

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
 Data File : BN010096.D
 Acq On : 05 Mar 2020 01:51
 Operator : CG/JU
 Sample : L1641-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB6Y5

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 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-BN021220MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	67	71	76	rVB2	12789	14899	0.54%	0.034%
2	3.493	81	86	95	rBV	47531	69631	2.51%	0.158%
3	3.687	115	119	123	rVB	41552	56887	2.05%	0.129%
4	3.905	153	156	161	rVB6	10067	16062	0.58%	0.036%
5	4.046	174	180	184	rBV5	11024	17176	0.62%	0.039%
6	4.216	205	209	217	rVB7	7435	13994	0.50%	0.032%
7	4.563	264	268	274	rBV	16033	22614	0.82%	0.051%
8	5.428	409	415	420	rVB5	14129	21722	0.78%	0.049%
9	5.481	420	424	430	rVV5	8447	16486	0.59%	0.037%
10	6.563	602	608	622	rVV	500492	872094	31.44%	1.977%
11	6.716	629	634	647	rVV	376232	613109	22.10%	1.390%
12	6.899	659	665	676	rVB	533340	858971	30.97%	1.947%
13	6.993	678	681	689	rVB4	9650	16873	0.61%	0.038%
14	7.075	689	695	701	rBV6	8427	17355	0.63%	0.039%
15	7.357	736	743	753	rBV	404383	667624	24.07%	1.514%
16	8.075	859	865	880	rBV	397338	724405	26.12%	1.642%
17	8.499	931	937	947	rBV	518102	919505	33.15%	2.085%
18	8.675	962	967	976	rVB3	29270	53012	1.91%	0.120%
19	8.934	1007	1011	1015	rVB	15487	20670	0.75%	0.047%
20	9.210	1052	1058	1076	rBV	453311	803974	28.98%	1.823%
21	9.746	1143	1149	1171	rVV	531528	1038371	37.44%	2.354%
22	10.098	1200	1209	1216	rBV	590148	982084	35.41%	2.226%
23	10.263	1229	1237	1260	rBV2	263166	652728	23.53%	1.480%
24	12.839	1670	1675	1680	rBV3	9992	16007	0.58%	0.036%
25	13.428	1769	1775	1781	rBV	1057335	1522848	54.90%	3.452%
26	13.481	1781	1784	1793	rVB2	46106	75924	2.74%	0.172%
27	13.687	1812	1819	1832	rBV	1259026	1880184	67.78%	4.262%
28	13.998	1866	1872	1879	rVV	843469	1232378	44.43%	2.794%
29	14.063	1879	1883	1890	rVB2	22777	34348	1.24%	0.078%
30	14.263	1911	1917	1932	rBV	318165	709600	25.58%	1.609%
31	14.422	1938	1944	1949	rBV5	23539	44067	1.59%	0.100%
32	15.010	2037	2044	2050	rBV	1501416	2177299	78.50%	4.936%
33	15.175	2066	2072	2082	rBV	476179	815110	29.39%	1.848%
34	15.251	2082	2085	2088	rVB4	17598	23516	0.85%	0.053%

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
 Title : SVOA CALIBRATION

35	15.510	2125	2129	2137	rVV4	8110	15195	0.55%	0.034%
36	15.763	2166	2172	2174	rBV6	14751	22651	0.82%	0.051%
37	15.792	2174	2177	2182	rVB4	11796	17821	0.64%	0.040%
38	16.322	2262	2267	2274	rVV2	46331	75267	2.71%	0.171%
39	16.433	2280	2286	2292	rVB5	15067	28792	1.04%	0.065%
40	16.580	2309	2311	2316	rVB	12814	15731	0.57%	0.036%
41	16.763	2336	2342	2346	rBV	984070	1388453	50.06%	3.148%
42	16.804	2346	2349	2353	rVV	333898	424673	15.31%	0.963%
43	16.863	2353	2359	2372	rVB	1470056	2170855	78.26%	4.921%
44	17.186	2406	2414	2422	rBV3	27002	60824	2.19%	0.138%
45	17.292	2429	2432	2436	rVB5	12724	15646	0.56%	0.035%
46	17.645	2486	2492	2495	rVV2	48281	66431	2.39%	0.151%
47	17.692	2495	2500	2506	rVB2	67000	97360	3.51%	0.221%
48	17.775	2508	2514	2518	rVB4	15832	27968	1.01%	0.063%
49	17.833	2519	2524	2528	rBV4	68437	107505	3.88%	0.244%
50	17.875	2528	2531	2538	rVB	32508	44842	1.62%	0.102%
51	17.986	2543	2550	2551	rBV3	15046	20949	0.76%	0.047%
52	18.151	2575	2578	2583	rBV2	31320	44683	1.61%	0.101%
53	18.204	2583	2587	2592	rVB5	20816	26513	0.96%	0.060%
54	18.392	2615	2619	2626	rVB3	27491	49703	1.79%	0.113%
55	18.457	2626	2630	2633	rBV4	20695	32798	1.18%	0.074%
56	18.492	2633	2636	2640	rVB4	13259	17244	0.62%	0.039%
57	18.580	2645	2651	2656	rBV5	57293	105874	3.82%	0.240%
58	18.633	2656	2660	2664	rVV3	36511	48672	1.75%	0.110%
59	18.669	2664	2666	2669	rVB3	14532	15315	0.55%	0.035%
60	18.727	2670	2676	2684	rBV6	32513	65070	2.35%	0.148%
61	18.833	2686	2694	2704	rBV	554716	819688	29.55%	1.858%
62	19.063	2728	2733	2735	rBV5	26674	44642	1.61%	0.101%
63	19.174	2746	2752	2755	rBV	1971218	2773771	100.00%	6.288%
64	19.286	2769	2771	2779	rVB4	34345	48540	1.75%	0.110%
65	19.398	2786	2790	2798	rVB	31425	44230	1.59%	0.100%
66	19.539	2809	2814	2815	rBV4	41436	55255	1.99%	0.125%
67	19.674	2833	2837	2840	rBV3	41663	74738	2.69%	0.169%
68	19.716	2840	2844	2849	rVB2	83920	141277	5.09%	0.320%
69	19.763	2849	2852	2856	rBV	56324	62481	2.25%	0.142%
70	19.816	2857	2861	2866	rBV	42361	55167	1.99%	0.125%
71	19.874	2866	2871	2873	rBV2	66305	91958	3.32%	0.208%

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 Title : SVOA CALIBRATION

72	20.010	2892	2894	2897	rVB	44030	43842	1.58%	0.099%
73	20.057	2899	2902	2906	rBV2	39926	60536	2.18%	0.137%
74	20.239	2931	2933	2934	rBV2	22383	19178	0.69%	0.043%
75	20.263	2935	2937	2941	rVV	49801	62788	2.26%	0.142%
76	20.474	2970	2973	2977	rVB	52050	72747	2.62%	0.165%
77	20.633	2996	3000	3003	rBV2	57616	81207	2.93%	0.184%
78	20.674	3003	3007	3011	rVV2	115558	204016	7.36%	0.463%
79	20.727	3011	3016	3021	rVV3	496198	796394	28.71%	1.805%
80	20.780	3021	3025	3034	rVB2	357840	513884	18.53%	1.165%
81	20.927	3048	3050	3053	rVB	52314	44539	1.61%	0.101%
82	20.986	3053	3060	3064	rBV3	1250090	2005712	72.31%	4.547%
83	21.021	3064	3066	3070	rVB	295363	325260	11.73%	0.737%
84	21.110	3078	3081	3084	rBV	87591	111051	4.00%	0.252%
85	21.174	3088	3092	3095	rVB	128408	163983	5.91%	0.372%
86	21.592	3158	3163	3170	rBV3	354678	591029	21.31%	1.340%
87	22.015	3232	3235	3243	rVB	224888	295221	10.64%	0.669%
88	22.227	3268	3271	3278	rVB3	123364	166438	6.00%	0.377%
89	22.480	3308	3314	3318	rBV2	902569	1592111	57.40%	3.609%
90	22.515	3318	3320	3329	rVB2	610258	813175	29.32%	1.843%
91	22.904	3382	3386	3389	rBV	148953	234692	8.46%	0.532%
92	22.951	3389	3394	3398	rVV	1377681	2231786	80.46%	5.060%
93	22.992	3398	3401	3407	rVB	565925	812324	29.29%	1.842%
94	23.080	3410	3416	3422	rBV	766690	1267845	45.71%	2.874%
95	23.574	3494	3500	3506	rVB	845209	1437128	51.81%	3.258%
96	23.645	3507	3512	3523	rVB2	405016	748479	26.98%	1.697%
97	24.227	3605	3611	3620	rVB2	396620	912147	32.88%	2.068%
98	25.004	3737	3743	3754	rVB2	304441	790392	28.50%	1.792%
99	25.121	3758	3763	3772	rVB3	301077	729679	26.31%	1.654%
100	25.780	3871	3875	3887	rVB3	285676	742878	26.78%	1.684%

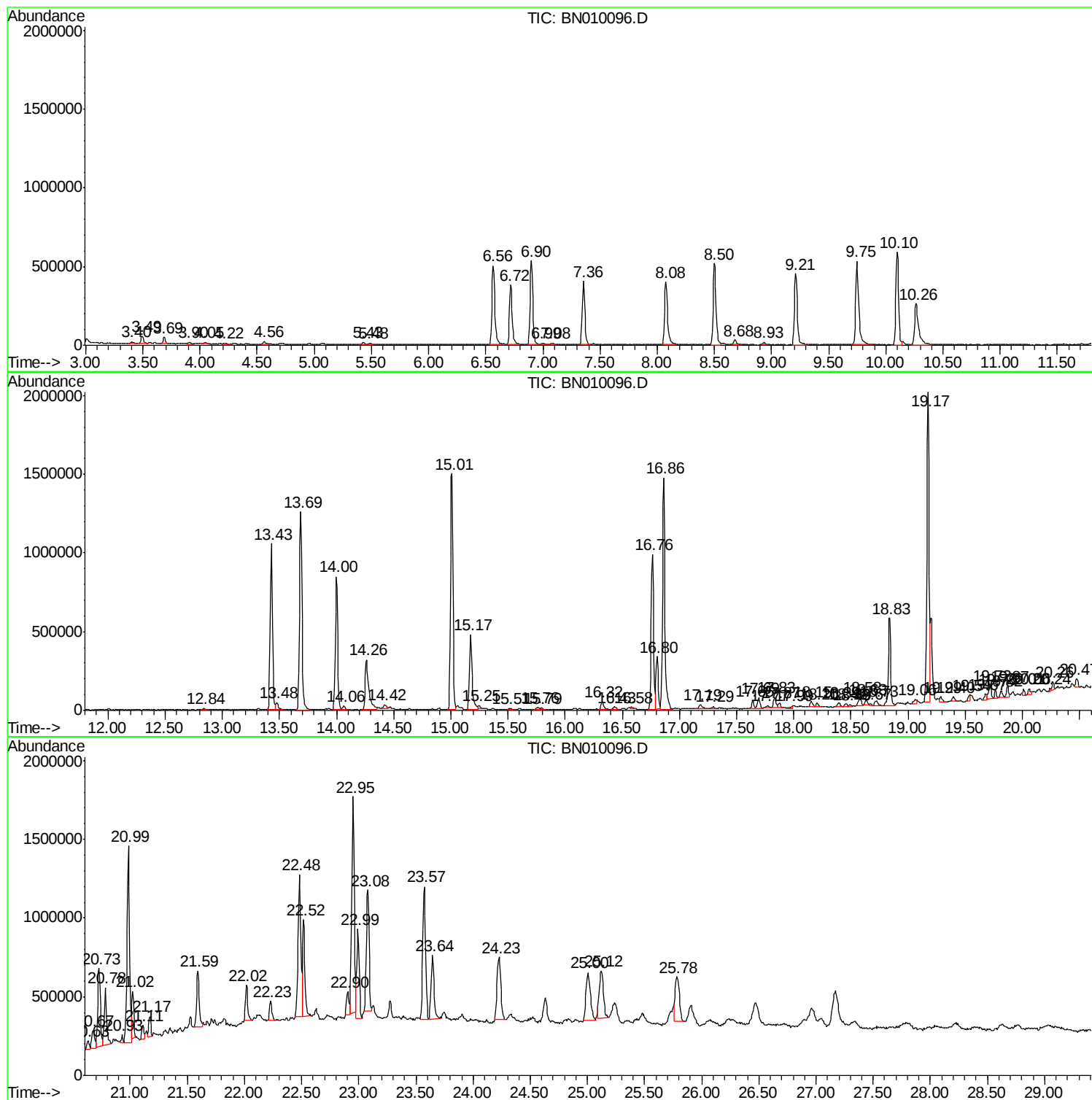
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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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



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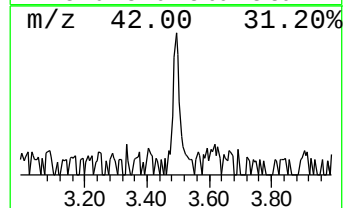
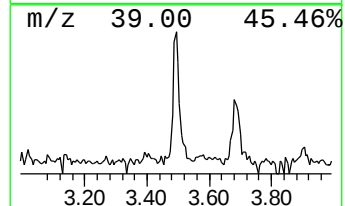
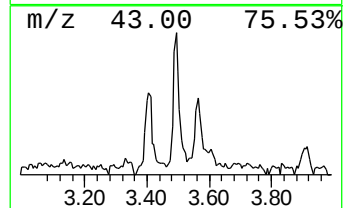
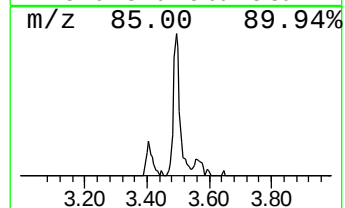
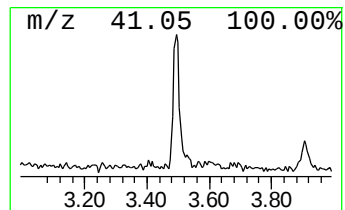
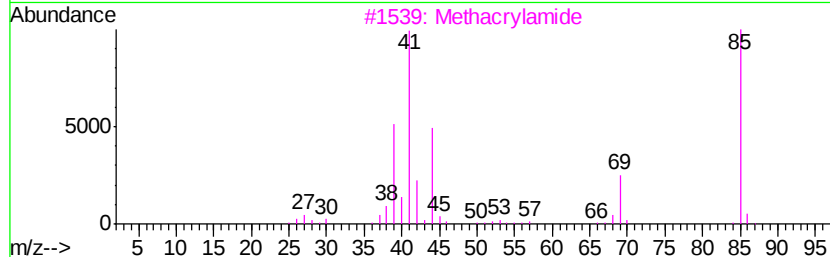
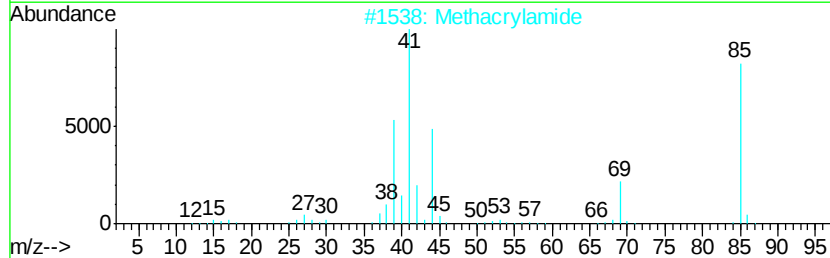
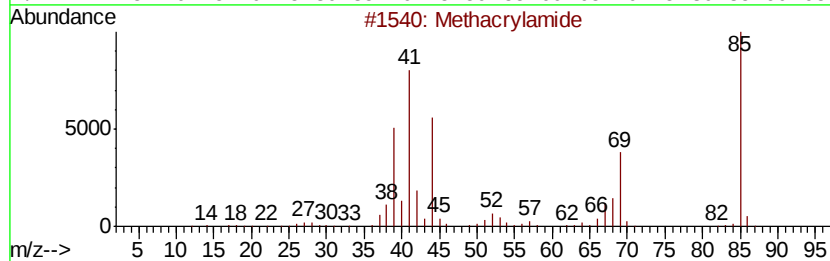
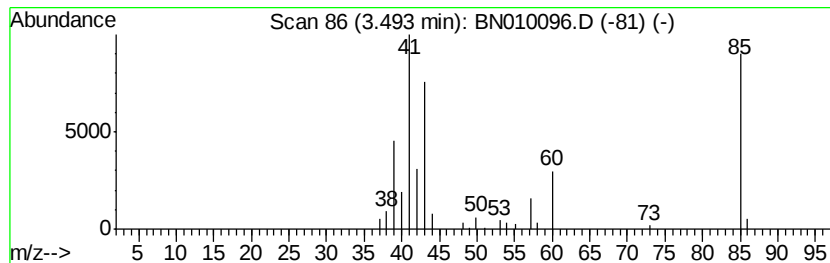
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Methacrylamide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.49	2.09 ng/ul	69631	1,4-Dichlorobenzene-d4	7.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	53
2			Methacrylamide	85	C4H7NO	000079-39-0	42
3			Methacrylamide	85	C4H7NO	000079-39-0	42
4			trans-Crotonamide	85	C4H7NO	000625-37-6	33
5			2-Pyrrolidinone	85	C4H7NO	000616-45-5	32



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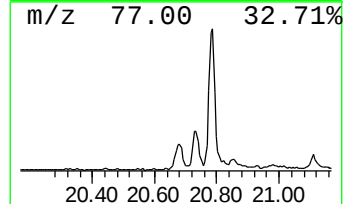
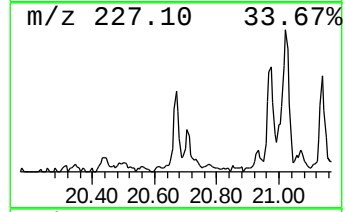
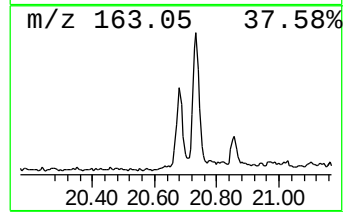
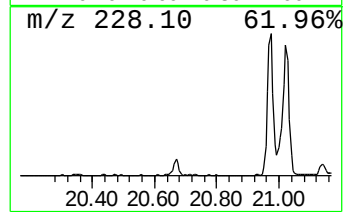
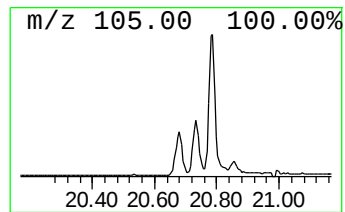
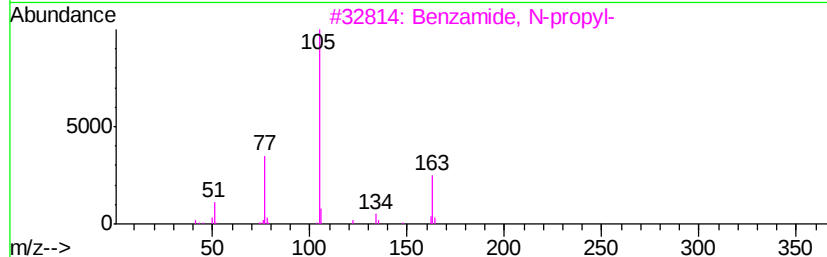
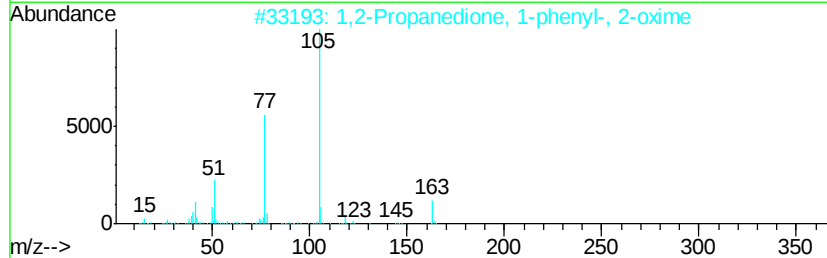
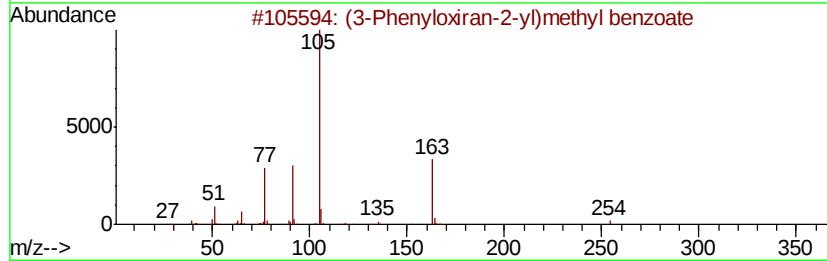
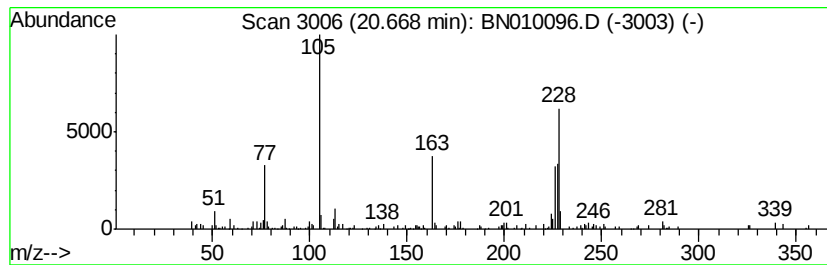
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.67	2.03 ng/ul	204016	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	47
2		1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	47
3		Benzamide, N-propyl-	163	C10H13NO	010546-70-0	47
4		Benzenepropanamide, .alpha.-(ben...	284	C16H16N2O3	058690-81-6	47
5		Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	47



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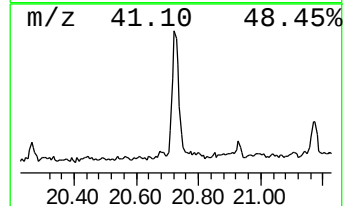
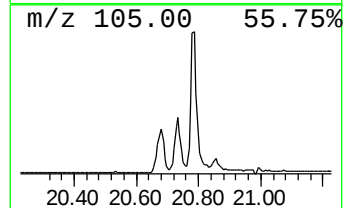
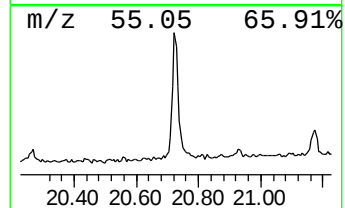
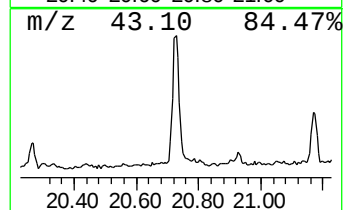
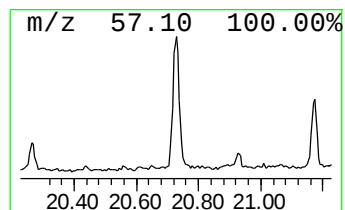
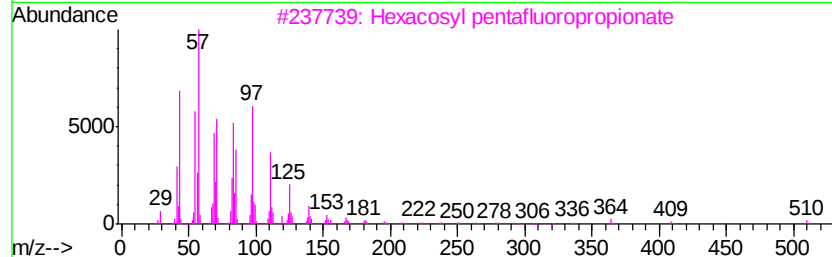
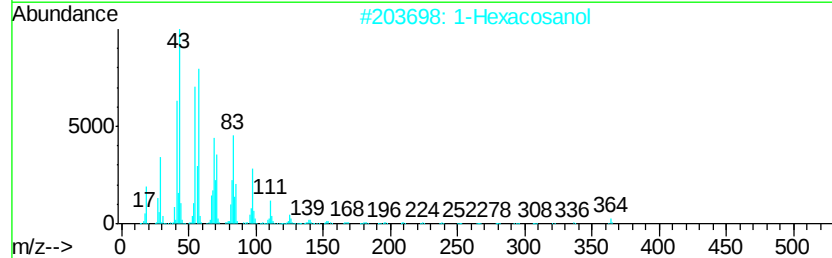
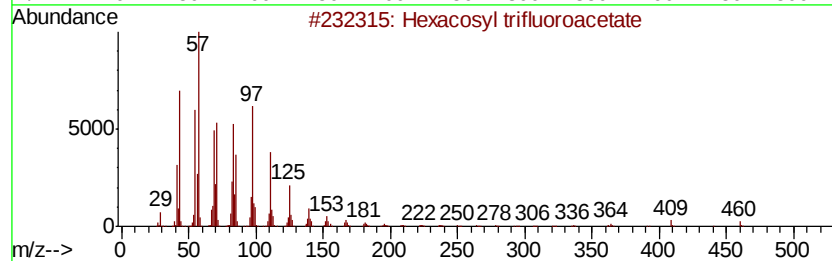
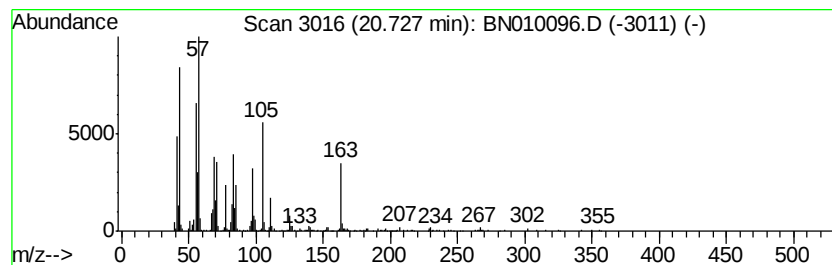
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Hexacosyl trifluoroacetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.73	7.94 ng/ul	796394	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	74
2		1-Hexacosanol	382	C26H54O	000506-52-5	72
3		Hexacosyl pentafluoropropionate	528	C29H53F5O2	1000351-81-2	72
4		Triaccontyl heptafluorobutyrte	634	C34H61F7O2	1000351-83-2	64
5		Oxalic acid, hexadecyl propyl ester	356	C21H40O4	1000309-26-9	62



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Data File : BN010096.D
Acq On : 05 Mar 2020 01:51
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Sample : L1641-02
Misc :
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ClientSampleId :
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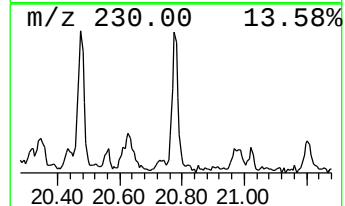
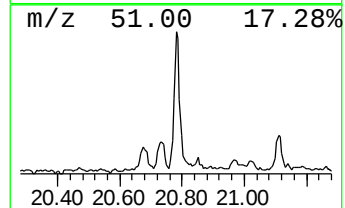
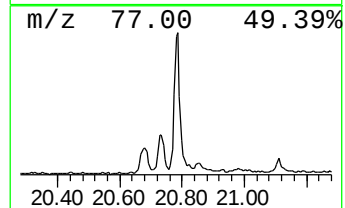
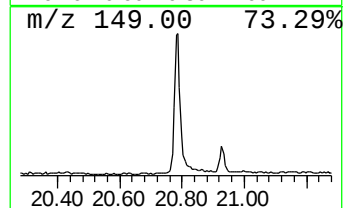
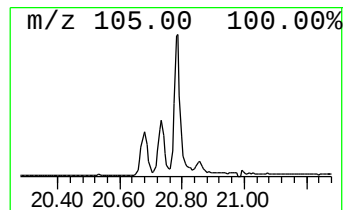
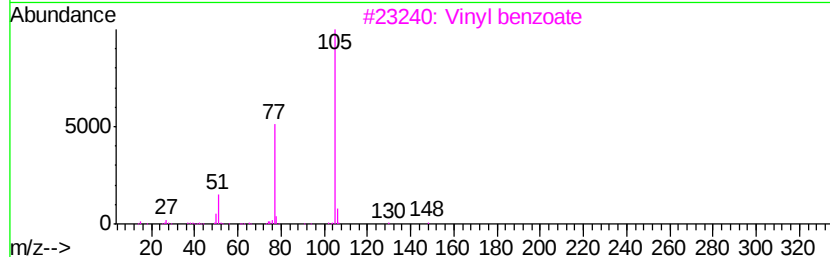
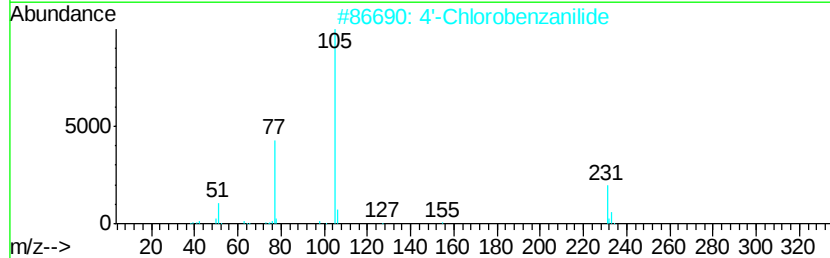
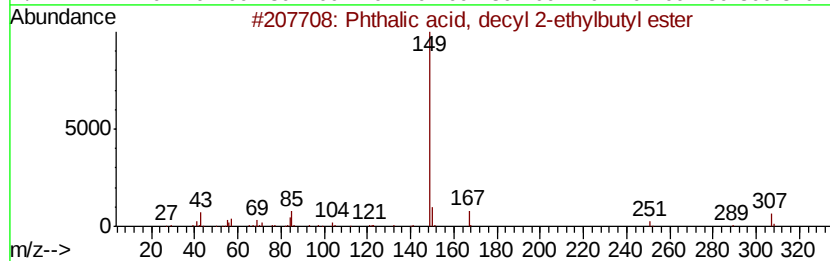
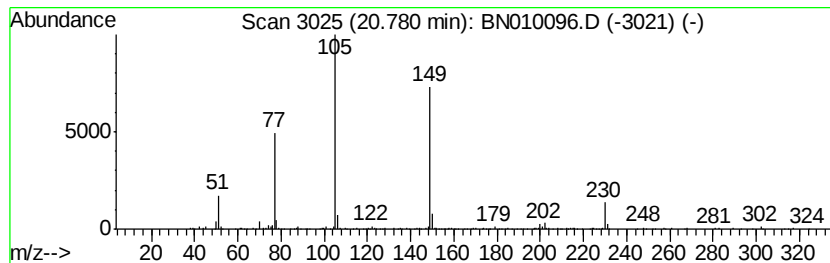
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.78	5.12 ng/ul	513884	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, decyl 2-ethylbutyl...	390	C24H38O4	1000356-89-5	35
2		4'-Chlorobenzanilide	231	C13H10ClNO	002866-82-2	30
3		Vinyl benzoate	148	C9H8O2	000769-78-8	30
4		Cyclobutyl phenyl ketone	160	C11H12O	005407-98-7	27
5		Benzenepentanoic acid, .delta.-oxo-	192	C11H12O3	001501-05-9	27



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010096.D
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Operator : CG/JU
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ClientSampleId :
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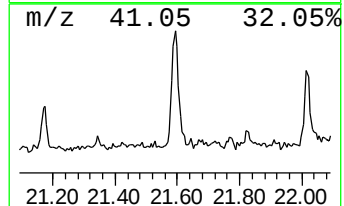
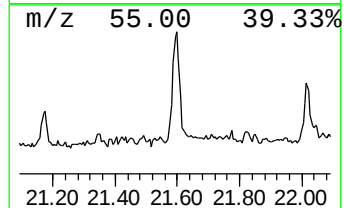
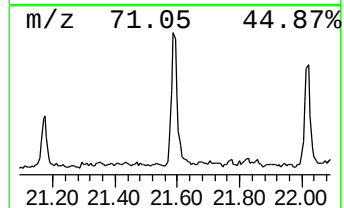
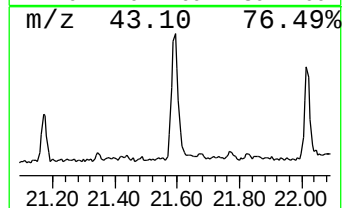
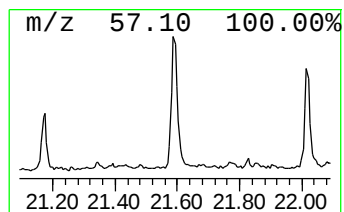
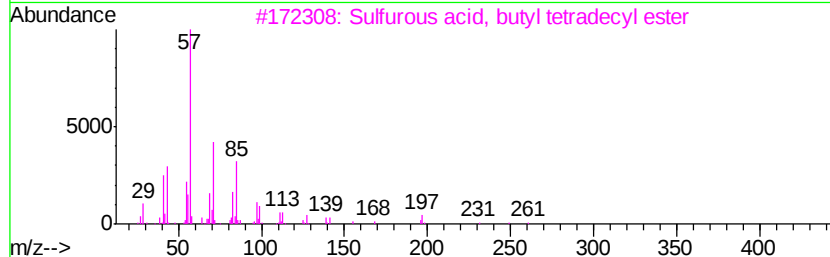
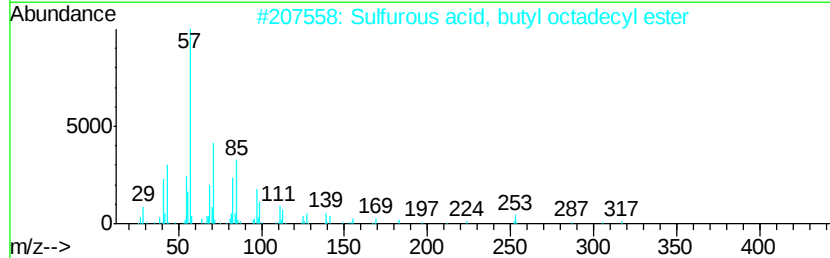
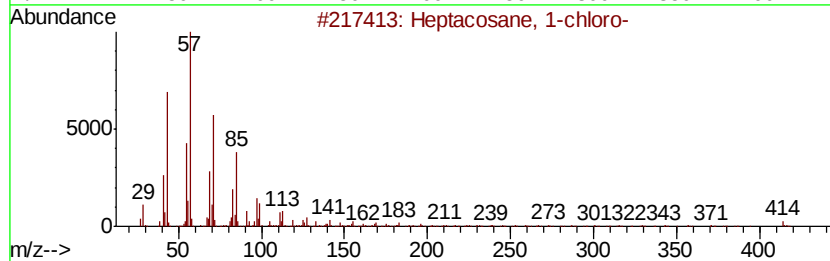
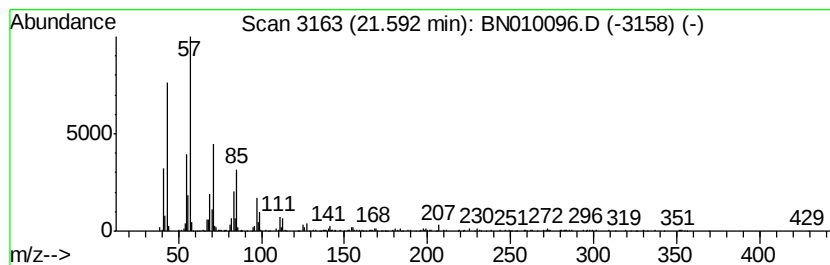
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Heptacosane, 1-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.59	5.89 ng/ul	591029	Chrysene-d12	20.99

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane, 1-chloro-		414	C27H55Cl	062016-79-9	91
2	Sulfurous acid, butyl octadecyl ...		390	C22H46O3S	1000309-18-5	91
3	Sulfurous acid, butyl tetradecyl...		334	C18H38O3S	1000309-18-1	87
4	Sulfurous acid, butyl dodecyl ester		306	C16H34O3S	1000309-17-9	83
5	Sulfurous acid, butyl tridecyl e...		320	C17H36O3S	1000309-18-0	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010096.D
Acq On : 05 Mar 2020 01:51
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Sample : L1641-02
Misc :
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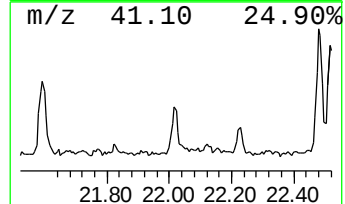
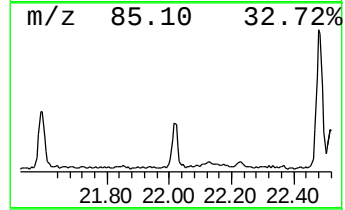
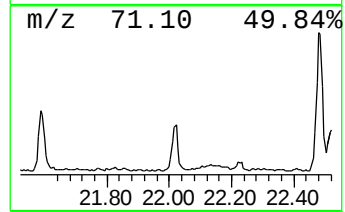
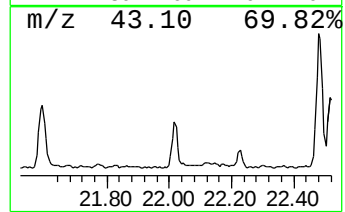
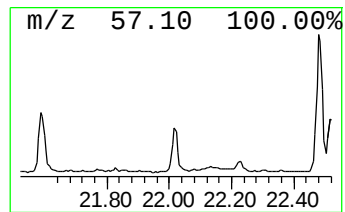
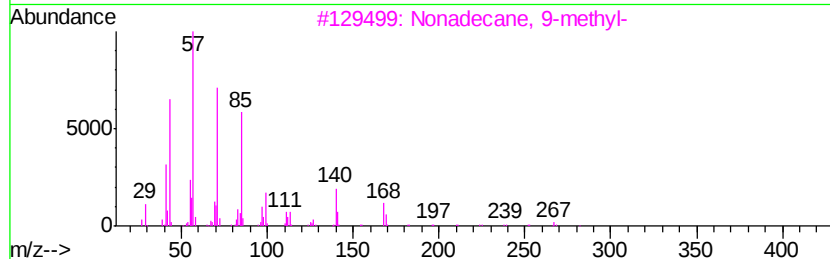
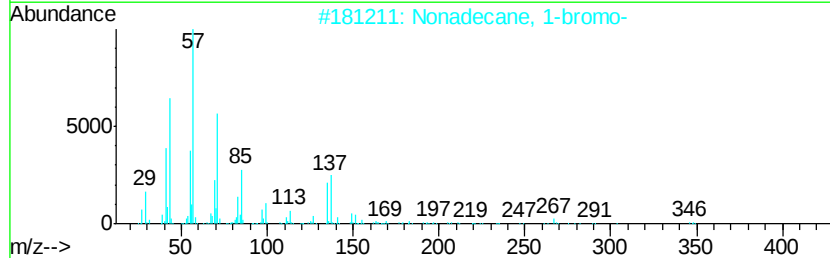
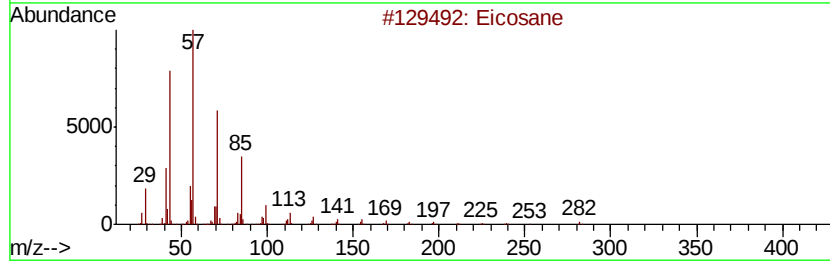
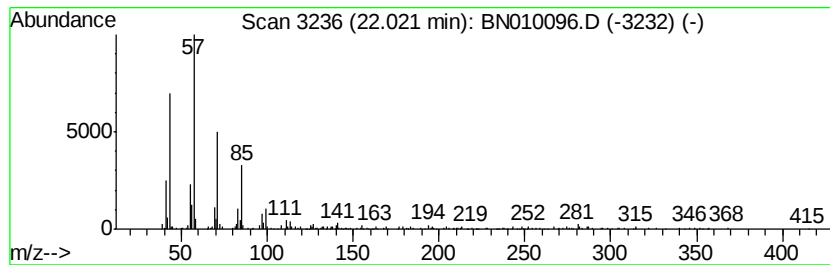
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.02	2.94 ng/ul	295221	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	92
2		Nonadecane, 1-bromo-	346	C19H39Br	004434-66-6	91
3		Nonadecane, 9-methyl-	282	C20H42	013287-24-6	87
4		Heptacosane	380	C27H56	000593-49-7	87
5		Eicosane, 3-methyl-	296	C21H44	006418-46-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010096.D
Acq On : 05 Mar 2020 01:51
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Sample : L1641-02
Misc :
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ClientSampleId :
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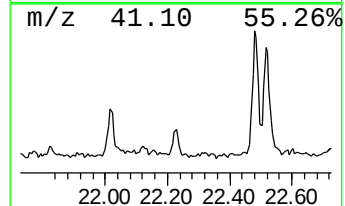
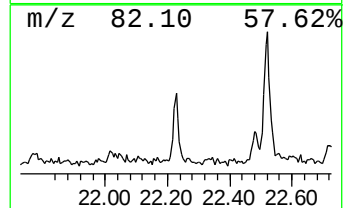
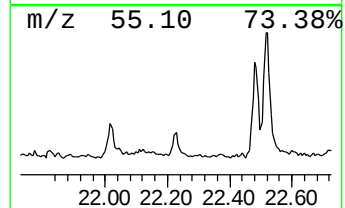
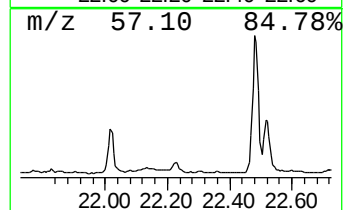
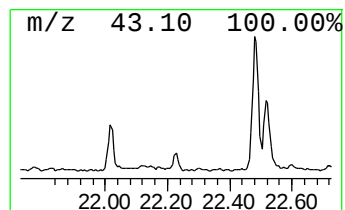
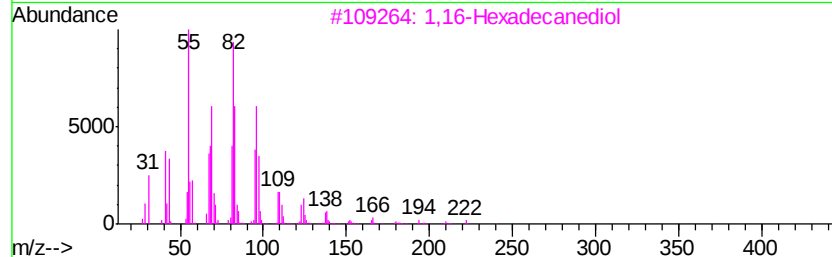
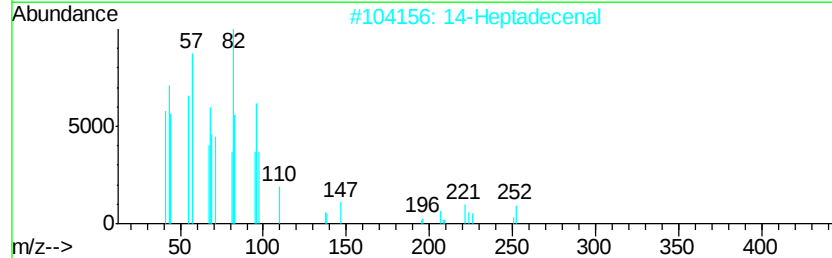
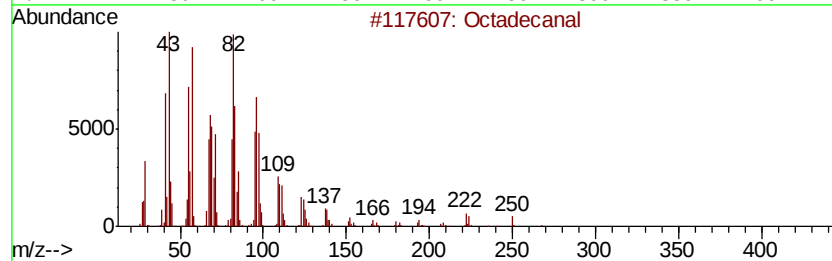
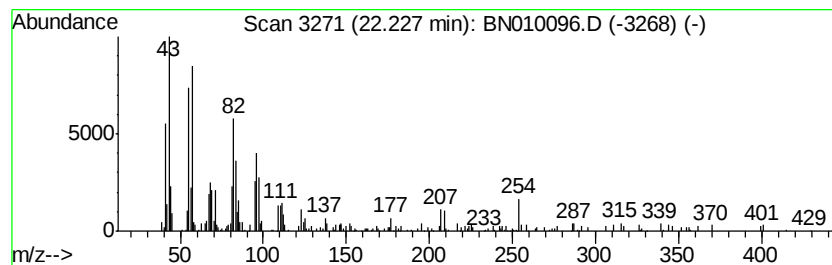
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.23	2.63 ng/ul	166438	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanal	268	C18H36O	000638-66-4	74
2		14-Heptadecenal	252	C17H32O	1000144-58-0	70
3		1,16-Hexadecanediol	258	C16H34O2	007735-42-4	62
4		E-3-Pentadecen-2-ol	226	C15H30O	1000130-83-8	58
5		1,15-Pentadecanediol	244	C15H32O2	014722-40-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010096.D
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Misc :
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ClientSampleId :
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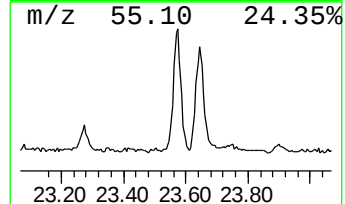
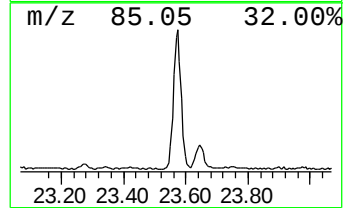
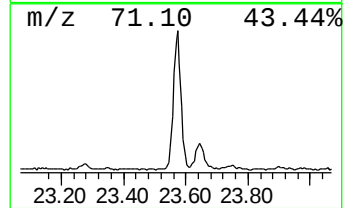
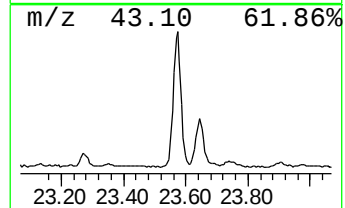
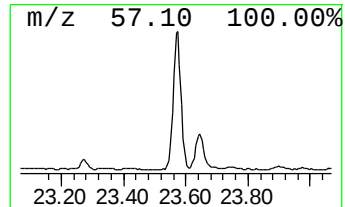
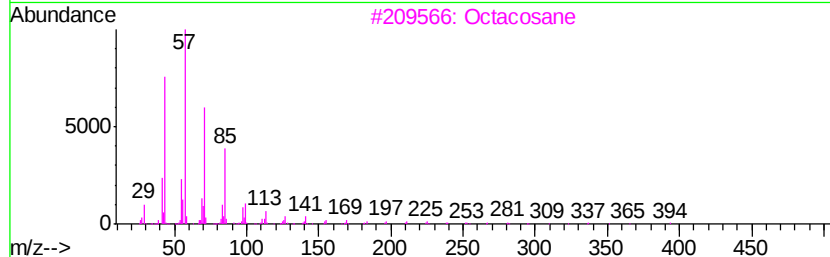
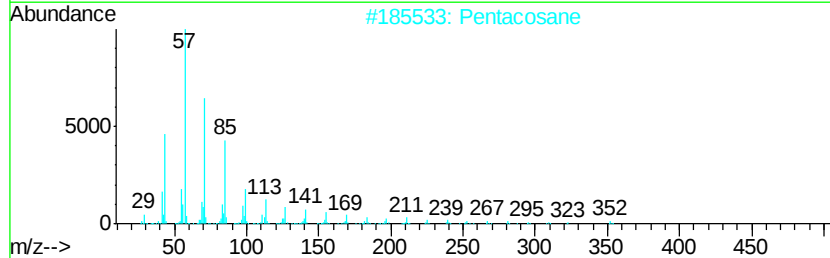
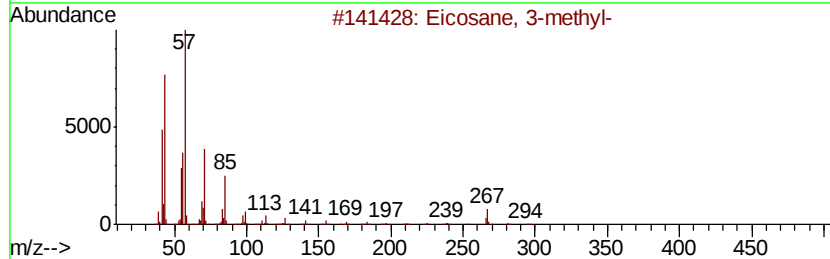
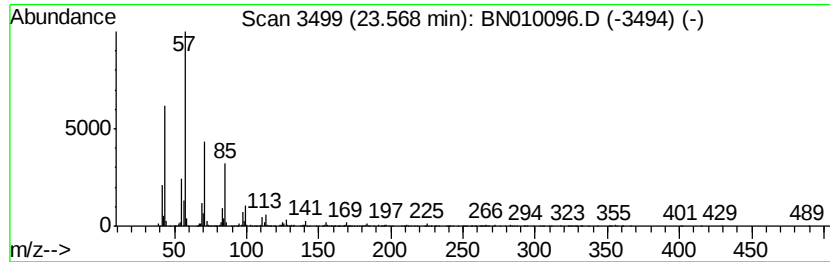
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.57	22.67 ng/ul	1437130	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 3-methyl-	296	C21H44	006418-46-8	91
2		Pentacosane	352	C25H52	000629-99-2	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
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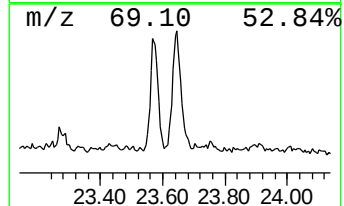
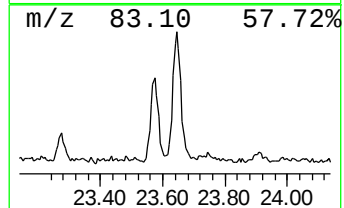
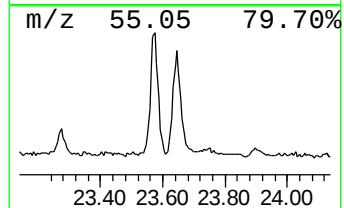
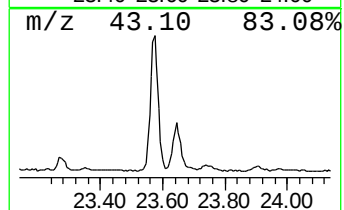
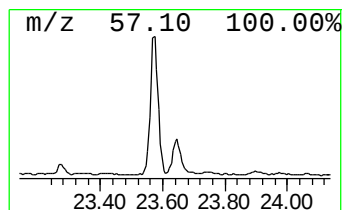
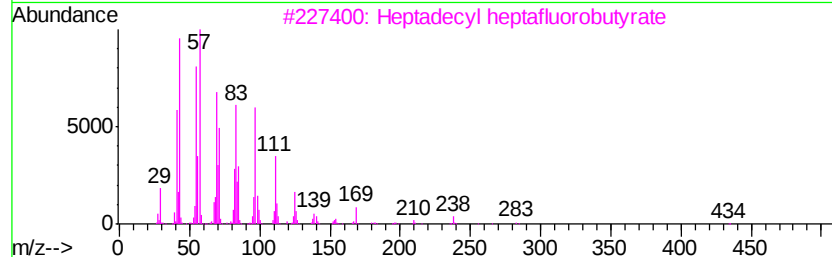
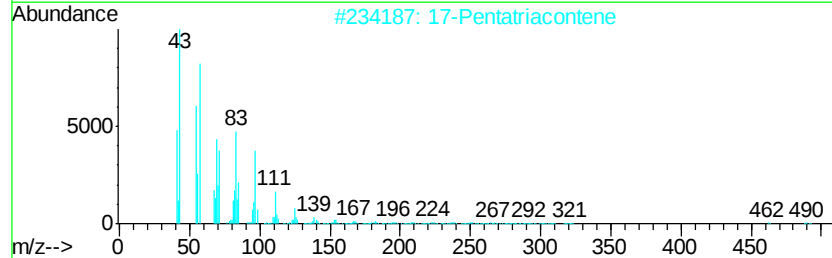
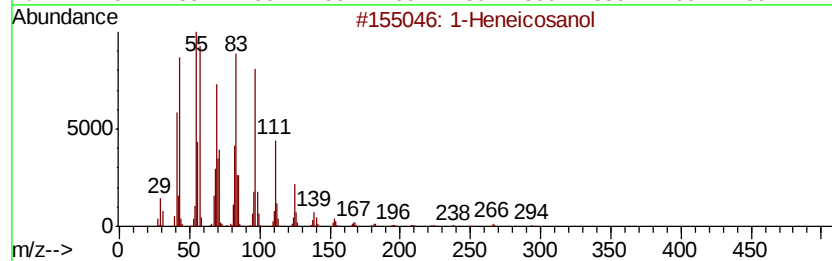
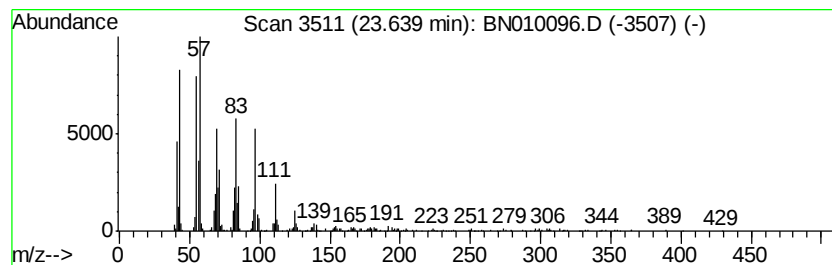
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.64	11.81 ng/ul	748479	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosanol	312	C21H44O	015594-90-8	91
2		17-Pentatriacontene	491	C35H70	006971-40-0	91
3		Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	91
4		Octacosanol	410	C28H58O	000557-61-9	90
5		1-Hexacosanol	382	C26H54O	000506-52-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010096.D
Acq On : 05 Mar 2020 01:51
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ClientSampleId :
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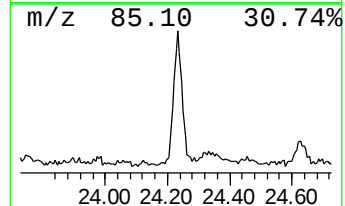
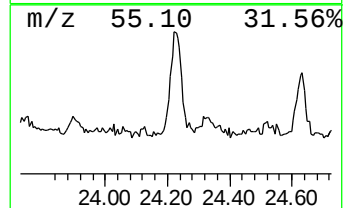
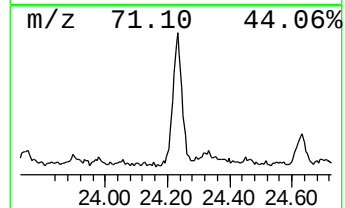
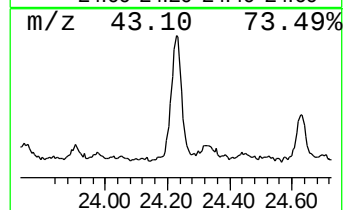
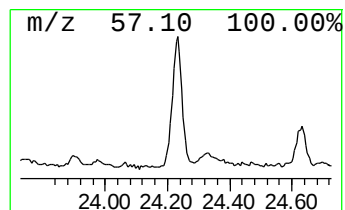
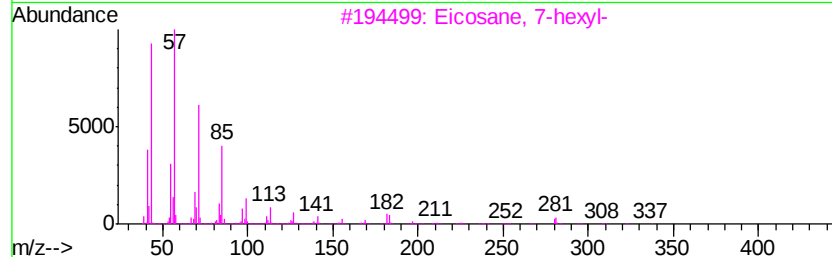
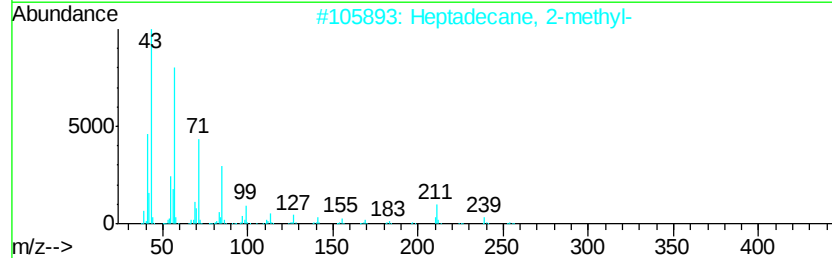
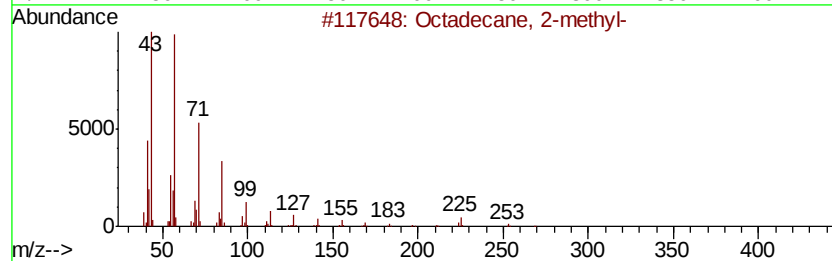
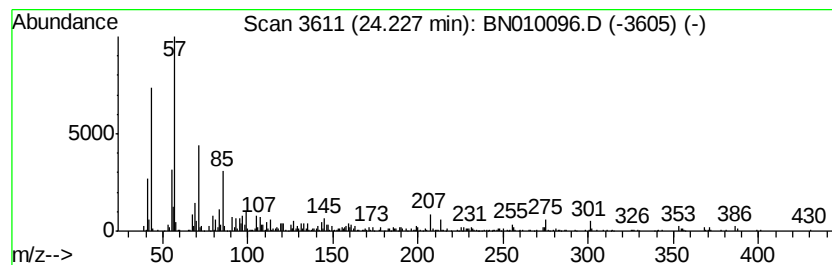
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.23	14.39 ng/ul	912147	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	86
2		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	86
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	81
4		Hexadecane	226	C16H34	000544-76-3	76
5		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030420\
Data File : BN010096.D
Acq On : 05 Mar 2020 01:51
Operator : CG/JU
Sample : L1641-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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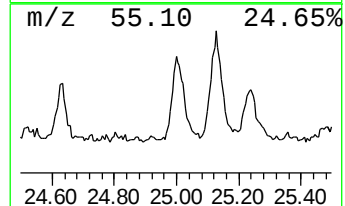
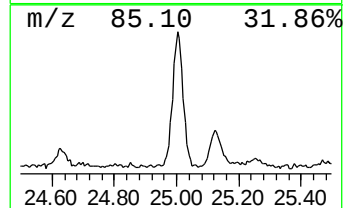
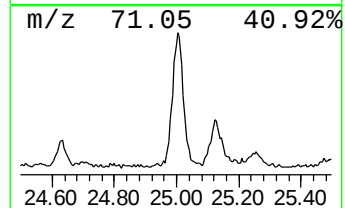
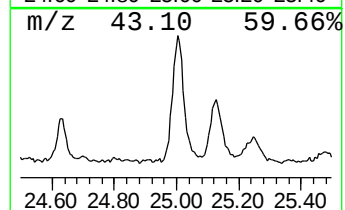
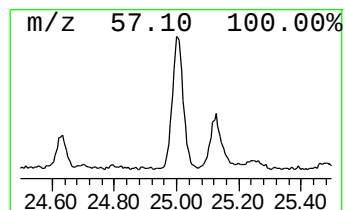
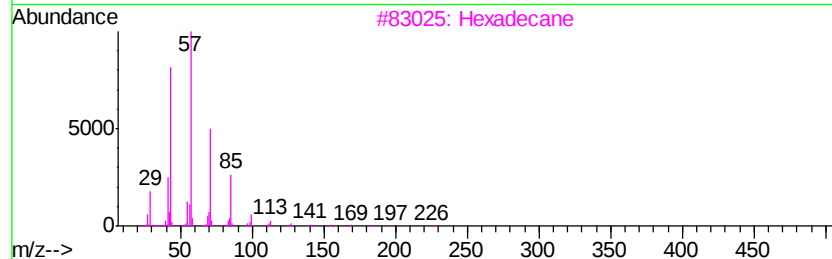
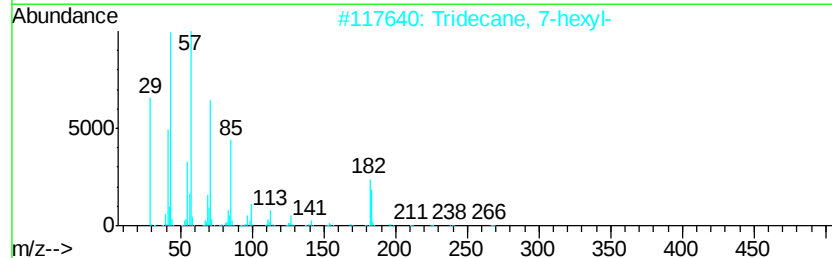
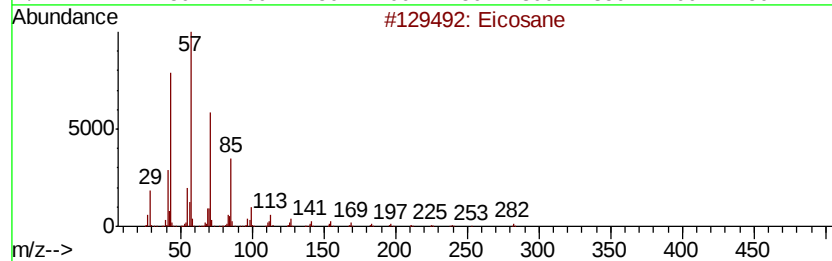
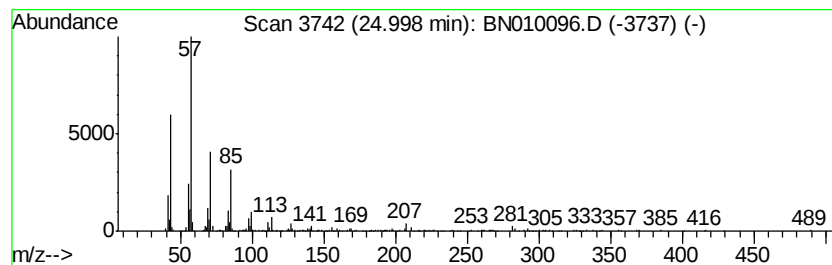
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.00	12.47 ng/ul	790392	Perylene-d12	23.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	94
3		Hexadecane	226	C16H34	000544-76-3	93
4		Hentriacontane	437	C31H64	000630-04-6	83
5		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030420\
 Data File : BN010096.D
 Acq On : 05 Mar 2020 01:51
 Operator : CG/JU
 Sample : L1641-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB6Y5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.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
Methacrylamide	3.49	2.1	ng/ul	69631	1	7.36	667624	20.0
unknown-01	20.67	2.0	ng/ul	204016	5	20.99	2005710	20.0
Hexacosyl trifluo...	20.73	7.9	ng/ul	796394	5	20.99	2005710	20.0
unknown-02	20.78	5.1	ng/ul	513884	5	20.99	2005710	20.0
Heptacosane, 1-ch...	21.59	5.9	ng/ul	591029	5	20.99	2005710	20.0
(DEL) Alkane: Str...	22.02	2.9	ng/ul	295221	5	20.99	2005710	20.0
Octadecanal	22.23	2.6	ng/ul	166438	6	23.08	1267850	20.0
(DEL) Alkane: Str...	23.57	22.7	ng/ul	1437130	6	23.08	1267850	20.0
1-Heneicosanol	23.64	11.8	ng/ul	748479	6	23.08	1267850	20.0
(DEL) Alkane: Str...	24.23	14.4	ng/ul	912147	6	23.08	1267850	20.0
(DEL) Alkane: Str...	25.00	12.5	ng/ul	790392	6	23.08	1267850	20.0