

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
 Data File : BN031930.D
 Acq On : 12 Jun 2024 04:09
 Operator : MA/JU
 Sample : P2659-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BHC33

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: BN031930.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	29	33	40	rVB	49719	65931	0.94%	0.069%
2	3.440	74	77	88	rBV	188328	265904	3.79%	0.280%
3	3.587	99	102	115	rBV	467491	730417	10.40%	0.768%
4	3.687	117	119	126	rVB2	14853	20690	0.29%	0.022%
5	3.899	151	155	163	rVB	37690	54070	0.77%	0.057%
6	4.793	302	307	313	rBV	29533	44998	0.64%	0.047%
7	5.763	463	472	483	rVB	32093	60092	0.86%	0.063%
8	6.834	642	654	667	rBV	1137926	1863491	26.53%	1.960%
9	7.004	675	683	693	rBV	897861	1421718	20.24%	1.495%
10	7.193	708	715	723	rBV	1119861	1789481	25.48%	1.882%
11	7.269	723	728	740	rVB	58266	111287	1.58%	0.117%
12	7.657	787	794	802	rBV	1248234	2017739	28.73%	2.122%
13	7.734	802	807	815	rVB2	29601	50924	0.73%	0.054%
14	8.369	907	915	925	rBV	1475123	2494647	35.52%	2.623%
15	8.816	983	991	1003	rBV	1178299	1902260	27.08%	2.000%
16	8.922	1004	1009	1014	rVV4	17011	29314	0.42%	0.031%
17	9.381	1080	1087	1100	rBV	159480	250811	3.57%	0.264%
18	9.534	1105	1113	1126	rBV	957004	1611842	22.95%	1.695%
19	10.063	1194	1203	1215	rBV	1512453	2687188	38.26%	2.826%
20	10.228	1226	1231	1240	rVB4	13426	25266	0.36%	0.027%
21	10.440	1258	1267	1275	rBV	1866234	3185685	45.36%	3.350%
22	10.587	1283	1292	1309	rBV	1331912	2422069	34.48%	2.547%
23	10.987	1350	1360	1368	rBV	55438	107188	1.53%	0.113%
24	11.187	1386	1394	1411	rBV	390701	808702	11.51%	0.850%
25	12.034	1533	1538	1546	rVB2	30204	52722	0.75%	0.055%
26	12.334	1584	1589	1595	rVB2	14149	21758	0.31%	0.023%
27	13.122	1719	1723	1729	rVB4	14170	22470	0.32%	0.024%
28	13.622	1802	1808	1813	rVB3	19975	36550	0.52%	0.038%
29	13.722	1818	1825	1839	rVB	3128163	4355339	62.01%	4.580%
30	13.998	1860	1872	1884	rBV	3389903	5463000	77.78%	5.745%
31	14.304	1914	1924	1931	rBV	2919998	4301941	61.25%	4.524%
32	14.369	1931	1935	1941	rVB3	17894	31488	0.45%	0.033%
33	14.522	1954	1961	1974	rBV	1290712	2261853	32.20%	2.378%
34	14.781	2001	2005	2009	rVB2	17683	23308	0.33%	0.025%
35	14.922	2022	2029	2034	rBV5	19637	45461	0.65%	0.048%
36	15.098	2052	2059	2063	rBV6	19946	43159	0.61%	0.045%

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Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
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37	15.163	2066	2070	2077	rVB2	18263	40475	0.58%	0.043%
38	15.298	2086	2093	2100	rBV	4410751	6353273	90.46%	6.681%
39	15.351	2100	2102	2109	rVB	43296	51065	0.73%	0.054%
40	15.428	2109	2115	2127	rBV	1191611	1635444	23.28%	1.720%
41	15.539	2130	2134	2137	rBV2	17528	28459	0.41%	0.030%
42	15.651	2148	2153	2159	rBV3	23446	44164	0.63%	0.046%
43	15.798	2174	2178	2185	rVB3	19711	40126	0.57%	0.042%
44	15.875	2185	2191	2197	rVB8	12615	28480	0.41%	0.030%
45	16.028	2211	2217	2224	rBV4	48691	99884	1.42%	0.105%
46	16.116	2228	2232	2238	rVV	28392	47575	0.68%	0.050%
47	16.363	2266	2274	2279	rBV4	23934	63103	0.90%	0.066%
48	16.410	2279	2282	2286	rVB	16812	22925	0.33%	0.024%
49	16.463	2286	2291	2296	rBV4	34956	60358	0.86%	0.063%
50	16.586	2309	2312	2316	rVV4	15396	21765	0.31%	0.023%
51	16.710	2328	2333	2340	rVB2	45733	86192	1.23%	0.091%
52	16.786	2341	2346	2348	rVV5	16728	27410	0.39%	0.029%
53	16.816	2348	2351	2354	rVV	26850	38500	0.55%	0.040%
54	16.869	2356	2360	2368	rVB	55919	98646	1.40%	0.104%
55	16.975	2371	2378	2382	rBV6	25674	58737	0.84%	0.062%
56	17.051	2382	2391	2395	rBV	3166925	4686717	66.73%	4.928%
57	17.092	2395	2398	2403	rVV	801672	1139683	16.23%	1.198%
58	17.151	2403	2408	2419	rVB	4423382	6368248	90.67%	6.696%
59	17.239	2419	2423	2428	rBV5	30459	61239	0.87%	0.064%
60	17.463	2453	2461	2471	rVB2	54259	114103	1.62%	0.120%
61	17.557	2473	2477	2478	rBV5	19250	25866	0.37%	0.027%
62	17.575	2478	2480	2485	rVV2	30358	35751	0.51%	0.038%
63	17.633	2486	2490	2497	rVB5	33661	64386	0.92%	0.068%
64	17.727	2503	2506	2510	rVB3	44322	59554	0.85%	0.063%
65	17.833	2520	2524	2530	rVB6	25437	54984	0.78%	0.058%
66	17.957	2535	2545	2559	rVV2	1325015	2589556	36.87%	2.723%
67	18.057	2559	2562	2567	rVV2	60507	108016	1.54%	0.114%
68	18.122	2567	2573	2577	rVV2	198099	392631	5.59%	0.413%
69	18.157	2577	2579	2586	rVB	116802	152424	2.17%	0.160%
70	18.245	2590	2594	2597	rBV5	50283	76037	1.08%	0.080%
71	18.280	2598	2600	2605	rVB2	44576	61020	0.87%	0.064%
72	18.386	2615	2618	2622	rVB6	22918	27906	0.40%	0.029%
73	18.433	2622	2626	2630	rVV	101575	141757	2.02%	0.149%
74	18.474	2630	2633	2639	rVB2	102472	136993	1.95%	0.144%
75	18.680	2665	2668	2671	rVB	36194	42797	0.61%	0.045%
76	18.733	2673	2677	2680	rVV	50892	73197	1.04%	0.077%

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 Title : SVOA CALIBRATION

77	18.769	2680	2683	2687	rVB	47125	56555	0.81%	0.059%
78	18.851	2692	2697	2701	rBV	129511	220883	3.14%	0.232%
79	18.910	2704	2707	2710	rVB3	64590	82740	1.18%	0.087%
80	18.939	2710	2712	2716	rVB	43502	51395	0.73%	0.054%
81	18.998	2717	2722	2723	rBV	90118	134810	1.92%	0.142%
82	19.116	2733	2742	2751	rBV	1565962	2372949	33.79%	2.495%
83	19.239	2757	2763	2772	rBV3	403927	761993	10.85%	0.801%
84	19.451	2793	2799	2803	rBV2	4626028	7023611	100.00%	7.386%
85	19.557	2813	2817	2822	rVB2	79591	118621	1.69%	0.125%
86	19.727	2842	2846	2851	rVB3	43953	78874	1.12%	0.083%
87	19.804	2855	2859	2861	rBV3	81164	131796	1.88%	0.139%
88	19.980	2885	2889	2891	rBV	135794	158426	2.26%	0.167%
89	20.080	2902	2906	2909	rBV2	104645	131801	1.88%	0.139%
90	20.133	2912	2915	2920	rVB2	131750	189813	2.70%	0.200%
91	20.498	2974	2977	2981	rBV2	124100	148363	2.11%	0.156%
92	20.727	3013	3016	3022	rVB	118250	167427	2.38%	0.176%
93	20.974	3054	3058	3065	rBV2	934575	1498100	21.33%	1.575%
94	21.286	3104	3111	3116	rBV3	2573800	4651233	66.22%	4.891%
95	21.327	3116	3118	3128	rVB	556976	801531	11.41%	0.843%
96	22.033	3235	3238	3240	rBV3	143400	197832	2.82%	0.208%
97	22.651	3338	3343	3351	rVB3	247434	530808	7.56%	0.558%
98	23.292	3447	3452	3459	rBV2	327521	836898	11.92%	0.880%
99	24.015	3568	3575	3582	rBV2	2047232	4753234	67.68%	4.998%
100	24.215	3601	3609	3618	rBV	1583653	4059221	57.79%	4.268%

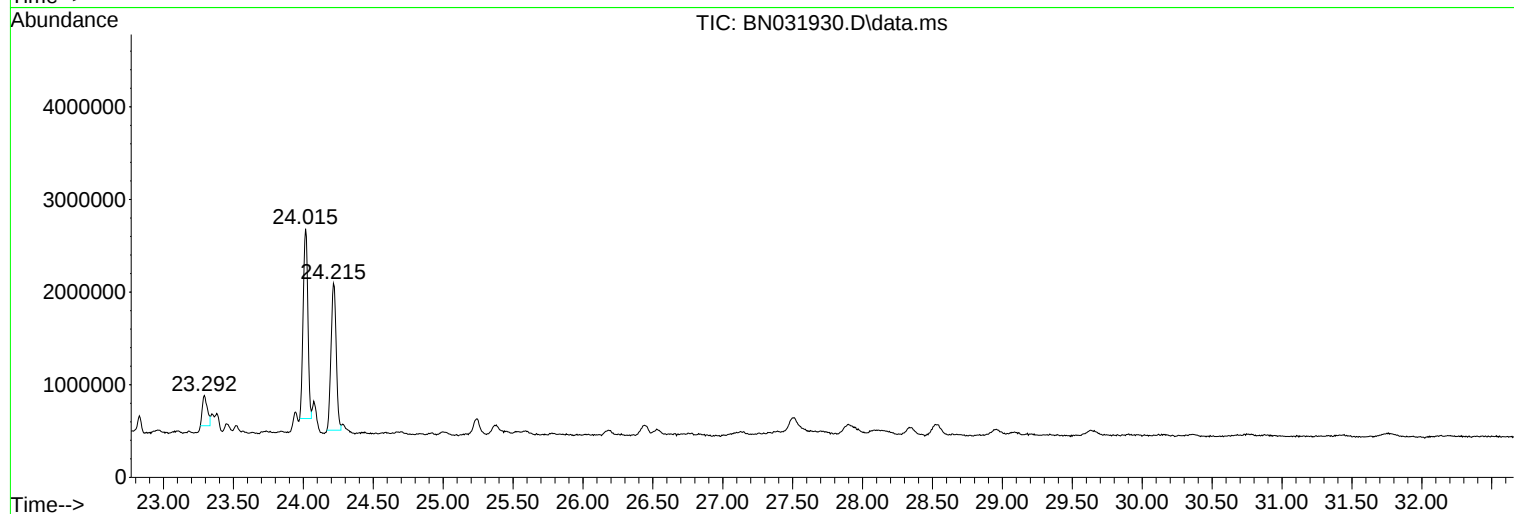
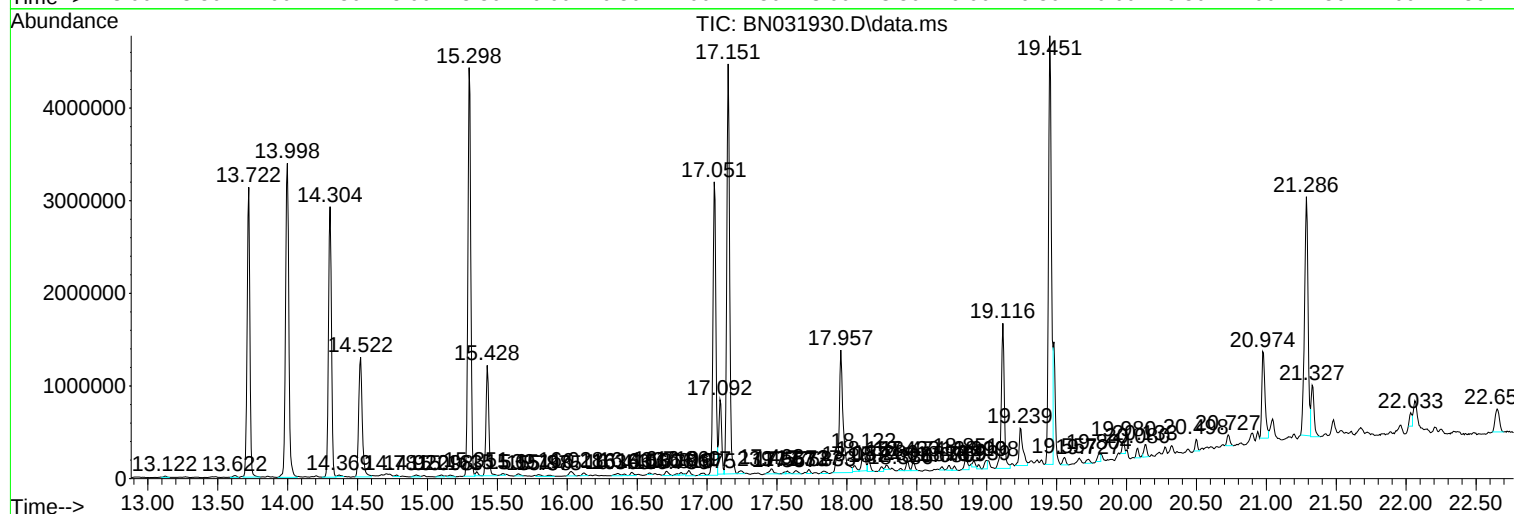
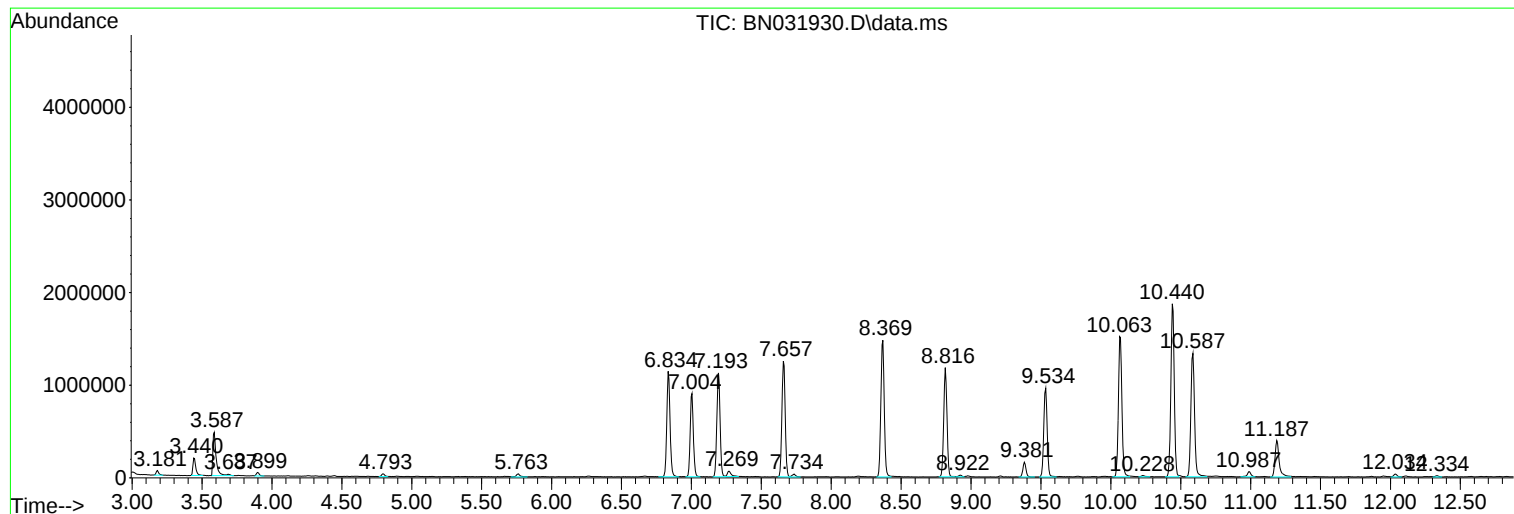
Sum of corrected areas: 95099123

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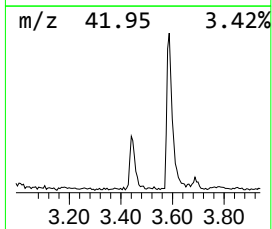
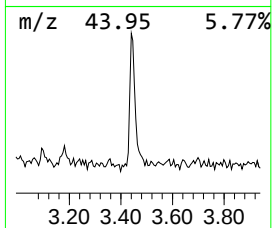
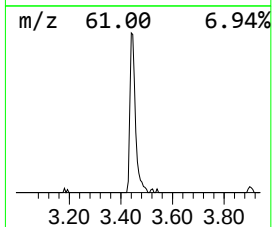
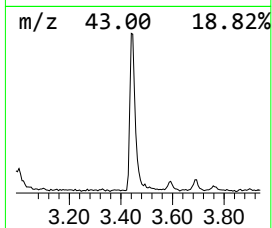
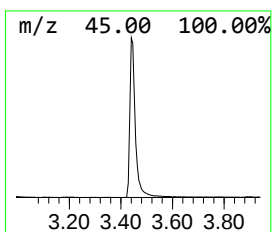
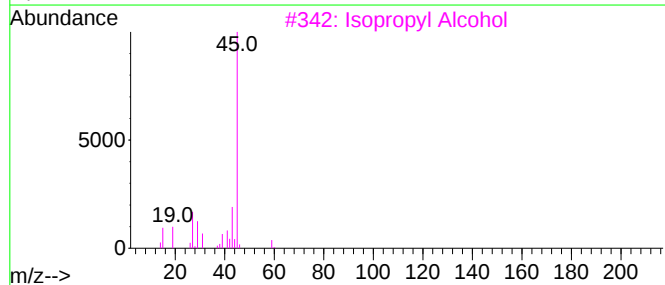
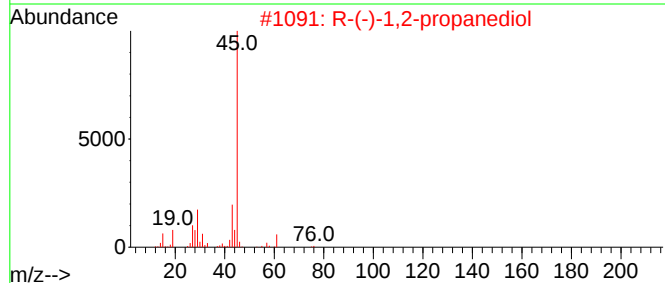
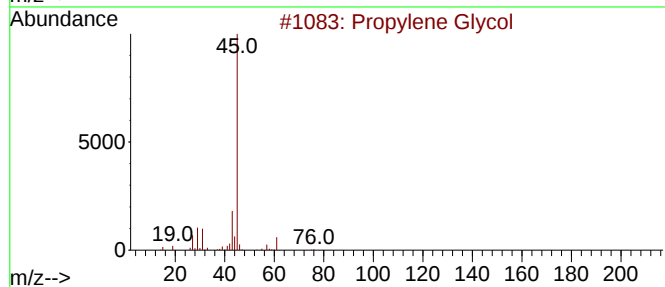
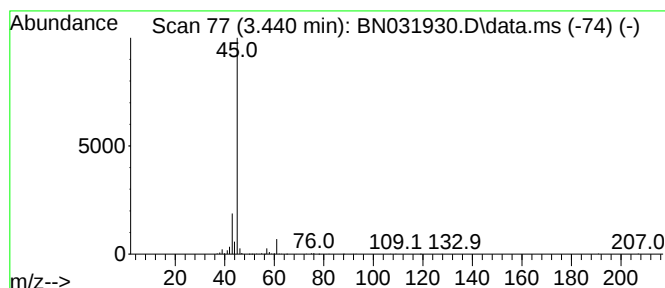
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.440	2.64 ng/ul	265904	1,4-Dichlorobenzene-d4	7.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	83
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
4			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40
5			(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	40



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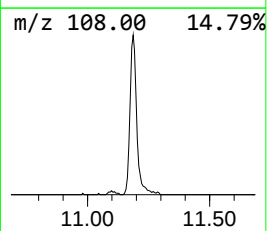
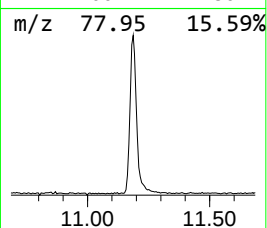
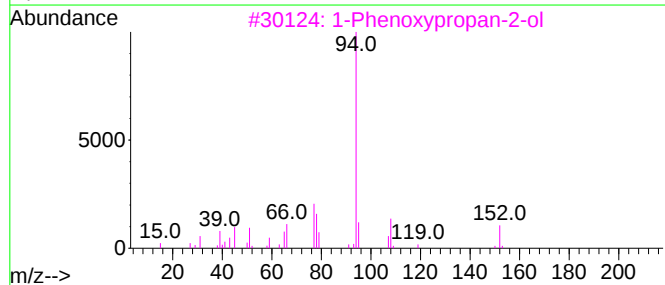
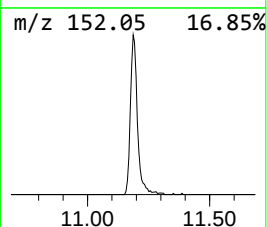
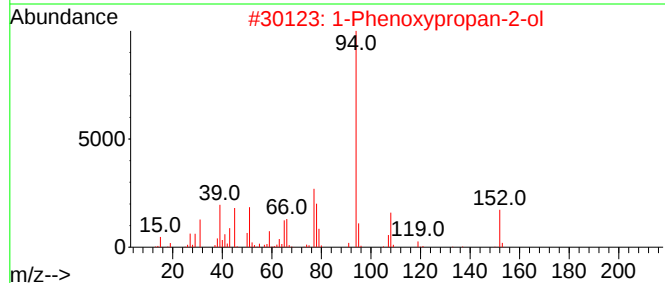
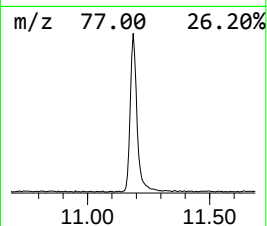
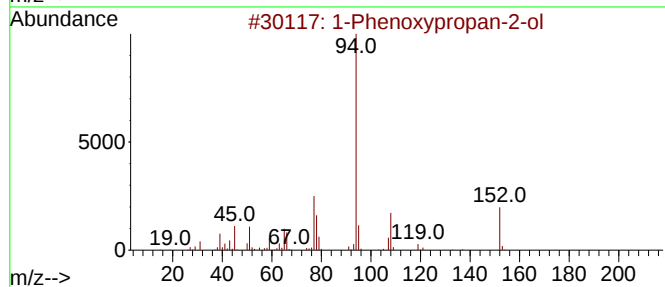
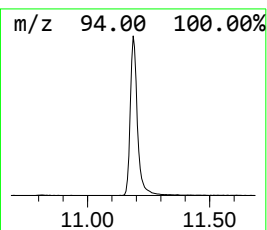
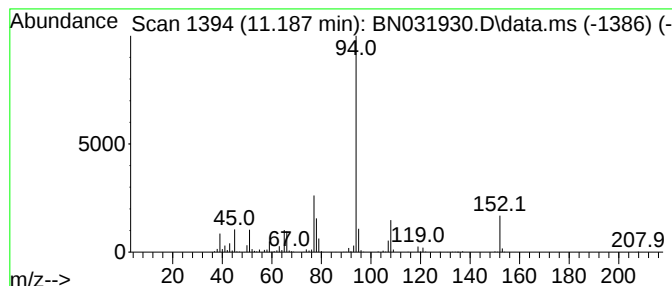
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN052924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	5.08 ng/ul	808702	Naphthalene-d8	10.440

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	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83
5			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	64



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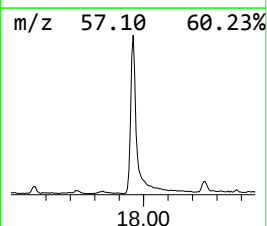
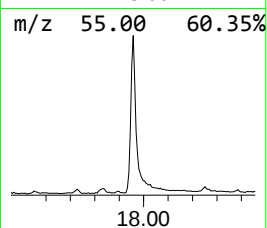
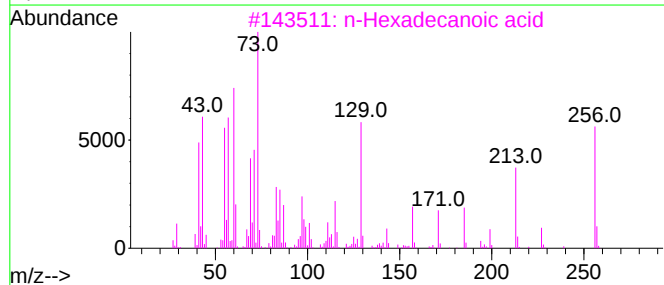
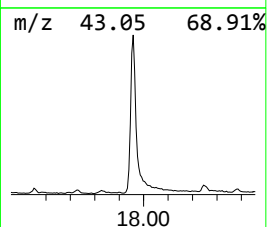
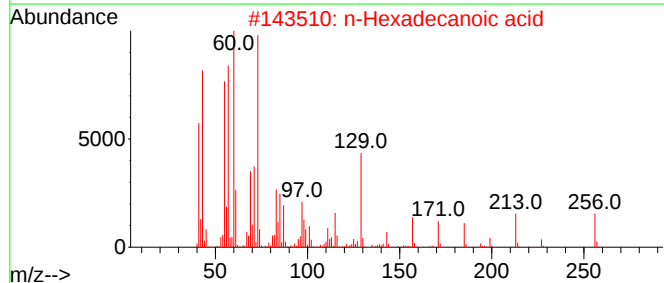
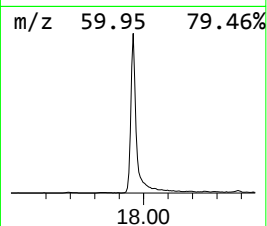
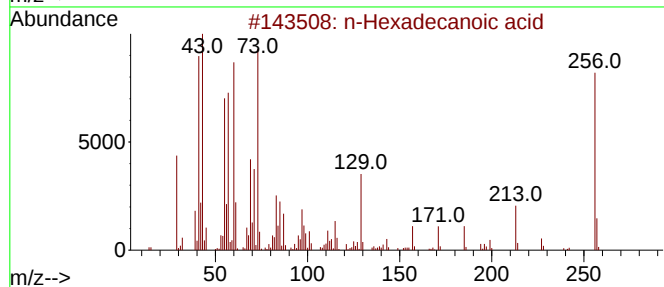
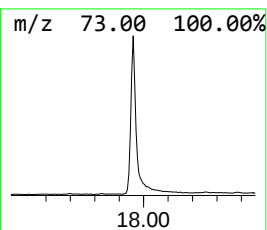
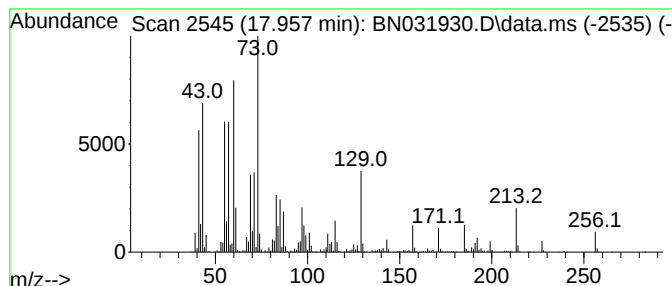
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.957	11.05 ng/ul	2589560	Phenanthrene-d10	17.051

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031930.D
Acq On : 12 Jun 2024 04:09
Operator : MA/JU
Sample : P2659-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHC33

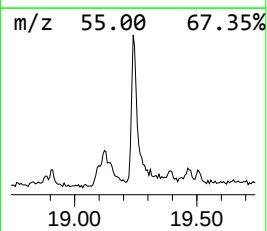
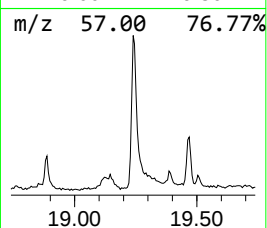
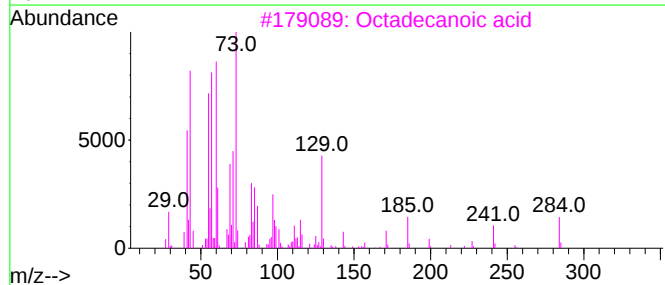
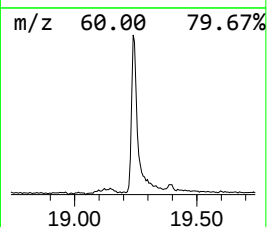
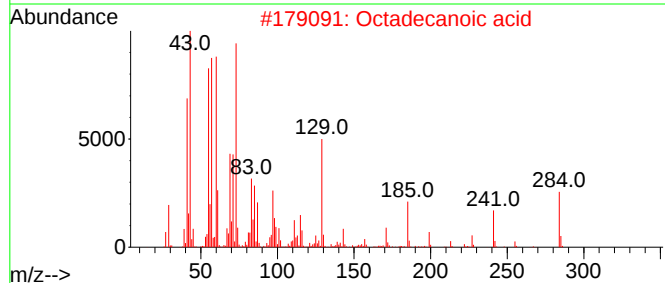
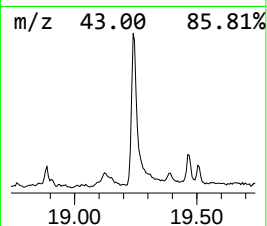
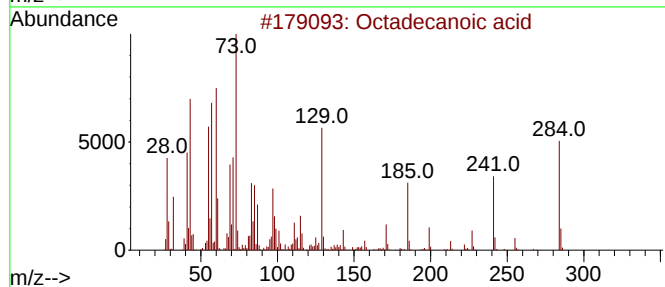
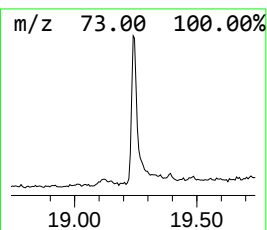
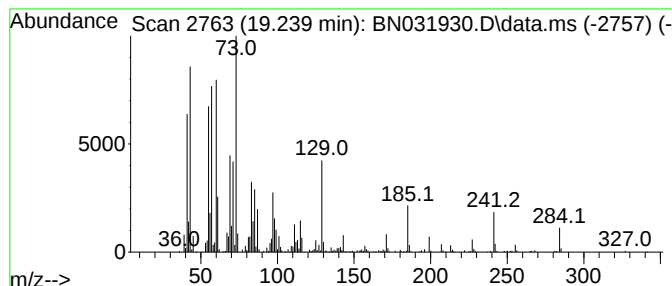
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.239	3.28 ng/ul	761993	Chrysene-d12	21.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031930.D
Acq On : 12 Jun 2024 04:09
Operator : MA/JU
Sample : P2659-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHC33

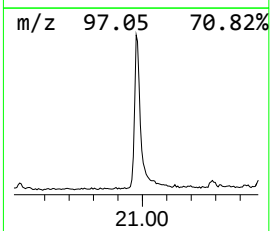
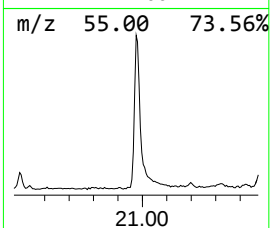
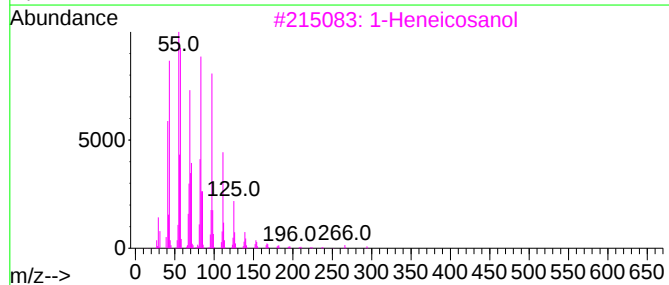
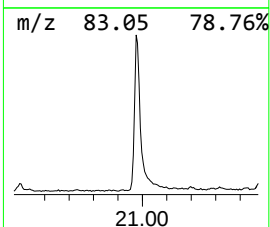
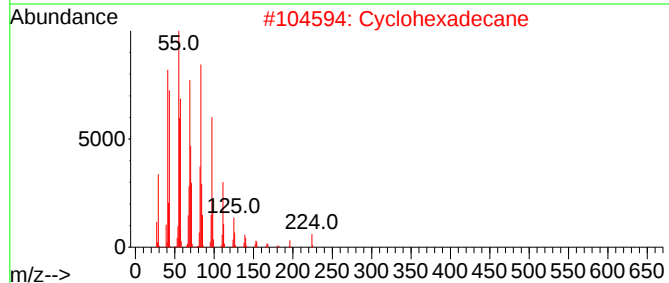
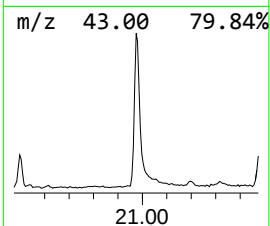
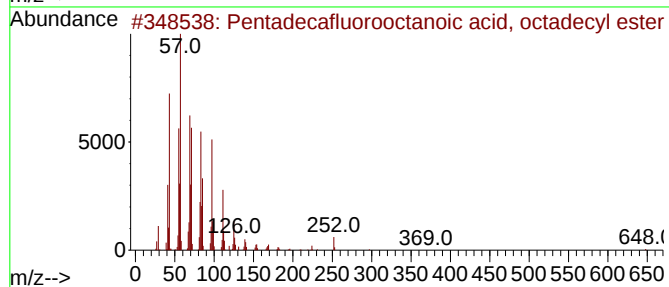
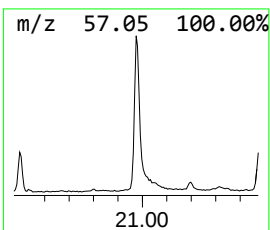
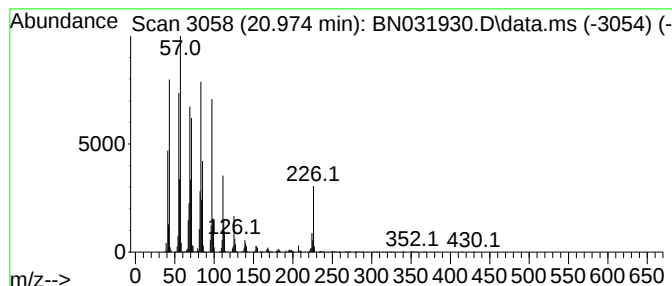
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecafluorooctanoic aci... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.974	6.44 ng/ul	1498100	Chrysene-d12	21.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2			Cyclohexadecane	224	C16H32	000295-65-8	93
3			1-Heneicosanol	312	C21H44O	015594-90-8	92
4			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	87
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031930.D
Acq On : 12 Jun 2024 04:09
Operator : MA/JU
Sample : P2659-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHC33

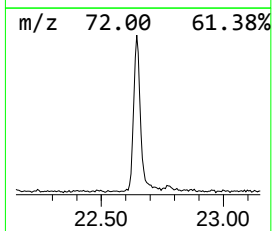
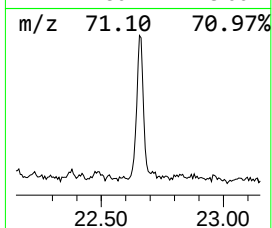
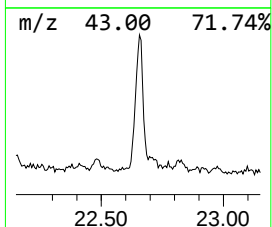
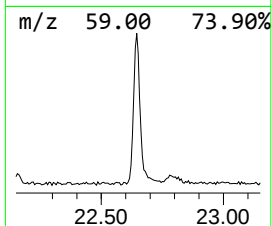
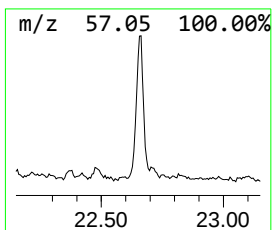
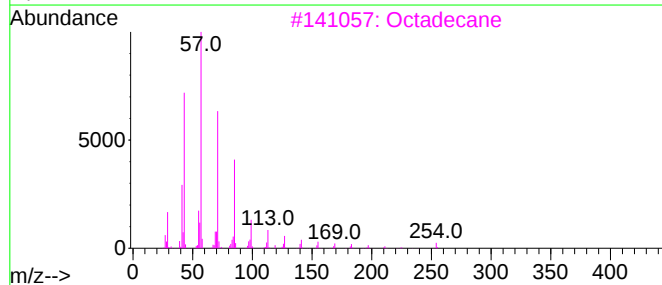
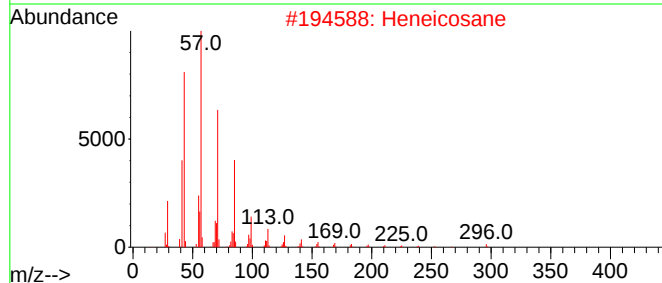
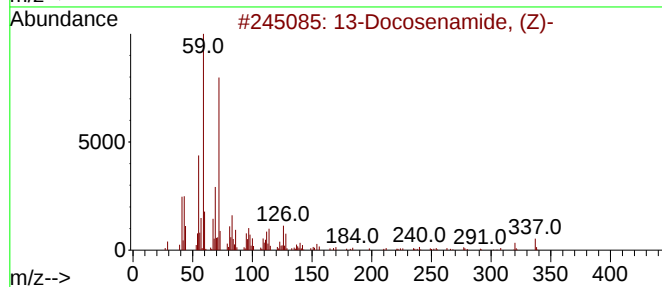
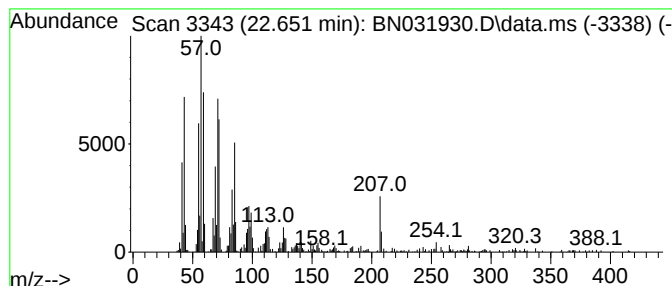
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 13-Docosenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.651	2.28 ng/ul	530808	Chrysene-d12	21.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	95
2		Heneicosane	296	C21H44	000629-94-7	66
3		Octadecane	254	C18H38	000593-45-3	64
4		Heptadecane	240	C17H36	000629-78-7	64
5		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061224\
Data File : BN031930.D
Acq On : 12 Jun 2024 04:09
Operator : MA/JU
Sample : P2659-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BHC33

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	3.440	2.6	ng/ul	265904	1	7.657	2017740	20.0
1-Phenoxypropan...	11.187	5.1	ng/ul	808702	2	10.440	3185690	20.0
n-Hexadecanoic ...	17.957	11.1	ng/ul	2589560	4	17.051	4686720	20.0
Octadecanoic acid	19.239	3.3	ng/ul	761993	5	21.286	4651230	20.0
Pentadecafluoro...	20.974	6.4	ng/ul	1498100	5	21.286	4651230	20.0
13-Docosenamide...	22.651	2.3	ng/ul	530808	5	21.286	4651230	20.0