LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

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_G\METHODS\SOM01.2-EPA-BG021015.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	2.812	17	21	34	rVB	669534	814867	14.39%	1.425%
2	2.953	40	45	54	rBV	203917	343306	6.06%	0.600%
3	3.030	54	58	74	rVB	107411	152535	2.69%	0.267%
4	4.340	276	281	289	rBV	127867	165069	2.91%	0.289%
5	4.933	377	382	391	rBV	57121	78731	1.39%	0.138%
6	6.954	719	726	741	rBV2	755609	1417158	25.03%	2.478%
7	7.148	751	759	767	rBV	203854	324626	5.73%	0.568%
8	7.342	785	792	794	rBV	251503	394261	6.96%	0.689%
9	7.371	794	797	804	rVV	496085	763123	13.48%	1.334%
10	7.441	804	809	815	rVB	16370	21395	0.38%	0.037%
11	7.812	866	872	878	rVB	204200	308282	5.44%	0.539%
12	8.511	984	991	995	rBV	314613	502867	8.88%	0.879%
13	8.563	995	1000	1006	rVV	521796	814649	14.39%	1.425%
14	8.628	1006	1011	1019	rVB	544854	865503	15.28%	1.513%
15	8.892	1048	1056	1062	rBV	304434	488493	8.63%	0.854%
16	8.963	1062	1068	1075	rVV	222864	370271	6.54%	0.647%
17	9.639	1176	1183	1184	rBV4	5813	10169	0.18%	0.018%
18	9.685	1184	1191	1199	rVB	168559	280327	4.95%	0.490%
19	10.244	1275	1286	1295	rBV2	674892	1513201	26.72%	2.646%
20	10.602	1336	1347	1351	rBV	315276	600403	10.60%	1.050%
21	10.655	1351	1356	1363	rVV	829578	1369648	24.19%	2.395%
22	10.737	1363	1370	1382	rVB	281253	482025	8.51%	0.843%
23	11.495	1492	1499	1501	rBV3	35705	66929	1.18%	0.117%
24	11.918	1564	1571	1578	rVB	40886	64985	1.15%	0.114%
25	12.129	1600	1607	1614	rVV3	14546	25469	0.45%	0.045%
26	12.188	1614	1617	1622	rVV4	5379	9295	0.16%	0.016%
27	12.264	1622	1630	1640	rVB	1081966	1641333	28.98%	2.870%
28	12.482	1663	1667	1675	rVB	10723	15788	0.28%	0.028%
29	12.711	1701	1706	1712	rVB5	6410	10629	0.19%	0.019%
30	12.916	1736	1741	1746	rBV3	15133	22127	0.39%	0.039%
31	13.187	1783	1787	1791	rBV3	8878	12661	0.22%	0.022%
32	13.281	1795	1803	1805	rBV	95008	149881	2.65%	0.262%
33	13.322	1805	1810	1817	rVB	1042474	1595877	28.18%	2.791%
34	13.856	1895	1901	1905	rBV	579735	757812	13.38%	1.325%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

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_G\METHODS\SOM01.2-EPA-BG021015.M
Title : SVOA CALIBRATION

35	13.897	1905	1908	1914	rVV	75385	103511	1.83%	0.181%
36	14.132	1936	1948	1955	rBV	638925	967368	17.08%	1.692%
37	14.356	1979	1986	1990	rBV2	10475	15949	0.28%	0.028%
38	14.444	1995	2001	2007	rVV2	484338	723110	12.77%	1.264%
39	14.508	2007	2012	2020	rVB	835455	1196500	21.13%	2.092%
40	14.626	2026	2032	2041	rBV	312499	452141	7.98%	0.791%
41	14.708	2043	2046	2052	rVB5	8130	13435	0.24%	0.023%
42	14.849	2065	2070	2075	rBV4	13005	18746	0.33%	0.033%
43	15.213	2128	2132	2143	rBV	52352	82082	1.45%	0.144%
44	15.301	2143	2147	2152	rVV	51384	62823	1.11%	0.110%
45	15.437	2163	2170	2174	rBV	881726	1255754	22.18%	2.196%
46	15.489	2174	2179	2184	rVV	953072	1322941	23.36%	2.313%
47	15.560	2184	2191	2204	rVB3	640590	1031110	18.21%	1.803%
48	15.677	2204	2211	2214	rBV3	9653	13766	0.24%	0.024%
49	15.801	2225	2232	2238	rVB	281048	372603	6.58%	0.652%
50	16.324	2317	2321	2329	rBV6	7886	15586	0.28%	0.027%
51	16.435	2336	2340	2345	rVB5	10416	18073	0.32%	0.032%
52	16.512	2345	2353	2354	rBV5	9132	15262	0.27%	0.027%
53	16.570	2359	2363	2365	rVV2	15618	23766	0.42%	0.042%
54	16.588	2365	2366	2371	rVV4	21282	22072	0.39%	0.039%
55	16.653	2371	2377	2382	rVV	1363429	1780410	31.44%	3.113%
56	16.688	2382	2383	2387	rVV3	19668	15285	0.27%	0.027%
57	16.952	2424	2428	2433	rVV2	22491	35212	0.62%	0.062%
58	16.999	2433	2436	2440	rVV2	22864	34184	0.60%	0.060%
59	17.187	2458	2468	2471	rBV2	584168	922064	16.28%	1.612%
60	17.228	2471	2475	2480	rVV	1430316	2084307	36.81%	3.645%
61	17.281	2480	2484	2490	rVB2	1085529	1480925	26.15%	2.590%
62	17.587	2528	2536	2545	rBV	2035597	2903644	51.28%	5.077%
63	17.822	2572	2576	2579	rBV3	8719	13741	0.24%	0.024%
64	17.857	2579	2582	2586	rVB5	10864	14978	0.26%	0.026%
65	17.980	2597	2603	2606	rBV6	5530	11955	0.21%	0.021%
66	18.045	2609	2614	2617	rBV2	12315	15034	0.27%	0.026%
67	18.086	2617	2621	2625	rVV4	31517	47733	0.84%	0.083%
68	18.156	2631	2633	2637	rVV	12778	15731	0.28%	0.028%
69	18.397	2669	2674	2683	rVB6	17128	39440	0.70%	0.069%
70	18.826	2739	2747	2749	rBV8	6556	13528	0.24%	0.024%
71	18.944	2763	2767	2773	rBV5	10595	18210	0.32%	0.032%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

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_G\METHODS\SOM01.2-EPA-BG021015.M
Title : SVOA CALIBRATION

72	19.208	2808	2812	2816	rVB2	16145	23105	0.41%	0.040%
73	19.361	2834	2838	2842	rVV6	10328	18080	0.32%	0.032%
74	19.460	2851	2855	2856	rBV	14860	16476	0.29%	0.029%
75	19.484	2857	2859	2863	rVB	19640	19805	0.35%	0.035%
76	19.572	2868	2874	2877	rBV	1335236	2054734	36.28%	3.593%
77	19.607	2877	2880	2886	rVB	2150607	2752211	48.60%	4.813%
78	20.095	2959	2963	2966	rBV5	14843	19824	0.35%	0.035%
79	20.130	2966	2969	2973	rVB4	10709	15194	0.27%	0.027%
80	20.301	2996	2998	3006	rVB8	11418	19495	0.34%	0.034%
81	20.506	3028	3033	3040	rVV	1997496	2248490	39.71%	3.932%
82	20.859	3090	3093	3096	rVB3	7709	9403	0.17%	0.016%
83	21.076	3125	3130	3141	rBV	849954	1055018	18.63%	1.845%
84	21.282	3161	3165	3172	rBV	2244611	2441184	43.11%	4.269%
85	21.358	3172	3178	3183	rVB	826663	1017751	17.97%	1.780%
86	21.522	3202	3206	3213	rBV3	23257	32215	0.57%	0.056%
87	21.981	3278	3284	3290	rBV3	50519	95766	1.69%	0.167%
88	22.233	3322	3327	3333	rVB	65297	87327	1.54%	0.153%
89	22.451	3360	3364	3367	rBV3	33012	51126	0.90%	0.089%
90	22.586	3382	3387	3397	rVB	174970	270800	4.78%	0.474%
91	22.979	3450	3454	3459	rBV	39844	56842	1.00%	0.099%
92	23.514	3537	3545	3552	rBV2	916323	1779728	31.43%	3.112%
93	23.578	3552	3556	3561	rVV3	37498	66439	1.17%	0.116%
94	23.661	3562	3570	3583	rVB	530651	1014427	17.91%	1.774%
95	24.266	3668	3673	3680	rVB3	41379	79995	1.41%	0.140%
96	24.348	3681	3687	3697	rBV2	88683	201982	3.57%	0.353%
97	25.071	3804	3810	3816	rVB3	28553	63860	1.13%	0.112%
98	26.040	3960	3975	3988	rBV2	1749574	5662794	100.00%	9.902%
99	26.756	4083	4097	4109	rBV	726330	2350730	41.51%	4.111%
100	27.115	4151	4158	4167	rVB8	12708	36890	0.65%	0.065%

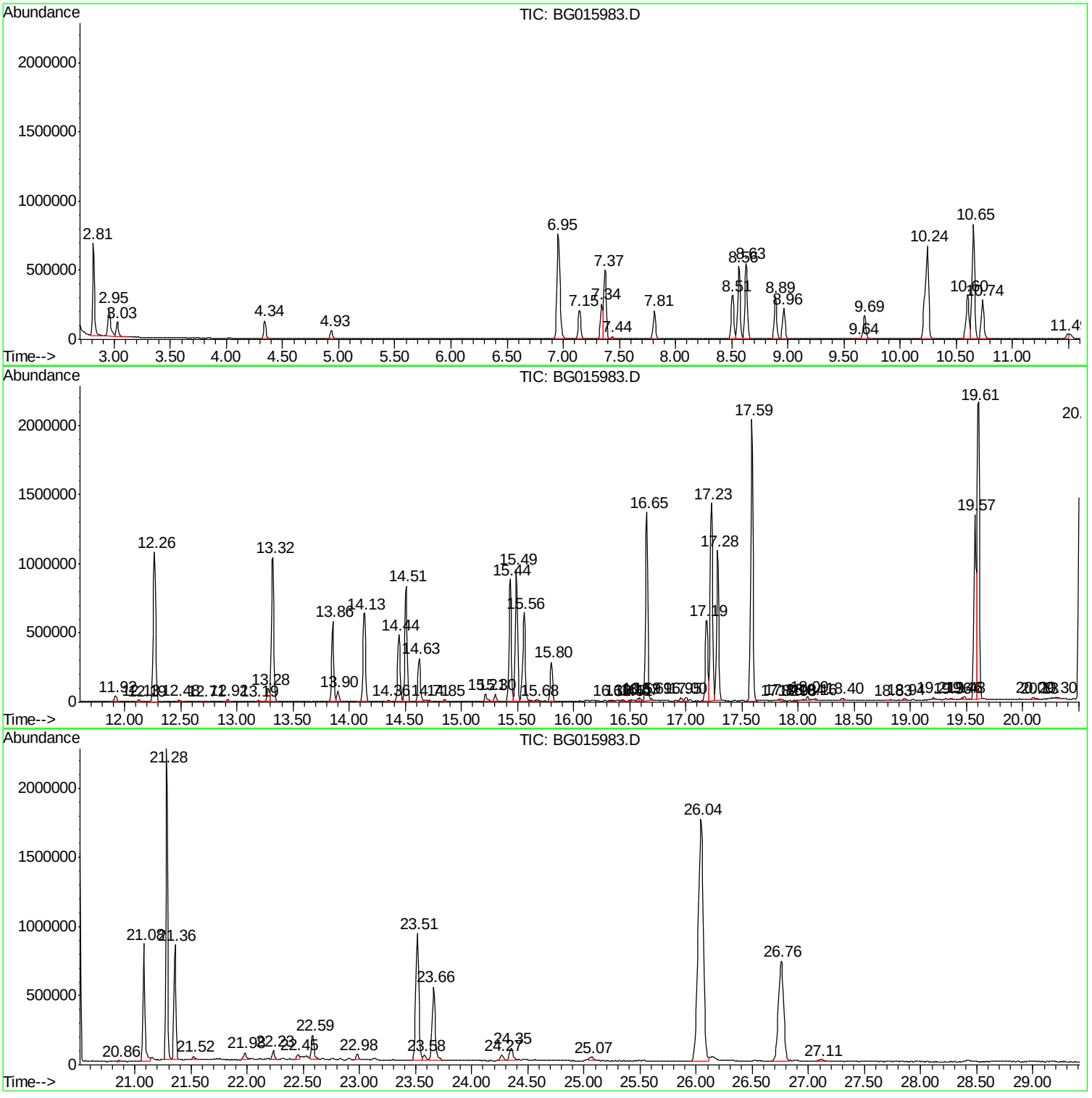
Sum of corrected areas: 57188365

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

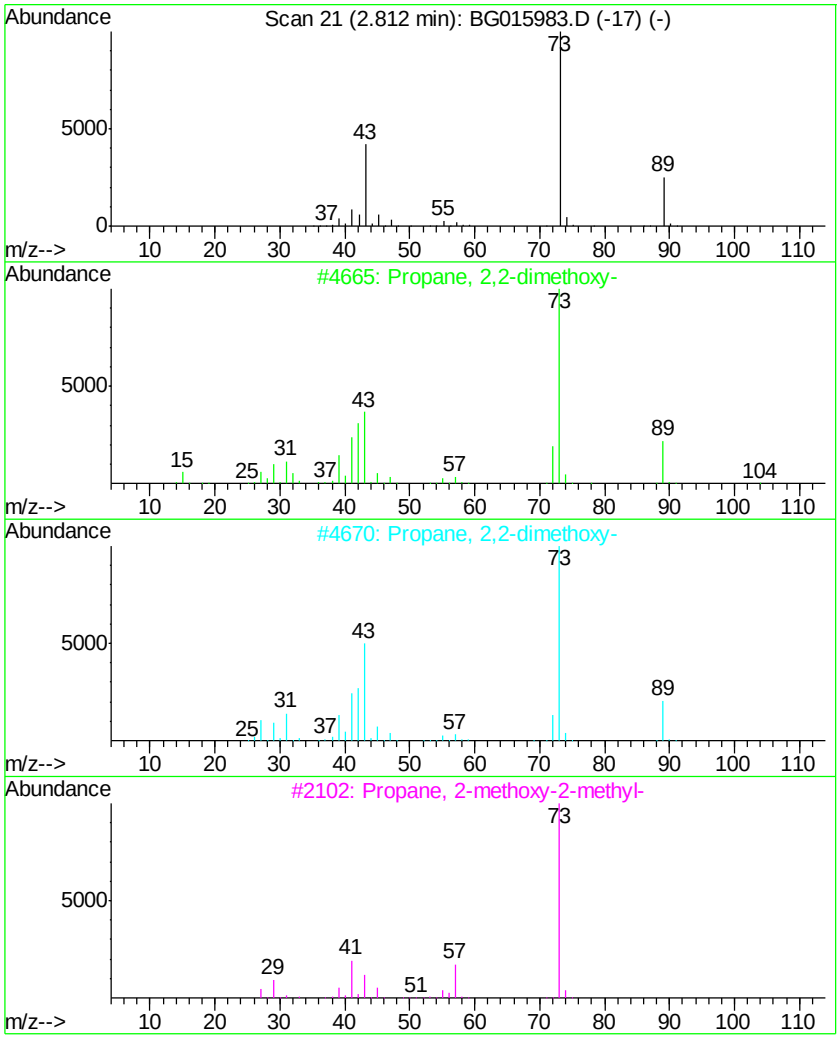
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

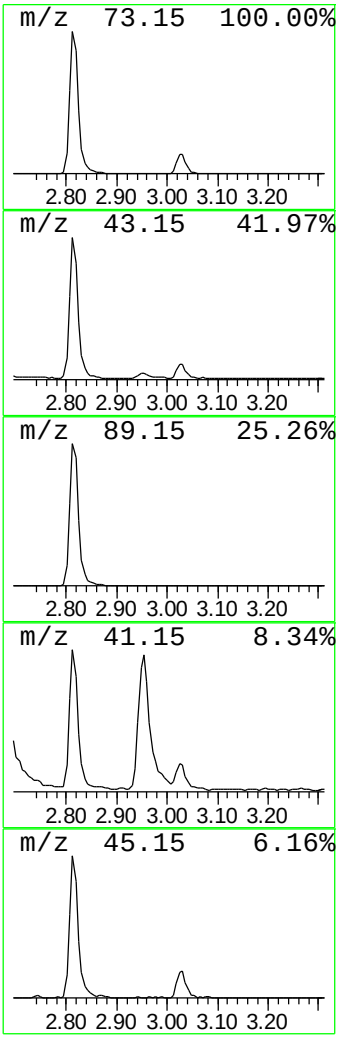
****

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.81	52.87 ng/ul	814867	1,4-Dichlorobenzene-d4	7.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	50
2	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	50
3	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	5





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

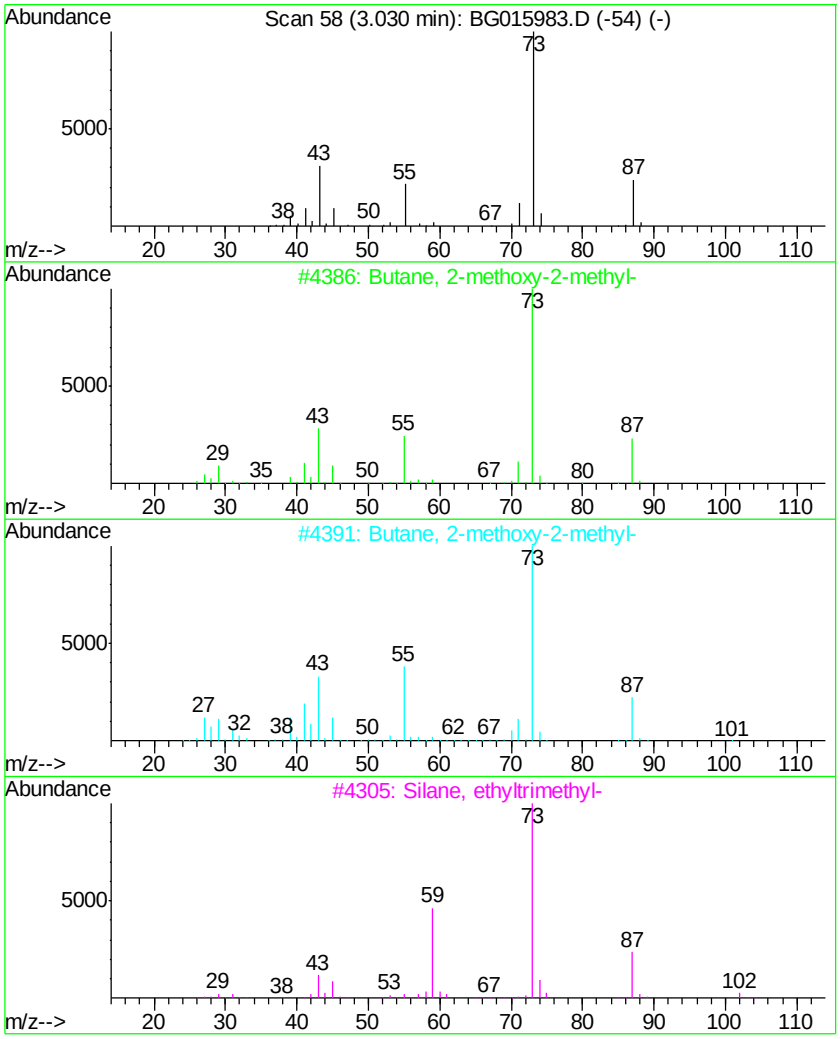
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

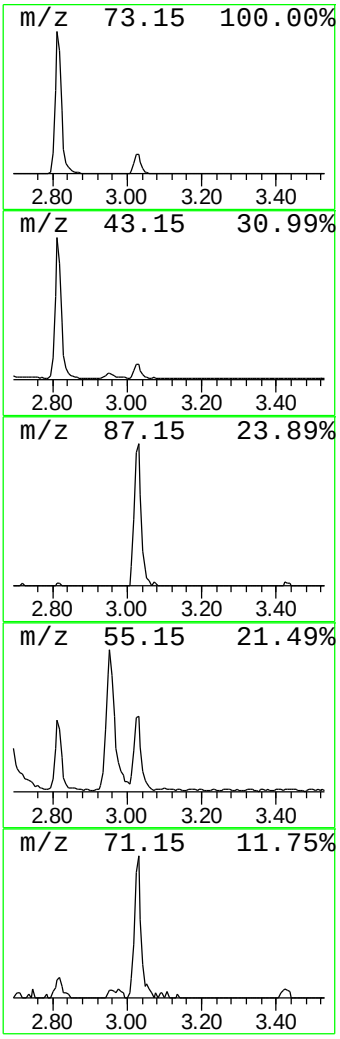
****

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.03	9.90 ng/ul	152535	1,4-Dichlorobenzene-d4	7.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

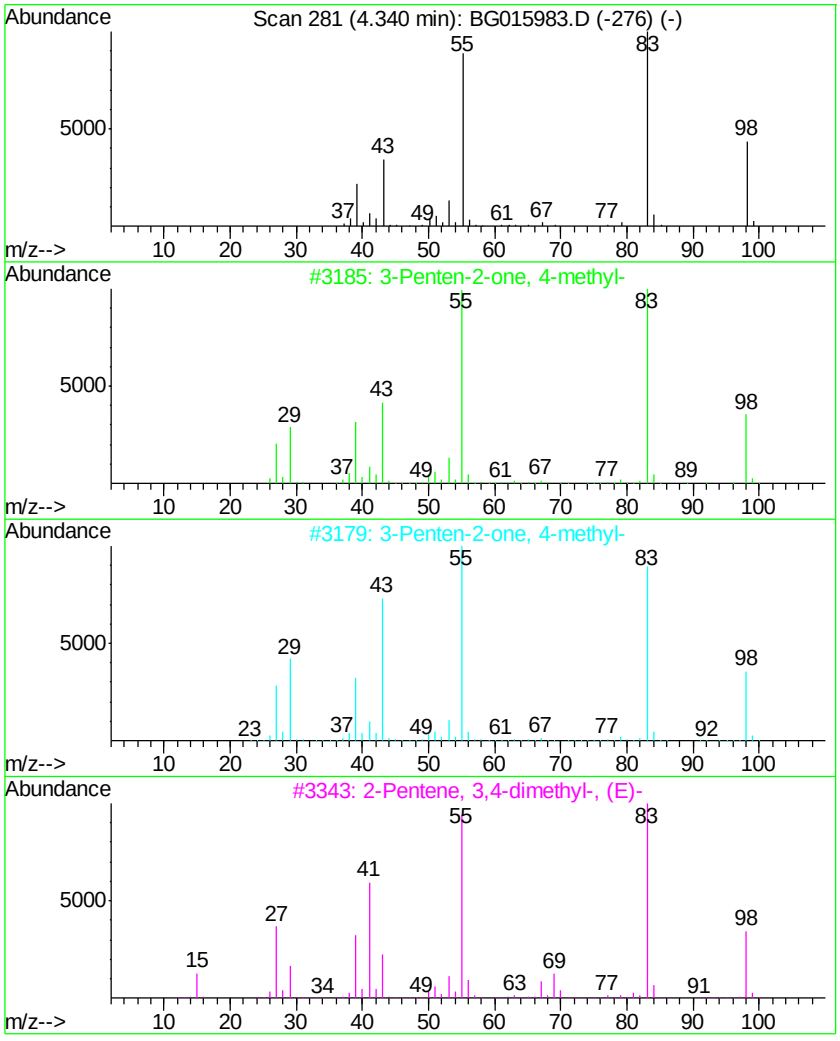
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

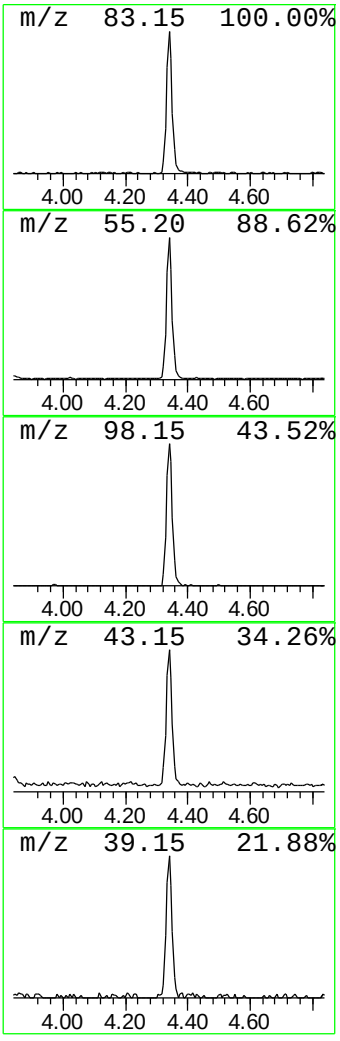
****

Peak Number 4 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.34	10.71 ng/ul	165069	1,4-Dichlorobenzene-d4	7.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
5	2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

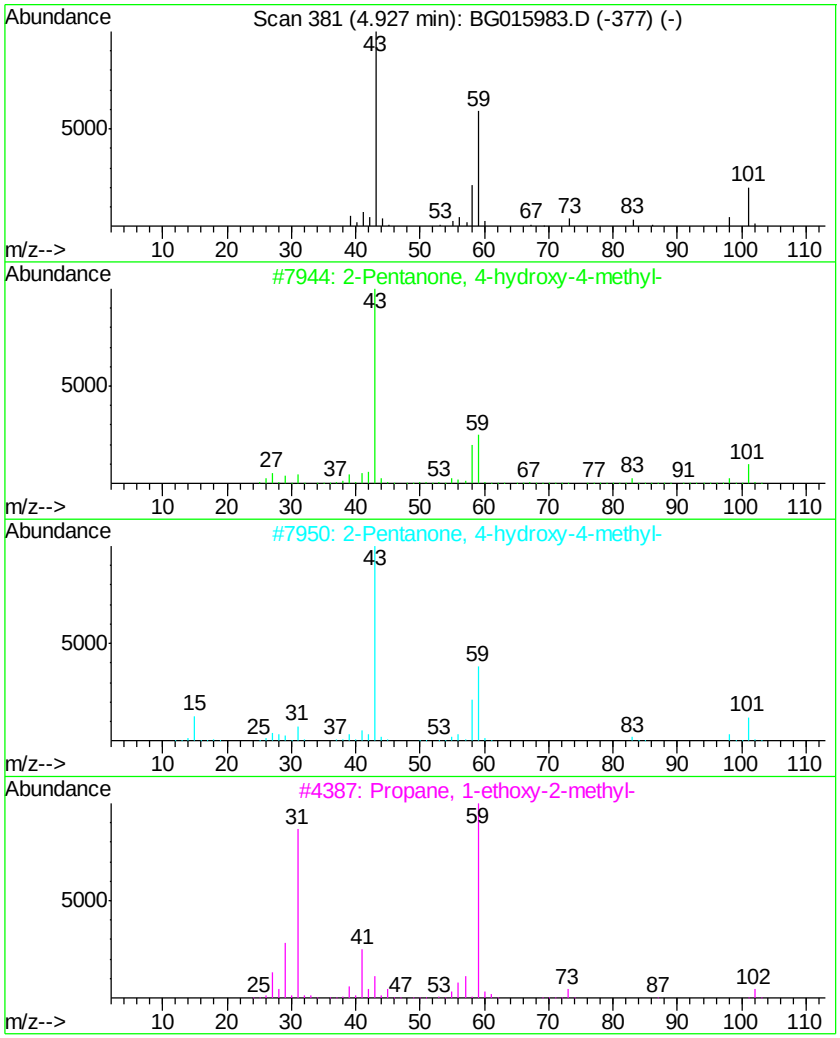
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

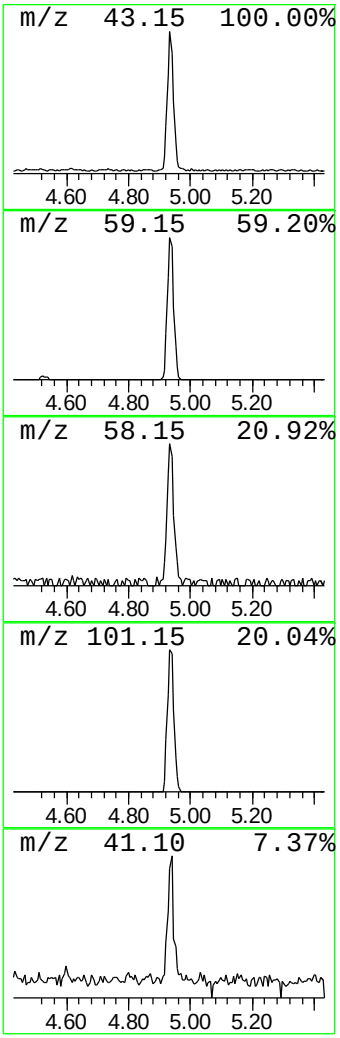
****

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	5.11 ng/ul	78731	1,4-Dichlorobenzene-d4	7.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
3	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	35
4	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	33
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

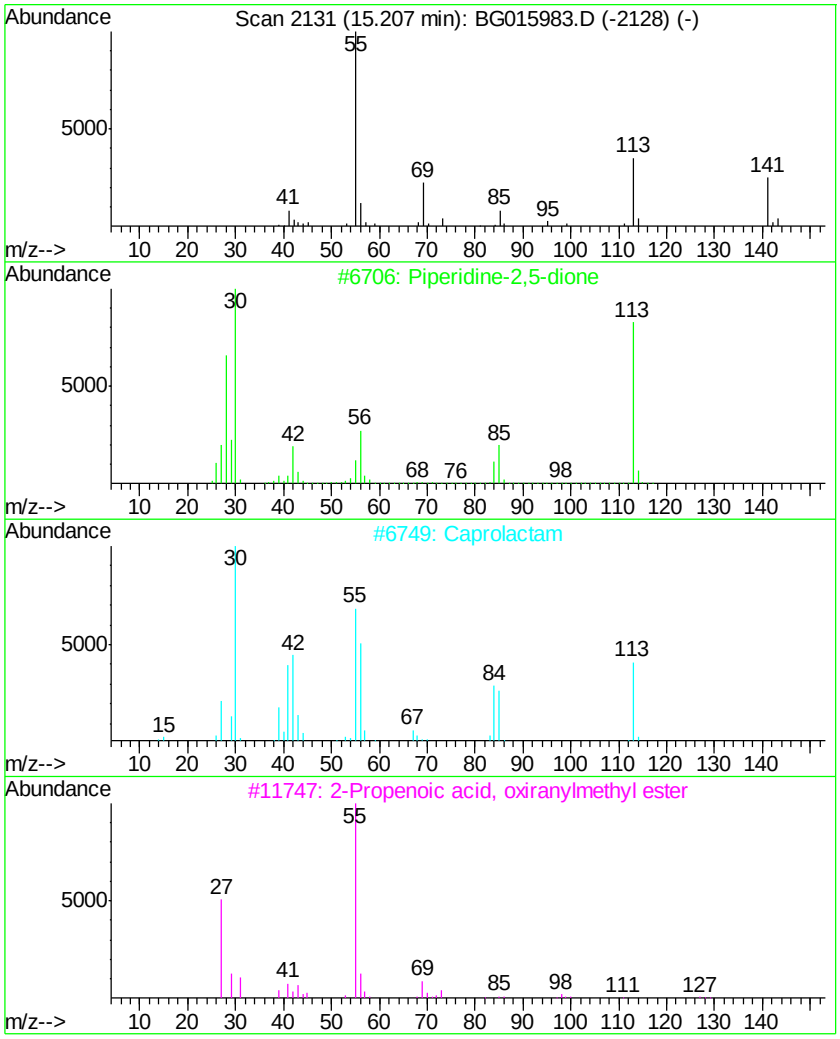
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

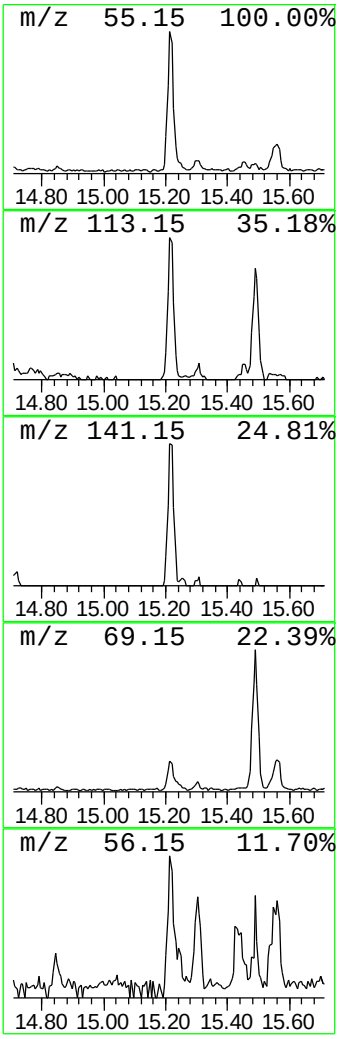
****

Peak Number 7 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	2.27 ng/ul	82082	Acenaphthene-d10	14.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Piperidine-2,5-dione	113	C5H7NO2	052065-78-8	35
2	Caprolactam	113	C6H11NO	000105-60-2	12
3	2-Propenoic acid, oxiranylmethyl...	128	C6H8O3	000106-90-1	10
4	2-Propenoic acid, butyl ester	128	C7H12O2	000141-32-2	9
5	2-Propenoic acid, butyl ester	128	C7H12O2	000141-32-2	9





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021115\
Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

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

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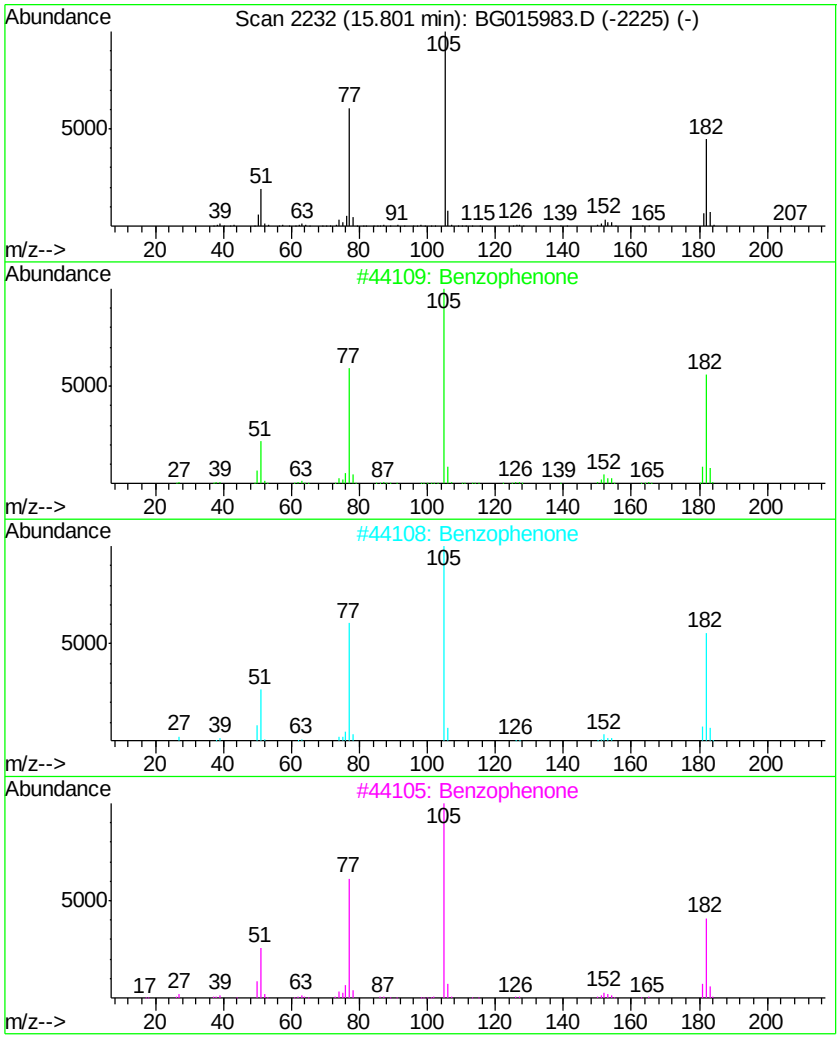
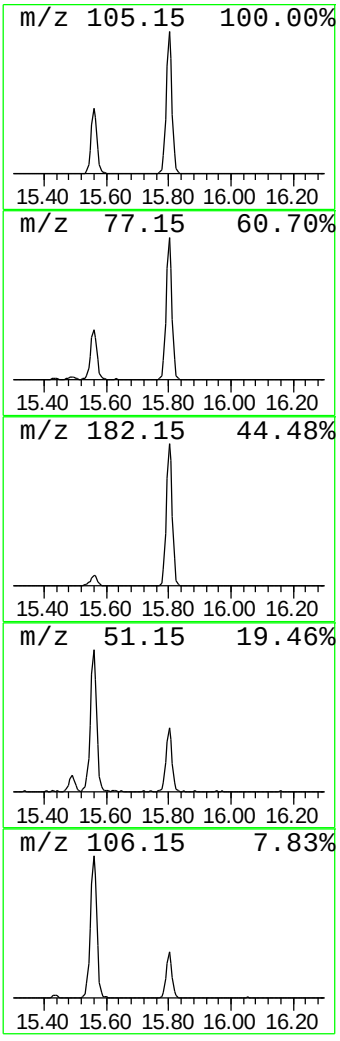
Peak Number 8 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	10.31 ng/ul	372603	Acenaphthene-d10	14.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzophenone	182	C13H10O	000119-61-9	96
3	Benzophenone	182	C13H10O	000119-61-9	95
4	Benzophenone	182	C13H10O	000119-61-9	93
5	Benzophenone	182	C13H10O	000119-61-9	91





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Operator : TP/IZ
Sample : G1227-07
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ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

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TIC Library : C:\DATABASE\NIST02.L
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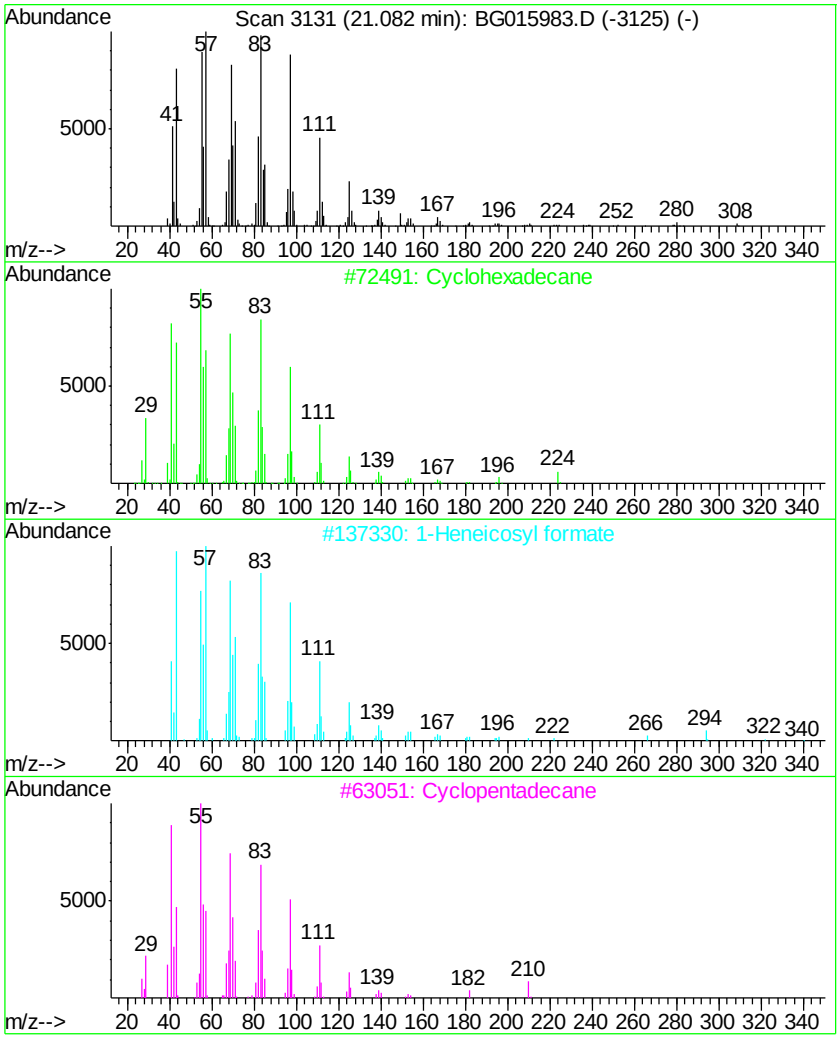
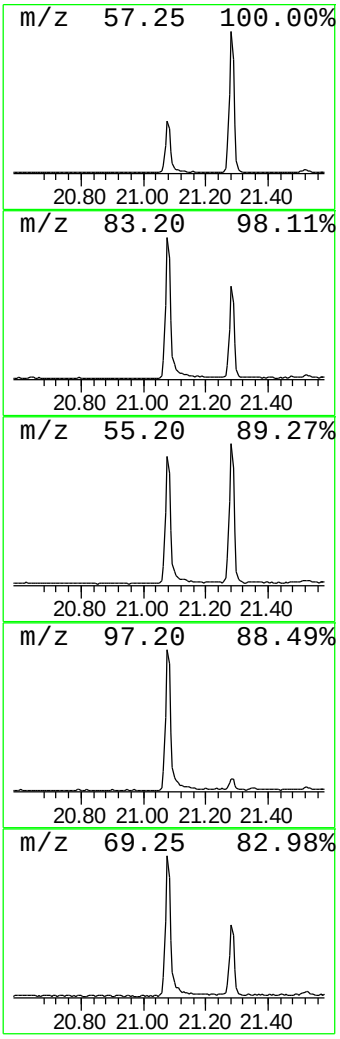
Peak Number 9 Cyclohexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.08	20.73 ng/ul	1055020	Chrysene-d12	21.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	94
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3	Cyclopentadecane	210	C15H30	000295-48-7	93
4	3-Eicosene, (E)-	280	C20H40	074685-33-9	90
5	1-Octadecanol	270	C18H38O	000112-92-5	90





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Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

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

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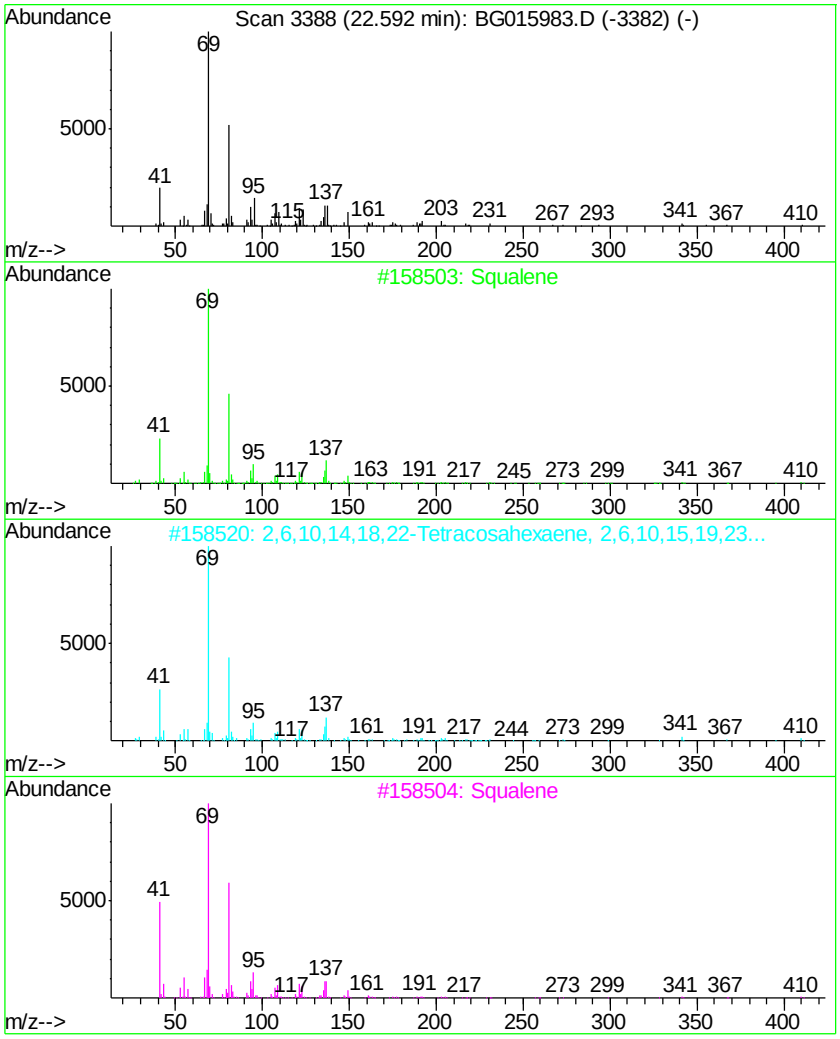
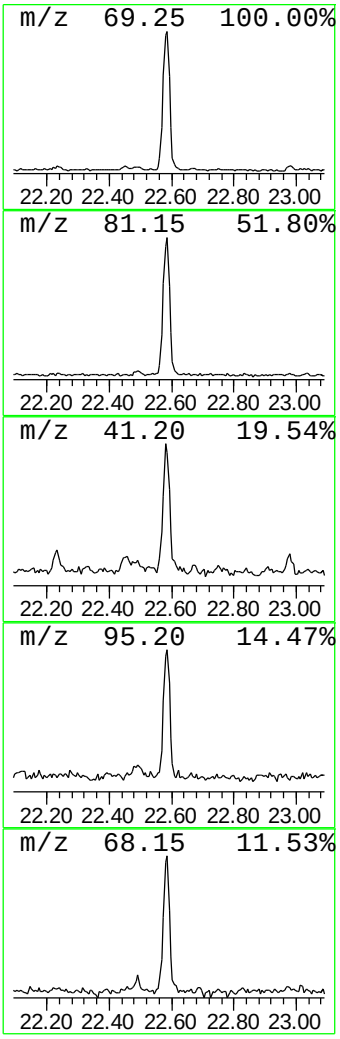
Peak Number 10 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.59	5.34 ng/ul	270800	Perylene-d12	23.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	007683-64-9	91
2	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	91
3	Squalene	410	C30H50	007683-64-9	91
4	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	91
5	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	87





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Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
Quant Title : SVOA CALIBRATION

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

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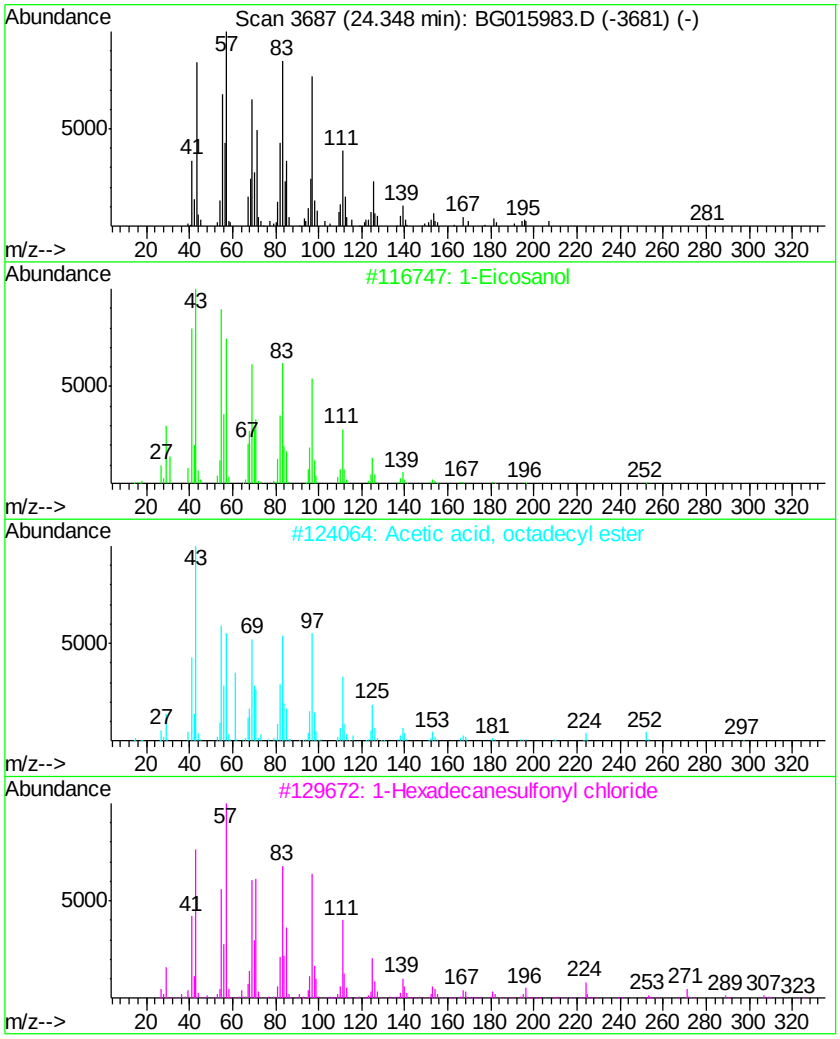
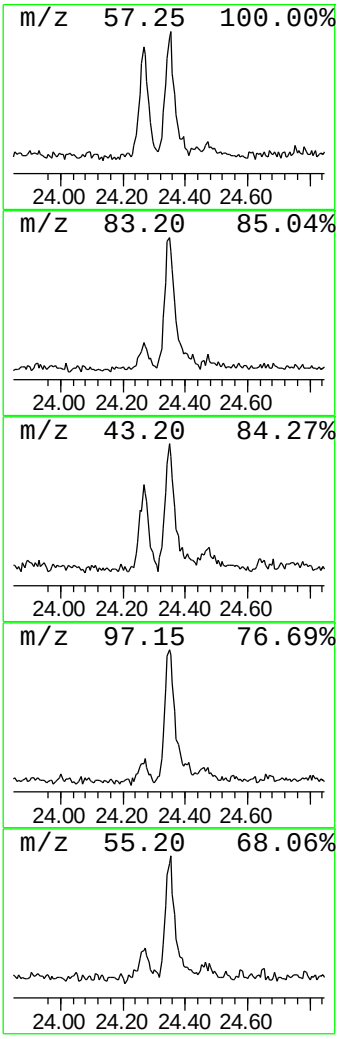
Peak Number 11 1-Eicosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.35	3.98 ng/ul	201982	Perylene-d12	23.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Eicosanol	298	C20H42O	000629-96-9	91
2	Acetic acid, octadecyl ester	312	C20H40O2	000822-23-1	86
3	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	68
4	13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	64
5	1-Octadecanol	270	C18H38O	000112-92-5	58





Tentatively Identified Compound (LSC) summary

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Data File : BG015983.D
Acq On : 11 Feb 2015 13:55
Operator : TP/IZ
Sample : G1227-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : X1H38

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-BG021015.M
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
Propane, 2,2-dime...	2.81	52.9	ng/ul	814867	1	7.81	308282	20.0
Butane, 2-methoxy...	3.03	9.9	ng/ul	152535	1	7.81	308282	20.0
3-Penten-2-one, 4...	4.34	10.7	ng/ul	165069	1	7.81	308282	20.0
2-Pentanone, 4-hy...	4.93	5.1	ng/ul	78731	1	7.81	308282	20.0
unknown-01	15.21	2.3	ng/ul	82082	3	14.44	723110	20.0
Benzophenone	15.80	10.3	ng/ul	372603	3	14.44	723110	20.0
Cyclohexadecane	21.08	20.7	ng/ul	1055020	5	21.36	1017750	20.0
Squalene	22.59	5.3	ng/ul	270800	6	23.66	1014430	20.0
1-Eicosanol	24.35	4.0	ng/ul	201982	6	23.66	1014430	20.0