

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
 Data File : BN032196.D
 Acq On : 23 Jun 2024 10:58
 Operator : MA/JU
 Sample : P2775-08
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4C86

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-BN052924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN032196.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	26	30	39	rVB	77354	107397	1.92%	0.158%
2	3.428	71	75	87	rBV	246625	349831	6.24%	0.514%
3	3.570	96	99	110	rBV	872336	1323813	23.63%	1.946%
4	3.875	147	151	156	rVB	25894	34672	0.62%	0.051%
5	4.375	233	236	239	rVV3	5193	6417	0.11%	0.009%
6	4.423	240	244	249	rVB	10351	16320	0.29%	0.024%
7	4.570	265	269	272	rBV6	3337	6133	0.11%	0.009%
8	4.658	281	284	291	rVB7	3880	5900	0.11%	0.009%
9	4.770	299	303	310	rVB	43291	64077	1.14%	0.094%
10	5.734	462	467	478	rBV	32583	58803	1.05%	0.086%
11	6.258	552	556	560	rBV6	3649	5825	0.10%	0.009%
12	6.658	621	624	630	rVB5	6815	10067	0.18%	0.015%
13	6.817	643	651	661	rBV	1130961	1847253	32.97%	2.715%
14	6.975	671	678	692	rVB	975490	1527423	27.26%	2.245%
15	7.169	703	711	719	rBV	1206302	2004479	35.77%	2.946%
16	7.240	719	723	735	rVB	81794	157541	2.81%	0.232%
17	7.628	781	789	797	rBV	1264520	2024654	36.13%	2.976%
18	7.705	797	802	809	rVB2	20688	45909	0.82%	0.067%
19	8.346	903	911	923	rBV	1388106	2344339	41.84%	3.445%
20	8.787	975	986	999	rBV	1061645	1789076	31.93%	2.629%
21	8.911	1003	1007	1011	rBV2	8904	15080	0.27%	0.022%
22	9.163	1045	1050	1054	rBV3	7010	11494	0.21%	0.017%
23	9.352	1076	1082	1090	rBV	88690	151706	2.71%	0.223%
24	9.505	1100	1108	1121	rBV	867346	1497652	26.73%	2.201%
25	9.752	1145	1150	1156	rBV8	5243	11717	0.21%	0.017%
26	10.040	1192	1199	1213	rBV	1275809	2282558	40.74%	3.355%
27	10.410	1254	1262	1274	rBV	1599372	2665617	47.57%	3.918%
28	10.557	1278	1287	1307	rVV	1296151	2296875	40.99%	3.376%
29	10.957	1349	1355	1367	rBV3	23948	48201	0.86%	0.071%
30	11.157	1382	1389	1407	rBV	237323	498006	8.89%	0.732%
31	11.922	1514	1519	1523	rBV6	5895	10127	0.18%	0.015%
32	12.004	1526	1533	1539	rVV3	25938	43603	0.78%	0.064%
33	12.622	1629	1638	1640	rBV8	4439	8355	0.15%	0.012%
34	12.746	1652	1659	1663	rBV9	3243	7416	0.13%	0.011%
35	12.887	1676	1683	1689	rBV9	5479	14988	0.27%	0.022%
36	13.240	1734	1743	1750	rBV9	8573	17881	0.32%	0.026%

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37	13.693	1813	1820	1828	rBV	2070462	2937298	52.42%	4.317%
38	13.969	1855	1867	1884	rBV2	2508755	3912223	69.82%	5.750%
39	14.275	1912	1919	1930	rBV	2043838	3102827	55.37%	4.560%
40	14.516	1944	1960	1978	rBV	666368	1428939	25.50%	2.100%
41	14.710	1989	1993	2003	rVB7	13886	28301	0.51%	0.042%
42	14.793	2004	2007	2014	rVB9	6530	9976	0.18%	0.015%
43	14.963	2031	2036	2039	rVV7	3870	6227	0.11%	0.009%
44	15.069	2050	2054	2058	rBV5	11544	18165	0.32%	0.027%
45	15.128	2060	2064	2067	rBV4	8847	11632	0.21%	0.017%
46	15.275	2082	2089	2102	rVB	3035920	4427007	79.01%	6.506%
47	15.410	2107	2112	2125	rVV	383439	605881	10.81%	0.890%
48	15.510	2125	2129	2136	rVV4	20463	37774	0.67%	0.056%
49	15.634	2147	2150	2156	rVB7	6233	9079	0.16%	0.013%
50	15.698	2156	2161	2164	rBV6	5468	7501	0.13%	0.011%
51	15.828	2179	2183	2191	rBV8	12783	25349	0.45%	0.037%
52	15.992	2207	2211	2219	rVB7	11546	17965	0.32%	0.026%
53	16.310	2257	2265	2266	rBV8	4224	8203	0.15%	0.012%
54	16.328	2266	2268	2276	rVV9	2997	6591	0.12%	0.010%
55	16.445	2284	2288	2293	rBV7	5281	10396	0.19%	0.015%
56	16.569	2306	2309	2315	rVB8	6446	9861	0.18%	0.014%
57	16.628	2315	2319	2322	rBV6	4365	5819	0.10%	0.009%
58	16.704	2326	2332	2335	rBV8	5934	10769	0.19%	0.016%
59	16.740	2335	2338	2342	rBV6	5050	6139	0.11%	0.009%
60	16.781	2342	2345	2348	rBV4	11661	15417	0.28%	0.023%
61	16.810	2348	2350	2362	rVB7	10878	27785	0.50%	0.041%
62	16.904	2363	2366	2369	rBV5	5350	7566	0.14%	0.011%
63	16.945	2370	2373	2378	rVB7	6446	9474	0.17%	0.014%
64	17.028	2380	2387	2393	rBV	2270360	3340851	59.62%	4.910%
65	17.128	2398	2404	2415	rVB	2985262	4312373	76.96%	6.338%
66	17.234	2419	2422	2429	rVB6	15940	27084	0.48%	0.040%
67	17.316	2430	2436	2446	rBV3	37082	108553	1.94%	0.160%
68	17.522	2467	2471	2475	rBV6	8357	14087	0.25%	0.021%
69	17.634	2485	2490	2494	rVB4	13302	24160	0.43%	0.036%
70	17.698	2498	2501	2506	rVB	35188	43296	0.77%	0.064%
71	17.928	2535	2540	2551	rBV2	298610	530300	9.46%	0.779%
72	18.098	2565	2569	2573	rBV3	27655	37791	0.67%	0.056%
73	18.198	2584	2586	2588	rBV2	9487	9064	0.16%	0.013%
74	18.263	2593	2597	2603	rBV4	38854	69055	1.23%	0.101%
75	18.381	2613	2617	2620	rBV6	26169	36303	0.65%	0.053%
76	18.992	2718	2721	2727	rVB3	42989	51274	0.92%	0.075%

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Title : SVOA CALIBRATION

77	19.063	2729	2733	2736	rBV	55386	95394	1.70%	0.140%
78	19.092	2736	2738	2746	rVB	80002	114417	2.04%	0.168%
79	19.216	2755	2759	2769	rBV4	87179	168835	3.01%	0.248%
80	19.428	2790	2795	2808	rBV	3519938	5080455	90.67%	7.467%
81	20.951	3050	3054	3063	rVB	479667	683109	12.19%	1.004%
82	21.263	3102	3107	3113	rBV	2319703	3544432	63.26%	5.209%
83	23.980	3562	3569	3587	rVB2	2253185	5603374	100.00%	8.235%
84	24.180	3595	3603	3615	rVB	1671366	4131127	73.73%	6.072%

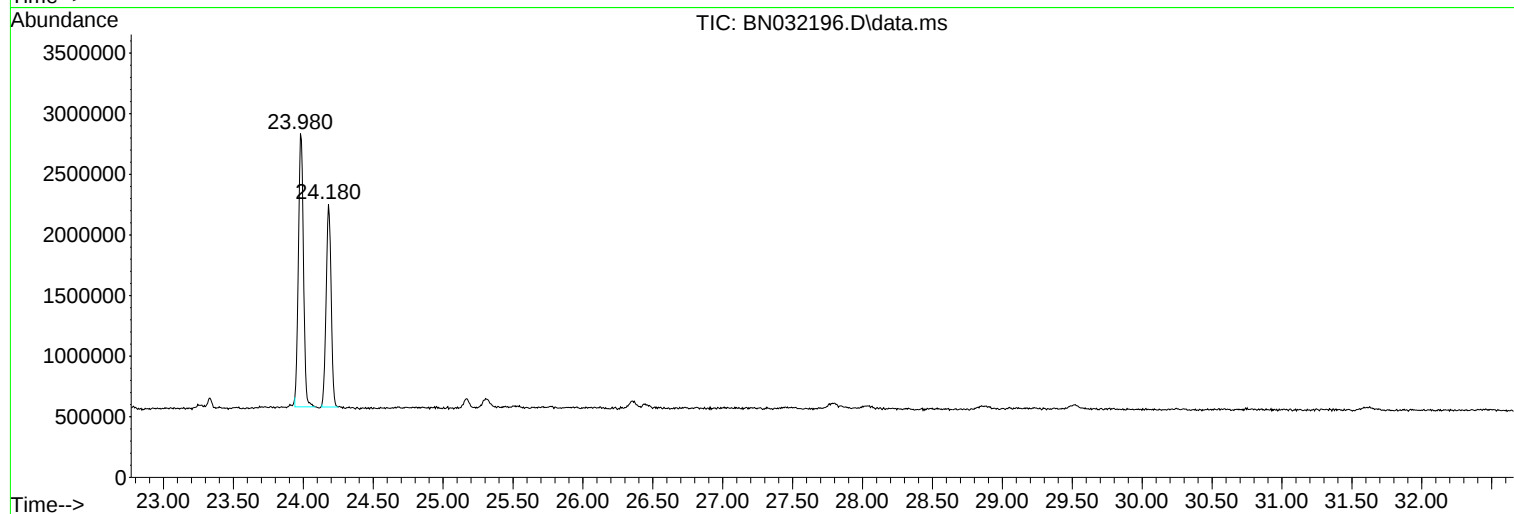
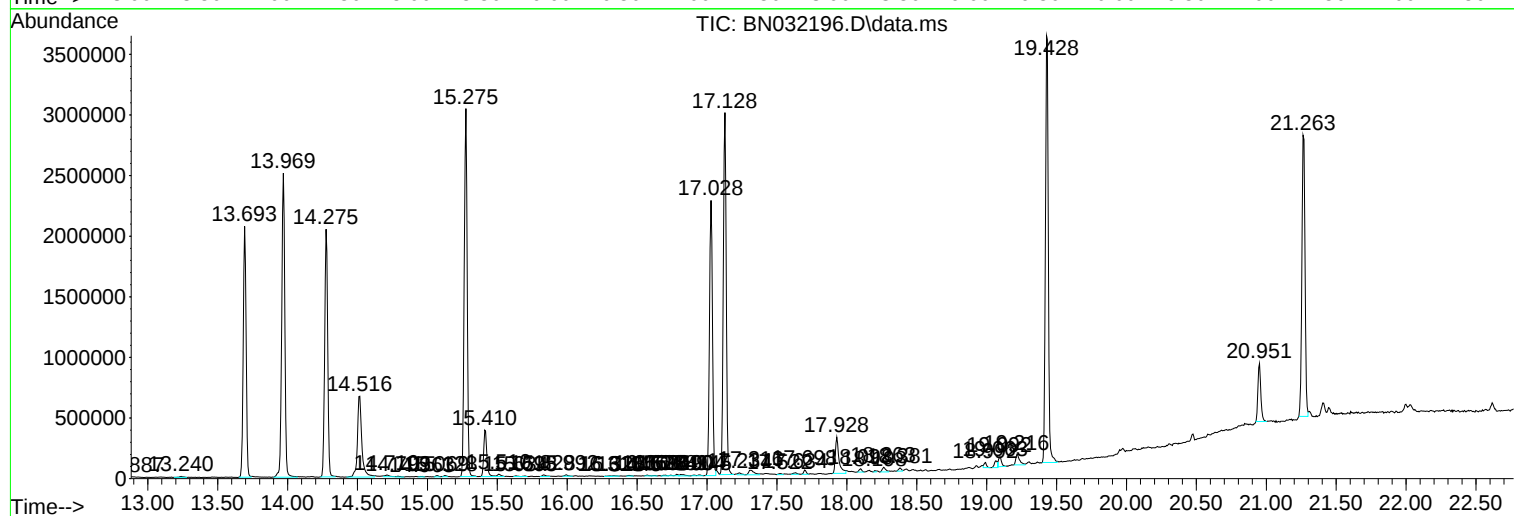
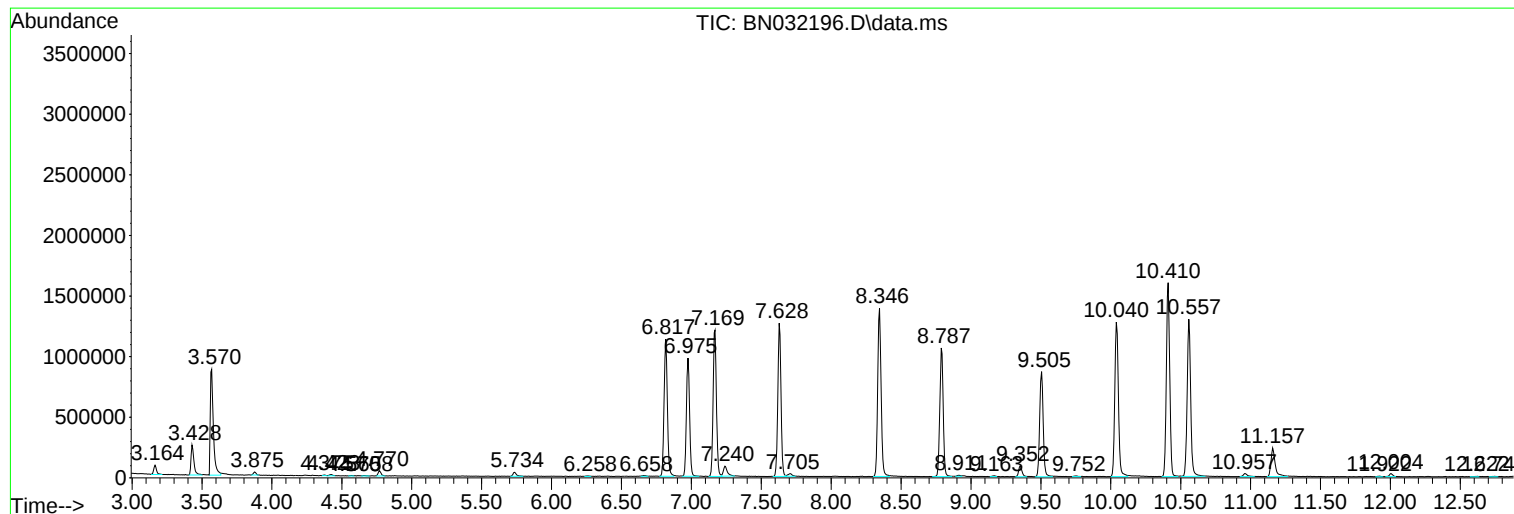
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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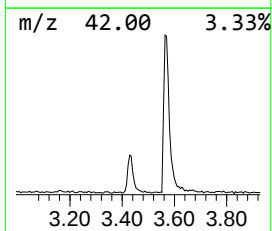
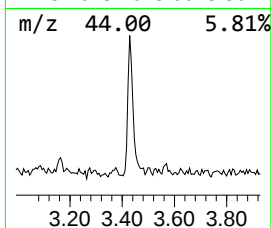
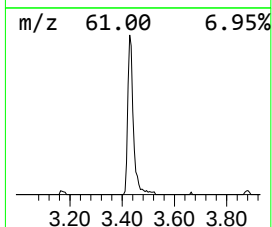
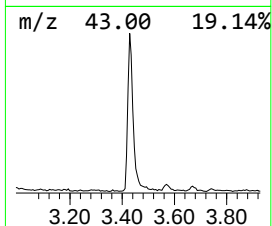
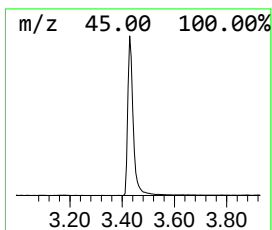
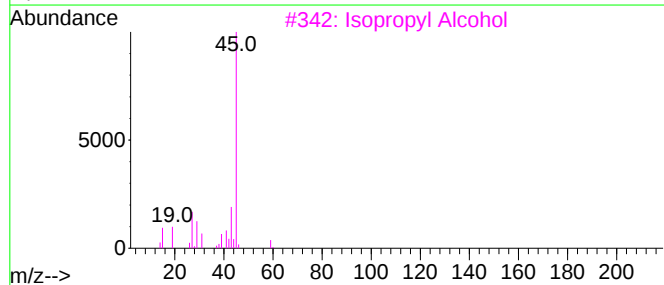
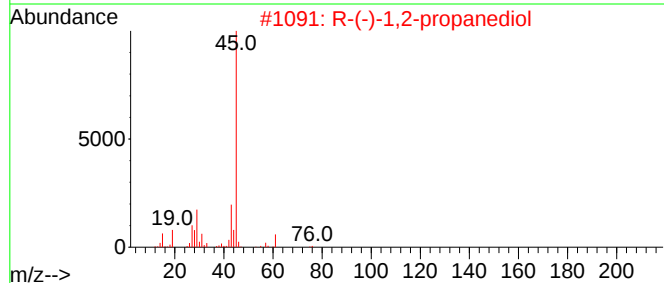
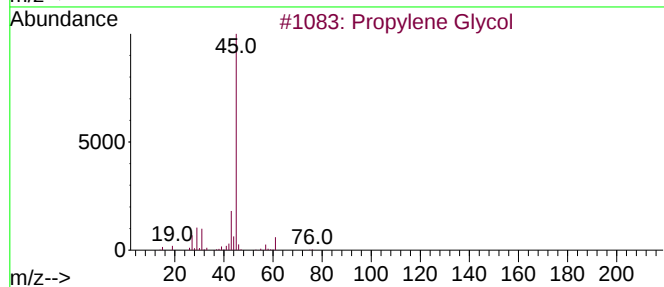
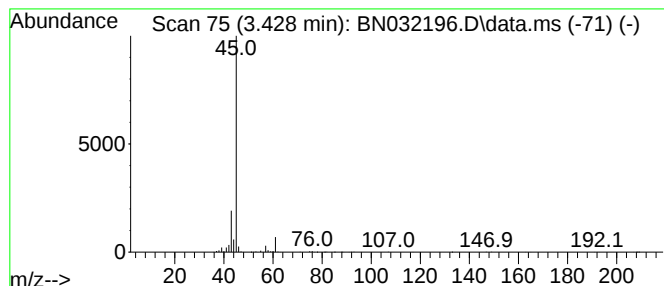
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.428	3.46 ng/ul	349831	1,4-Dichlorobenzene-d4	7.628

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			Isopropyl Alcohol	60	C3H8O	000067-63-0	50
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			2-Butanol	74	C4H10O	000078-92-2	40



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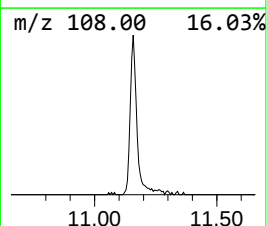
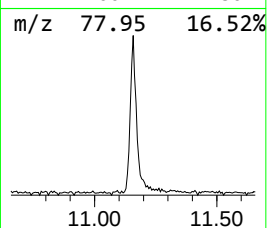
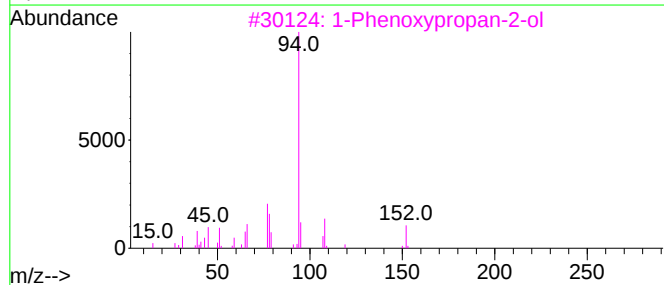
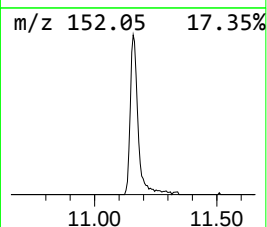
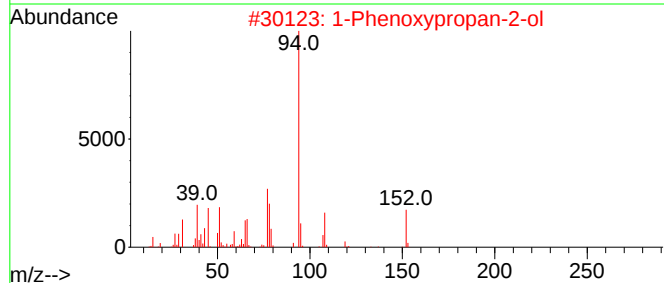
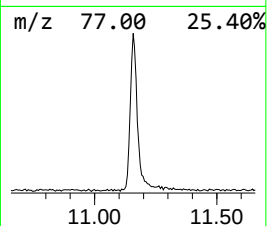
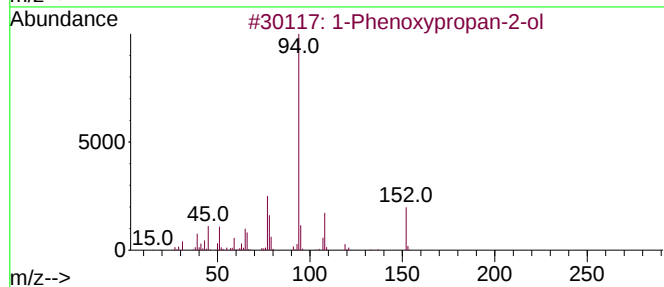
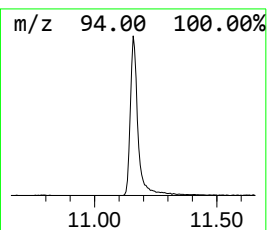
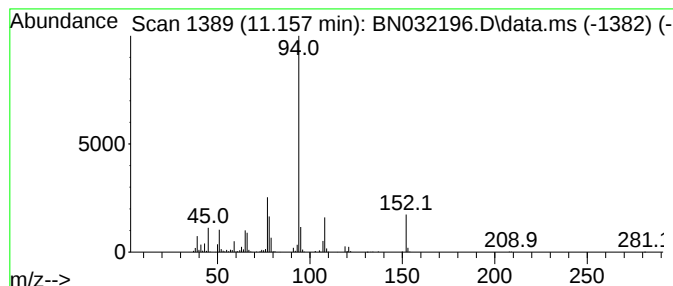
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.157	3.74 ng/ul	498006	Naphthalene-d8	10.410

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	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
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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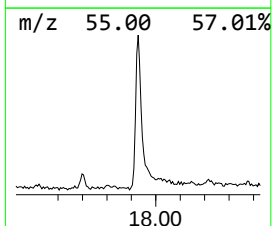
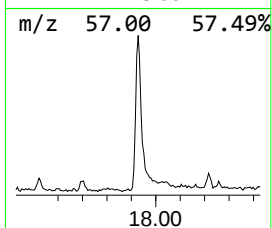
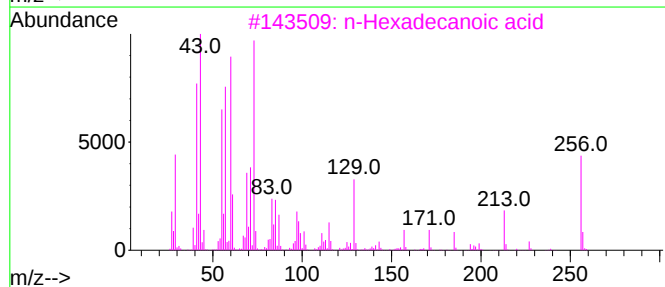
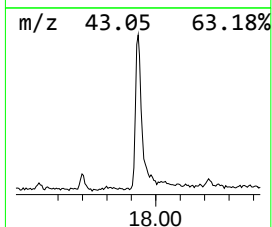
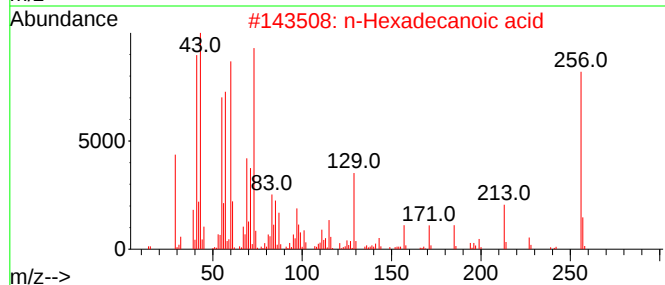
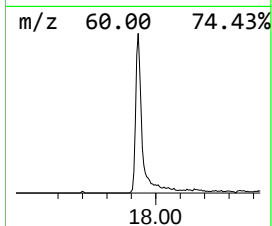
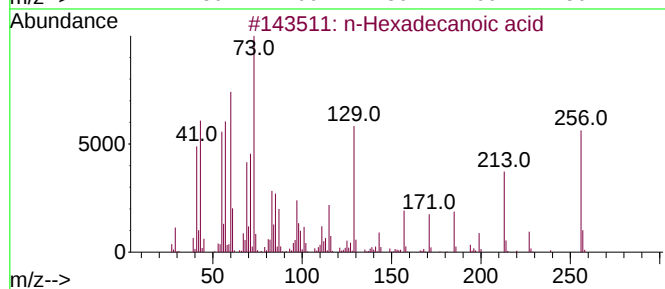
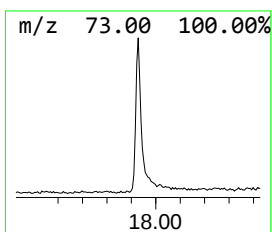
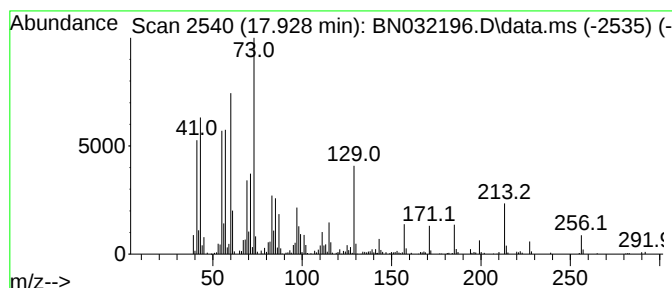
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.928	3.17 ng/ul	530300	Phenanthrene-d10	17.028	
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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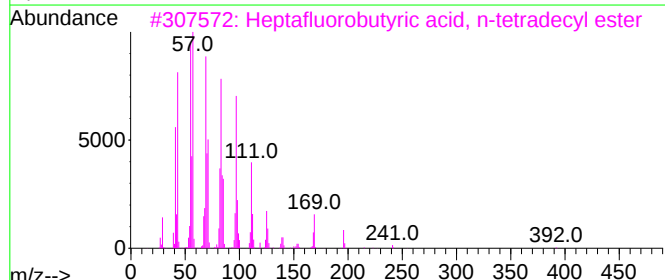
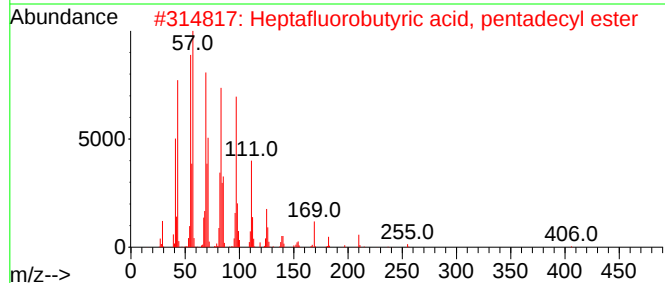
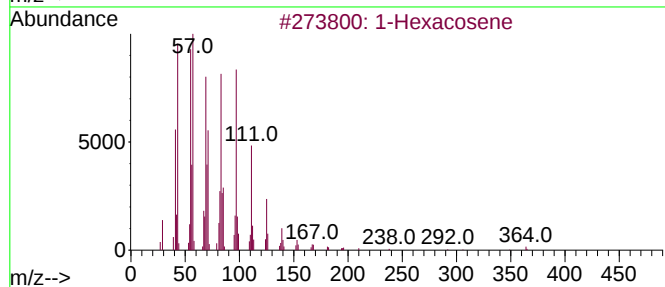
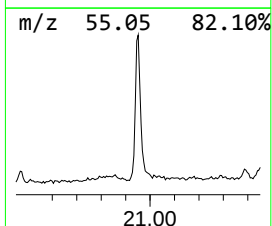
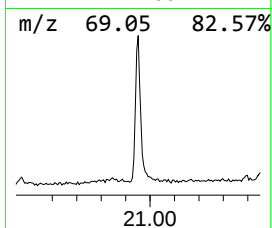
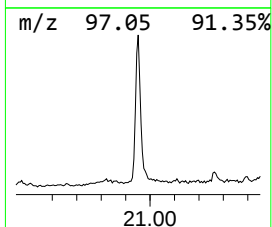
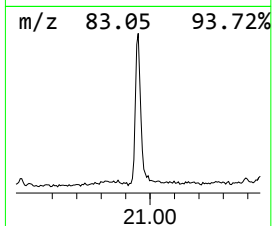
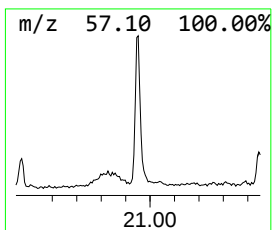
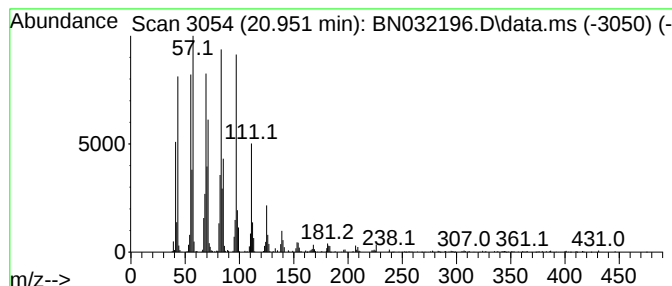
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.951	3.85 ng/ul	683109	Chrysene-d12	21.269

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene			364	C26H52	018835-33-1	95
2	Heptafluorobutyric acid, pentade...			424	C19H31F7O2	959261-23-5	94
3	Heptafluorobutyric acid, n-tetra...			410	C18H29F7O2	007365-36-8	94
4	1-Heneicosanol			312	C21H44O	015594-90-8	93
5	Nonadecyl trifluoroacetate			380	C21H39F3O2	1000351-76-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062224\
Data File : BN032196.D
Acq On : 23 Jun 2024 10:58
Operator : MA/JU
Sample : P2775-08
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A4C86

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.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	3.428	3.5	ng/u1	349831	1	7.628	2024650	20.0
1-Phenoxypropan...	11.157	3.7	ng/u1	498006	2	10.410	2665620	20.0
n-Hexadecanoic ...	17.928	3.2	ng/u1	530300	4	17.028	3340850	20.0
(DEL) Alkane: S...	20.951	3.9	ng/u1	683109	5	21.269	3544430	20.0