

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042523\
 Data File : BN025126.D
 Acq On : 25 Apr 2023 12:18
 Operator : CG/JU
 Sample : 02400-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 COML6

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-BN041023.M
 Title : SVOA CALIBRATION

Signal : TIC: BN025126.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.293	14	18	29	rVB	76332	112718	1.06%	0.079%
2	3.723	88	91	114	rBV	414386	749596	7.04%	0.528%
3	3.952	128	130	137	rVB6	11949	18146	0.17%	0.013%
4	4.052	143	147	154	rVB2	27752	45728	0.43%	0.032%
5	4.993	302	307	310	rBV4	7111	13368	0.13%	0.009%
6	5.199	335	342	356	rBV	5651428	9006030	84.54%	6.349%
7	6.587	573	578	583	rVB2	15802	25522	0.24%	0.018%
8	6.993	644	647	654	rBV6	9716	16690	0.16%	0.012%
9	7.075	654	661	673	rVB	393447	650417	6.11%	0.459%
10	7.240	683	689	702	rBV	1419696	2272701	21.33%	1.602%
11	7.440	716	723	734	rBV	1394156	2192709	20.58%	1.546%
12	7.528	734	738	740	rBV2	9504	11584	0.11%	0.008%
13	7.799	777	784	795	rVB	2275415	3600333	33.80%	2.538%
14	7.916	795	804	805	rBV	1384210	1982510	18.61%	1.398%
15	7.952	805	810	820	rVB	6150982	9896806	92.90%	6.977%
16	8.269	855	864	876	rBV	6701246	10653199	100.00%	7.510%
17	8.622	917	924	935	rBV	1099307	1769099	16.61%	1.247%
18	9.081	995	1002	1007	rBV	1826522	2956049	27.75%	2.084%
19	9.122	1007	1009	1017	rVB	175232	249271	2.34%	0.176%
20	9.240	1026	1029	1034	rVB6	7265	11673	0.11%	0.008%
21	9.316	1036	1042	1048	rVB3	5963	15470	0.15%	0.011%
22	9.416	1055	1059	1066	rBV4	14301	27149	0.25%	0.019%
23	9.487	1066	1071	1076	rVV	14352	28913	0.27%	0.020%
24	9.805	1118	1125	1136	rBV	1472282	2394129	22.47%	1.688%
25	10.346	1209	1217	1228	rBV	1945816	3244162	30.45%	2.287%
26	10.587	1250	1258	1274	rBV	1504653	2529075	23.74%	1.783%
27	10.722	1274	1281	1293	rVB	1934224	3317536	31.14%	2.339%
28	10.863	1297	1305	1323	rBV	1850493	3178503	29.84%	2.241%
29	11.110	1340	1347	1357	rVB	628221	1045272	9.81%	0.737%
30	11.334	1379	1385	1392	rBV	79453	143555	1.35%	0.101%
31	11.552	1412	1422	1433	rVB7	24159	69036	0.65%	0.049%
32	12.140	1515	1522	1527	rVB8	10502	20589	0.19%	0.015%
33	12.310	1546	1551	1557	rBV	34285	55139	0.52%	0.039%
34	12.728	1617	1622	1623	rBV5	11871	16474	0.15%	0.012%
35	12.757	1624	1627	1632	rVB4	12462	19061	0.18%	0.013%
36	13.087	1676	1683	1687	rBV9	6403	12516	0.12%	0.009%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
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37	13.157	1691	1695	1702	rVB10	6084	11420	0.11%	0.008%
38	13.369	1724	1731	1739	rBV2	66169	141581	1.33%	0.100%
39	13.451	1740	1745	1750	rVB4	8617	14733	0.14%	0.010%
40	13.528	1750	1758	1761	rBV8	12578	27083	0.25%	0.019%
41	13.581	1763	1767	1773	rVB6	16326	28226	0.26%	0.020%
42	13.969	1825	1833	1841	rBV	4104091	5572936	52.31%	3.929%
43	14.087	1849	1853	1859	rVB8	13696	21848	0.21%	0.015%
44	14.151	1859	1864	1868	rBV	31667	46664	0.44%	0.033%
45	14.251	1875	1881	1890	rBV	4249372	6261558	58.78%	4.414%
46	14.446	1908	1914	1918	rVB8	11504	22776	0.21%	0.016%
47	14.563	1925	1934	1941	rBV	3054305	4459132	41.86%	3.144%
48	14.663	1947	1951	1955	rVB7	7651	12285	0.12%	0.009%
49	14.757	1961	1967	1983	rBV	317230	553462	5.20%	0.390%
50	14.881	1983	1988	1991	rVB6	9675	14369	0.13%	0.010%
51	14.969	1999	2003	2010	rVB7	16025	28868	0.27%	0.020%
52	15.551	2095	2102	2111	rBV	5604113	7944189	74.57%	5.600%
53	15.675	2115	2123	2135	rVV	2150730	2939497	27.59%	2.072%
54	15.793	2138	2143	2148	rVB4	28811	44648	0.42%	0.031%
55	15.916	2159	2164	2172	rVB	1442703	1923760	18.06%	1.356%
56	16.481	2255	2260	2266	rBV	229539	298812	2.80%	0.211%
57	16.987	2342	2346	2350	rBV5	12154	16534	0.16%	0.012%
58	17.310	2394	2401	2410	rBV	3770283	5282055	49.58%	3.724%
59	17.410	2411	2418	2433	rVV	6382156	9035966	84.82%	6.370%
60	17.587	2445	2448	2455	rVB7	13859	21435	0.20%	0.015%
61	17.692	2462	2466	2469	rBV2	9423	13250	0.12%	0.009%
62	17.975	2510	2514	2518	rBV5	12254	15604	0.15%	0.011%
63	18.022	2518	2522	2527	rVV7	13888	20641	0.19%	0.015%
64	18.198	2547	2552	2561	rBV	296728	431935	4.05%	0.305%
65	18.281	2562	2566	2572	rVB4	25783	59478	0.56%	0.042%
66	18.492	2599	2602	2607	rBV5	17758	30056	0.28%	0.021%
67	19.051	2694	2697	2700	rBV5	8242	11864	0.11%	0.008%
68	19.116	2703	2708	2712	rBV8	18115	42228	0.40%	0.030%
69	19.263	2729	2733	2739	rBV2	87017	119101	1.12%	0.084%
70	19.334	2741	2745	2751	rVV	82889	130669	1.23%	0.092%
71	19.475	2764	2769	2774	rBV	137701	191629	1.80%	0.135%
72	19.622	2791	2794	2799	rVB2	43502	51638	0.48%	0.036%
73	19.704	2802	2808	2821	rVV	8066747	10620866	99.70%	7.487%
74	20.675	2970	2973	2976	rBV4	34808	37126	0.35%	0.026%
75	20.733	2981	2983	2986	rVB3	27624	26771	0.25%	0.019%
76	21.204	3058	3063	3072	rBV	423214	711319	6.68%	0.501%

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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

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Title : SVOA CALIBRATION

77	21.504	3109	3114	3121	rBV	4364666	5386243	50.56%	3.797%
78	22.116	3215	3218	3225	rVB2	81634	120417	1.13%	0.085%
79	22.616	3300	3303	3308	rVB	166513	212062	1.99%	0.149%
80	23.174	3394	3398	3403	rVB	229001	318233	2.99%	0.224%
81	23.774	3492	3500	3517	rBV2	4889327	10131692	95.10%	7.143%
82	23.927	3519	3526	3537	rVB2	2470636	5037050	47.28%	3.551%
83	24.527	3623	3628	3635	rVB	268919	539659	5.07%	0.380%
84	25.374	3768	3772	3779	rVB	237254	515302	4.84%	0.363%

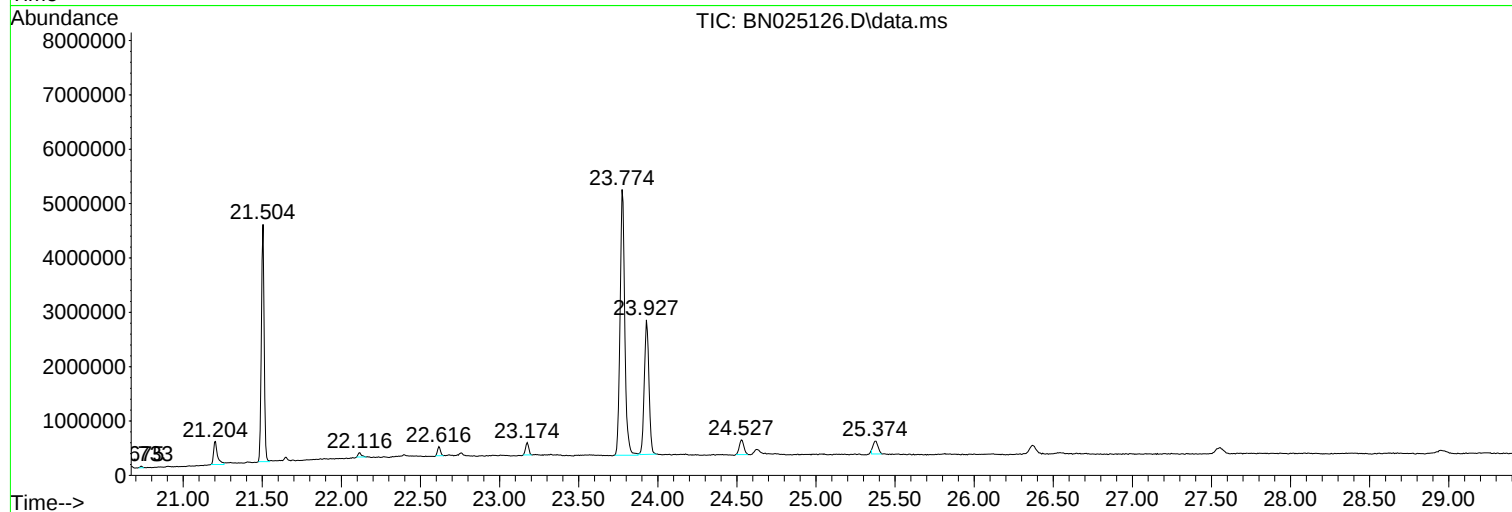
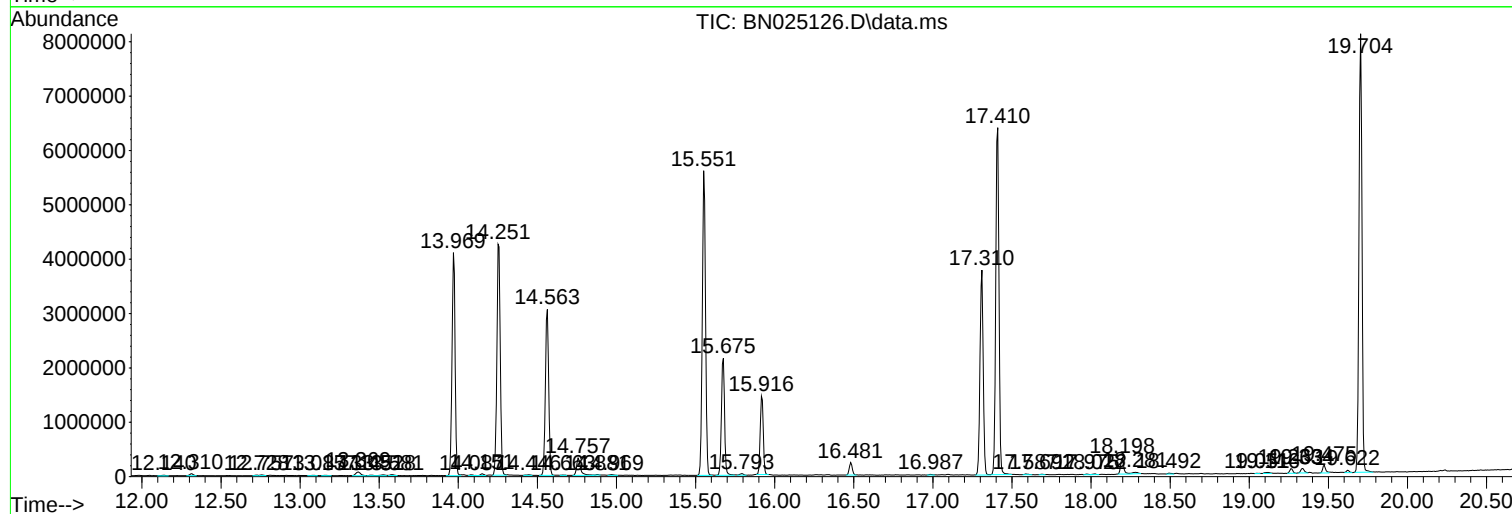
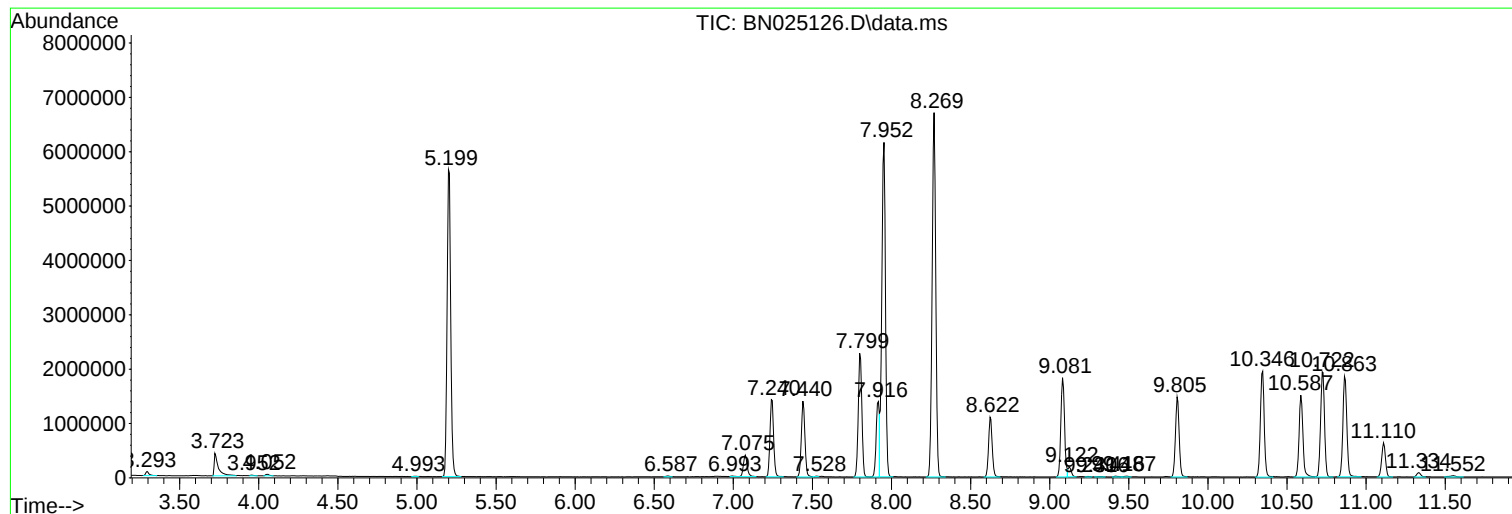
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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



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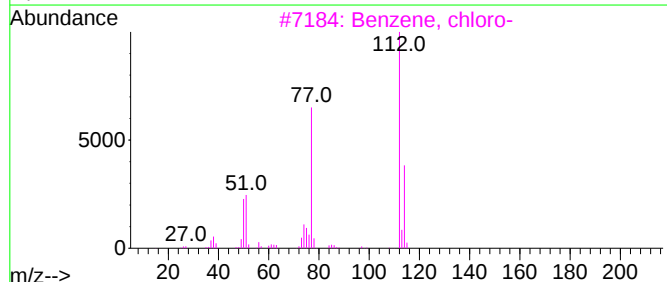
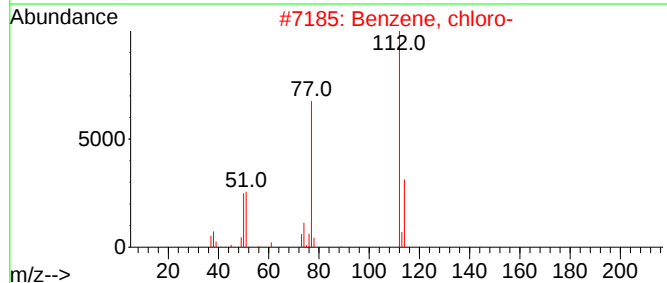
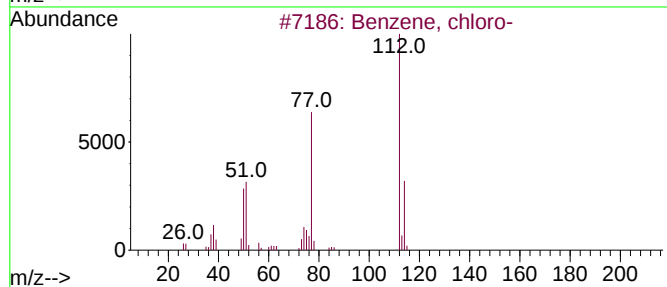
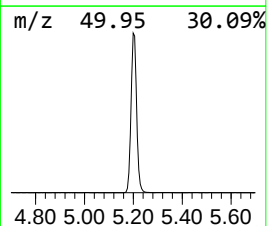
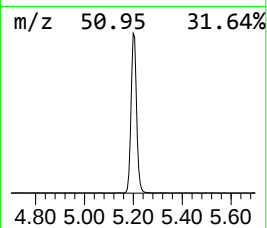
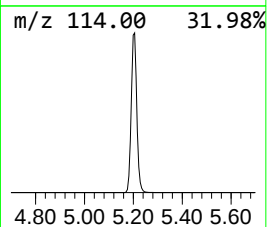
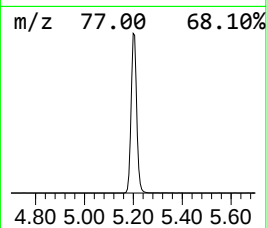
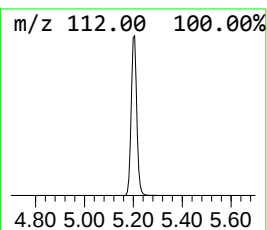
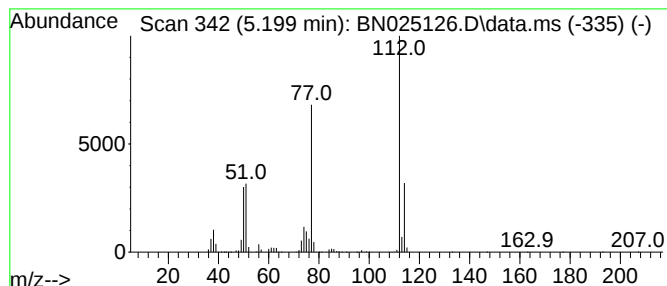
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.199	90.85 ng/ul	9006030	1,4-Dichlorobenzene-d4	7.916

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	95
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	94



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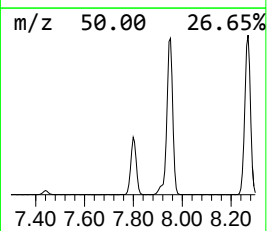
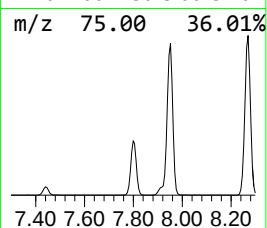
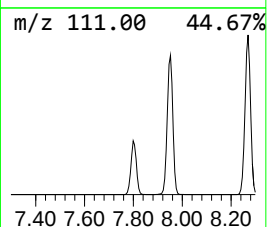
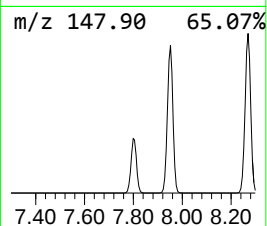
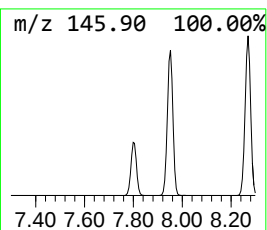
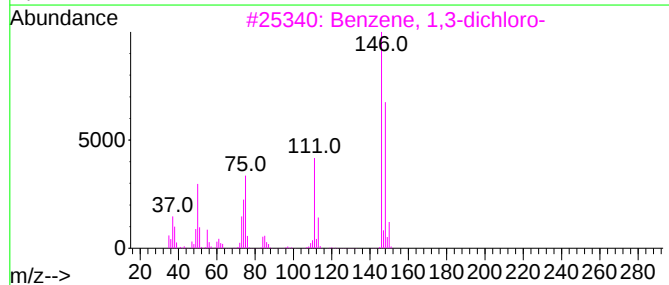
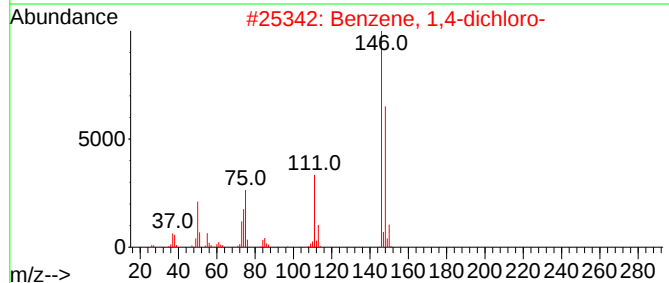
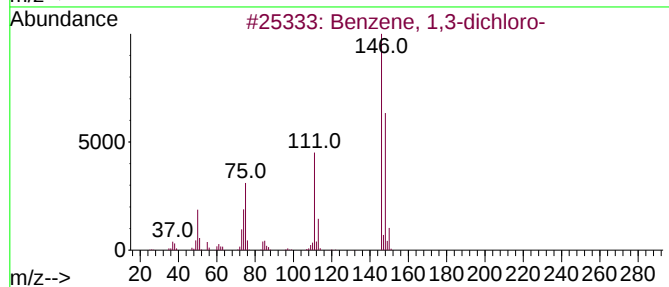
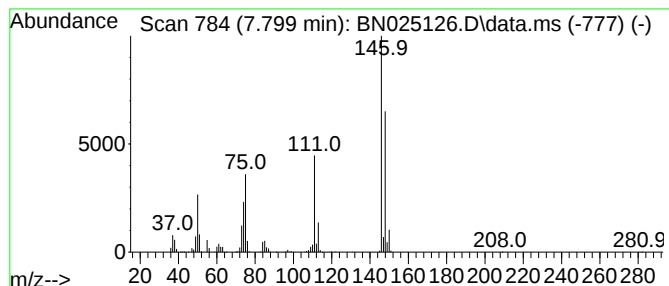
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.799	36.32 ng/ul	3600330	1,4-Dichlorobenzene-d4	7.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
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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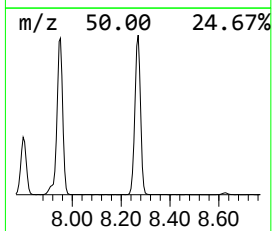
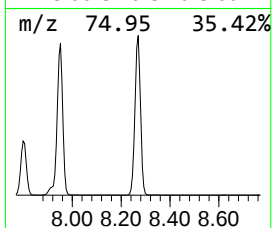
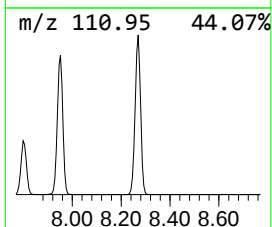
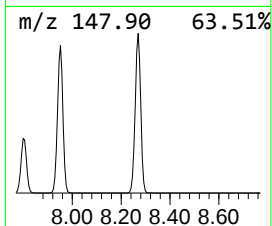
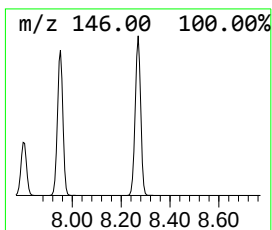
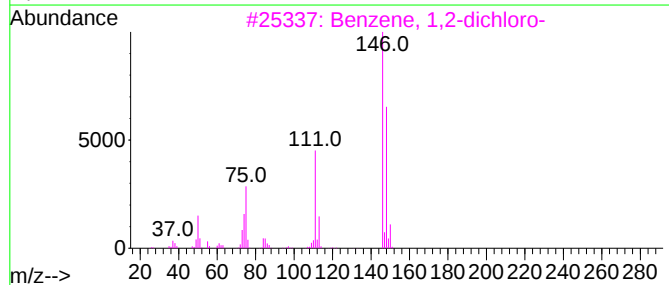
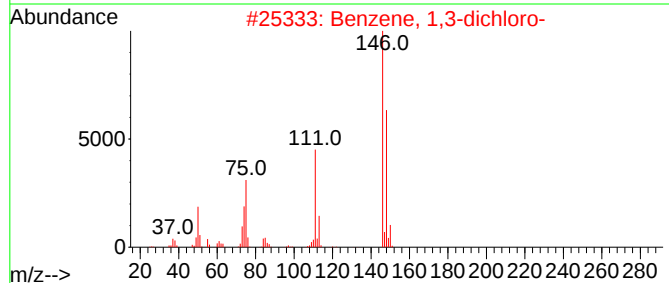
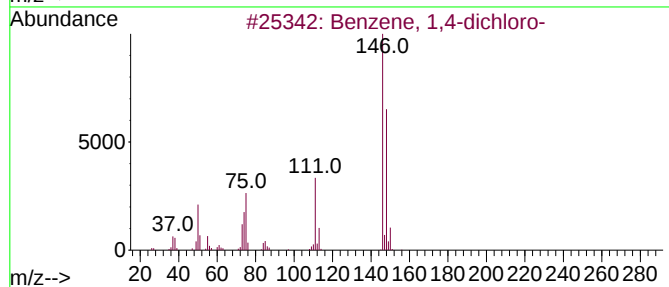
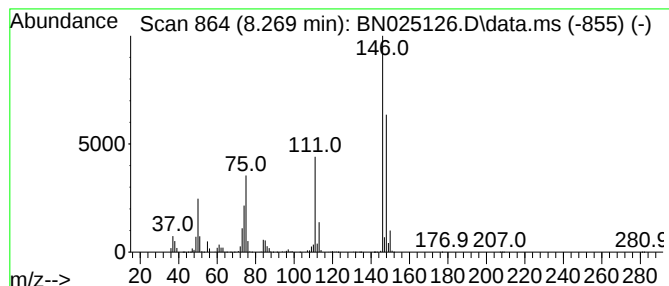
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.269	107.47 ng/ul	10653200	1,4-Dichlorobenzene-d4	7.916

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	97



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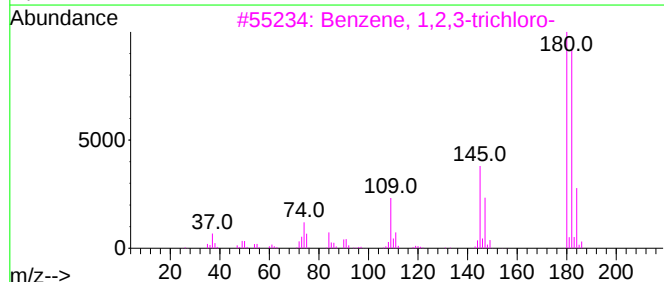
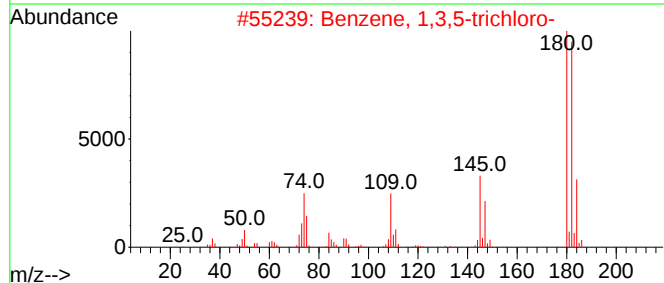
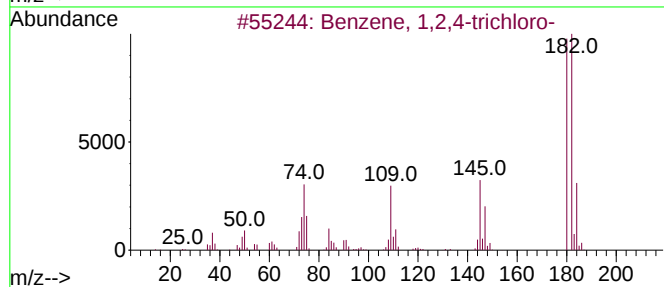
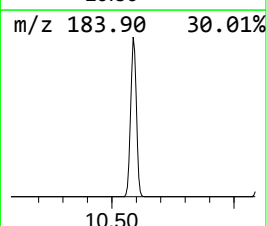
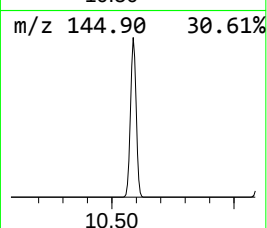
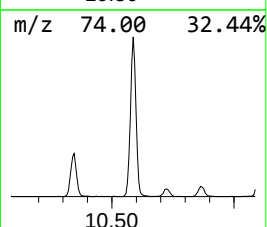
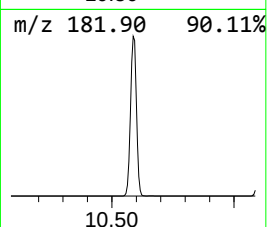
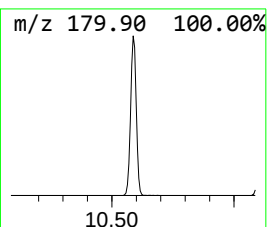
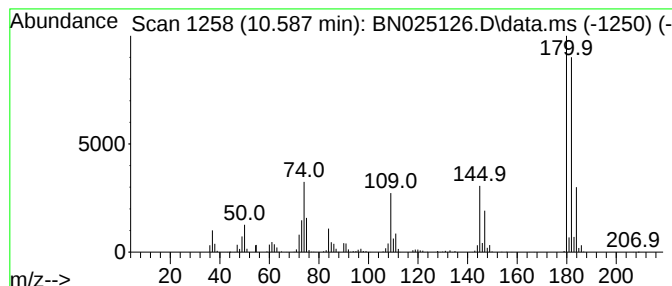
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.587	15.25 ng/ul	2529080	Naphthalene-d8	10.728

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\BN042523\
Data File : BN025126.D
Acq On : 25 Apr 2023 12:18
Operator : CG/JU
Sample : 02400-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
COML6

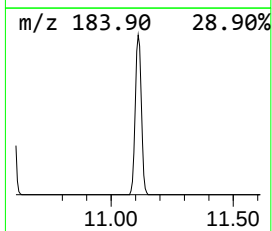
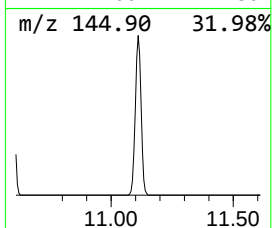
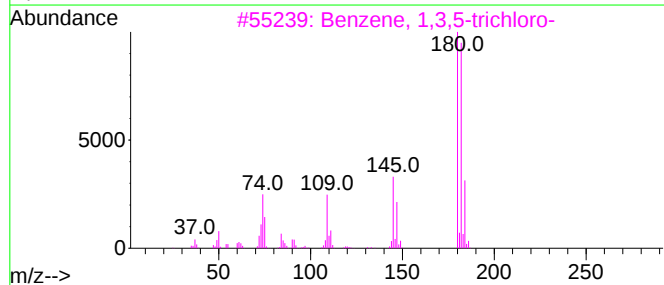
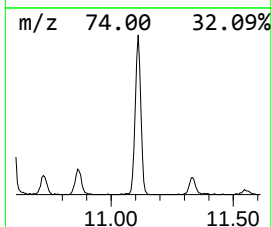
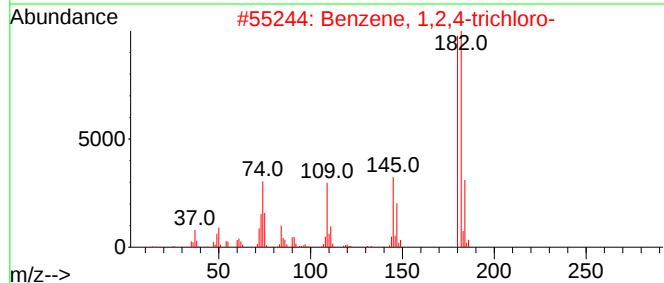
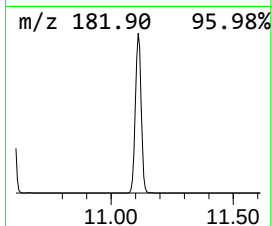
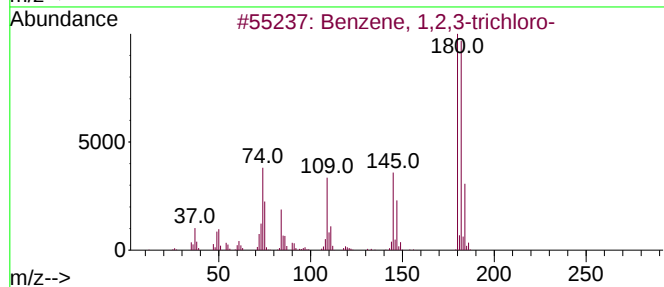
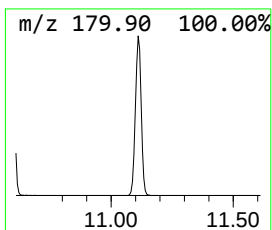
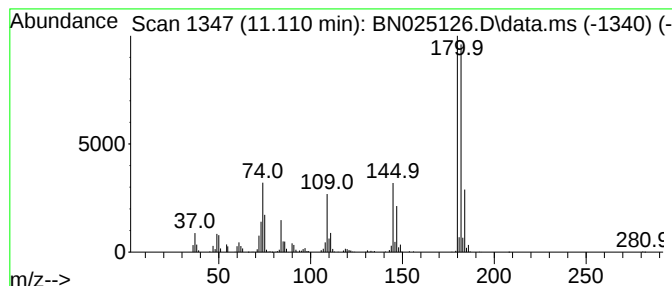
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2,3-trichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.110	6.30 ng/ul	1045270	Naphthalene-d8	10.728

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	98
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, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042523\
Data File : BN025126.D
Acq On : 25 Apr 2023 12:18
Operator : CG/JU
Sample : 02400-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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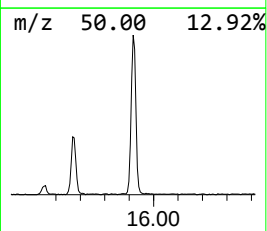
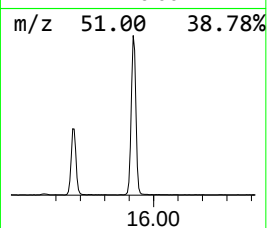
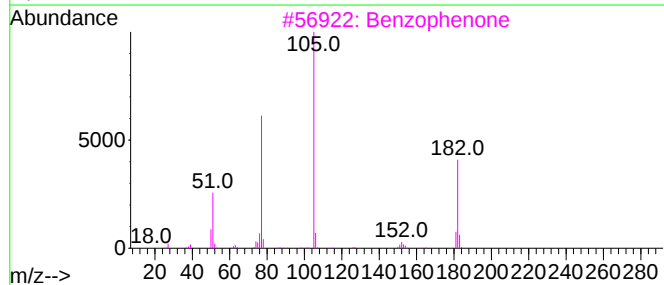
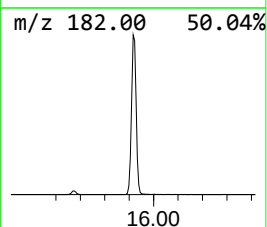
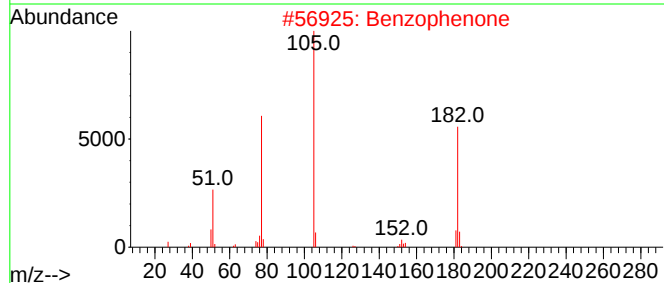
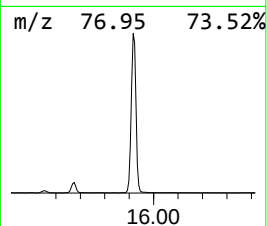
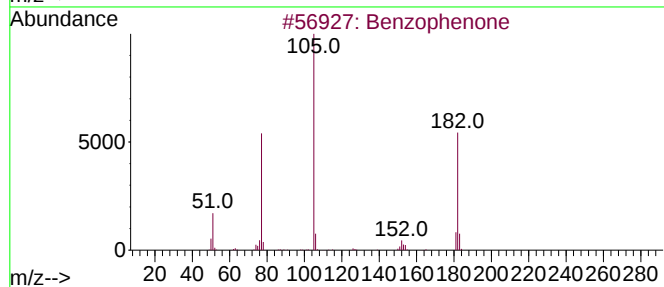
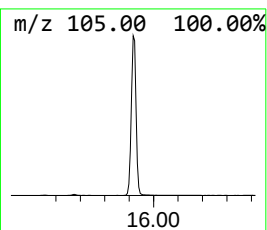
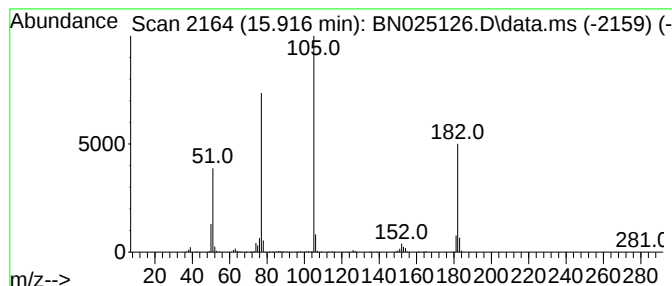
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.916	8.63 ng/ul	1923760	Acenaphthene-d10	14.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042523\
Data File : BN025126.D
Acq On : 25 Apr 2023 12:18
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Sample : 02400-10
Misc :
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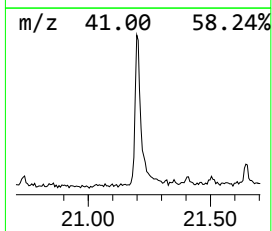
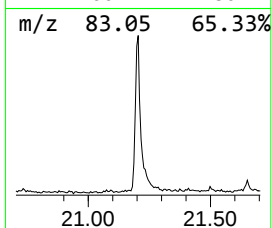
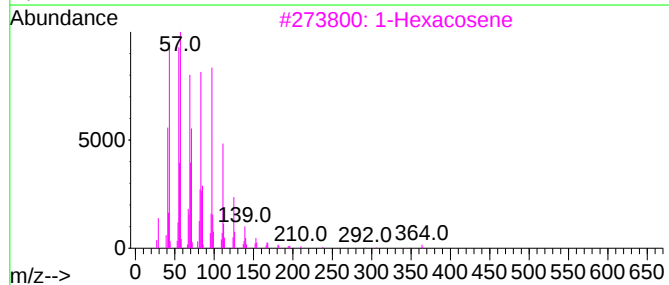
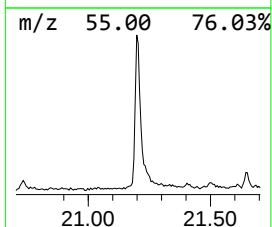
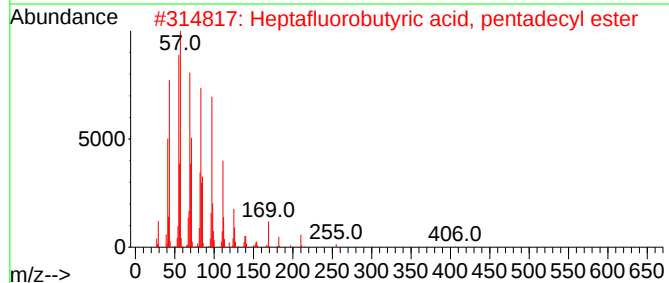
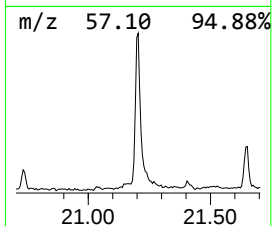
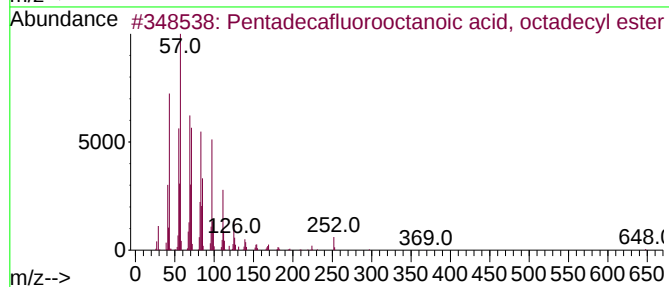
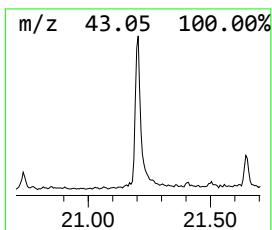
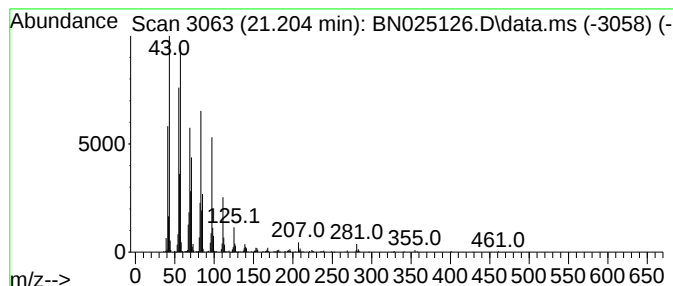
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Pentadecafluorooctanoic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.204	2.64 ng/ul	711319	Chrysene-d12		21.504	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...		666	C26H37F15O2	1000406-04-8	94
2	Heptafluorobutyric acid, pentade...		424	C19H31F7O2	959261-23-5	94
3	1-Hexacosene		364	C26H52	018835-33-1	93
4	1-Eicosene		280	C20H40	003452-07-1	91
5	Eicosyl trifluoroacetate		394	C22H41F3O2	1000351-75-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042523\
Data File : BN025126.D
Acq On : 25 Apr 2023 12:18
Operator : CG/JU
Sample : 02400-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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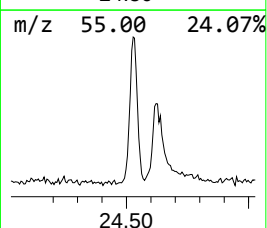
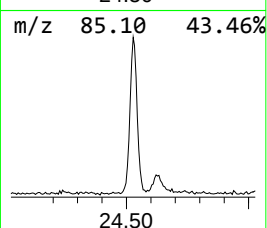
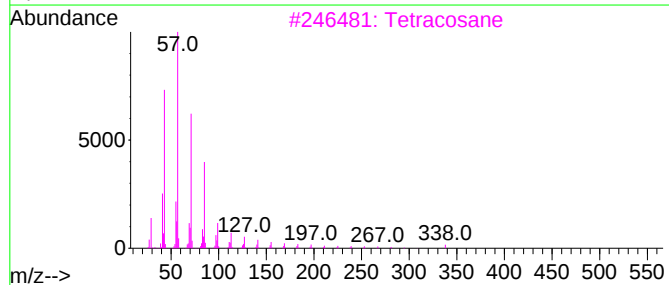
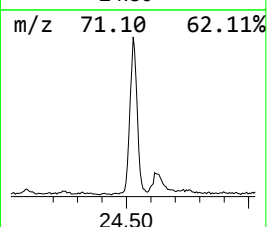
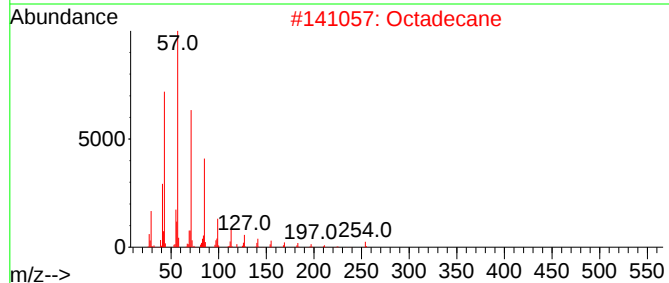
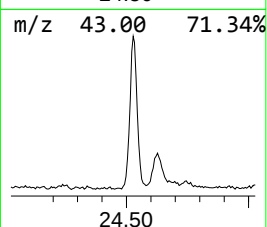
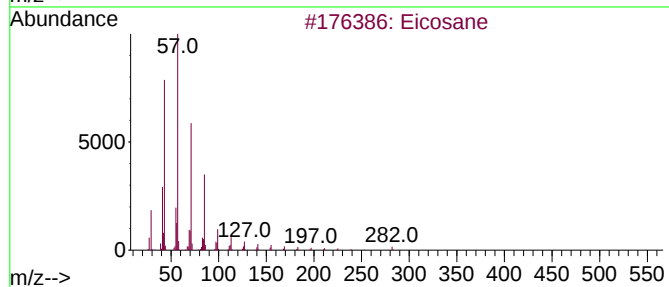
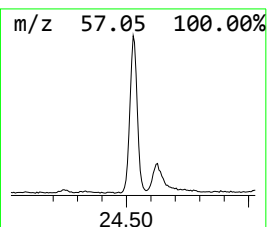
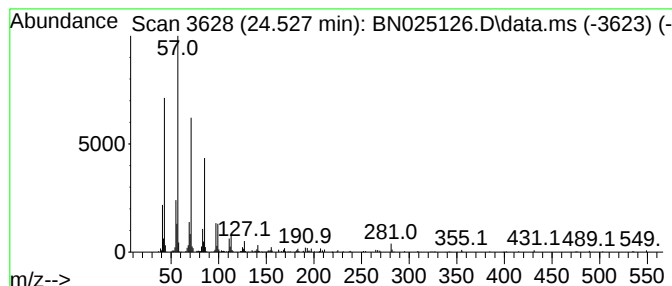
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	2.14 ng/ul	539659	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Octadecane	254	C18H38	000593-45-3	95
3			Tetracosane	338	C24H50	000646-31-1	87
4			Octacosane	394	C28H58	000630-02-4	87
5			Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042523\
Data File : BN025126.D
Acq On : 25 Apr 2023 12:18
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Sample : 02400-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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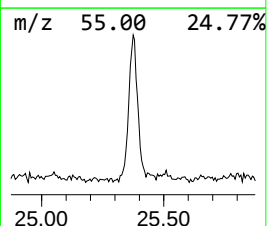
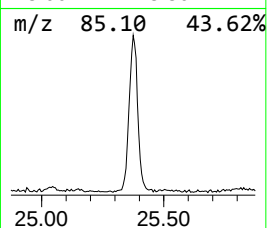
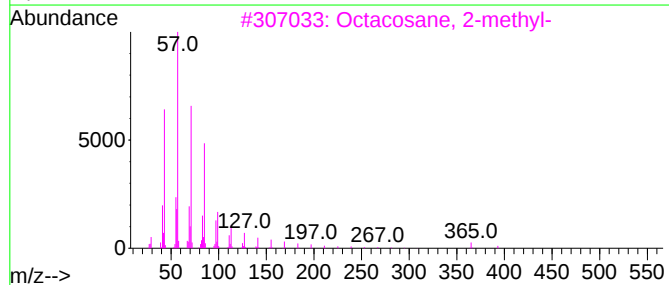
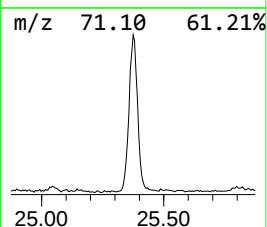
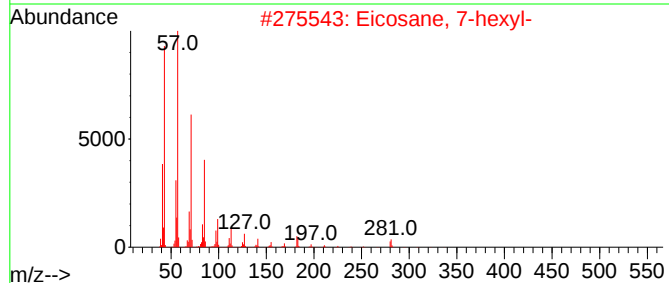
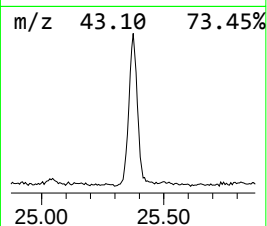
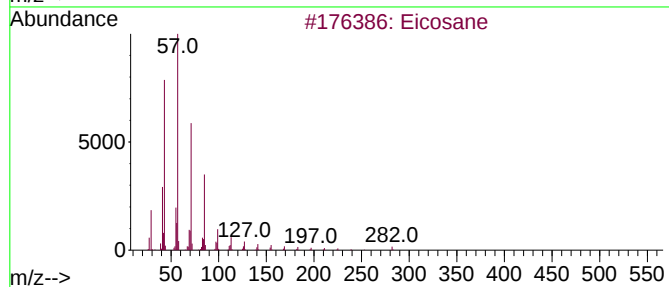
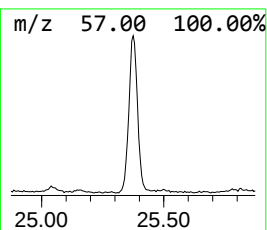
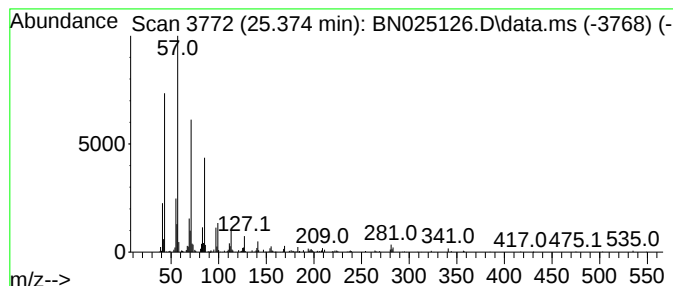
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.374	2.05 ng/ul	515302	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
4			Hexacosane	366	C26H54	000630-01-3	90
5			Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042523\
Data File : BN025126.D
Acq On : 25 Apr 2023 12:18
Operator : CG/JU
Sample : 02400-10
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
COML6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.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
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Benzene, 1,3-di...	7.799	36.3	ng/ul	3600330	1	7.916	1982510	20.0
Benzene, 1,4-di...	8.269	107.5	ng/ul	10653200	1	7.916	1982510	20.0
Benzene, 1,2,4-...	10.587	15.3	ng/ul	2529080	2	10.728	3317540	20.0
Benzene, 1,2,3-...	11.110	6.3	ng/ul	1045270	2	10.728	3317540	20.0
Benzophenone	15.916	8.6	ng/ul	1923760	3	14.563	4459130	20.0
Pentadecafluoro...	21.204	2.6	ng/ul	711319	5	21.504	5386240	20.0
(DEL) Alkane: S...	24.527	2.1	ng/ul	539659	6	23.927	5037050	20.0
(DEL) Alkane: S...	25.374	2.0	ng/ul	515302	6	23.927	5037050	20.0