

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
 Data File : BM026086.D
 Acq On : 22 May 2020 19:08
 Operator : MA/SJ
 Sample : L2737-18
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9B40

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-BM051320MA.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	3.069	28	31	38	rVB	46872	58517	1.28%	0.122%
2	3.687	133	136	141	rVB4	9095	15477	0.34%	0.032%
3	3.775	147	151	155	rVB2	15273	22841	0.50%	0.048%
4	4.281	232	237	241	rVV	11660	18484	0.40%	0.038%
5	5.581	451	458	465	rVB2	21977	39397	0.86%	0.082%
6	5.645	465	469	474	rBV2	11427	22402	0.49%	0.047%
7	6.534	614	620	624	rBV6	7772	16962	0.37%	0.035%
8	6.634	632	637	640	rBV5	10704	14888	0.32%	0.031%
9	6.686	640	646	657	rVB	176304	299452	6.53%	0.623%
10	6.839	661	672	687	rBV	639966	1008563	21.99%	2.099%
11	7.034	698	705	713	rVB	660844	1035511	22.58%	2.155%
12	7.492	774	783	788	rBV	689028	1061834	23.16%	2.210%
13	7.539	788	791	793	rVV	119944	164740	3.59%	0.343%
14	7.575	793	797	807	rVB	324429	476982	10.40%	0.993%
15	7.745	822	826	831	rBV7	6991	13344	0.29%	0.028%
16	7.798	831	835	842	rVB7	13651	23232	0.51%	0.048%
17	8.204	898	904	915	rVV	506312	773759	16.87%	1.610%
18	8.310	915	922	929	rVB2	46900	80849	1.76%	0.168%
19	8.580	962	968	972	rBV	170916	274813	5.99%	0.572%
20	8.633	972	977	983	rVV	818296	1268570	27.66%	2.640%
21	8.686	983	986	993	rVB	38684	61476	1.34%	0.128%
22	8.828	1005	1010	1015	rVB5	11562	22917	0.50%	0.048%
23	9.110	1053	1058	1065	rVV6	10062	20693	0.45%	0.043%
24	9.186	1067	1071	1080	rVB10	11562	24700	0.54%	0.051%
25	9.351	1092	1099	1106	rBV	636727	1013939	22.11%	2.110%
26	9.416	1106	1110	1117	rVB3	32527	59403	1.30%	0.124%
27	9.533	1122	1130	1135	rVV8	13051	27061	0.59%	0.056%
28	9.616	1141	1144	1151	rVV7	8429	17409	0.38%	0.036%
29	9.680	1151	1155	1161	rVB7	10510	16810	0.37%	0.035%
30	9.822	1172	1179	1182	rBV7	10658	19677	0.43%	0.041%
31	9.886	1182	1190	1198	rVV	903634	1450650	31.63%	3.019%
32	10.045	1212	1217	1223	rVB9	11276	23703	0.52%	0.049%
33	10.122	1223	1230	1232	rBV8	13593	25542	0.56%	0.053%
34	10.169	1233	1238	1246	rBV	143241	261165	5.70%	0.544%

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35	10.251	1246	1252	1258	rBV	893474	1409136	30.73%	2.933%
36	10.304	1258	1261	1263	rBV	56213	75937	1.66%	0.158%
37	10.398	1271	1277	1285	rVB2	46358	91654	2.00%	0.191%
38	10.498	1288	1294	1301	rVB5	9467	19030	0.41%	0.040%
39	10.569	1301	1306	1311	rVB8	8075	13543	0.30%	0.028%
40	10.804	1341	1346	1354	rVB10	6940	14682	0.32%	0.031%
41	10.904	1356	1363	1372	rVB	37980	72868	1.59%	0.152%
42	11.074	1385	1392	1397	rVV7	13304	25326	0.55%	0.053%
43	11.439	1449	1454	1463	rBV4	17581	34031	0.74%	0.071%
44	11.557	1469	1474	1484	rVV	71300	127828	2.79%	0.266%
45	11.645	1484	1489	1495	rVB9	8189	18430	0.40%	0.038%
46	11.857	1518	1525	1531	rVV6	20464	45553	0.99%	0.095%
47	11.927	1531	1537	1544	rVB	13558	25607	0.56%	0.053%
48	12.398	1609	1617	1620	rBV7	11933	22431	0.49%	0.047%
49	12.486	1625	1632	1639	rBV	102465	165609	3.61%	0.345%
50	12.745	1671	1676	1682	rBV	178461	250544	5.46%	0.521%
51	13.051	1722	1728	1737	rBV	203053	322078	7.02%	0.670%
52	13.121	1737	1740	1744	rVB5	14172	22681	0.49%	0.047%
53	13.557	1808	1814	1819	rVV	1725494	2268267	49.46%	4.721%
54	13.604	1819	1822	1829	rVB	143615	187009	4.08%	0.389%
55	13.821	1852	1859	1869	rBV	1850028	2617378	57.08%	5.447%
56	14.004	1884	1890	1895	rVV3	18147	30810	0.67%	0.064%
57	14.133	1905	1912	1920	rBV	1245141	1798649	39.22%	3.743%
58	14.198	1920	1923	1926	rVV2	25349	37452	0.82%	0.078%
59	14.233	1926	1929	1934	rVB4	22666	33476	0.73%	0.070%
60	14.357	1944	1950	1959	rBV	129494	223582	4.88%	0.465%
61	14.539	1976	1981	1987	rVV	16793	26220	0.57%	0.055%
62	14.833	2023	2031	2038	rBV	55012	91556	2.00%	0.191%
63	14.974	2048	2055	2062	rVV	487409	638441	13.92%	1.329%
64	15.033	2062	2065	2069	rVB2	17469	21889	0.48%	0.046%
65	15.133	2075	2082	2088	rVV	2424957	3412180	74.41%	7.101%
66	15.192	2088	2092	2098	rVB2	27134	36807	0.80%	0.077%
67	15.262	2098	2104	2118	rBV	814370	1100663	24.00%	2.291%
68	15.374	2119	2123	2129	rVB2	30352	46426	1.01%	0.097%
69	15.509	2139	2146	2154	rBV	44972	76149	1.66%	0.158%
70	15.809	2194	2197	2204	rBV6	16464	26212	0.57%	0.055%
71	15.921	2209	2216	2220	rVB7	11220	18107	0.39%	0.038%

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72	16.080	2240	2243	2250	rVB4	9955	19229	0.42%	0.040%
73	16.633	2331	2337	2340	rBV5	13876	21742	0.47%	0.045%
74	16.886	2373	2380	2384	rVV	1483281	2057499	44.87%	4.282%
75	16.921	2384	2386	2390	rVV	60957	75994	1.66%	0.158%
76	16.980	2390	2396	2410	rVB	2768958	3910065	85.27%	8.137%
77	17.292	2445	2449	2453	rVB	13401	16765	0.37%	0.035%
78	17.386	2460	2465	2470	rBV7	10818	18932	0.41%	0.039%
79	17.574	2492	2497	2505	rBV	742714	930532	20.29%	1.937%
80	17.739	2523	2525	2532	rVB6	10101	15866	0.35%	0.033%
81	17.809	2534	2537	2542	rBV4	27062	39659	0.86%	0.083%
82	17.886	2545	2550	2555	rVV3	24703	44867	0.98%	0.093%
83	18.050	2574	2578	2583	rBV4	15435	20670	0.45%	0.043%
84	18.562	2659	2665	2673	rBV7	15186	27656	0.60%	0.058%
85	18.915	2721	2725	2728	rBV	29784	43946	0.96%	0.091%
86	18.950	2728	2731	2737	rVB3	23733	29072	0.63%	0.061%
87	19.174	2764	2769	2775	rBV2	36005	52586	1.15%	0.109%
88	19.286	2781	2788	2796	rBV	3427537	4585701	100.00%	9.543%
89	20.103	2924	2927	2932	rBV5	16635	24840	0.54%	0.052%
90	20.780	3038	3042	3046	rBV	42150	62615	1.37%	0.130%
91	20.833	3046	3051	3055	rBV2	170361	265071	5.78%	0.552%
92	20.880	3055	3059	3065	rVB	201631	224993	4.91%	0.468%
93	21.080	3088	3093	3103	rBV	1934506	2395061	52.23%	4.984%
94	21.691	3195	3197	3201	rVB2	81501	85050	1.85%	0.177%
95	22.133	3266	3272	3279	rBV4	169560	332688	7.25%	0.692%
96	22.609	3350	3353	3357	rVB	177252	216960	4.73%	0.452%
97	23.068	3425	3431	3439	rBV	2573413	4358718	95.05%	9.071%
98	23.138	3439	3443	3447	rVV	217545	317028	6.91%	0.660%
99	23.197	3448	3453	3464	rVB2	1367292	2403513	52.41%	5.002%
100	23.738	3541	3545	3552	rVB	165816	269813	5.88%	0.562%

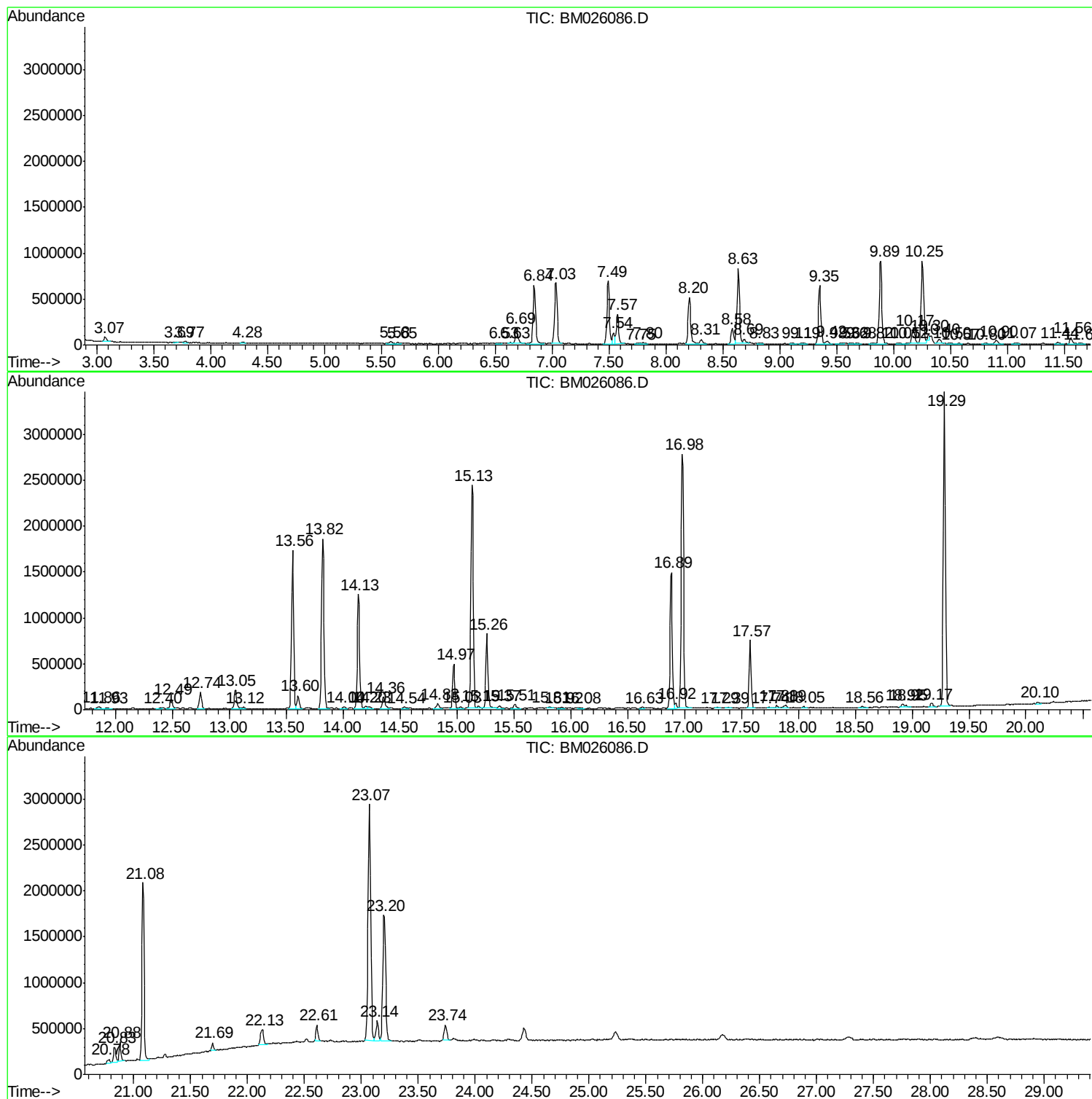
Sum of corrected areas: 48051134

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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



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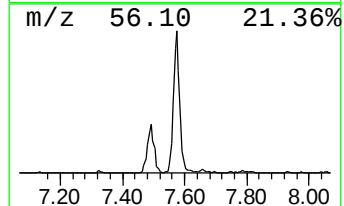
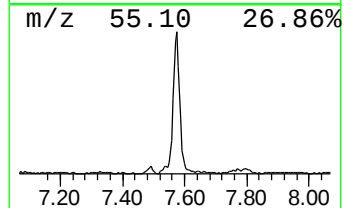
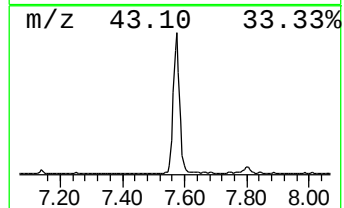
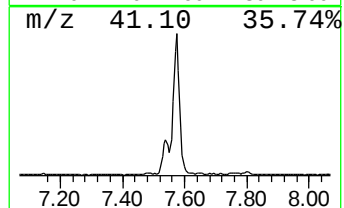
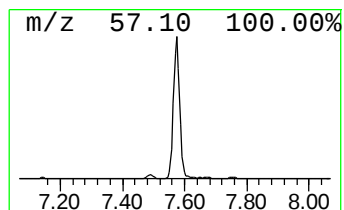
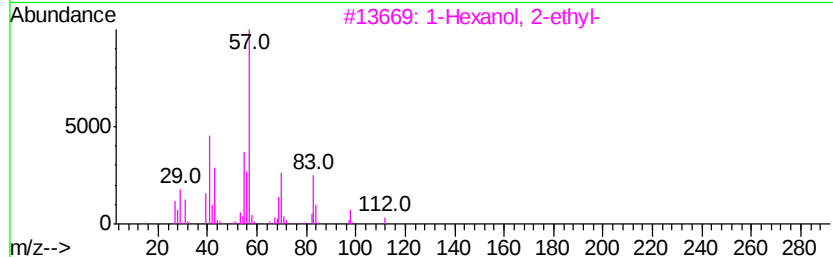
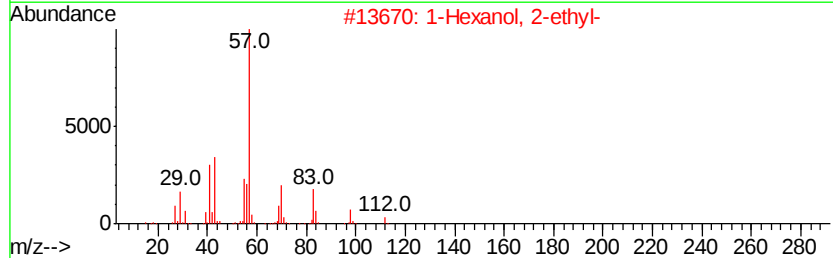
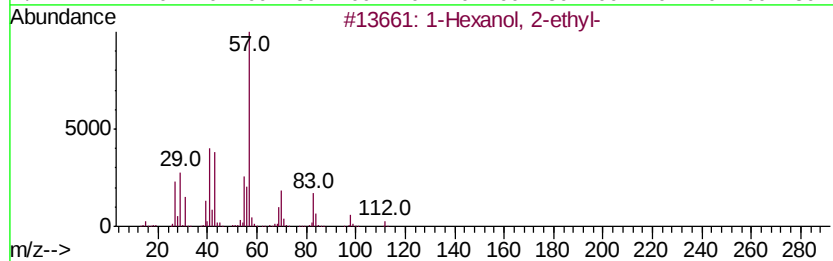
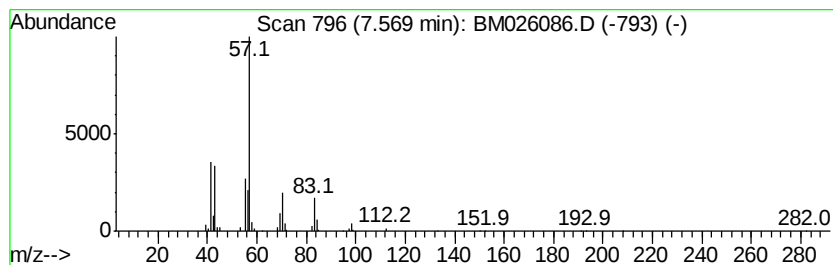
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Hexanol, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	8.98 ng/ul	476982	1,4-Dichlorobenzene-d4	7.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	74
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
4			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
5			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50



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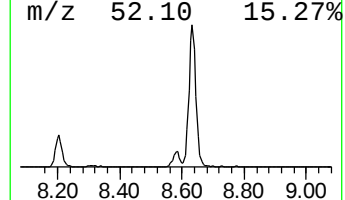
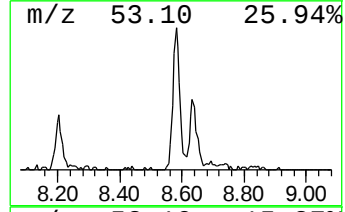
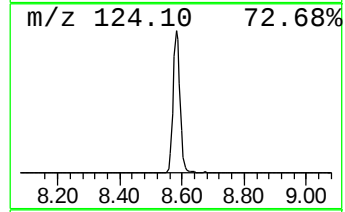
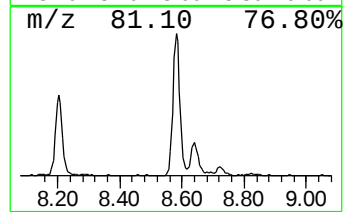
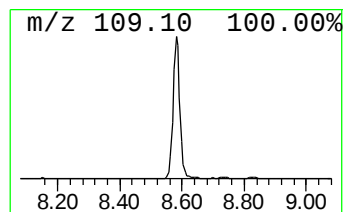
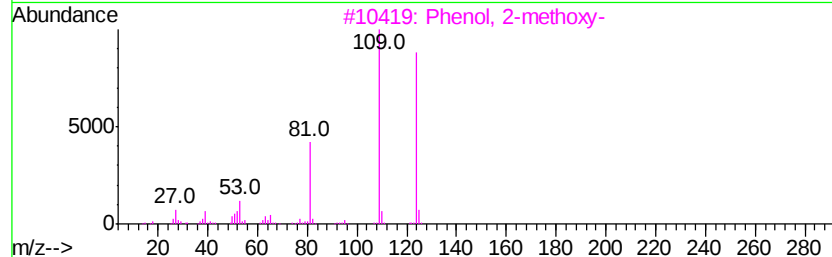
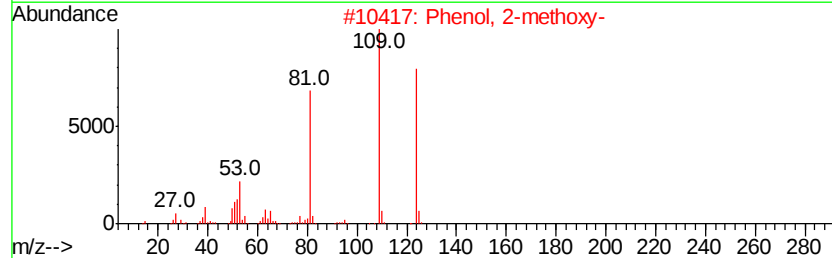
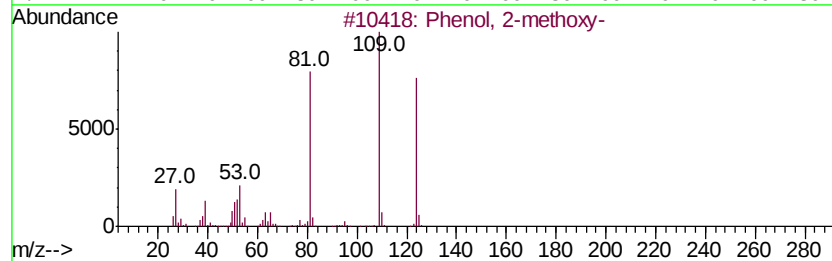
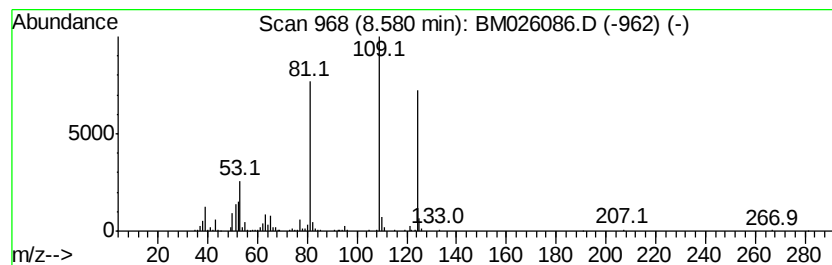
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenol, 2-methoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	5.18 ng/ul	274813	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	97
2		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	96
3		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	95
4		Mequinol	124	C7H8O2	000150-76-5	91
5		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	90



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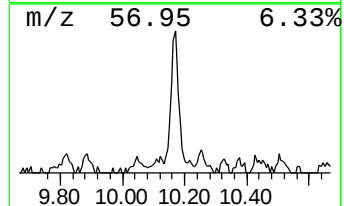
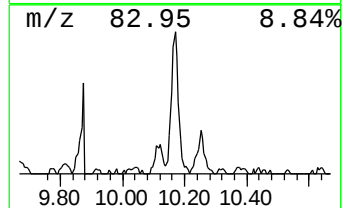
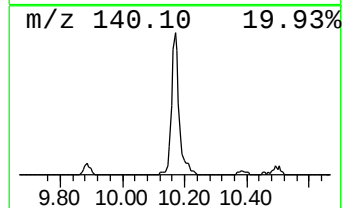
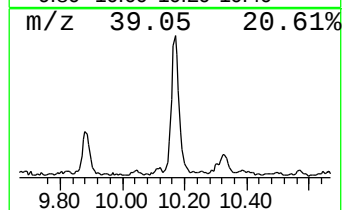
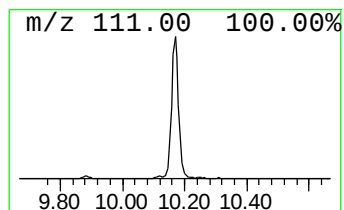
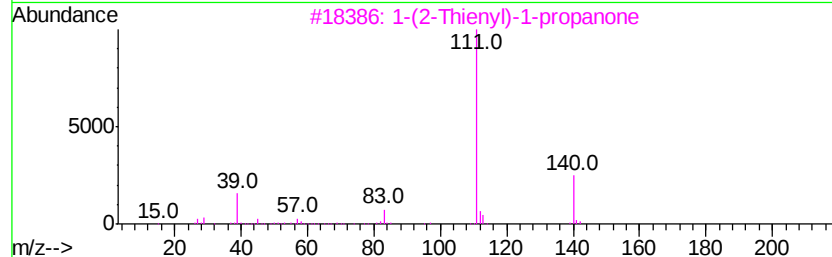
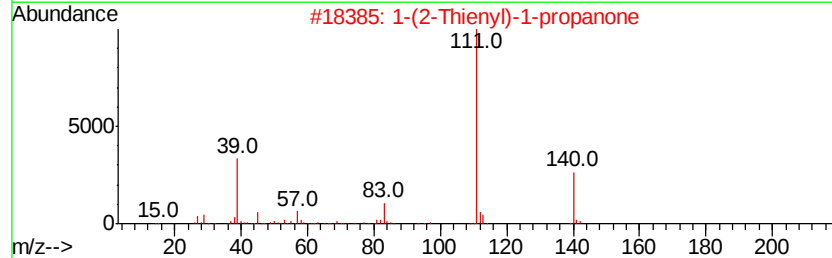
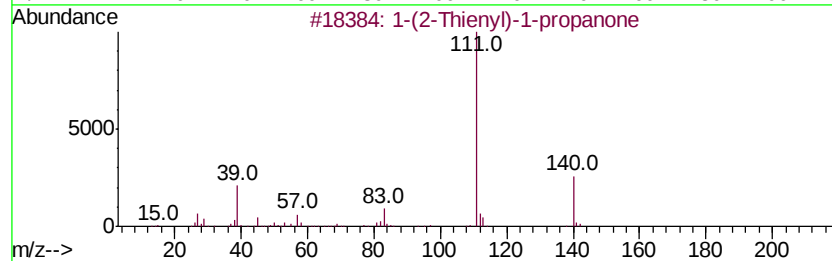
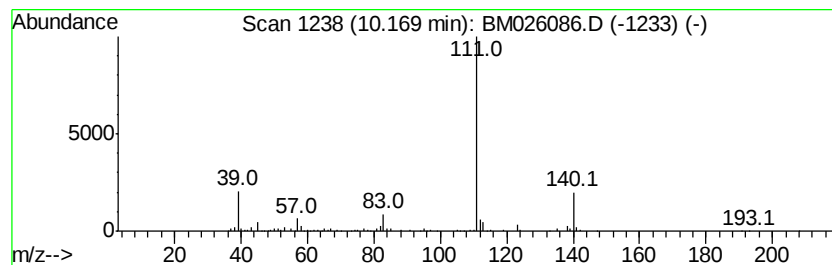
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-(2-Thienyl)-1-propanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.17	3.71 ng/ul	261165	Naphthalene-d8	10.25

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(2-Thienyl)-1-propanone	140	C7H8OS	013679-75-9	93
2		1-(2-Thienyl)-1-propanone	140	C7H8OS	013679-75-9	90
3		1-(2-Thienyl)-1-propanone	140	C7H8OS	013679-75-9	86
4		Glycine, N-(2-thienylcarbonyl)-,...	199	C8H9NO3S	1000299-60-2	78
5		Thiophene-2-carboxylic acid [1,2...	194	C7H6N4OS	349127-13-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026086.D
Acq On : 22 May 2020 19:08
Operator : MA/SJ
Sample : L2737-18
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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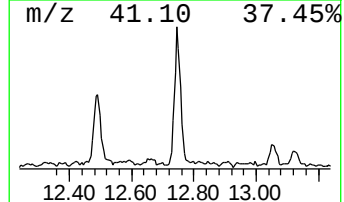
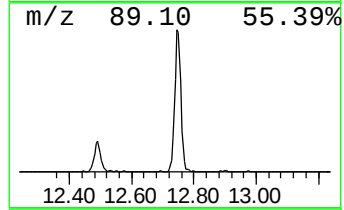
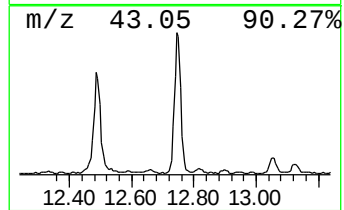
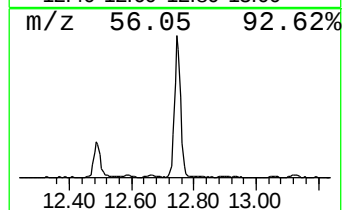
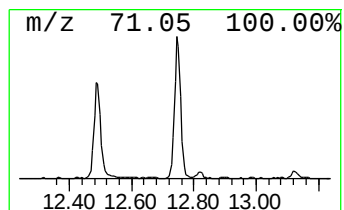
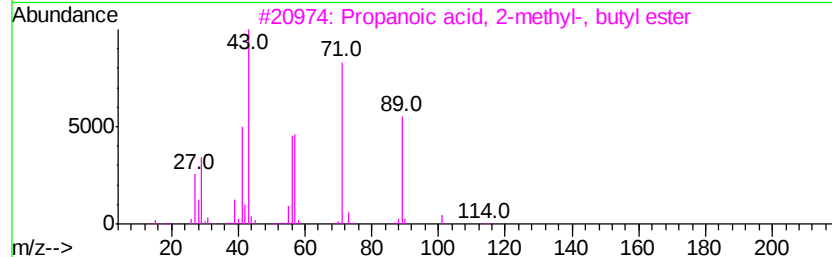
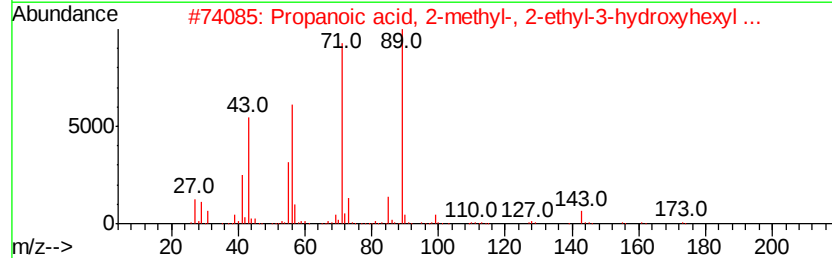
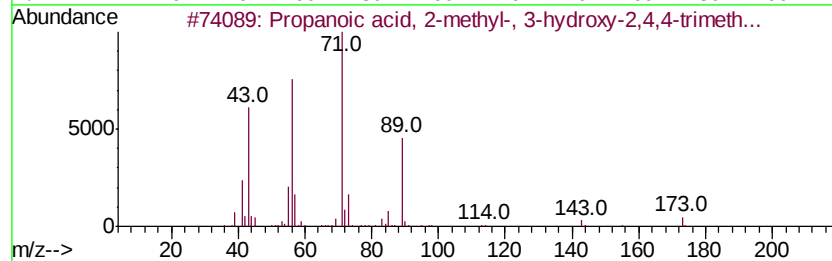
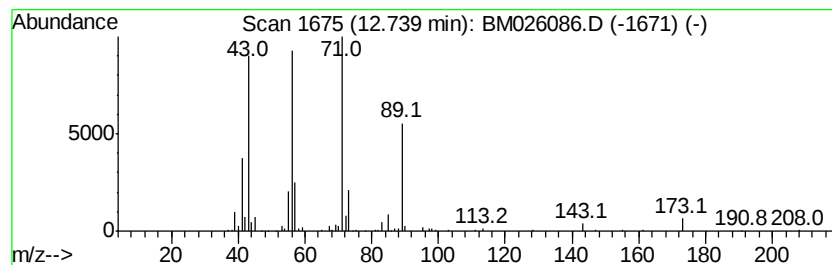
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Propanoic acid, 2-methyl-, ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	2.79 ng/ul	250544	Acenaphthene-d10	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, 3-hyd...	216	C12H24O3	074367-34-3	91
2		Propanoic acid, 2-methyl-, 2-eth...	216	C12H24O3	074367-31-0	83
3		Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	53
4		cis-2,3-Epoxyoctane	128	C8H16O	023024-54-6	50
5		Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026086.D
Acq On : 22 May 2020 19:08
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Sample : L2737-18
Misc :
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ClientSampleId :
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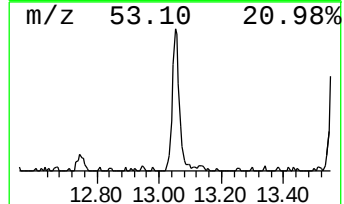
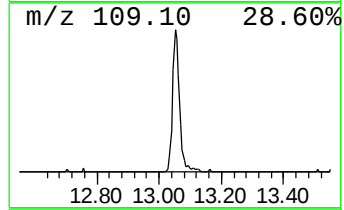
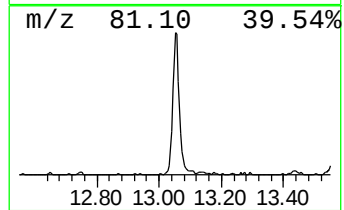
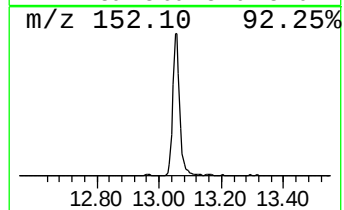
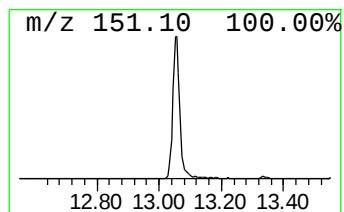
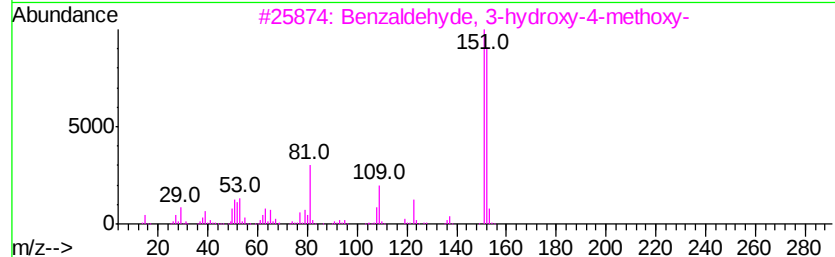
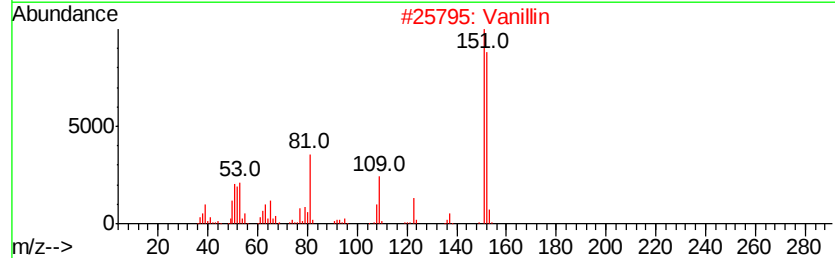
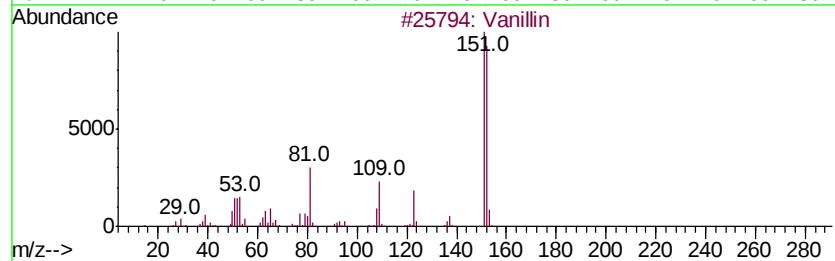
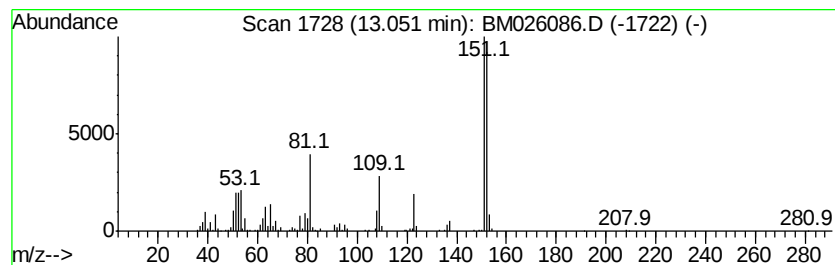
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Vanillin Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.05	3.58 ng/ul	322078	Acenaphthene-d10	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vanillin	152	C8H8O3	000121-33-5	97
2		Vanillin	152	C8H8O3	000121-33-5	97
3		Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96
4		Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	95
5		Vanillin	152	C8H8O3	000121-33-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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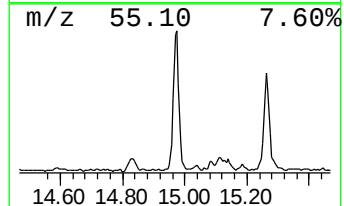
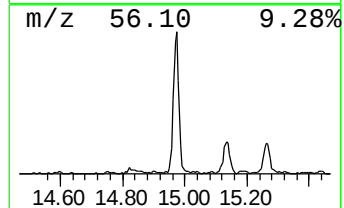
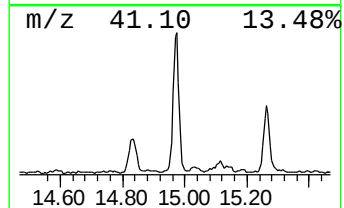
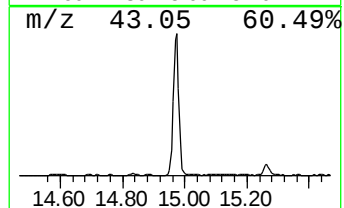
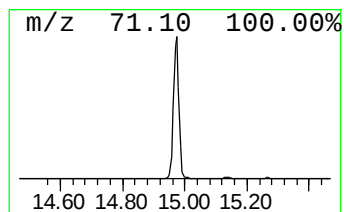
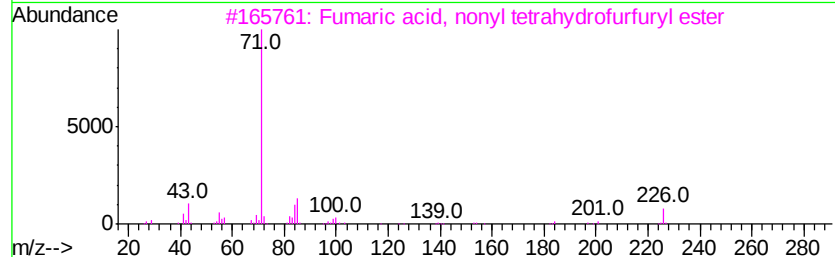
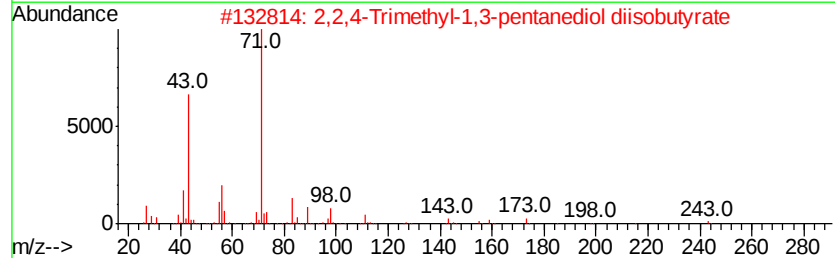
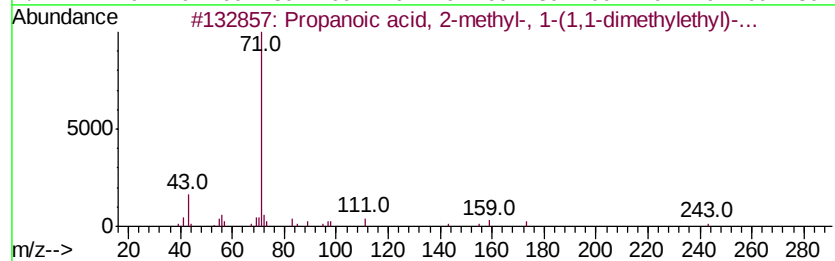
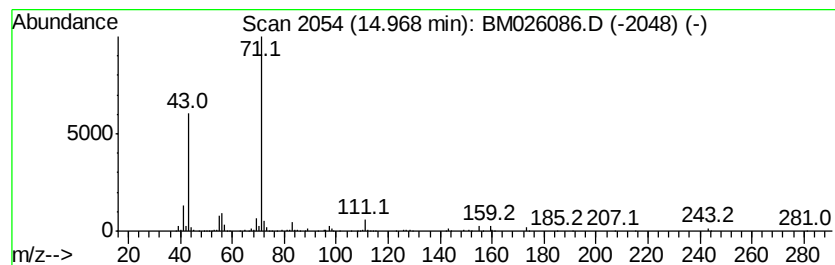
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Propanoic acid, 2-methyl-, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	7.10 ng/ul	638441	Acenaphthene-d10	14.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 1-(1,...	286	C16H30O4	074381-40-1	78	
2	2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	56	
3	Fumaric acid, nonyl tetrahydrofu...	326	C18H30O5	1000330-58-3	45	
4	Oxalic acid, decyl neopentyl ester	300	C17H32O4	1000309-73-4	45	
5	Diethylmalonic acid, monochlorid...	262	C12H19ClO4	1000370-65-0	42	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026086.D
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Sample : L2737-18
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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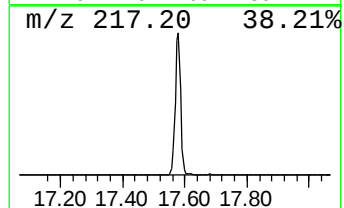
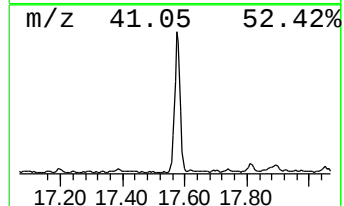
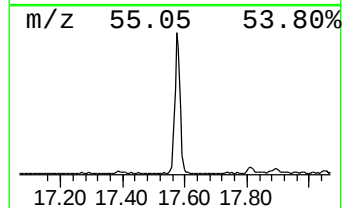
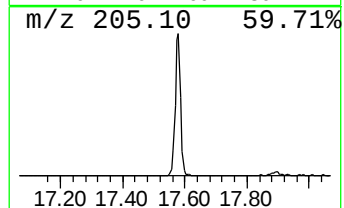
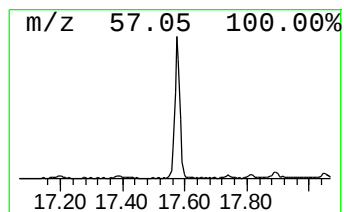
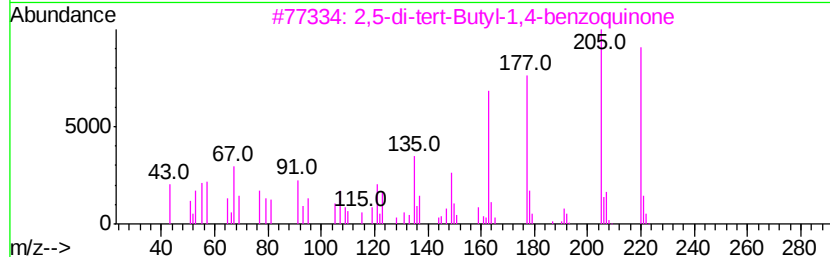
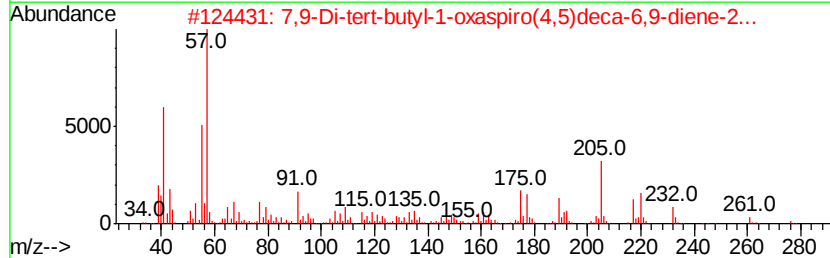
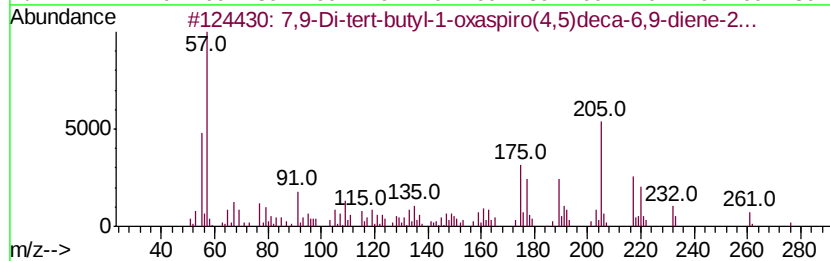
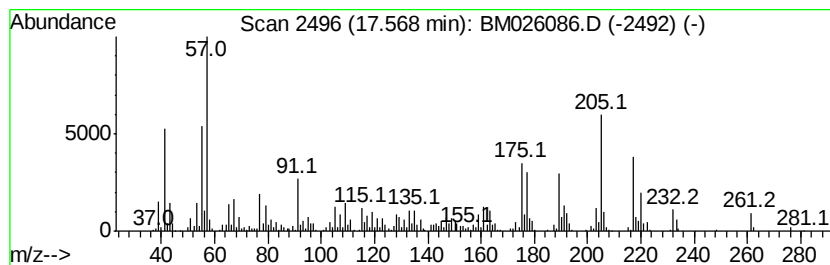
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.57	9.05 ng/ul	930532	Phenanthrene-d10	16.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,9-Di-tert-butyl-1-oxaspiro(4,5)deca-6,9-diene-2-one	276	C17H24O3	082304-66-3	99
2			7,9-Di-tert-butyl-1-oxaspiro(4,5)deca-6,9-diene-2-one	276	C17H24O3	082304-66-3	98
3			2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H20O2	002460-77-7	48
4			2-Methyl-4,5-dimethoxycrotonophenone	220	C13H16O3	207233-94-1	30
5			1-Ethyl-2-methylphenanthrene	220	C17H16	061983-53-7	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026086.D
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Operator : MA/SJ
Sample : L2737-18
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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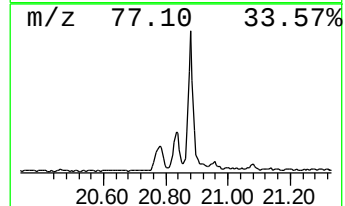
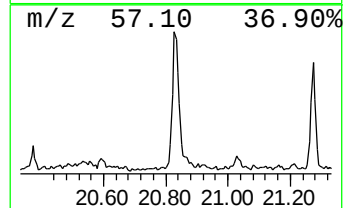
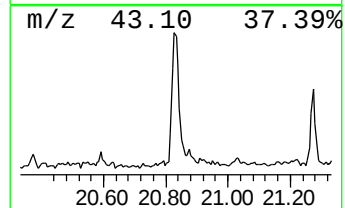
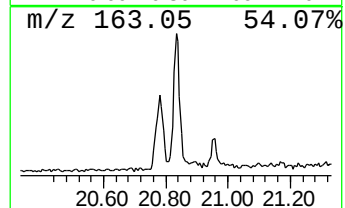
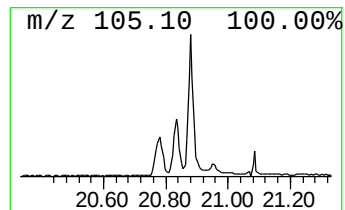
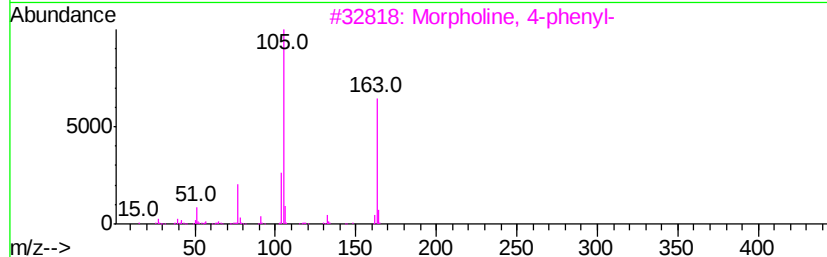
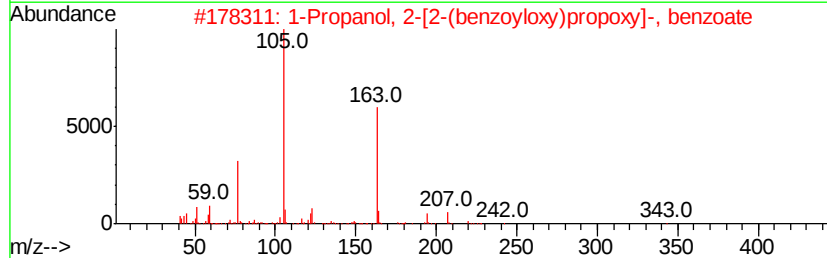
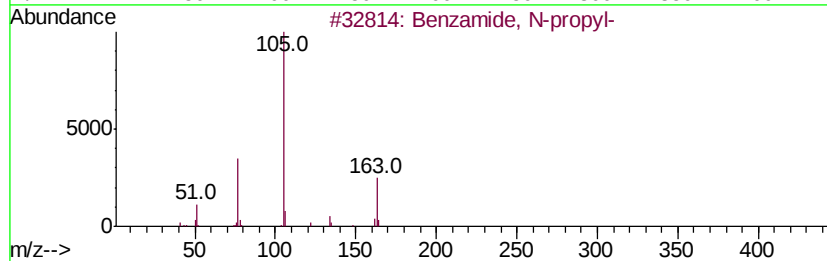
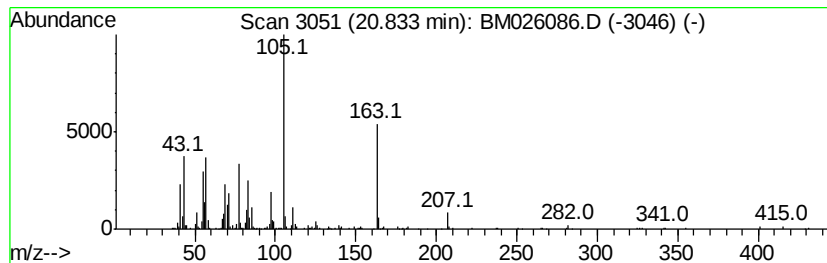
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	2.21 ng/ul	265071	Chrysene-d12	21.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, N-propyl-	163	C10H13NO	010546-70-0	43
2		1-Propanol, 2-[2-(benzoyloxy)propoxy]-, benzoate	342	C20H22O5	020109-39-1	43
3		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	38
4		Benzenepropanamide, .alpha.-(benzoyloxy)-, benzoyl ester	284	C16H16N2O3	058690-81-6	37
5		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	27



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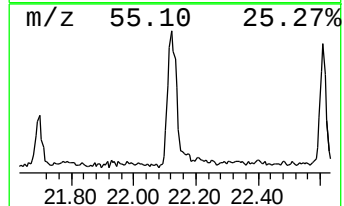
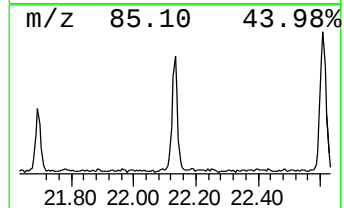
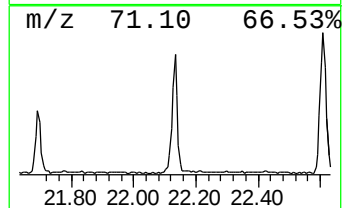
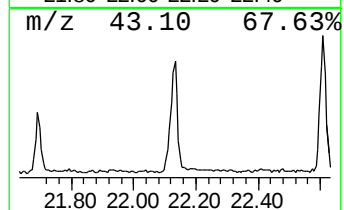
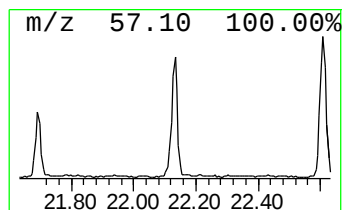
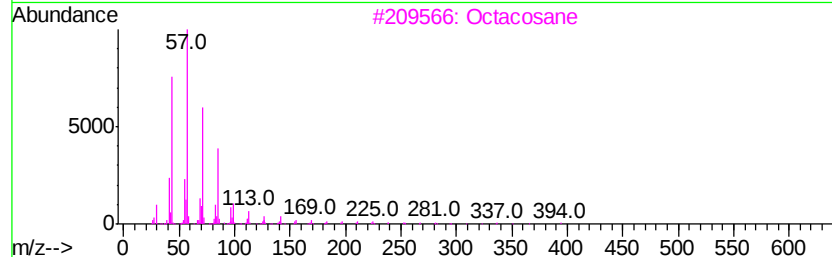
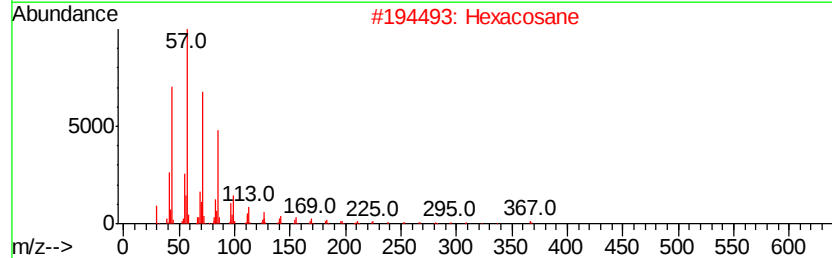
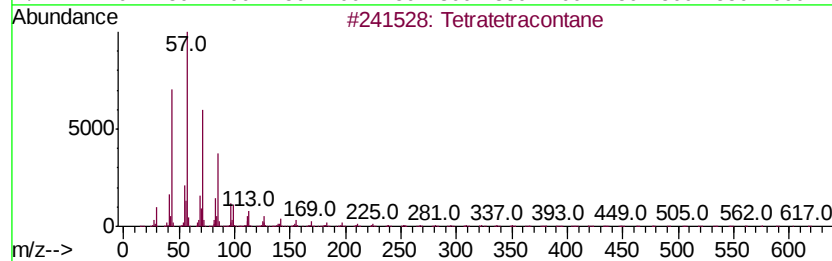
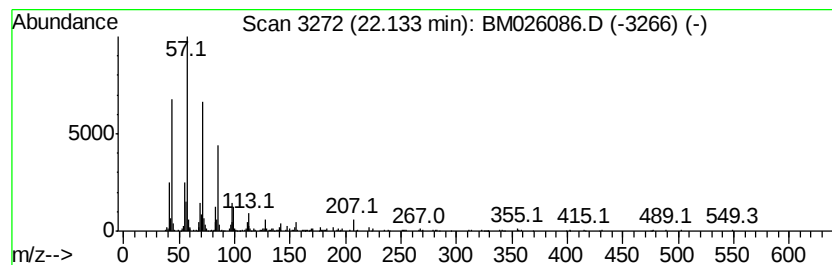
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.13	2.78 ng/ul	332688	Chrysene-d12	21.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	90
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Eicosane, 9-octyl-	394	C28H58	013475-77-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026086.D
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Sample : L2737-18
Misc :
ALS Vial : 11 Sample Multiplier: 1

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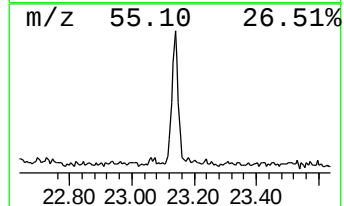
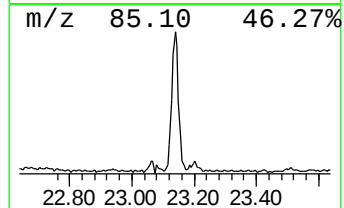
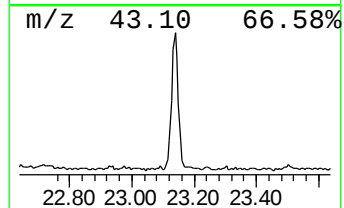
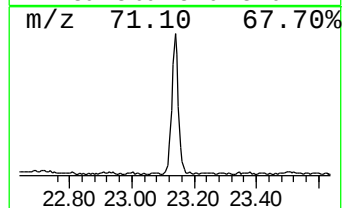
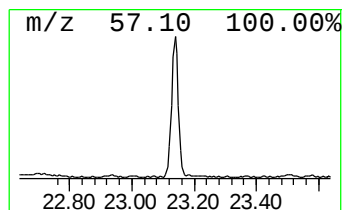
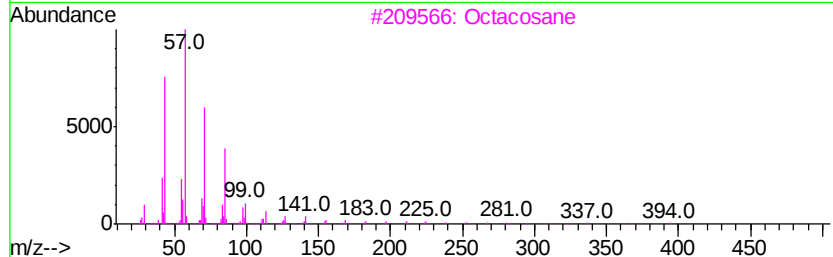
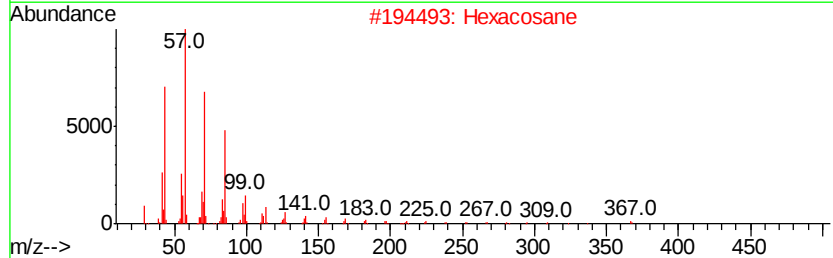
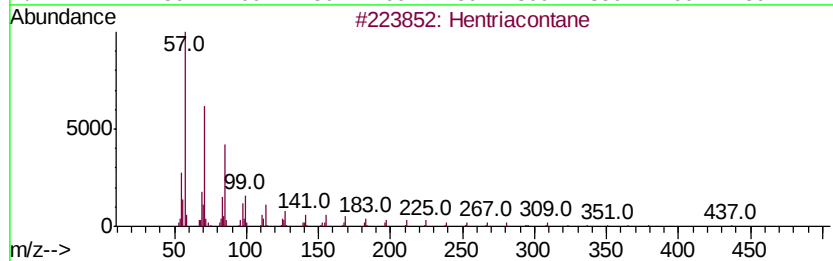
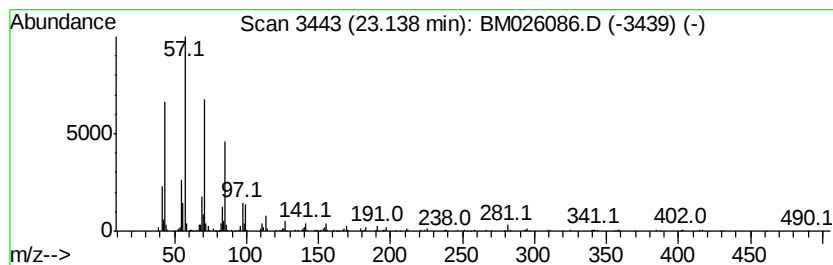
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.14	2.64 ng/ul	317028	Perylene-d12	23.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	90
2			Hexacosane	366	C26H54	000630-01-3	90
3			Octacosane	394	C28H58	000630-02-4	90
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	87
5			Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026086.D
Acq On : 22 May 2020 19:08
Operator : MA/SJ
Sample : L2737-18
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F9B40

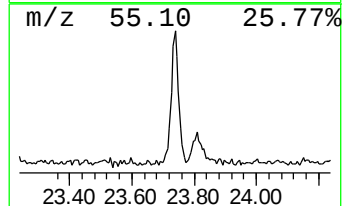
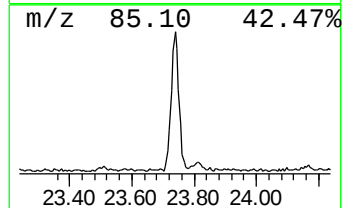
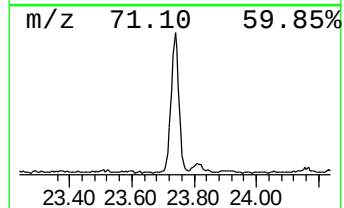
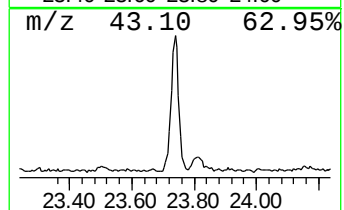
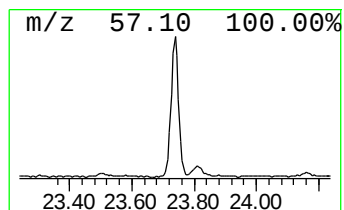
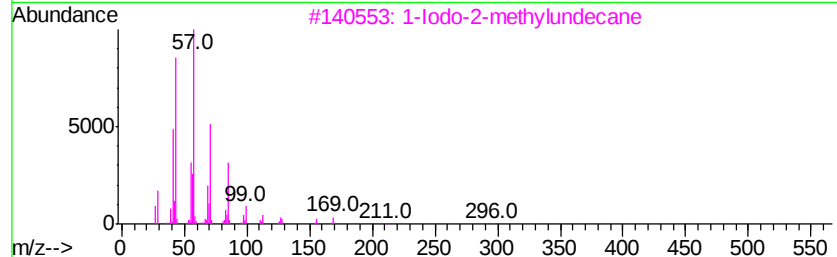
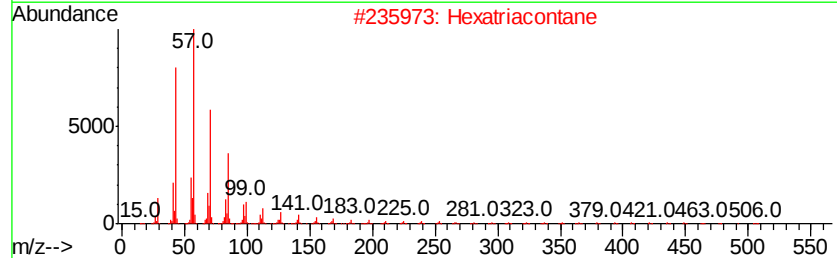
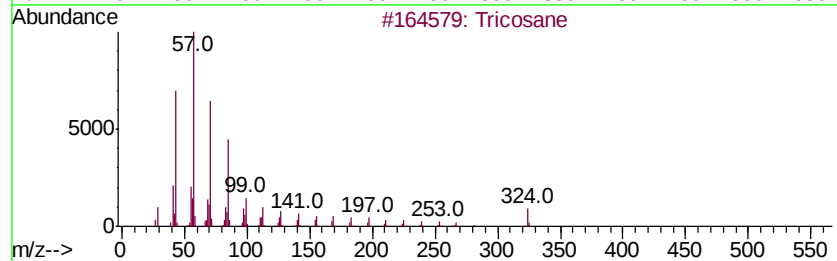
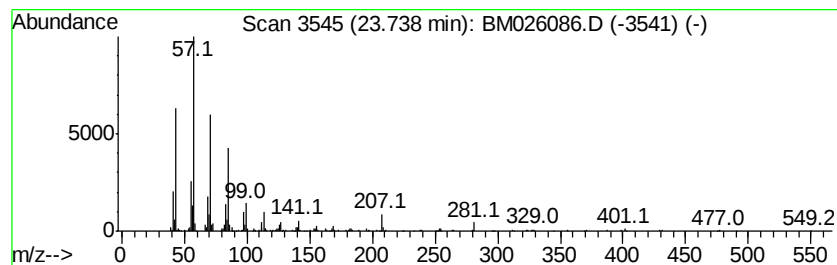
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.74	2.25 ng/ul	269813	Perylene-d12	23.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	93
2			Hexatriacontane	507	C36H74	000630-06-8	90
3			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
4			Heneicosane	296	C21H44	000629-94-7	90
5			Tricosane	324	C23H48	000638-67-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052220\
 Data File : BM026086.D
 Acq On : 22 May 2020 19:08
 Operator : MA/SJ
 Sample : L2737-18
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F9B40

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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
1-Hexanol, 2-ethyl-	7.57	9.0	ng/ul	476982	1	7.49	1061830	20.0
Phenol, 2-methoxy-	8.58	5.2	ng/ul	274813	1	7.49	1061830	20.0
1-(2-Thienyl)-1-p...	10.17	3.7	ng/ul	261165	2	10.25	1409140	20.0
Propanoic acid, 2...	12.74	2.8	ng/ul	250544	3	14.13	1798650	20.0
Vanillin	13.05	3.6	ng/ul	322078	3	14.13	1798650	20.0
Propanoic acid, 2...	14.97	7.1	ng/ul	638441	3	14.13	1798650	20.0
7,9-Di-tert-butyl...	17.57	9.1	ng/ul	930532	4	16.89	2057500	20.0
unknown-01	20.83	2.2	ng/ul	265071	5	21.08	2395060	20.0
(DEL) Alkane: Str...	22.13	2.8	ng/ul	332688	5	21.08	2395060	20.0
(DEL) Alkane: Str...	23.14	2.6	ng/ul	317028	6	23.20	2403510	20.0
(DEL) Alkane: Str...	23.74	2.3	ng/ul	269813	6	23.20	2403510	20.0