LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.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	2.775	12	15	22	rVB2	31080	43967	1.37%	0.094%
2	2.987	45	51	59	rBV	440352	701384	21.85%	1.494%
3	3.058	59	63	79	rVB	1639619	1998076	62.24%	4.256%
4	3.316	103	107	116	rVB	42010	60173	1.87%	0.128%
5	3.452	127	130	135	rVB6	8829	9201	0.29%	0.020%
6	3.581	149	152	156	rBV	13936	21270	0.66%	0.045%
7	3.863	197	200	205	rBV6	8854	11372	0.35%	0.024%
8	4.046	228	231	236	rVB7	5062	7311	0.23%	0.016%
9	4.951	380	385	392	rBV	92300	140561	4.38%	0.299%
10	6.046	567	571	578	rBV5	5009	8603	0.27%	0.018%
11	6.810	696	701	702	rBV2	5022	7323	0.23%	0.016%
12	6.987	724	731	743	rVV	753193	1306886	40.71%	2.783%
13	7.157	750	760	769	rBV	537987	817119	25.45%	1.740%
14	7.357	787	794	801	rBV	735341	1166236	36.33%	2.484%
15	7.445	804	809	817	rVB2	34983	55781	1.74%	0.119%
16	7.734	852	858	862	rBV3	6052	10520	0.33%	0.022%
17	7.822	866	873	880	rVV	489592	798030	24.86%	1.700%
18	7.881	880	883	888	rVV3	9530	14361	0.45%	0.031%
19	8.192	931	936	942	rBV4	4956	8598	0.27%	0.018%
20	8.522	985	992	1001	rBV	880803	1390263	43.31%	2.961%
21	8.981	1063	1070	1082	rVV	725928	1176896	36.66%	2.507%
22	9.392	1135	1140	1146	rVB4	9437	16469	0.51%	0.035%
23	9.698	1183	1192	1199	rBV	618758	993016	30.93%	2.115%
24	10.233	1275	1283	1295	rBV	867087	1455613	45.34%	3.100%
25	10.616	1341	1348	1355	rBV	660949	1107702	34.50%	2.359%
26	10.751	1363	1371	1379	rBV	725514	1261564	39.30%	2.687%
27	11.928	1564	1571	1580	rVB	166596	265416	8.27%	0.565%
28	12.210	1612	1619	1623	rBV3	10714	17741	0.55%	0.038%
29	13.063	1759	1764	1769	rVB4	5972	9782	0.30%	0.021%
30	13.280	1796	1801	1806	rVB4	8089	14062	0.44%	0.030%
31	13.351	1806	1813	1821	rVB	80659	129789	4.04%	0.276%
32	13.869	1894	1901	1905	rBV	1455690	2045518	63.72%	4.357%
33	13.910	1905	1908	1914	rVV	592251	816803	25.44%	1.740%
34	14.098	1936	1940	1941	rBV4	5603	6992	0.22%	0.015%

LSC Area Percent Report

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Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Title : SVOA CALIBRATION

35	14.151	1943	1949	1959	rVV	1453341	2157488	67.21%	4.595%
36	14.363	1979	1985	1989	rVB5	7969	14924	0.46%	0.032%
37	14.463	1995	2002	2008	rBV	929387	1414325	44.06%	3.012%
38	14.651	2027	2034	2042	rBV	670490	1079863	33.64%	2.300%
39	14.733	2042	2048	2053	rVV2	114622	194185	6.05%	0.414%
40	14.780	2053	2056	2057	rVV3	6113	6527	0.20%	0.014%
41	14.880	2068	2073	2078	rBV7	18163	32755	1.02%	0.070%
42	15.098	2107	2110	2115	rVB4	9299	12816	0.40%	0.027%
43	15.221	2125	2131	2135	rBV	62426	80223	2.50%	0.171%
44	15.310	2143	2146	2150	rVV2	5667	7444	0.23%	0.016%
45	15.451	2162	2170	2177	rBV	1887607	2724554	84.87%	5.803%
46	15.563	2183	2189	2200	rBV	602910	853122	26.57%	1.817%
47	15.639	2200	2202	2208	rVV5	5061	7747	0.24%	0.016%
48	15.692	2208	2211	2218	rVB4	18110	28018	0.87%	0.060%
49	15.774	2218	2225	2227	rBV5	16947	25866	0.81%	0.055%
50	15.815	2227	2232	2242	rVV	1003843	1372631	42.76%	2.923%
51	15.992	2258	2262	2266	rBV3	18730	26900	0.84%	0.057%
52	16.027	2266	2268	2276	rVB4	12698	21268	0.66%	0.045%
53	16.110	2279	2282	2288	rBV6	4060	6956	0.22%	0.015%
54	16.168	2288	2292	2299	rBV6	5573	13714	0.43%	0.029%
55	16.339	2317	2321	2326	rBV3	13902	19772	0.62%	0.042%
56	16.439	2333	2338	2343	rVB	52741	70470	2.20%	0.150%
57	16.527	2344	2353	2358	rBV	244052	398733	12.42%	0.849%
58	16.586	2358	2363	2364	rVV2	40256	69857	2.18%	0.149%
59	16.604	2364	2366	2370	rVV	53373	68557	2.14%	0.146%
60	16.657	2374	2375	2381	rVB4	11063	12065	0.38%	0.026%
61	16.815	2397	2402	2408	rBV4	18881	32688	1.02%	0.070%
62	17.045	2437	2441	2446	rVB	68098	84929	2.65%	0.181%
63	17.157	2453	2460	2463	rBV	117916	174305	5.43%	0.371%
64	17.204	2463	2468	2477	rVB	1156818	1673484	52.13%	3.564%
65	17.298	2478	2484	2493	rVB	1984738	2906995	90.55%	6.191%
66	17.386	2493	2499	2500	rBV5	6347	6915	0.22%	0.015%
67	17.462	2505	2512	2519	rBV	152947	268523	8.36%	0.572%
68	17.580	2528	2532	2535	rBV4	17675	26564	0.83%	0.057%
69	17.886	2577	2584	2590	rBV3	18761	30711	0.96%	0.065%
70	18.092	2614	2619	2628	rBV2	390635	587642	18.31%	1.252%
71	18.315	2655	2657	2662	rVB5	4382	7121	0.22%	0.015%

LSC Area Percent Report

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Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Title : SVOA CALIBRATION

72	18.386	2665	2669	2673	rVV5	13528	19874	0.62%	0.042%
73	18.521	2688	2692	2695	rBV3	12129	14888	0.46%	0.032%
74	18.586	2698	2703	2706	rVB6	8457	16463	0.51%	0.035%
75	18.627	2706	2710	2713	rBV5	18604	28445	0.89%	0.061%
76	18.968	2764	2768	2770	rBV5	6853	9801	0.31%	0.021%
77	19.021	2774	2777	2781	rBV5	11049	17102	0.53%	0.036%
78	19.180	2802	2804	2808	rBV5	7937	9816	0.31%	0.021%
79	19.227	2808	2812	2818	rBV5	44611	95448	2.97%	0.203%
80	19.374	2832	2837	2848	rBV	255374	428430	13.35%	0.912%
81	19.480	2853	2855	2859	rVB4	21628	26962	0.84%	0.057%
82	19.598	2869	2875	2879	rBV	2262531	3103596	96.68%	6.610%
83	19.639	2879	2882	2890	rVB	144650	228582	7.12%	0.487%
84	19.798	2907	2909	2918	rVB9	27119	47485	1.48%	0.101%
85	19.909	2924	2928	2931	rBV6	15264	26269	0.82%	0.056%
86	20.133	2963	2966	2971	rVB5	28039	35633	1.11%	0.076%
87	20.633	3046	3051	3054	rBV5	34513	51958	1.62%	0.111%
88	20.745	3068	3070	3074	rBV3	18316	18680	0.58%	0.040%
89	21.092	3124	3129	3140	rBV	511235	744624	23.20%	1.586%
90	21.297	3160	3164	3171	rVB	230301	259322	8.08%	0.552%
91	21.386	3174	3179	3185	rVV	1453391	1864829	58.09%	3.972%
92	21.497	3195	3198	3201	rBV5	68502	79703	2.48%	0.170%
93	21.533	3201	3204	3208	rVB5	28404	36905	1.15%	0.079%
94	22.021	3284	3287	3291	rVB5	29432	36041	1.12%	0.077%
95	22.444	3355	3359	3363	rVB4	49817	70056	2.18%	0.149%
96	22.956	3444	3446	3452	rVB4	27858	34821	1.08%	0.074%
97	23.533	3536	3544	3555	rVB2	1670212	3210274	100.00%	6.837%
98	23.674	3562	3568	3578	rVB	938376	1809985	56.38%	3.855%
99	24.297	3669	3674	3685	rVB9	68387	168036	5.23%	0.358%
100	24.544	3713	3716	3718	rBV4	30017	42598	1.33%	0.091%

Sum of corrected areas: 46951681

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
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Sample : G1065-14
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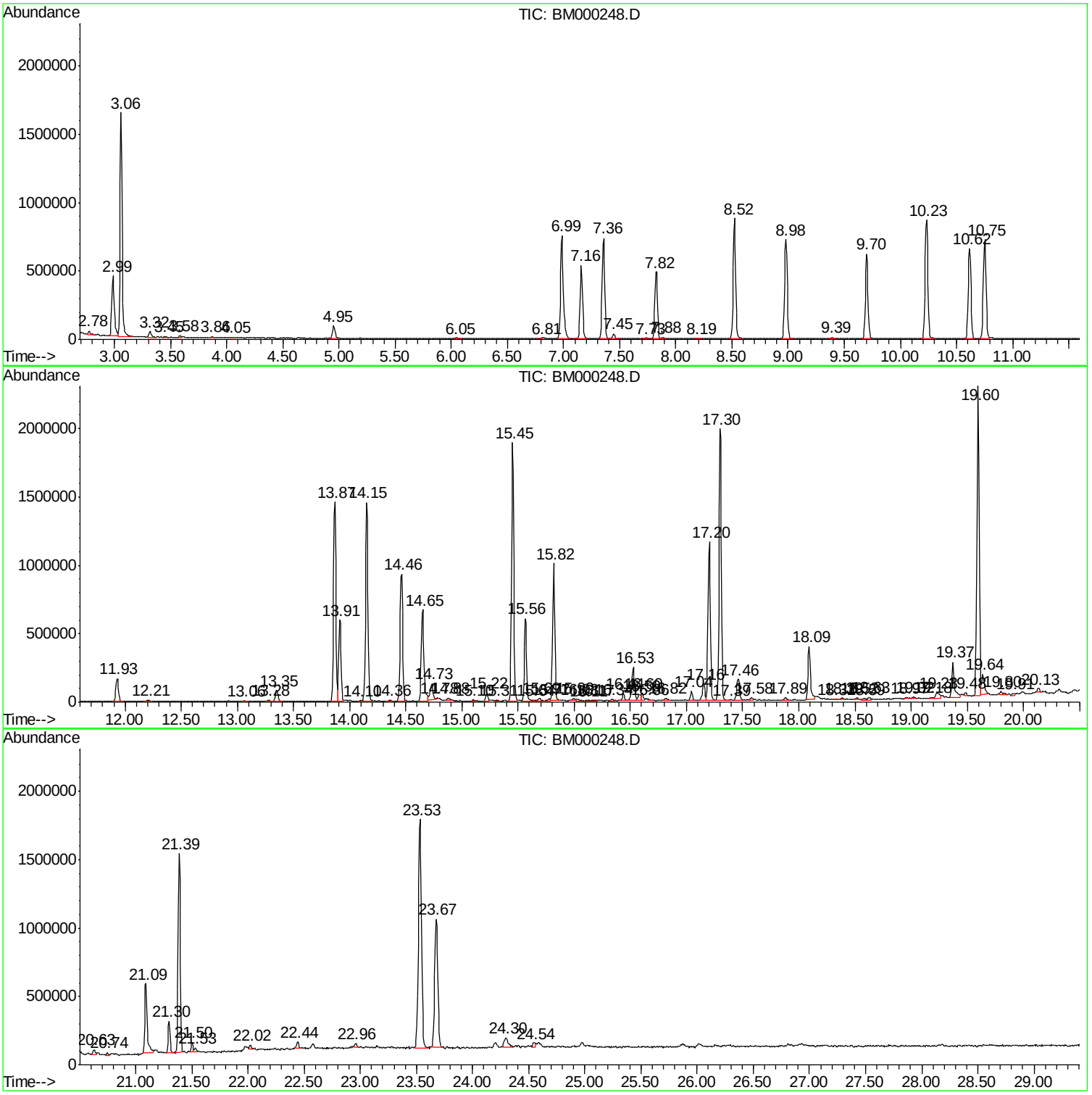
Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
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Operator : TP/IZ
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

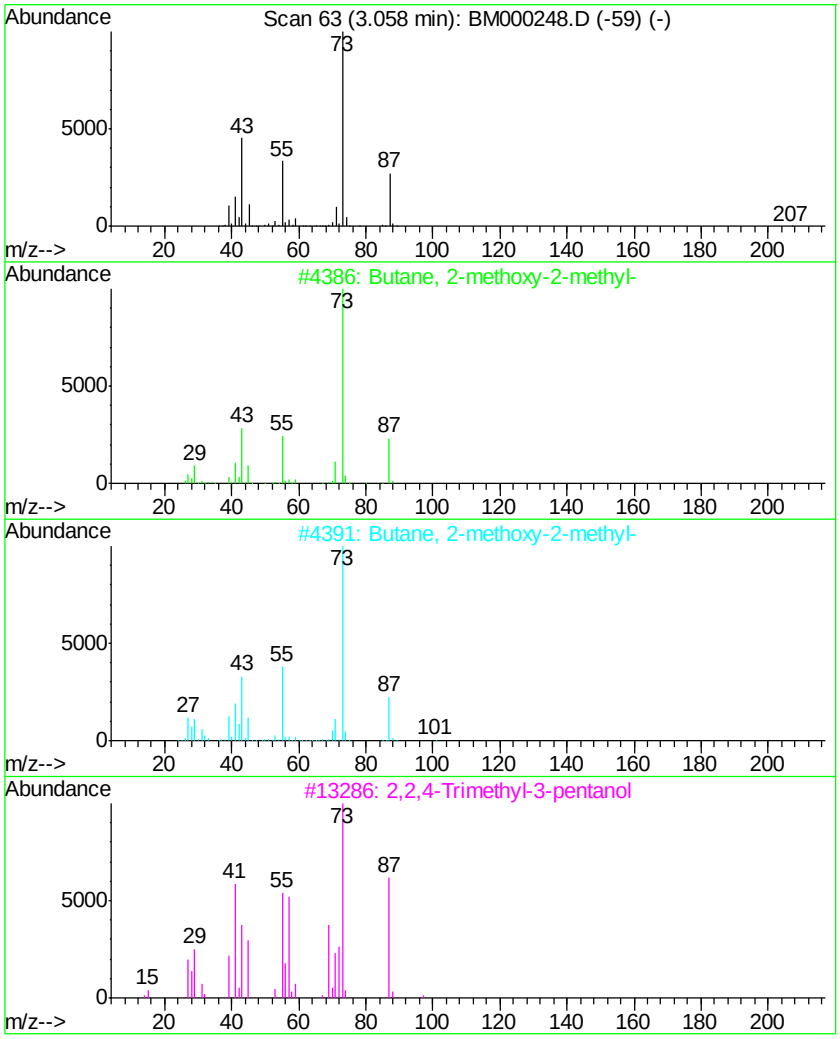
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

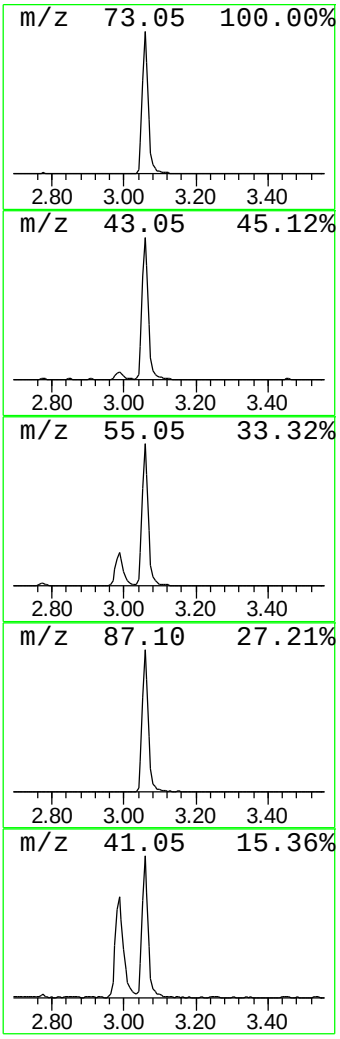
****

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	50.08 ng/ul	1998080	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

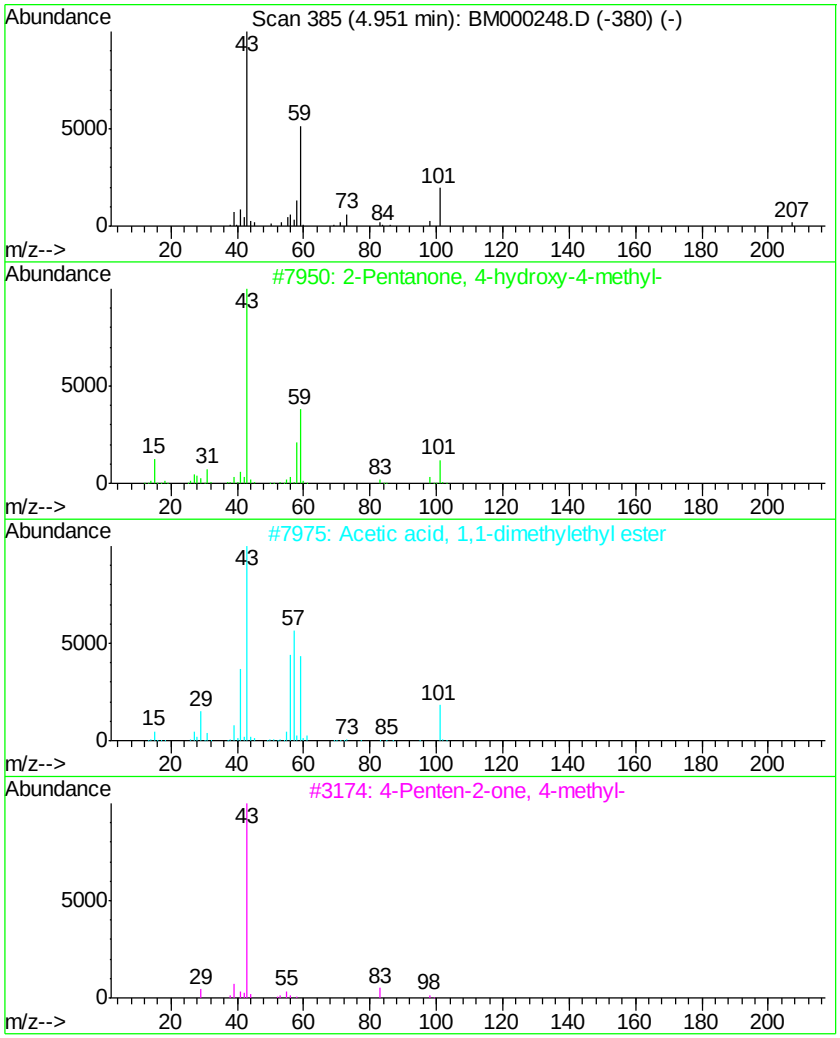
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

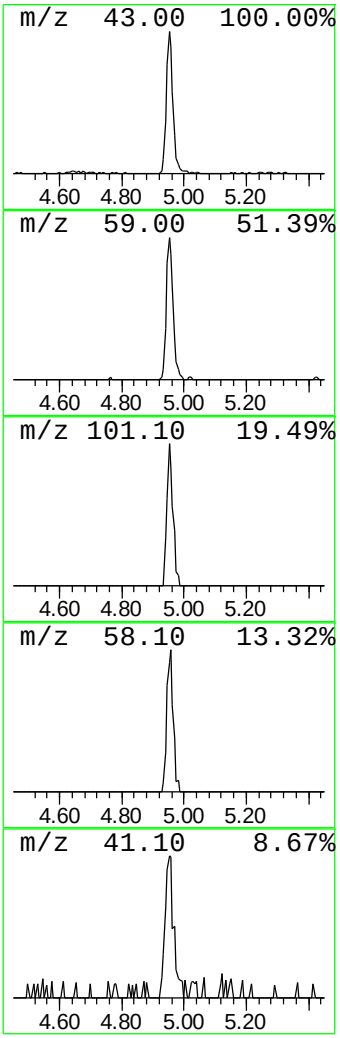
****

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.95	3.52 ng/ul	140561	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	10
4	2-Propyn-1-ol, acetate	98	C5H6O2	000627-09-8	9
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

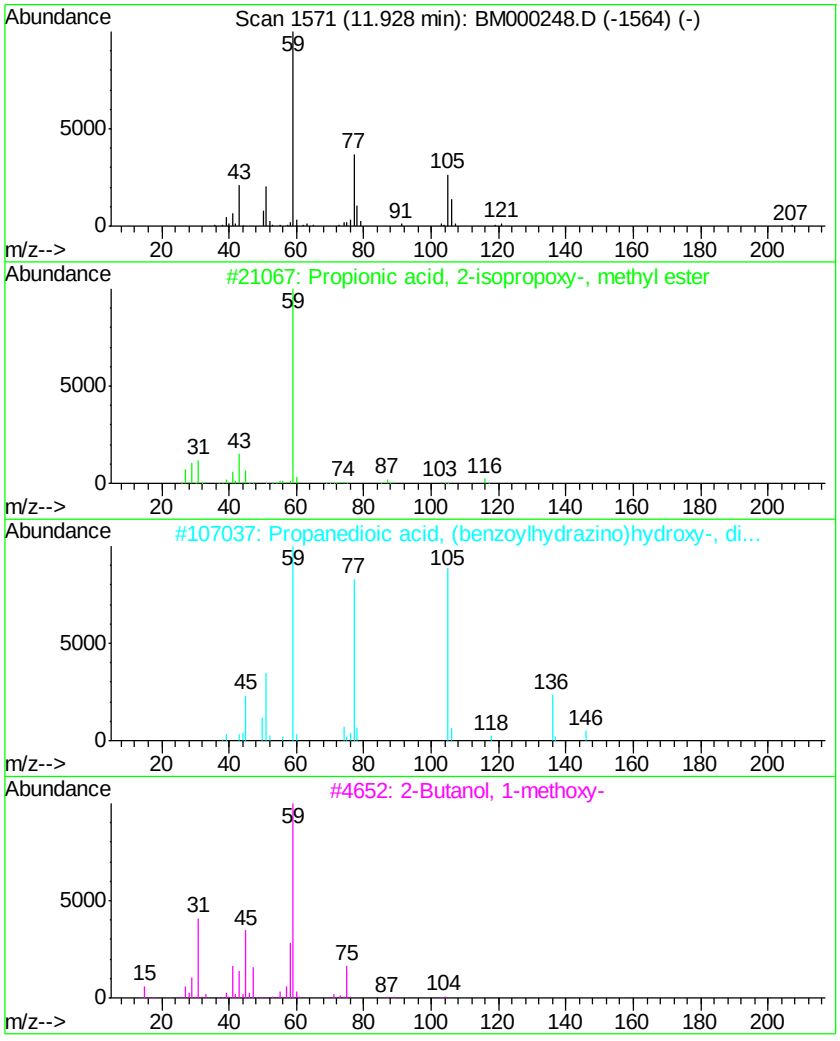
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

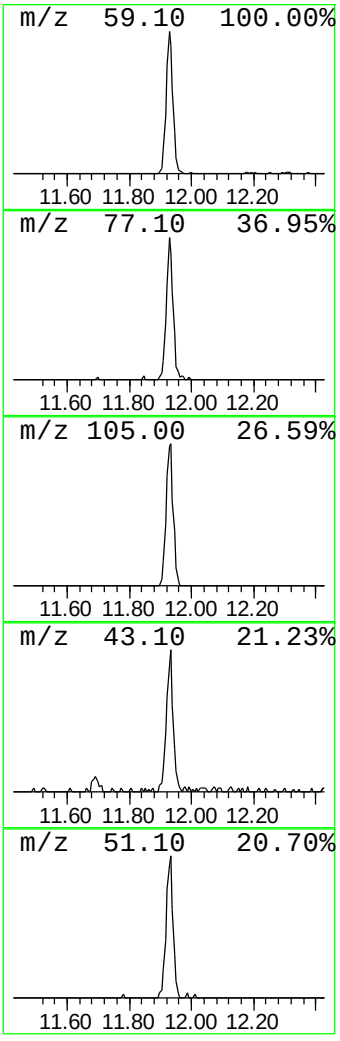
****

Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	4.79 ng/ul	265416	Naphthalene-d8	10.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propionic acid, 2-isopropoxy-, m...	146	C7H14O3	1000151-15-1	23
2	Propanedioic acid, (benzoylhydra...	282	C12H14N2O6	1000142-99-9	23
3	2-Butanol, 1-methoxy-	104	C5H12O2	053778-73-7	9
4	Guanidine	59	CH5N3	000113-00-8	9
5	Ethanol, 1-methoxy-, benzoate	180	C10H12O3	051835-44-0	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

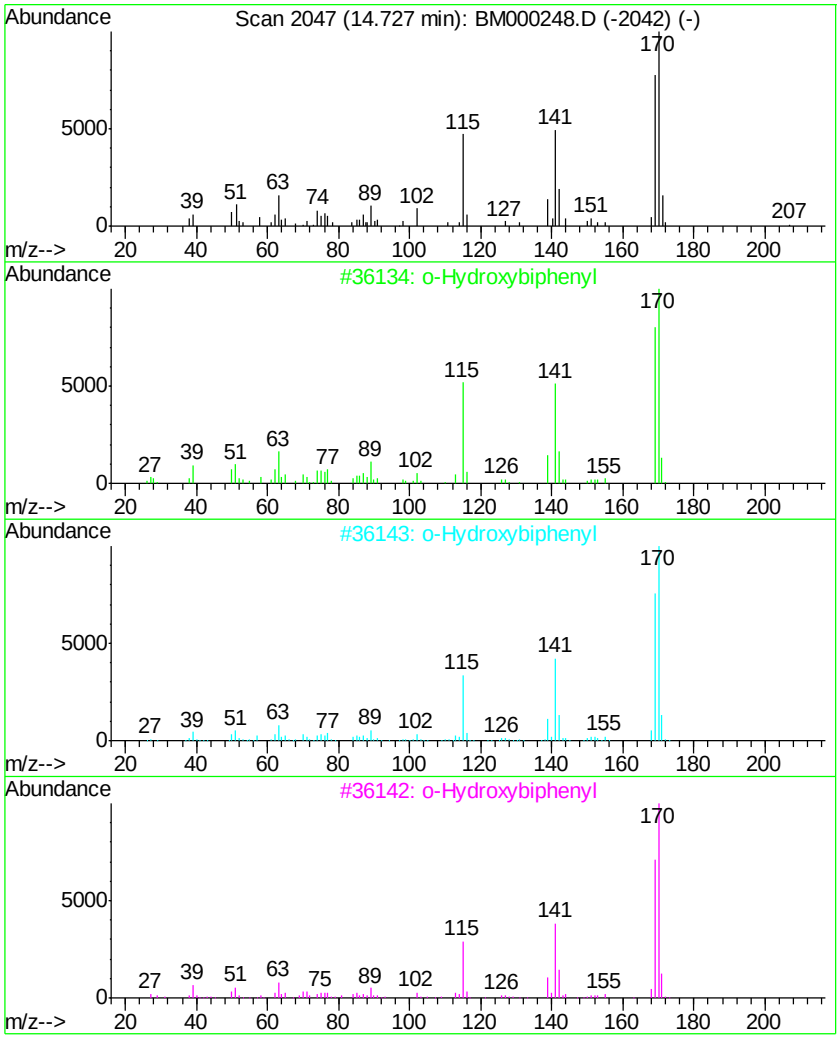
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

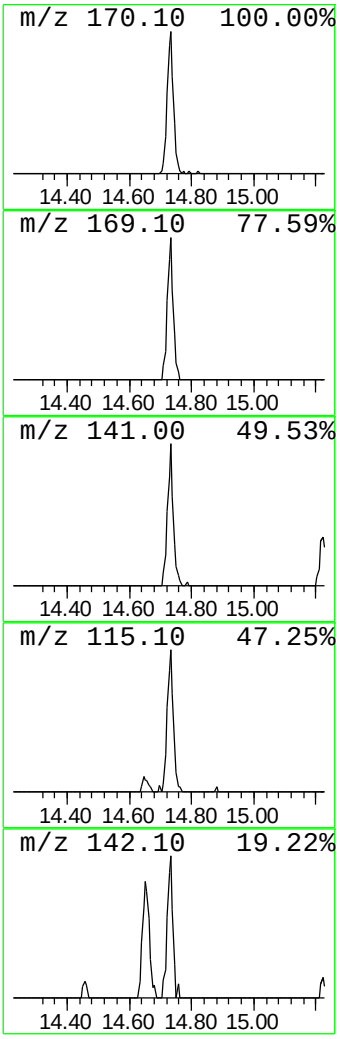
****

Peak Number 5 o-Hydroxybiphenyl Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	2.75 ng/ul	194185	Acenaphthene-d10	14.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
2	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	94
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	91
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	91
5	1-Naphthalenecarboxaldehyde, 4-m...	170	C12H10O	033738-48-6	87





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

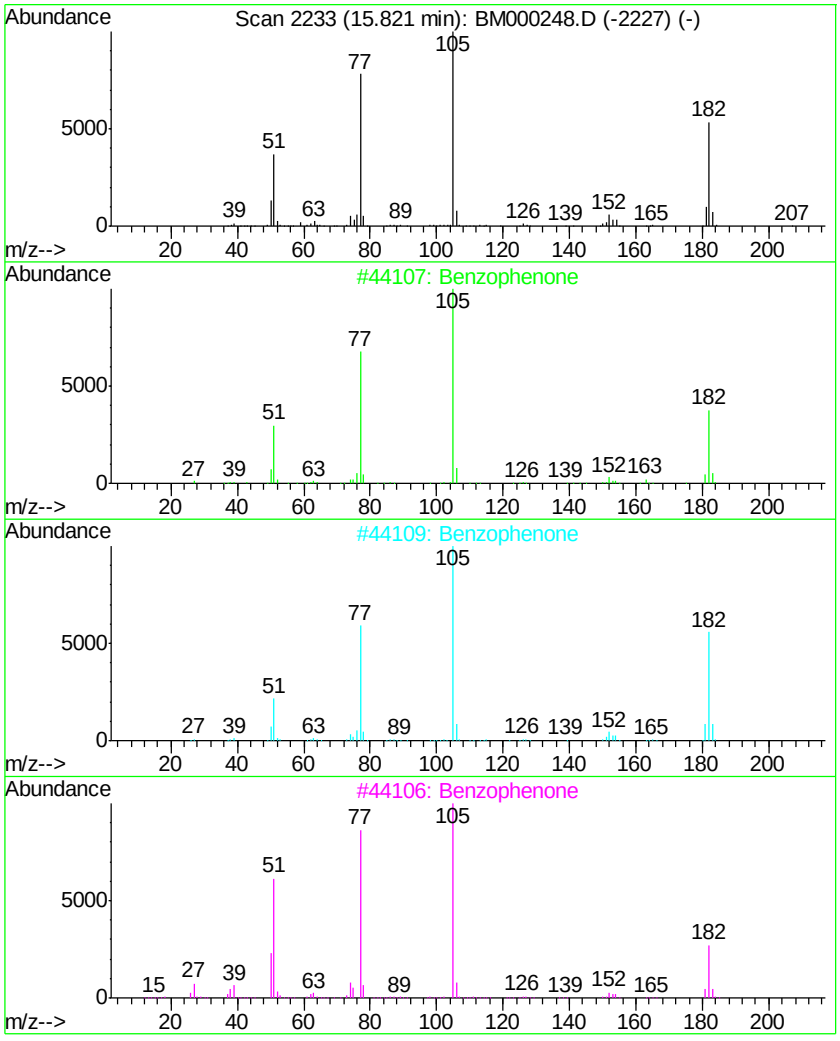
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

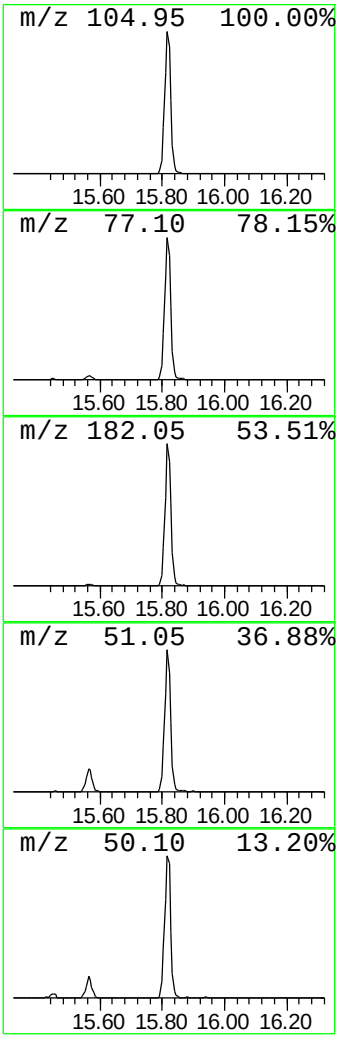
****

Peak Number 6 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	19.41 ng/ul	1372630	Acenaphthene-d10	14.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Benzophenone	182	C13H10O	000119-61-9	97
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	95
5	Benzophenone	182	C13H10O	000119-61-9	91





Data Path : Z:\HPCHEM1\BNA M\DATA\BM011615\
Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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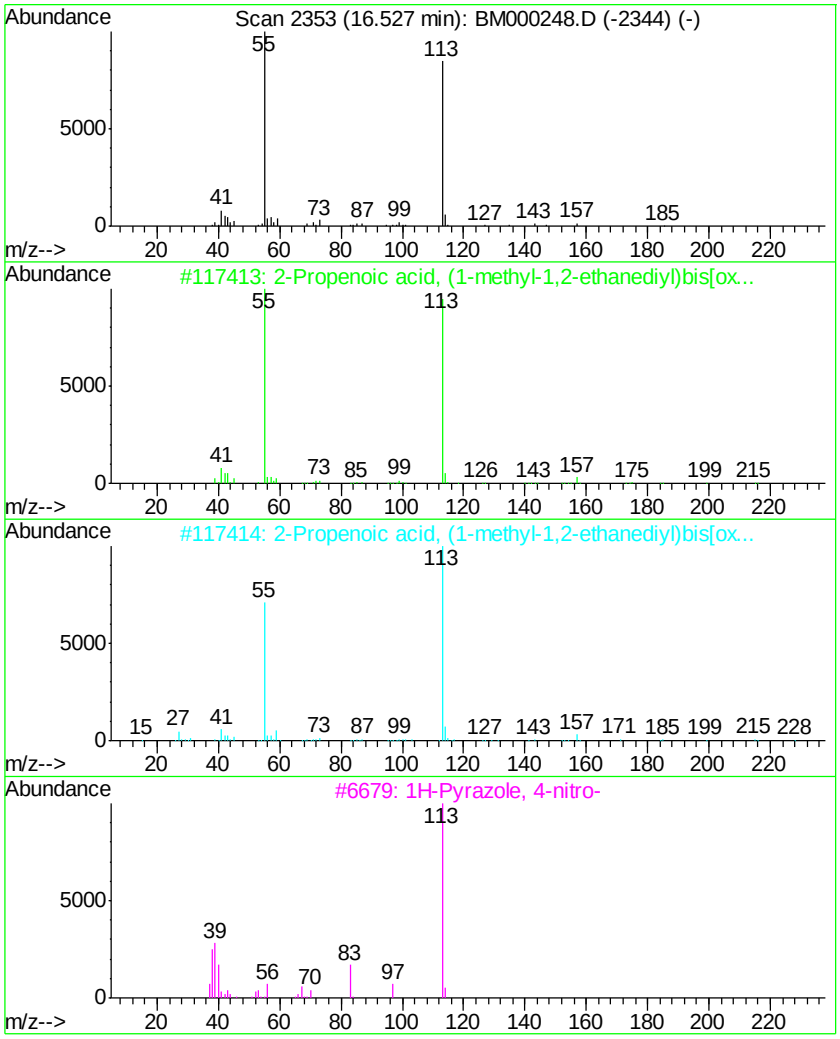
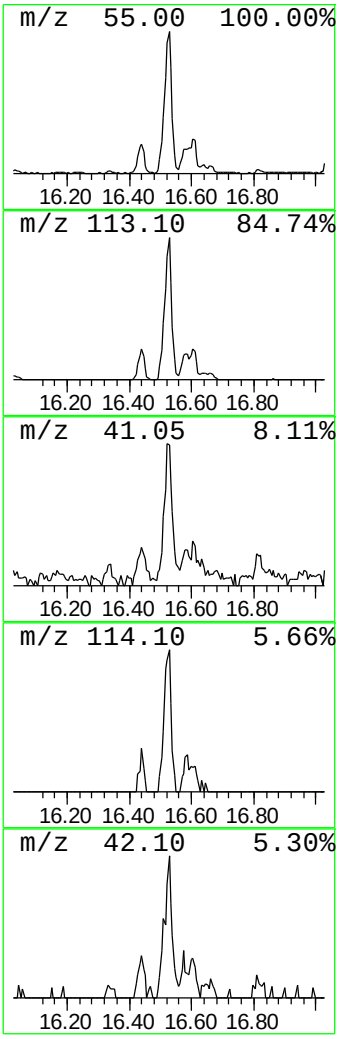
Peak Number 7 2-Propenoic acid, (1-methyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	4.77 ng/ul	398733	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	78
2	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	64
3	1H-Pyrazole, 4-nitro-	113	C3H3N3O2	002075-46-9	40
4	Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	36
5	2-Propenoic acid, 2-methyl-, 1,2...	286	C14H22O6	000109-16-0	36





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Data File : BM000248.D
Acq On : 17 Jan 2015 00:31
Operator : TP/IZ
Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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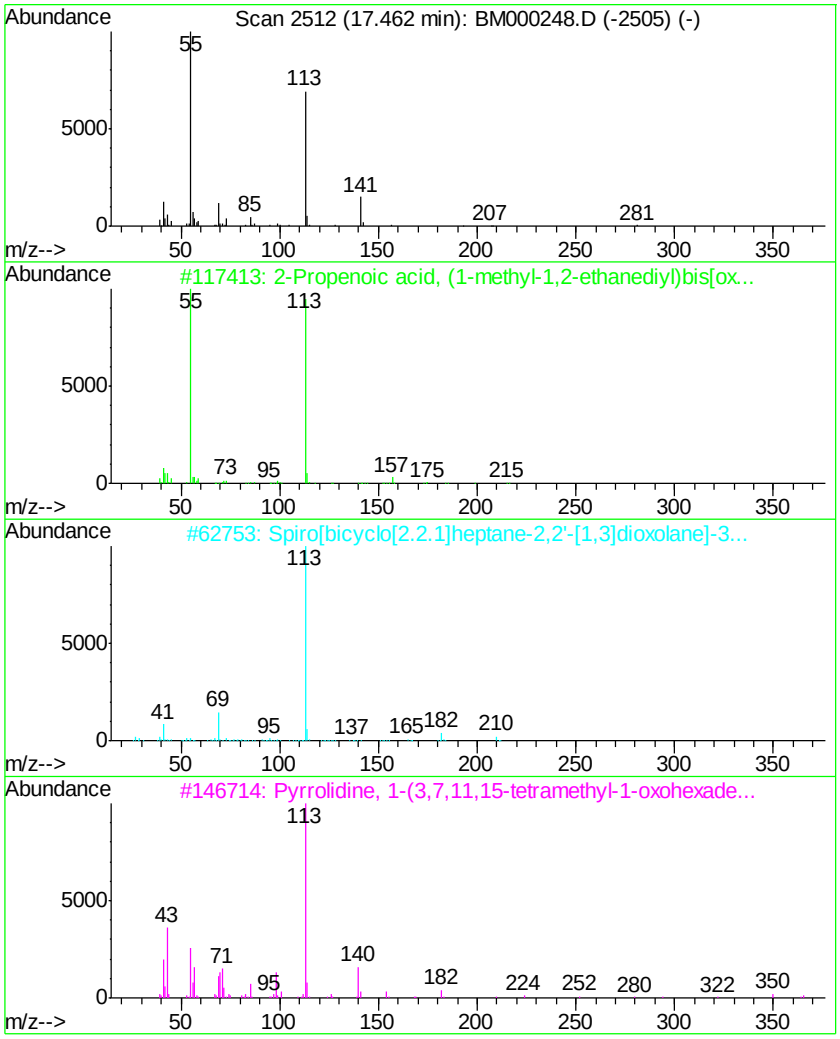
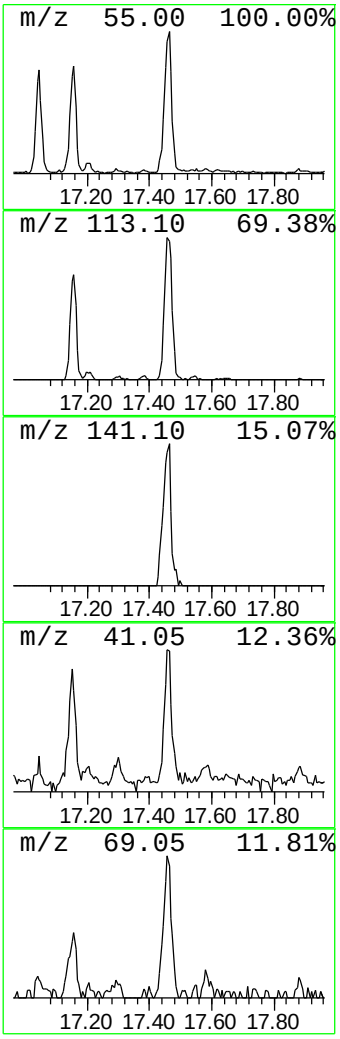
Peak Number 8 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.46	3.21 ng/ul	268523	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	40
2	Spiro[bicyclo[2.2.1]heptane-2,2'...	210	C12H18O3	018501-56-9	37
3	Pyrrolidine, 1-(3,7,11,15-tetram...	365	C24H47NO	056630-63-8	33
4	(1-Propoxy-pentyl)-cyclopropane	170	C11H22O	094883-97-3	28
5	Pyrrolidine, 1-(3-methyl-1-oxohe...	337	C22H43NO	1000036-69-5	28





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Instrument :
ClientSampleId :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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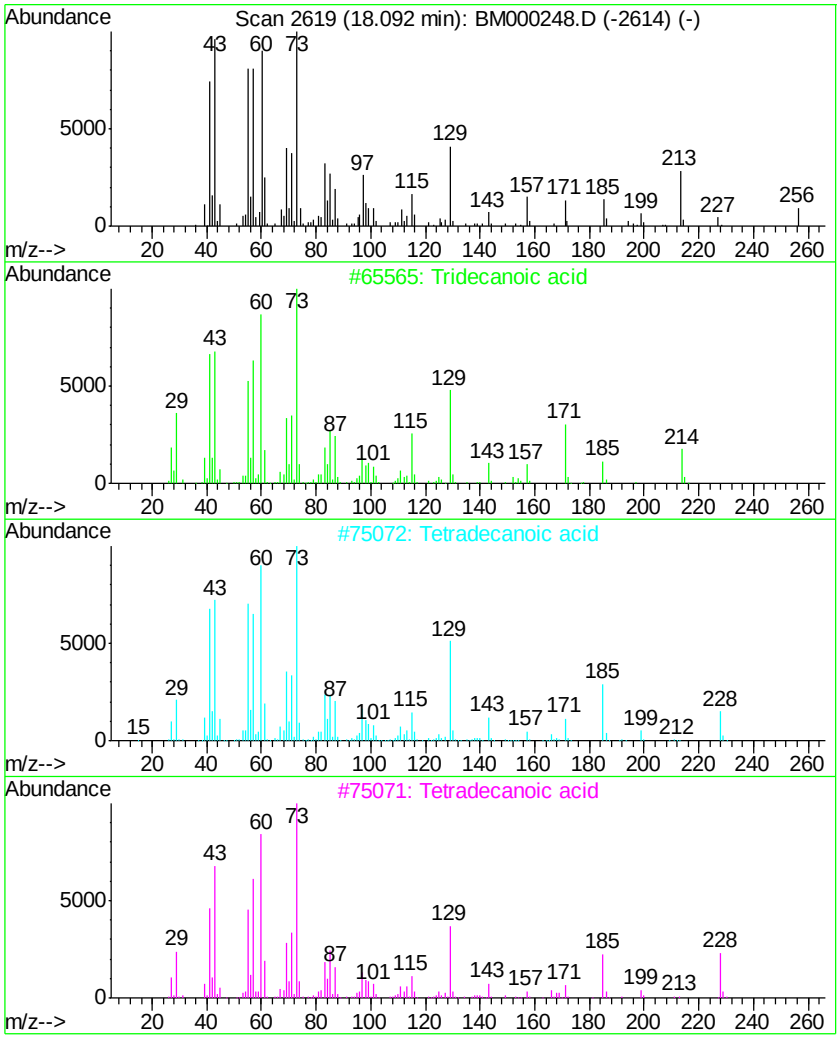
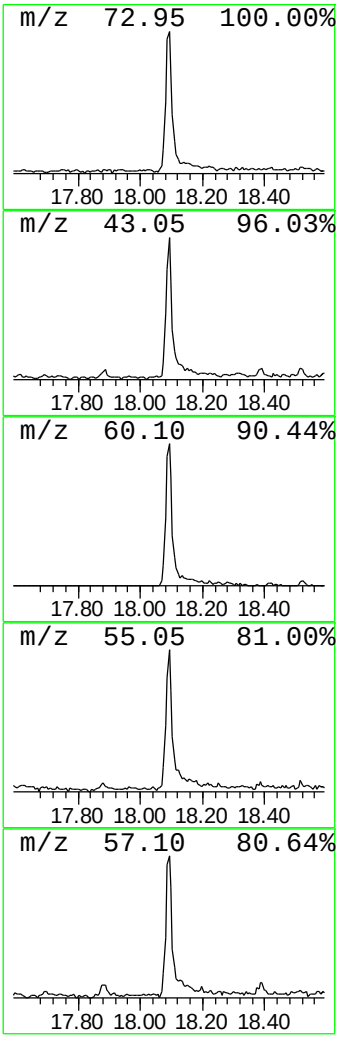
Peak Number 9 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	7.02 ng/ul	587642	Phenanthrene-d10	17.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72





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Sample : G1065-14
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
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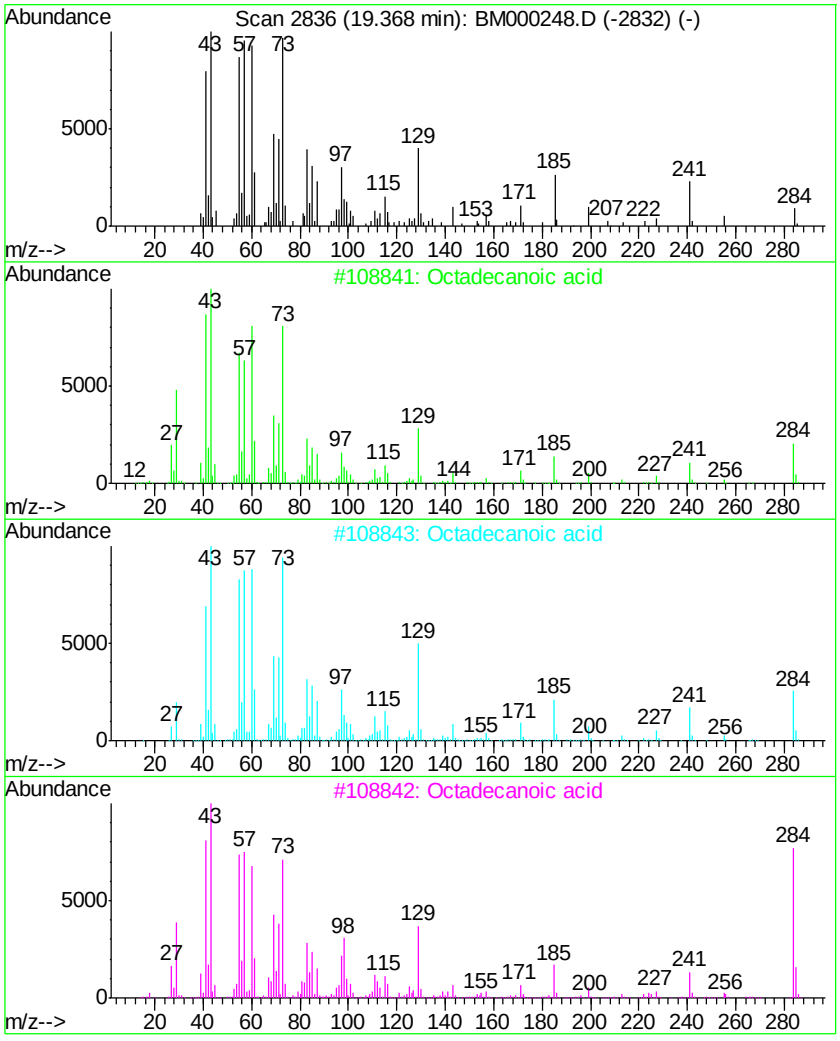
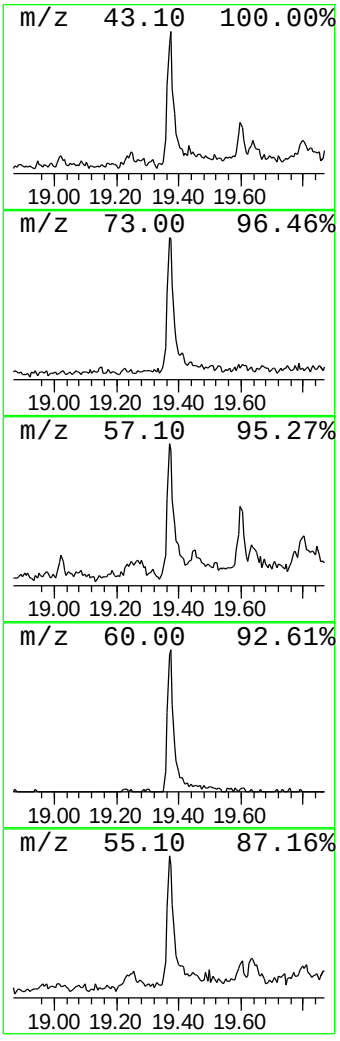
Peak Number 10 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.37	4.59 ng/ul	428430	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	96
2	Octadecanoic acid	284	C18H36O2	000057-11-4	90
3	Octadecanoic acid	284	C18H36O2	000057-11-4	87
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5	Octadecanoic acid	284	C18H36O2	000057-11-4	68





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Sample : G1065-14
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
ClientSampleId :
C0AG1

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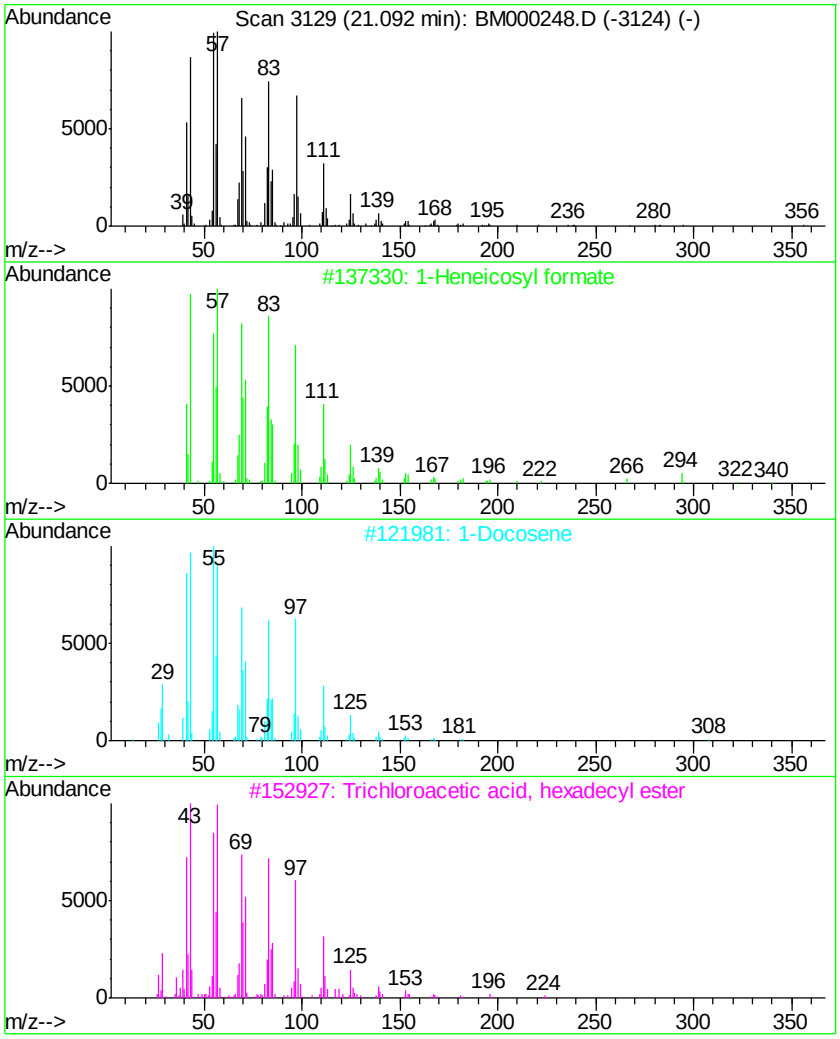
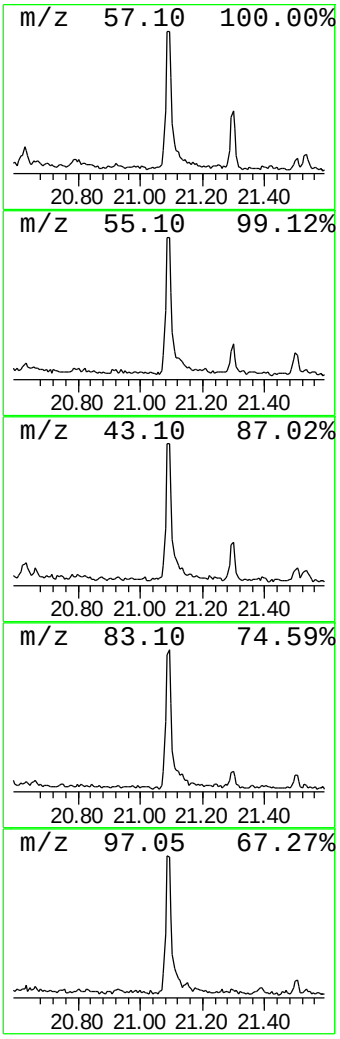
Peak Number 11 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.09	7.99 ng/ul	744624	Chrysene-d12	21.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2	1-Docosene	308	C22H44	001599-67-3	94
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
5	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91





Tentatively Identified Compound (LSC) summary

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	3.06	50.1	ng/ul	1998080	1	7.82	798030	20.0
2-Pentanone, 4-hy...	4.95	3.5	ng/ul	140561	1	7.82	798030	20.0
unknown-01	11.93	4.8	ng/ul	265416	2	10.62	1107700	20.0
o-Hydroxybiphenyl	14.73	2.8	ng/ul	194185	3	14.46	1414330	20.0
Benzophenone	15.82	19.4	ng/ul	1372630	3	14.46	1414330	20.0
2-Propenoic acid,...	16.53	4.8	ng/ul	398733	4	17.20	1673480	20.0
unknown-02	17.46	3.2	ng/ul	268523	4	17.20	1673480	20.0
Tridecanoic acid	18.09	7.0	ng/ul	587642	4	17.20	1673480	20.0
Octadecanoic acid	19.37	4.6	ng/ul	428430	5	21.39	1864830	20.0
1-Heneicosyl formate	21.09	8.0	ng/ul	744624	5	21.39	1864830	20.0