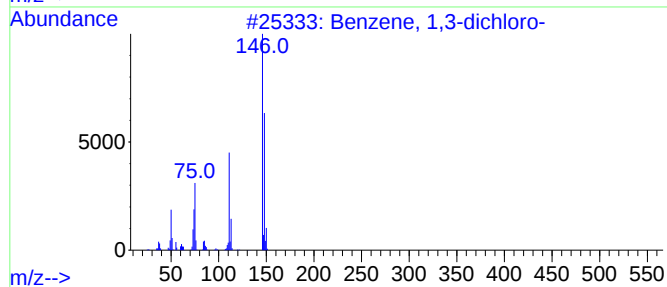
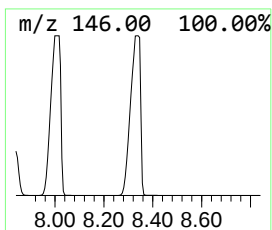
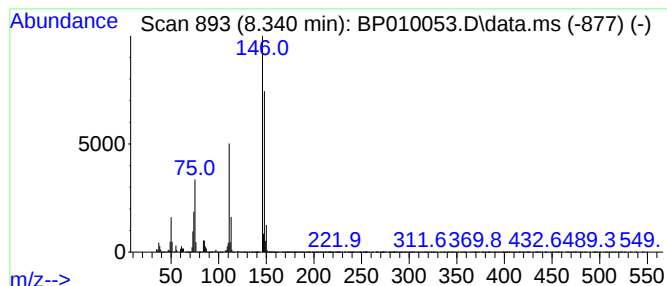
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.340	19.46 ng/ul	96225500	1,4-Dichlorobenzene-d4	7.958

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



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Instrument :
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ClientSampleId :
C0HZ1DL

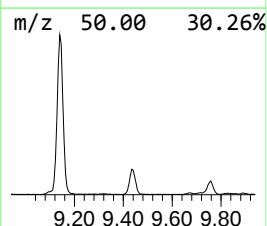
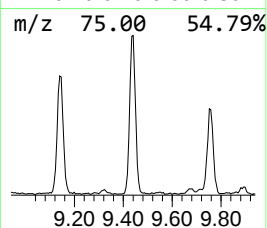
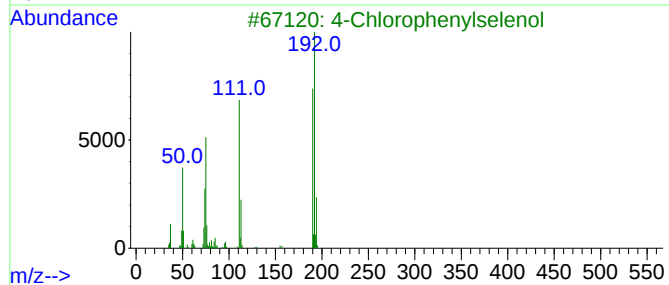
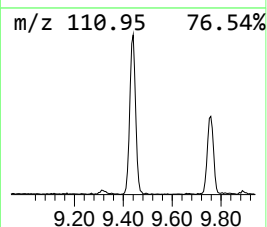
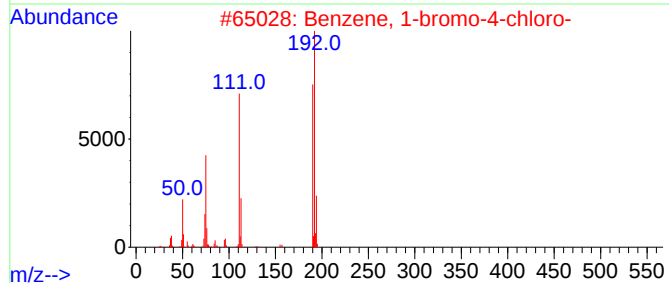
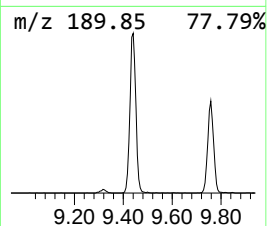
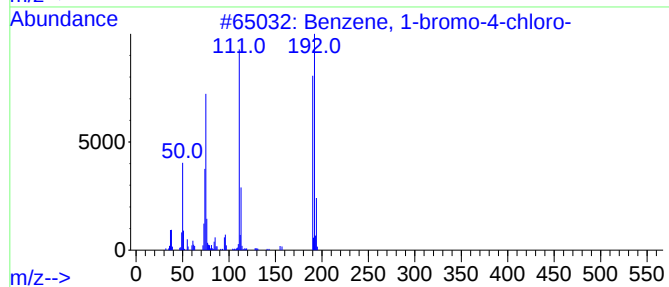
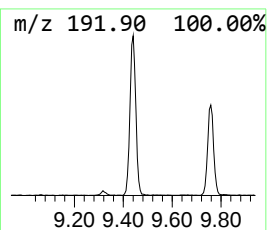
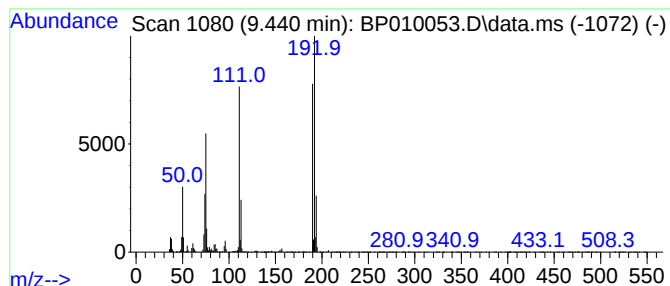
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.440	4.88 ng/ul	424816	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	99
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
4			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	96
5			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HZ1DL

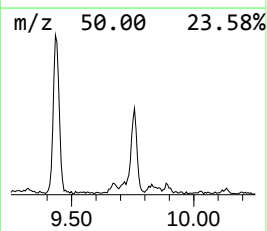
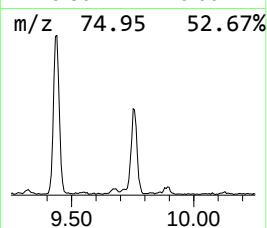
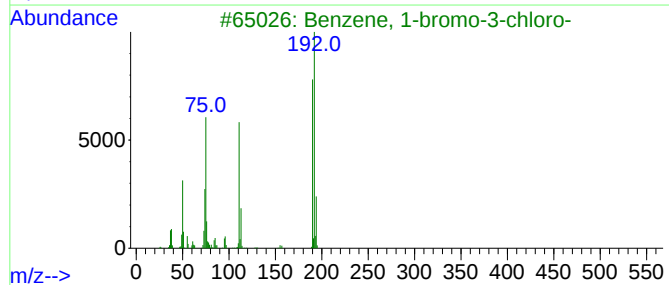
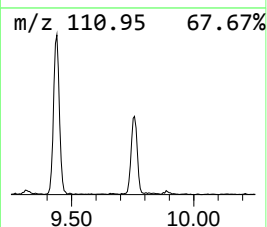
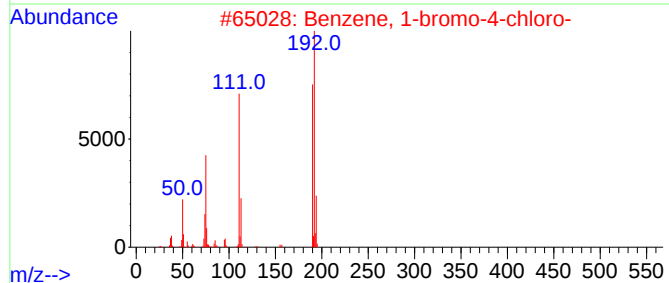
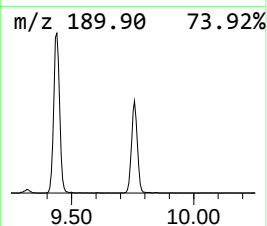
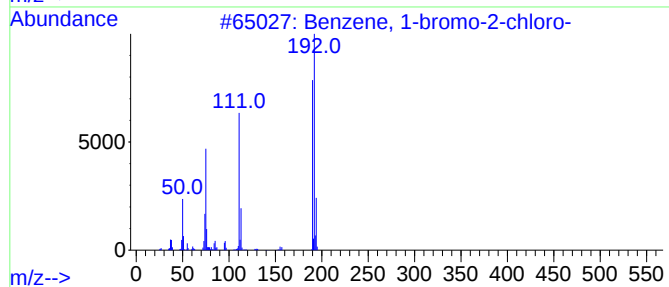
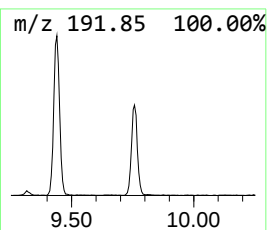
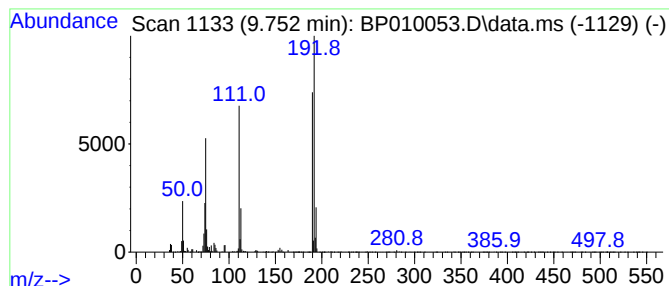
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.752	3.14 ng/ul	273460	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
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-2-chloro-	190	C6H4BrCl	000694-80-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HZ1DL

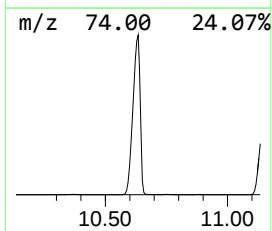
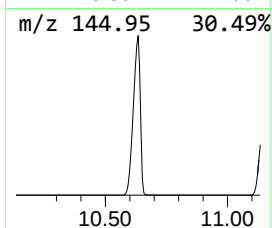
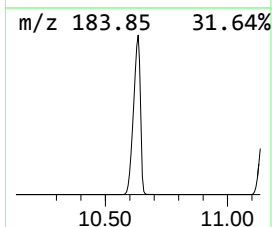
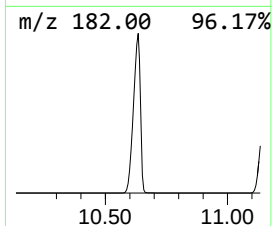
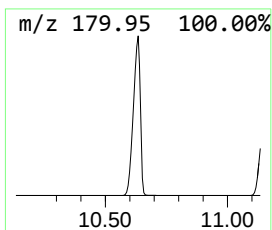
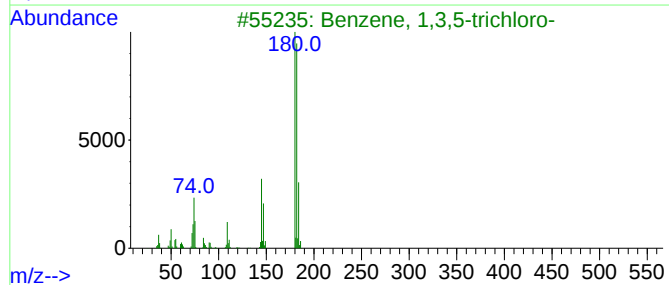
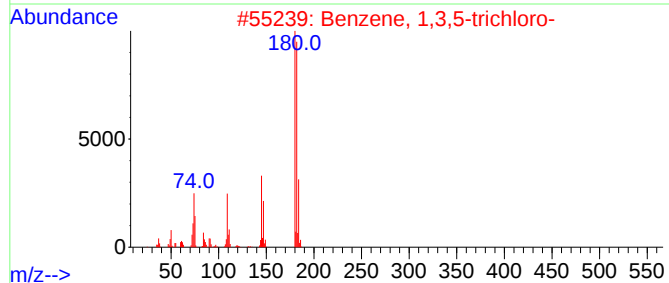
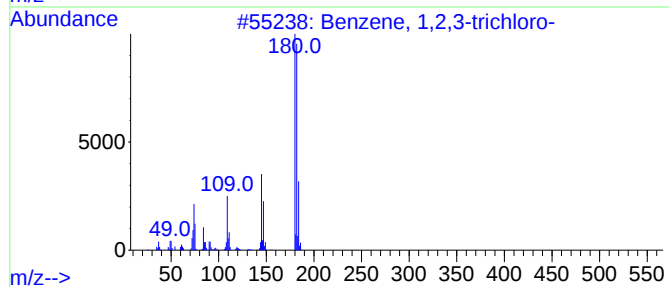
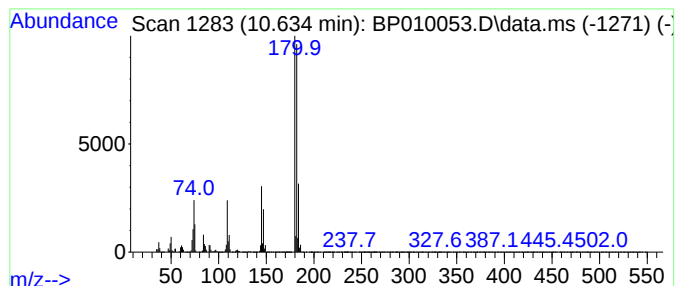
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.634	570.80 ng/ul	49728300	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
4			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HZ1DL

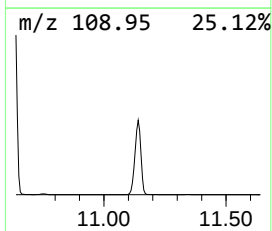
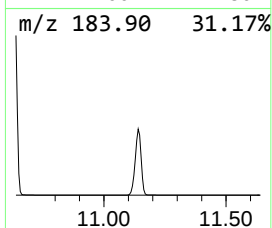
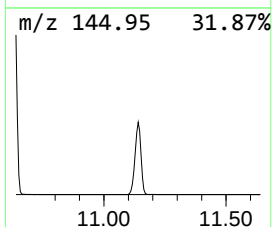
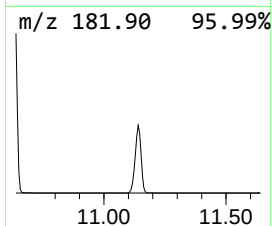
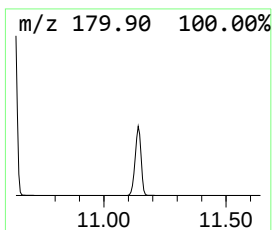
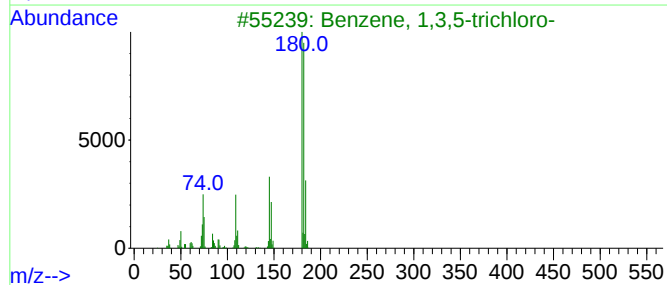
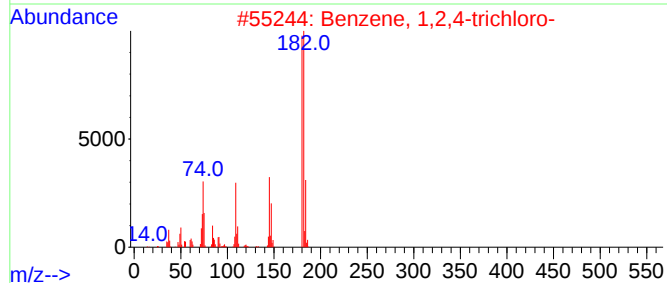
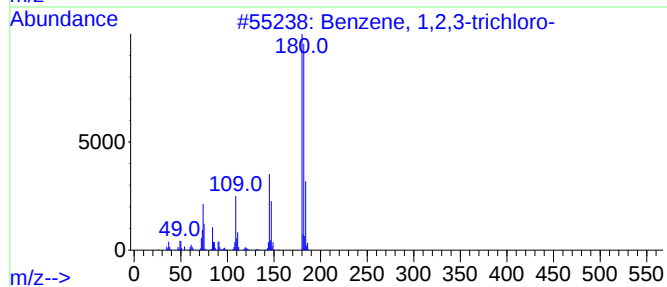
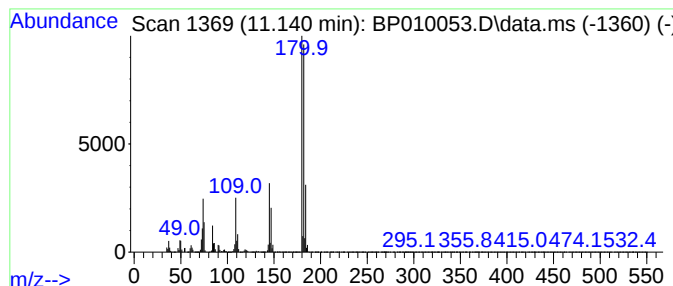
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Quant Title : SVOA CALIBRATION

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.
11.140	174.30 ng/ul	15185400	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	99
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, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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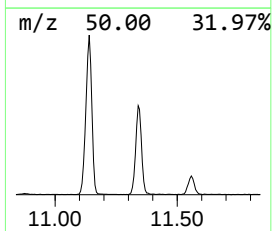
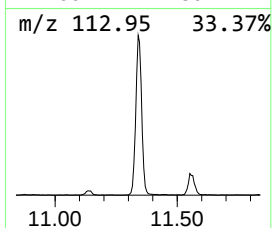
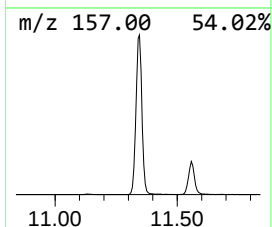
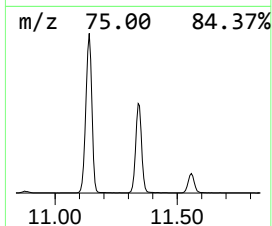
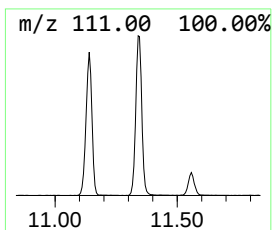
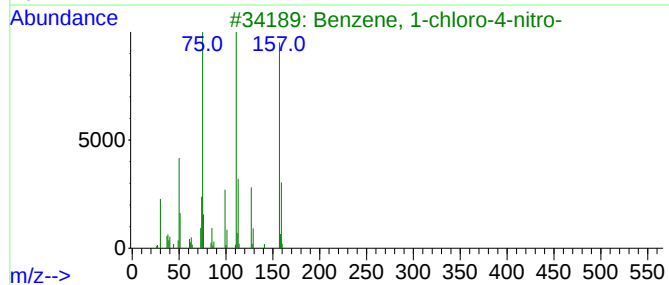
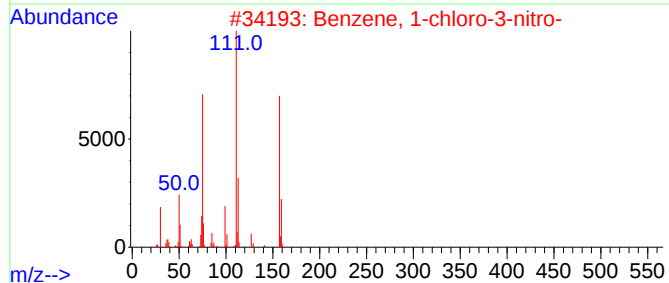
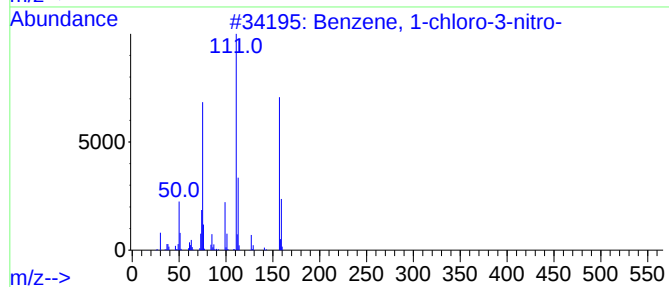
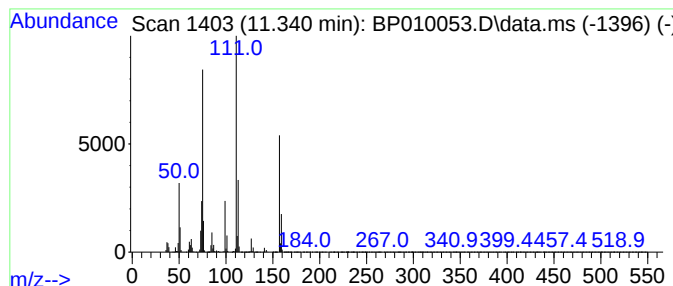
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.340	16.68 ng/ul	1453040	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	94
3			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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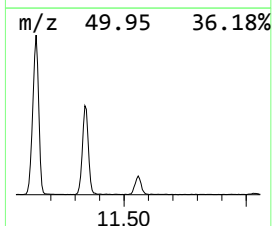
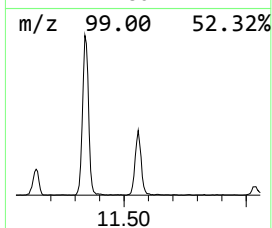
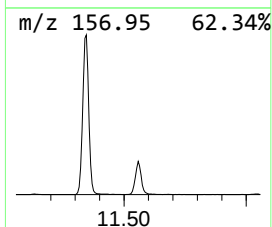
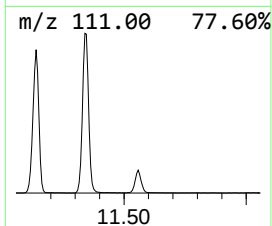
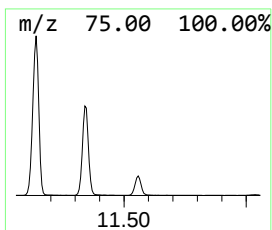
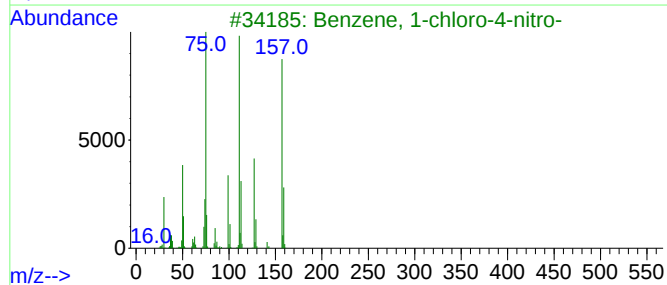
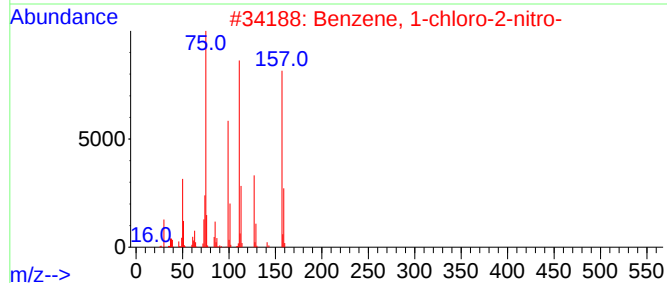
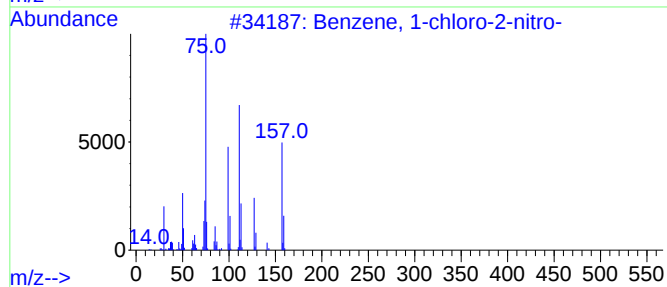
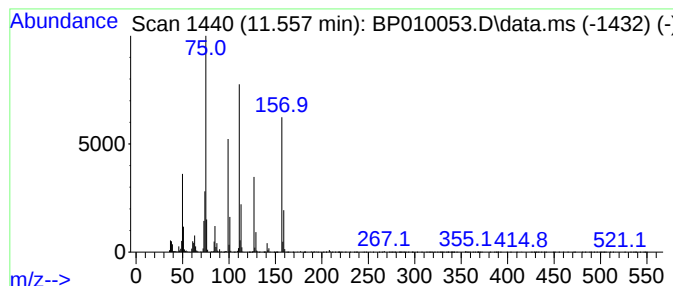
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.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.
11.557	3.82 ng/ul	332585	Naphthalene-d8	10.752

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
2			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96
5			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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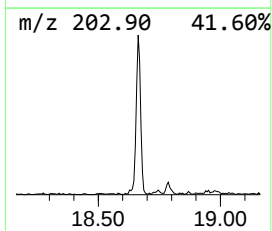
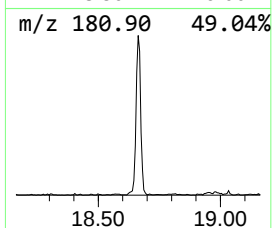
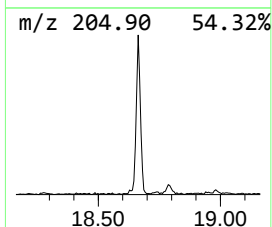
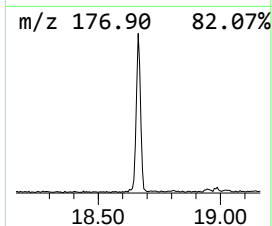
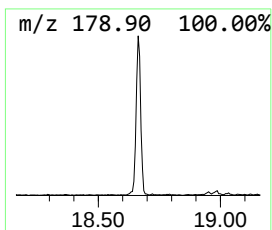
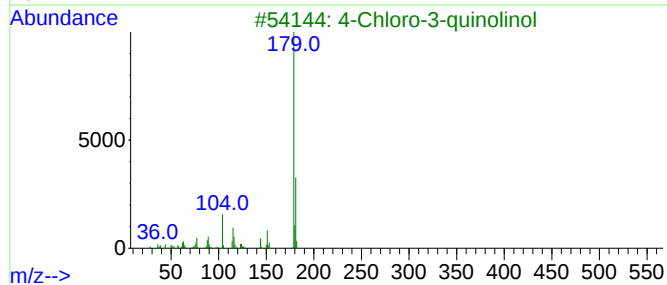
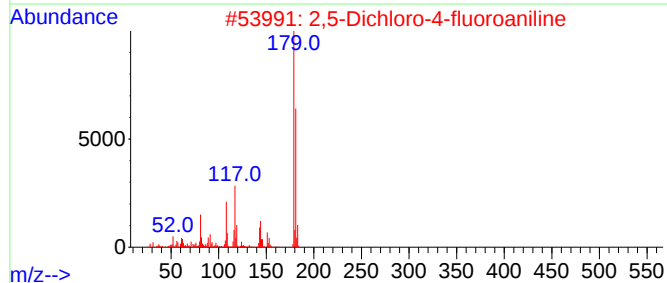
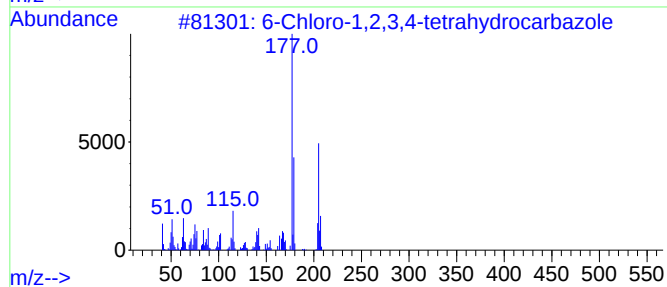
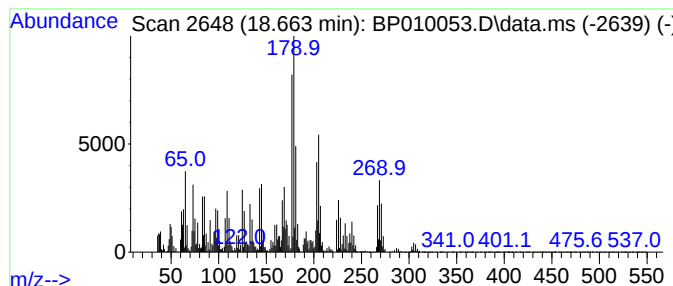
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.663	5.34 ng/ul	856080	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	30
2			2,5-Dichloro-4-fluoroaniline	179	C6H4Cl2FN	002729-37-5	18
3			4-Chloro-3-quinolinol	179	C9H6ClNO	032435-60-2	18
4			4,5-Dichloro-2-fluoroaniline	179	C6H4Cl2FN	002729-36-4	18
5			trans-.beta.-Ionone	192	C13H20O	000079-77-6	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010053.D
Acq On : 22 Apr 2022 13:30
Operator : CG/JU
Sample : N2473-09DL 5X
Misc :
ALS Vial : 49 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0HZ1DL

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.281	21.5	ng/ul	106478000	1	7.958	98910100	20.0
Benzene, 1,2-di...	7.840	3.3	ng/ul	16326000	1	7.958	98910100	20.0
Benzene, 1,4-di...	8.011	20.0	ng/ul	98910100	1	7.958	98910100	20.0
Benzene, 1,3-di...	8.340	19.5	ng/ul	96225500	1	7.958	98910100	20.0
Benzene, 1-brom...	9.440	4.9	ng/ul	424816	2	10.752	1742410	20.0
Benzene, 1-brom...	9.752	3.1	ng/ul	273460	2	10.752	1742410	20.0
Benzene, 1,2,3-...	10.634	570.8	ng/ul	49728300	2	10.752	1742410	20.0
Benzene, 1,2,4-...	11.140	174.3	ng/ul	15185400	2	10.752	1742410	20.0
Benzene, 1-chlo...	11.340	16.7	ng/ul	1453040	2	10.752	1742410	20.0
Benzene, 1-chlo...	11.557	3.8	ng/ul	332585	2	10.752	1742410	20.0
unknown-01	18.663	5.3	ng/ul	856080	4	17.322	3203920	20.0