

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
 Data File : BP010920.D
 Acq On : 30 Jun 2022 15:34
 Operator : CG/JU
 Sample : N3523-09DL 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0JY0DL

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-BP062322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010920.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.616	87	90	103	rBV	245496	360833	0.17%	0.041%
2	5.087	325	340	344	rBV4	53391239	199567703	92.82%	22.802%
3	6.393	553	562	573	rVB	226615	366818	0.17%	0.042%
4	6.893	637	647	660	rVB2	116567	239657	0.11%	0.027%
5	7.057	668	675	687	rBV	725032	1167991	0.54%	0.133%
6	7.252	701	708	719	rBV	451015	750328	0.35%	0.086%
7	7.610	760	769	779	rVB	29777909	55238818	25.69%	6.312%
8	7.793	779	800	805	rBV4	51979443	188072367	87.47%	21.489%
9	8.122	839	856	861	rBV5	52760991	215014455	100.00%	24.567%
10	8.428	902	908	917	rVB	376199	572079	0.27%	0.065%
11	8.875	977	984	987	rBV	784621	1272278	0.59%	0.145%
12	8.916	987	991	1001	rVB	2165834	3526675	1.64%	0.403%
13	9.204	1033	1040	1050	rBV	677686	1088608	0.51%	0.124%
14	9.434	1073	1079	1083	rBV	159388	260753	0.12%	0.030%
15	9.481	1083	1087	1089	rVV	154647	248411	0.12%	0.028%
16	9.522	1089	1094	1100	rVV	339210	618501	0.29%	0.071%
17	9.593	1100	1106	1112	rVV	358038	599048	0.28%	0.068%
18	10.134	1191	1198	1206	rBV	550654	906137	0.42%	0.104%
19	10.398	1231	1243	1249	rBV2	48934894	114603110	53.30%	13.094%
20	10.516	1256	1263	1270	rBV	2070023	3184955	1.48%	0.364%
21	10.651	1277	1286	1298	rVB	575950	957527	0.45%	0.109%
22	10.898	1318	1328	1338	rBV	19278357	33326230	15.50%	3.808%
23	11.116	1354	1365	1372	rBV	1145151	1834063	0.85%	0.210%
24	11.328	1394	1401	1409	rBV	246929	438073	0.20%	0.050%
25	12.045	1516	1523	1537	rBV2	121189	269951	0.13%	0.031%
26	12.510	1593	1602	1603	rBV3	248074	348290	0.16%	0.040%
27	12.545	1603	1608	1617	rVB	2000157	2974714	1.38%	0.340%
28	13.157	1705	1712	1724	rVB	5159099	7743852	3.60%	0.885%
29	13.786	1808	1819	1835	rBV	1443740	2068914	0.96%	0.236%
30	14.063	1859	1866	1875	rVB	1578709	2378775	1.11%	0.272%
31	14.369	1909	1918	1934	rBV	2658894	3860531	1.80%	0.441%
32	14.686	1966	1972	1979	rVB	166888	245020	0.11%	0.028%
33	15.369	2081	2088	2093	rBV	2168281	3096583	1.44%	0.354%
34	15.486	2102	2108	2118	rBV	193354	293177	0.14%	0.033%
35	17.133	2381	2388	2398	rBV	3024515	4290351	2.00%	0.490%
36	17.233	2399	2405	2418	rVV	2288209	3211239	1.49%	0.367%

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Instrument :
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ClientSampleId :
C0JY0DL

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-BP062322.M
Title : SVOA CALIBRATION

37	18.498	2612	2620	2628	rBV	508888	692783	0.32%	0.079%
38	19.486	2782	2788	2795	rBV	3090236	4030544	1.87%	0.461%
39	20.974	3038	3041	3046	rBV4	162907	216348	0.10%	0.025%
40	21.245	3082	3087	3094	rBV	4170662	5315867	2.47%	0.607%
41	23.462	3457	3464	3474	rBV	2242059	4392132	2.04%	0.502%
42	23.615	3483	3490	3501	rVB2	2772457	5561003	2.59%	0.635%

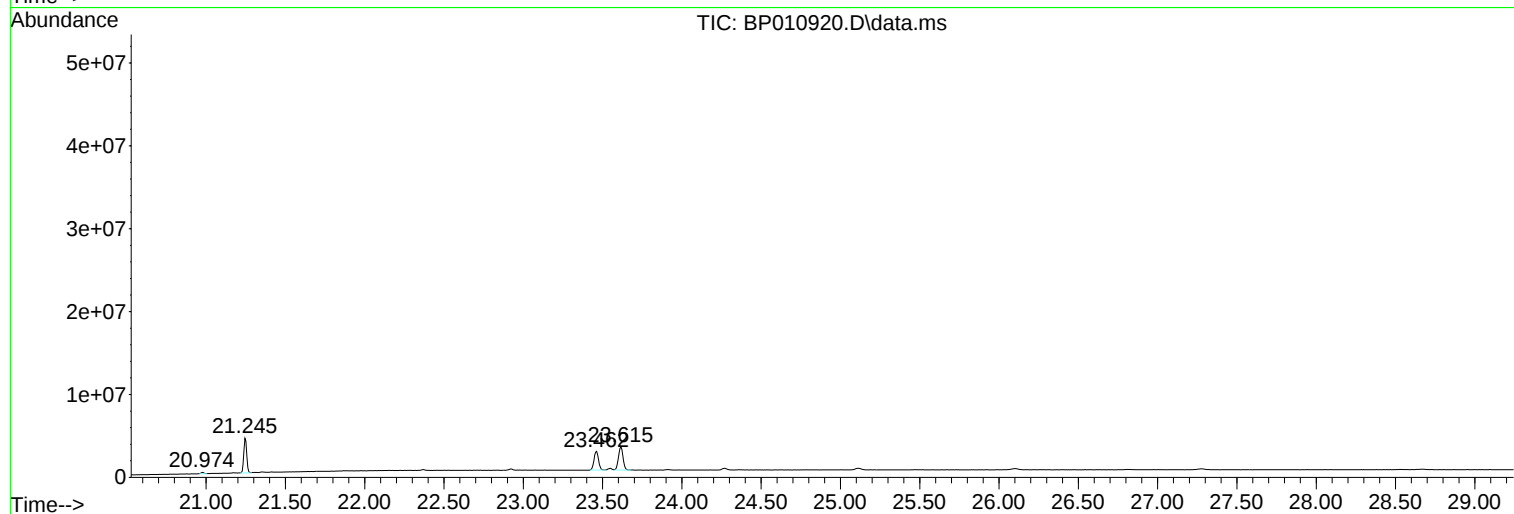
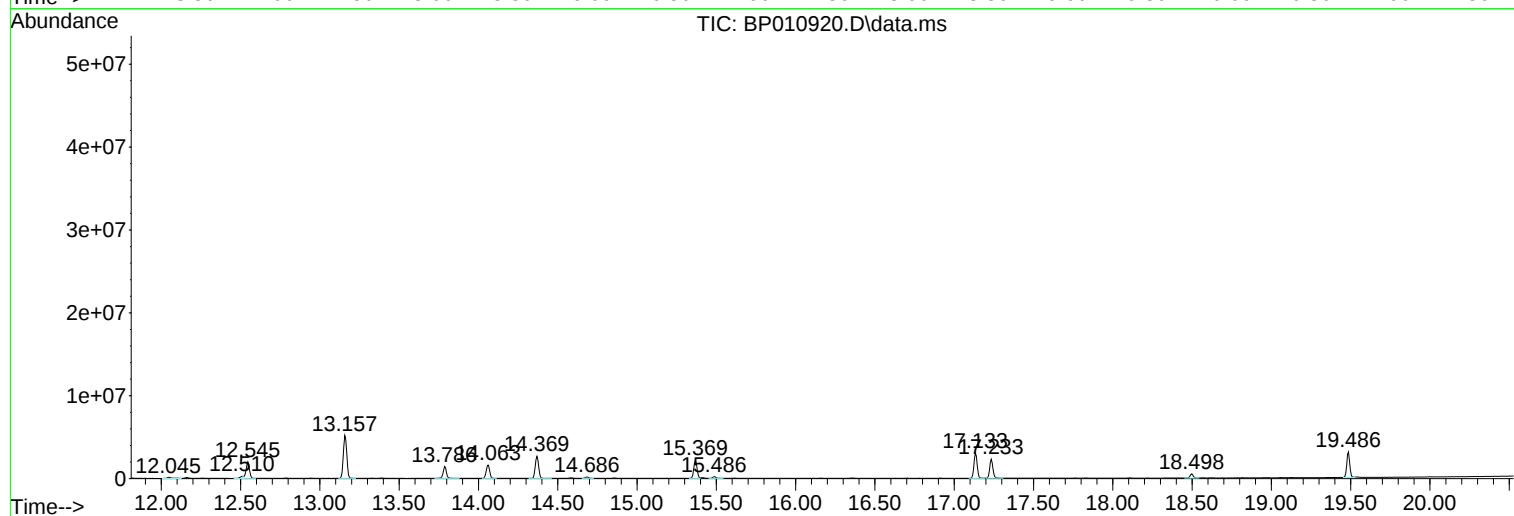
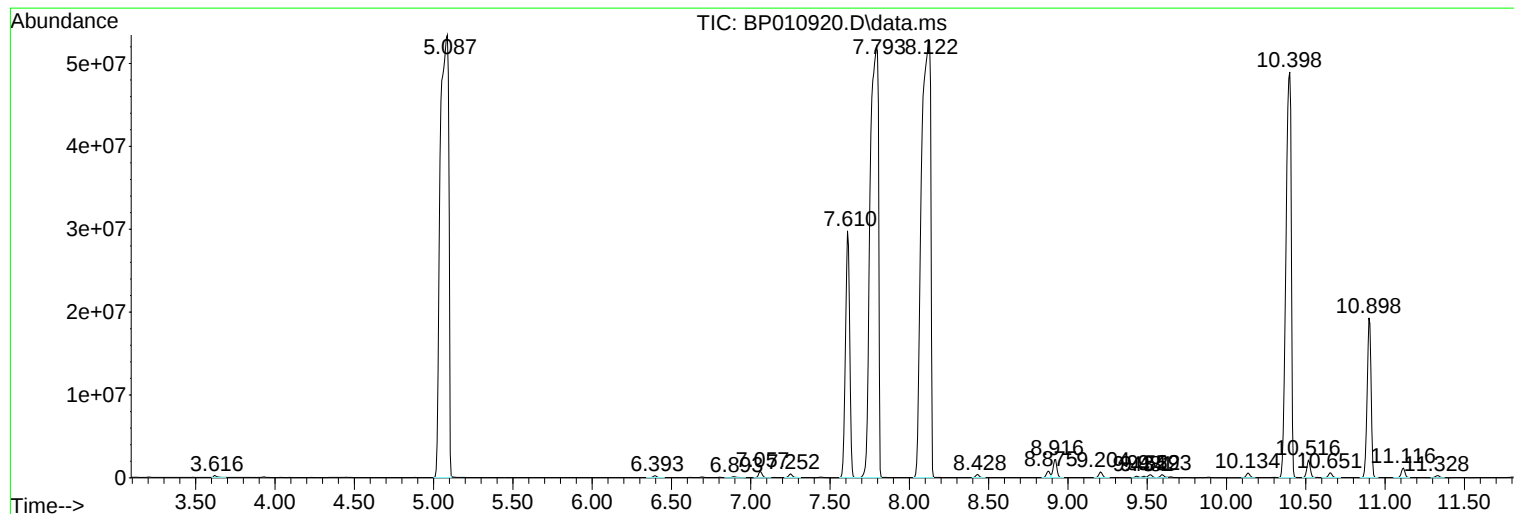
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ClientSampleId :
C0JY0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_P
ClientSampleId :
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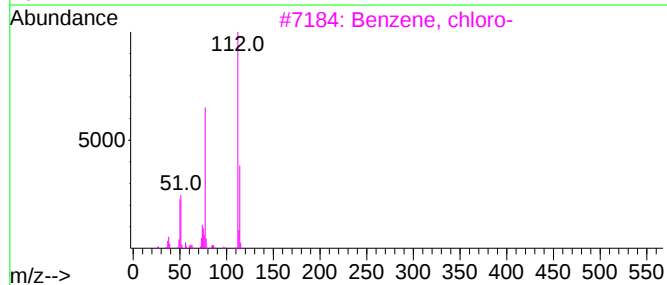
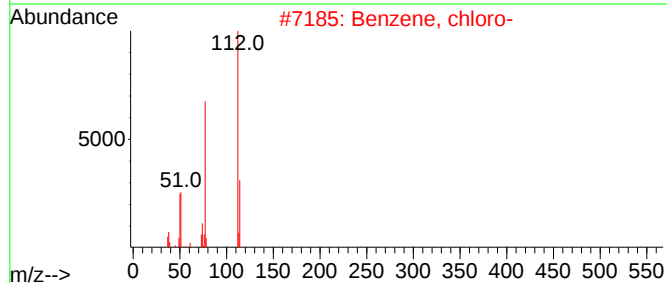
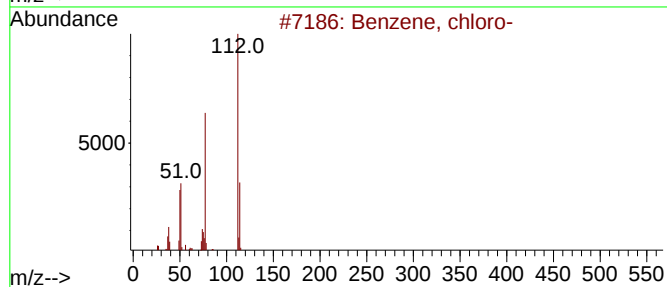
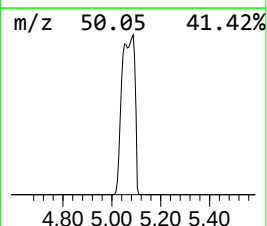
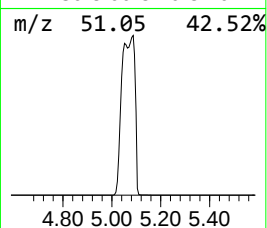
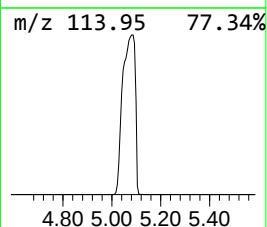
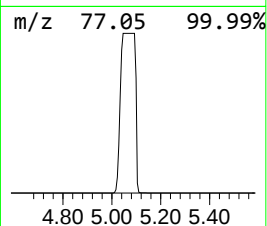
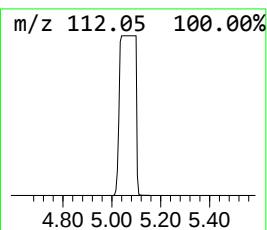
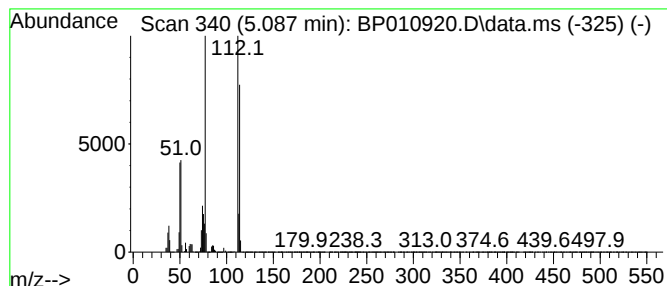
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.087	21.22 ng/ul	199568000	1,4-Dichlorobenzene-d4	7.728

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	91
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	80
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	43



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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

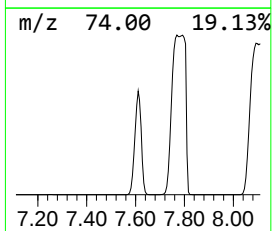
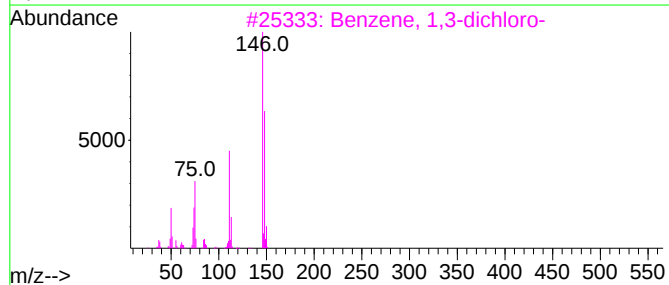
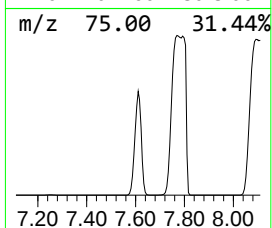
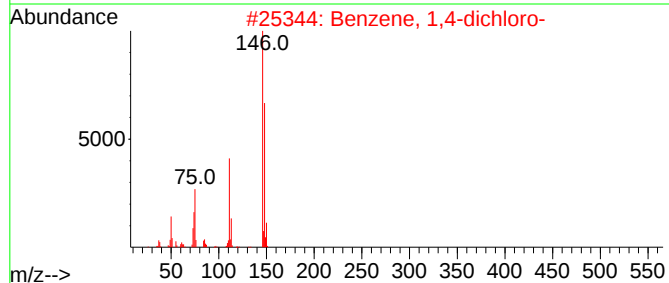
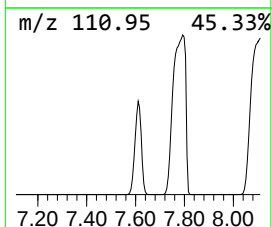
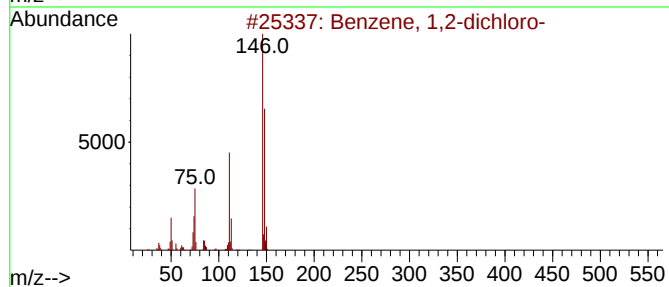
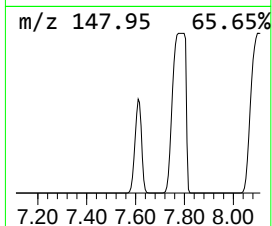
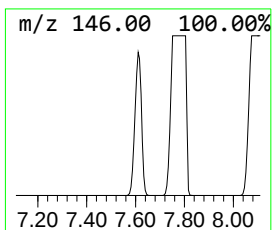
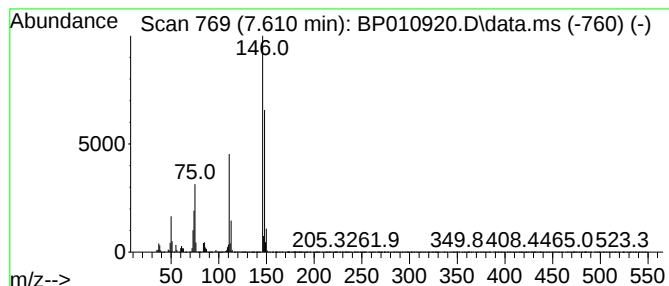
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.610	5.87 ng/ul	55238800	1,4-Dichlorobenzene-d4	7.728

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, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

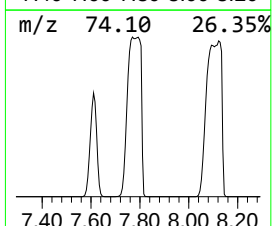
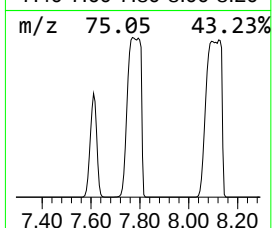
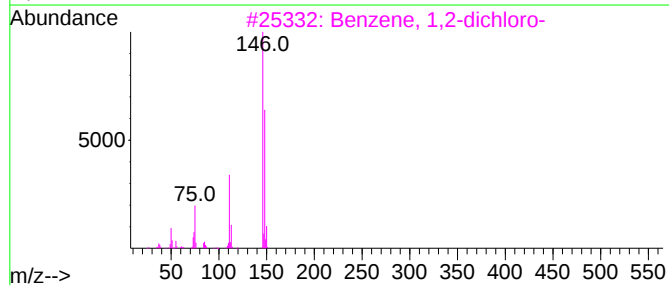
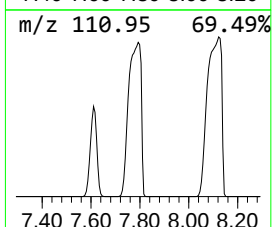
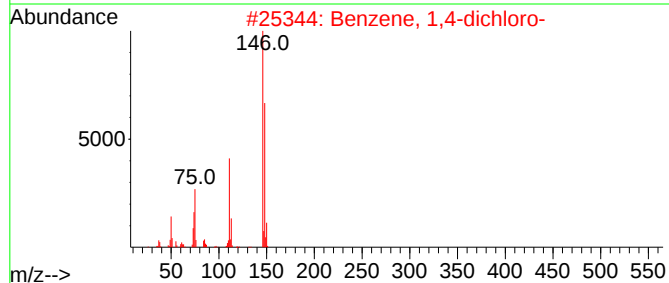
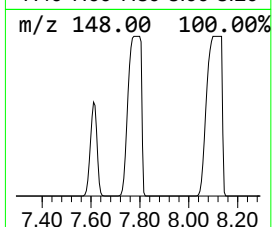
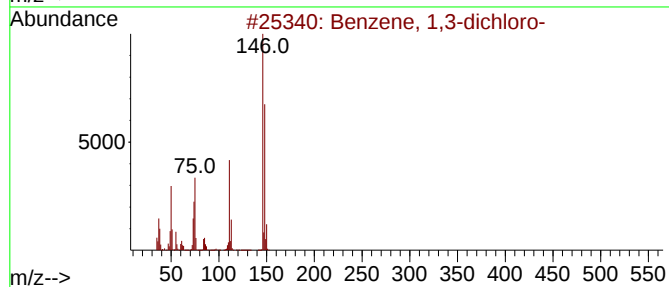
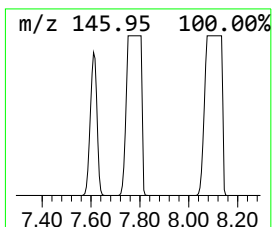
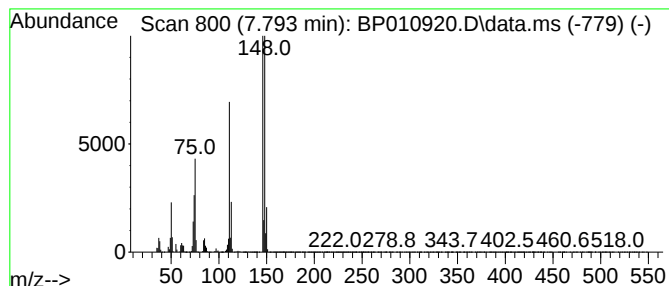
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.793	20.00 ng/ul	188072000	1,4-Dichlorobenzene-d4	7.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	89
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	86
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	81



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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

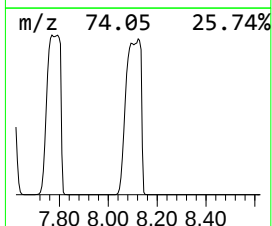
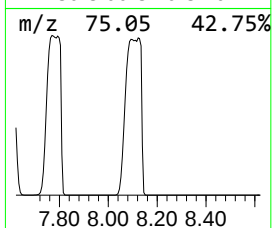
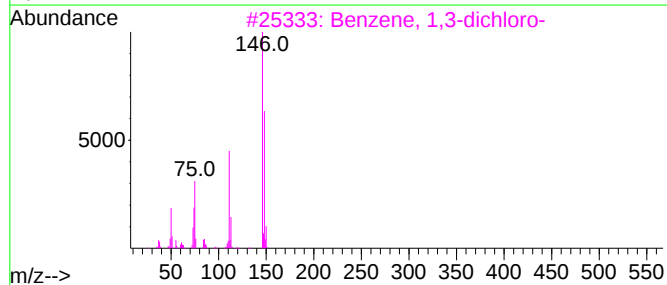
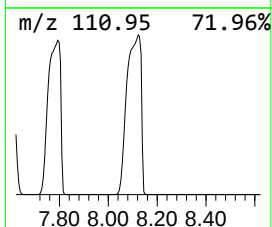
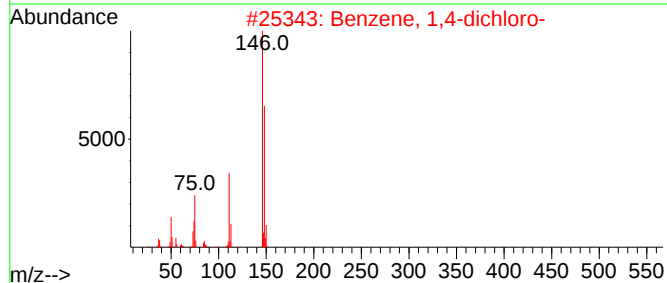
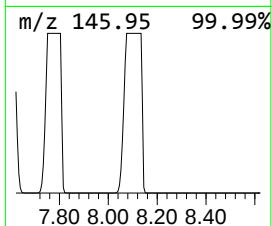
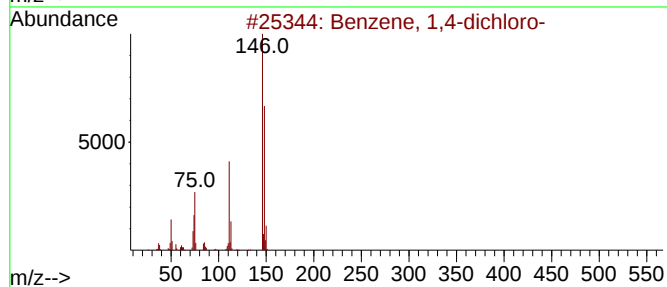
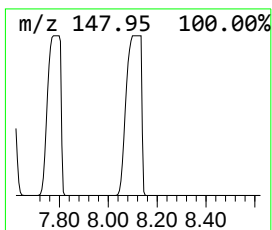
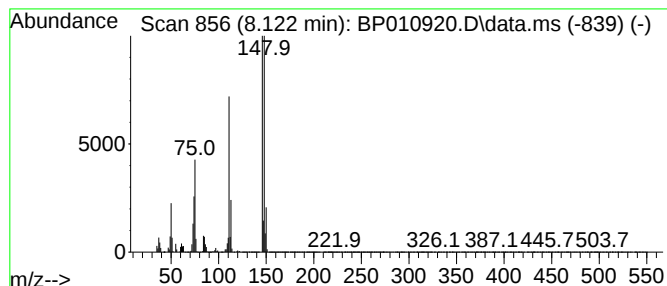
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	22.87 ng/ul	215014000	1,4-Dichlorobenzene-d4	7.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	86
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	81
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	81



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Instrument :
BNA_P
ClientSampleId :
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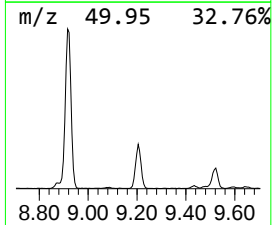
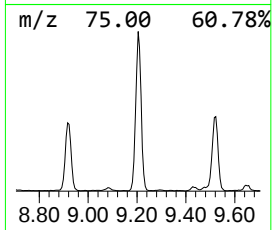
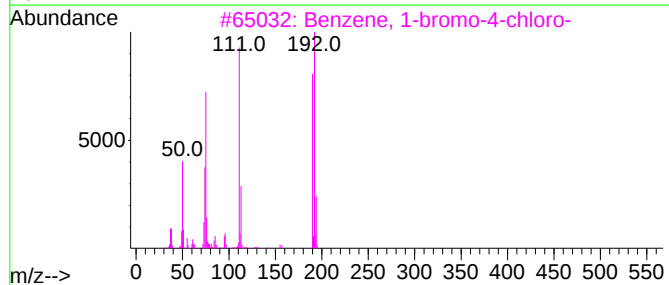
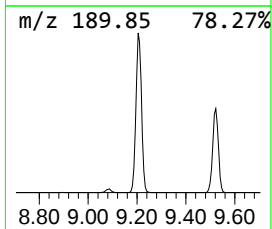
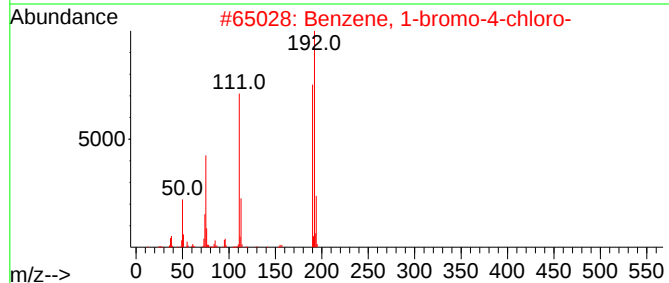
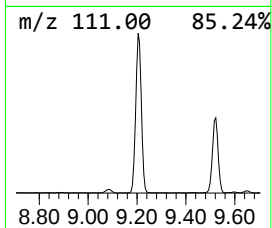
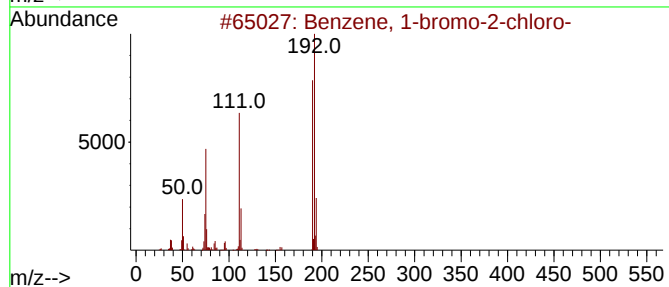
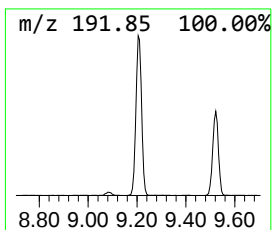
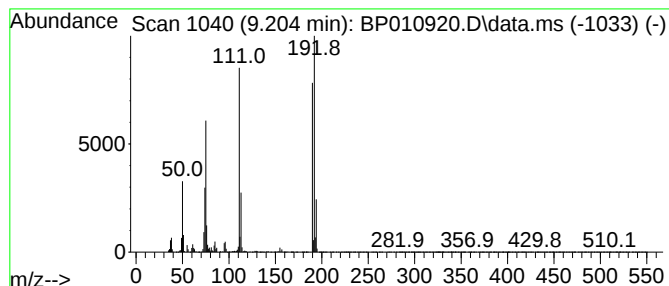
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.204	6.84 ng/ul	1088610	Naphthalene-d8	10.516

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	98
3			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

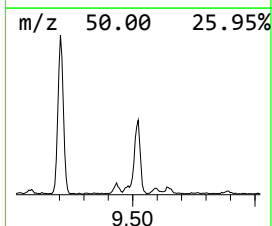
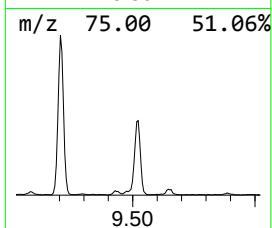
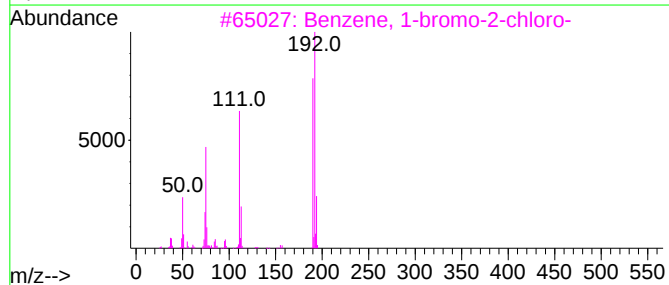
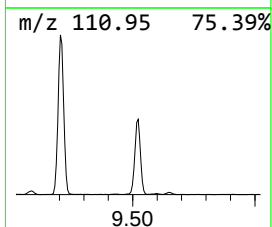
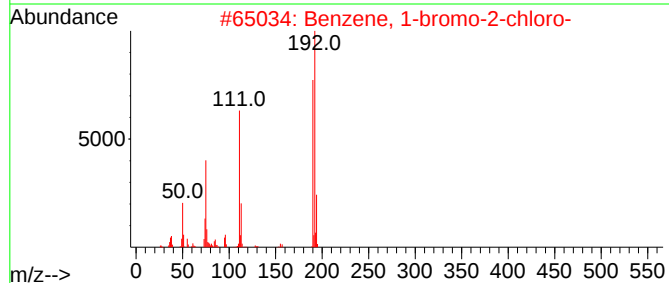
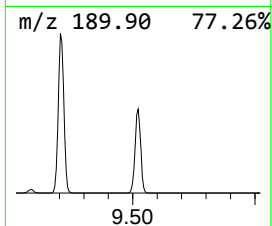
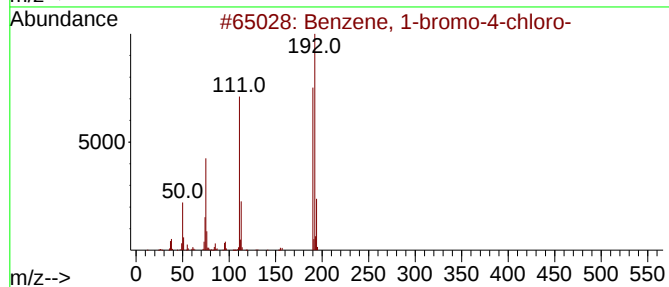
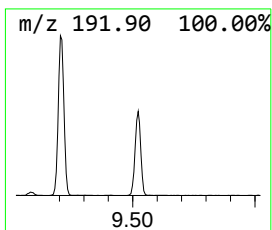
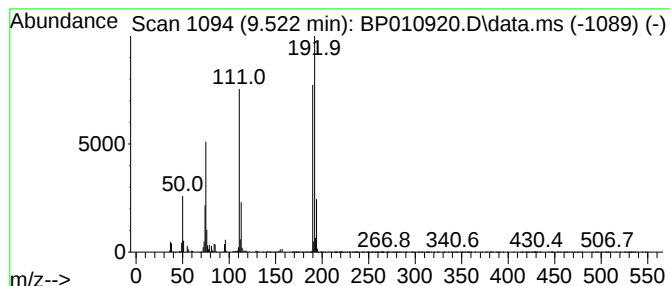
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.522	3.88 ng/ul	618501	Naphthalene-d8	10.516

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-2-chloro-	190	C6H4BrCl	000694-80-4	98
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

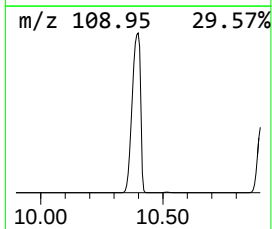
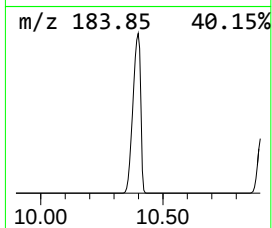
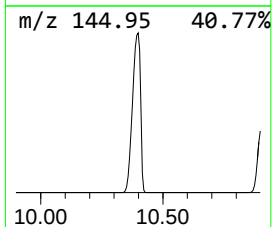
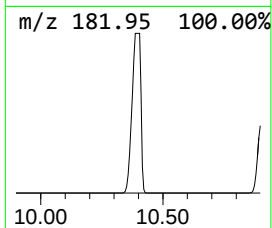
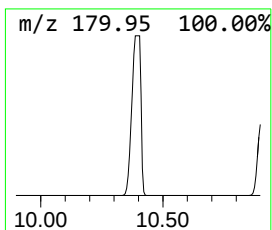
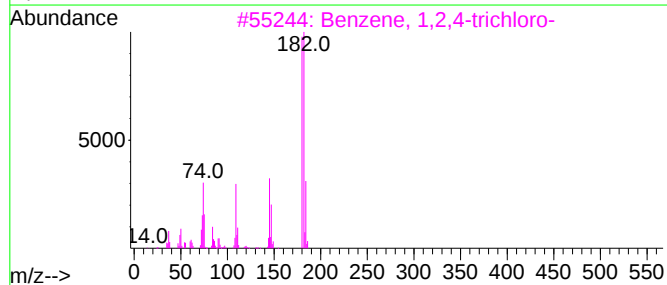
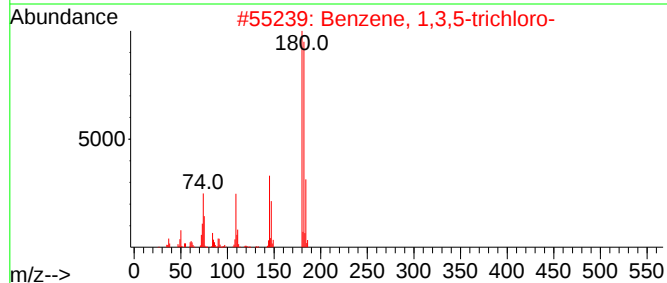
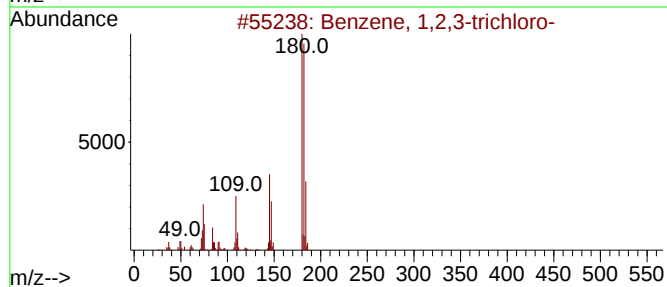
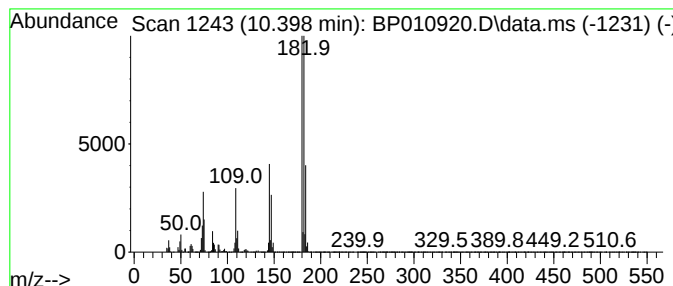
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Quant Title : SVOA CALIBRATION

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.
10.398	719.65 ng/ul	114603000	Naphthalene-d8	10.516

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, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
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_P\Data\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

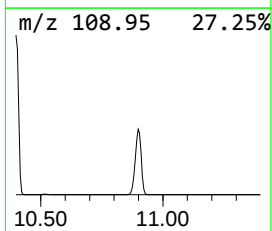
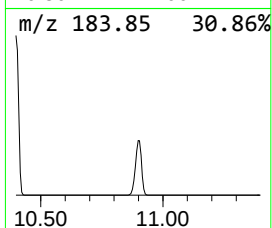
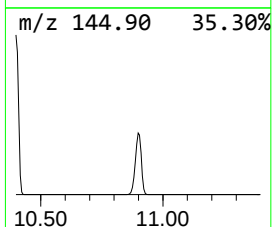
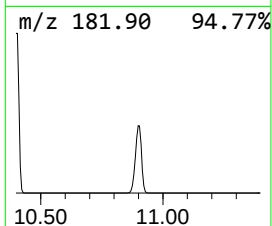
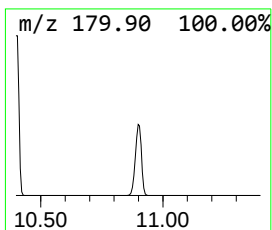
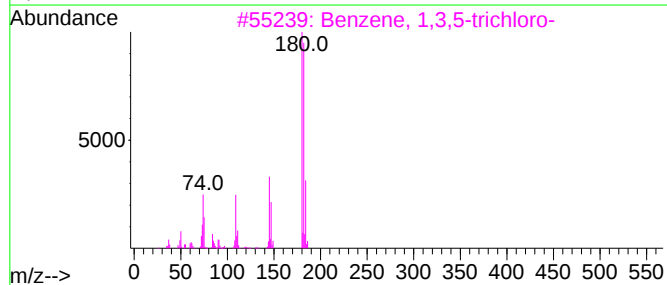
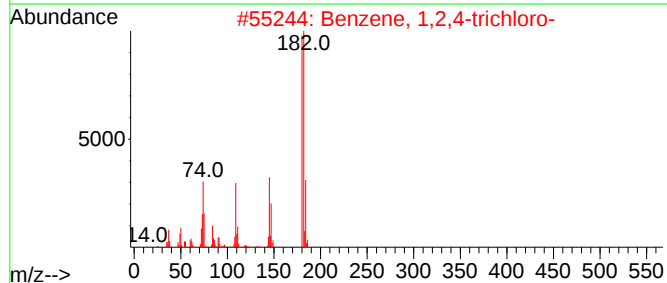
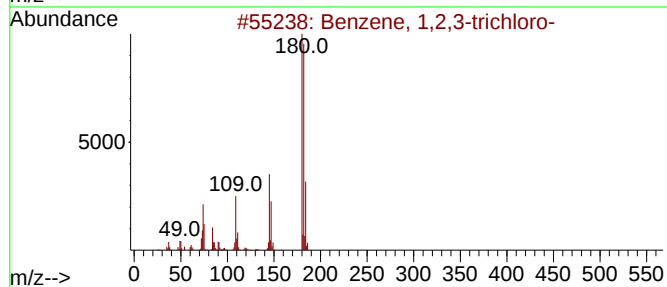
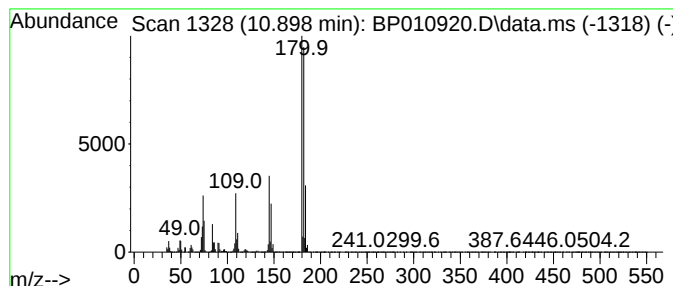
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Quant Title : SVOA CALIBRATION

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.
10.898	209.27 ng/ul	33326200	Naphthalene-d8	10.516

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\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

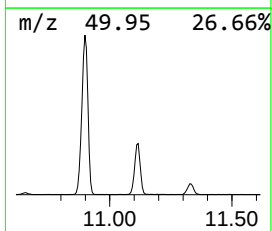
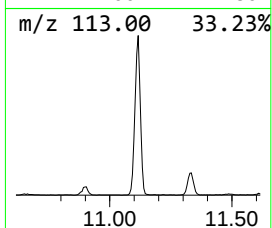
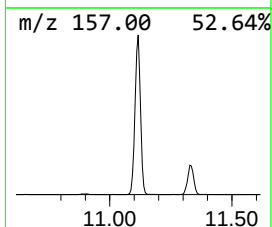
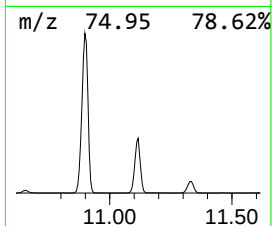
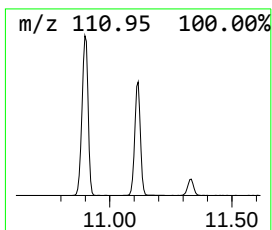
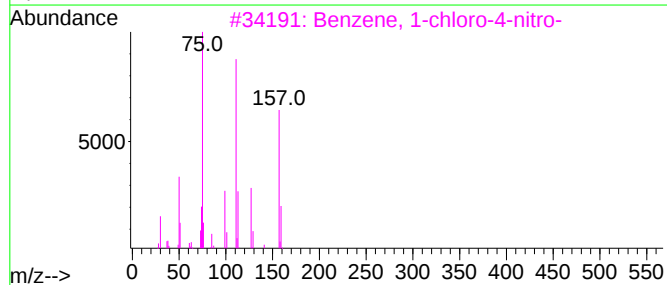
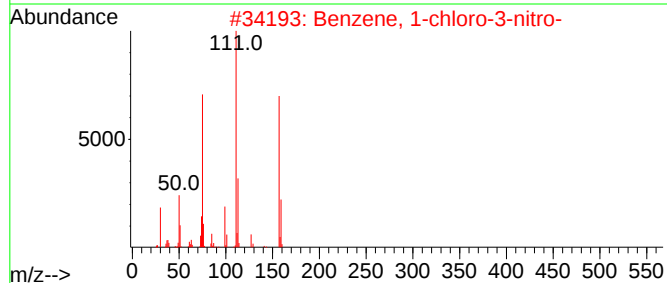
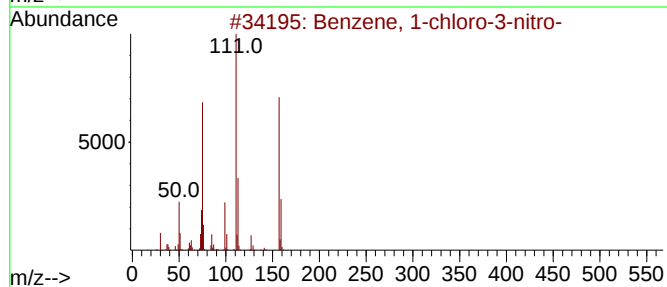
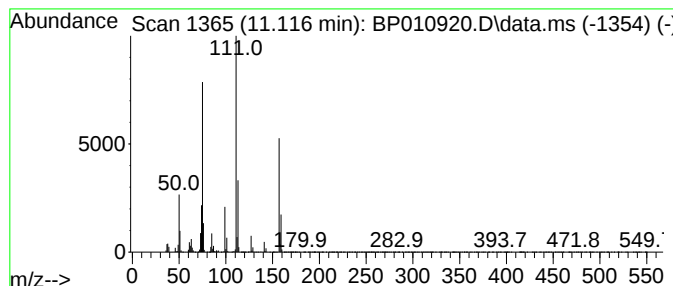
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Quant Title : SVOA CALIBRATION

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.
11.116	11.52 ng/ul	1834060	Naphthalene-d8	10.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
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-3-nitro-	157	C6H4ClNO2	000121-73-3	92
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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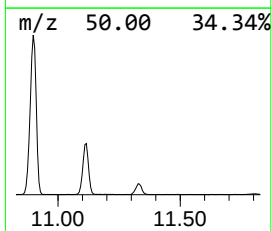
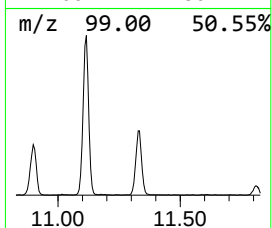
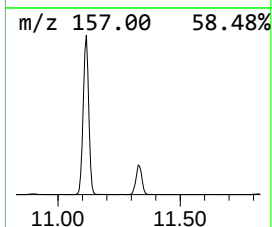
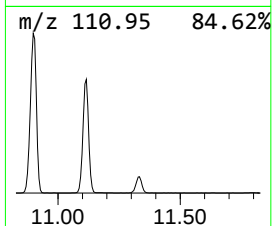
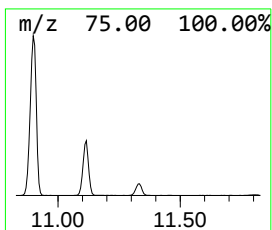
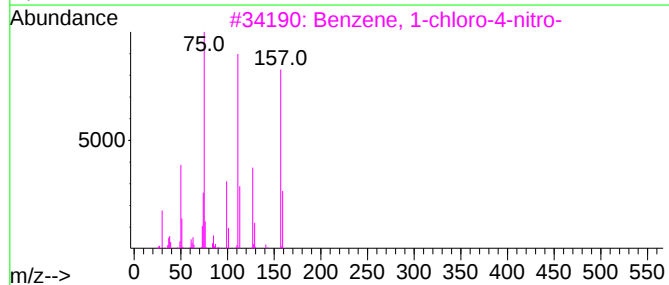
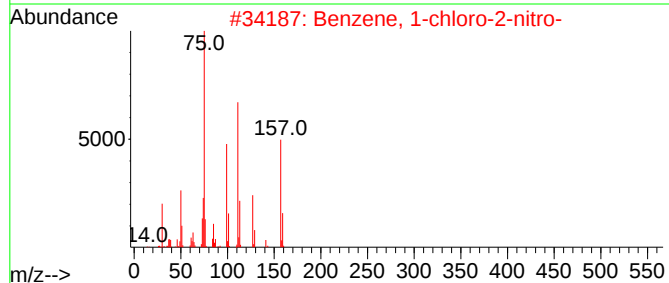
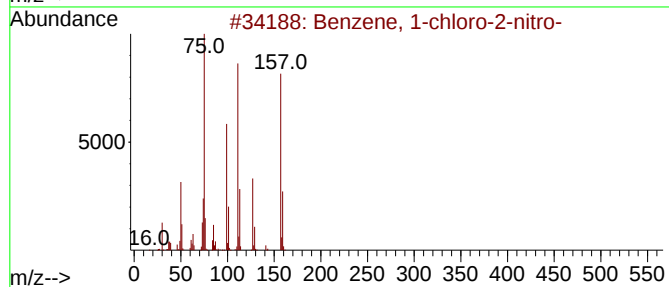
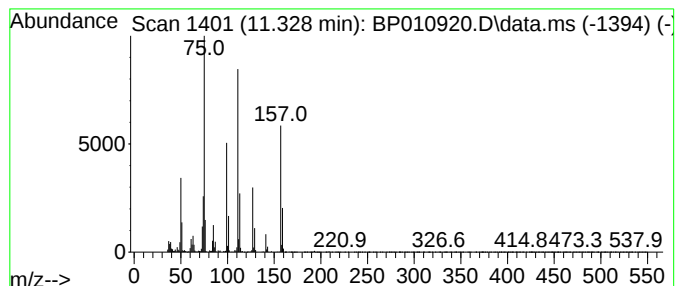
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.328	2.75 ng/ul	438073	Naphthalene-d8	10.516

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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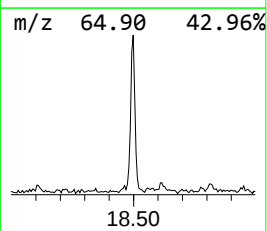
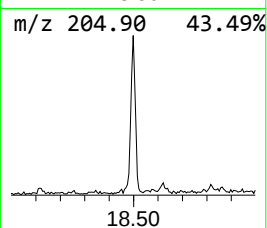
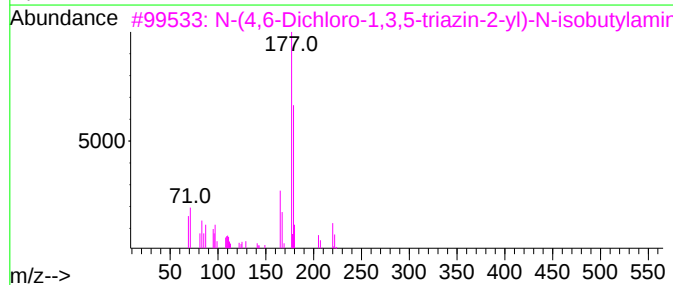
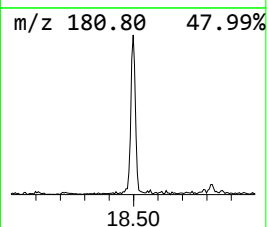
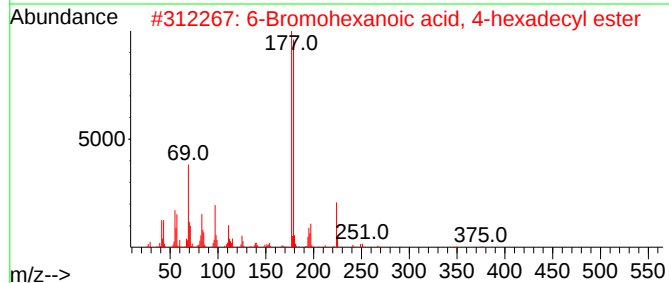
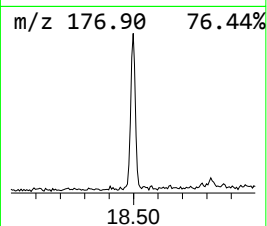
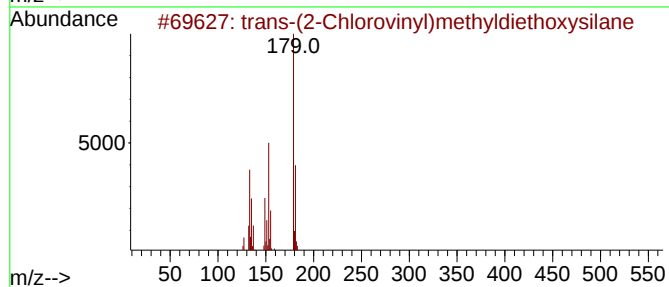
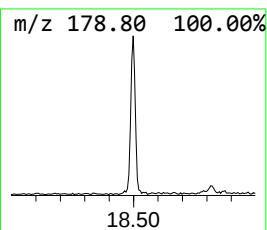
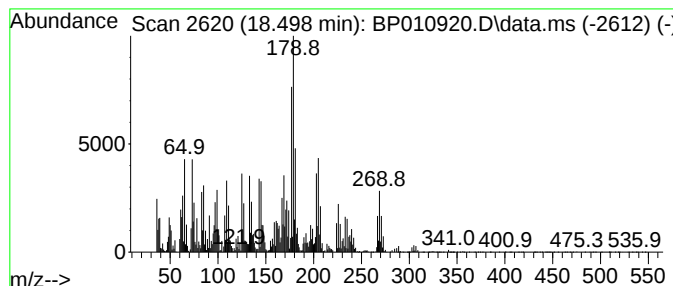
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.498	3.23 ng/ul	692783	Phenanthrene-d10	17.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-(2-Chlorovinyl)methyldieth...	194	C7H15ClO2Si	1000139-96-9	27
2			6-Bromohexanoic acid, 4-hexadecy...	418	C22H43BrO2	1000282-90-6	18
3			N-(4,6-Dichloro-1,3,5-triazin-2-...	220	C7H10Cl2N4	1000326-88-2	16
4			3-Chloro-1H-indole-2-carbaldehyde	179	C9H6ClNO	110912-15-7	14
5			2-Chloro-6-quinolinol	179	C9H6ClNO	577967-89-6	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063022\
Data File : BP010920.D
Acq On : 30 Jun 2022 15:34
Operator : CG/JU
Sample : N3523-09DL 2X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0JY0DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.087	21.2	ng/ul	199568000	1	7.728	188072000	20.0
Benzene, 1,2-di...	7.610	5.9	ng/ul	55238800	1	7.728	188072000	20.0
Benzene, 1,3-di...	7.793	20.0	ng/ul	188072000	1	7.728	188072000	20.0
Benzene, 1,4-di...	8.122	22.9	ng/ul	215014000	1	7.728	188072000	20.0
Benzene, 1-brom...	9.204	6.8	ng/ul	1088610	2	10.516	3184960	20.0
Benzene, 1-brom...	9.522	3.9	ng/ul	618501	2	10.516	3184960	20.0
Benzene, 1,2,3-...	10.398	719.6	ng/ul	114603000	2	10.516	3184960	20.0
Benzene, 1,2,4-...	10.898	209.3	ng/ul	33326200	2	10.516	3184960	20.0
Benzene, 1-chlo...	11.116	11.5	ng/ul	1834060	2	10.516	3184960	20.0
Benzene, 1-chlo...	11.328	2.8	ng/ul	438073	2	10.516	3184960	20.0
unknown-01	18.498	3.2	ng/ul	692783	4	17.133	4290350	20.0