LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 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-BG011817.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	3.534	115	118	121	rBV3	8417	12103	0.53%	0.084%
2	3.869	173	175	183	rBV7	7511	12101	0.53%	0.084%
3	4.186	226	229	232	rBV4	6760	8243	0.36%	0.057%
4	4.786	328	331	335	rVB5	7350	9516	0.42%	0.066%
5	5.761	492	497	501	rBV5	6801	8713	0.38%	0.060%
6	6.713	656	659	666	rVB7	4369	10761	0.47%	0.074%
7	6.789	666	672	674	rBV4	6470	9977	0.44%	0.069%
8	7.147	729	733	736	rBV4	7188	9273	0.41%	0.064%
9	7.288	750	757	763	rVB7	17962	37748	1.65%	0.261%
10	7.512	787	795	799	rBV7	13752	30287	1.32%	0.209%
11	7.565	799	804	815	rVB2	48694	94010	4.11%	0.650%
12	7.717	825	830	832	rVV5	7197	9604	0.42%	0.066%
13	7.811	842	846	849	rVB5	9763	12344	0.54%	0.085%
14	8.176	900	908	912	rBV2	275490	525654	22.99%	3.634%
15	8.417	941	949	954	rBV2	6108	14224	0.62%	0.098%
16	8.469	954	958	961	rVV6	6011	9288	0.41%	0.064%
17	8.552	966	972	974	rBV6	10559	14766	0.65%	0.102%
18	8.599	974	980	991	rVB2	48953	106058	4.64%	0.733%
19	8.722	994	1001	1008	rBV7	15639	28966	1.27%	0.200%
20	8.793	1008	1013	1020	rVB7	7806	17327	0.76%	0.120%
21	8.840	1020	1021	1027	rVB4	8431	12042	0.53%	0.083%
22	8.969	1037	1043	1053	rVB9	9490	22619	0.99%	0.156%
23	9.286	1091	1097	1102	rBV8	7979	20202	0.88%	0.140%
24	9.357	1106	1109	1113	rVB4	6791	9790	0.43%	0.068%
25	9.392	1113	1115	1119	rBV4	7616	11597	0.51%	0.080%
26	9.462	1119	1127	1136	rBV2	62411	144504	6.32%	0.999%
27	9.774	1176	1180	1185	rVB8	7444	14703	0.64%	0.102%
28	9.850	1189	1193	1199	rBV7	5879	15517	0.68%	0.107%
29	9.915	1199	1204	1206	rVV5	12227	18101	0.79%	0.125%
30	9.944	1206	1209	1215	rVB5	12812	19663	0.86%	0.136%
31	10.015	1218	1221	1224	rBV4	8949	12840	0.56%	0.089%
32	10.085	1231	1233	1242	rVB7	11563	26208	1.15%	0.181%
33	10.156	1242	1245	1253	rBV10	9070	14682	0.64%	0.102%
34	10.385	1280	1284	1288	rVB6	10573	14507	0.63%	0.100%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 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-BG011817.M
Title : SVOA CALIBRATION

35	10.502	1298	1304	1314	rBV9	10533	32585	1.42%	0.225%
36	10.626	1320	1325	1327	rBV5	6271	9058	0.40%	0.063%
37	10.878	1361	1368	1375	rVB9	4554	13143	0.57%	0.091%
38	11.002	1375	1389	1394	rBV	402203	809484	35.40%	5.596%
39	11.049	1394	1397	1407	rVB2	215337	409181	17.89%	2.829%
40	11.131	1408	1411	1414	rVV5	9051	11199	0.49%	0.077%
41	11.184	1414	1420	1424	rVB6	9257	20774	0.91%	0.144%
42	11.260	1427	1433	1440	rBV7	22320	49727	2.17%	0.344%
43	11.331	1440	1445	1458	rVB5	29183	59205	2.59%	0.409%
44	11.472	1464	1469	1475	rBV6	18016	33214	1.45%	0.230%
45	11.713	1504	1510	1513	rVB7	7274	9935	0.43%	0.069%
46	11.836	1523	1531	1536	rBV7	27535	54781	2.40%	0.379%
47	11.883	1536	1539	1549	rVV7	20631	48431	2.12%	0.335%
48	12.001	1556	1559	1563	rVB4	8122	9001	0.39%	0.062%
49	12.418	1625	1630	1635	rBV6	7215	9818	0.43%	0.068%
50	12.559	1650	1654	1657	rBV4	4851	9323	0.41%	0.064%
51	12.647	1662	1669	1678	rVV3	53992	106477	4.66%	0.736%
52	12.747	1678	1686	1695	rVB2	24714	54106	2.37%	0.374%
53	12.870	1698	1707	1714	rVB3	39875	85042	3.72%	0.588%
54	13.076	1740	1742	1747	rBV4	6857	9052	0.40%	0.063%
55	13.311	1777	1782	1784	rBV4	5932	9235	0.40%	0.064%
56	13.358	1784	1790	1801	rVV7	59062	174662	7.64%	1.208%
57	13.581	1822	1828	1835	rBV7	13408	28703	1.26%	0.198%
58	13.634	1835	1837	1848	rVB8	15136	33681	1.47%	0.233%
59	13.834	1868	1871	1876	rVB5	12063	15318	0.67%	0.106%
60	13.898	1876	1882	1888	rBV10	14808	35665	1.56%	0.247%
61	13.975	1890	1895	1905	rVB10	10091	26193	1.15%	0.181%
62	14.133	1914	1922	1928	rBV9	15424	41656	1.82%	0.288%
63	14.198	1928	1933	1941	rVB7	14087	38603	1.69%	0.267%
64	14.363	1956	1961	1965	rBV4	11325	20124	0.88%	0.139%
65	14.504	1979	1985	1991	rBV6	17007	32956	1.44%	0.228%
66	14.592	1995	2000	2004	rBV7	8090	15136	0.66%	0.105%
67	14.662	2004	2012	2022	rBV5	87958	182864	8.00%	1.264%
68	14.803	2027	2036	2044	rBV	689121	1170614	51.19%	8.093%
69	14.868	2044	2047	2052	rVB6	20371	30877	1.35%	0.213%
70	14.968	2059	2064	2076	rVB9	23305	52279	2.29%	0.361%
71	15.085	2076	2084	2090	rBV2	98317	195363	8.54%	1.351%

LSC Area Percent Report

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Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 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-BG011817.M
Title : SVOA CALIBRATION

72	15.150	2090	2095	2102	rVV6	42281	75797	3.31%	0.524%
73	15.244	2103	2111	2119	rVB6	43024	118001	5.16%	0.816%
74	15.320	2121	2124	2130	rVB5	23568	38936	1.70%	0.269%
75	15.402	2130	2138	2140	rBV4	48997	75062	3.28%	0.519%
76	15.444	2141	2145	2156	rVB7	65254	149510	6.54%	1.034%
77	15.567	2161	2166	2170	rBV6	6813	15736	0.69%	0.109%
78	15.673	2182	2184	2189	rVB6	9110	11951	0.52%	0.083%
79	15.796	2201	2205	2211	rBV2	21618	50789	2.22%	0.351%
80	15.855	2212	2215	2221	rVB5	23292	38074	1.66%	0.263%
81	15.949	2223	2231	2253	rBV	1154381	2286876	100.00%	15.810%
82	16.390	2305	2306	2311	rBV4	7745	12258	0.54%	0.085%
83	16.484	2320	2322	2327	rBV5	8723	11791	0.52%	0.082%
84	16.542	2327	2332	2336	rBV7	15954	32097	1.40%	0.222%
85	16.748	2364	2367	2375	rVB9	7419	14300	0.63%	0.099%
86	16.830	2378	2381	2386	rVB6	6892	10796	0.47%	0.075%
87	16.912	2391	2395	2403	rVB10	10449	24828	1.09%	0.172%
88	17.212	2434	2446	2456	rBV	710501	1337192	58.47%	9.245%
89	17.435	2478	2484	2489	rVB10	23547	40048	1.75%	0.277%
90	17.553	2497	2504	2517	rBV2	891901	1474714	64.49%	10.195%
91	17.659	2518	2522	2526	rVB4	16283	23871	1.04%	0.165%
92	17.764	2537	2540	2545	rBV2	34907	45061	1.97%	0.312%
93	17.976	2570	2576	2582	rVB	43933	72156	3.16%	0.499%
94	18.105	2594	2598	2601	rBV6	11767	15550	0.68%	0.108%
95	18.346	2638	2639	2641	rBV2	12133	9583	0.42%	0.066%
96	18.428	2652	2653	2657	rBV4	10213	12392	0.54%	0.086%
97	18.869	2727	2728	2735	rVB7	10866	18366	0.80%	0.127%
98	19.938	2906	2910	2917	rBV3	28620	52278	2.29%	0.361%
99	21.848	3228	3235	3243	rVB	888497	1578853	69.04%	10.915%
100	25.238	3802	3812	3829	rVB2	462132	1532236	67.00%	10.593%

Sum of corrected areas: 14464505

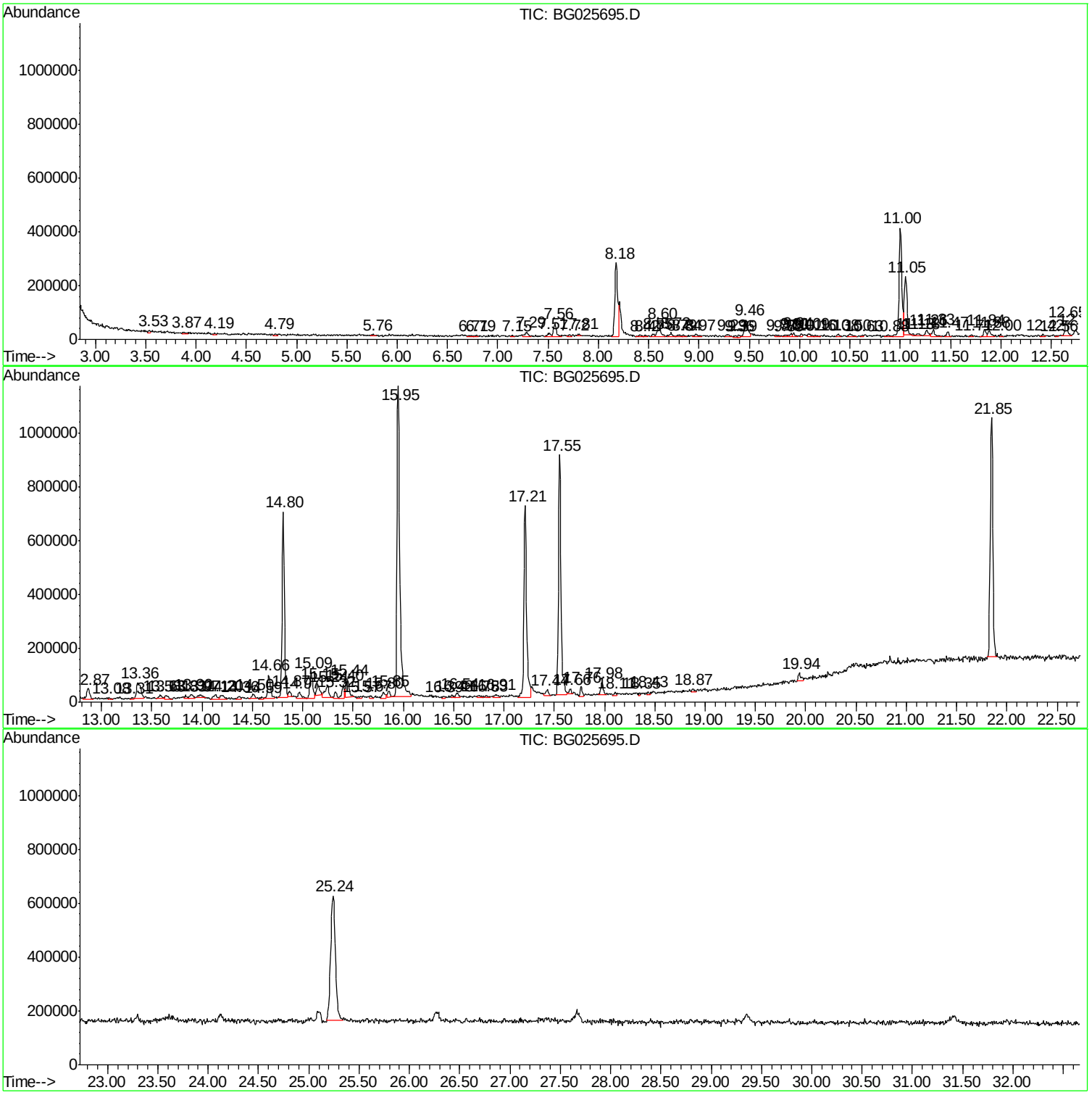
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
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Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

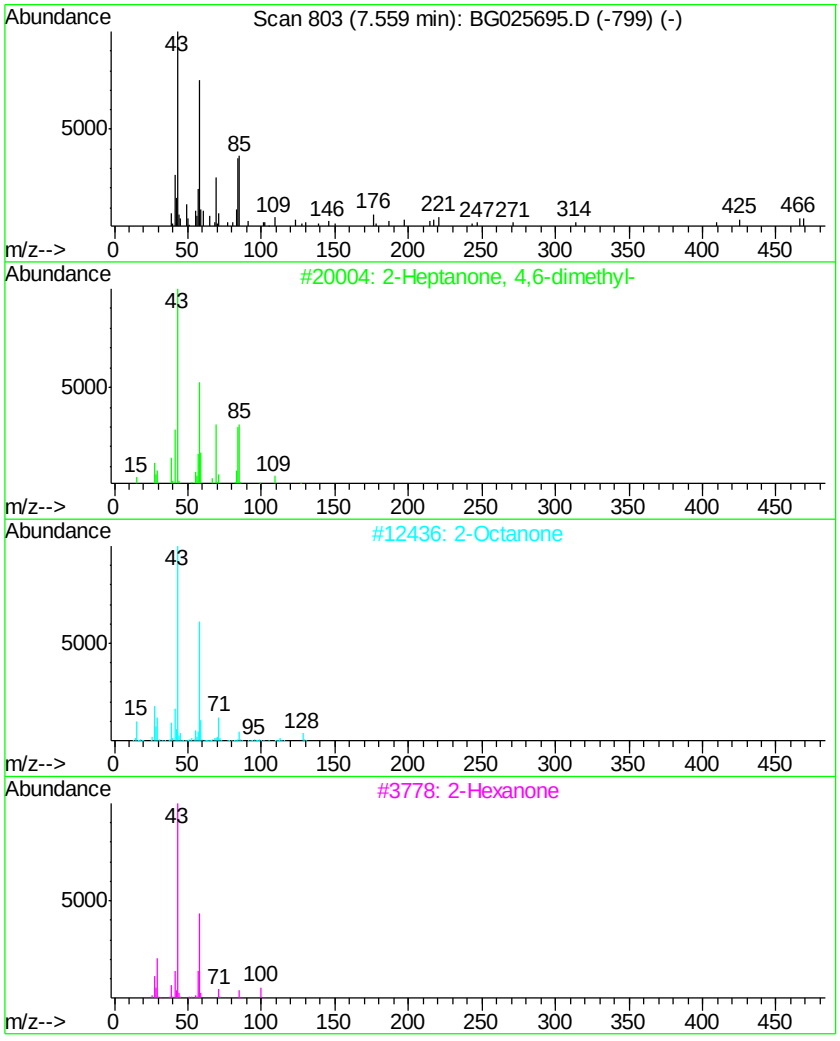
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

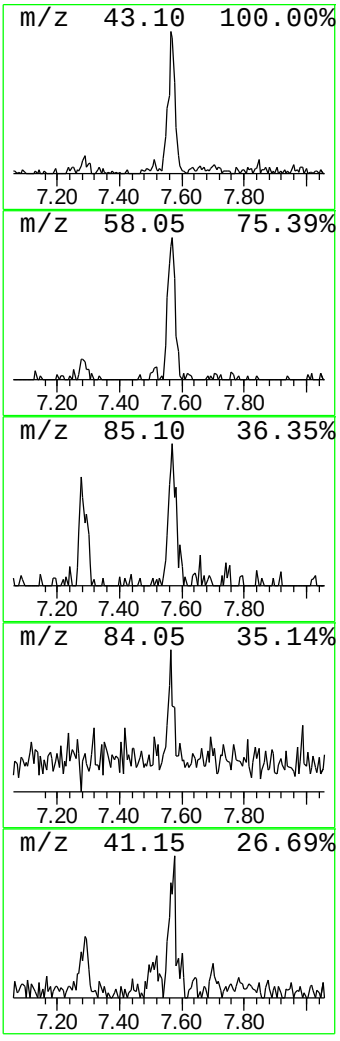
****

Peak Number 1 2-Heptanone, 4,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	3.58 ng/ul	94010	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	59
2	2-Octanone	128	C8H16O	000111-13-7	32
3	2-Hexanone	100	C6H12O	000591-78-6	30
4	Heptanoic acid, 6-oxo-	144	C7H12O3	003128-07-2	28
5	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	25





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

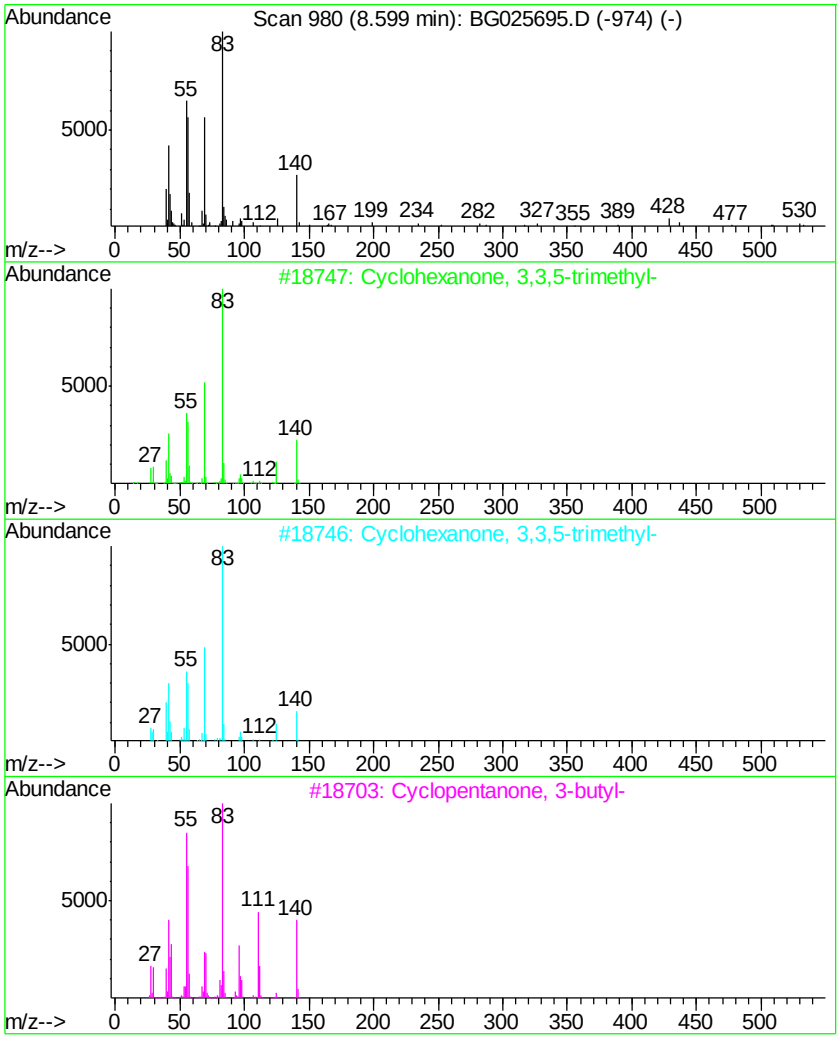
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

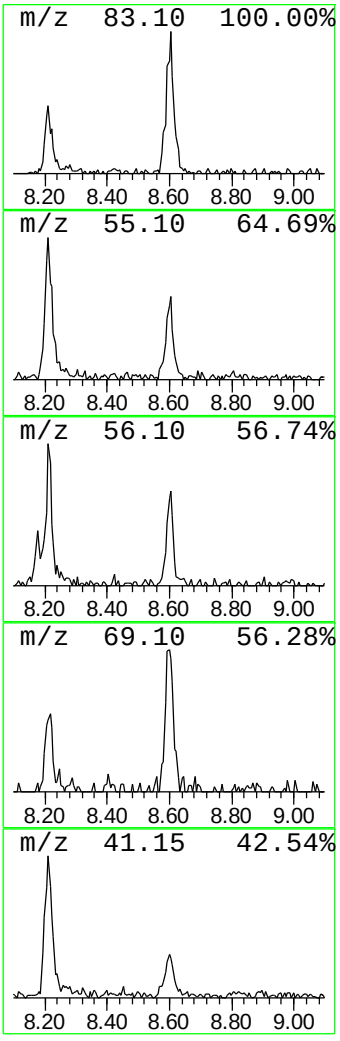
****

Peak Number 2 Cyclohexanone, 3,3,5-trimet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	4.04 ng/ul	106058	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	83
3	Cyclopentanone, 3-butyl-	140	C9H16O	057283-81-5	72
4	1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	59
5	1-Hexyn-3-ol	98	C6H10O	000105-31-7	59





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
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Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

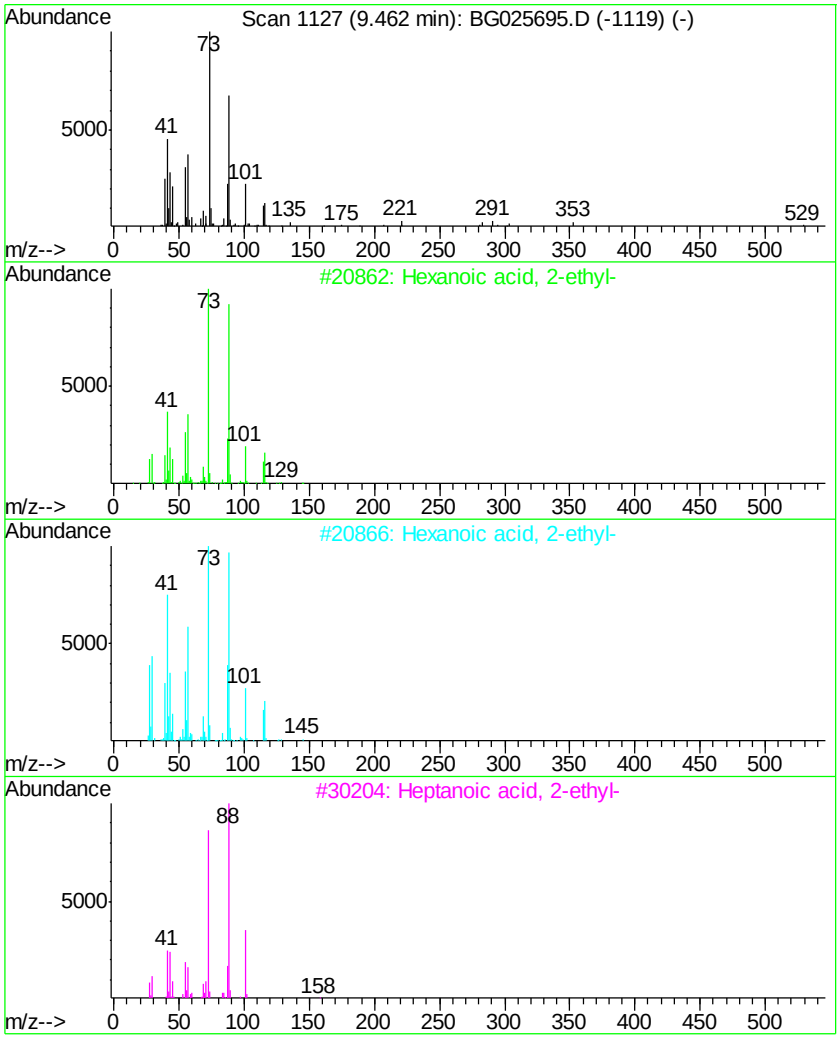
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

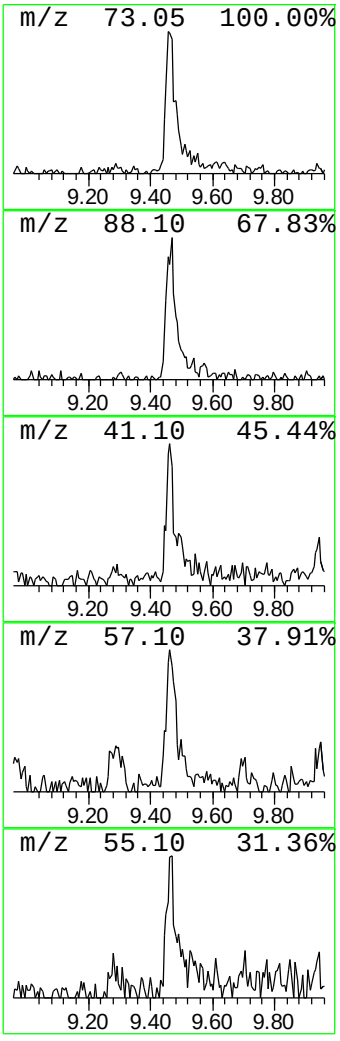
****

Peak Number 3 Hexanoic acid, 2-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.46	5.50 ng/ul	144504	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	59
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	50
3	Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	47
4	.beta.-D-Gluco-Hexodialdo-1,5-py...	414	C16H22N4O9	041545-31-7	38
5	Carbonimidodithioic acid, methyl...	135	C4H9NS2	018805-25-9	38





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

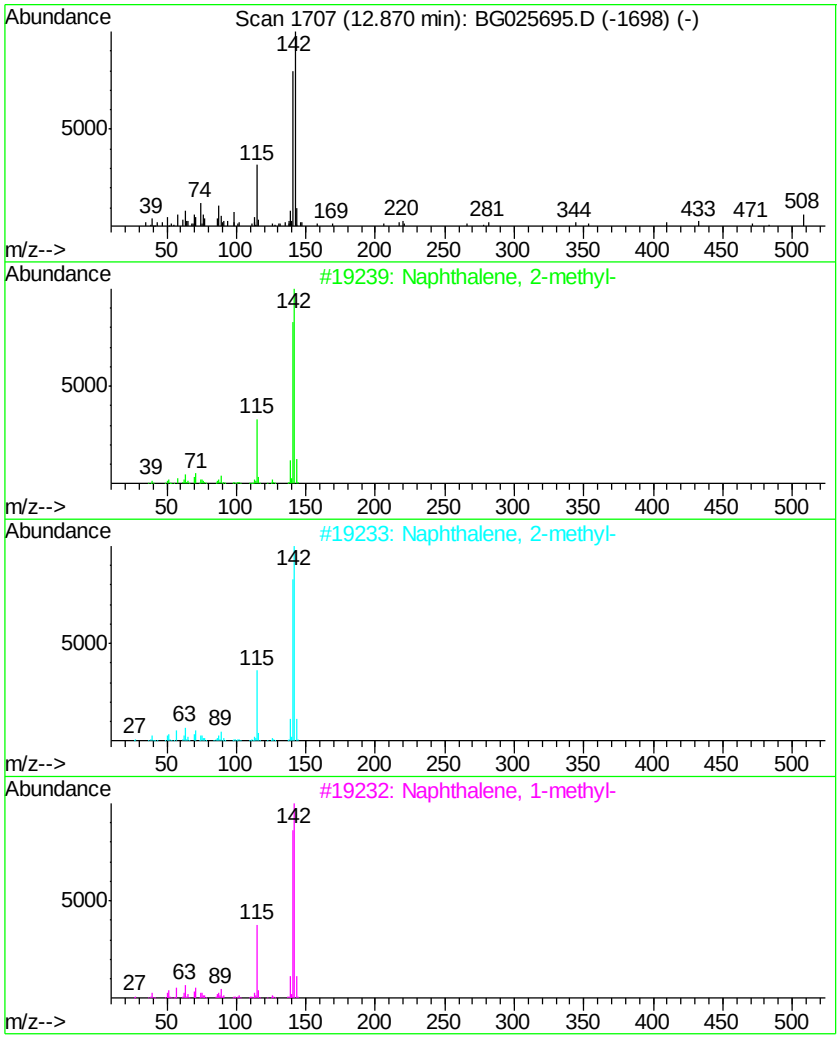
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

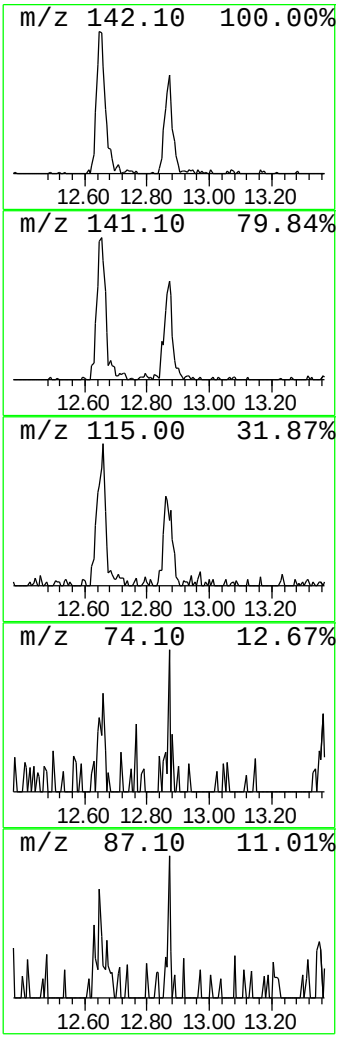
****

Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	2.10 ng/ul	85042	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	90
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	90
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	90
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	87





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

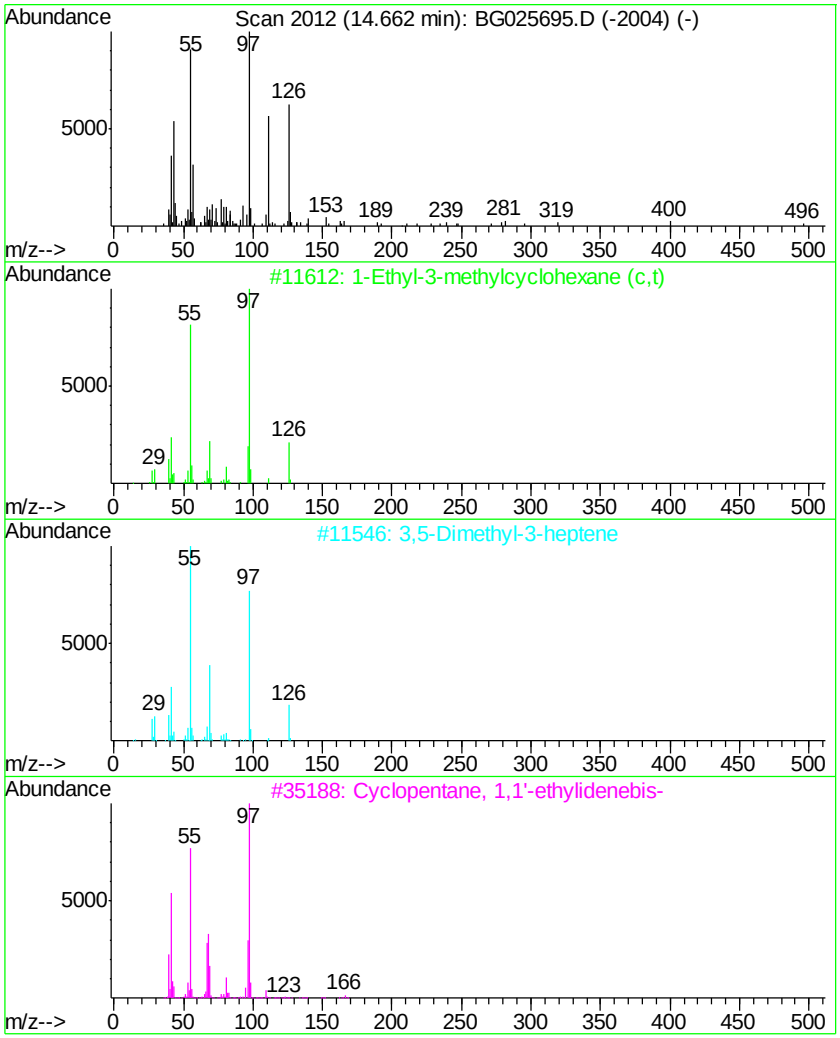
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

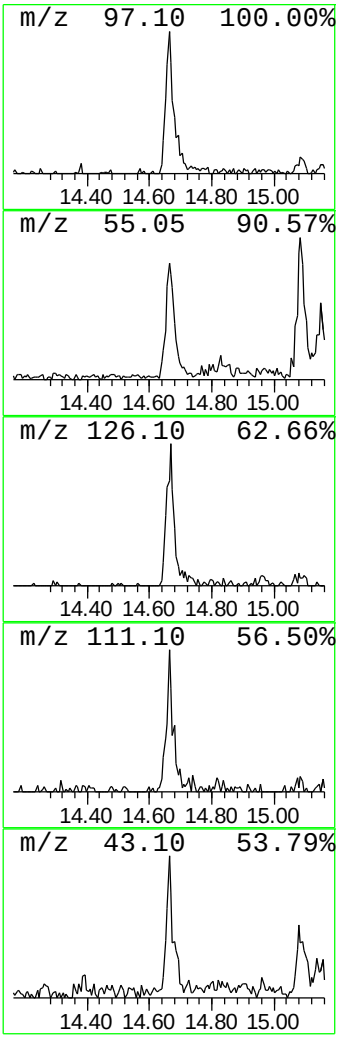
****

Peak Number 5 (DEL) Alkane: Cyclic14.66 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	3.12 ng/ul	182864	Acenaphthene-d10	14.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	58
2	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	43
3	Cyclopentane, 1,1'-ethylidenebis-	166	C12H22	004413-21-2	43
4	2-Pentene, 2,3,4-trimethyl-	112	C8H16	000565-77-5	38
5	2-Pentene, 2,4,4-trimethyl-	112	C8H16	000107-40-4	38





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Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

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

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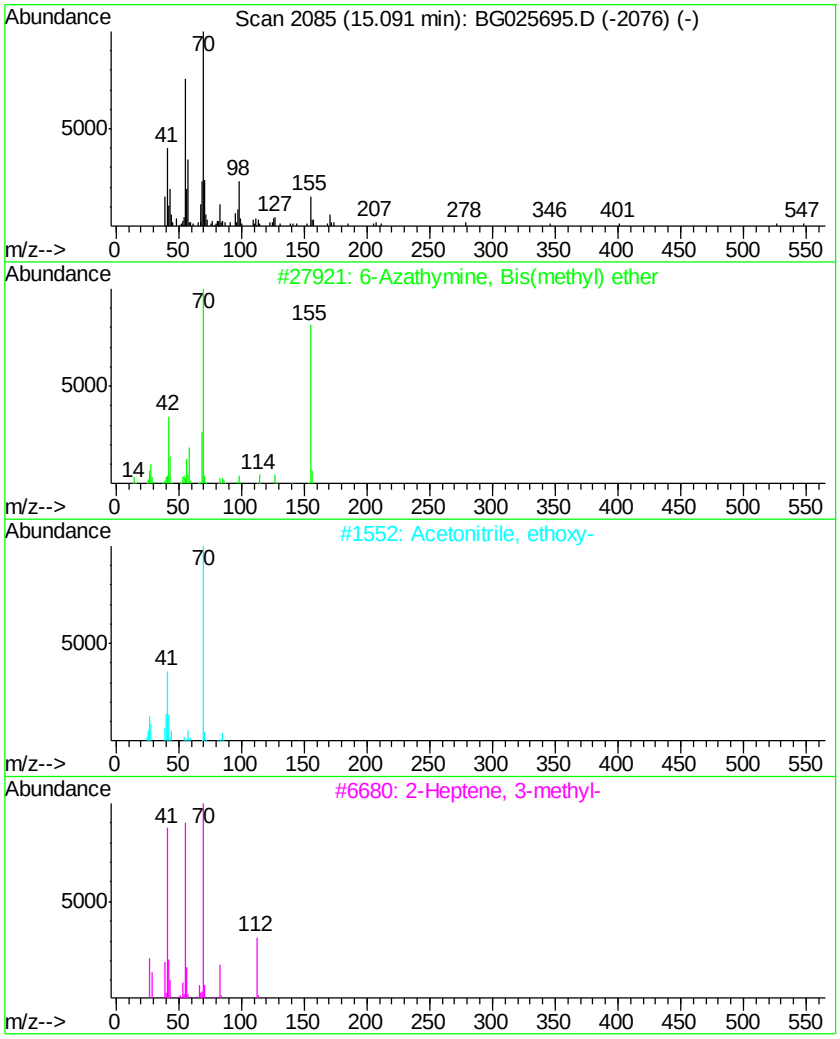
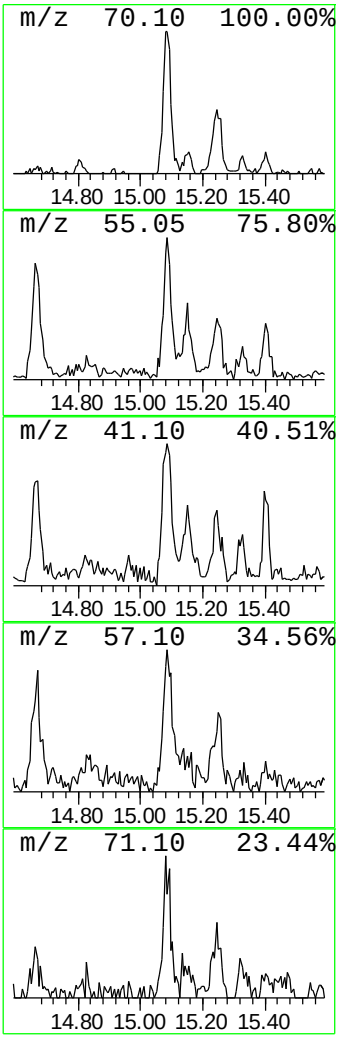
Peak Number 6 6-Azathymine, Bis(methyl) e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	3.34 ng/ul	195363	Acenaphthene-d10	14.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Azathymine, Bis(methyl) ether	155	C6H9N3O2	1000333-80-4	52
2	Acetonitrile, ethoxy-	85	C4H7NO	062957-60-2	47
3	2-Heptene, 3-methyl-	112	C8H16	003404-75-9	47
4	Pentanoic acid, 3-methylbutyl ester	172	C10H20O2	002050-09-1	47
5	2-Nonene, 3-methyl-, (E)-	140	C10H20	017003-99-5	47





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Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

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

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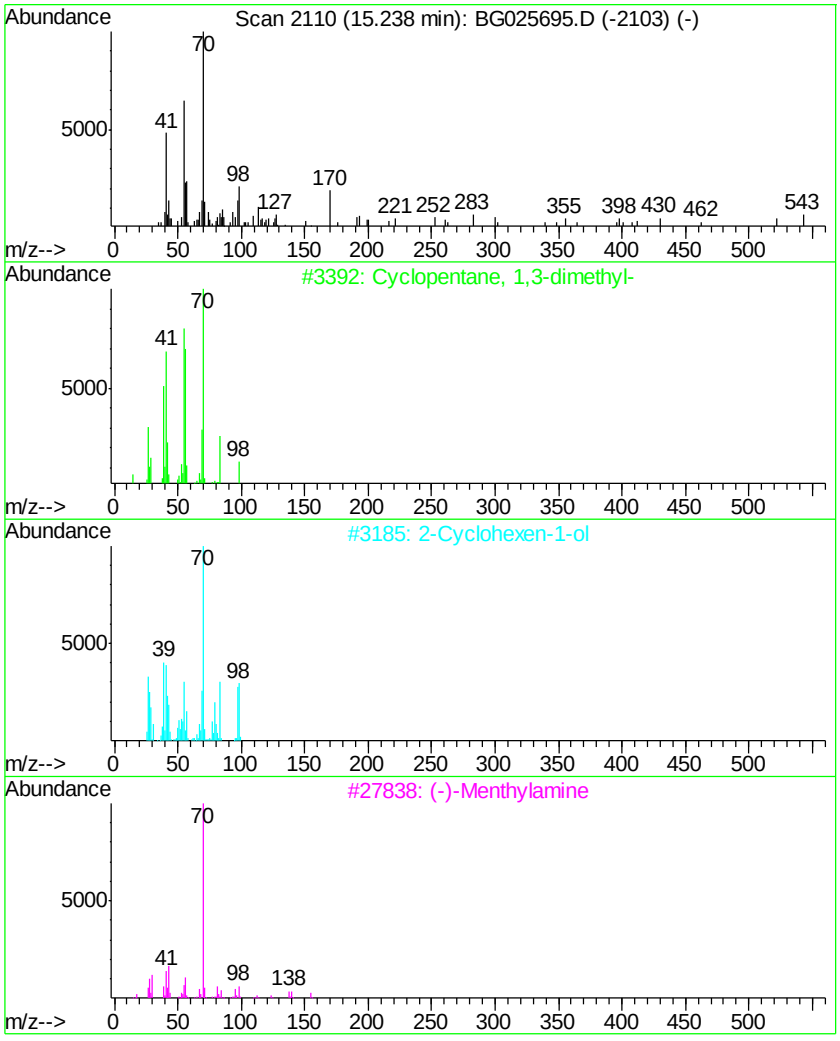
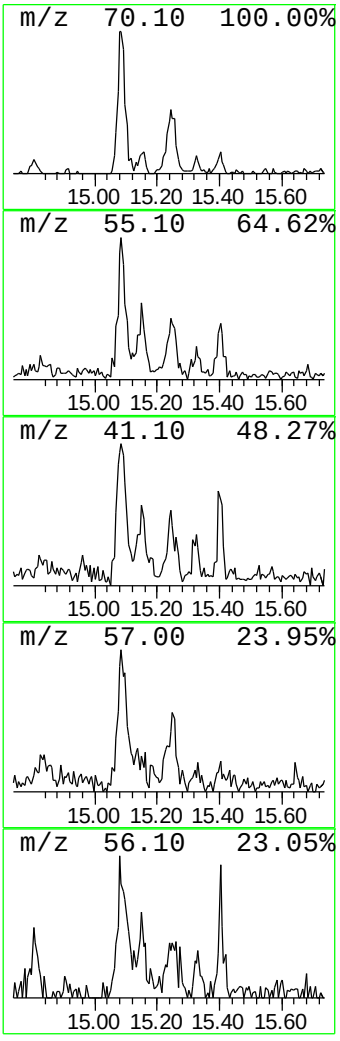
Peak Number 7 (DEL) Alkane: Cyclic15.24 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.24	2.02 ng/ul	118001	Acenaphthene-d10	14.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	52
2	2-Cyclohexen-1-ol	98	C6H10O	000822-67-3	50
3	(-)-Menthylamine	155	C10H21N	021411-81-4	46
4	Cyclopentane, 1,2,3-trimethyl-	112	C8H16	002815-57-8	43
5	Cyclopropane, 1-methyl-1-(1-meth...	224	C16H32	041977-40-6	42





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Data File : BG025695.D
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Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

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

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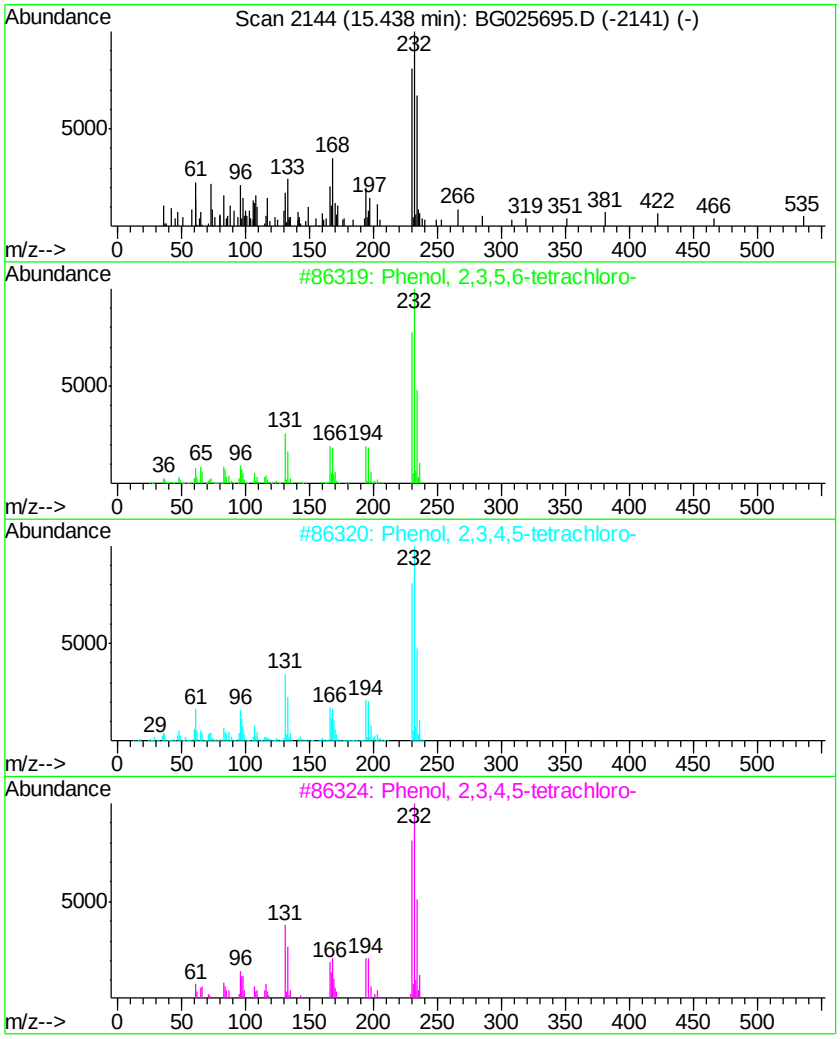
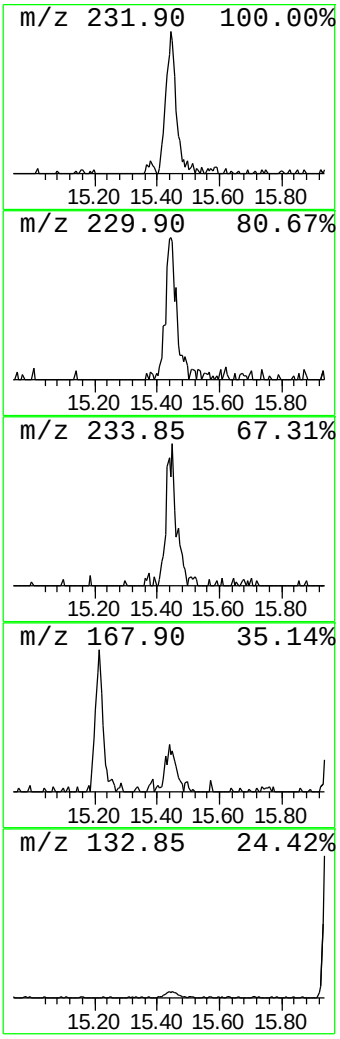
Peak Number 8 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.44	2.55 ng/ul	149510	Acenaphthene-d10	14.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	93
2	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	90
3	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	89
4	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	87
5	2,4,5,6-Tetrachloro-pyridin-3-yl...	230	C5H2Cl4N2	1000277-70-7	76





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Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

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

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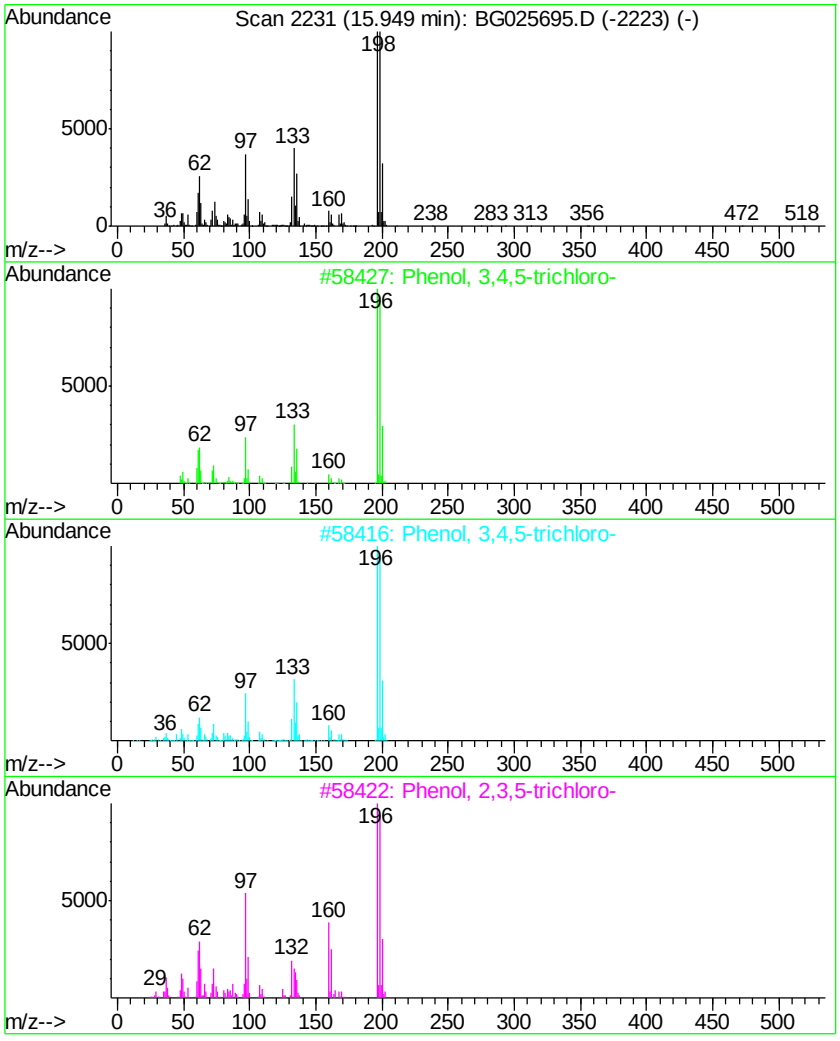
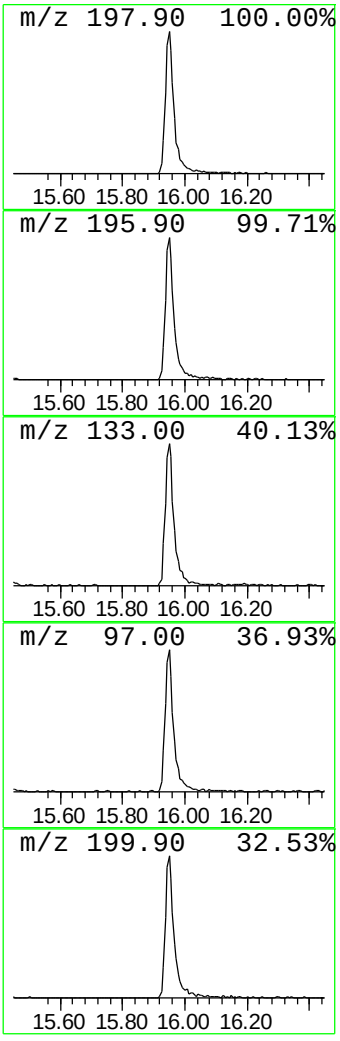
Peak Number 9 Phenol, 3,4,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.95	39.07 ng/ul	2286880	Acenaphthene-d10	14.80

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4,5-trichloro-	196	C6H3Cl3O	000609-19-8	95
2	Phenol, 3,4,5-trichloro-	196	C6H3Cl3O	000609-19-8	94
3	Phenol, 2,3,5-trichloro-	196	C6H3Cl3O	000933-78-8	87
4	Phenol, 2,3,4-trichloro-	196	C6H3Cl3O	015950-66-0	87
5	Phenol, 2,3,4-trichloro-	196	C6H3Cl3O	015950-66-0	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025695.D
Acq On : 27 Jan 2017 9:39
Operator : SJ/MA
Sample : I1387-03DL2 100X
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9Q7DL2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Heptanone, 4,6-...	7.56	3.6	ng/ul	94010	1	8.18	525654	20.0
Cyclohexanone, 3,...	8.60	4.0	ng/ul	106058	1	8.18	525654	20.0
Hexanoic acid, 2-...	9.46	5.5	ng/ul	144504	1	8.18	525654	20.0
Naphthalene, 1-me...	12.87	2.1	ng/ul	85042	2	11.00	809484	20.0
(DEL) Alkane: Cyc...	14.66	3.1	ng/ul	182864	3	14.80	1170610	20.0
6-Azathymine, Bis...	15.09	3.3	ng/ul	195363	3	14.80	1170610	20.0
(DEL) Alkane: Cyc...	15.24	2.0	ng/ul	118001	3	14.80	1170610	20.0
Phenol, 2,3,5,6-t...	15.44	2.5	ng/ul	149510	3	14.80	1170610	20.0
Phenol, 3,4,5-tri...	15.95	39.1	ng/ul	2286880	3	14.80	1170610	20.0