

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
 Data File : BN031103.D
 Acq On : 21 Apr 2024 11:12
 Operator : MA/JU
 Sample : P2159-19
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CORE0

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_N\Methods\SFAM-EPA-BN041924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031103.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.558	94	97	120	rBV	113651	226446	0.58%	0.117%
2	4.987	330	340	359	rBV	18677028	39280475	100.00%	20.355%
3	6.805	644	649	658	rVB	84432	139043	0.35%	0.072%
4	6.975	671	678	689	rBV	460833	719702	1.83%	0.373%
5	7.164	701	710	718	rBV	398900	638733	1.63%	0.331%
6	7.522	762	771	784	rBV	3677787	6018681	15.32%	3.119%
7	7.681	784	798	812	rVV	15335068	30473725	77.58%	15.791%
8	7.993	841	851	858	rBV	14799707	29317246	74.64%	15.192%
9	8.340	903	910	921	rBV	296548	481133	1.22%	0.249%
10	8.787	979	986	990	rBV	558069	914465	2.33%	0.474%
11	8.828	990	993	1004	rVB	422638	667910	1.70%	0.346%
12	9.116	1035	1042	1051	rBV	242185	386668	0.98%	0.200%
13	9.340	1070	1080	1084	rBV2	70828	119363	0.30%	0.062%
14	9.387	1084	1088	1090	rVV	70314	113817	0.29%	0.059%
15	9.428	1091	1095	1101	rVV	144073	248620	0.63%	0.129%
16	9.505	1101	1108	1114	rVV	422915	705969	1.80%	0.366%
17	9.558	1114	1117	1123	rVB2	53111	81177	0.21%	0.042%
18	9.793	1151	1157	1162	rVB2	26842	45232	0.12%	0.023%
19	10.040	1190	1199	1215	rBV	630151	1083668	2.76%	0.562%
20	10.287	1231	1241	1256	rBV	13798137	28277766	71.99%	14.653%
21	10.416	1256	1263	1270	rBV	752908	1160277	2.95%	0.601%
22	10.557	1279	1287	1300	rBV	472685	796180	2.03%	0.413%
23	10.799	1319	1328	1342	rBV	6296754	10915538	27.79%	5.656%
24	11.016	1358	1365	1373	rBV	168926	279294	0.71%	0.145%
25	11.228	1394	1401	1410	rVB	38062	72457	0.18%	0.038%
26	11.710	1470	1483	1492	rBV	135359	249094	0.63%	0.129%
27	12.416	1597	1603	1605	rBV2	118597	189815	0.48%	0.098%
28	12.451	1605	1609	1617	rVB	420038	646165	1.65%	0.335%
29	12.699	1646	1651	1658	rBV2	48697	79174	0.20%	0.041%
30	13.069	1706	1714	1726	rBV	1753058	2684917	6.84%	1.391%
31	13.298	1746	1753	1760	rBV	150005	231021	0.59%	0.120%
32	13.693	1808	1820	1825	rBV	1431291	2106857	5.36%	1.092%
33	13.740	1825	1828	1840	rVB2	123844	180438	0.46%	0.094%
34	13.969	1859	1867	1879	rBV	1693116	2431699	6.19%	1.260%
35	14.069	1879	1884	1892	rVB4	30389	47785	0.12%	0.025%
36	14.275	1913	1919	1933	rVB	1180223	1744050	4.44%	0.904%

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Filtering: 5

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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

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37	14.493	1950	1956	1967	rBV	94807	168286	0.43%	0.087%
38	14.593	1967	1973	1978	rVB	82817	115185	0.29%	0.060%
39	14.769	1998	2003	2009	rBV2	51251	84109	0.21%	0.044%
40	15.275	2082	2089	2095	rBV	2212152	3155330	8.03%	1.635%
41	15.322	2095	2097	2103	rVV	58839	86816	0.22%	0.045%
42	15.398	2104	2110	2123	rVV	646384	913648	2.33%	0.473%
43	16.057	2218	2222	2227	rVB3	27387	39489	0.10%	0.020%
44	17.028	2381	2387	2397	rBV	1627566	2265874	5.77%	1.174%
45	17.128	2397	2404	2418	rVV2	2688406	3736457	9.51%	1.936%
46	17.928	2535	2540	2547	rBV	79864	111652	0.28%	0.058%
47	18.404	2617	2621	2627	rVB2	151462	208606	0.53%	0.108%
48	18.992	2716	2721	2725	rBV2	38353	55377	0.14%	0.029%
49	19.063	2728	2733	2738	rVV	35337	53137	0.14%	0.028%
50	19.216	2755	2759	2769	rBV2	85066	122116	0.31%	0.063%
51	19.428	2789	2795	2807	rBV	3480039	4889338	12.45%	2.534%
52	19.951	2880	2884	2888	rBV4	42754	61210	0.16%	0.032%
53	21.263	3101	3107	3114	rBV	2212381	3101735	7.90%	1.607%
54	23.969	3557	3567	3587	rBV2	2680233	6324752	16.10%	3.277%
55	24.163	3591	3600	3613	rVB2	1553010	3730423	9.50%	1.933%

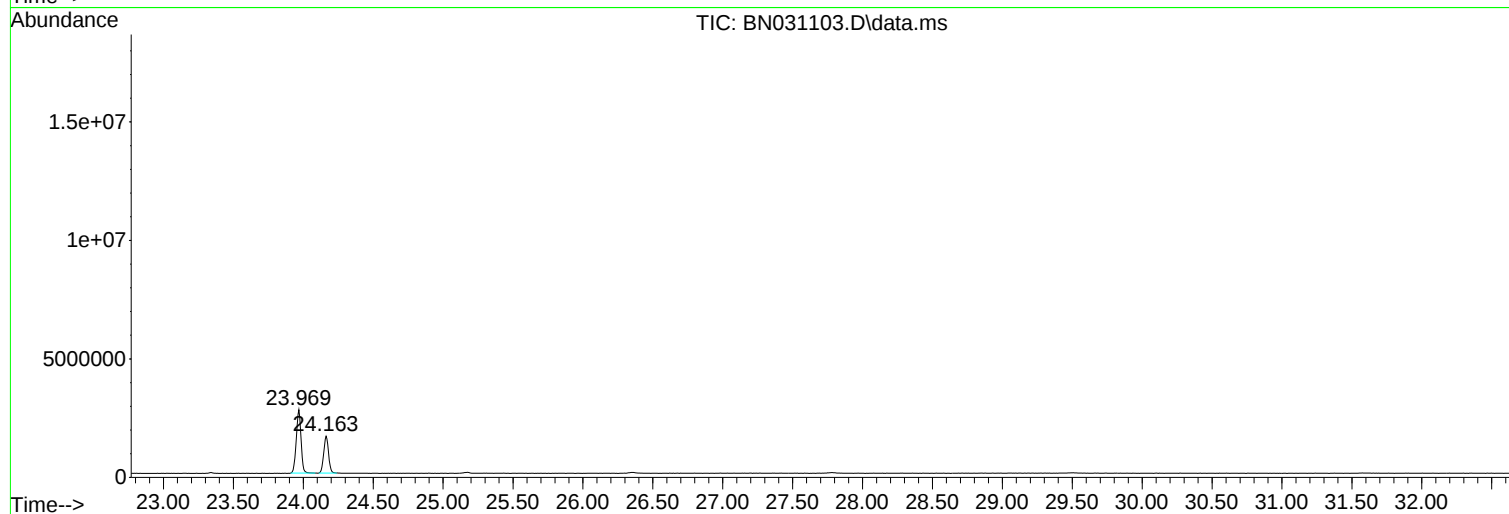
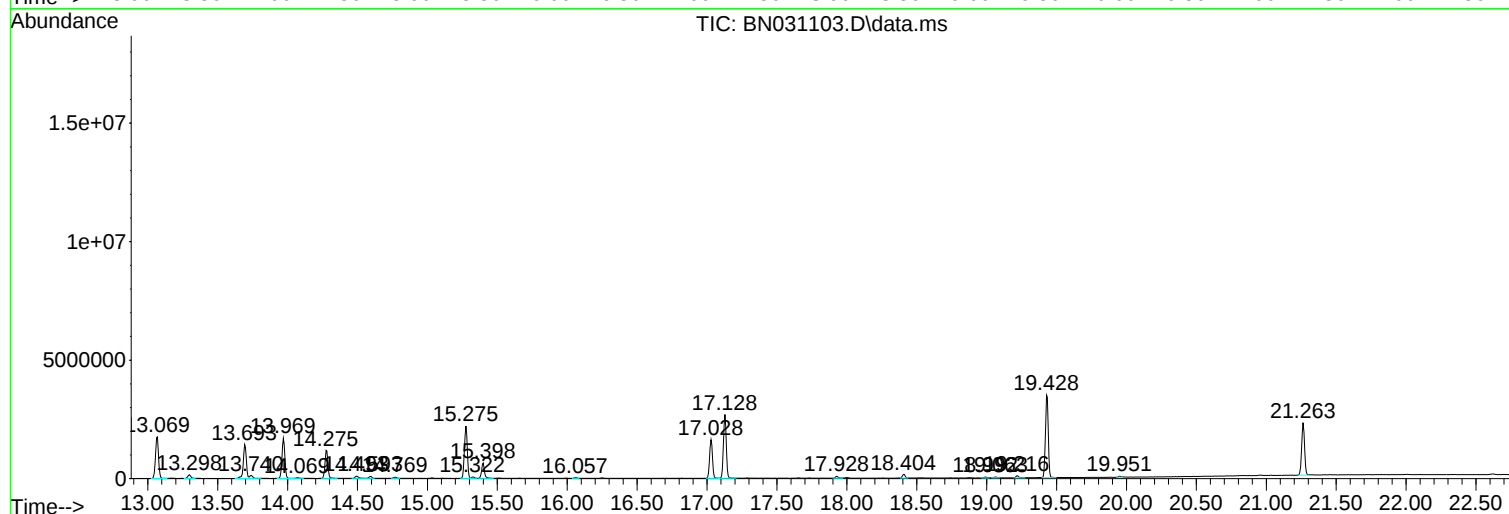
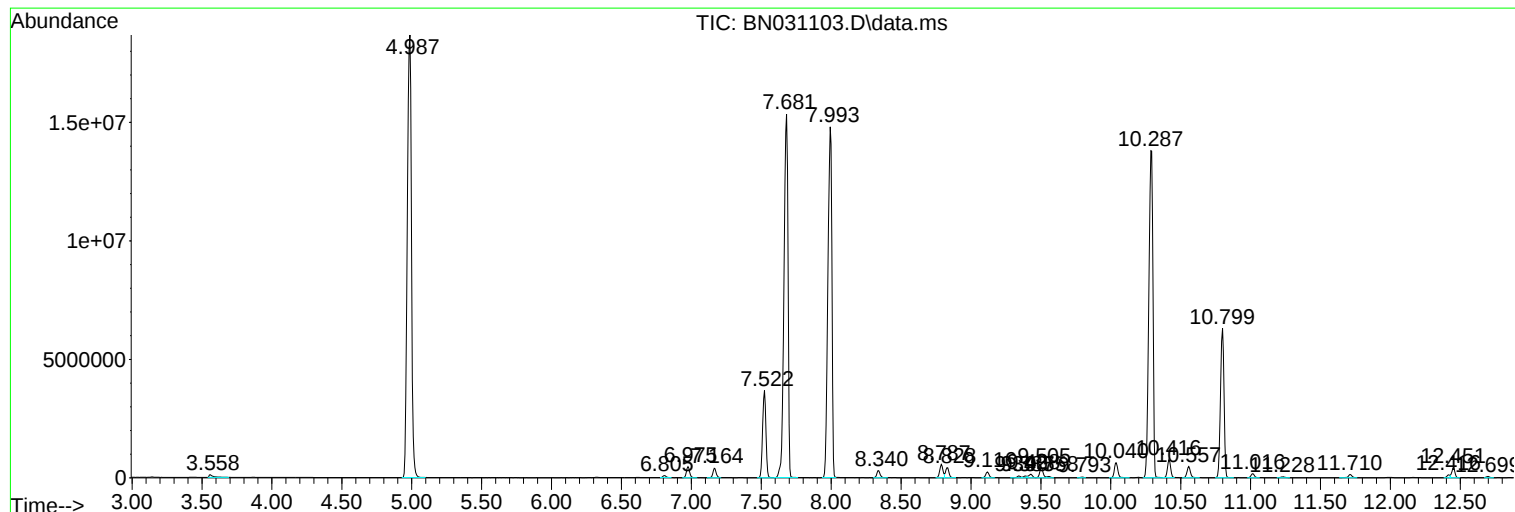
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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



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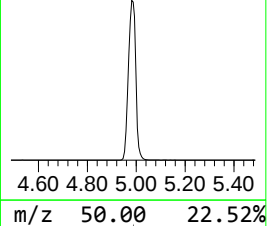
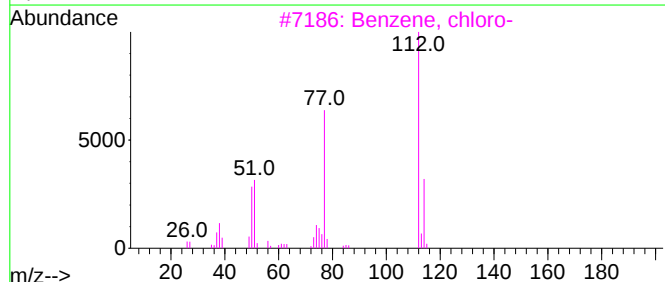
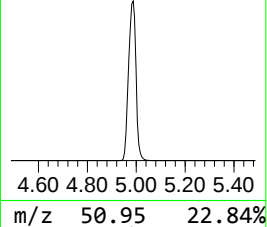
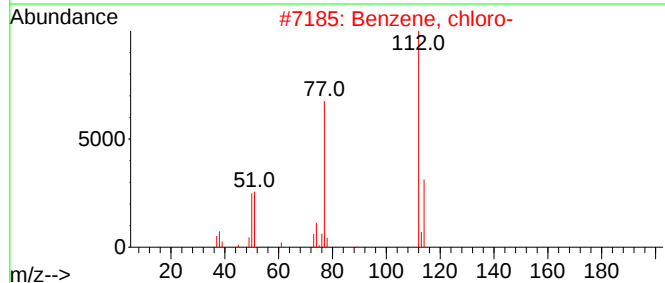
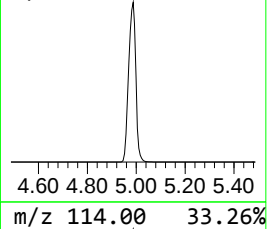
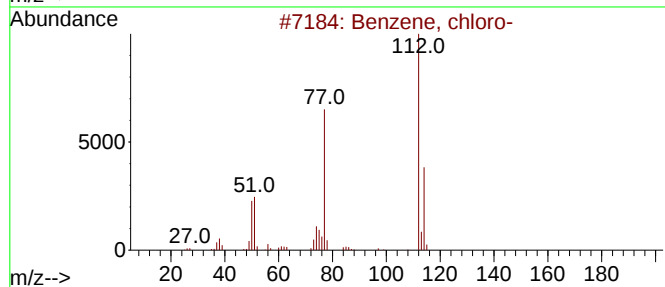
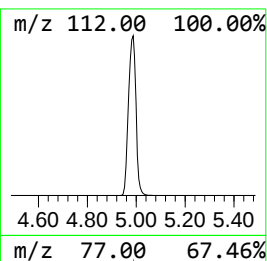
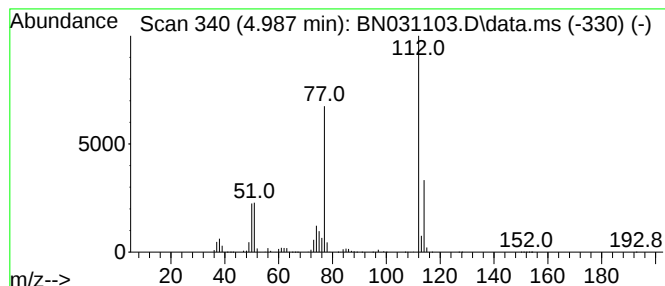
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.987	25.78 ng/ul	39280500	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	91



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ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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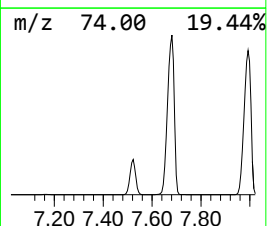
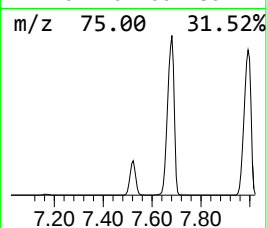
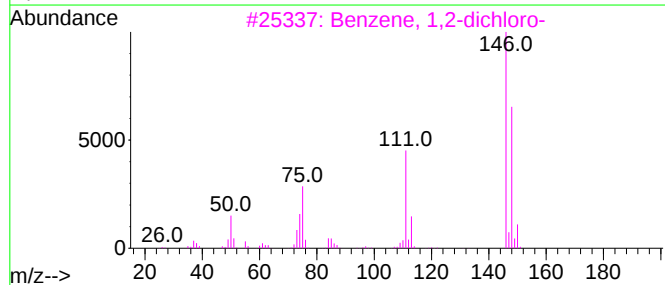
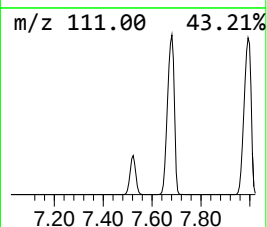
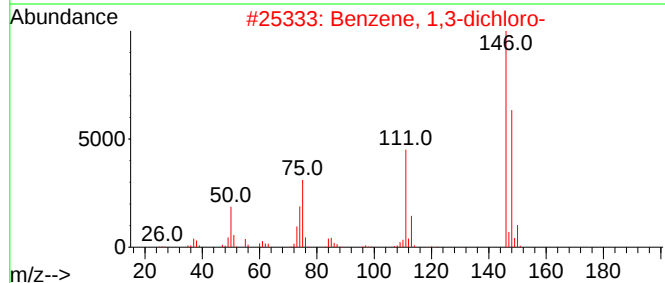
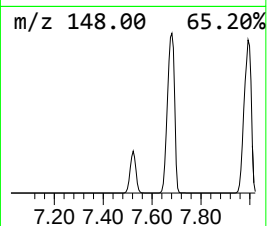
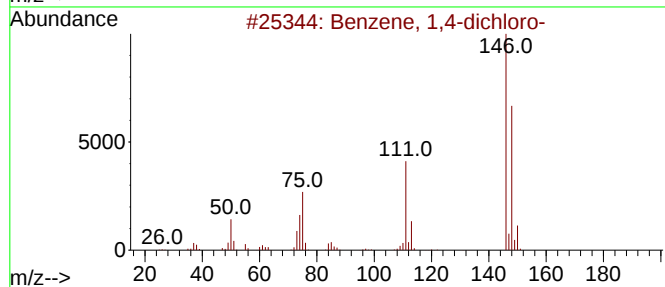
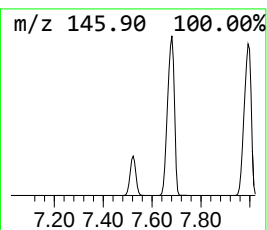
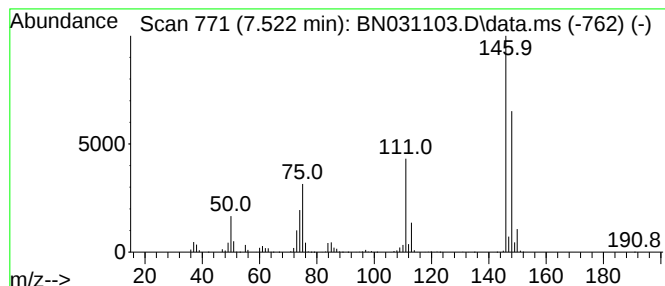
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.522	3.95 ng/ul	6018680	1,4-Dichlorobenzene-d4	7.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
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Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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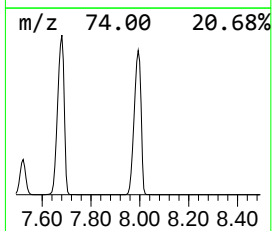
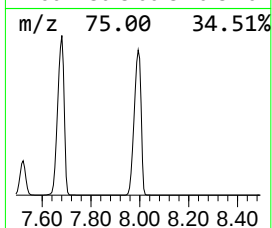
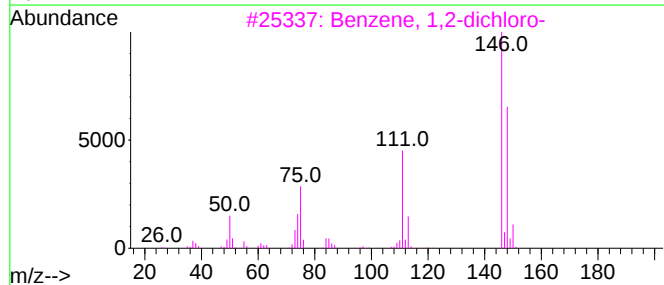
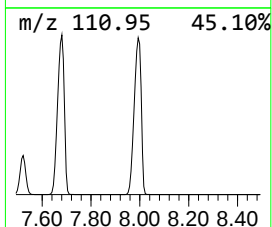
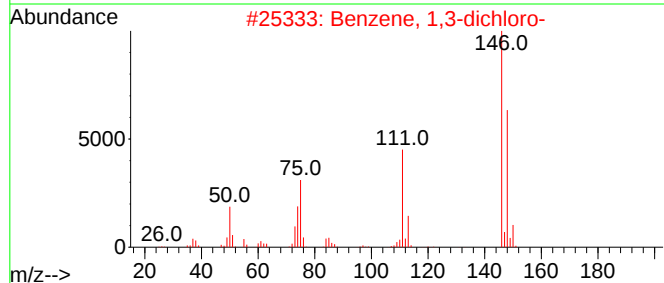
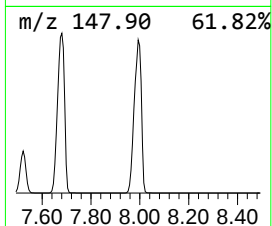
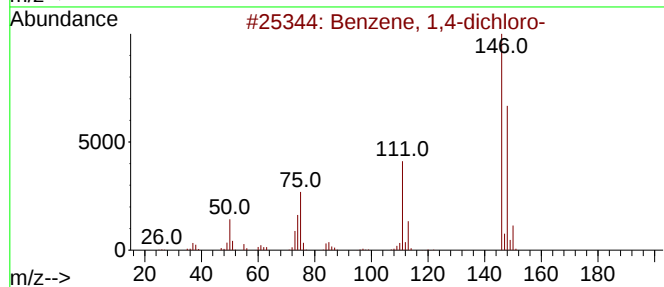
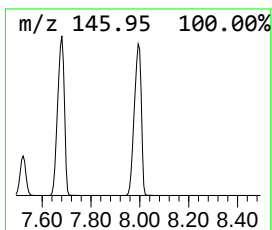
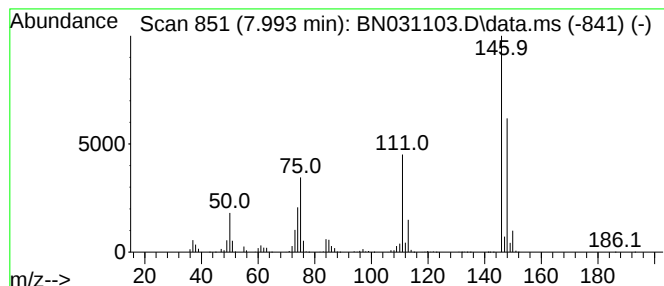
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	19.24 ng/ul	29317200	1,4-Dichlorobenzene-d4	7.634

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



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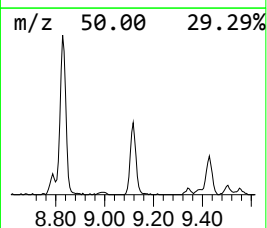
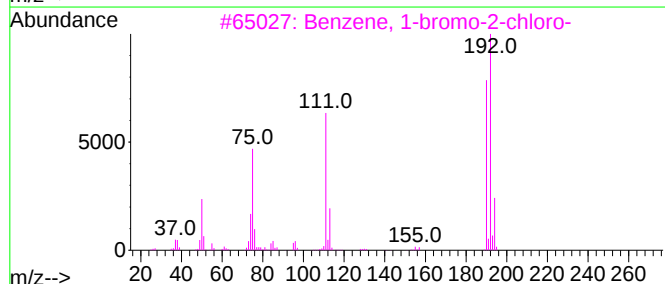
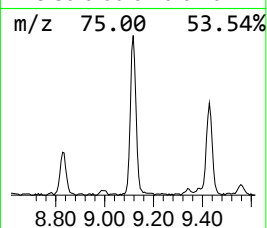
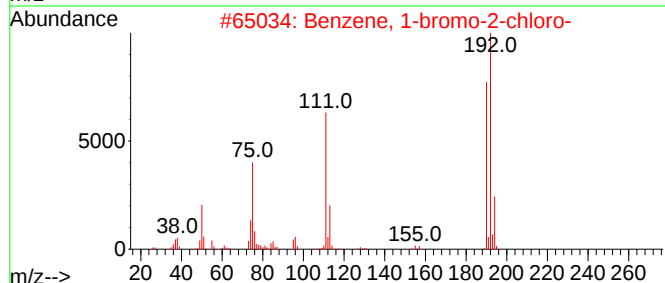
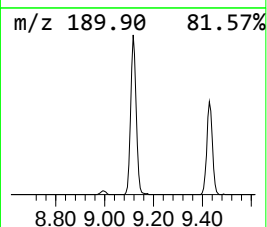
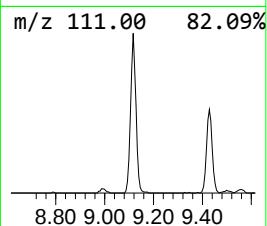
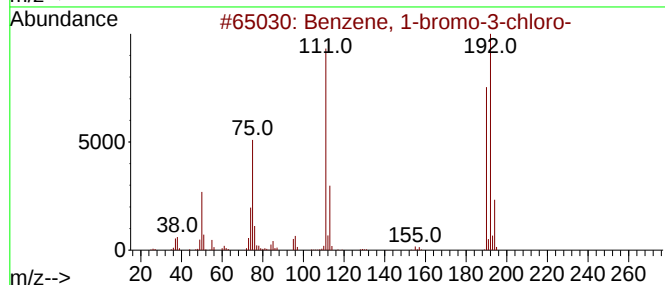
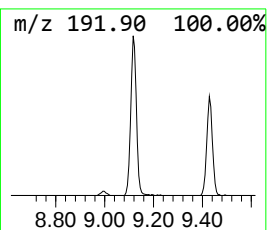
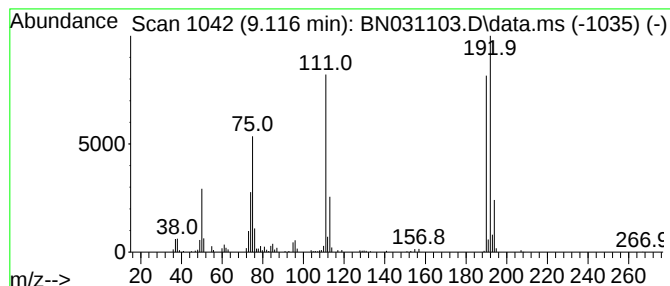
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.116	6.67 ng/ul	386668	Naphthalene-d8	10.416

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



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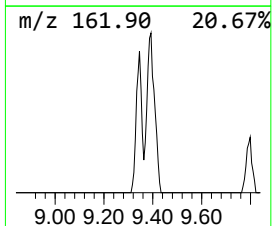
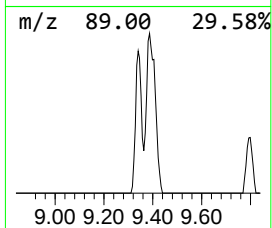
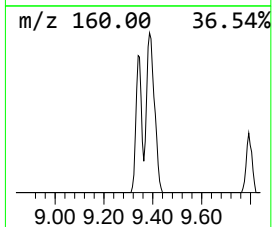
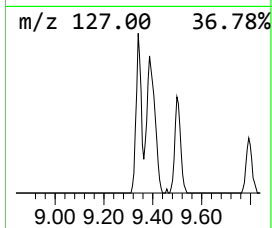
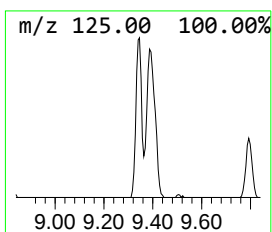
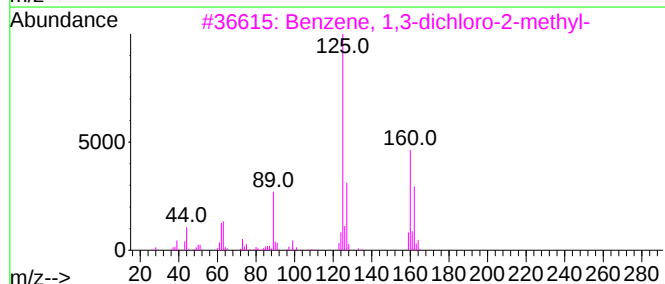
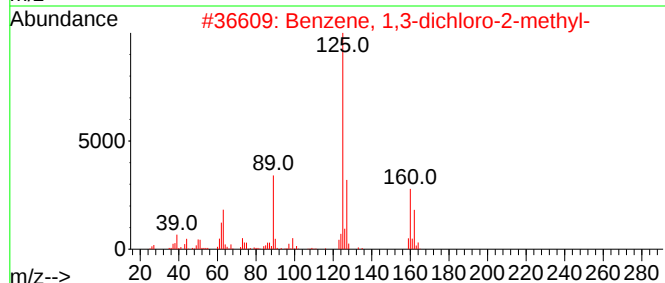
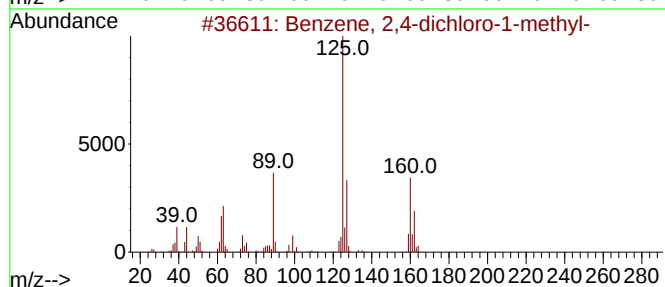
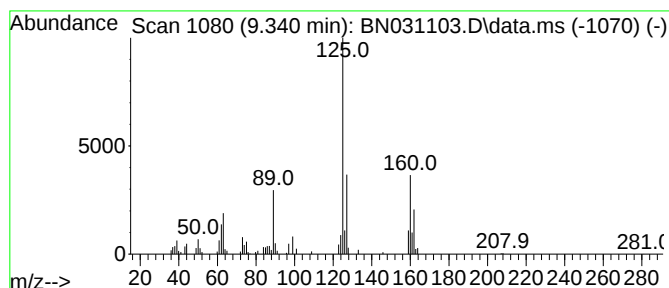
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 2,4-dichloro-1-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.340	2.06 ng/ul	119363	Naphthalene-d8	10.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2,4-dichloro-1-methyl-	160	C7H6Cl2	000095-73-8	98
2			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
3			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
4			Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97
5			Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031103.D
Acq On : 21 Apr 2024 11:12
Operator : MA/JU
Sample : P2159-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORE0

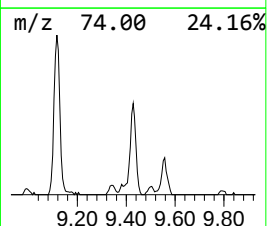
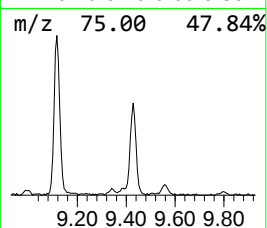
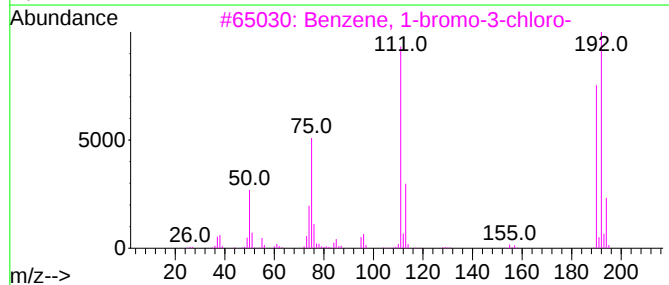
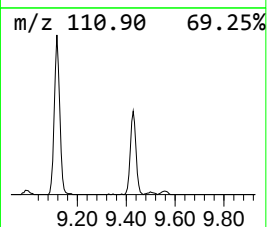
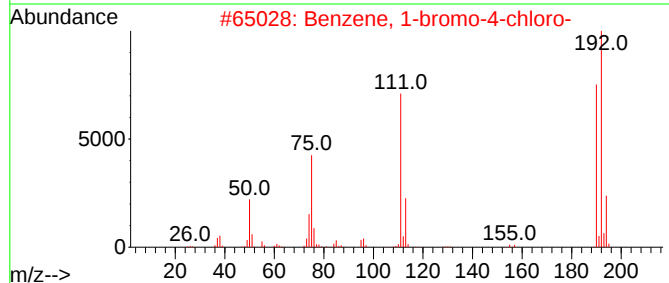
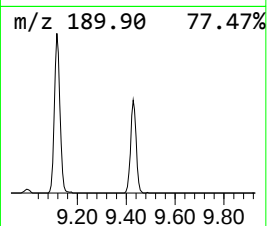
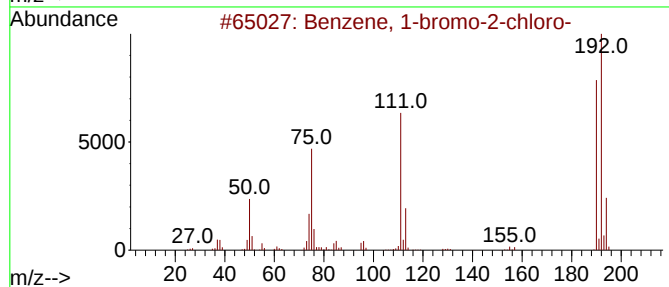
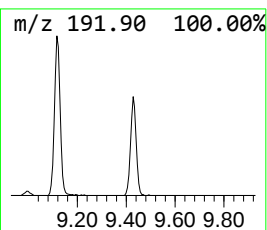
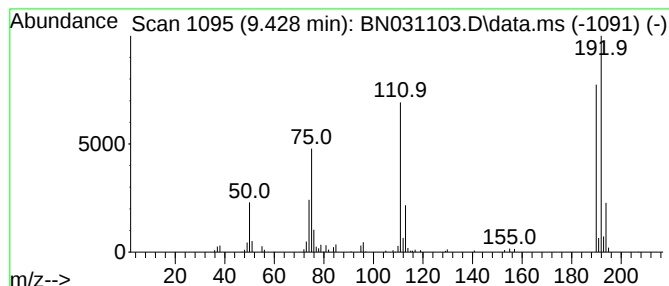
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.428	4.29 ng/ul	248620	Naphthalene-d8	10.416

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031103.D
Acq On : 21 Apr 2024 11:12
Operator : MA/JU
Sample : P2159-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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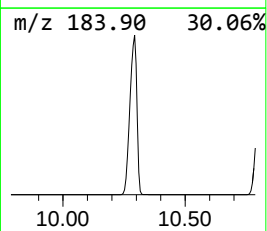
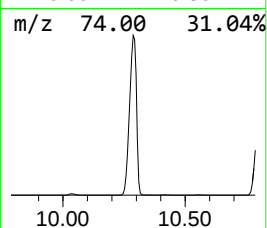
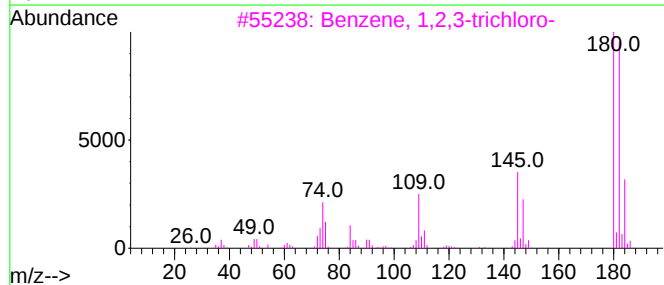
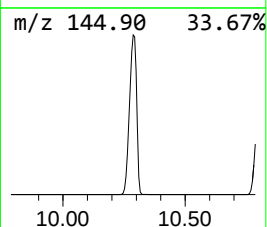
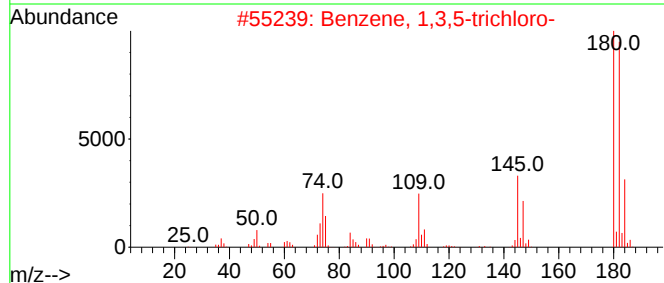
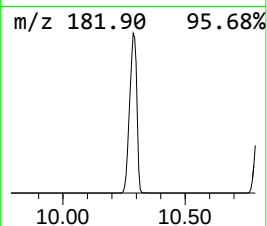
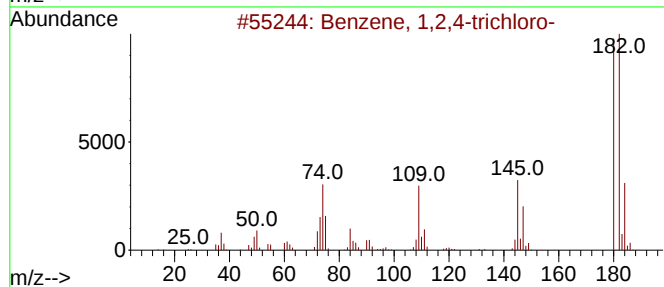
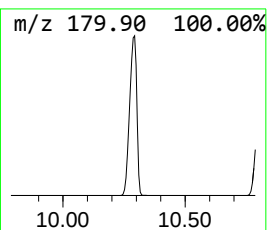
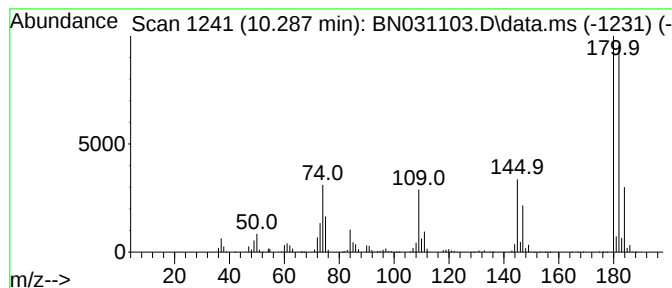
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,2,4-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.287	487.43 ng/ul	28277800	Naphthalene-d8	10.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	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	97
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031103.D
Acq On : 21 Apr 2024 11:12
Operator : MA/JU
Sample : P2159-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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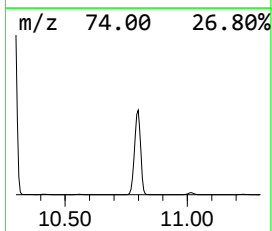
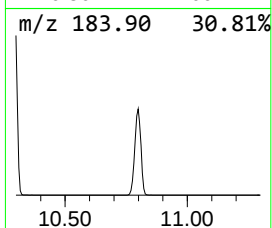
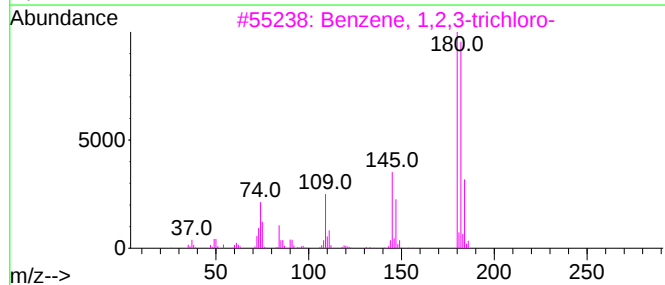
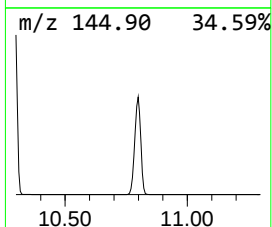
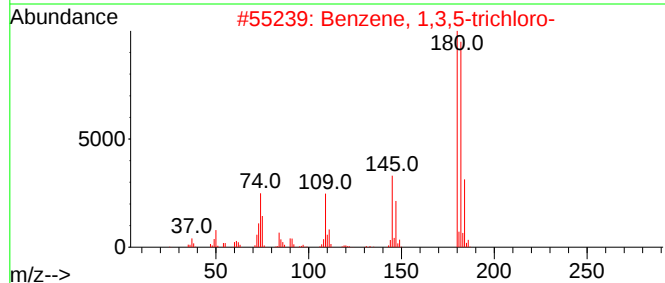
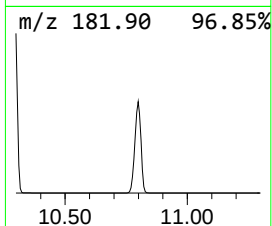
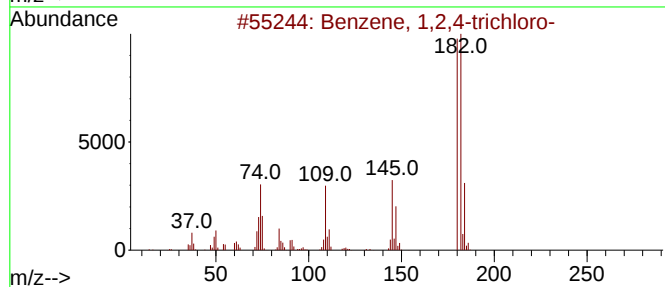
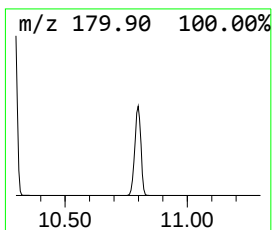
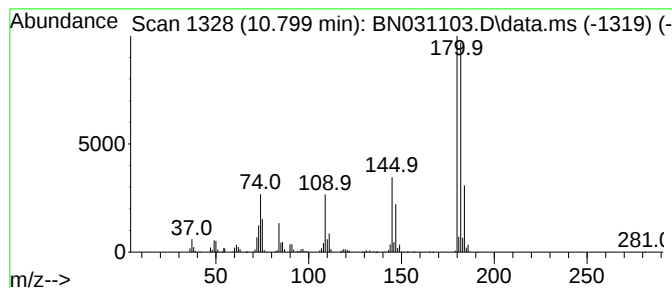
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,3,5-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.799	188.15 ng/ul	10915500	Naphthalene-d8	10.416

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031103.D
Acq On : 21 Apr 2024 11:12
Operator : MA/JU
Sample : P2159-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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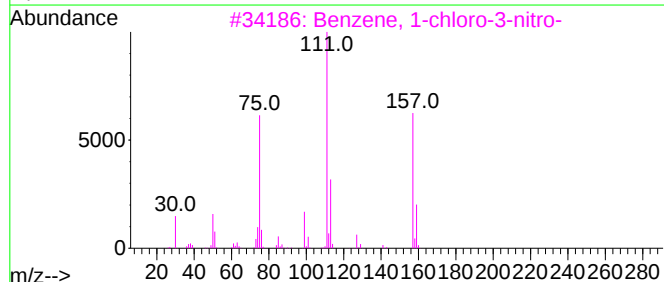
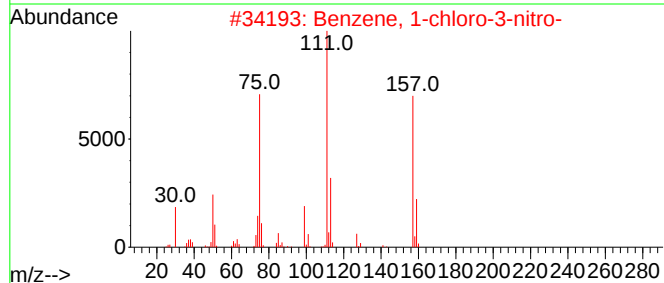
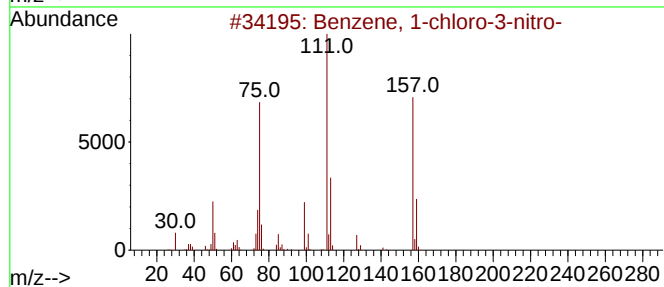
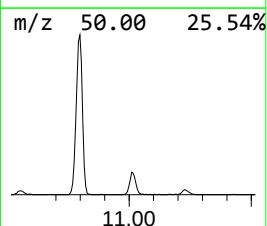
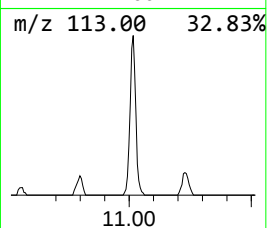
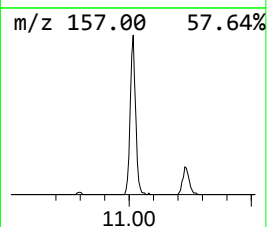
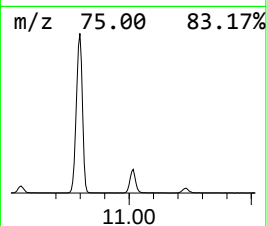
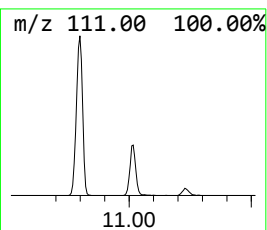
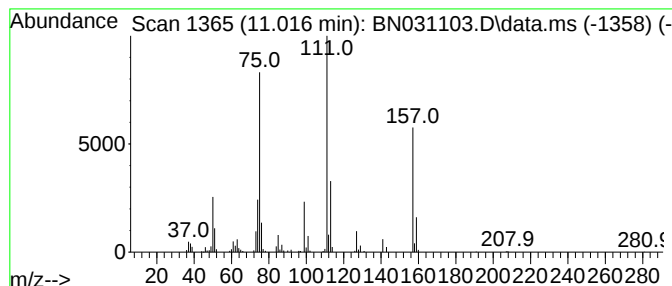
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Quant Title : SVOA CALIBRATION

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.
11.016	4.81 ng/ul	279294	Naphthalene-d8	10.416

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	95
3			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031103.D
Acq On : 21 Apr 2024 11:12
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Sample : P2159-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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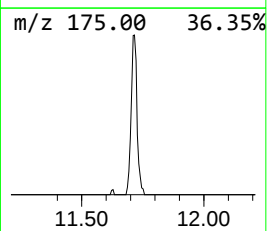
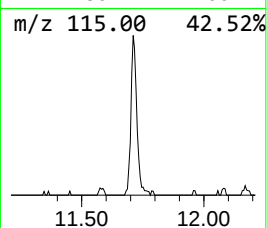
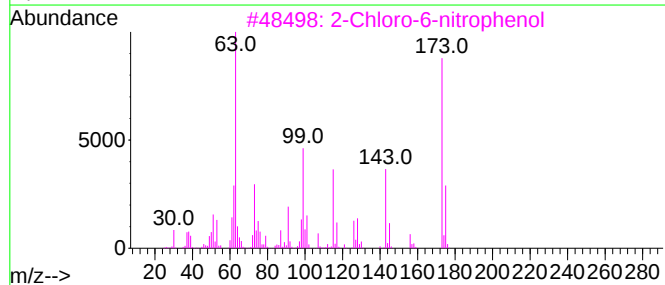
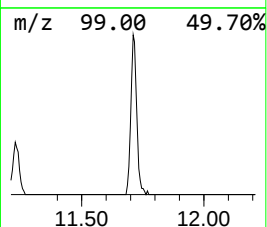
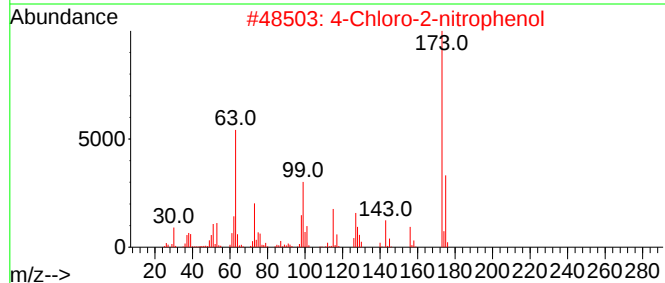
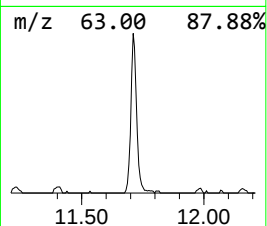
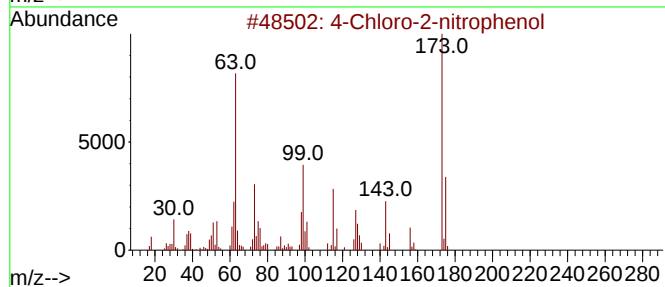
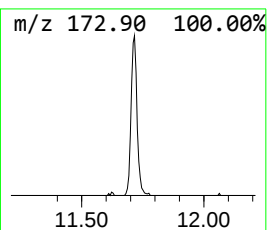
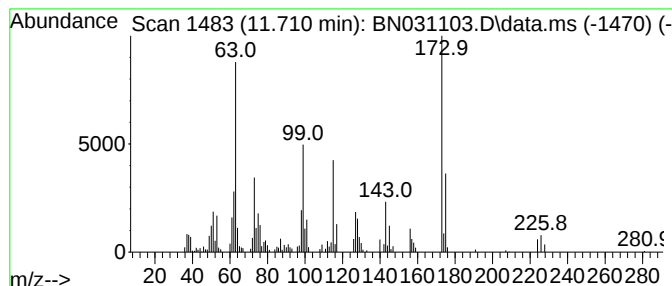
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 4-Chloro-2-nitrophenol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.710	4.29 ng/ul	249094	Naphthalene-d8	10.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	94
3			2-Chloro-6-nitrophenol	173	C6H4ClNO3	000603-86-1	94
4			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	90
5			Phenol, 2-chloro-4-nitro-	173	C6H4ClNO3	000619-08-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
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Acq On : 21 Apr 2024 11:12
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Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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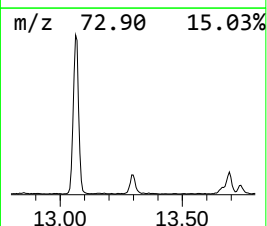
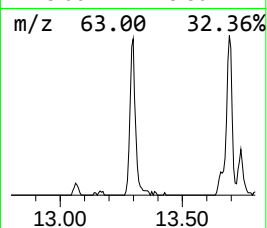
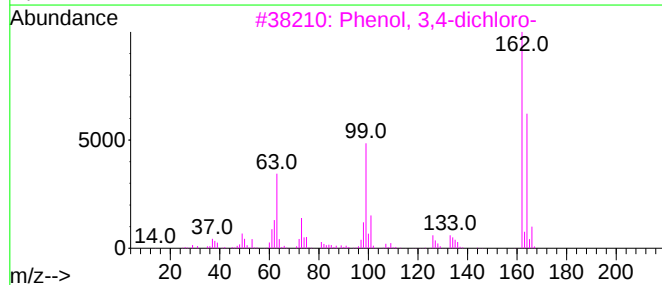
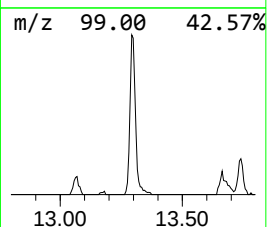
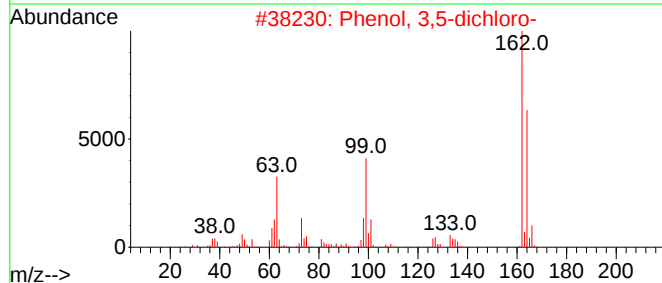
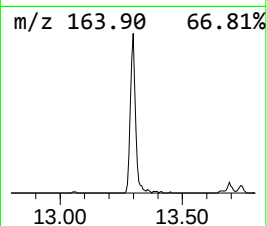
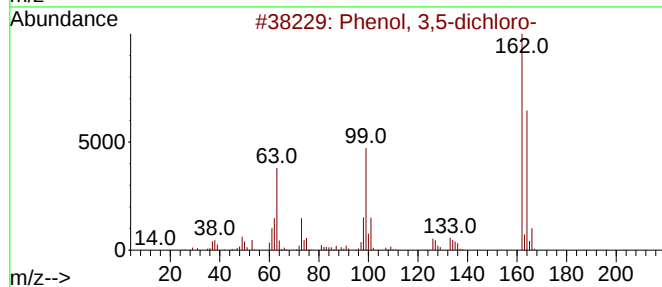
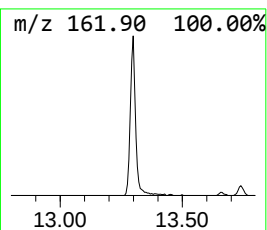
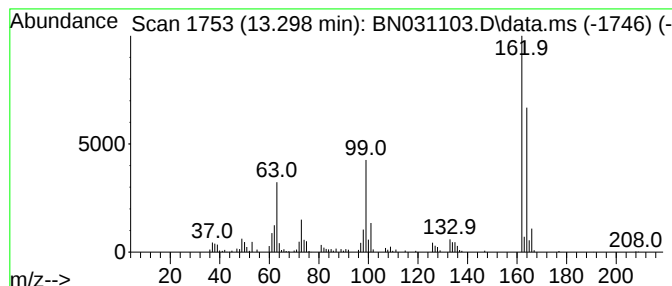
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenol, 3,5-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.299	2.65 ng/ul	231021	Acenaphthene-d10	14.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
2			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
4			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
5			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031103.D
Acq On : 21 Apr 2024 11:12
Operator : MA/JU
Sample : P2159-19
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORE0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.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-	4.987	25.8	ng/ul	39280500	1	7.634	30473700	20.0
Benzene, 1,4-di...	7.522	4.0	ng/ul	6018680	1	7.634	30473700	20.0
Benzene, 1,3-di...	7.993	19.2	ng/ul	29317200	1	7.634	30473700	20.0
Benzene, 1-brom...	9.116	6.7	ng/ul	386668	2	10.416	1160280	20.0
Benzene, 2,4-di...	9.340	2.1	ng/ul	119363	2	10.416	1160280	20.0
Benzene, 1-brom...	9.428	4.3	ng/ul	248620	2	10.416	1160280	20.0
Benzene, 1,2,4-...	10.287	487.4	ng/ul	28277800	2	10.416	1160280	20.0
Benzene, 1,3,5-...	10.799	188.2	ng/ul	10915500	2	10.416	1160280	20.0
Benzene, 1-chlo...	11.016	4.8	ng/ul	279294	2	10.416	1160280	20.0
4-Chloro-2-nitr...	11.710	4.3	ng/ul	249094	2	10.416	1160280	20.0
Phenol, 3,5-dic...	13.299	2.6	ng/ul	231021	3	14.275	1744050	20.0