

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM090420\
 Data File : BM027344.D
 Acq On : 04 Sep 2020 23:35
 Operator : JU/CG
 Sample : L3843-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC1

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_M\METHODS\SOM-EPA-BM090420MA.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	3.134	21	25	31	rVB	32001	40745	0.82%	0.090%
2	3.404	69	71	72	rBV2	8028	5517	0.11%	0.012%
3	6.757	634	641	650	rBV	539300	864584	17.32%	1.907%
4	6.916	662	668	678	rVB	447894	686276	13.75%	1.514%
5	7.110	694	701	710	rBV	600581	929763	18.63%	2.051%
6	7.186	710	714	728	rVV	84278	153900	3.08%	0.339%
7	7.569	773	779	789	rVV	509127	782204	15.67%	1.725%
8	7.651	789	793	801	rVB2	24629	39483	0.79%	0.087%
9	8.281	893	900	916	rBV	696728	1118647	22.42%	2.468%
10	8.716	967	974	982	rBV	562495	924419	18.52%	2.039%
11	8.875	995	1001	1013	rBV	43565	78726	1.58%	0.174%
12	9.433	1089	1096	1108	rBV	485969	813601	16.30%	1.795%
13	9.851	1160	1167	1174	rVB5	4271	7967	0.16%	0.018%
14	9.969	1179	1187	1197	rBV	796114	1286991	25.79%	2.839%
15	10.339	1239	1250	1263	rBV	679959	1110284	22.25%	2.449%
16	10.475	1266	1273	1290	rBV	706023	1202060	24.09%	2.652%
17	11.933	1515	1521	1531	rBV3	9358	19238	0.39%	0.042%
18	13.127	1719	1724	1732	rBV2	104149	161408	3.23%	0.356%
19	13.627	1803	1809	1814	rBV	1486361	2133180	42.74%	4.706%
20	13.674	1814	1817	1826	rVB	585497	780266	15.63%	1.721%
21	13.898	1848	1855	1868	rBV	1624774	2365143	47.39%	5.217%
22	14.210	1902	1908	1917	rBV	1172434	1735118	34.77%	3.828%
23	14.421	1938	1944	1964	rBV	507327	814272	16.32%	1.796%
24	14.657	1981	1984	1993	rVB6	4142	7710	0.15%	0.017%
25	15.210	2071	2078	2087	rBV	2245141	3023776	60.59%	6.670%
26	15.333	2093	2099	2110	rBV	563203	764244	15.31%	1.686%
27	15.451	2115	2119	2124	rVB2	25853	35650	0.71%	0.079%
28	15.780	2171	2175	2179	rBV4	4800	8039	0.16%	0.018%
29	15.998	2208	2212	2219	rVB7	8906	15228	0.31%	0.034%
30	16.745	2337	2339	2345	rVB3	6986	9585	0.19%	0.021%
31	16.956	2369	2375	2386	rBV	1672458	2260181	45.29%	4.986%
32	17.056	2386	2392	2401	rBV	2537758	3436862	68.87%	7.581%
33	17.486	2463	2465	2471	rBV7	3440	5626	0.11%	0.012%
34	17.886	2526	2533	2541	rBV	218053	342864	6.87%	0.756%

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35	17.945	2541	2543	2549	rVB2	16246	24421	0.49%	0.054%
36	18.750	2678	2680	2686	rVB7	8007	12285	0.25%	0.027%
37	18.821	2689	2692	2695	rVV4	11917	14415	0.29%	0.032%
38	18.992	2717	2721	2724	rBV	27289	44482	0.89%	0.098%
39	19.021	2724	2726	2733	rVB4	19422	31477	0.63%	0.069%
40	19.168	2747	2751	2759	rBV2	70678	108368	2.17%	0.239%
41	19.245	2760	2764	2771	rVB	32790	53579	1.07%	0.118%
42	19.356	2777	2783	2791	rBV	3426253	4477005	89.71%	9.876%
43	19.439	2795	2797	2803	rVB7	14822	19126	0.38%	0.042%
44	19.944	2880	2883	2886	rBV3	23816	25294	0.51%	0.056%
45	20.397	2956	2960	2964	rBV	142108	159533	3.20%	0.352%
46	20.439	2965	2967	2971	rVB2	44889	47552	0.95%	0.105%
47	20.897	3041	3045	3055	rBV	411438	565743	11.34%	1.248%
48	21.156	3084	3089	3100	rBV	2423634	3120521	62.53%	6.884%
49	21.344	3118	3121	3127	rVB2	92217	104040	2.08%	0.230%
50	21.768	3190	3193	3200	rBV4	111122	172030	3.45%	0.379%
51	22.221	3267	3270	3274	rVB4	96964	110548	2.22%	0.244%
52	23.191	3428	3435	3442	rBV	2843568	4990540	100.00%	11.009%
53	23.327	3452	3458	3472	rVB	1820871	3287921	65.88%	7.253%

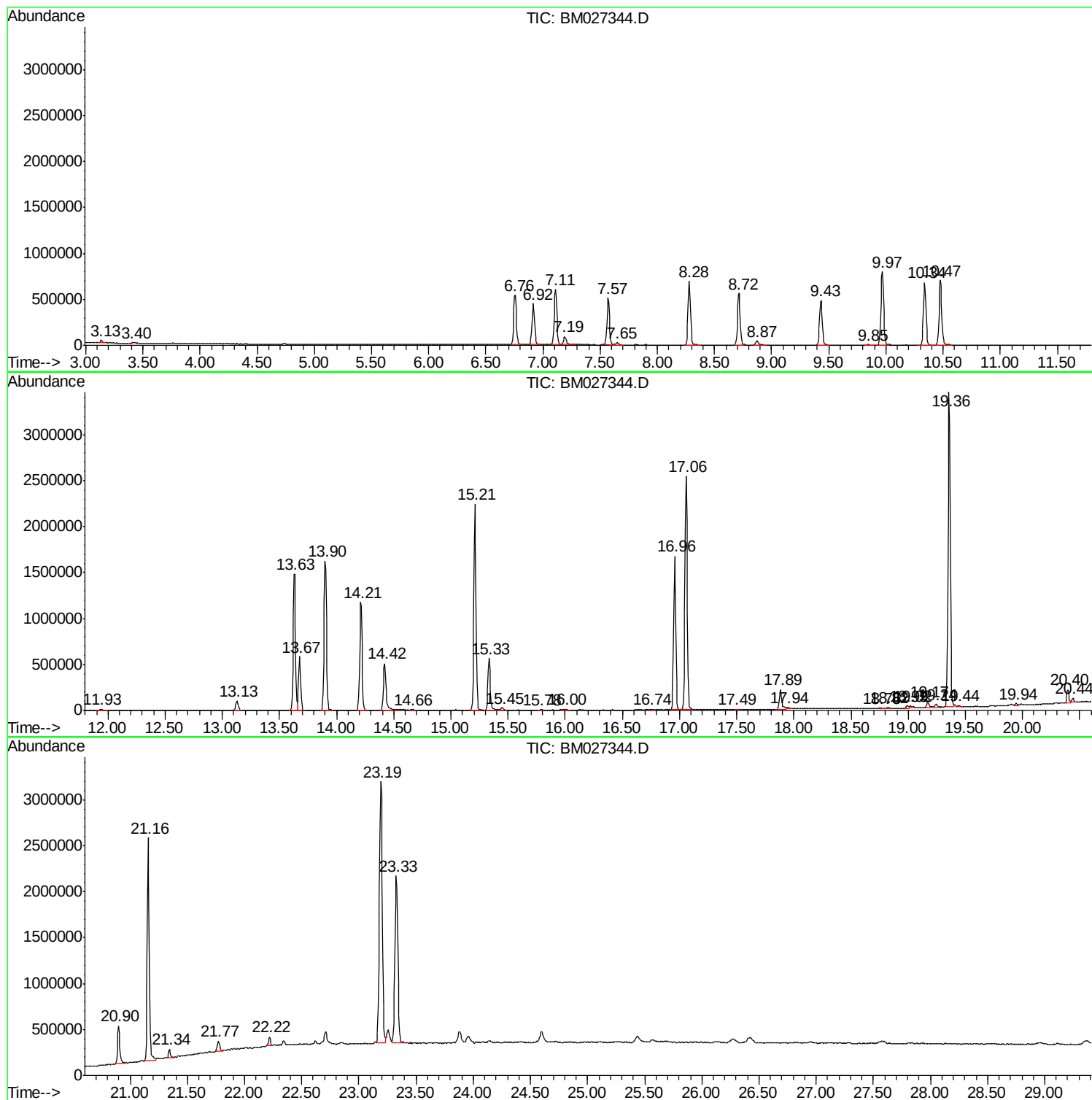
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090420MA.M
Quant Title : SVOA CALIBRATION

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



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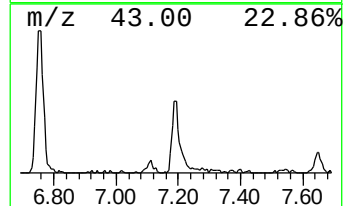
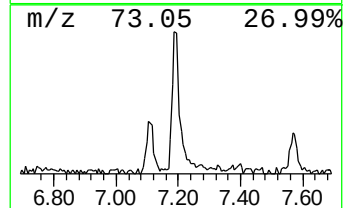
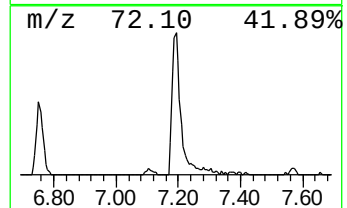
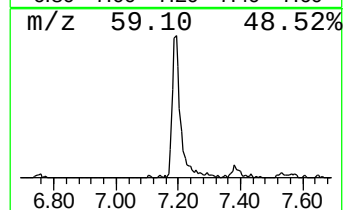
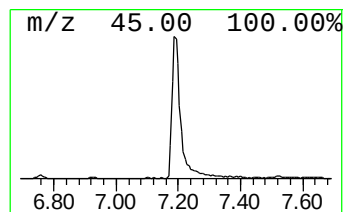
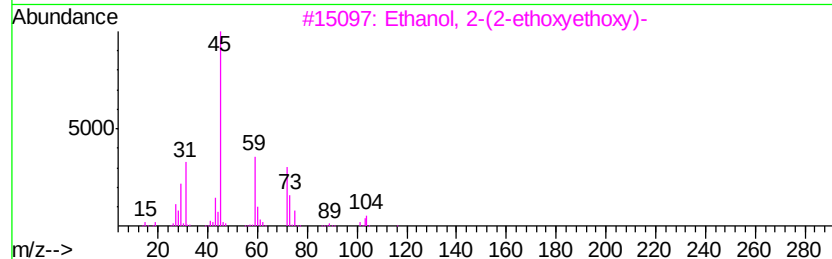
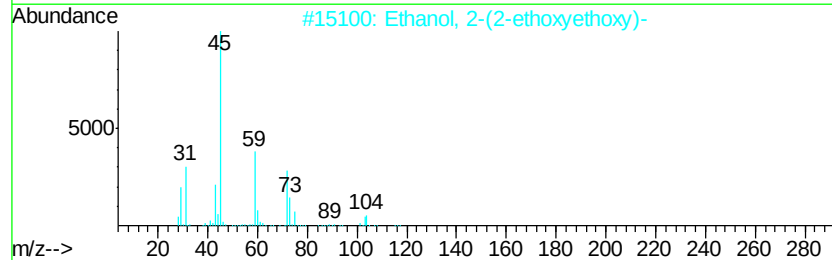
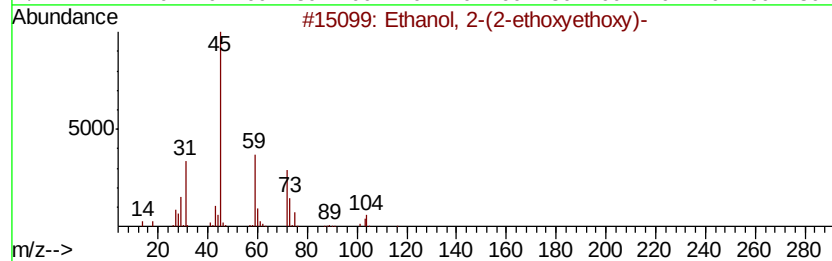
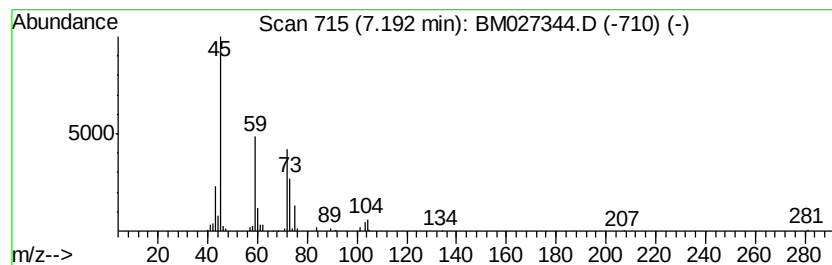
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090420MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	3.94 ng/ul	153900	1,4-Dichlorobenzene-d4	7.57

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-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
4			Diethyl carbitol	162	C8H18O3	000112-36-7	72
5			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72



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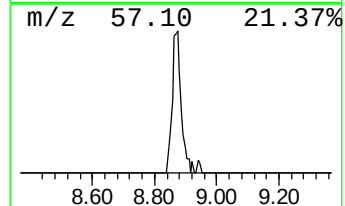
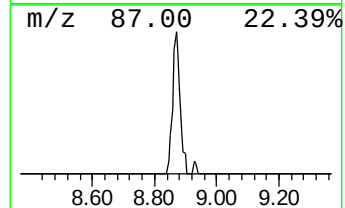
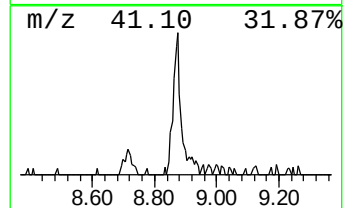
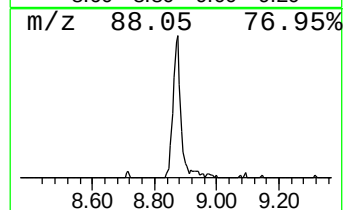
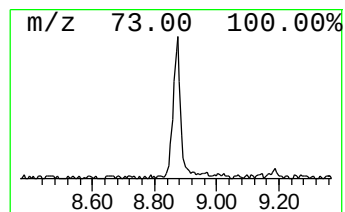
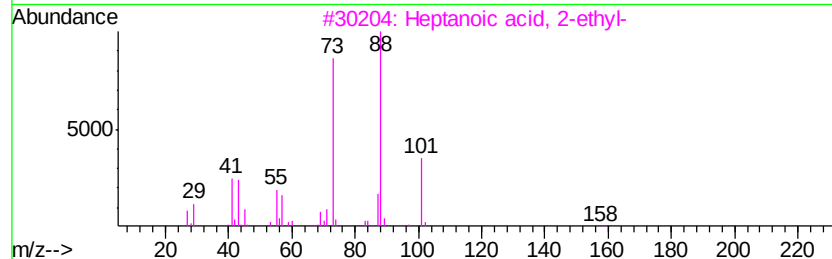
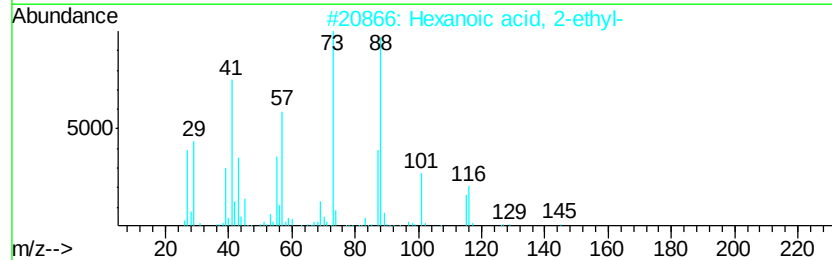
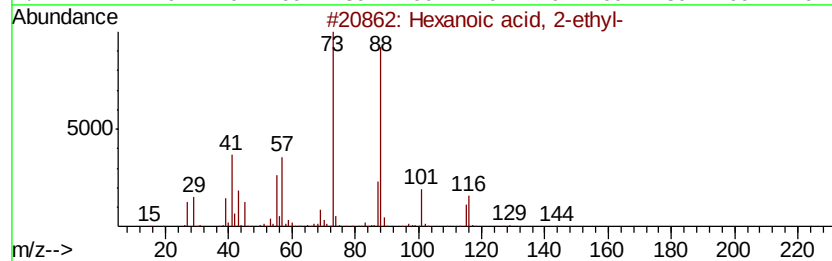
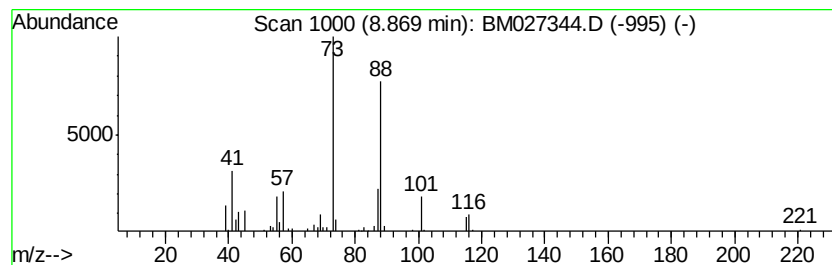
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexanoic acid, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.87	2.01 ng/ul	78726	1,4-Dichlorobenzene-d4	7.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	64
2		Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	59
3		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	40
4		Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	40
5		Methyl 2,3,4-tri-O-methyl-6-deox...	220	C10H20O5	072983-16-5	38



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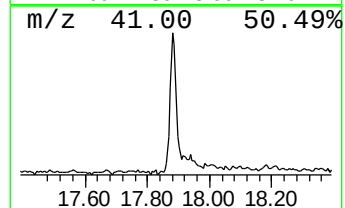
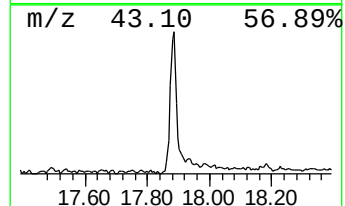
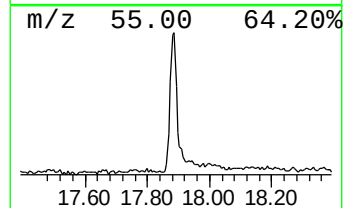
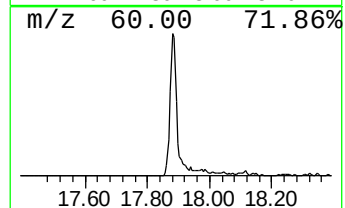
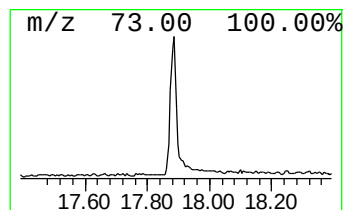
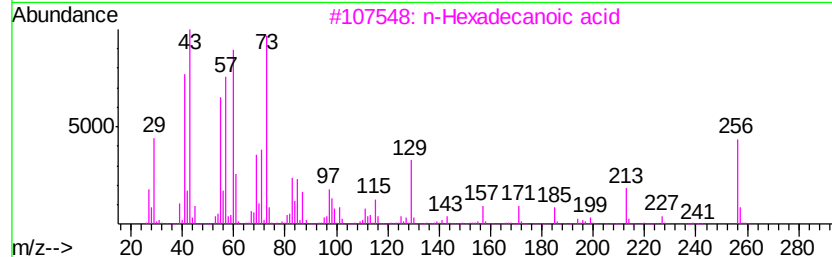
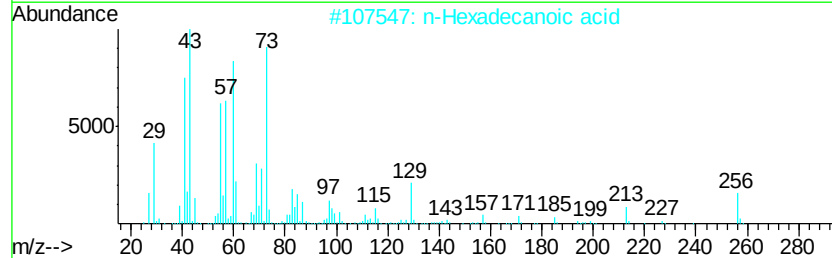
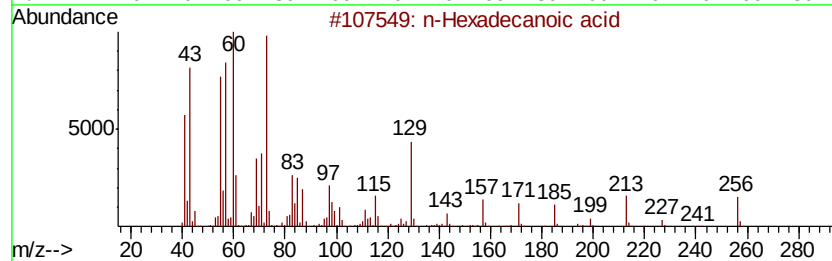
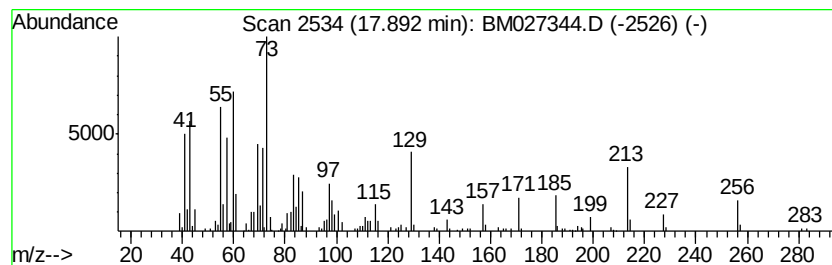
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.89	3.03 ng/ul	342864	Phenanthrene-d10	16.96
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	n-Hexadecanoic acid	256 C16H32O2	000057-10-3	93
2	n-Hexadecanoic acid	256 C16H32O2	000057-10-3	93
3	n-Hexadecanoic acid	256 C16H32O2	000057-10-3	90
4	Tridecanoic acid	214 C13H26O2	000638-53-9	64
5	Pentadecanoic acid	242 C15H30O2	001002-84-2	53



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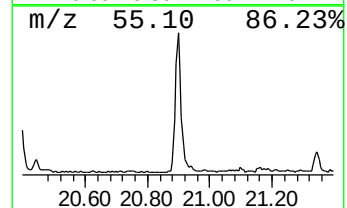
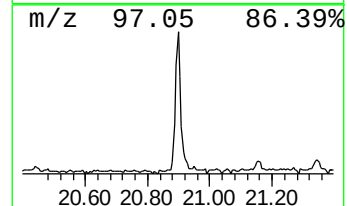
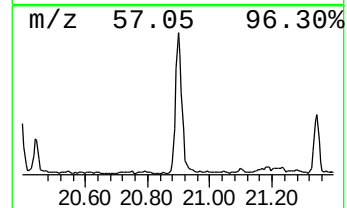
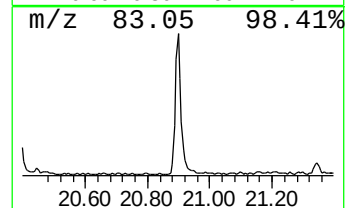
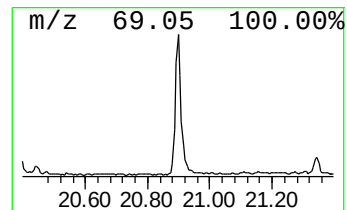
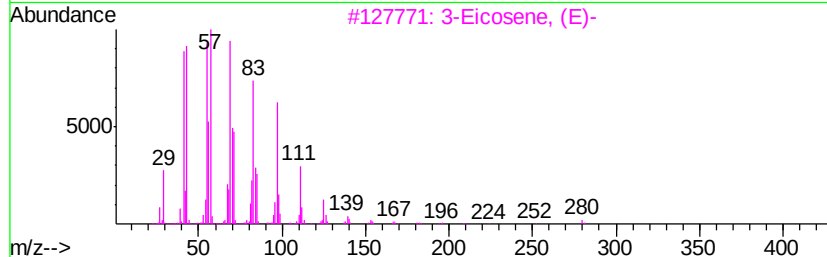
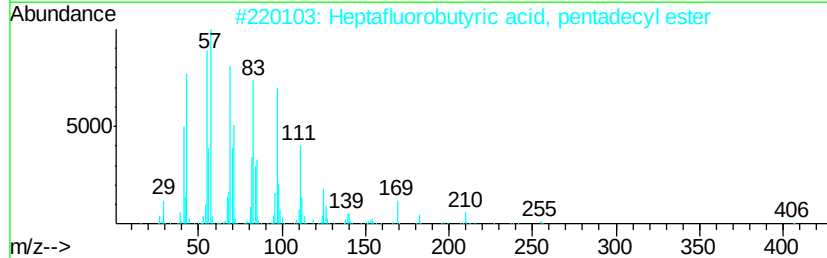
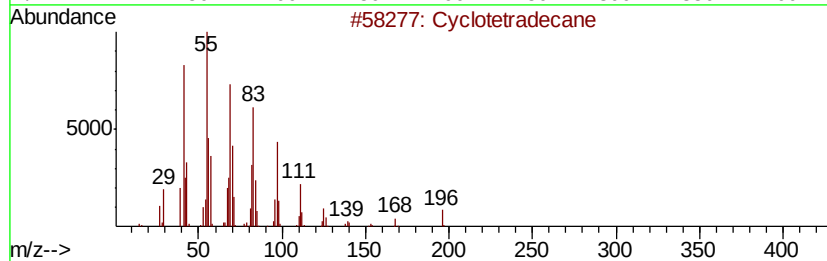
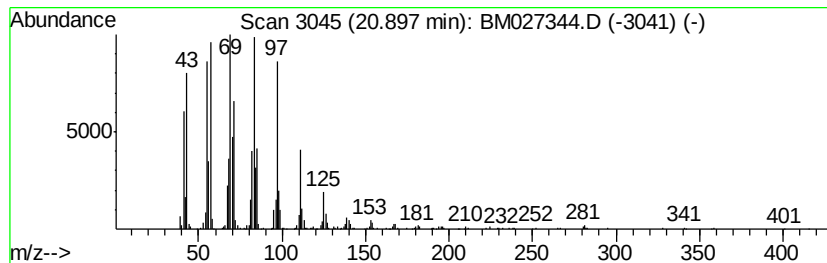
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 (DEL) Alkane: Cyclic20.90 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.90	3.63 ng/ul	565743	Chrysene-d12	21.16
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclotetradecane	196 C14H28	000295-17-0 94
2		Heptafluorobutyric acid, pentade...	424 C19H31F7O2	959261-23-5 94
3		3-Eicosene, (E)-	280 C20H40	074685-33-9 90
4		Heptafluorobutyric acid, n-tetra...	410 C18H29F7O2	007365-36-8 87
5		1-Heptacosanol	396 C27H56O	002004-39-9 87



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-eth...	7.19	3.9	ng/ul	153900	1	7.57	782204	20.0
Hexanoic acid, 2-...	8.87	2.0	ng/ul	78726	1	7.57	782204	20.0
n-Hexadecanoic acid	17.89	3.0	ng/ul	342864	4	16.96	2260180	20.0
(DEL) Alkane: Cyc...	20.90	3.6	ng/ul	565743	5	21.16	3120520	20.0