LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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	3.866	169	175	184	rVB7	16373	44048	0.19%	0.086%
2	5.916	519	524	532	rVB7	13951	34759	0.15%	0.068%
3	6.645	643	648	656	rVB5	17315	37185	0.16%	0.073%
4	7.285	744	757	765	rBV	121596	244132	1.03%	0.477%
5	7.503	785	794	799	rBV3	66449	126932	0.53%	0.248%
6	7.567	799	805	815	rVB	461704	760824	3.20%	1.486%
7	7.714	824	830	838	rBV7	28909	60287	0.25%	0.118%
8	8.178	896	909	911	rBV	270609	452385	1.91%	0.884%
9	8.214	911	915	922	rVB	386009	655378	2.76%	1.280%
10	8.601	967	981	992	rVV	573855	1055770	4.45%	2.062%
11	8.801	1008	1015	1026	rVB6	23543	65932	0.28%	0.129%
12	8.913	1026	1034	1037	rVV9	19894	51260	0.22%	0.100%
13	8.966	1037	1043	1054	rVB2	55014	130374	0.55%	0.255%
14	9.283	1088	1097	1105	rVB	61411	113043	0.48%	0.221%
15	9.377	1105	1113	1123	rBV7	45076	121669	0.51%	0.238%
16	9.612	1149	1153	1159	rVB7	19876	36637	0.15%	0.072%
17	9.682	1159	1165	1173	rVB4	40179	89406	0.38%	0.175%
18	9.771	1173	1180	1192	rBV3	71830	136510	0.57%	0.267%
19	9.941	1200	1209	1216	rVB2	100276	216801	0.91%	0.424%
20	10.011	1216	1221	1231	rBV4	57876	156748	0.66%	0.306%
21	10.100	1231	1236	1243	rVB3	105977	184307	0.78%	0.360%
22	10.505	1295	1305	1312	rVV7	80516	186662	0.79%	0.365%
23	10.658	1323	1331	1341	rVV9	31902	109894	0.46%	0.215%
24	10.728	1341	1343	1352	rVB7	20363	52241	0.22%	0.102%
25	10.887	1361	1370	1377	rVV10	19212	47786	0.20%	0.093%
26	11.004	1377	1390	1394	rVV	454180	928227	3.91%	1.813%
27	11.051	1394	1398	1406	rVV	585599	1121220	4.72%	2.190%
28	11.128	1406	1411	1414	rVV6	44940	92442	0.39%	0.181%
29	11.175	1414	1419	1426	rVV4	50807	136122	0.57%	0.266%
30	11.257	1426	1433	1439	rVV3	225698	458833	1.93%	0.896%
31	11.328	1439	1445	1456	rVB4	169918	381647	1.61%	0.746%
32	11.469	1463	1469	1478	rBV2	66110	131131	0.55%	0.256%
33	11.721	1506	1512	1517	rBV10	19536	46062	0.19%	0.090%
34	11.886	1525	1540	1552	rBV8	80503	289472	1.22%	0.565%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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	12.009	1556	1561	1571	rBV10	13301	41183	0.17%	0.080%
36	12.209	1584	1595	1606	rVB10	30701	87159	0.37%	0.170%
37	12.397	1621	1627	1636	rVB10	52191	124985	0.53%	0.244%
38	12.514	1641	1647	1652	rBV10	16174	43211	0.18%	0.084%
39	12.603	1657	1662	1665	rBV2	36897	51053	0.22%	0.100%
40	12.650	1665	1670	1676	rVB2	131912	225285	0.95%	0.440%
41	12.796	1685	1695	1700	rBV2	20597	60471	0.25%	0.118%
42	12.861	1700	1706	1713	rVV	99871	185004	0.78%	0.361%
43	12.967	1713	1724	1734	rVV10	38963	110324	0.46%	0.216%
44	13.084	1740	1744	1750	rVB5	23662	34220	0.14%	0.067%
45	13.178	1756	1760	1765	rVB7	18535	36264	0.15%	0.071%
46	13.313	1773	1783	1788	rBV5	26456	74168	0.31%	0.145%
47	13.372	1788	1793	1800	rVB6	44787	89651	0.38%	0.175%
48	13.631	1824	1837	1844	rBV6	56821	168210	0.71%	0.329%
49	13.707	1844	1850	1858	rVV4	80511	163844	0.69%	0.320%
50	13.819	1858	1869	1879	rVV4	133972	352093	1.48%	0.688%
51	13.960	1885	1893	1909	rBV4	67901	261549	1.10%	0.511%
52	14.142	1909	1924	1929	rBV4	40802	133662	0.56%	0.261%
53	14.201	1929	1934	1946	rVB2	128823	238527	1.00%	0.466%
54	14.377	1958	1964	1969	rBV9	21780	49775	0.21%	0.097%
55	14.506	1973	1986	1995	rVB2	118553	257518	1.08%	0.503%
56	14.665	2004	2013	2021	rBV2	348482	683605	2.88%	1.335%
57	14.806	2027	2037	2048	rBV	723572	1252089	5.27%	2.446%
58	14.900	2048	2053	2060	rVV	40700	109844	0.46%	0.215%
59	14.959	2060	2063	2075	rVB2	49295	111740	0.47%	0.218%
60	15.088	2076	2085	2092	rBV2	366243	858538	3.62%	1.677%
61	15.153	2092	2096	2100	rVV3	152318	271114	1.14%	0.530%
62	15.205	2100	2105	2108	rVV	98544	155353	0.65%	0.303%
63	15.252	2108	2113	2121	rVB4	113106	226901	0.96%	0.443%
64	15.329	2121	2126	2134	rBV7	71715	117775	0.50%	0.230%
65	15.417	2134	2141	2152	rVB2	539641	1243242	5.24%	2.429%
66	15.546	2162	2163	2174	rVB2	18450	45137	0.19%	0.088%
67	15.746	2191	2197	2200	rBV8	26364	55910	0.24%	0.109%
68	15.799	2200	2206	2221	rVB2	152841	380179	1.60%	0.743%
69	15.946	2224	2231	2250	rBV	1351639	2491071	10.49%	4.866%
70	16.087	2252	2255	2261	rVB8	22130	36825	0.16%	0.072%
71	16.251	2277	2283	2292	rVB10	30731	74849	0.32%	0.146%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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	16.492	2318	2324	2326	rBV7	22237	39725	0.17%	0.078%
73	16.533	2326	2331	2339	rVB3	101259	188022	0.79%	0.367%
74	16.704	2358	2360	2370	rVB3	17854	42525	0.18%	0.083%
75	16.827	2376	2381	2386	rBV2	54187	80543	0.34%	0.157%
76	16.892	2386	2392	2396	rBV9	16114	35555	0.15%	0.069%
77	17.221	2438	2448	2465	rBV	13982121	23743883	100.00%	46.384%
78	17.344	2465	2469	2473	rVB7	56261	92248	0.39%	0.180%
79	17.556	2499	2505	2511	rBV2	919814	1464722	6.17%	2.861%
80	17.661	2517	2523	2531	rVB	171091	296854	1.25%	0.580%
81	17.767	2539	2541	2546	rVB6	23301	34831	0.15%	0.068%
82	17.873	2555	2559	2567	rVB10	23431	53955	0.23%	0.105%
83	17.967	2567	2575	2581	rBV	110287	196978	0.83%	0.385%
84	18.360	2637	2642	2648	rBV10	27240	64173	0.27%	0.125%
85	18.983	2742	2748	2754	rVB5	38765	67968	0.29%	0.133%
86	19.400	2813	2819	2825	rBV3	49960	94902	0.40%	0.185%
87	19.718	2870	2873	2882	rVB10	31889	69216	0.29%	0.135%
88	19.935	2904	2910	2917	rBV2	196941	319000	1.34%	0.623%
89	20.417	2985	2992	3000	rBV4	68428	188550	0.79%	0.368%
90	21.845	3228	3235	3245	rVB	990037	1722347	7.25%	3.365%
91	21.974	3252	3257	3265	rBV	27696	71281	0.30%	0.139%
92	23.290	3478	3481	3488	rVB9	22578	40847	0.17%	0.080%
93	24.113	3617	3621	3631	rVB9	35169	96652	0.41%	0.189%
94	25.000	3763	3772	3783	rBV2	94342	312764	1.32%	0.611%
95	25.111	3783	3791	3798	rVB9	45843	124745	0.53%	0.244%
96	25.241	3802	3813	3831	rBV	516497	1653308	6.96%	3.230%
97	26.269	3978	3988	3999	rBV6	54729	171191	0.72%	0.334%
98	27.673	4221	4227	4238	rVB6	48210	138280	0.58%	0.270%
99	29.348	4506	4512	4514	rBV7	26188	49519	0.21%	0.097%
100	31.410	4856	4863	4873	rVB7	22805	77323	0.33%	0.151%

Sum of corrected areas: 51190361

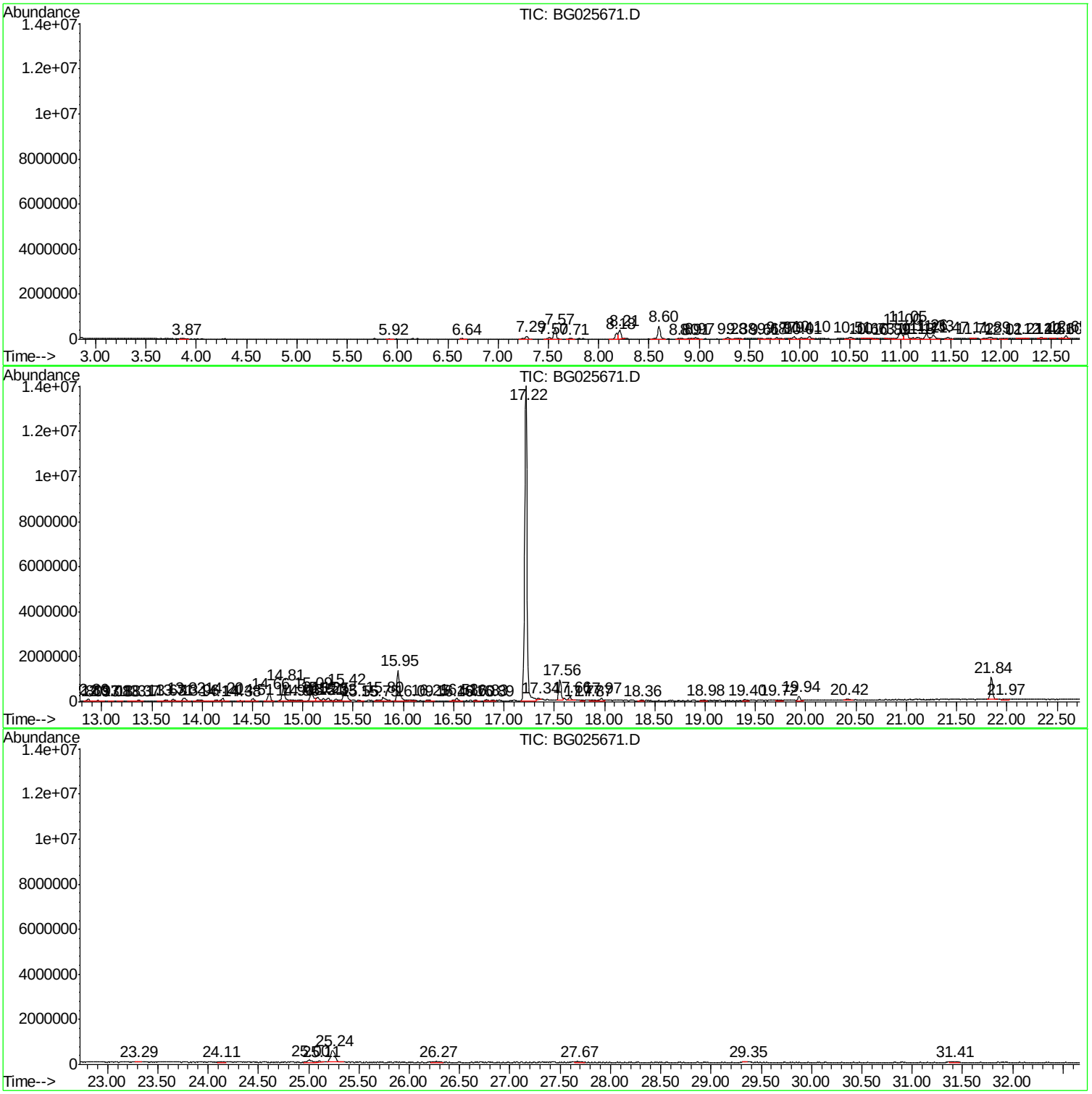
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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 : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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

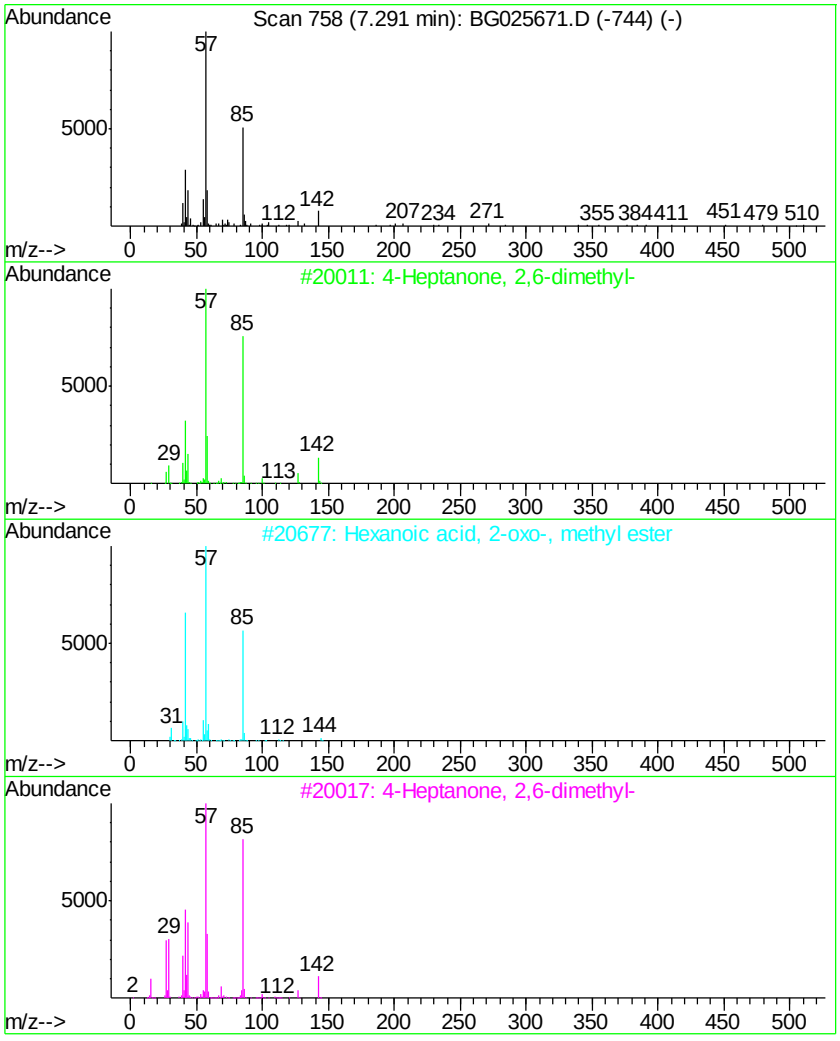
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

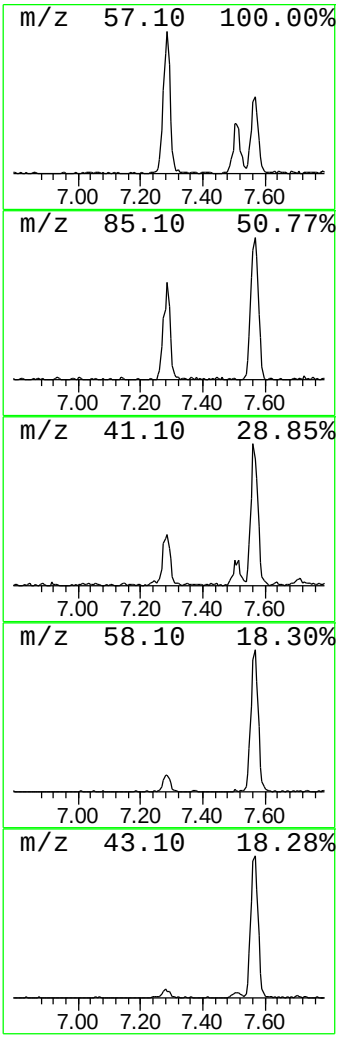
****

Peak Number 1 4-Heptanone, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.29	10.79 ng/ul	244132	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	64
2	Hexanoic acid, 2-oxo-, methyl ester	144	C7H12O3	006395-83-1	59
3	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	50
4	5-Tridecanone	198	C13H26O	030692-16-1	39
5	Pentanethioic acid, S-ethyl ester	146	C7H14OS	002432-92-0	38





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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

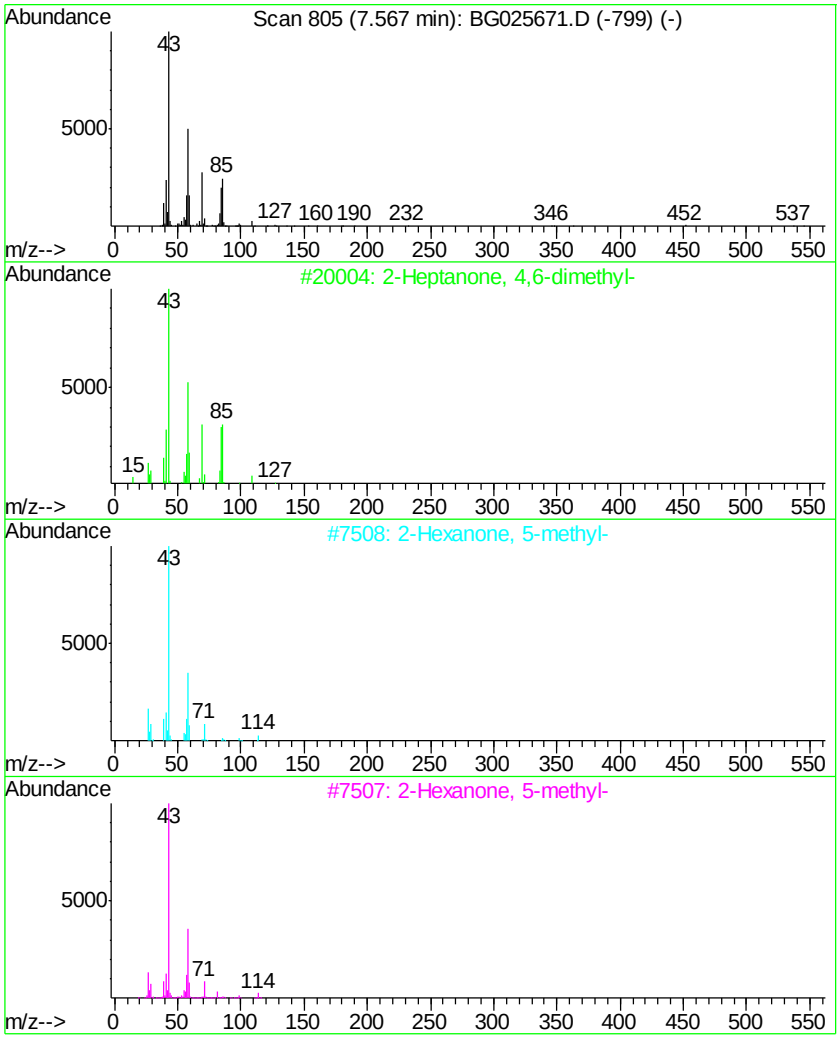
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

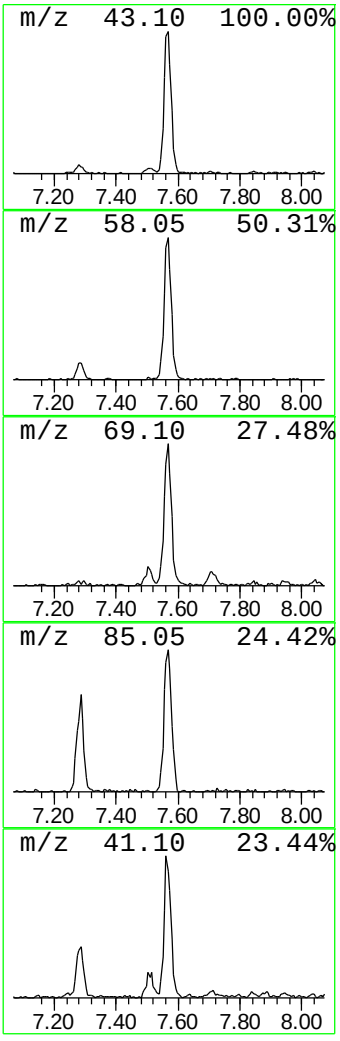
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	33.64 ng/ul	760824	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	83
2	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	50
3	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	47
4	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	40
5	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	38





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

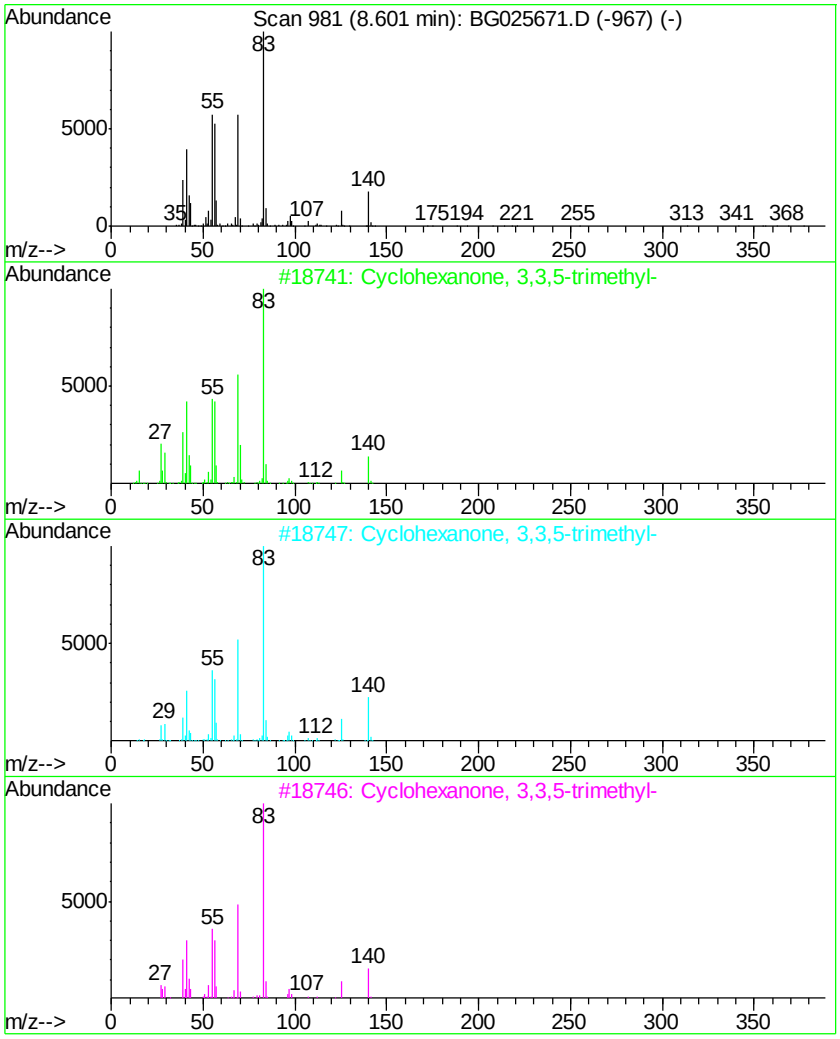
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

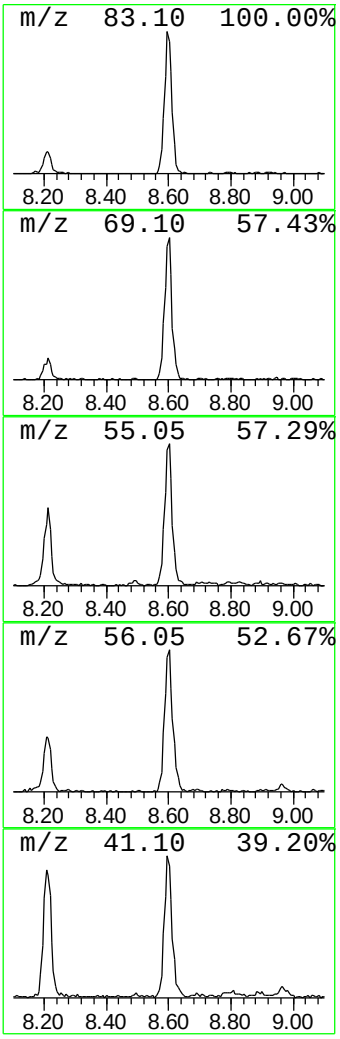
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclohexanone, 3,3,5-trimet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	46.68 ng/ul	1055770	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	94
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	76
4	1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	53
5	2,4,6-Trimethyl-3-heptene	140	C10H20	1000113-56-9	52





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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

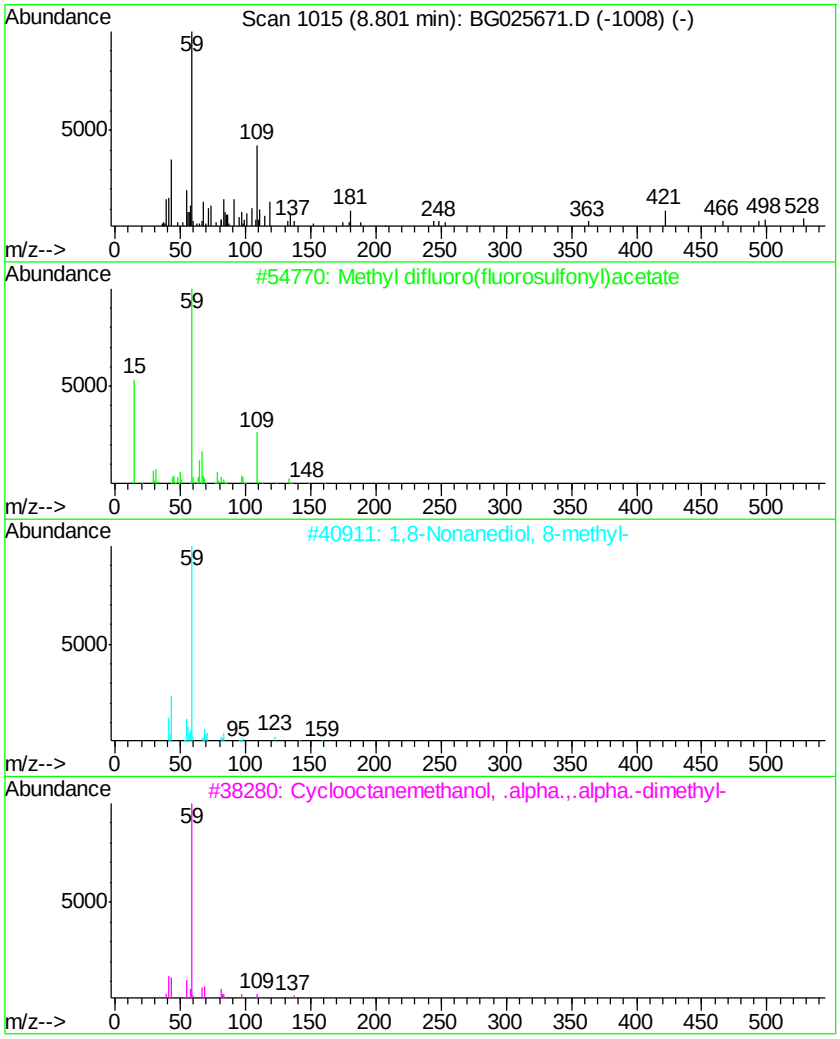
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

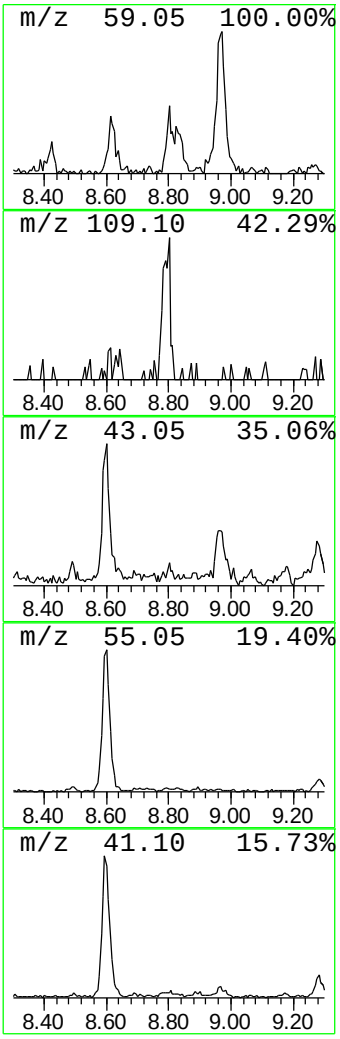
****

Peak Number 4 unknown-01 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	2.91 ng/ul	65932	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl difluoro(fluorosulfonyl)a...	192	C3H3F3O4S	000680-15-9	36
2	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	35
3	Cyclooctanemethanol, .alpha.,.al...	170	C11H22O	016624-06-9	32
4	Butanamide	87	C4H9NO	000541-35-5	25
5	Cyclohexane, (ethoxymethoxy)-	158	C9H18O2	054699-29-5	25





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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

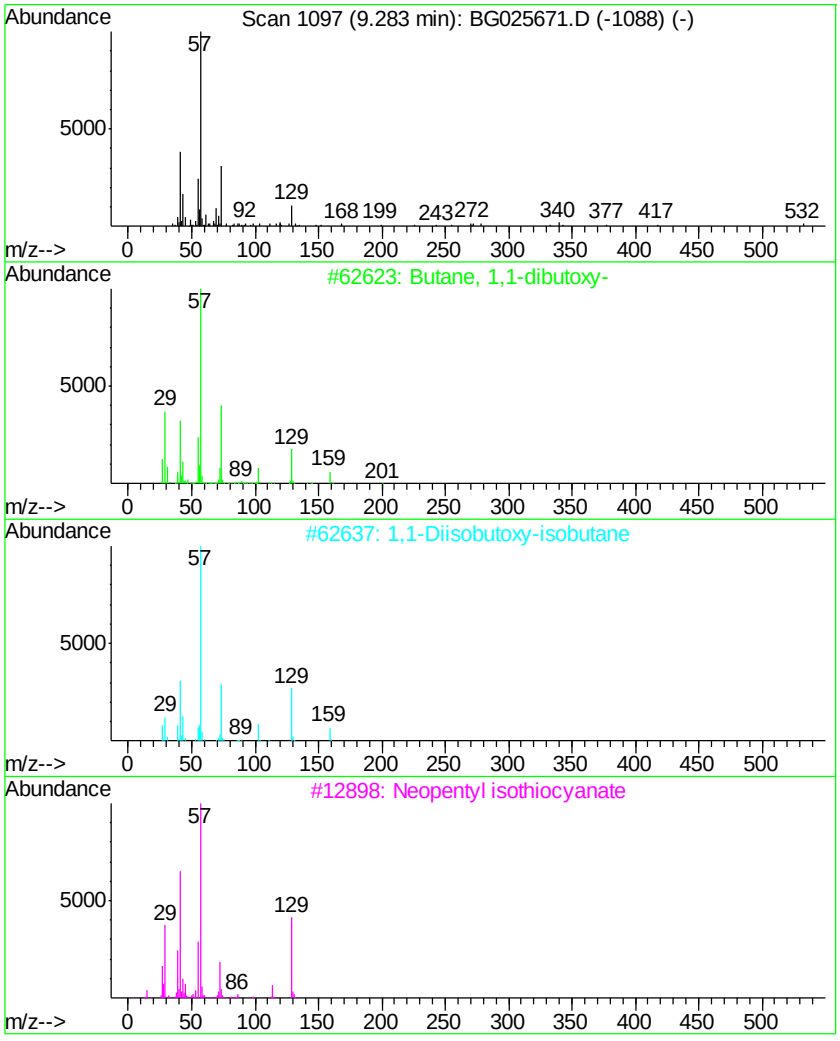
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

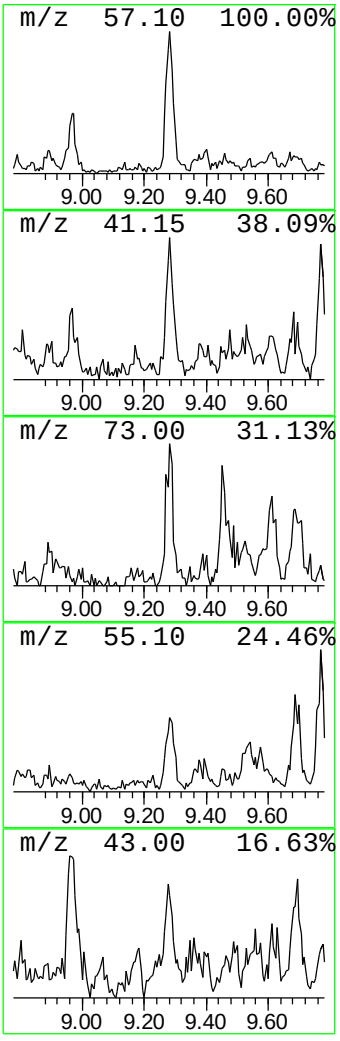
****

Peak Number 5 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	5.00 ng/ul	113043	1,4-Dichlorobenzene-d4	8.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1,1-dibutoxy-	202	C12H26O2	005921-80-2	42
2	1,1-Diisobutoxy-isobutane	202	C12H26O2	013262-24-3	42
3	Neopentyl isothiocyanate	129	C6H11NS	000597-97-7	38
4	Formic acid, neopentyl ester	116	C6H12O2	1000367-90-3	37
5	1-Butoxy-1-isobutoxy-butane	202	C12H26O2	020266-12-0	28





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Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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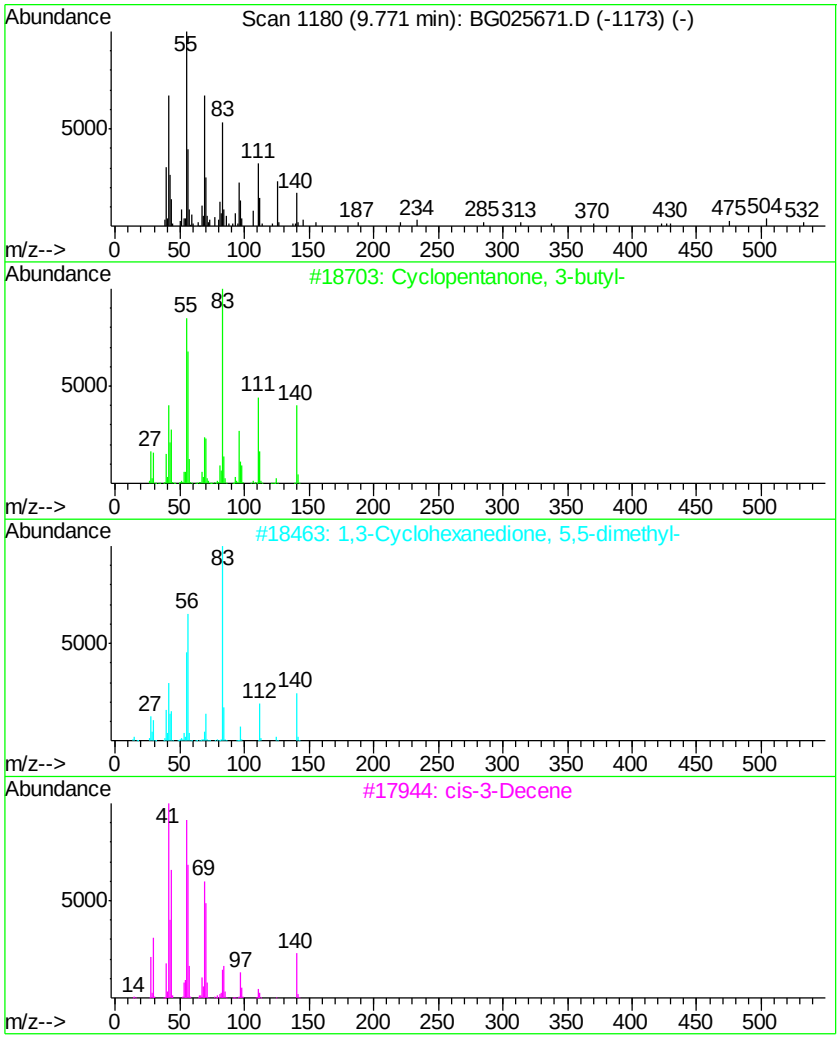
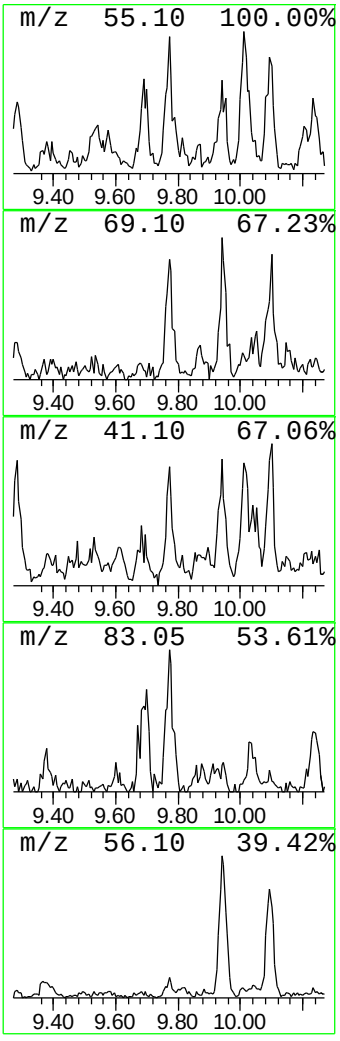
Peak Number 6 Cyclopentanone, 3-butyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.77	2.94 ng/ul	136510	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanone, 3-butyl-	140	C9H16O	057283-81-5	52
2	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	38
3	cis-3-Decene	140	C10H20	019398-86-8	35
4	Cyclopentane, propyl-	112	C8H16	002040-96-2	30
5	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	27





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Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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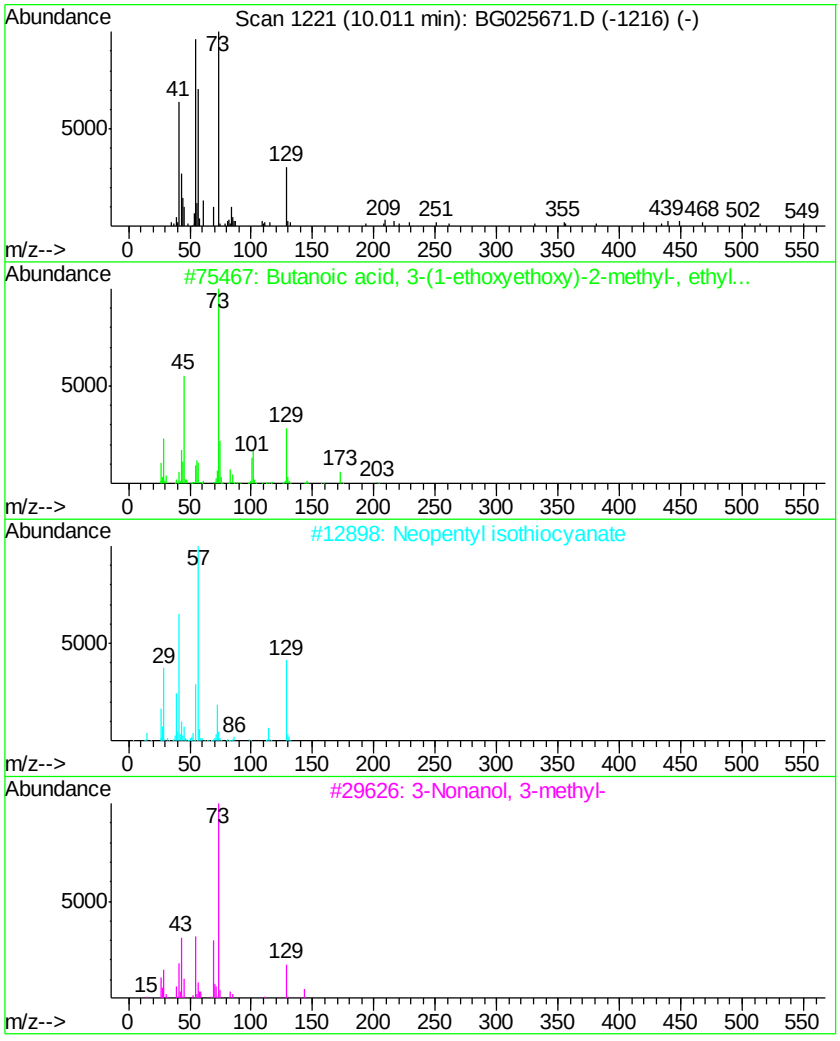
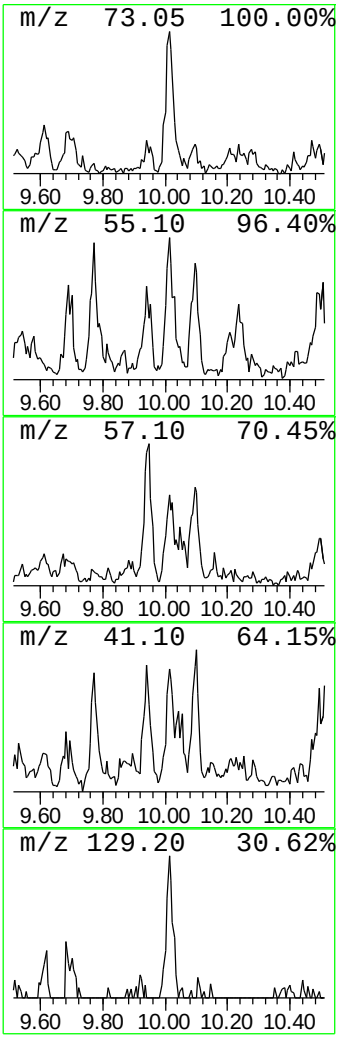
Peak Number 7 unknown-03 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.01	3.38 ng/ul	156748	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 3-(1-ethoxyethoxy...	218	C11H22O4	086845-49-0	33
2	Neopentyl isothiocyanate	129	C6H11NS	000597-97-7	12
3	3-Nonanol, 3-methyl-	158	C10H22O	021078-72-8	9
4	1,4-Anhydro-d-galactitol	164	C6H12O5	1000129-01-0	9
5	1,1-Dibutoxy-isobutane	202	C12H26O2	013112-62-4	9





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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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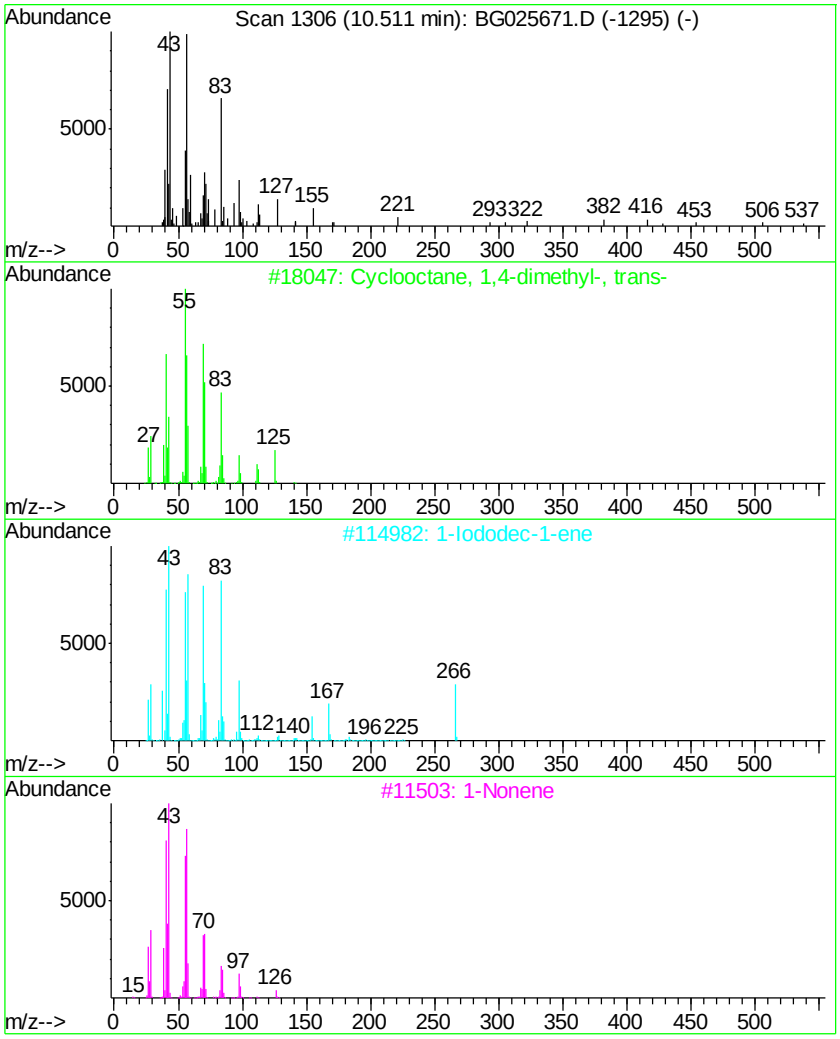
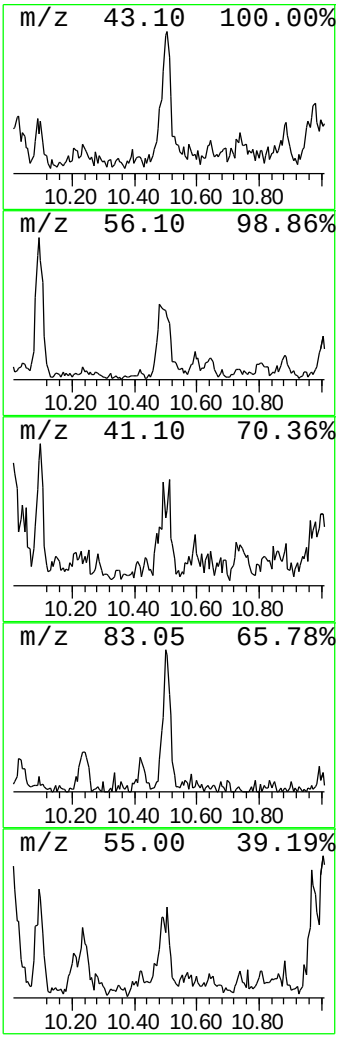
Peak Number 8 (DEL) Alkane: Cyclic10.51 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	4.02 ng/ul	186662	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctane, 1,4-dimethyl-, trans-	140	C10H20	013151-98-9	72
2	1-Iododec-1-ene	266	C10H19I	1000192-68-7	43
3	1-Nonene	126	C9H18	000124-11-8	38
4	Cyclohexane, ethyl-	112	C8H16	001678-91-7	30
5	1-Heptene, 2-methyl-	112	C8H16	015870-10-7	30





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Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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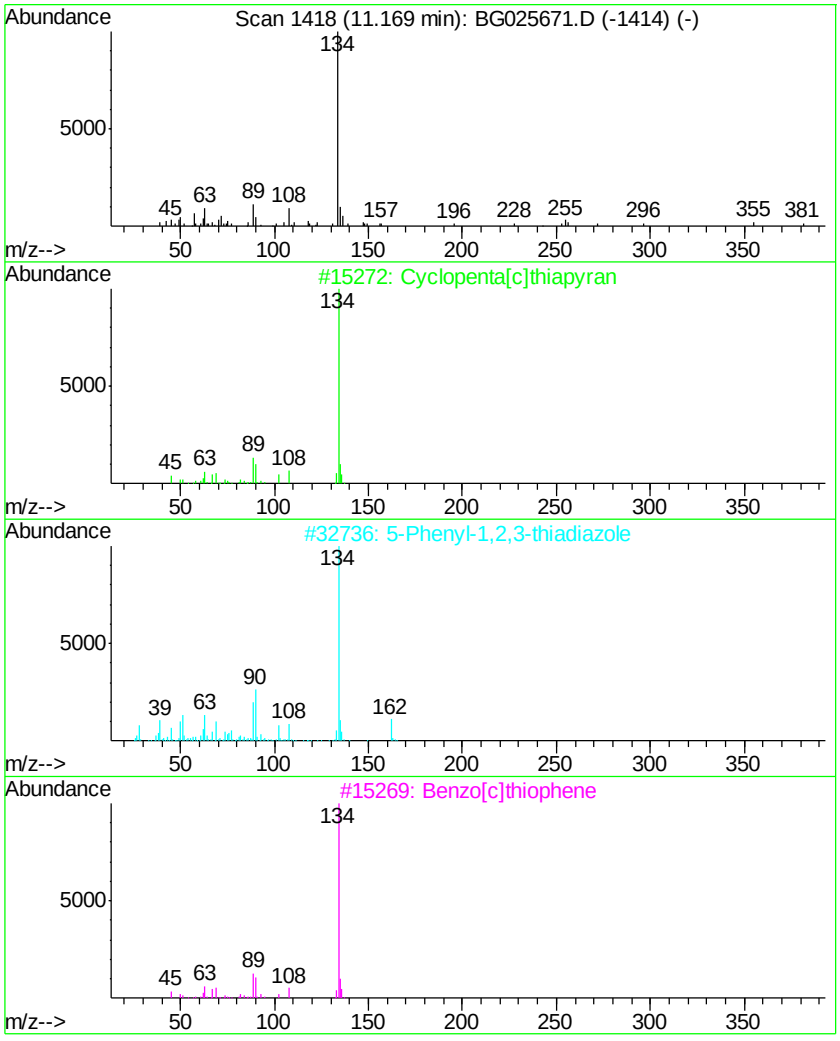
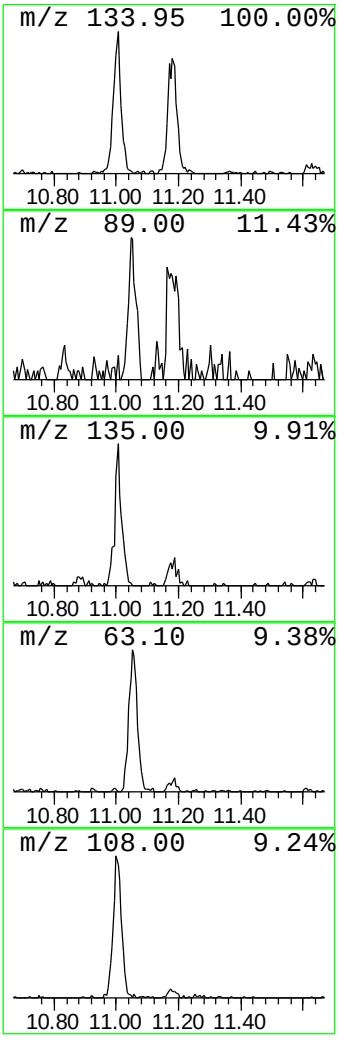
Peak Number 9 Cyclopenta[c]thiapyran Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.17	2.93 ng/ul	136122	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	80
2	5-Phenyl-1,2,3-thiadiazole	162	C8H6N2S	018212-29-8	72
3	Benzo[c]thiophene	134	C8H6S	000270-82-6	72
4	Benzo[b]thiophene	134	C8H6S	000095-15-8	72
5	3-Bromomethylbenzamide	213	C8H8BrNO	1000211-93-8	64





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Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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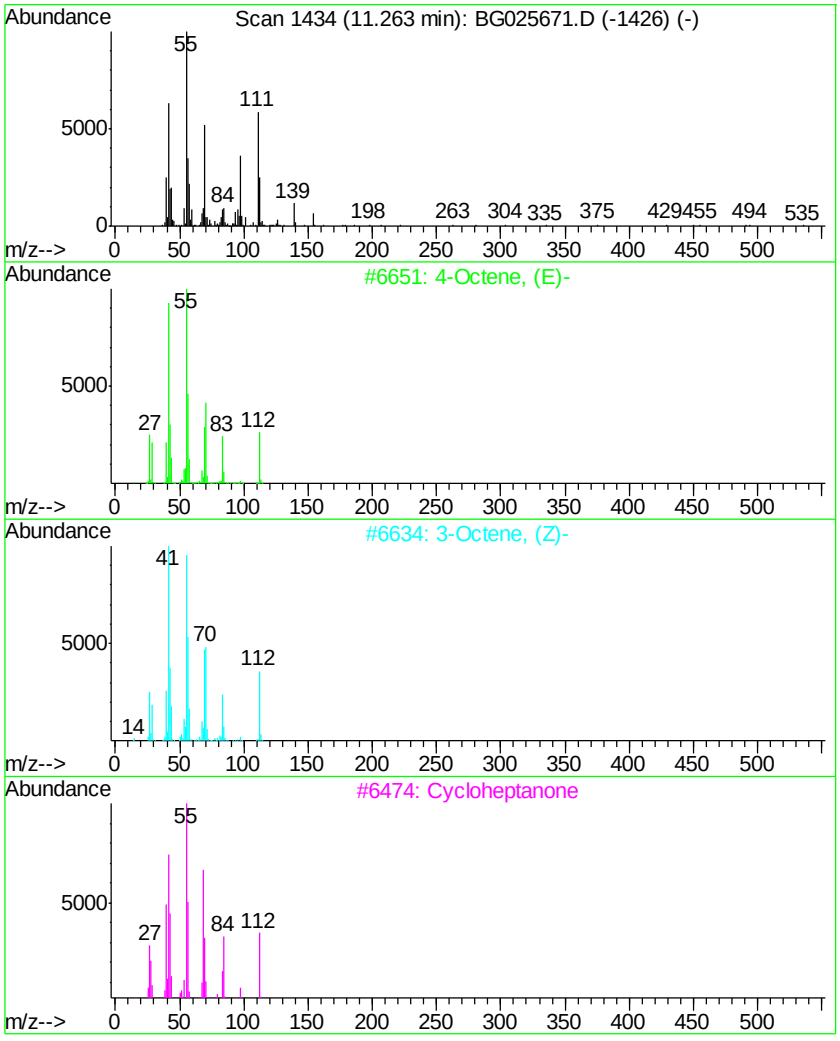
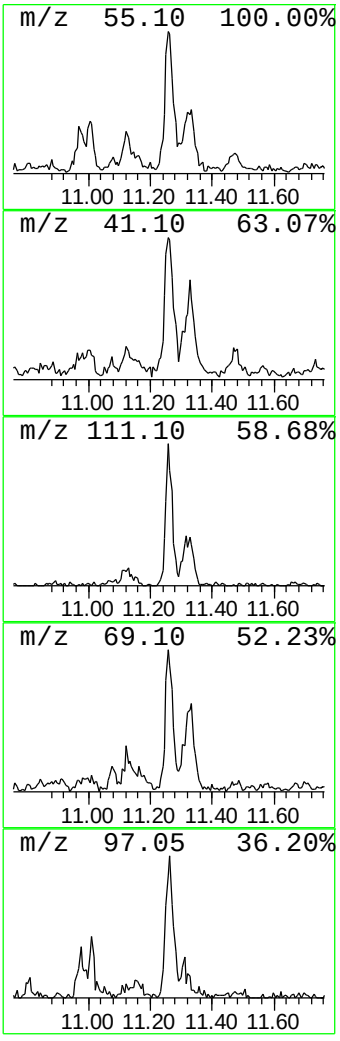
Peak Number 10 (DEL) Alkane: Cyclic11.26 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	9.89 ng/ul	458833	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Octene, (E)-	112	C8H16	014850-23-8	43
2	3-Octene, (Z)-	112	C8H16	014850-22-7	38
3	Cycloheptanone	112	C7H12O	000502-42-1	35
4	4-Decene, 6-methyl-, (E)-	154	C11H22	036229-57-9	30
5	2-Octene, (Z)-	112	C8H16	007642-04-8	27





Library Search Compound Report

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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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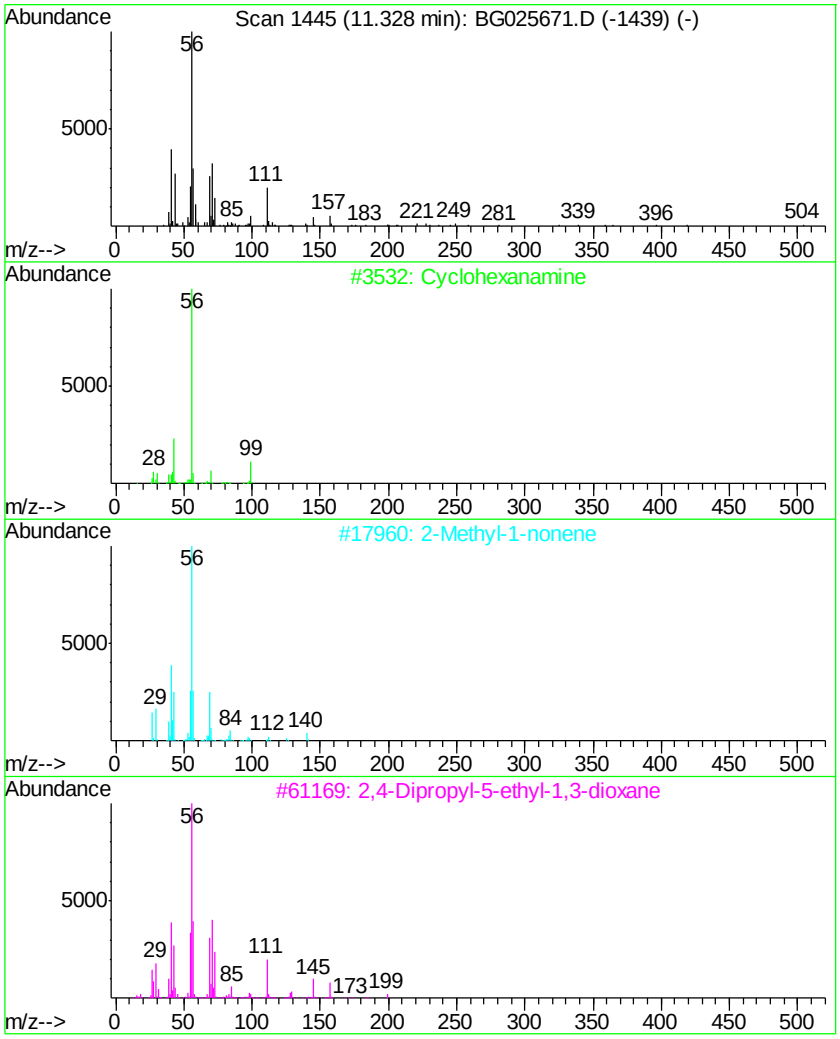
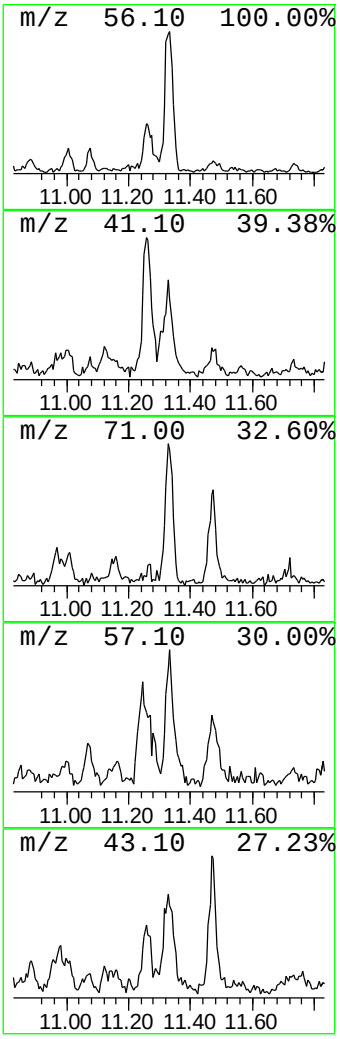
Peak Number 11 unknown-04 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.33	8.22 ng/ul	381647	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanamine	99	C6H13N	000108-91-8	38
2	2-Methyl-1-nonene	140	C10H20	002980-71-4	38
3	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	37
4	Cyclopentanone, 2,2,5,5-tetramet...	140	C9H16O	004541-35-9	32
5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	28





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Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

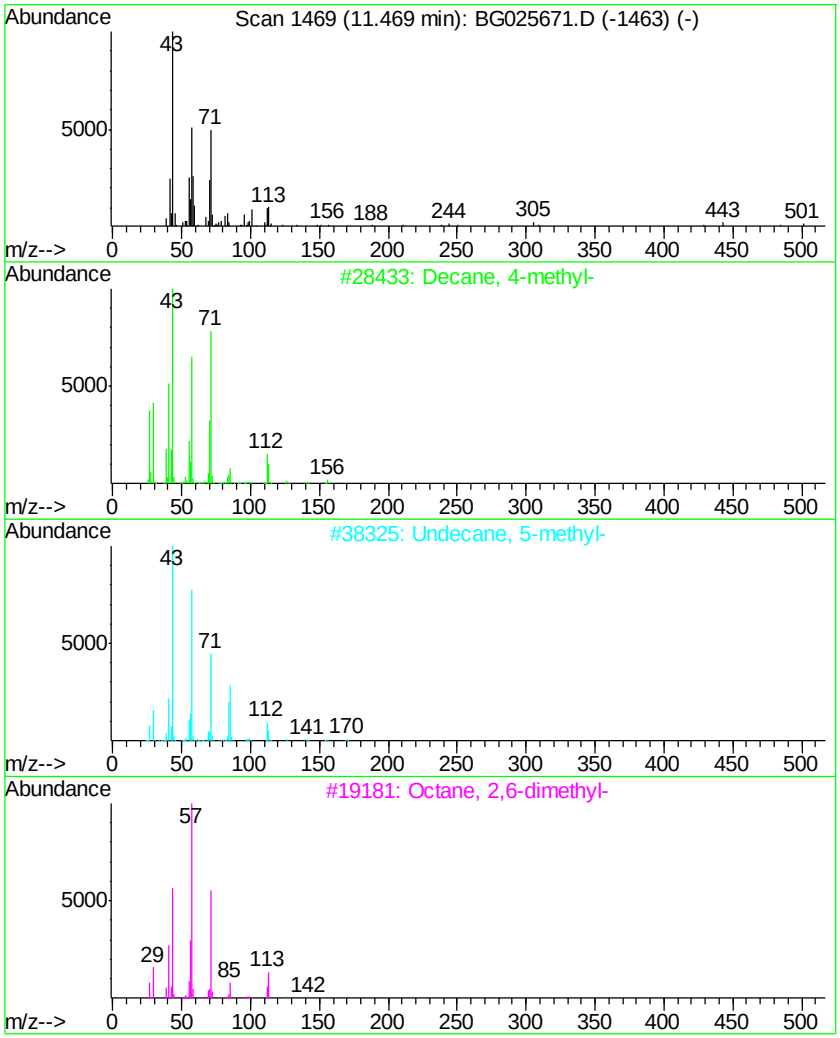
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG011817.M
Quant Title : SVOA CALIBRATION

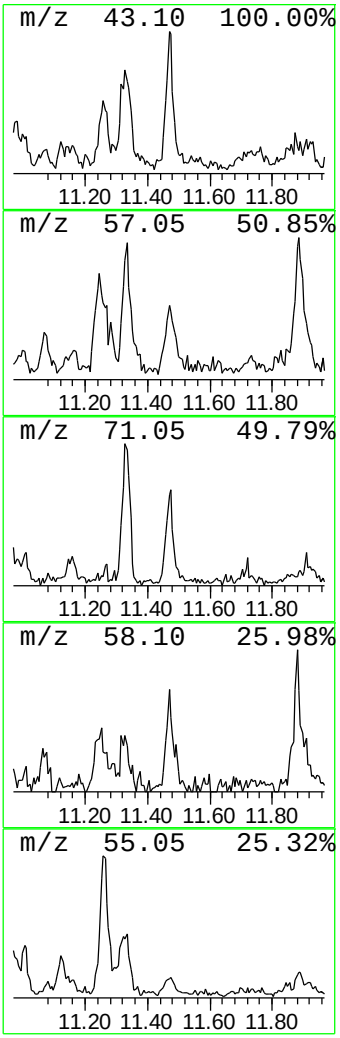
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	2.83 ng/ul	131131	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-	156	C11H24	002847-72-5	58
2	Undecane, 5-methyl-	170	C12H26	001632-70-8	47
3	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	47
4	Heptane, 5-ethyl-2-methyl-	142	C10H22	013475-78-0	47
5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	47





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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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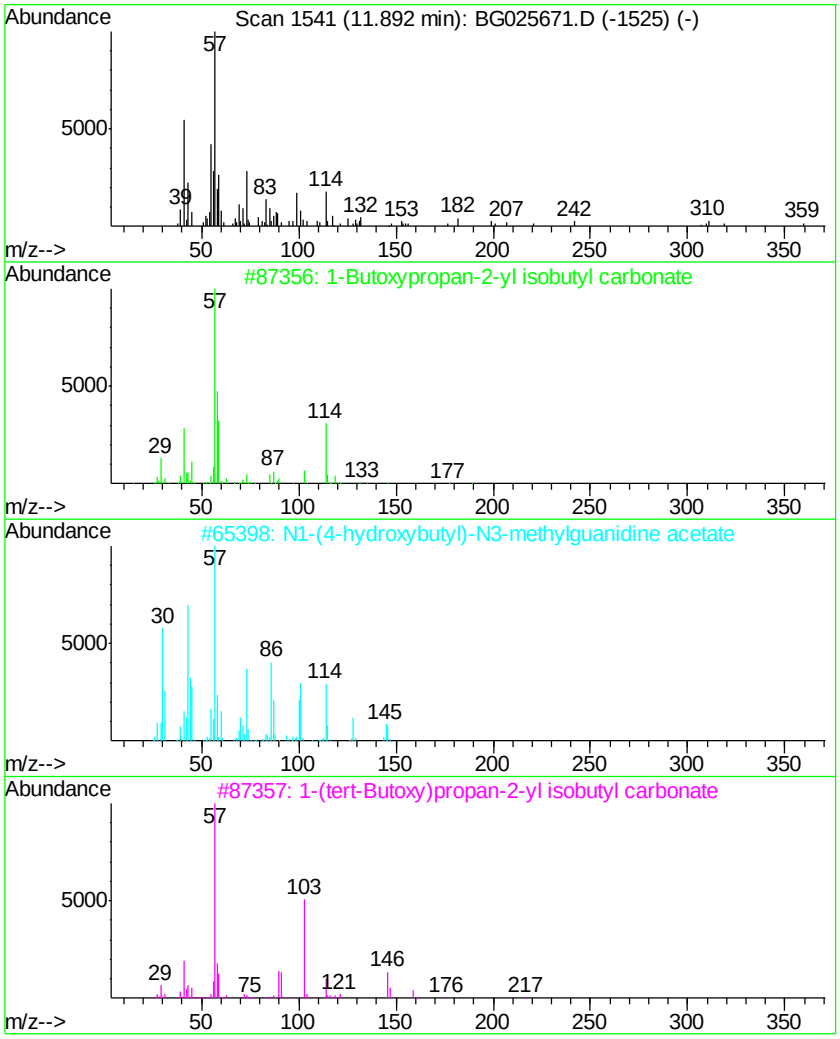
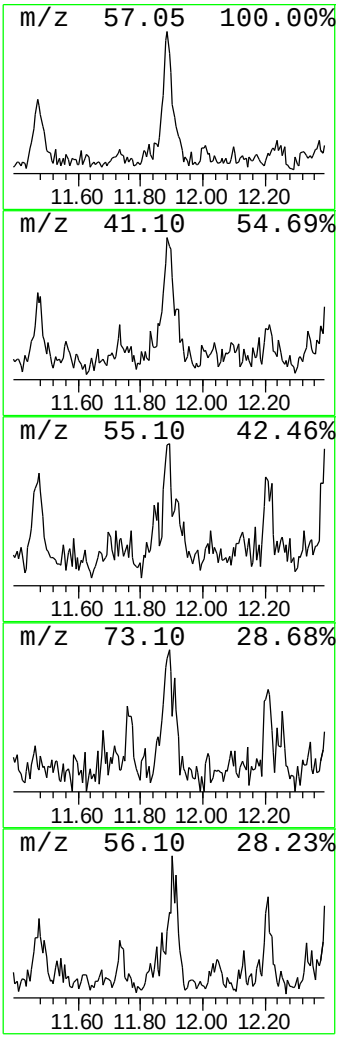
Peak Number 13 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	6.24 ng/ul	289472	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butoxypropan-2-yl isobutyl car...	232	C12H24O4	1000378-27-6	37
2	N1-(4-hydroxybutyl)-N3-methylgua...	205	C8H19N3O3	1000188-13-5	25
3	1-(tert-Butoxy)propan-2-yl isobu...	232	C12H24O4	1000378-31-8	25
4	2-(2-Ethoxyethoxy)ethyl 2-methyl...	218	C11H22O4	1000366-97-5	17
5	Butanoic acid, 2-methyl-, methyl...	116	C6H12O2	000868-57-5	12





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Misc :
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Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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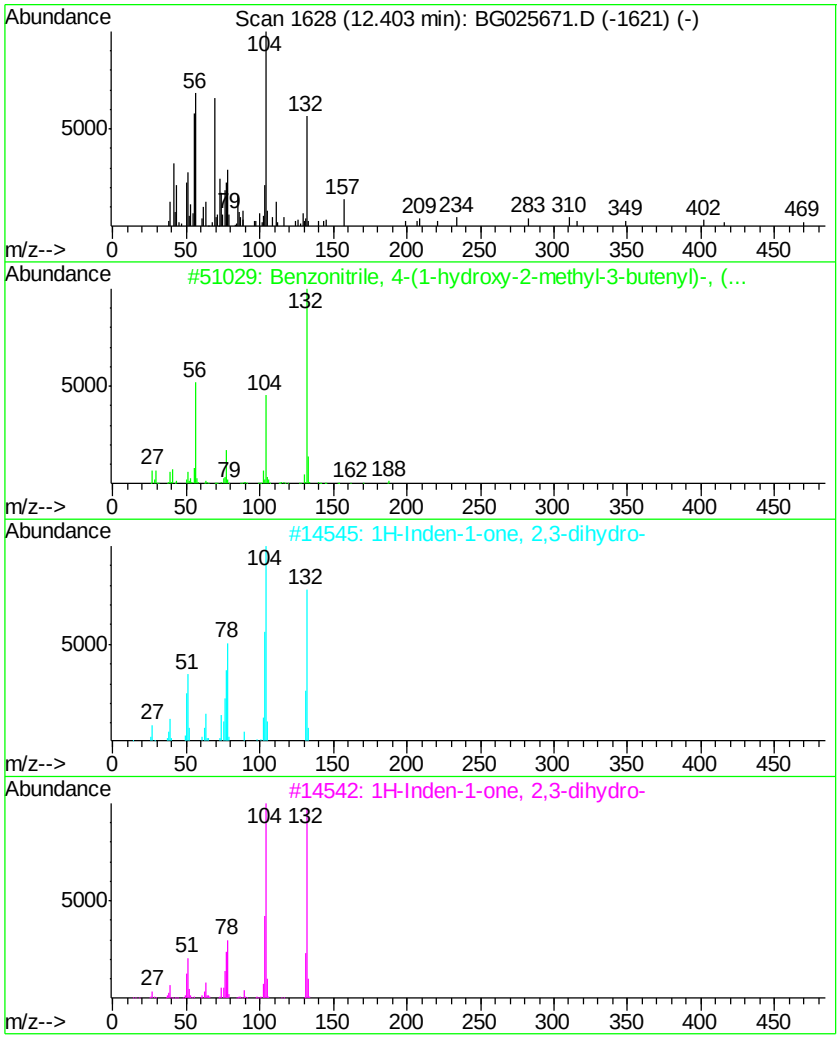
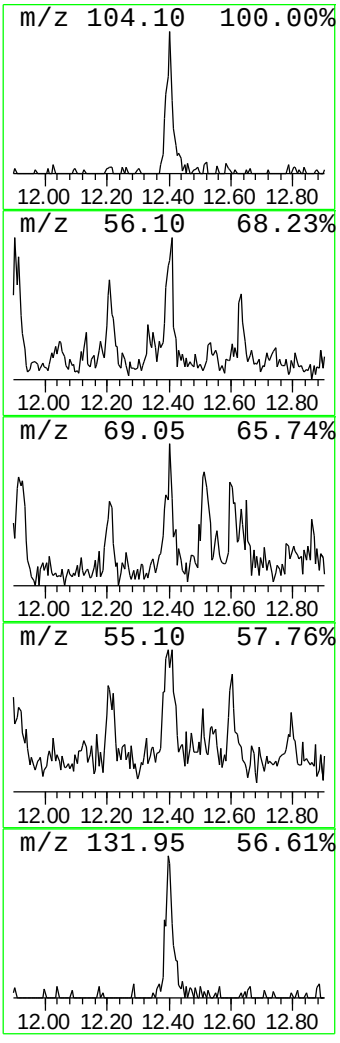
Peak Number 14 unknown-06 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	2.69 ng/ul	124985	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzonitrile, 4-(1-hydroxy-2-met...	187	C12H13NO	083173-80-2	43
2	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	38
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	38
4	Propanedinitrile, 2-(1-cycloprop...	132	C8H8N2	017407-30-6	38
5	Indan, epoxide	132	C9H8O	000768-22-9	38





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Sample : I1387-07DL 10X
Misc :
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Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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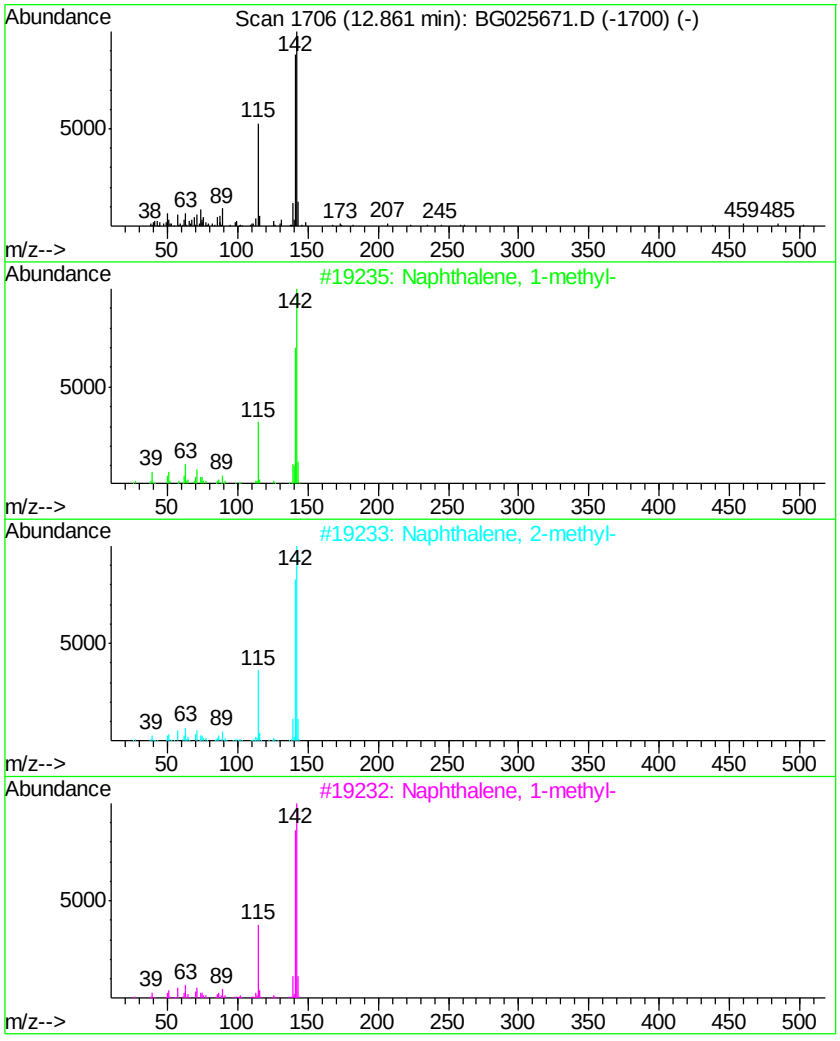
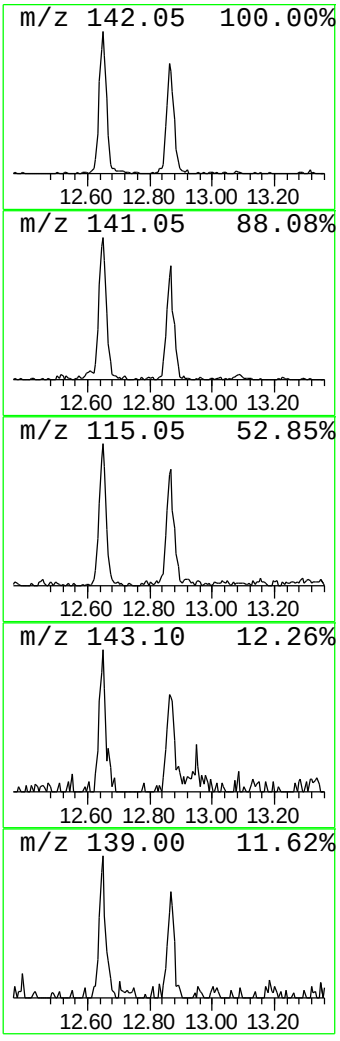
Peak Number 15 Naphthalene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.86	3.99 ng/ul	185004	Naphthalene-d8	11.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91





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Quant Title : SVOA CALIBRATION

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

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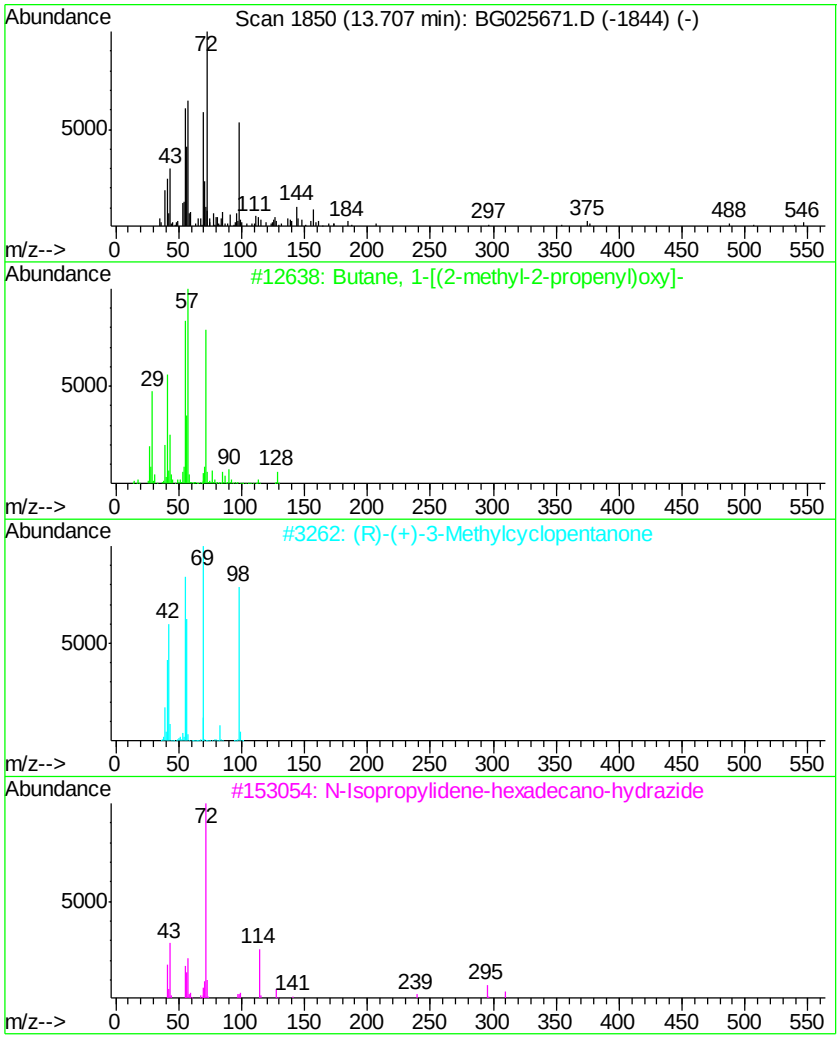
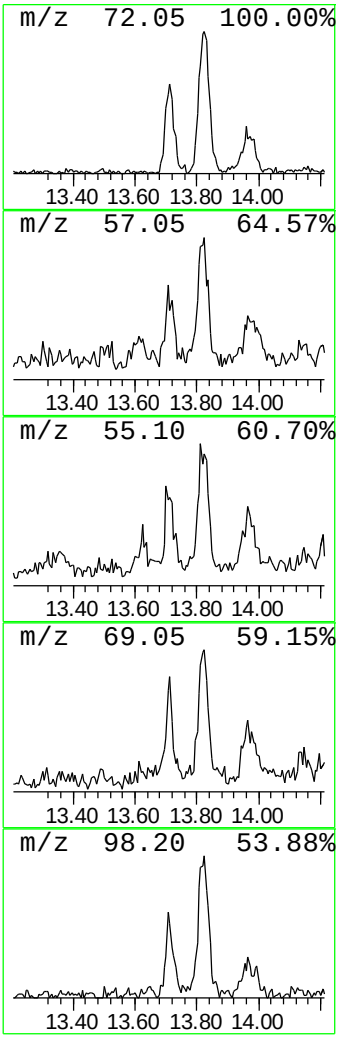
Peak Number 16 Butane, 1-[(2-methyl-2-prop... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	2.62 ng/ul	163844	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1-[(2-methyl-2-propenyl)...	128	C8H16O	053907-74-7	50
2	(R)-(+)-3-Methylcyclopentanone	98	C6H10O	006672-30-6	35
3	N-Isopropylidene-hexadecano-hydr...	310	C19H38N2O	093612-49-8	32
4	Cyclopentanone, 3-methyl-	98	C6H10O	001757-42-2	27
5	Cyclopentanone, 3-methyl-	98	C6H10O	001757-42-2	27





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Data File : BG025671.D
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Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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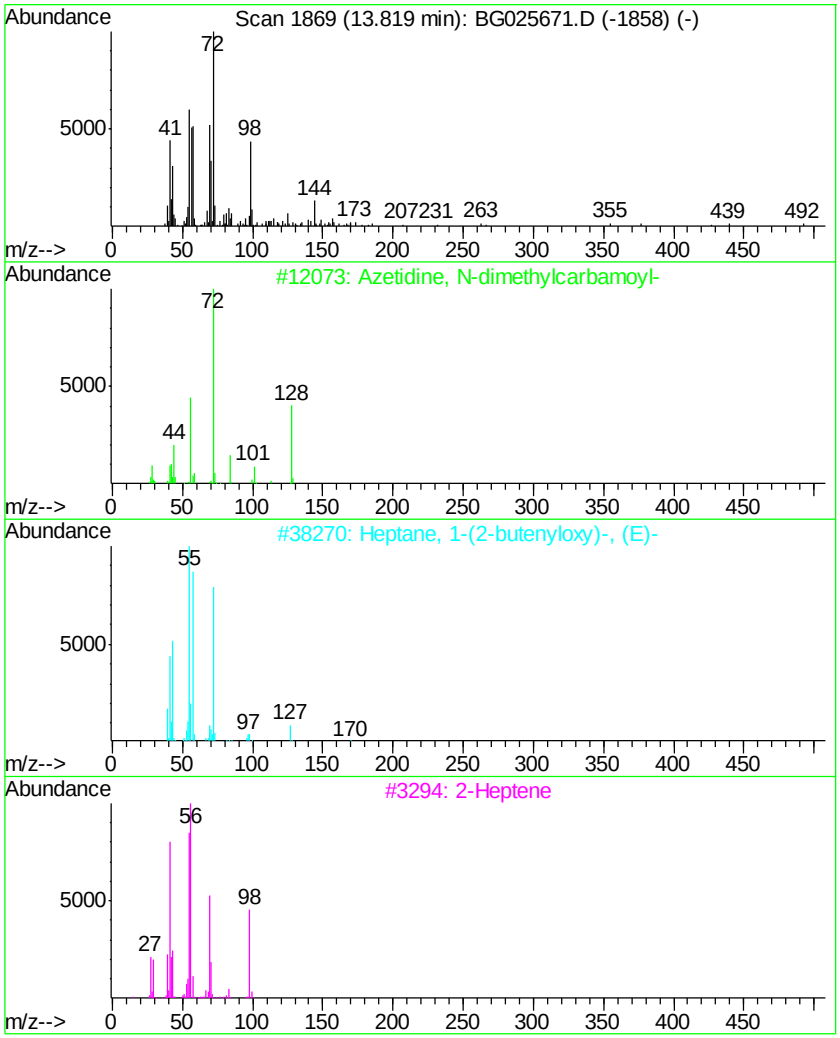
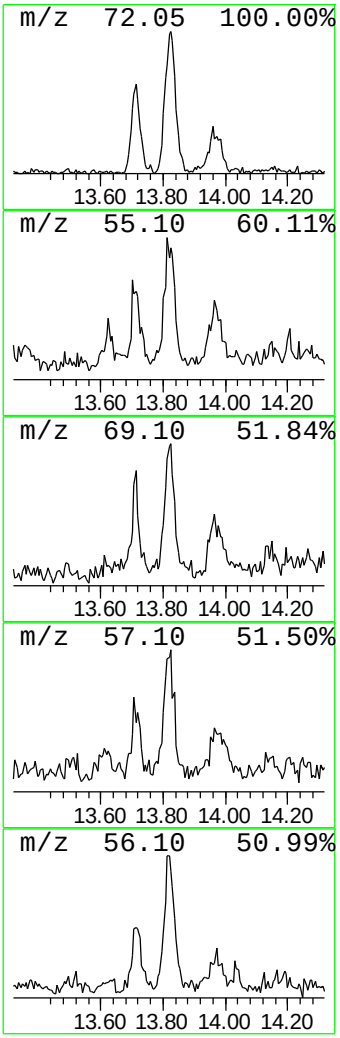
Peak Number 17 unknown-07 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	5.62 ng/ul	352093	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Azetidine, N-dimethylcarbamoyl-	128	C6H12N2O	057031-53-5	49
2	Heptane, 1-(2-butenyloxy)-, (E)-	170	C11H22O	056052-77-8	43
3	2-Heptene	98	C7H14	000592-77-8	30
4	2-Heptene, (E)-	98	C7H14	014686-13-6	30
5	2-Hexene, 2-methyl-	98	C7H14	002738-19-4	27





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Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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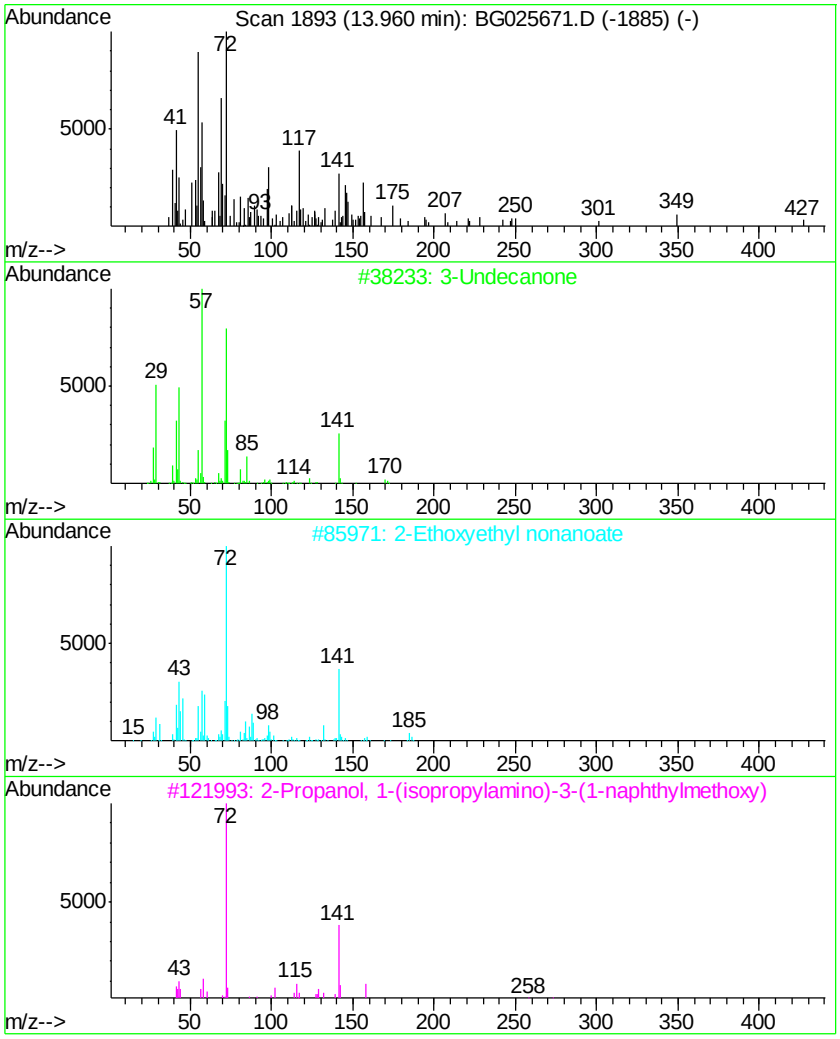
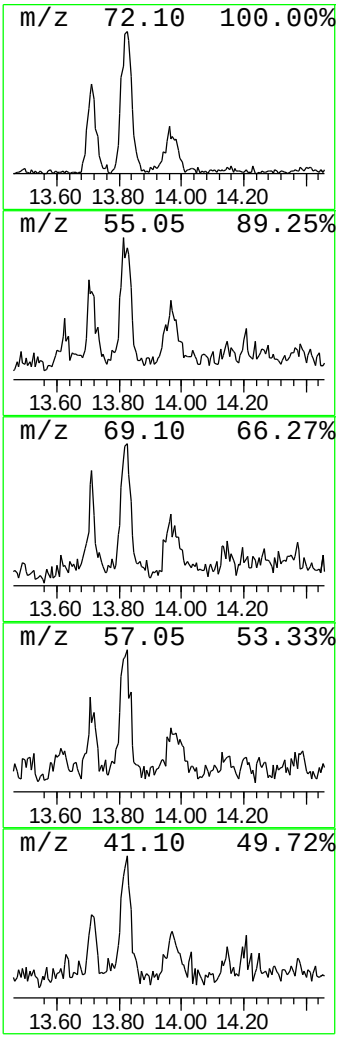
Peak Number 18 unknown-08 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	4.18 ng/ul	261549	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Undecanone	170	C11H22O	002216-87-7	27
2	2-Ethoxyethyl nonanoate	230	C13H26O3	1000378-25-2	16
3	2-Propanol, 1-(isopropylamino)-3...	273	C17H23NO2	019343-21-6	16
4	3,cis-(1,1-dimethylethyl)-4,cis-...	186	C11H22O2	1000215-62-0	16
5	2-Acrylamido-2-methyl-1-propanes...	207	C7H13NO4S	015214-89-8	16





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Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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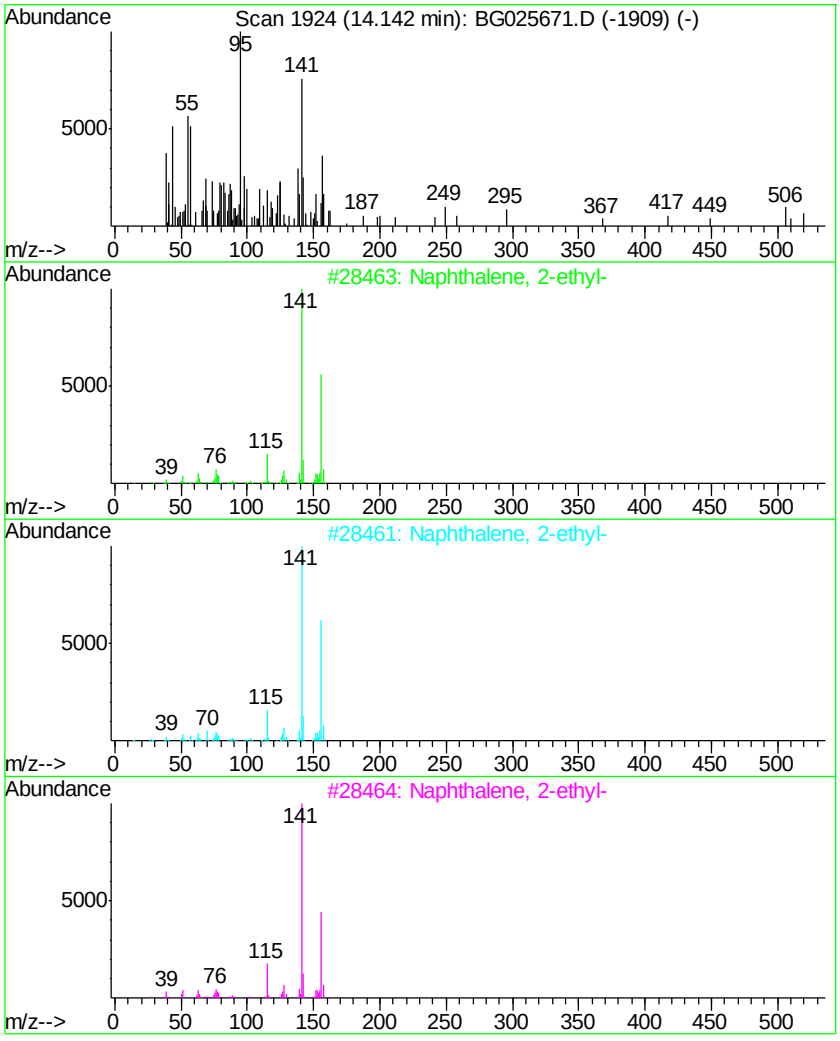
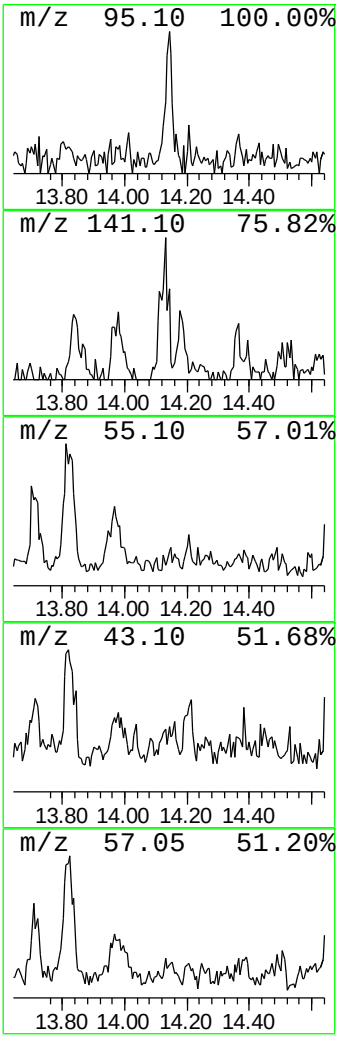
Peak Number 19 unknown-09 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	2.14 ng/ul	133662	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	38
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	35
3	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	35
4	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	30
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	27





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Acq On : 26 Jan 2017 17:55
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Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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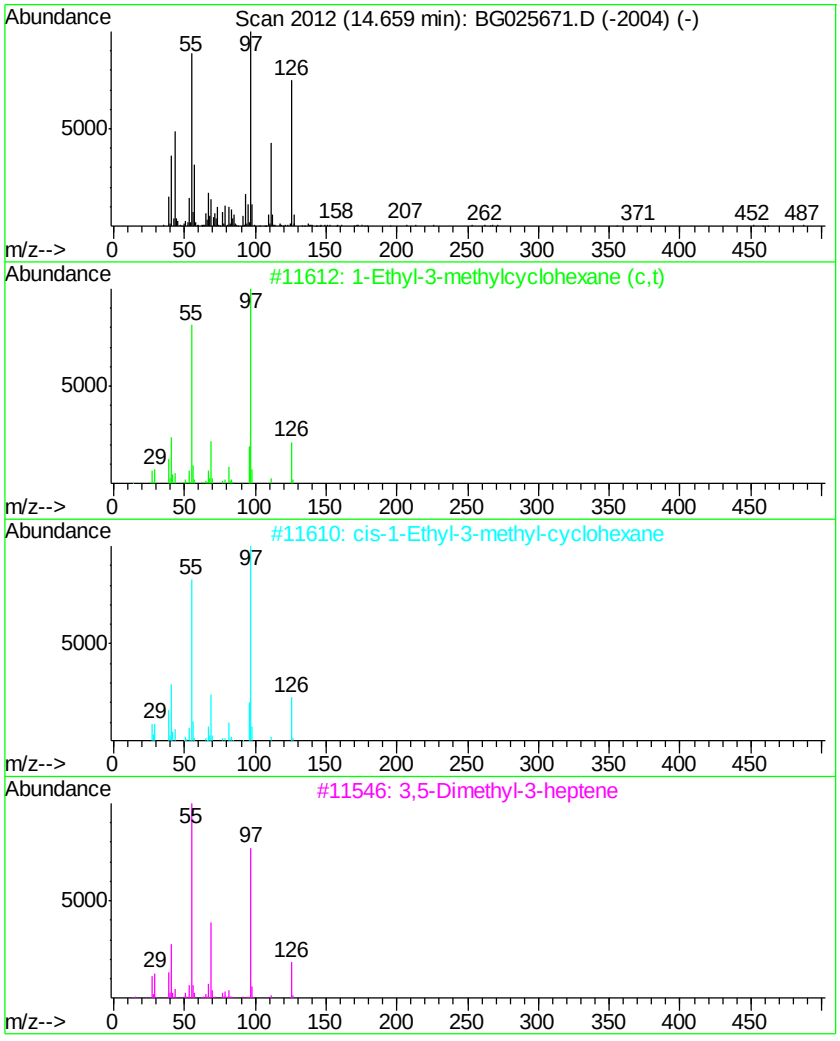
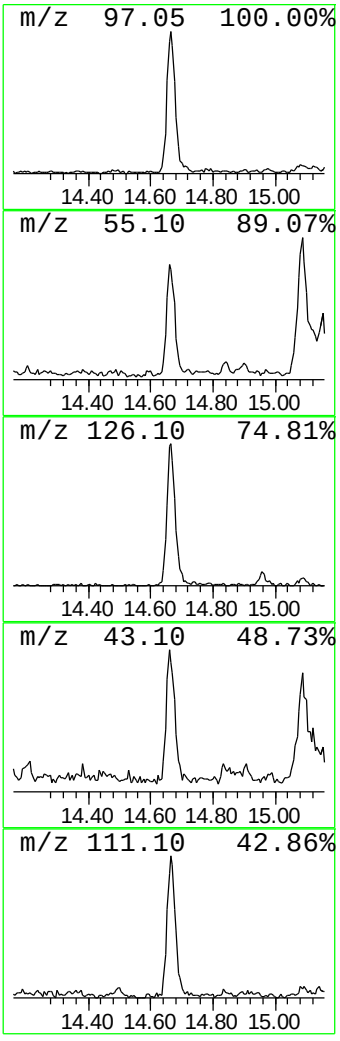
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	10.92 ng/ul	683605	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	64
2	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	59
3	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	59
4	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	59
5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	59





Library Search Compound Report

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Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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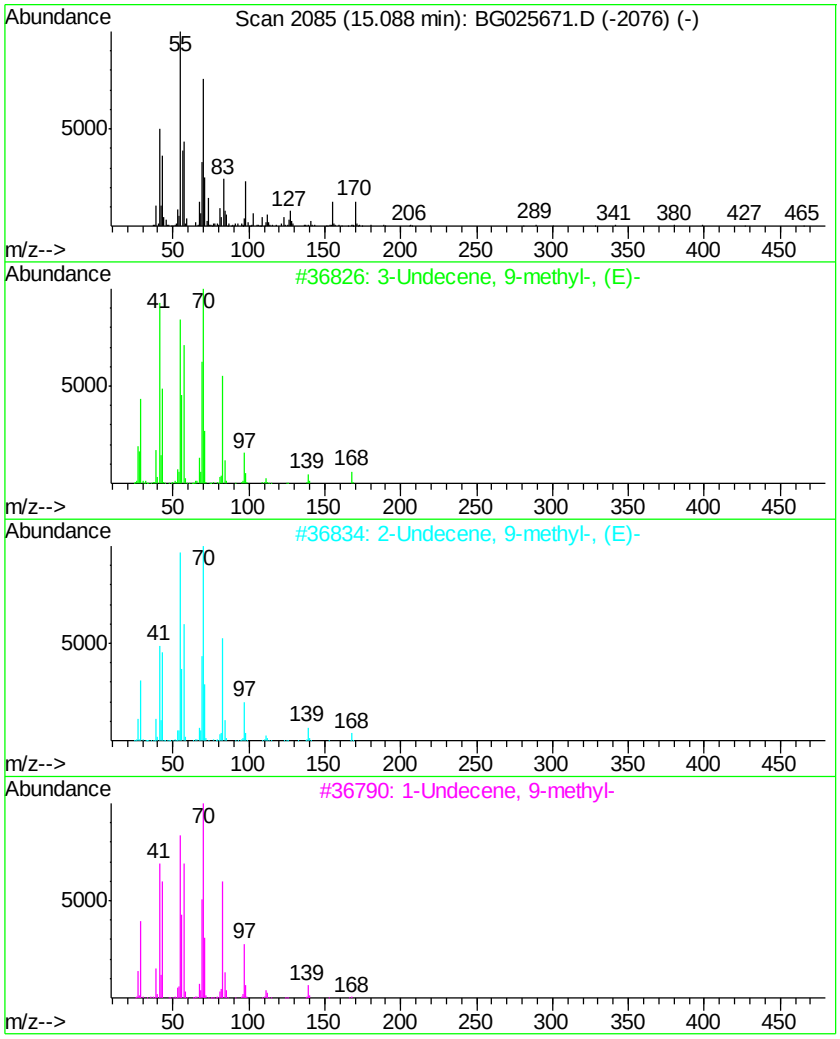
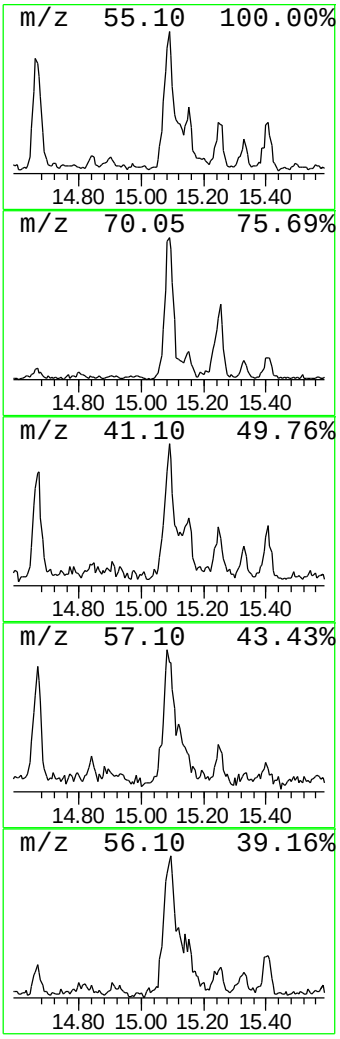
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	13.71 ng/ul	858538	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Undecene, 9-methyl-, (E)-	168	C12H24	074630-54-9	53
2	2-Undecene, 9-methyl-, (E)-	168	C12H24	074630-46-9	53
3	1-Undecene, 9-methyl-	168	C12H24	074630-41-4	50
4	4-Undecene, 9-methyl-, (Z)-	168	C12H24	074630-56-1	50
5	1,2-Cyclohexanediol, trans-	116	C6H12O2	001460-57-7	43





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Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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Quant Title : SVOA CALIBRATION

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

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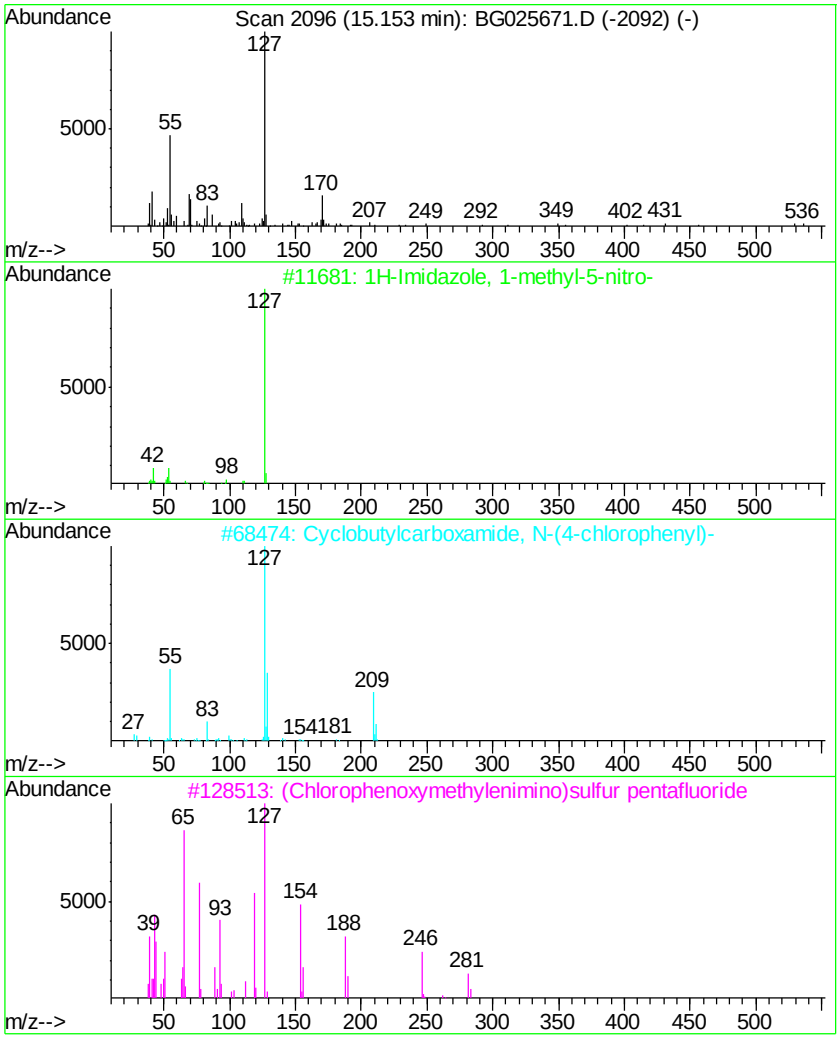
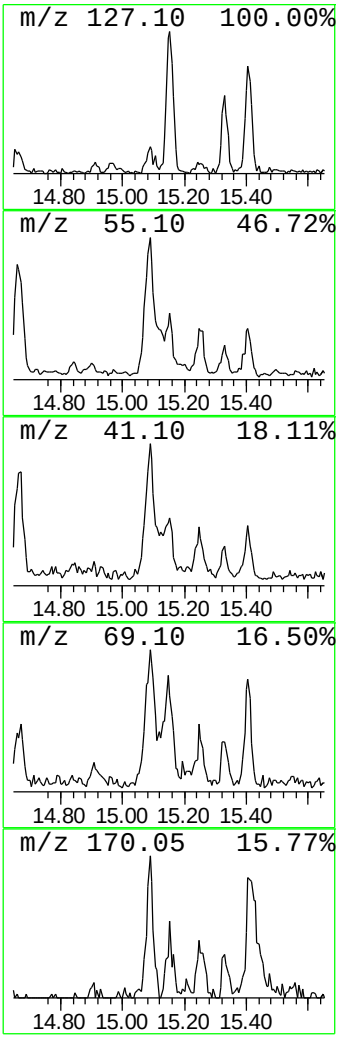
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.15	4.33 ng/ul	271114	Acenaphthene-d10	14.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Imidazole, 1-methyl-5-nitro-	127	C4H5N3O2	003034-42-2	52
2	Cyclobutylcarboxamide, N-(4-chlo...	209	C11H12ClNO	1000340-37-4	50
3	(Chlorophenoxymethylenimino)sulf...	281	C7H5ClF5NOS	081439-21-6	50
4	3,5,5-Trimethylcyclohexyl isopro...	250	C12H24FO2P	1000298-27-4	50
5	2-Isopropyl-5-methylcyclohexanyl...	264	C13H26FO2P	1000298-30-9	45





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Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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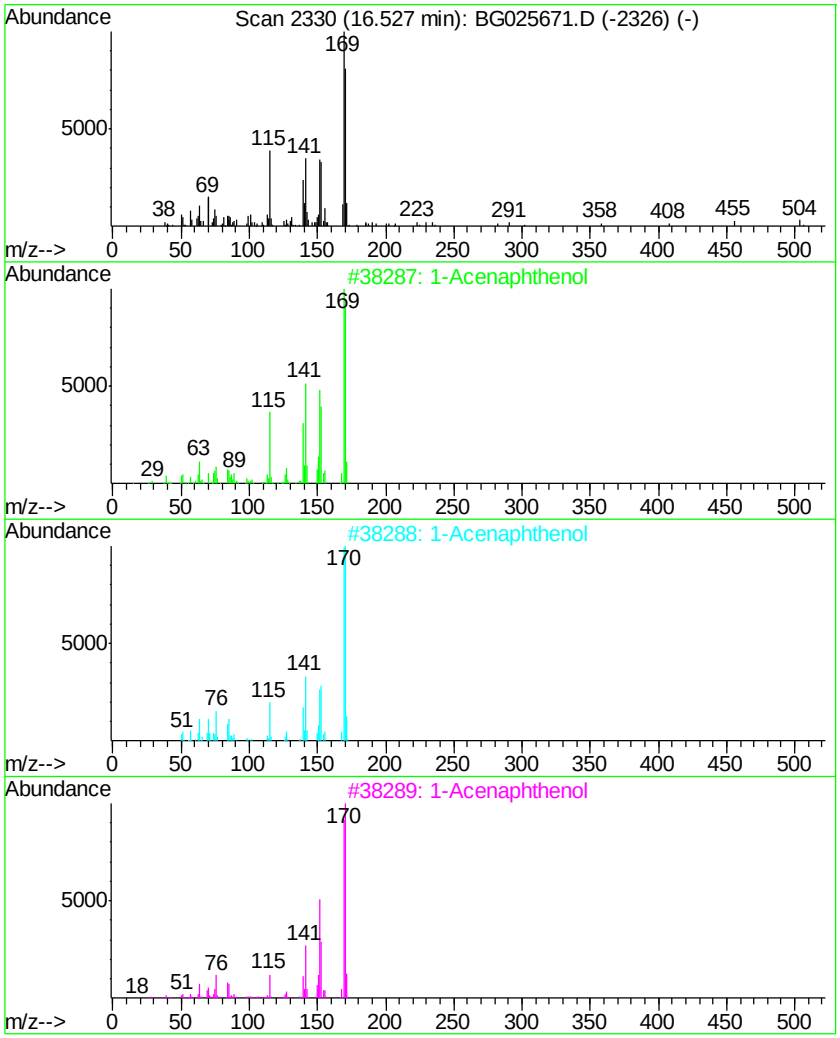
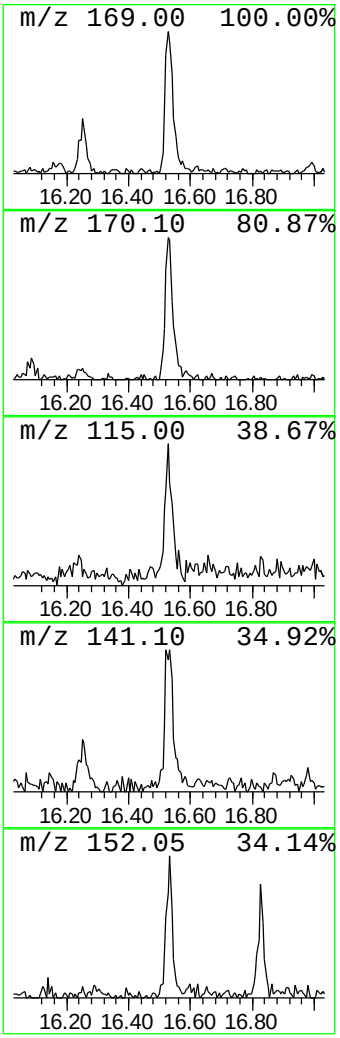
Peak Number 23 1-Acenaphthenol Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	2.57 ng/ul	188022	Phenanthrene-d10	17.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Acenaphthenol	170	C12H10O	006306-07-6	96
2	1-Acenaphthenol	170	C12H10O	006306-07-6	83
3	1-Acenaphthenol	170	C12H10O	006306-07-6	80
4	1,8-Naphthalenedimethanol	188	C12H12O2	002026-08-6	74
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	62





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Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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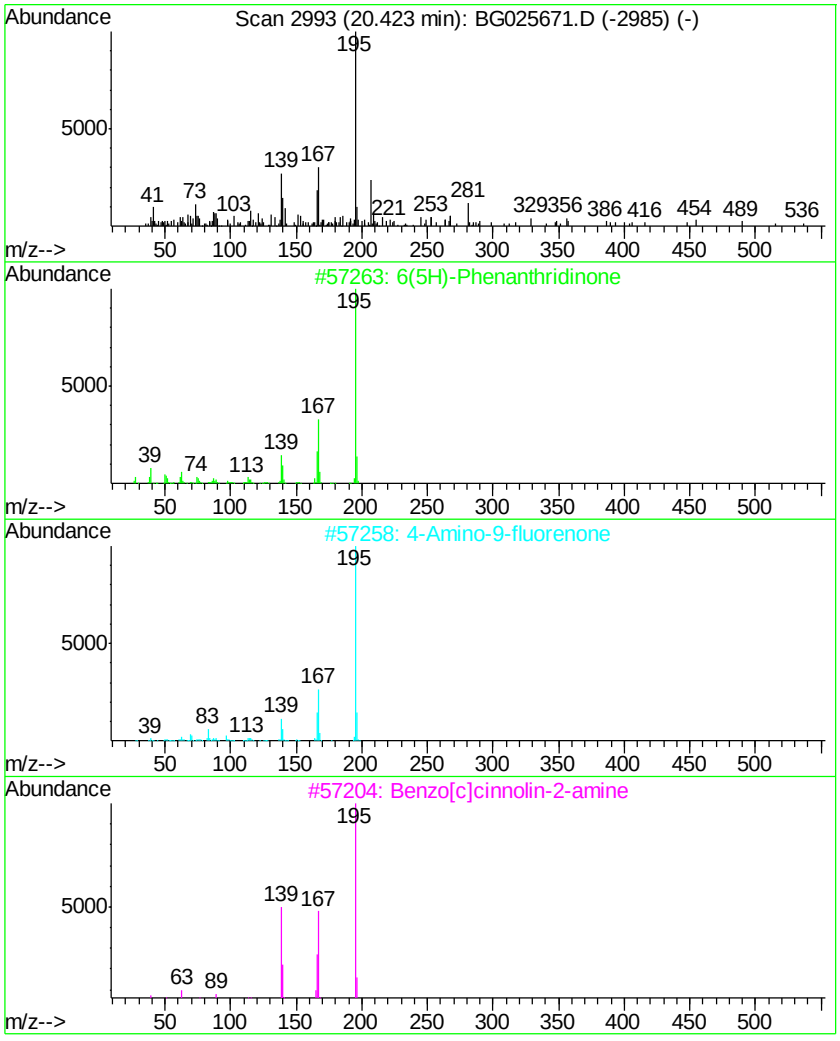
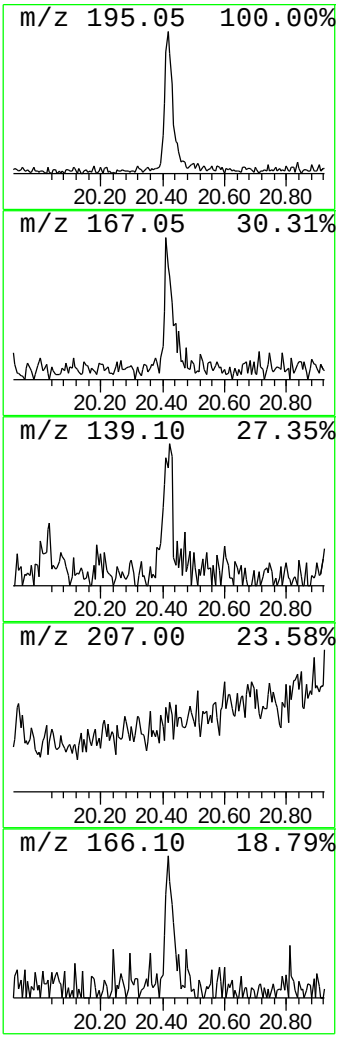
Peak Number 24 6(5H)-Phenanthridinone Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.42	2.19 ng/ul	188550	Chrysene-d12	21.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6(5H)-Phenanthridinone	195	C13H9NO	001015-89-0	93
2	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	90
3	Benzo[c]cinnolin-2-amine	195	C12H9N3	004827-07-0	90
4	4-Amino-9-fluorenone	195	C13H9NO	004269-15-2	90
5	3-Acridinol	195	C13H9NO	007132-70-9	90





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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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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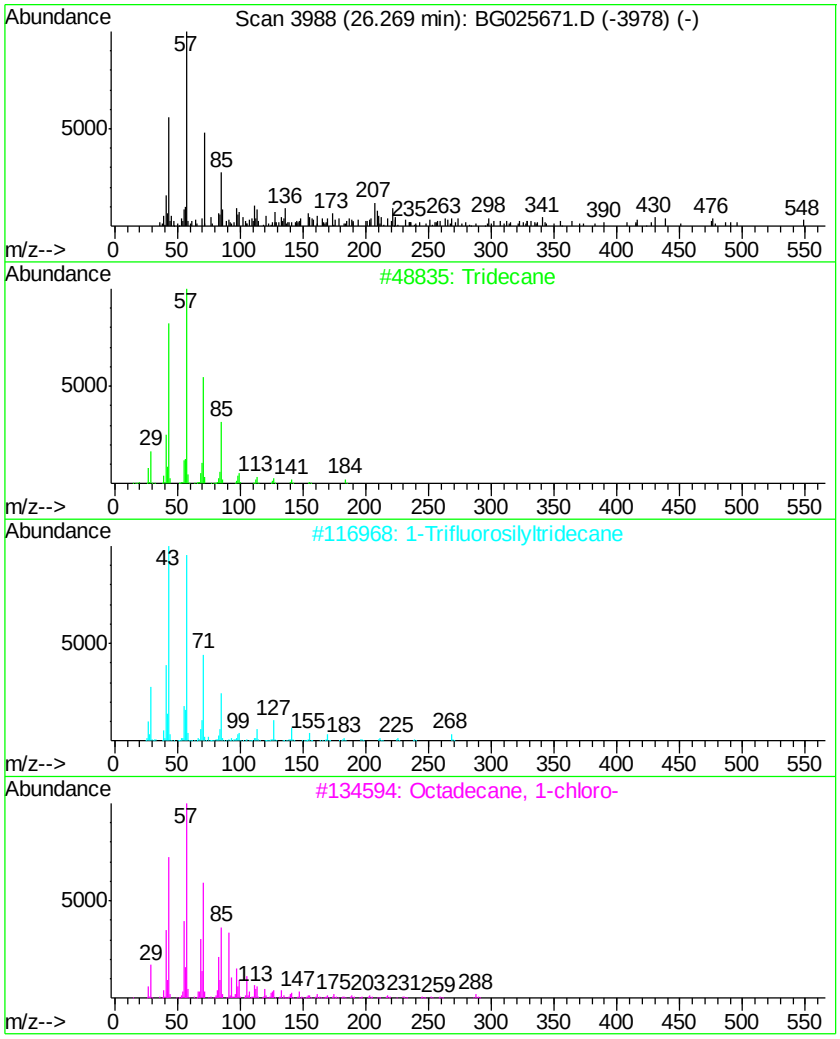
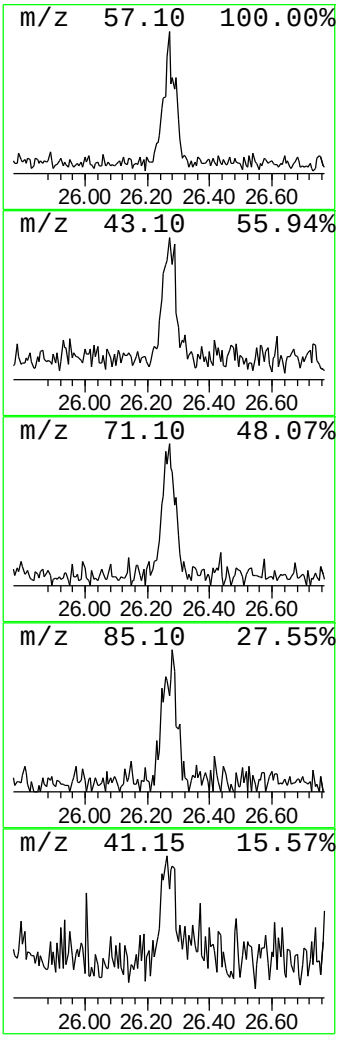
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.27	2.07 ng/ul	171191	Perylene-d12	25.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	64
2	1-Trifluorosilyltridecane	268	C13H27F3Si	1000217-08-2	58
3	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	58
4	Docosane	310	C22H46	000629-97-0	52
5	Nonacosane	408	C29H60	000630-03-5	52





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012617\
Data File : BG025671.D
Acq On : 26 Jan 2017 17:55
Operator : SJ/MA
Sample : I1387-07DL 10X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DA9R3DL

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
4-Heptanone, 2,6-...	7.29	10.8	ng/ul	244132	1	8.18	452385	20.0
2-Heptanone, 4,6-...	7.57	33.6	ng/ul	760824	1	8.18	452385	20.0
Cyclohexanone, 3,...	8.60	46.7	ng/ul	1055770	1	8.18	452385	20.0
unknown-01	8.80	2.9	ng/ul	65932	1	8.18	452385	20.0
unknown-02	9.28	5.0	ng/ul	113043	1	8.18	452385	20.0
Cyclopentanone, 3...	9.77	2.9	ng/ul	136510	2	11.00	928227	20.0
unknown-03	10.01	3.4	ng/ul	156748	2	11.00	928227	20.0
(DEL) Alkane: Cyc...	10.51	4.0	ng/ul	186662	2	11.00	928227	20.0
Cyclopenta[c]thia...	11.17	2.9	ng/ul	136122	2	11.00	928227	20.0
(DEL) Alkane: Cyc...	11.26	9.9	ng/ul	458833	2	11.00	928227	20.0
unknown-04	11.33	8.2	ng/ul	381647	2	11.00	928227	20.0
(DEL) Alkane: Str...	11.47	2.8	ng/ul	131131	2	11.00	928227	20.0
unknown-05	11.89	6.2	ng/ul	289472	2	11.00	928227	20.0
unknown-06	12.40	2.7	ng/ul	124985	2	11.00	928227	20.0
Naphthalene, 1-me...	12.86	4.0	ng/ul	185004	2	11.00	928227	20.0
Butane, 1-[(2-met...	13.71	2.6	ng/ul	163844	3	14.81	1252090	20.0
unknown-07	13.82	5.6	ng/ul	352093	3	14.81	1252090	20.0
unknown-08	13.96	4.2	ng/ul	261549	3	14.81	1252090	20.0
unknown-09	14.14	2.1	ng/ul	133662	3	14.81	1252090	20.0
(DEL) Alkane: Str...	14.66	10.9	ng/ul	683605	3	14.81	1252090	20.0
(DEL) Alkane: Str...	15.09	13.7	ng/ul	858538	3	14.81	1252090	20.0
(DEL) Alkane: Str...	15.15	4.3	ng/ul	271114	3	14.81	1252090	20.0
1-Acenaphthenol	16.53	2.6	ng/ul	188022	4	17.56	1464720	20.0
6(5H)-Phenanthrid...	20.42	2.2	ng/ul	188550	5	21.84	1722350	20.0
(DEL) Alkane: Str...	26.27	2.1	ng/ul	171191	6	25.23	1653310	20.0