

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
 Data File : BG025073.D
 Acq On : 10 Dec 2016 17:58
 Operator : UM/SJ
 Sample : H5861-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DA8D8

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 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: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.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	3.595	126	129	132	rVB4	13297	17101	0.46%	0.053%
2	3.695	142	146	148	rBV5	7055	12060	0.33%	0.037%
3	4.136	218	221	225	rVB5	9015	11515	0.31%	0.036%
4	4.359	253	259	263	rBV7	8629	16179	0.44%	0.050%
5	4.677	306	313	315	rBV6	10717	22333	0.61%	0.069%
6	4.900	344	351	360	rVB5	56947	104433	2.83%	0.325%
7	5.299	415	419	423	rBV6	8120	13934	0.38%	0.043%
8	5.646	471	478	480	rBV5	5497	12982	0.35%	0.040%
9	6.439	607	613	616	rBV6	7403	13456	0.36%	0.042%
10	6.739	658	664	667	rBV5	6230	11520	0.31%	0.036%
11	7.191	733	741	752	rBV2	52201	116327	3.16%	0.362%
12	7.291	752	758	764	rVB2	26089	46033	1.25%	0.143%
13	7.379	764	773	784	rVB3	95341	192351	5.22%	0.598%
14	7.550	794	802	815	rVB2	282858	471972	12.80%	1.467%
15	7.761	829	838	849	rBV	315496	577876	15.67%	1.796%
16	7.902	853	862	865	rBV8	6972	19398	0.53%	0.060%
17	8.237	912	919	929	rVB	259262	484224	13.13%	1.505%
18	8.519	965	967	973	rVB4	6600	11244	0.30%	0.035%
19	8.784	1007	1012	1016	rVB6	18029	33020	0.90%	0.103%
20	8.936	1029	1038	1047	rVV2	253794	479909	13.02%	1.492%
21	9.412	1111	1119	1127	rBV	399792	724311	19.65%	2.251%
22	9.489	1127	1132	1139	rVV4	46652	85451	2.32%	0.266%
23	9.553	1139	1143	1148	rVB6	14659	25402	0.69%	0.079%
24	9.735	1170	1174	1181	rBV4	15571	31575	0.86%	0.098%
25	9.964	1207	1213	1218	rVB8	7422	15582	0.42%	0.048%
26	10.135	1231	1242	1250	rBV2	358676	687735	18.65%	2.138%
27	10.335	1270	1276	1287	rVB6	39630	95355	2.59%	0.296%
28	10.493	1301	1303	1309	rVB5	7174	12436	0.34%	0.039%
29	10.540	1309	1311	1319	rVB7	7609	18135	0.49%	0.056%
30	10.675	1324	1334	1349	rVV	468722	902989	24.49%	2.807%
31	10.805	1349	1356	1367	rBV6	34298	84046	2.28%	0.261%
32	11.063	1392	1400	1412	rVB2	358408	660003	17.90%	2.051%
33	11.563	1478	1485	1490	rVB7	13573	28354	0.77%	0.088%
34	11.627	1490	1496	1501	rBV7	12817	21531	0.58%	0.067%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	11.815	1521	1528	1538	rBV3	100028	205665	5.58%	0.639%
36	11.956	1549	1552	1557	rVB6	8011	11299	0.31%	0.035%
37	12.232	1591	1599	1603	rBV6	20441	42321	1.15%	0.132%
38	12.385	1621	1625	1632	rVV9	6811	13003	0.35%	0.040%
39	12.526	1643	1649	1655	rVV7	9975	13031	0.35%	0.041%
40	12.573	1655	1657	1661	rVB3	9125	11066	0.30%	0.034%
41	12.632	1661	1667	1674	rBV8	14220	39924	1.08%	0.124%
42	12.749	1680	1687	1692	rBV9	10805	21939	0.60%	0.068%
43	12.826	1693	1700	1704	rVV6	6709	10659	0.29%	0.033%
44	12.937	1716	1719	1723	rBV4	6535	10465	0.28%	0.033%
45	13.108	1743	1748	1758	rVB10	21690	54453	1.48%	0.169%
46	13.202	1758	1764	1779	rBV2	169571	320422	8.69%	0.996%
47	13.313	1779	1783	1787	rBV5	7863	11538	0.31%	0.036%
48	13.449	1797	1806	1820	rVB	226343	382469	10.37%	1.189%
49	13.607	1832	1833	1841	rVB6	12894	17190	0.47%	0.053%
50	13.725	1845	1853	1857	rBV7	19241	42816	1.16%	0.133%
51	13.784	1857	1863	1877	rVB4	53698	118736	3.22%	0.369%
52	14.118	1917	1920	1923	rBV5	7834	11189	0.30%	0.035%
53	14.248	1928	1942	1954	rBV	1094524	1650676	44.77%	5.130%
54	14.353	1954	1960	1968	rBV10	28010	46537	1.26%	0.145%
55	14.459	1974	1978	1986	rBV8	34497	72024	1.95%	0.224%
56	14.553	1986	1994	2003	rVB	1019191	1637049	44.40%	5.088%
57	14.624	2004	2006	2014	rVB8	12992	21917	0.59%	0.068%
58	14.718	2017	2022	2025	rBV6	8712	11414	0.31%	0.035%
59	14.859	2038	2046	2056	rVB	628028	1011320	27.43%	3.143%
60	15.053	2074	2079	2090	rBV4	127389	250391	6.79%	0.778%
61	15.223	2102	2108	2115	rBV2	147507	233515	6.33%	0.726%
62	15.429	2140	2143	2149	rVB7	9183	16373	0.44%	0.051%
63	15.517	2155	2158	2161	rBV5	6889	11215	0.30%	0.035%
64	15.605	2165	2173	2177	rBV	170361	231557	6.28%	0.720%
65	15.646	2178	2180	2185	rVB3	33056	32744	0.89%	0.102%
66	15.769	2199	2201	2205	rVB4	11523	12085	0.33%	0.038%
67	15.846	2205	2214	2223	rBV	1434940	2273089	61.66%	7.065%
68	15.957	2226	2233	2247	rBV	674645	1070720	29.04%	3.328%
69	16.075	2250	2253	2261	rVB7	24339	43735	1.19%	0.136%
70	16.187	2268	2272	2276	rBV7	9787	15306	0.42%	0.048%
71	16.404	2303	2309	2315	rBV8	27767	57275	1.55%	0.178%

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Integration Parameters: LSCINT.P

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72	16.686	2350	2357	2364	rBV9	28495	66065	1.79%	0.205%
73	16.950	2395	2402	2406	rBV7	30753	55701	1.51%	0.173%
74	17.062	2417	2421	2428	rVB5	21465	39219	1.06%	0.122%
75	17.121	2429	2431	2436	rVB6	11925	14502	0.39%	0.045%
76	17.256	2449	2454	2457	rBV7	16325	24322	0.66%	0.076%
77	17.321	2461	2465	2470	rVB5	27290	45118	1.22%	0.140%
78	17.379	2473	2475	2481	rVV7	9354	16846	0.46%	0.052%
79	17.420	2481	2482	2486	rVB4	10623	12234	0.33%	0.038%
80	17.597	2505	2512	2521	rVB	916206	1439378	39.04%	4.474%
81	17.697	2521	2529	2540	rBV2	1803029	2852203	77.37%	8.865%
82	17.843	2549	2554	2560	rVV2	53409	81829	2.22%	0.254%
83	17.890	2560	2562	2565	rVB4	11667	11506	0.31%	0.036%
84	18.020	2583	2584	2587	rBV3	11127	11148	0.30%	0.035%
85	18.231	2617	2620	2629	rBV8	19617	24758	0.67%	0.077%
86	18.443	2652	2656	2660	rBV6	21661	32587	0.88%	0.101%
87	18.525	2666	2670	2676	rBV	59276	80451	2.18%	0.250%
88	18.737	2704	2706	2709	rBV4	9591	10421	0.28%	0.032%
89	18.972	2739	2746	2752	rBV4	22856	53282	1.45%	0.166%
90	19.165	2776	2779	2782	rVB4	10942	12286	0.33%	0.038%
91	19.336	2806	2808	2810	rBV3	15385	13579	0.37%	0.042%
92	19.377	2812	2815	2820	rBV6	35229	44228	1.20%	0.137%
93	19.606	2849	2854	2859	rBV	36585	43417	1.18%	0.135%
94	19.694	2867	2869	2874	rVB6	26065	34844	0.95%	0.108%
95	19.865	2893	2898	2904	rVB5	28847	54445	1.48%	0.169%
96	19.970	2909	2916	2926	rVB	2547518	3686665	100.00%	11.459%
97	21.886	3235	3242	3254	rVB	1077751	1939262	52.60%	6.027%
98	25.064	3772	3783	3797	rVB2	1200800	3478718	94.36%	10.812%
99	25.217	3806	3809	3813	rVB6	37676	51981	1.41%	0.162%
100	25.305	3814	3824	3834	rVB	598381	1845112	50.05%	5.735%

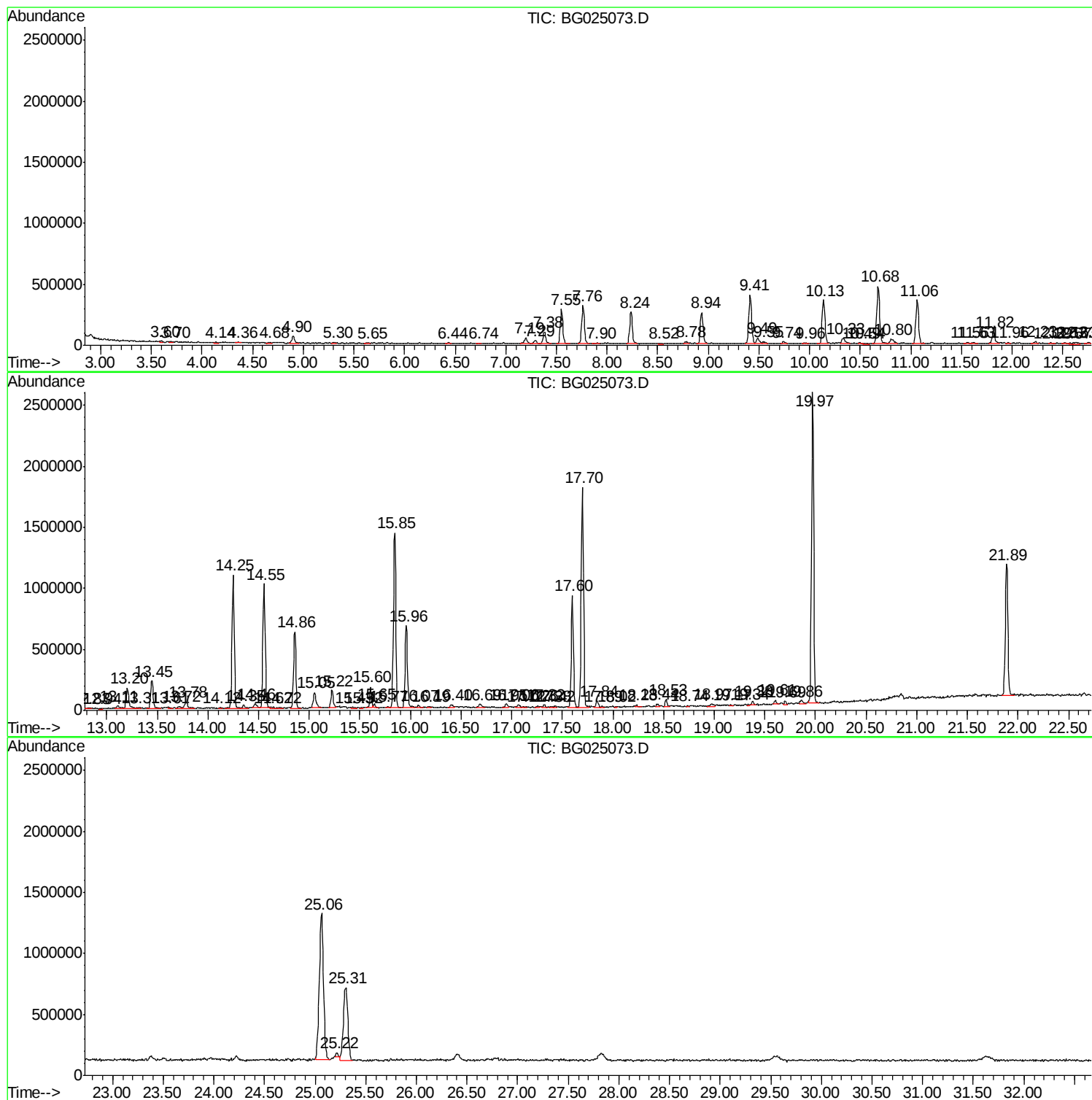
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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



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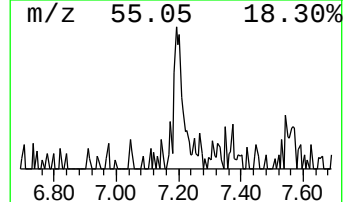
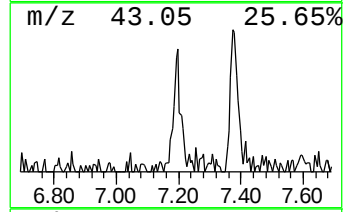
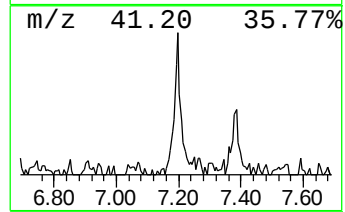
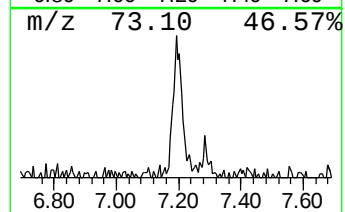
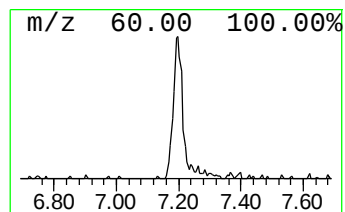
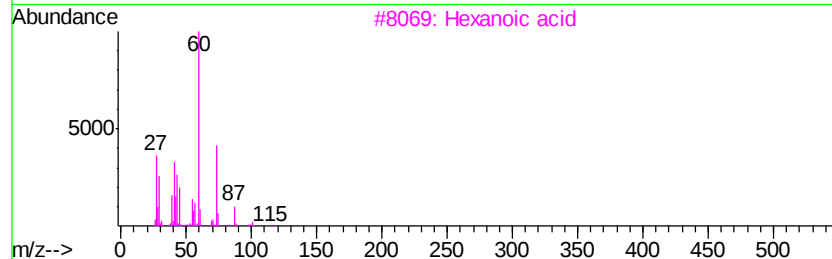
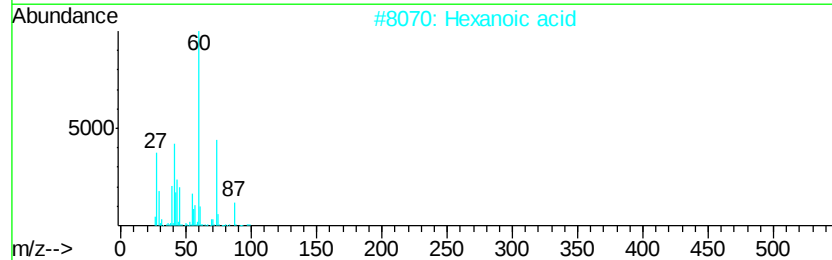
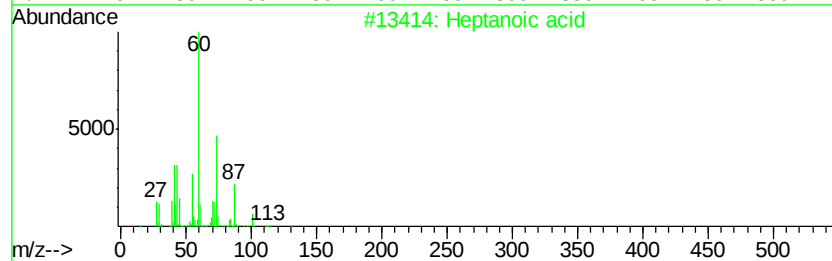
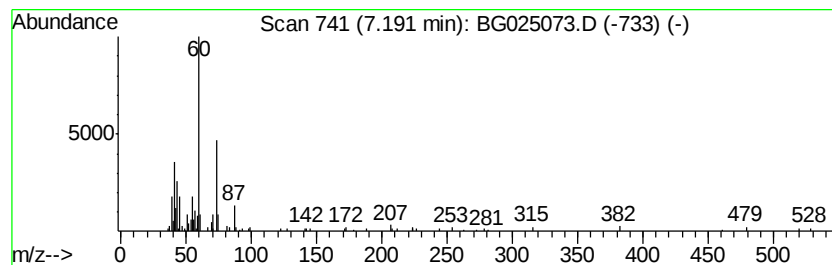
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Heptanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.19	4.80 ng/ul	116327	1,4-Dichlorobenzene-d4	8.24

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid	130	C7H14O2	000111-14-8	78	
2	Hexanoic acid	116	C6H12O2	000142-62-1	64	
3	Hexanoic acid	116	C6H12O2	000142-62-1	64	
4	8-Chlorocapric acid	178	C8H15ClO2	001795-62-6	64	
5	3-Methyl-hexanoic acid	130	C7H14O2	1000365-65-4	64	



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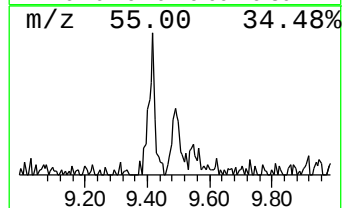
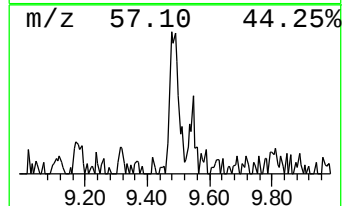
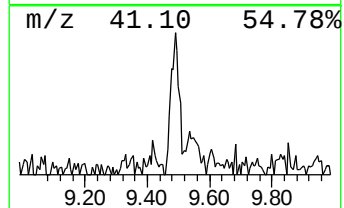
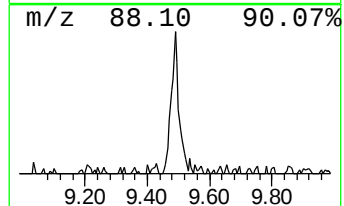
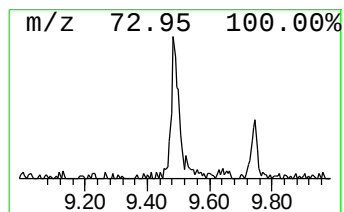
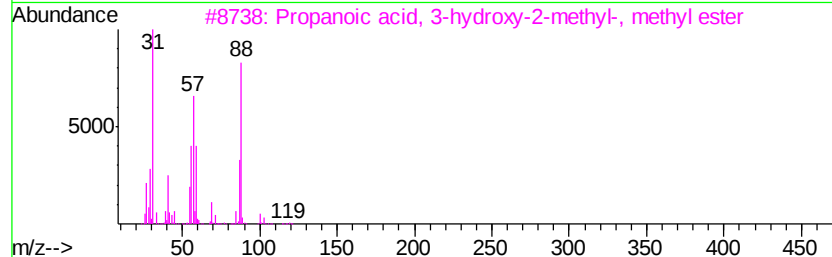
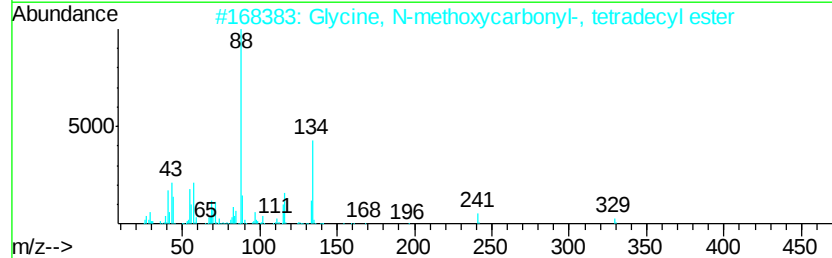
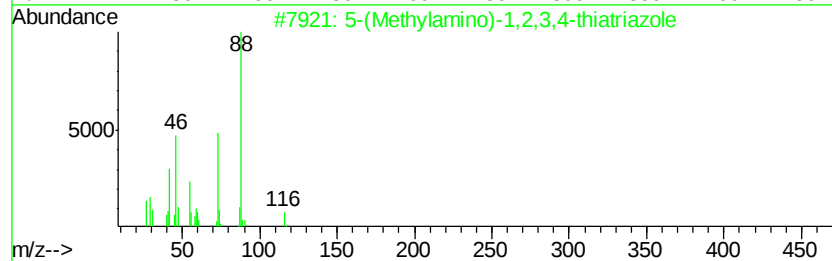
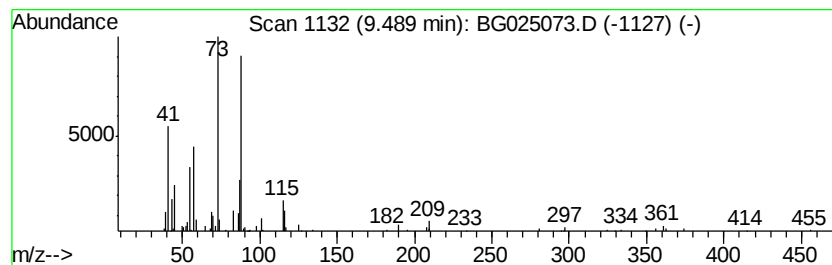
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 5-(Methylamino)-1,2,3,4-thi... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	3.53 ng/ul	85451	1,4-Dichlorobenzene-d4	8.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-(Methylamino)-1,2,3,4-thiatr...	116	C2H4N4S	052098-72-3	53
2			Glycine, N-methoxycarbonyl-, tet...	329	C18H35N04	1000313-49-4	43
3			Propanoic acid, 3-hydroxy-2-meth...	118	C5H10O3	042998-03-8	33
4			Acetic acid, diethyl-	116	C6H12O2	000088-09-5	33
5			Methyl (R)-(-)-3-hydroxy-2-methy...	118	C5H10O3	072657-23-9	33



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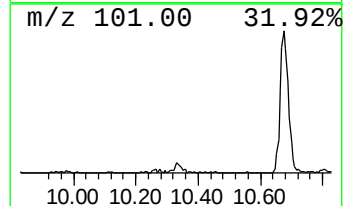
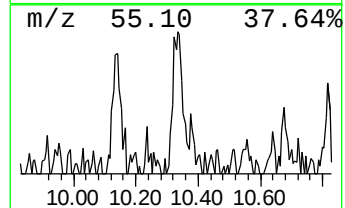
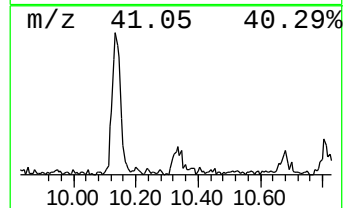
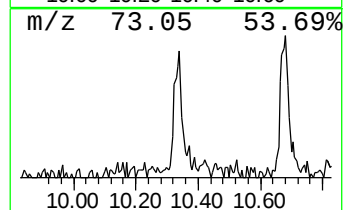
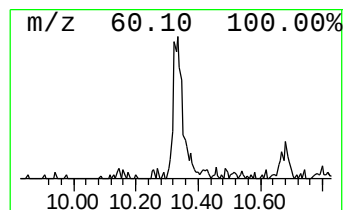
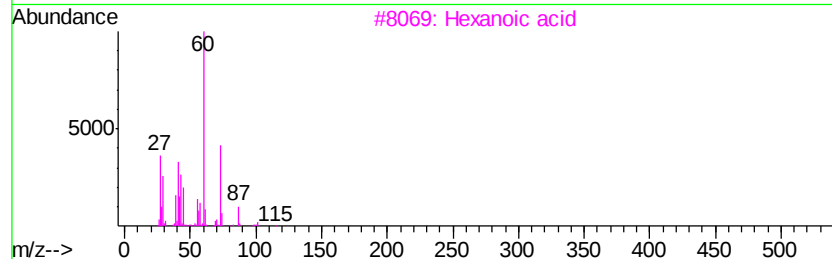
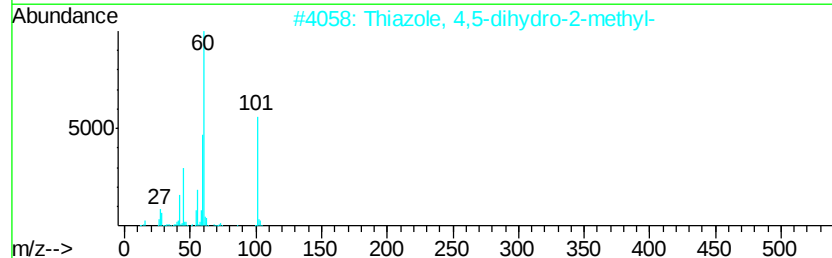
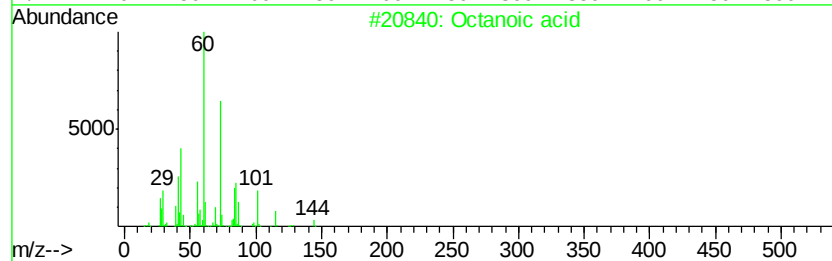
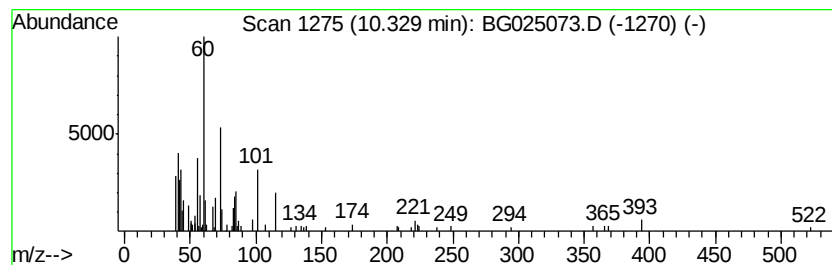
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Peak Number 4 Octanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	2.89 ng/ul	95355	Naphthalene-d8	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid		144	C8H16O2	000124-07-2	64
2	Thiazole, 4,5-dihydro-2-methyl-		101	C4H7NS	002346-00-1	59
3	Hexanoic acid		116	C6H12O2	000142-62-1	53
4	Hexanoic acid		116	C6H12O2	000142-62-1	50
5	.beta.-D-Ribopyranoside, methyl		164	C6H12O5	017289-61-1	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
Operator : UM/SJ
Sample : H5861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleID :
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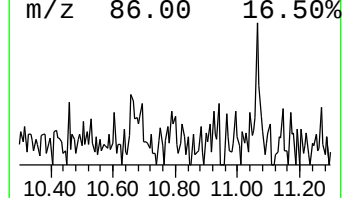
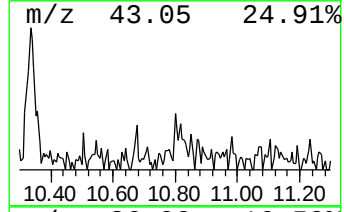
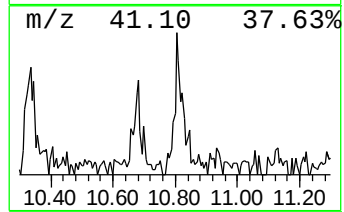
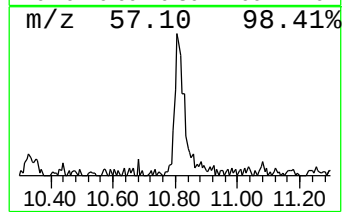
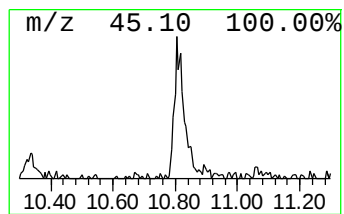
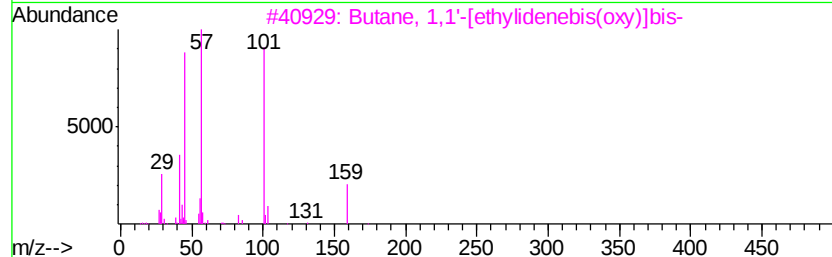
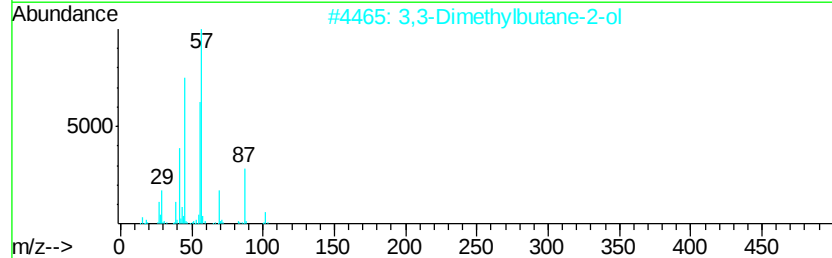
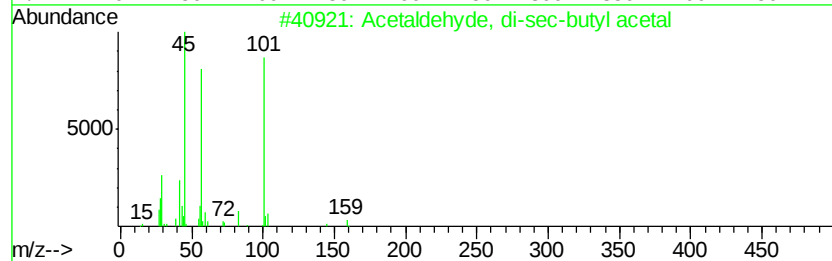
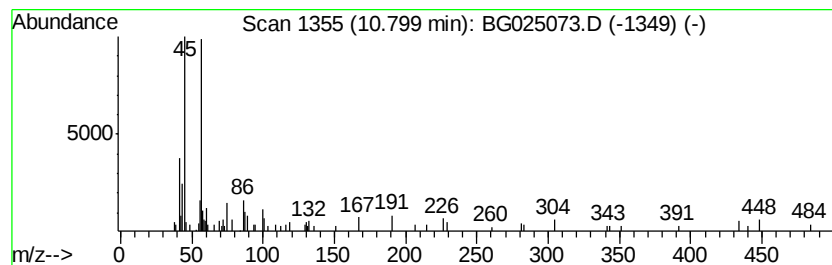
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Acetaldehyde, di-sec-butyl ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	2.55 ng/ul	84046	Naphthalene-d8	11.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetaldehyde, di-sec-butyl acetal	174	C10H22O2	005314-41-0	64
2		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	59
3		Butane, 1,1'-[ethylidenebis(oxy)]...	174	C10H22O2	000871-22-7	59
4		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	50
5		Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
Operator : UM/SJ
Sample : H5861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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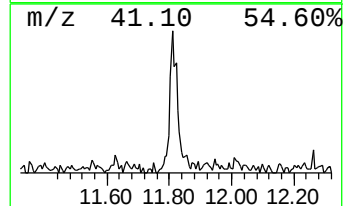
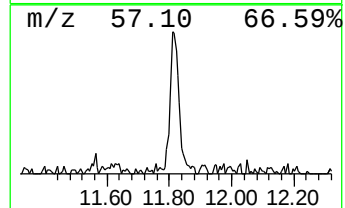
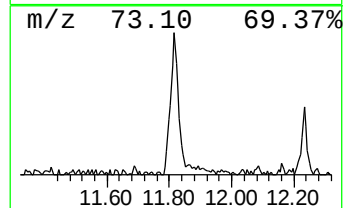
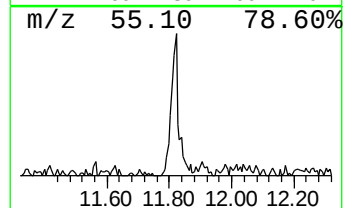
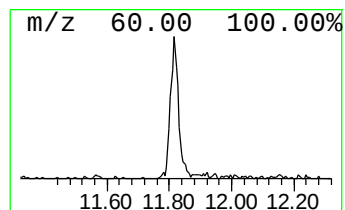
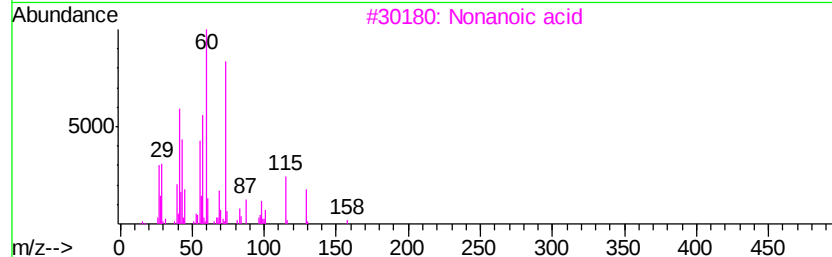
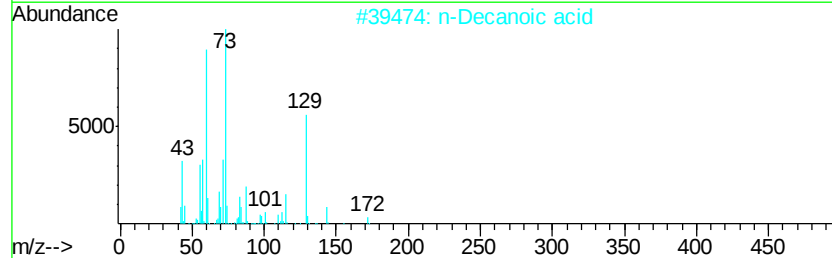
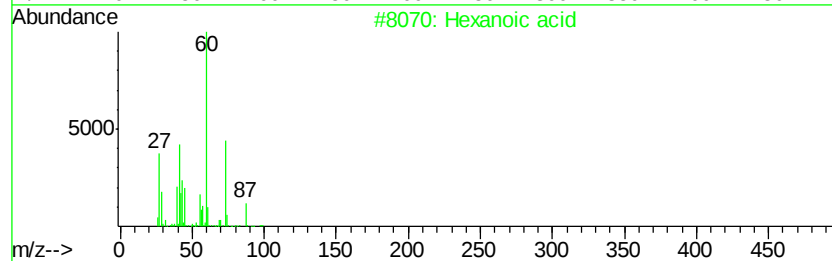
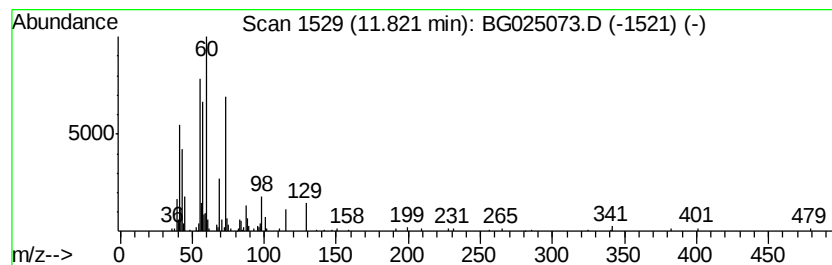
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	6.23 ng/ul	205665	Naphthalene-d8	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid		116	C6H12O2	000142-62-1	72
2	n-Decanoic acid		172	C10H20O2	000334-48-5	72
3	Nonanoic acid		158	C9H18O2	000112-05-0	64
4	Undecanoic acid, 10-bromo-		264	C11H21BrO2	018294-93-4	64
5	Pentanoic acid		102	C5H10O2	000109-52-4	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
Operator : UM/SJ
Sample : H5861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

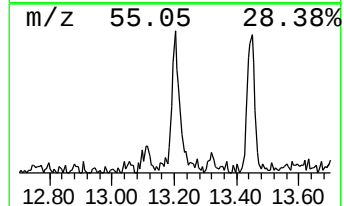
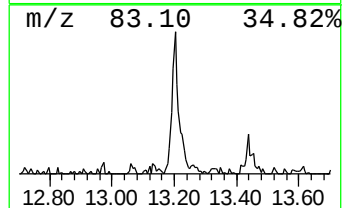
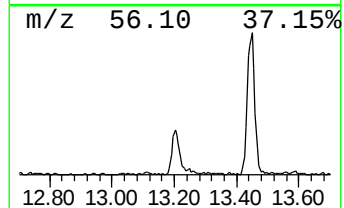
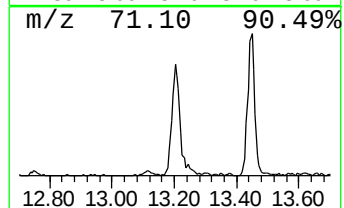
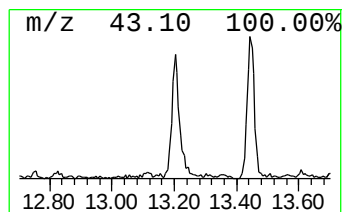
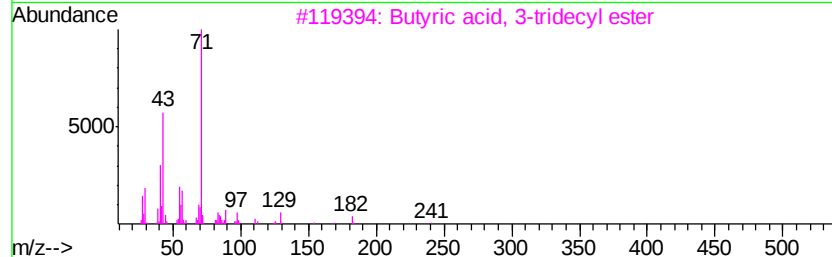
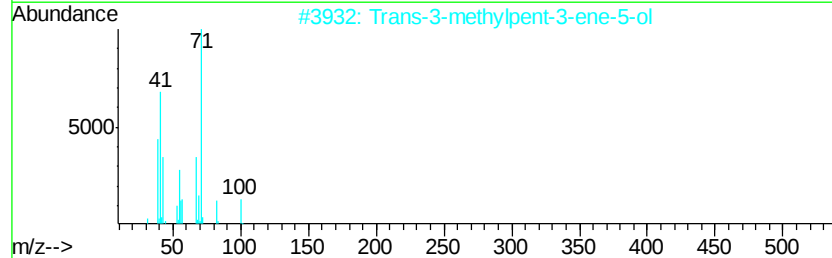
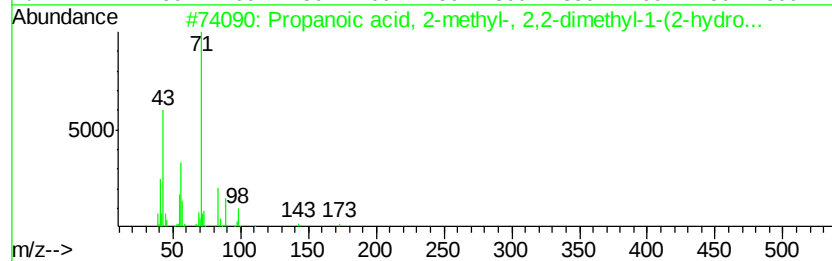
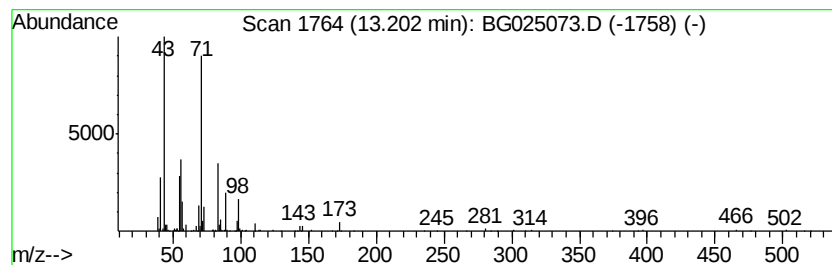
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	6.34 ng/ul	320422	Acenaphthene-d10	14.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Propanoic acid, 2-methyl-, 2,2-d...	216 C12H24O3	074367-33-2	72
2	Trans-3-methylpent-3-ene-5-ol	100 C6H12O	030801-95-7	43
3	Butyric acid, 3-tridecyl ester	270 C17H34O2	1000280-56-8	43
4	Sulfurous acid, 2-pentyl undecyl...	306 C16H34O3S	1000309-16-0	43
5	Propanoic acid, 2-methyl-, 2-met...	144 C8H16O2	000097-85-8	42



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
Operator : UM/SJ
Sample : H5861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

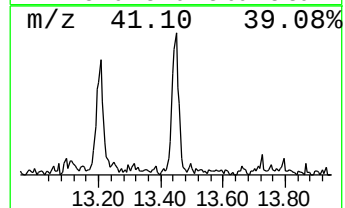
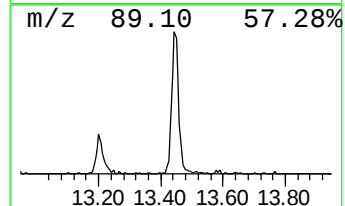
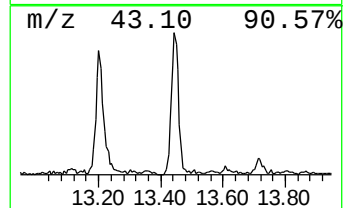
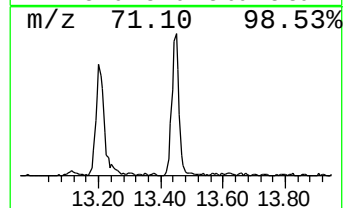
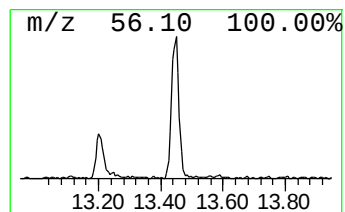
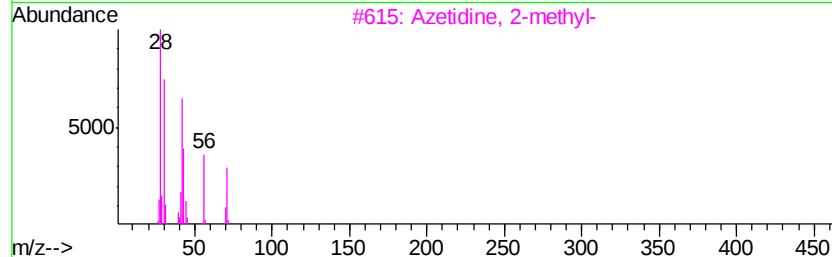
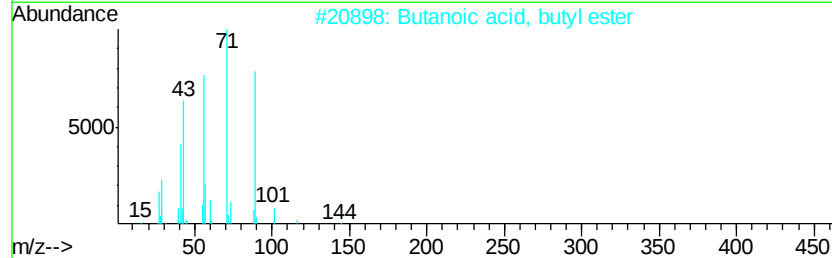
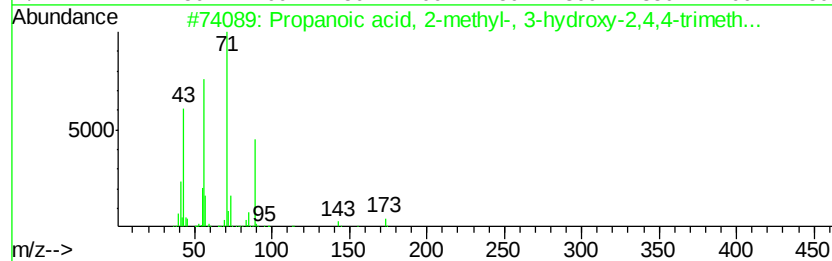
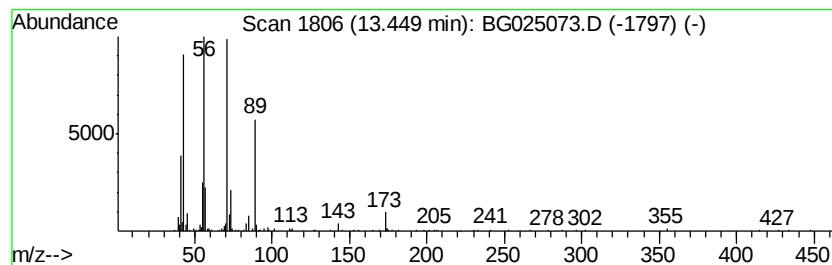
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Propanoic acid, 2-methyl-, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	7.56 ng/ul	382469	Acenaphthene-d10	14.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Propanoic acid, 2-methyl-, 3-hyd...	216 C12H24O3	074367-34-3	90
2	Butanoic acid, butyl ester	144 C8H16O2	000109-21-7	56
3	Azetidine, 2-methyl-	71 C4H9N	019812-49-8	50
4	Propanoic acid, 2-methyl-, butyl...	144 C8H16O2	000097-87-0	50
5	Butanoic acid, butyl ester	144 C8H16O2	000109-21-7	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
Operator : UM/SJ
Sample : H5861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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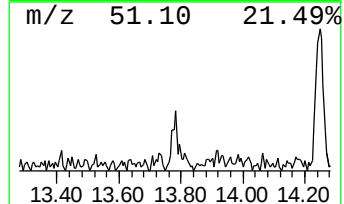
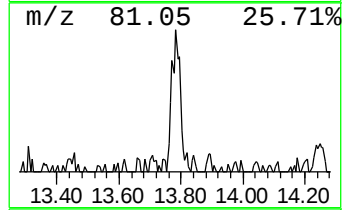
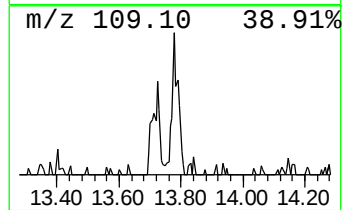
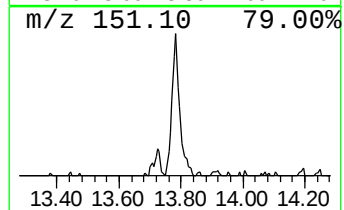
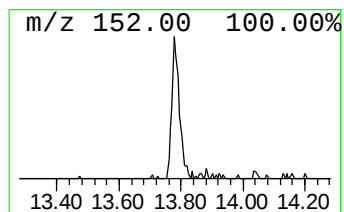
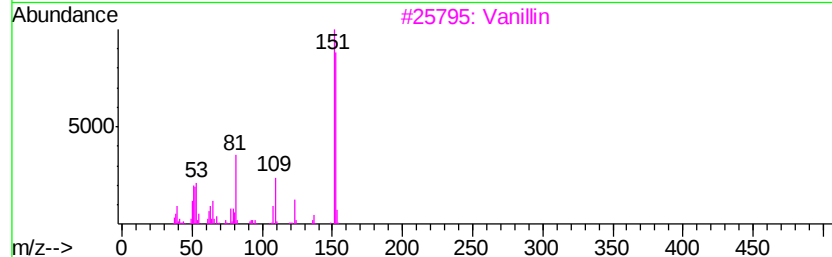
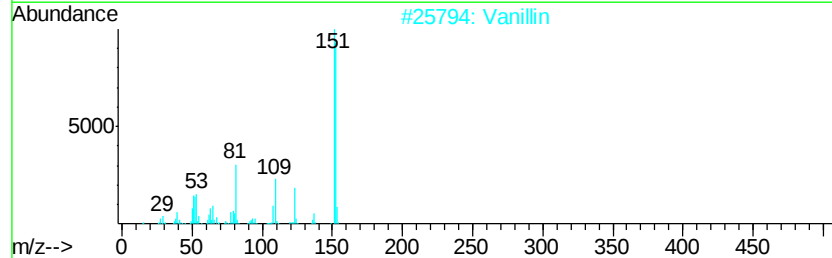
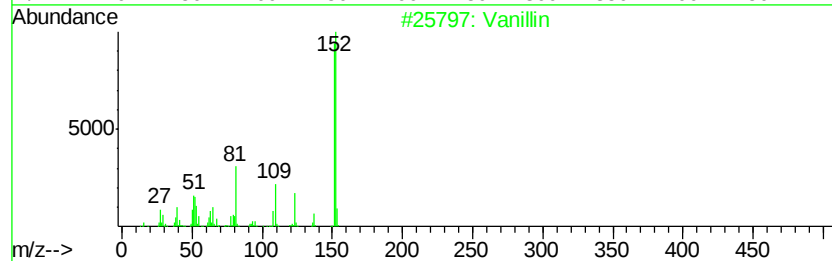
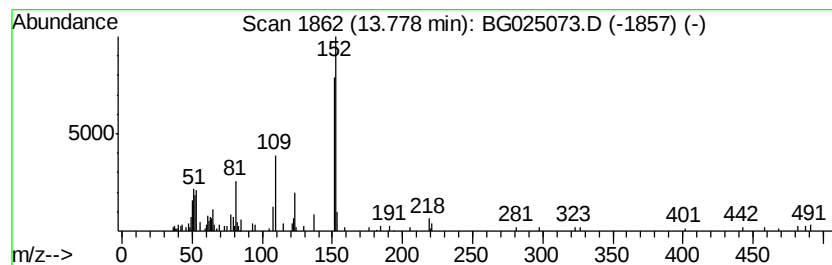
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Vanillin Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	2.35 ng/ul	118736	Acenaphthene-d10	14.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Vanillin	152	C8H8O3	000121-33-5	93
2		Vanillin	152	C8H8O3	000121-33-5	90
3		Vanillin	152	C8H8O3	000121-33-5	90
4		Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	64
5		Vanillin	152	C8H8O3	000121-33-5	60



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
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Misc :
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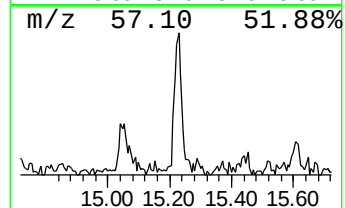
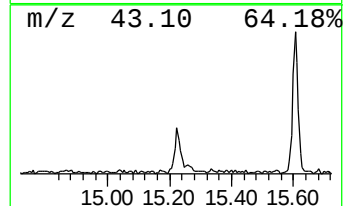
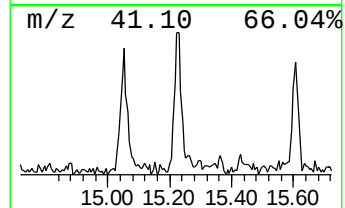
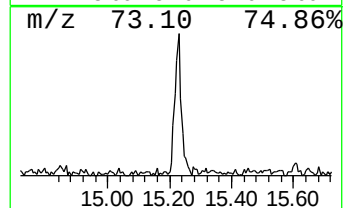
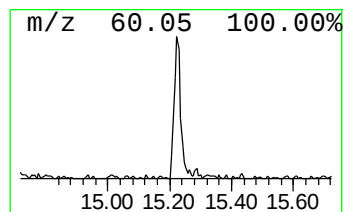
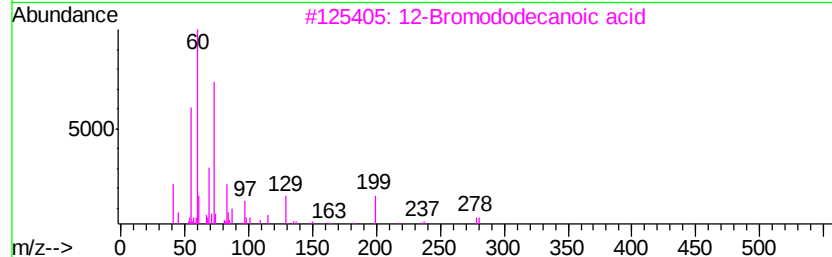
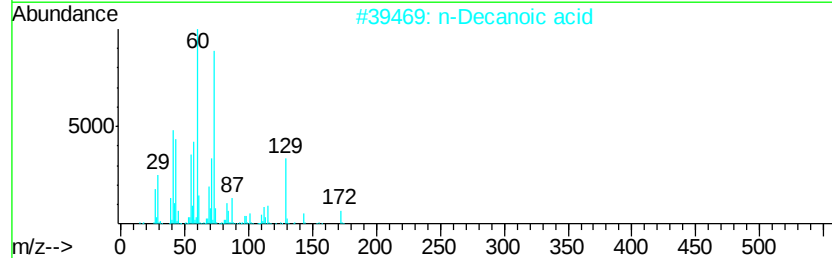
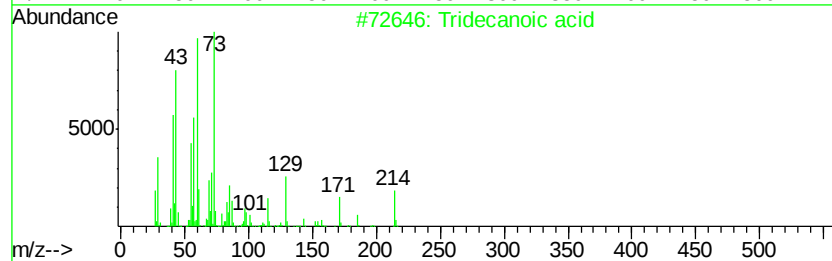
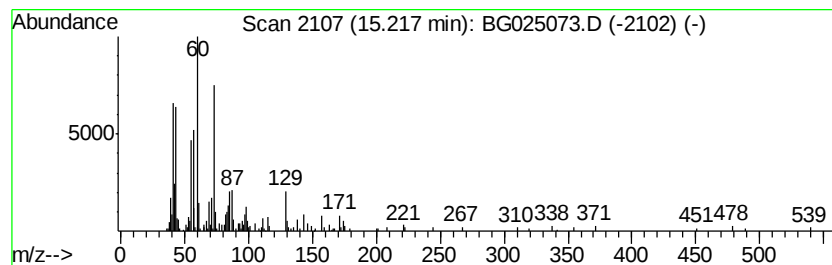
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	4.62 ng/ul	233515	Acenaphthene-d10	14.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	64
2		n-Decanoic acid	172	C10H20O2	000334-48-5	53
3		12-Bromododecanoic acid	278	C12H23BrO2	073367-80-3	50
4		Dodecanoic acid	200	C12H24O2	000143-07-7	49
5		Dodecanoic acid	200	C12H24O2	000143-07-7	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
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Misc :
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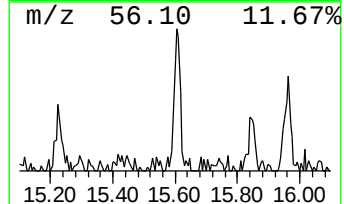
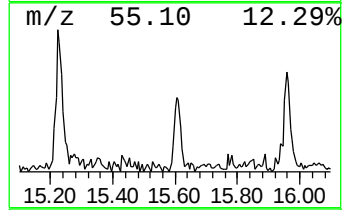
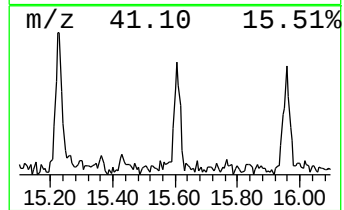
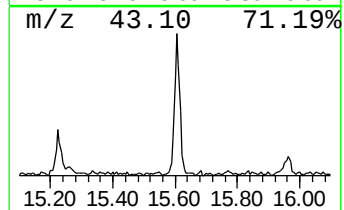
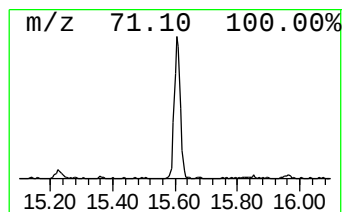
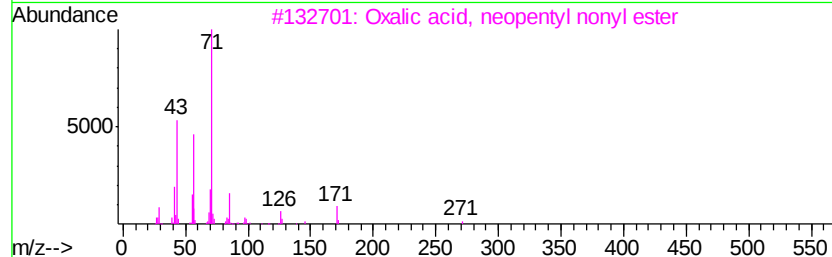
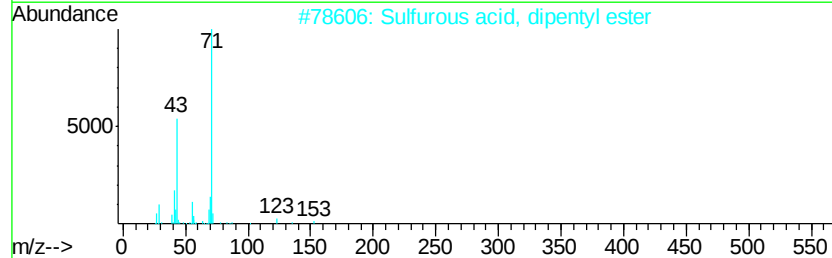
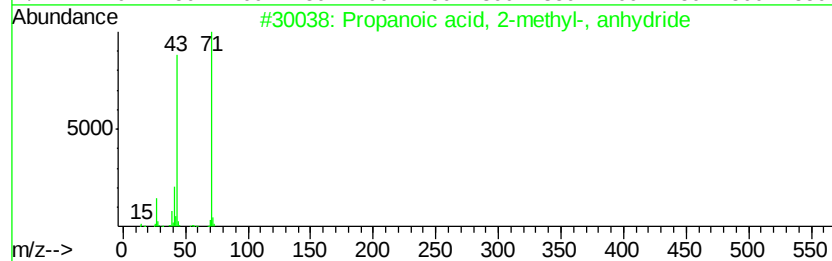
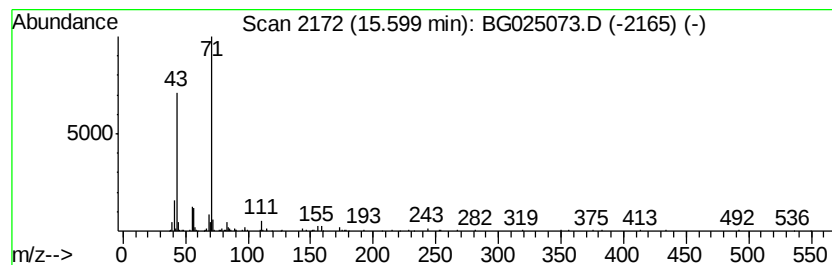
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Propanoic acid, 2-methyl-, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.60	4.58 ng/ul	231557	Acenaphthene-d10	14.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	53
2			Sulfurous acid, dipentyl ester	222	C10H22O3S	1000309-13-9	53
3			Oxalic acid, neopentyl nonyl ester	286	C16H30O4	1000309-73-3	45
4			Butane, 2-iodo-3-methyl-	198	C5H11I	018295-27-7	42
5			Propanoic acid, 2-methyl-, 1,2,3...	302	C15H26O6	014295-64-8	42



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025073.D
Acq On : 10 Dec 2016 17:58
Operator : UM/SJ
Sample : H5861-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DA8D8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.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
Heptanoic acid	7.19	4.8	ng/ul	116327	1	8.24	484224	20.0
5-(Methylamino)-1...	9.49	3.5	ng/ul	85451	1	8.24	484224	20.0
Octanoic acid	10.33	2.9	ng/ul	95355	2	11.06	660003	20.0
Acetaldehyde, di-...	10.80	2.5	ng/ul	84046	2	11.06	660003	20.0
Hexanoic acid	11.82	6.2	ng/ul	205665	2	11.06	660003	20.0
Propanoic acid, 2...	13.20	6.3	ng/ul	320422	3	14.86	1011320	20.0
Propanoic acid, 2...	13.45	7.6	ng/ul	382469	3	14.86	1011320	20.0
Vanillin	13.78	2.4	ng/ul	118736	3	14.86	1011320	20.0
Tridecanoic acid	15.22	4.6	ng/ul	233515	3	14.86	1011320	20.0
Propanoic acid, 2...	15.60	4.6	ng/ul	231557	3	14.86	1011320	20.0