

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
 Data File : BN012079.D
 Acq On : 06 Oct 2020 05:39
 Operator : CG/JU
 Sample : L4184-06
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AM1

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\SOM-EPA-BN100220.M
 Title : SVOA CALIBRATION

Signal : TIC: BN012079.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.193	32	35	46	rVB	87393	118001	1.86%	0.151%
2	3.469	79	82	88	rBV	40508	58616	0.92%	0.075%
3	3.911	153	157	160	rBV5	15642	22684	0.36%	0.029%
4	4.640	275	281	284	rBV5	5444	10347	0.16%	0.013%
5	6.652	619	623	626	rBV6	7265	12116	0.19%	0.016%
6	6.828	646	653	666	rBV	1343256	2096981	33.07%	2.690%
7	6.999	676	682	691	rBV	968003	1474991	23.26%	1.892%
8	7.193	708	715	724	rBV	1387998	2147401	33.86%	2.755%
9	7.269	724	728	741	rVV	156897	290892	4.59%	0.373%
10	7.375	744	746	753	rVB8	5538	11749	0.19%	0.015%
11	7.658	786	794	801	rBV	1129049	1738505	27.42%	2.230%
12	7.716	802	804	811	rVB6	16189	25571	0.40%	0.033%
13	7.969	841	847	848	rBV4	5276	9222	0.15%	0.012%
14	8.357	906	913	923	rBV	1560591	2443928	38.54%	3.135%
15	8.805	982	989	1001	rVB	1218421	1924220	30.34%	2.469%
16	8.928	1006	1010	1013	rVB6	8171	12230	0.19%	0.016%
17	8.975	1013	1018	1023	rBV8	9851	18383	0.29%	0.024%
18	9.234	1055	1062	1067	rVB5	14740	22091	0.35%	0.028%
19	9.522	1102	1111	1120	rBV	1007942	1648633	26.00%	2.115%
20	9.740	1147	1148	1155	rVB6	6313	10189	0.16%	0.013%
21	10.051	1194	1201	1210	rBV	1598102	2508490	39.56%	3.218%
22	10.428	1258	1265	1274	rBV	1383581	2324670	36.66%	2.982%
23	10.569	1281	1289	1301	rBV	1270395	2174413	34.29%	2.790%
24	10.946	1349	1353	1354	rBV2	7032	6986	0.11%	0.009%
25	10.957	1354	1355	1360	rVB4	7974	9058	0.14%	0.012%
26	12.022	1531	1536	1541	rBV4	23526	34979	0.55%	0.045%
27	12.645	1638	1642	1644	rBV4	7135	9284	0.15%	0.012%
28	12.798	1664	1668	1671	rBV4	4706	7643	0.12%	0.010%
29	12.904	1680	1686	1690	rVB6	10343	12939	0.20%	0.017%
30	13.193	1731	1735	1744	rVV2	47154	74618	1.18%	0.096%
31	13.710	1816	1823	1827	rBV	2410793	3382037	53.33%	4.339%
32	13.751	1827	1830	1840	rVB	1028950	1419167	22.38%	1.821%
33	13.987	1860	1870	1882	rBV	2837912	4110888	64.83%	5.274%
34	14.204	1904	1907	1912	rVB4	8316	12997	0.20%	0.017%
35	14.292	1916	1922	1930	rVV	2042357	2980558	47.00%	3.824%
36	14.492	1948	1956	1970	rBV	1111838	1601870	25.26%	2.055%

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Integration Parameters: LSCINT.P

Integrator: RTE
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37	14.716	1992	1994	2001	rVB7	14332	20466	0.32%	0.026%
38	15.128	2054	2064	2067	rBV2	54338	83631	1.32%	0.107%
39	15.222	2077	2080	2083	rVB4	5286	7002	0.11%	0.009%
40	15.292	2085	2092	2106	rVV	3533453	4951253	78.08%	6.352%
41	15.410	2106	2112	2124	rVV	1111817	1547002	24.40%	1.985%
42	15.528	2127	2132	2138	rBV2	42629	62914	0.99%	0.081%
43	15.657	2153	2154	2158	rVB4	7359	6868	0.11%	0.009%
44	15.851	2183	2187	2188	rBV4	9701	9371	0.15%	0.012%
45	15.869	2188	2190	2194	rVB5	9221	10334	0.16%	0.013%
46	15.922	2194	2199	2202	rBV4	9512	13230	0.21%	0.017%
47	16.028	2213	2217	2219	rBV5	6785	8359	0.13%	0.011%
48	16.075	2219	2225	2231	rVB7	16554	27372	0.43%	0.035%
49	16.192	2242	2245	2249	rVB6	8307	13611	0.21%	0.017%
50	16.316	2260	2266	2267	rBV4	5082	8966	0.14%	0.012%
51	16.451	2286	2289	2298	rVB7	44892	64014	1.01%	0.082%
52	16.545	2303	2305	2312	rVB5	8050	15055	0.24%	0.019%
53	16.828	2347	2353	2357	rBV6	12807	32619	0.51%	0.042%
54	17.039	2382	2389	2395	rBV	2422228	3346018	52.77%	4.293%
55	17.139	2400	2406	2414	rVV	3722014	5087632	80.23%	6.527%
56	17.222	2417	2420	2425	rVV4	44273	58764	0.93%	0.075%
57	17.339	2435	2440	2444	rBV	58158	78417	1.24%	0.101%
58	17.386	2444	2448	2454	rBV	115895	133032	2.10%	0.171%
59	17.557	2474	2477	2480	rVB5	7755	10124	0.16%	0.013%
60	17.716	2502	2504	2510	rBV7	9162	12775	0.20%	0.016%
61	17.816	2516	2521	2528	rBV3	79974	122322	1.93%	0.157%
62	17.945	2538	2543	2553	rBV	674264	1029859	16.24%	1.321%
63	18.369	2612	2615	2617	rBV4	9167	7254	0.11%	0.009%
64	18.622	2655	2658	2661	rBV4	16367	17479	0.28%	0.022%
65	18.663	2661	2665	2669	rVV	47285	61066	0.96%	0.078%
66	18.704	2669	2672	2678	rVB	114288	129045	2.04%	0.166%
67	18.886	2700	2703	2708	rVB5	23683	31907	0.50%	0.041%
68	19.080	2731	2736	2738	rBV3	48802	72715	1.15%	0.093%
69	19.104	2738	2740	2747	rVV8	41049	72228	1.14%	0.093%
70	19.233	2757	2762	2772	rBV2	300703	455952	7.19%	0.585%
71	19.327	2775	2778	2783	rVB	43533	61732	0.97%	0.079%
72	19.439	2791	2797	2806	rBV	4232054	5733489	90.42%	7.356%
73	19.510	2806	2809	2815	rVB4	51962	74013	1.17%	0.095%
74	19.792	2853	2857	2860	rBV	561153	617206	9.73%	0.792%
75	19.833	2860	2864	2869	rVB	1217259	1296865	20.45%	1.664%
76	19.969	2884	2887	2889	rBV4	33241	39914	0.63%	0.051%

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77	20.004	2891	2893	2897	rVB2	43649	46828	0.74%	0.060%
78	20.380	2951	2957	2963	rBV	325893	424905	6.70%	0.545%
79	20.504	2976	2978	2982	rVB2	79811	70701	1.11%	0.091%
80	20.774	3020	3024	3026	rBV	204667	210005	3.31%	0.269%
81	20.804	3026	3029	3034	rVB	440324	474551	7.48%	0.609%
82	20.963	3051	3056	3064	rBV	926037	1333471	21.03%	1.711%
83	21.239	3098	3103	3112	rBV	2915380	3693079	58.24%	4.738%
84	21.404	3128	3131	3134	rBV	143927	154753	2.44%	0.199%
85	21.674	3174	3177	3180	rVB2	207279	214816	3.39%	0.276%
86	21.833	3201	3204	3211	rBV4	161915	258784	4.08%	0.332%
87	22.416	3299	3303	3310	rVB	534551	688879	10.86%	0.884%
88	23.304	3448	3454	3467	rVB	3378775	6341239	100.00%	8.136%
89	23.445	3472	3478	3485	rVB	2108660	3834045	60.46%	4.919%

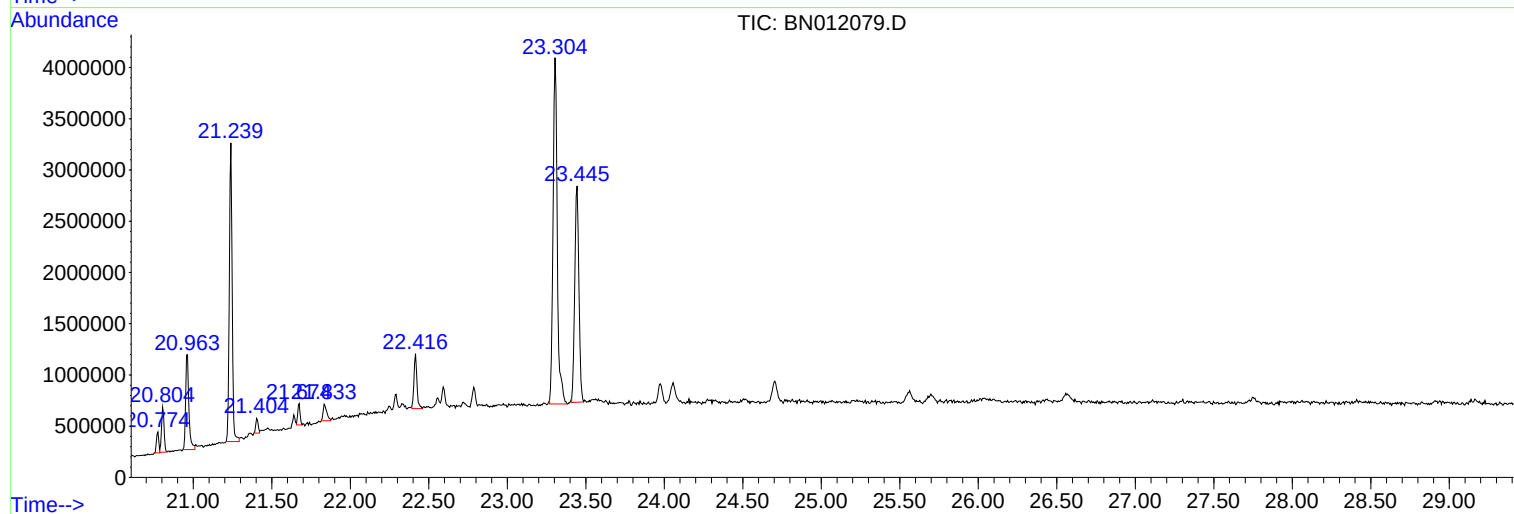
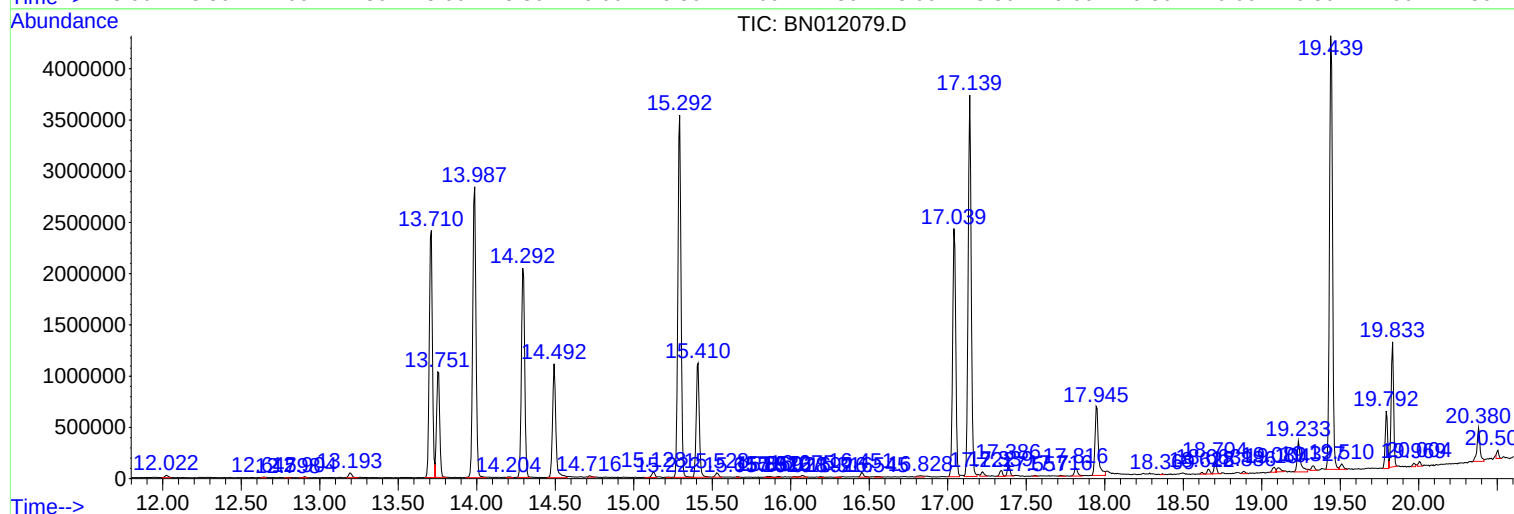
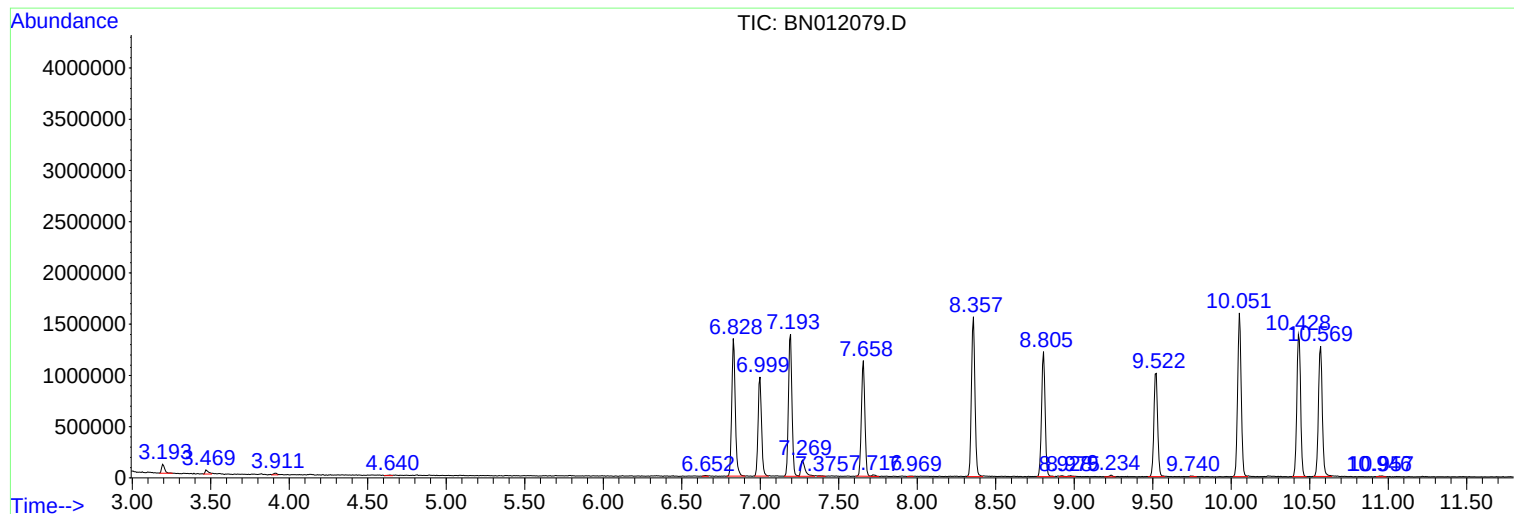
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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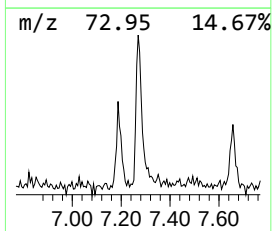
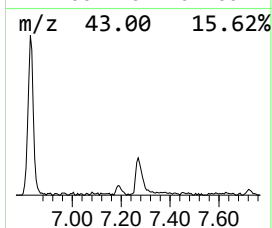
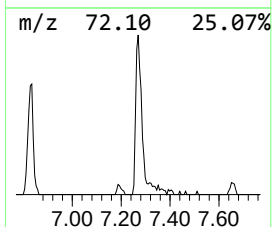
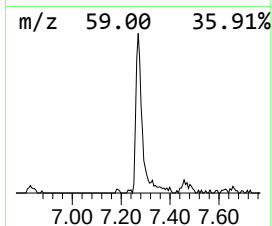
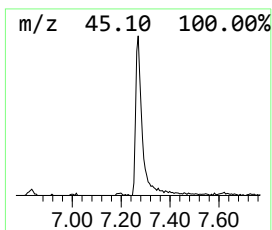
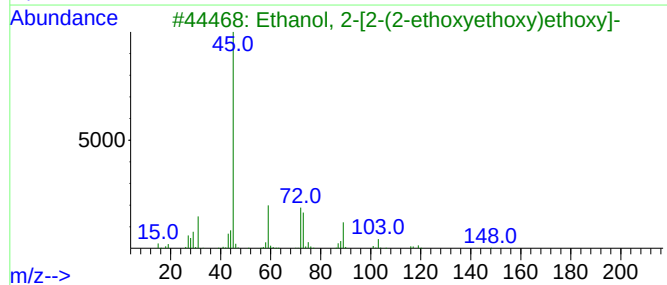
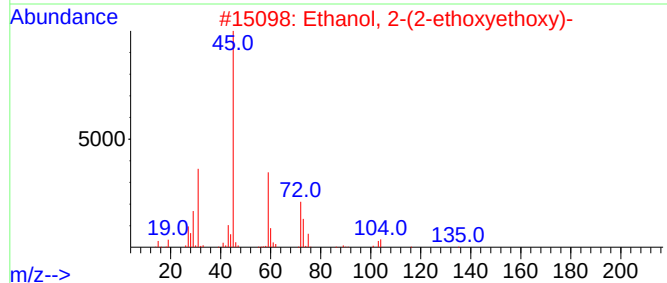
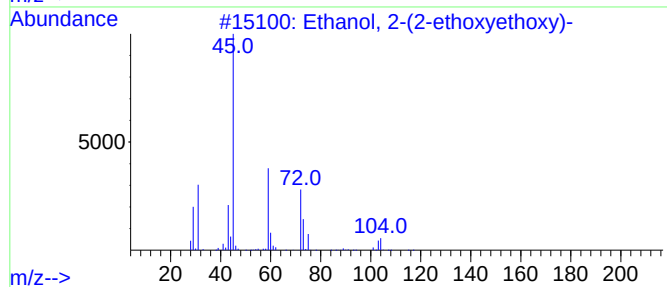
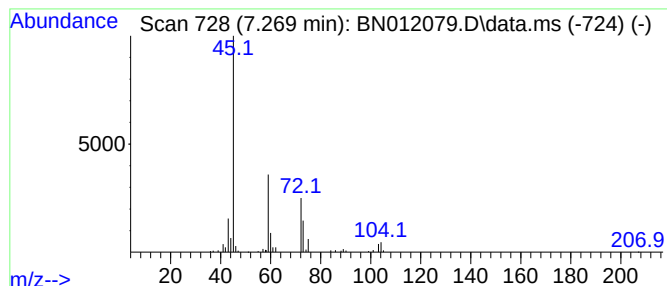
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.270	3.35 ng/ul	290892	1,4-Dichlorobenzene-d4	7.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	90
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
3			Ethanol, 2-[2-(2-ethoxyethoxy)et...	178	C8H18O4	000112-50-5	56
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	47
5			2-Butanol, 3-methyl-, (S)-	88	C5H12O	001517-66-4	47



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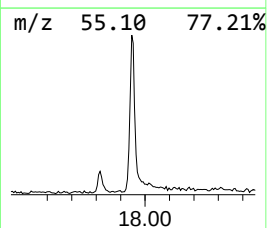
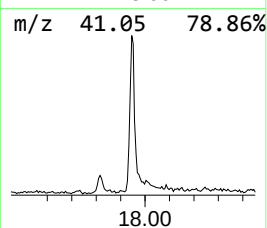
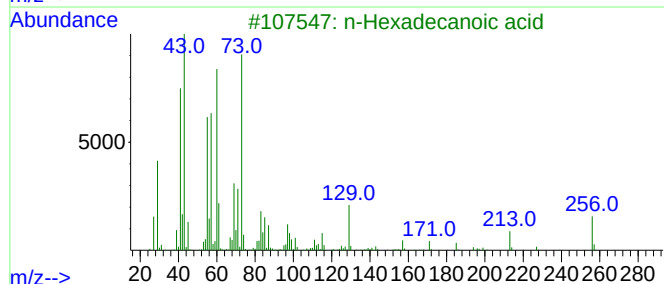
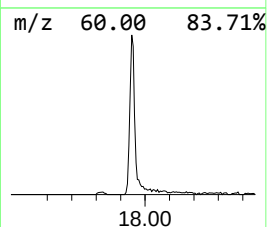
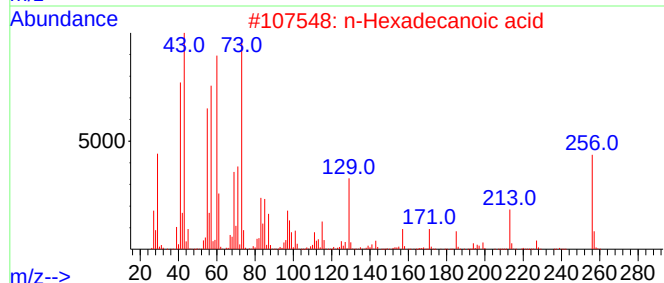
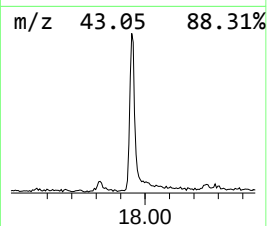
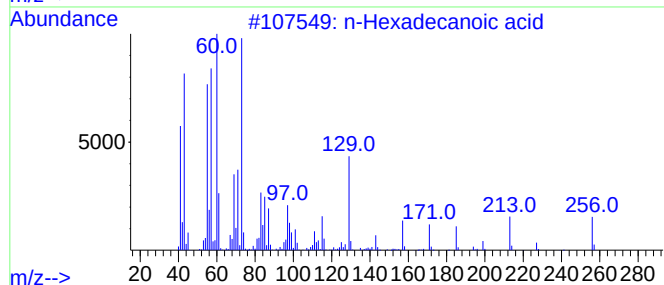
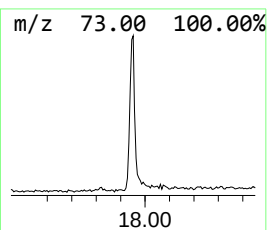
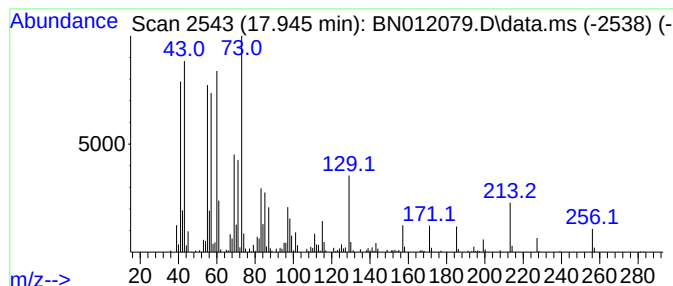
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.950	6.16 ng/ul	1029860	Phenanthrene-d10	17.039

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	87
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	83



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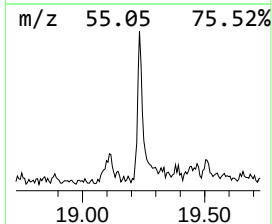
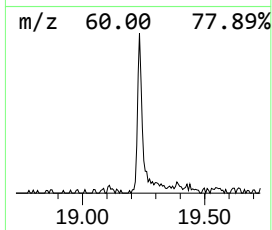
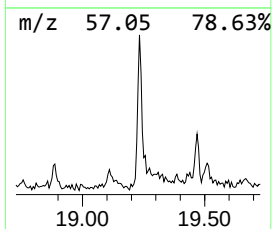
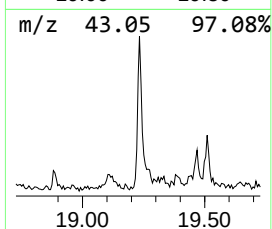
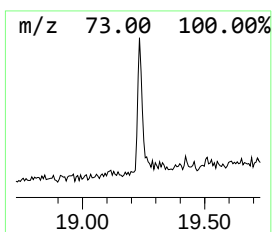
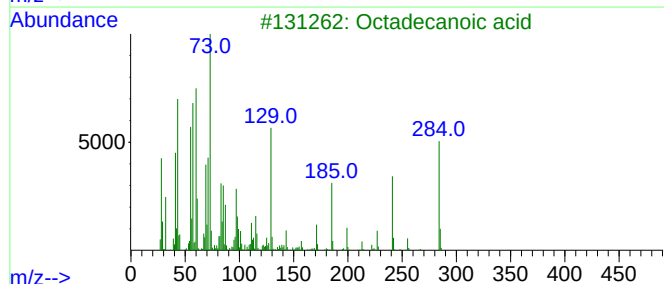
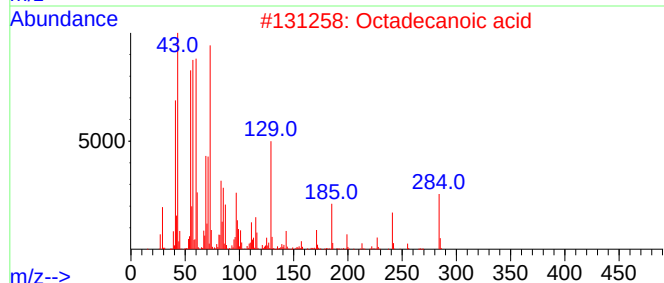
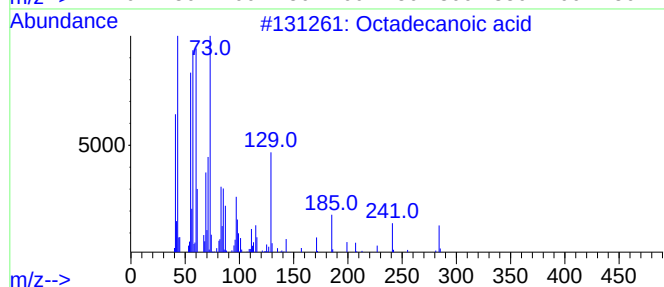
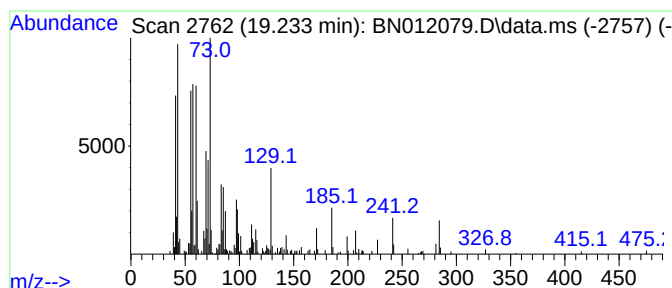
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.230	2.47 ng/ul	455952	Chrysene-d12	21.239

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	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



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C0AM1

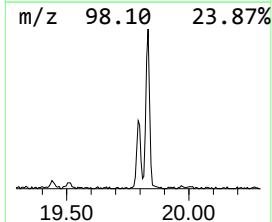
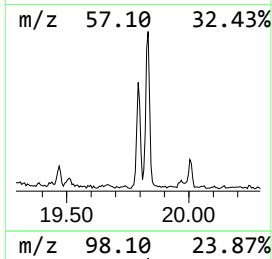
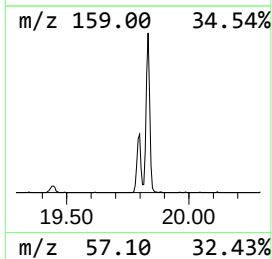
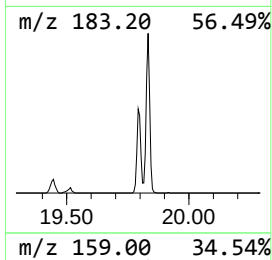
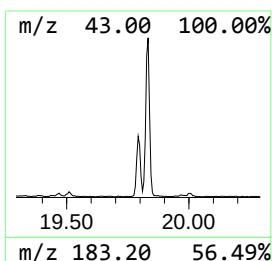
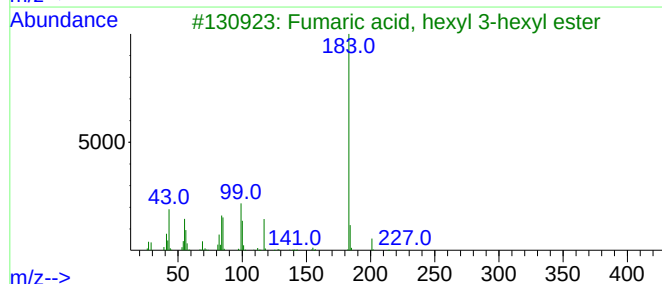
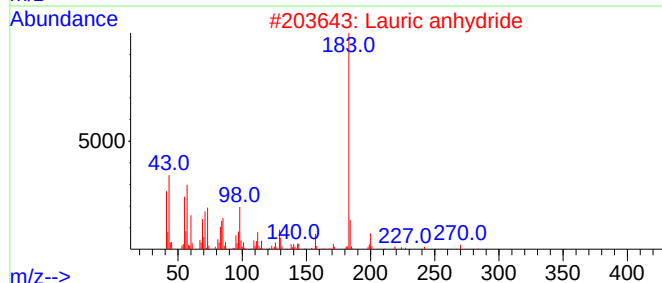
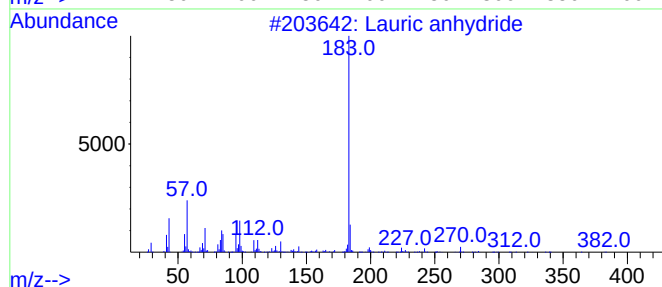
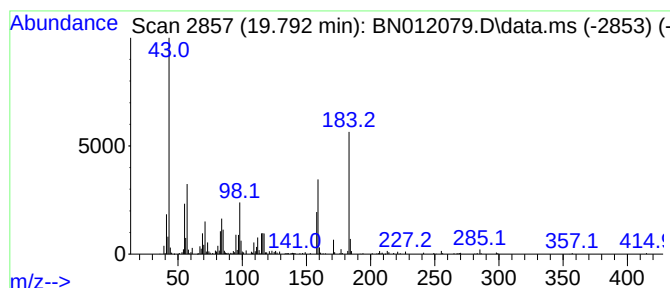
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.790	3.34 ng/ul	617206	Chrysene-d12	21.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lauric anhydride	382	C24H46O3	000645-66-9	38
2			Lauric anhydride	382	C24H46O3	000645-66-9	32
3			Fumaric acid, hexyl 3-hexyl ester	284	C16H28O4	1000348-60-5	32
4			Lauric acid, 3,4-dichlorophenyl ...	344	C18H26Cl2O2	1000357-92-7	27
5			Fumaric acid, cyclobutyl isohexy...	254	C14H22O4	1000345-03-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
Data File : BN012079.D
Acq On : 06 Oct 2020 05:39
Operator : CG/JU
Sample : L4184-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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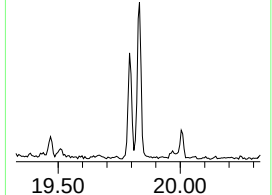
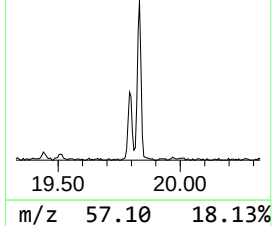
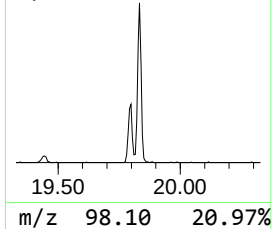
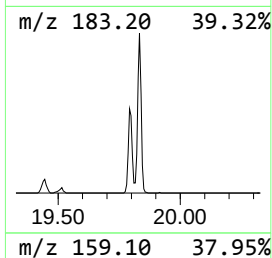
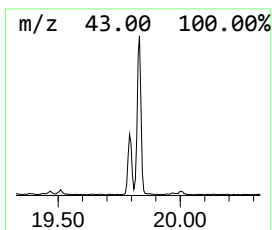
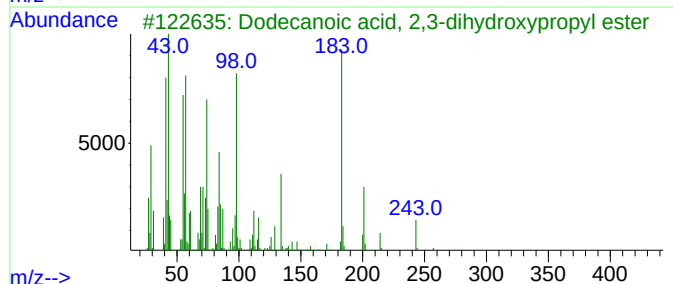
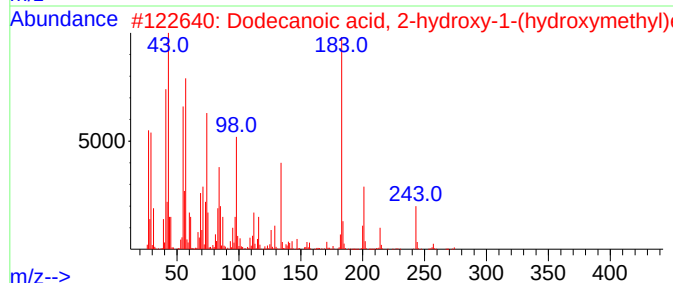
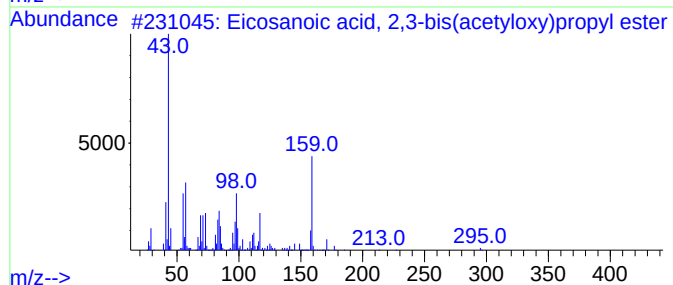
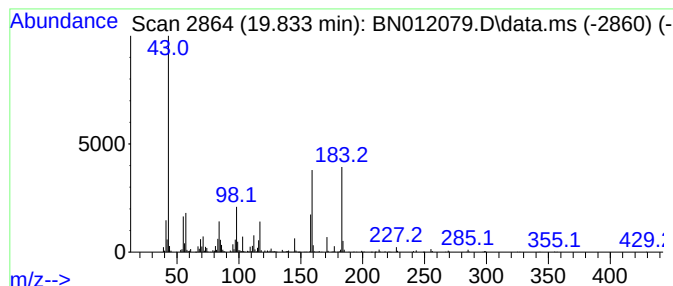
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Eicosanoic acid, 2,3-bis(ac... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.830	7.02 ng/ul	1296870	Chrysene-d12	21.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	50
2			Dodecanoic acid, 2-hydroxy-1-(hy...	274	C15H30O4	001678-45-1	47
3			Dodecanoic acid, 2,3-dihydroxypr...	274	C15H30O4	000142-18-7	35
4			Dodecanoyl chloride	218	C12H23ClO	000112-16-3	32
5			Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
Data File : BN012079.D
Acq On : 06 Oct 2020 05:39
Operator : CG/JU
Sample : L4184-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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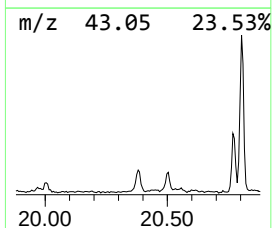
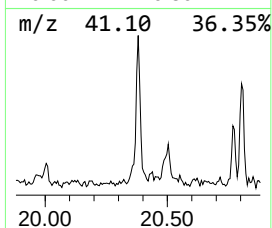
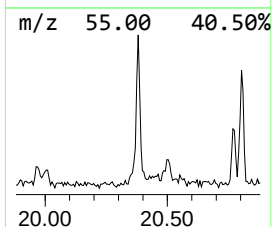
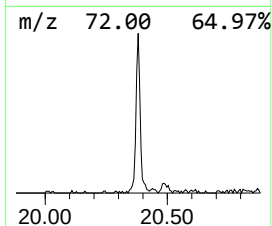
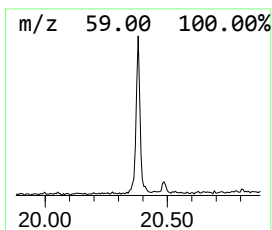
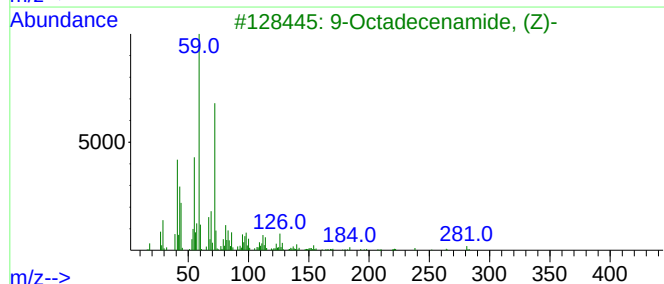
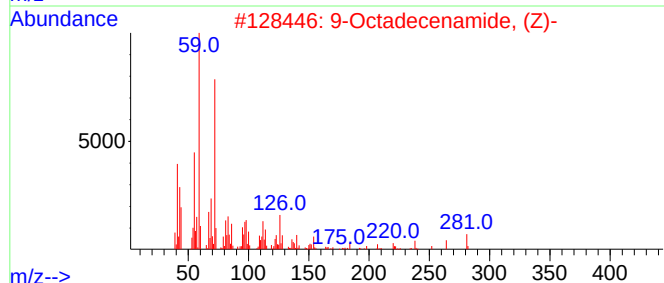
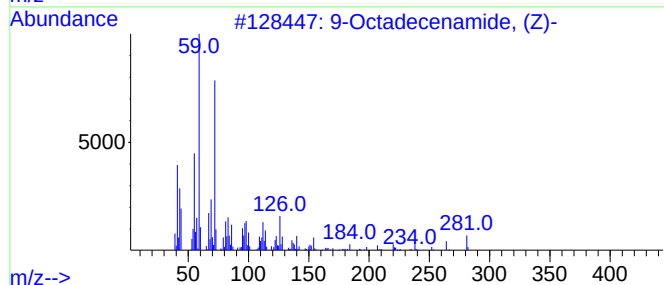
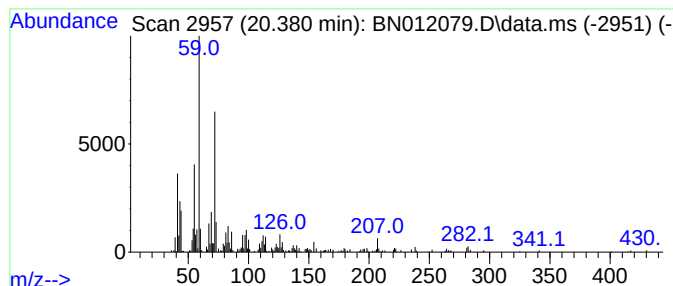
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.380	2.30 ng/ul	424905	Chrysene-d12	21.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	97
2	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	97
3	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	95
4	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	94
5	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
Data File : BN012079.D
Acq On : 06 Oct 2020 05:39
Operator : CG/JU
Sample : L4184-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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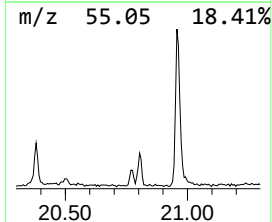
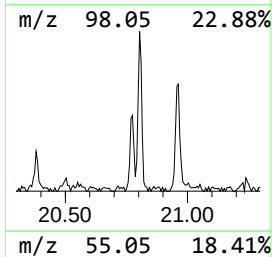
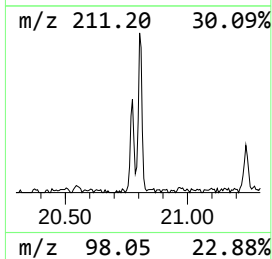
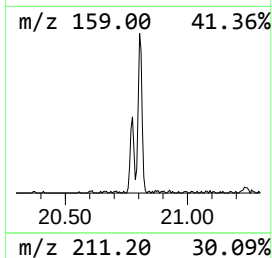
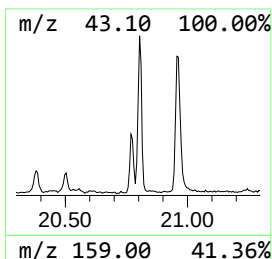
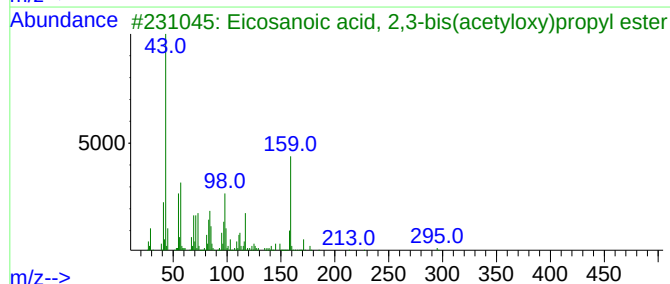
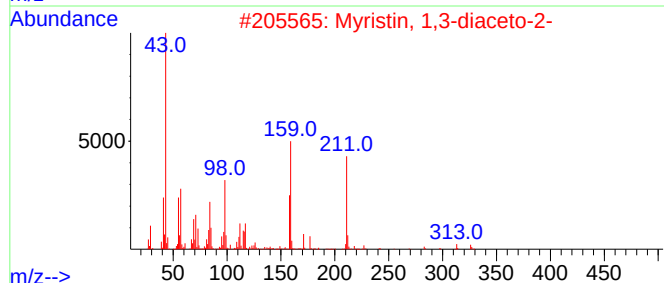
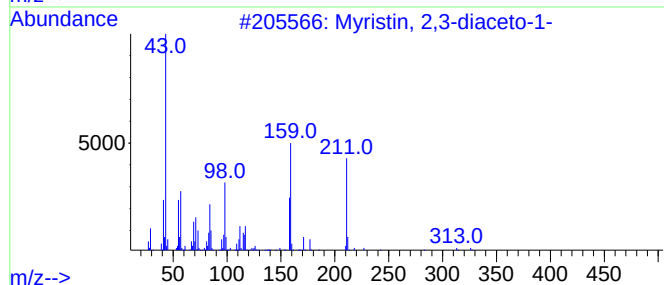
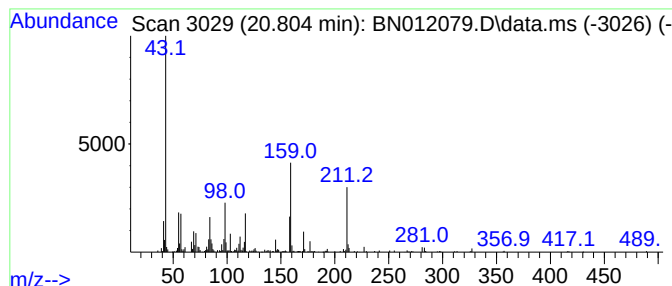
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Myristin, 2,3-diaceto-1- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.800	2.57 ng/ul	474551	Chrysene-d12	21.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	72
2			Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	38
3			Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	16
4			Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	10
5			9-Octadecenoic acid (Z)-, 2-(ace...	440	C25H44O6	055401-63-3	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
Data File : BN012079.D
Acq On : 06 Oct 2020 05:39
Operator : CG/JU
Sample : L4184-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AM1

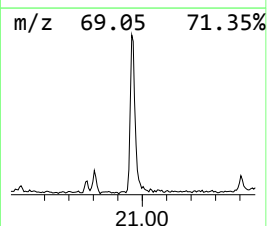
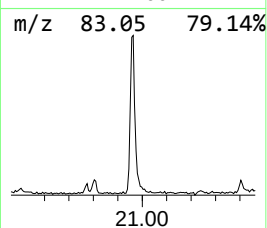
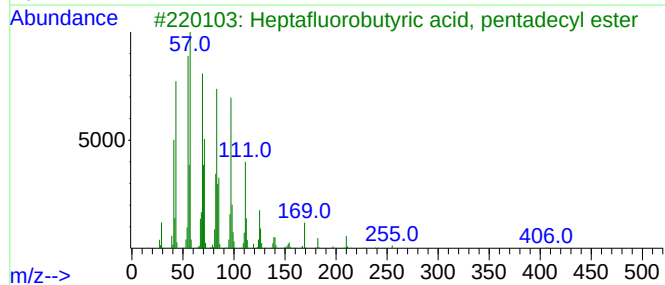
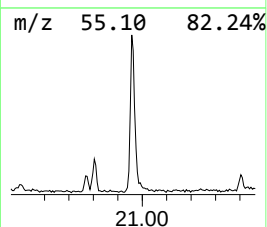
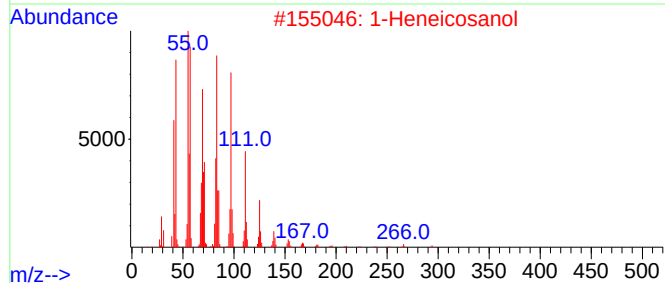
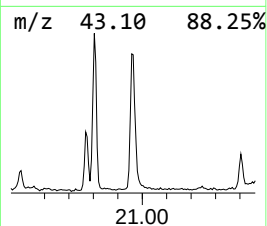
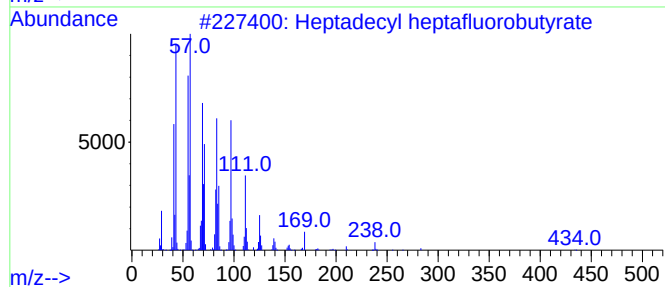
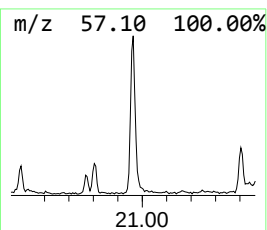
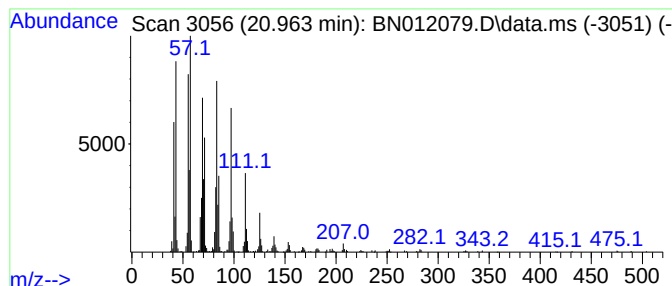
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.960	7.22 ng/ul	1333470	Chrysene-d12	21.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	93
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91
4			Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	91
5			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
Data File : BN012079.D
Acq On : 06 Oct 2020 05:39
Operator : CG/JU
Sample : L4184-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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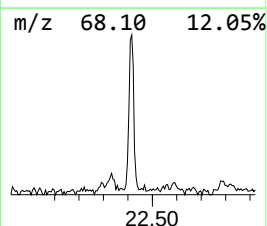
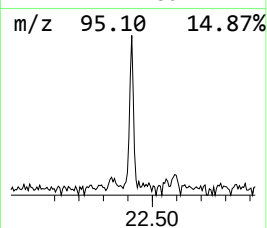
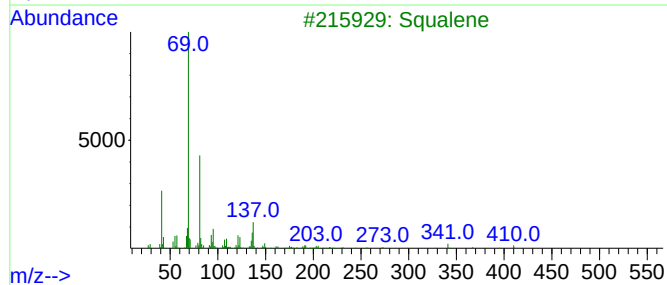
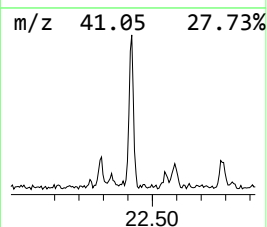
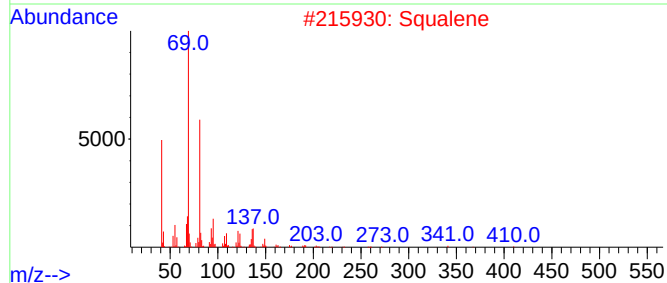
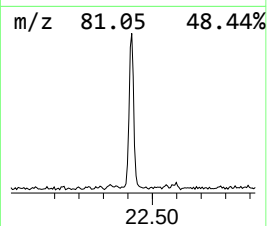
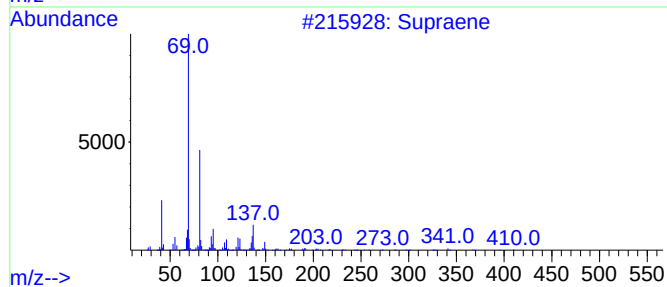
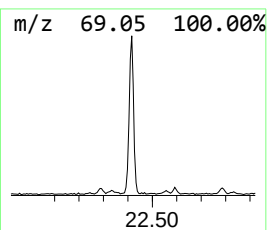
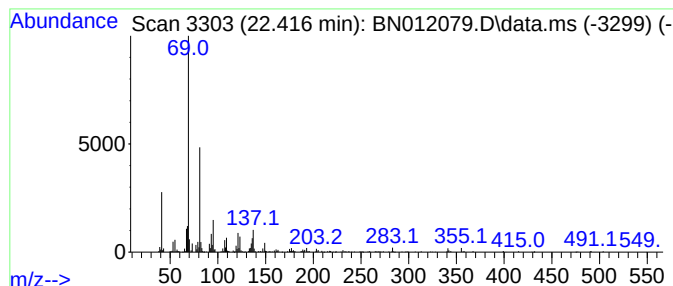
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Supraene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.420	3.59 ng/ul	688879	Perylene-d12	23.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	87
2			Squalene	410	C30H50	000111-02-4	87
3			Squalene	410	C30H50	000111-02-4	83
4			(.+/-)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	72
5			Squalene	410	C30H50	000111-02-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100520\
Data File : BN012079.D
Acq On : 06 Oct 2020 05:39
Operator : CG/JU
Sample : L4184-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN100220.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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n-Hexadecanoic ...	17.950	6.2	ng/ul	1029860	4	17.039	3346020	20.0
Octadecanoic acid	19.230	2.5	ng/ul	455952	5	21.239	3693080	20.0
unknown-01	19.790	3.3	ng/ul	617206	5	21.239	3693080	20.0
Eicosanoic acid...	19.830	7.0	ng/ul	1296870	5	21.239	3693080	20.0
9-Octadecenamid...	20.380	2.3	ng/ul	424905	5	21.239	3693080	20.0
Myristin, 2,3-d...	20.800	2.6	ng/ul	474551	5	21.239	3693080	20.0
Heptadecyl hept...	20.960	7.2	ng/ul	1333470	5	21.239	3693080	20.0
Supraene	22.420	3.6	ng/ul	688879	6	23.445	3834050	20.0