

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019065.D
 Acq On : 23 Dec 2023 06:47
 Operator : MA/JU
 Sample : 05835-04
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B34

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019065.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	24	28	37	rVB	45099	59291	1.23%	0.132%
2	3.528	72	75	84	rBV	187546	259181	5.37%	0.579%
3	3.675	96	100	120	rVB	192352	308476	6.39%	0.689%
4	3.993	151	154	160	rVB	10463	14074	0.29%	0.031%
5	4.705	270	275	276	rBV5	4318	6833	0.14%	0.015%
6	4.916	309	311	318	rVB2	5914	6982	0.14%	0.016%
7	5.899	476	478	485	rVB4	5036	8017	0.17%	0.018%
8	6.822	631	635	641	rBV6	3148	6680	0.14%	0.015%
9	6.993	655	664	676	rBV	638304	1008580	20.90%	2.253%
10	7.152	684	691	701	rBV	512383	803350	16.65%	1.794%
11	7.346	717	724	734	rBV	649027	1004590	20.82%	2.244%
12	7.428	734	738	754	rVB	31059	61056	1.27%	0.136%
13	7.816	796	804	813	rBV	573752	884063	18.32%	1.975%
14	7.893	813	817	825	rVB4	6903	14601	0.30%	0.033%
15	8.534	918	926	936	rBV	749941	1230533	25.50%	2.749%
16	8.975	993	1001	1011	rBV	585520	958475	19.87%	2.141%
17	9.140	1026	1029	1036	rVB8	4334	7771	0.16%	0.017%
18	9.381	1065	1070	1076	rVB8	4106	7242	0.15%	0.016%
19	9.551	1091	1099	1104	rBV	25686	41384	0.86%	0.092%
20	9.698	1116	1124	1138	rBV	452602	755446	15.66%	1.687%
21	10.234	1207	1215	1236	rBV	747397	1319613	27.35%	2.948%
22	10.610	1271	1279	1292	rVB	712890	1205121	24.98%	2.692%
23	10.751	1295	1303	1325	rBV	594885	1082571	22.44%	2.418%
24	11.175	1368	1375	1383	rBV	7892	23150	0.48%	0.052%
25	11.245	1383	1387	1394	rVB8	8234	13562	0.28%	0.030%
26	11.357	1400	1406	1424	rBV	75000	170223	3.53%	0.380%
27	12.198	1545	1549	1557	rVB2	13715	24244	0.50%	0.054%
28	13.063	1691	1696	1700	rBV8	3646	7316	0.15%	0.016%
29	13.275	1728	1732	1739	rVB5	4602	6948	0.14%	0.016%
30	13.351	1739	1745	1753	rVV3	15153	27278	0.57%	0.061%
31	13.410	1753	1755	1764	rVB9	3831	6914	0.14%	0.015%
32	13.675	1794	1800	1804	rVB8	10437	16731	0.35%	0.037%
33	13.869	1826	1833	1842	rBV	1378324	1919304	39.78%	4.287%
34	14.139	1869	1879	1893	rBV	1601989	2442660	50.63%	5.456%
35	14.451	1925	1932	1940	rBV	1119058	1684263	34.91%	3.762%
36	14.669	1961	1969	1985	rBV2	601050	1198450	24.84%	2.677%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Title : SVOA CALIBRATION

37	14.875	1998	2004	2012	rVB9	12477	26904	0.56%	0.060%
38	14.963	2015	2019	2026	rVB9	4884	8767	0.18%	0.020%
39	15.445	2094	2101	2114	rBV	2174120	3141546	65.11%	7.017%
40	15.569	2116	2122	2133	rVV	537562	747954	15.50%	1.671%
41	15.680	2137	2141	2147	rVB5	10788	20716	0.43%	0.046%
42	16.010	2191	2197	2202	rBV8	13491	26542	0.55%	0.059%
43	16.157	2219	2222	2226	rBV3	8335	10552	0.22%	0.024%
44	16.228	2231	2234	2240	rVB4	9138	14588	0.30%	0.033%
45	16.339	2250	2253	2257	rVB4	5372	6520	0.14%	0.015%
46	16.963	2356	2359	2361	rBV3	7065	9002	0.19%	0.020%
47	17.198	2393	2399	2409	rBV	1591261	2192142	45.44%	4.896%
48	17.298	2410	2416	2427	rVV2	2447073	3508796	72.73%	7.837%
49	17.404	2431	2434	2442	rVV8	19389	34053	0.71%	0.076%
50	17.492	2446	2449	2456	rVB7	4823	8627	0.18%	0.019%
51	17.704	2481	2485	2489	rBV5	7150	10103	0.21%	0.023%
52	17.874	2510	2514	2518	rVB	78394	96716	2.00%	0.216%
53	18.092	2546	2551	2560	rBV	329250	491740	10.19%	1.098%
54	18.374	2596	2599	2603	rVB4	11615	13189	0.27%	0.029%
55	18.986	2700	2703	2705	rBV4	14441	17104	0.35%	0.038%
56	19.116	2721	2725	2733	rBV4	41267	64651	1.34%	0.144%
57	19.186	2733	2737	2740	rBV	60778	82033	1.70%	0.183%
58	19.327	2757	2761	2771	rBV	130430	184810	3.83%	0.413%
59	19.539	2791	2797	2803	rBV	3415846	4410406	91.41%	9.851%
60	20.551	2967	2969	2973	rVB2	49799	49998	1.04%	0.112%
61	21.010	3043	3047	3056	rBV	374901	620801	12.87%	1.387%
62	21.292	3090	3095	3104	rBV	2104171	2705311	56.07%	6.043%
63	23.498	3463	3470	3485	rVB2	2422111	4824703	100.00%	10.777%
64	23.651	3490	3496	3505	rVB	1464277	2847309	59.02%	6.360%

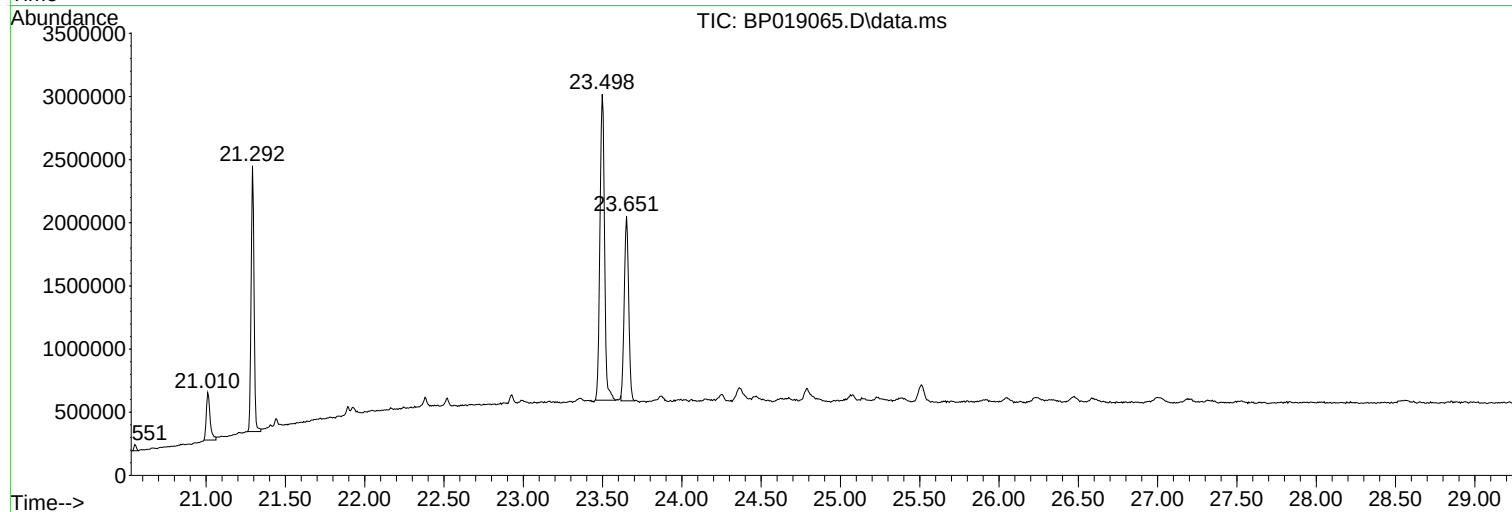
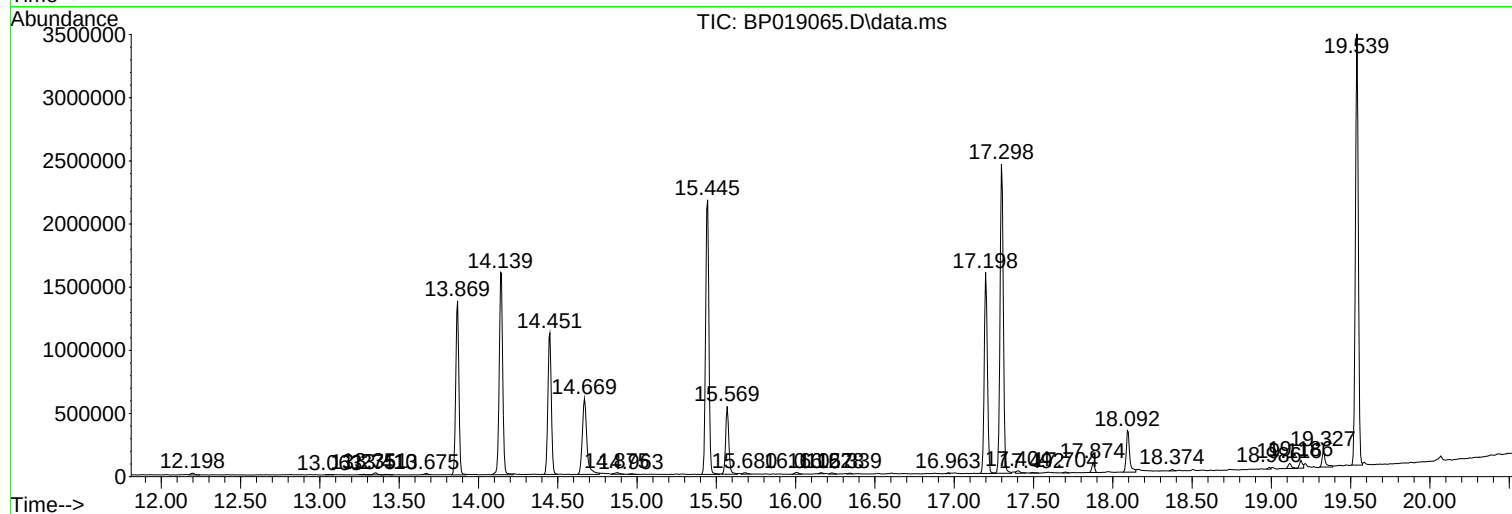
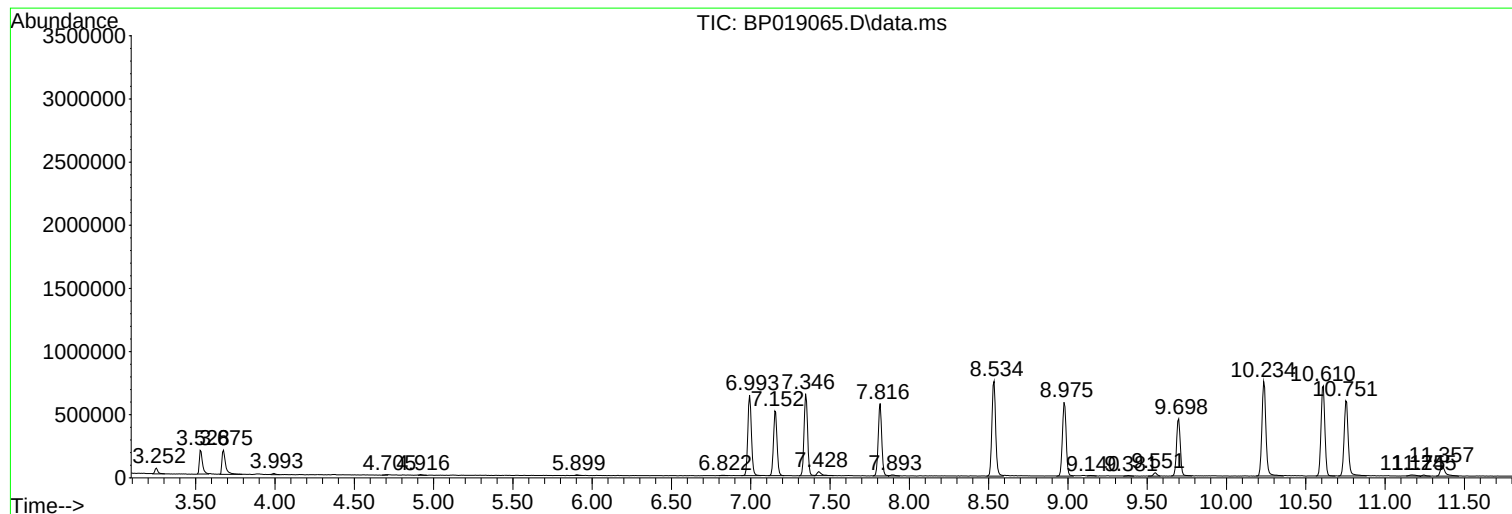
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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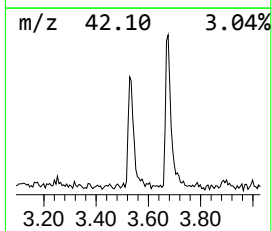
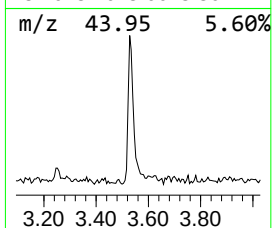
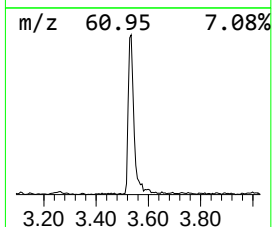
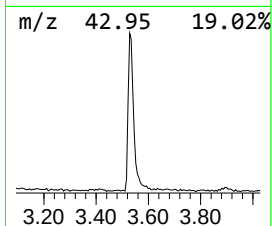
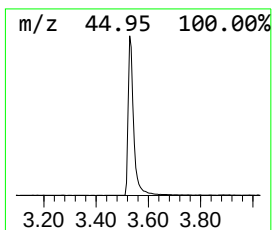
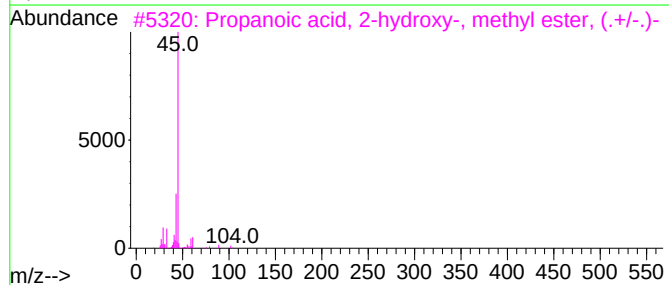
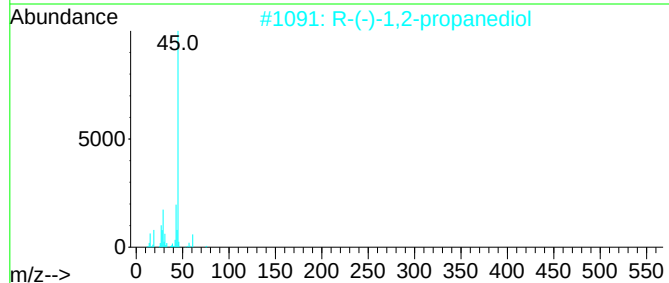
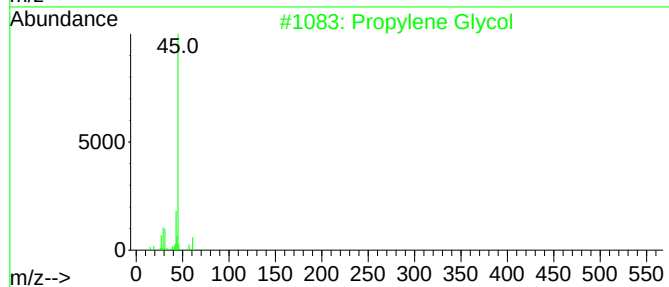
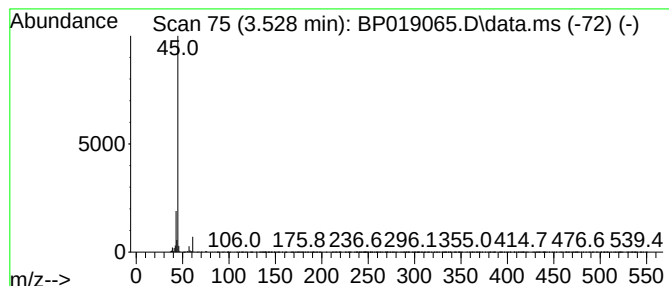
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.528	5.86 ng/ul	259181	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
4			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40
5			Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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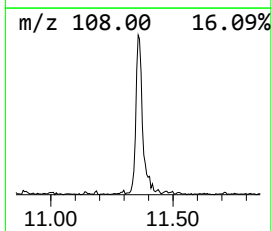
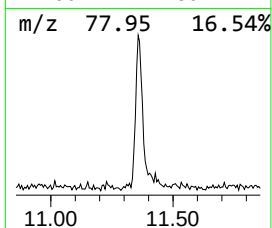
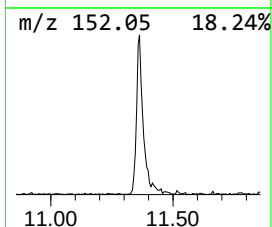
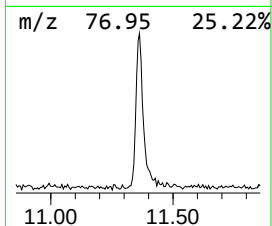
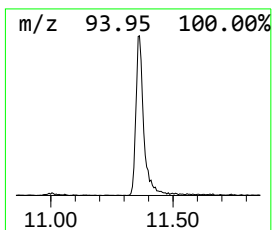
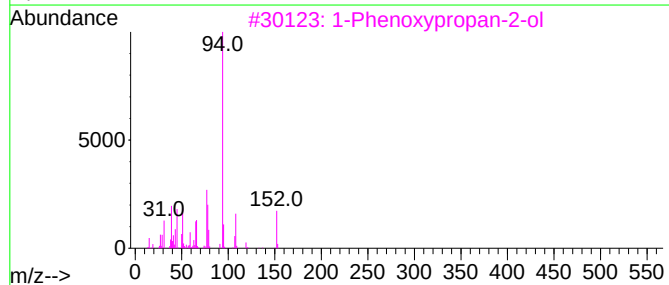
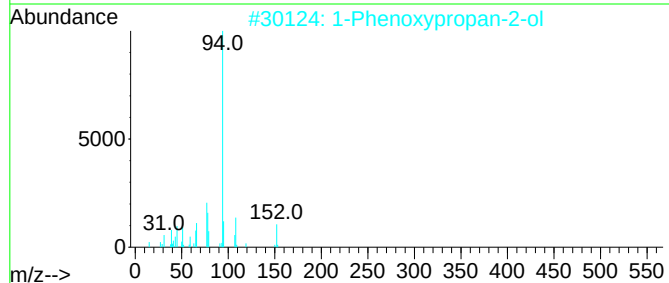
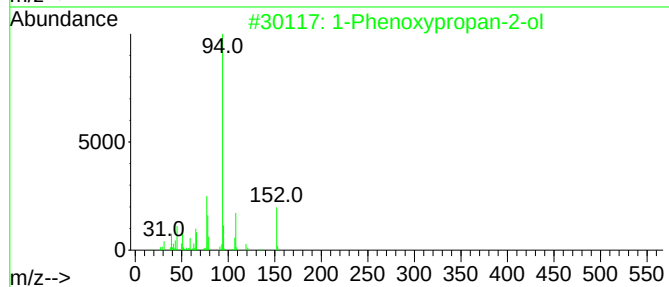
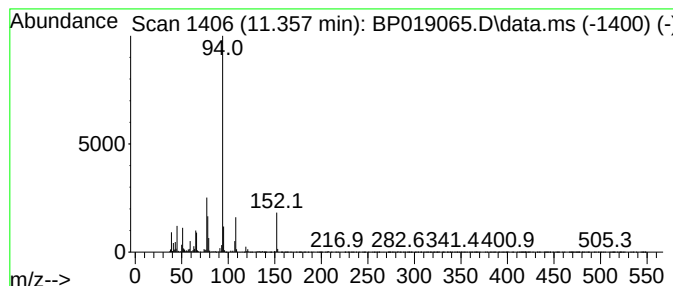
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.357	2.82 ng/ul	170223	Naphthalene-d8	10.610

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4		1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	90
5		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87



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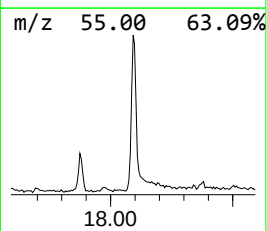
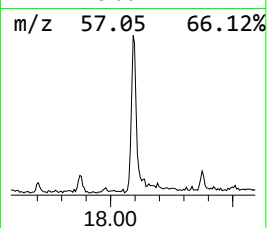
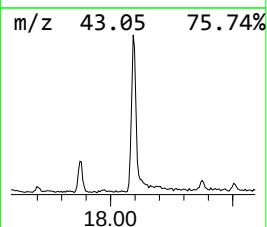
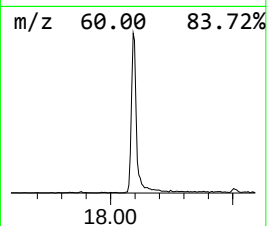
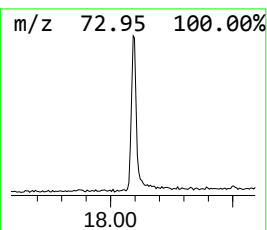
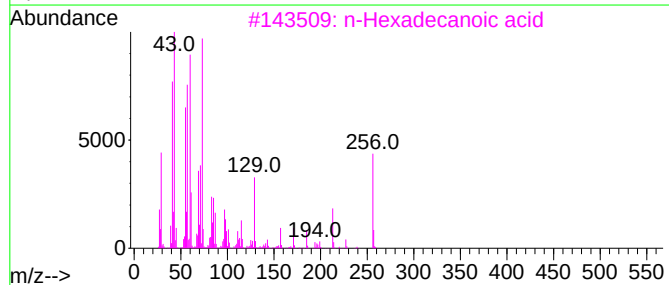
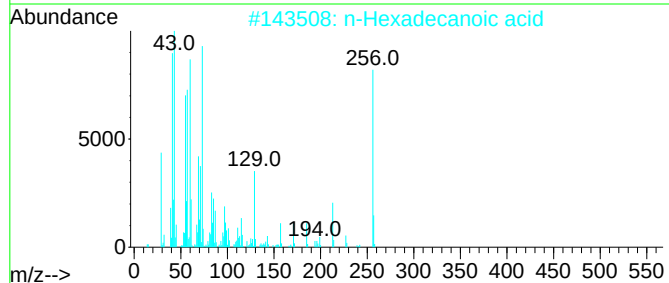
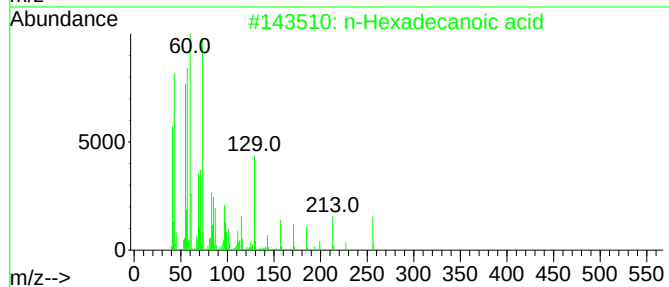
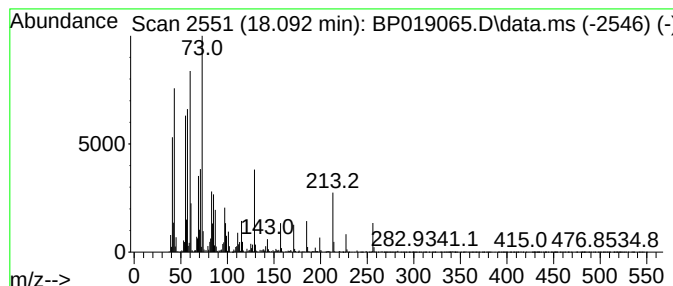
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	4.49 ng/ul	491740	Phenanthrene-d10	17.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



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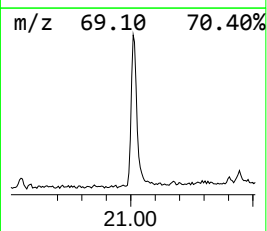
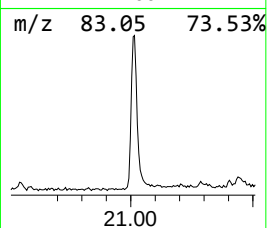
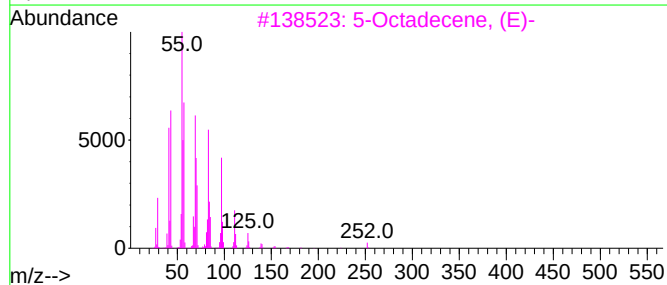
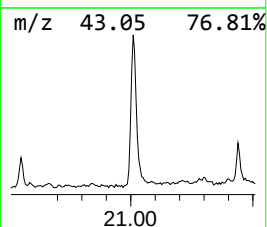
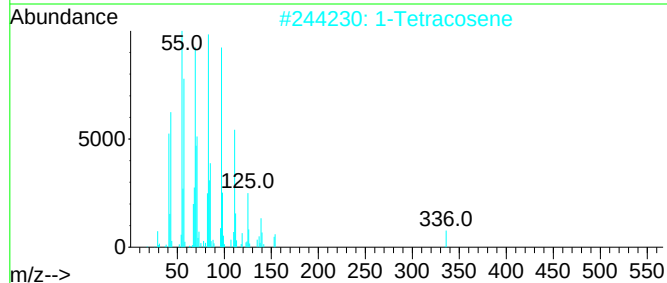
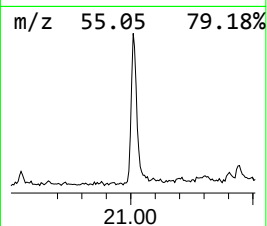
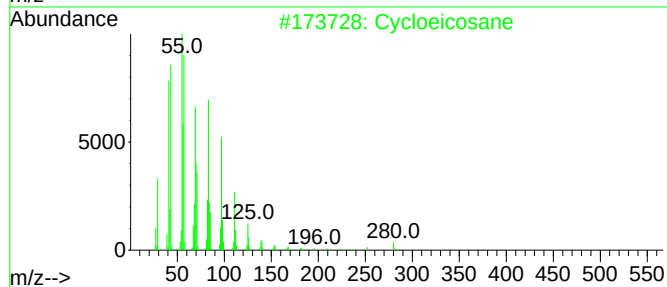
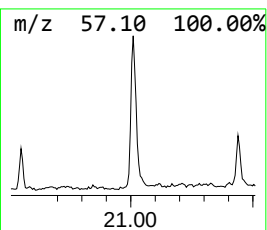
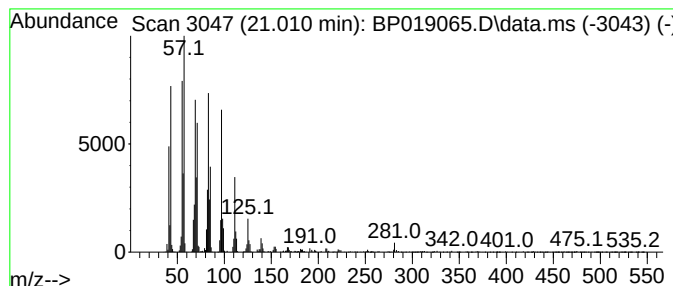
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic21.010 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.010	4.59 ng/ul	620801	Chrysene-d12	21.292

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloeicosane	280	C20H40	000296-56-0	98
2			1-Tetracosene	336	C24H48	010192-32-2	97
3			5-Octadecene, (E)-	252	C18H36	007206-21-5	95
4			1-Hexacosanol	382	C26H54O	000506-52-5	94
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94



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ClientSampleId :
C0B34

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.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
Propylene Glycol	3.528	5.9	ng/ul	259181	1	7.816	884063	20.0
1-Phenoxypropan...	11.357	2.8	ng/ul	170223	2	10.610	1205120	20.0
n-Hexadecanoic ...	18.092	4.5	ng/ul	491740	4	17.198	2192140	20.0
(DEL) Alkane: C...	21.010	4.6	ng/ul	620801	5	21.292	2705310	20.0