LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

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\SOM02.2-EPA-BG080915.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	2.890	7	9	14	rVB	36210	36377	1.13%	0.094%
2	3.113	42	47	54	rBV	68351	114809	3.57%	0.297%
3	3.184	55	59	73	rVB	694628	924627	28.74%	2.396%
4	3.442	100	103	110	rVB3	17667	30694	0.95%	0.080%
5	3.713	147	149	155	rBV3	7928	12044	0.37%	0.031%
6	4.565	290	294	296	rBV5	4152	6644	0.21%	0.017%
7	5.710	487	489	496	rVB6	4439	8057	0.25%	0.021%
8	5.910	521	523	526	rBV4	5774	7852	0.24%	0.020%
9	6.580	636	637	644	rVB4	6840	9353	0.29%	0.024%
10	6.703	653	658	664	rVB4	16137	27764	0.86%	0.072%
11	7.003	704	709	714	rVB6	16694	28640	0.89%	0.074%
12	7.167	728	737	746	rBV	336660	572067	17.78%	1.482%
13	7.332	757	765	772	rBV	272772	448044	13.92%	1.161%
14	7.461	781	787	795	rVB3	20627	37333	1.16%	0.097%
15	7.543	795	801	809	rBV	313742	527686	16.40%	1.367%
16	7.784	838	842	844	rBV4	6610	7896	0.25%	0.020%
17	7.914	859	864	871	rVB10	10032	20104	0.62%	0.052%
18	8.013	871	881	887	rBV	249192	429188	13.34%	1.112%
19	8.231	914	918	919	rBV3	6902	7739	0.24%	0.020%
20	8.713	993	1000	1008	rBV2	425458	740973	23.03%	1.920%
21	9.165	1069	1077	1086	rBV	319514	565314	17.57%	1.465%
22	9.494	1127	1133	1137	rVB5	4219	8556	0.27%	0.022%
23	9.535	1137	1140	1143	rBV5	4788	6445	0.20%	0.017%
24	9.894	1193	1201	1210	rBV	245500	450518	14.00%	1.167%
25	10.228	1251	1258	1262	rBV8	6462	14081	0.44%	0.036%
26	10.434	1286	1293	1303	rVV	476428	828567	25.75%	2.147%
27	10.581	1315	1318	1324	rBV7	5975	12865	0.40%	0.033%
28	10.663	1327	1332	1336	rBV6	5162	9166	0.28%	0.024%
29	10.816	1351	1358	1366	rBV	426164	739637	22.99%	1.917%
30	10.945	1372	1380	1388	rBV	381195	688540	21.40%	1.784%
31	11.462	1465	1468	1473	rBV6	4335	8337	0.26%	0.022%
32	11.686	1503	1506	1509	rBV5	4840	6802	0.21%	0.018%
33	12.167	1585	1588	1594	rVB6	11838	19303	0.60%	0.050%
34	12.414	1625	1630	1633	rVB6	5425	8423	0.26%	0.022%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

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\SOM02.2-EPA-BG080915.M
Title : SVOA CALIBRATION

35	13.002	1725	1730	1737	rBV6	14429	28078	0.87%	0.073%
36	13.243	1766	1771	1776	rVB3	26501	38569	1.20%	0.100%
37	13.389	1790	1796	1799	rBV7	7140	15673	0.49%	0.041%
38	13.525	1814	1819	1823	rBV7	14078	24739	0.77%	0.064%
39	13.583	1826	1829	1833	rVB6	10274	15164	0.47%	0.039%
40	13.965	1888	1894	1898	rBV8	8289	14393	0.45%	0.037%
41	14.036	1900	1906	1911	rVV	925038	1411855	43.88%	3.658%
42	14.083	1911	1914	1921	rVV	394736	562015	17.47%	1.456%
43	14.142	1921	1924	1925	rVV3	10810	11307	0.35%	0.029%
44	14.171	1925	1929	1933	rVB2	39952	53183	1.65%	0.138%
45	14.330	1945	1956	1962	rVV	1102188	1743567	54.19%	4.518%
46	14.535	1986	1991	1994	rBV5	9954	13736	0.43%	0.036%
47	14.600	1997	2002	2004	rVV2	105469	157542	4.90%	0.408%
48	14.635	2004	2008	2016	rVB2	765494	1267061	39.38%	3.283%
49	14.711	2016	2021	2024	rBV5	19509	30695	0.95%	0.080%
50	14.823	2033	2040	2048	rBV	483293	774204	24.06%	2.006%
51	15.035	2074	2076	2078	rBV3	6294	6958	0.22%	0.018%
52	15.170	2093	2099	2107	rVB4	50479	87892	2.73%	0.228%
53	15.434	2141	2144	2147	rVB3	9306	9529	0.30%	0.025%
54	15.481	2148	2152	2156	rVB5	22164	26584	0.83%	0.069%
55	15.534	2156	2161	2162	rBV5	5505	7382	0.23%	0.019%
56	15.628	2170	2177	2184	rBV	1581373	2245285	69.78%	5.818%
57	15.740	2189	2196	2207	rBV	363141	573539	17.82%	1.486%
58	15.869	2213	2218	2225	rBV9	18297	37638	1.17%	0.098%
59	15.916	2225	2226	2232	rVB6	5197	6283	0.20%	0.016%
60	16.028	2244	2245	2247	rBV2	6688	6819	0.21%	0.018%
61	16.110	2254	2259	2263	rBV7	14585	24082	0.75%	0.062%
62	16.345	2294	2299	2302	rBV4	34592	52251	1.62%	0.135%
63	16.686	2354	2357	2359	rBV4	6883	6767	0.21%	0.018%
64	16.768	2369	2371	2374	rBV4	7613	8107	0.25%	0.021%
65	16.909	2393	2395	2398	rVB4	6090	6612	0.21%	0.017%
66	17.032	2412	2416	2418	rBV5	8324	10810	0.34%	0.028%
67	17.132	2428	2433	2436	rBV	43685	61174	1.90%	0.159%
68	17.267	2452	2456	2460	rBV7	16009	22514	0.70%	0.058%
69	17.326	2464	2466	2468	rBV3	9014	8568	0.27%	0.022%
70	17.379	2469	2475	2481	rVV	1185757	1769198	54.98%	4.584%
71	17.420	2481	2482	2486	rVV	31984	28525	0.89%	0.074%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

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\SOM02.2-EPA-BG080915.M
Title : SVOA CALIBRATION

72	17.479	2486	2492	2499	rVB2	1720213	2594379	80.63%	6.722%
73	17.872	2555	2559	2563	rVB4	23750	29547	0.92%	0.077%
74	18.137	2600	2604	2612	rBV9	23501	44663	1.39%	0.116%
75	18.207	2612	2616	2620	rBV3	47580	72037	2.24%	0.187%
76	18.266	2621	2626	2636	rBV2	498777	763601	23.73%	1.979%
77	18.436	2652	2655	2656	rBV2	12461	13372	0.42%	0.035%
78	18.560	2673	2676	2681	rBV7	23219	42282	1.31%	0.110%
79	18.789	2713	2715	2719	rVB5	14756	12099	0.38%	0.031%
80	18.924	2736	2738	2742	rBV5	8695	10788	0.34%	0.028%
81	19.153	2775	2777	2778	rBV2	10654	8031	0.25%	0.021%
82	19.424	2815	2823	2830	rBV2	108540	230792	7.17%	0.598%
83	19.535	2838	2842	2847	rBV2	138331	202740	6.30%	0.525%
84	19.688	2865	2868	2874	rVB	85140	105653	3.28%	0.274%
85	19.758	2874	2880	2892	rVB2	2165030	3217680	100.00%	8.337%
86	20.628	3025	3028	3029	rBV3	29088	33479	1.04%	0.087%
87	20.669	3031	3035	3043	rVB	364609	494113	15.36%	1.280%
88	20.775	3049	3053	3058	rBV2	123756	190748	5.93%	0.494%
89	21.292	3137	3141	3153	rBV2	194362	461897	14.35%	1.197%
90	21.639	3193	3200	3206	rBV	1281923	2075466	64.50%	5.378%
91	21.833	3231	3233	3238	rVB	48344	56719	1.76%	0.147%
92	22.373	3320	3325	3332	rVB2	386240	726337	22.57%	1.882%
93	22.438	3333	3336	3339	rBV2	64657	90088	2.80%	0.233%
94	22.549	3351	3355	3363	rVB3	189989	371308	11.54%	0.962%
95	22.696	3375	3380	3384	rVB5	162840	268991	8.36%	0.697%
96	23.037	3435	3438	3443	rBV6	98048	182702	5.68%	0.473%
97	23.102	3444	3449	3464	rVB	981772	2042610	63.48%	5.293%
98	23.636	3535	3540	3550	rVB4	374739	881882	27.41%	2.285%
99	24.617	3697	3707	3720	rVB	1025000	2784707	86.54%	7.216%
100	24.835	3735	3744	3760	rVB	672077	2051568	63.76%	5.316%

Sum of corrected areas: 38593022

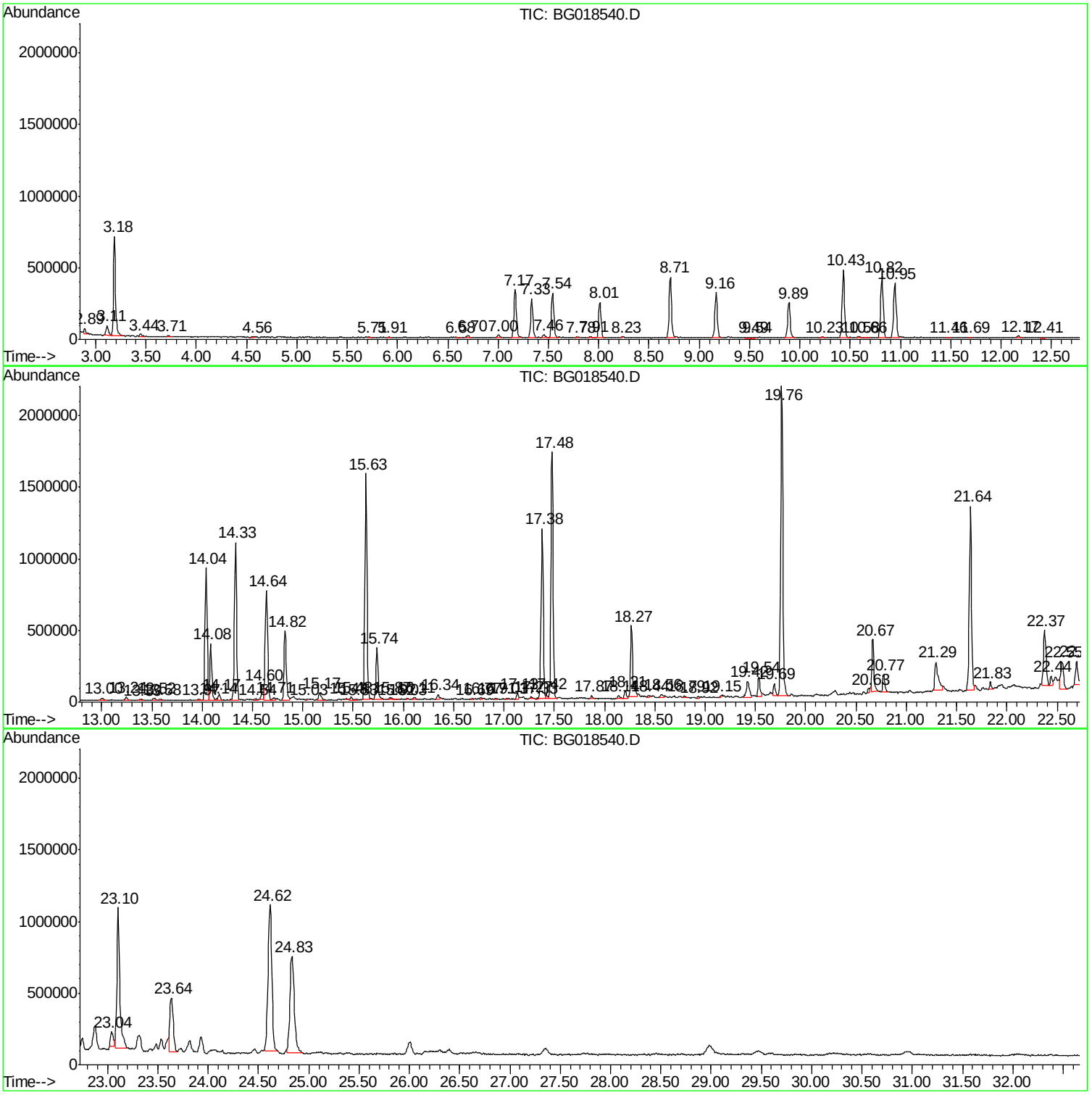
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.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\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

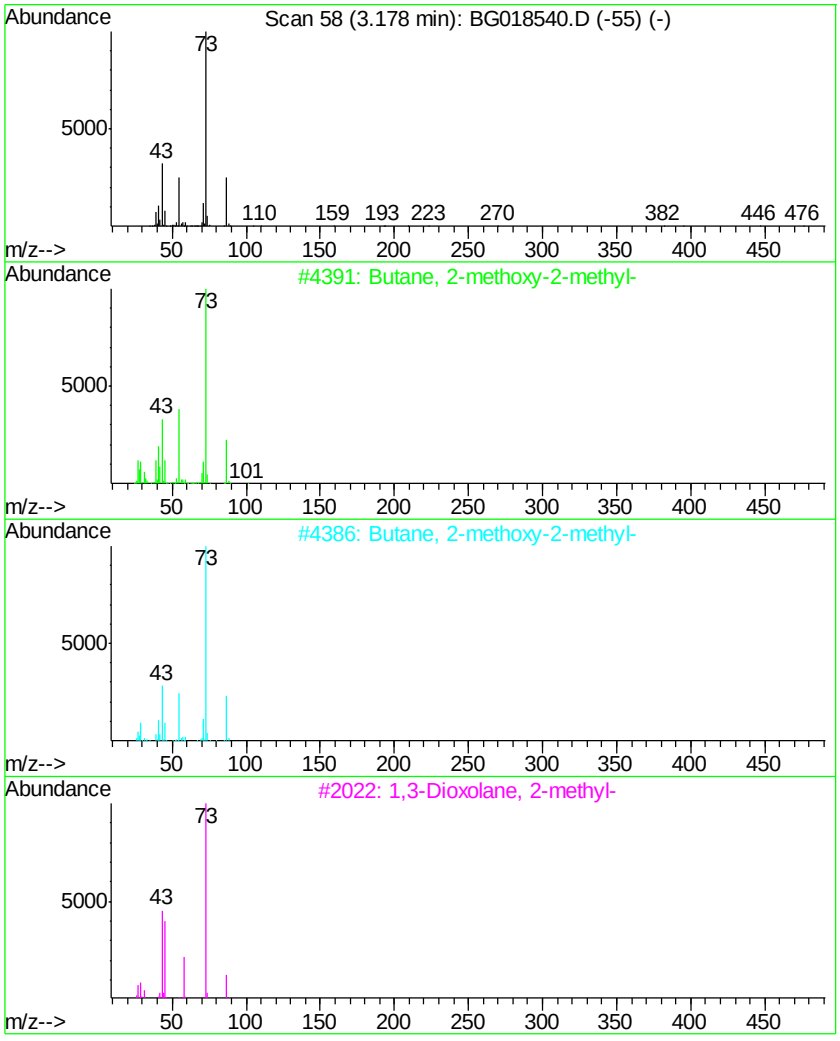
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

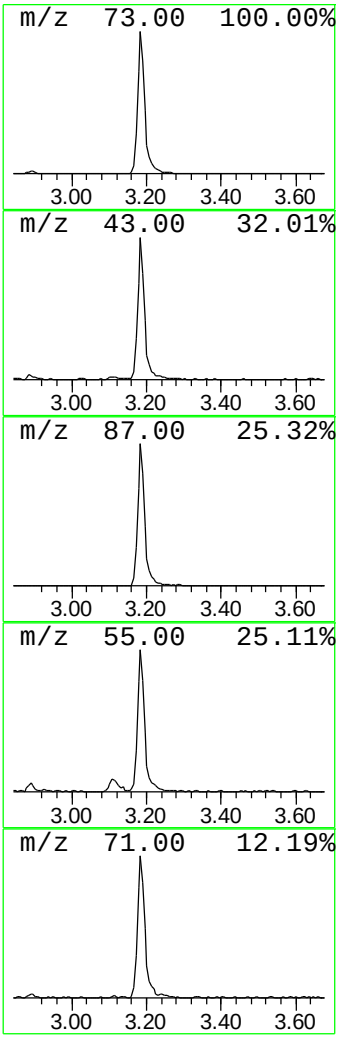
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	43.09 ng/ul	924627	1,4-Dichlorobenzene-d4	8.01

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	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

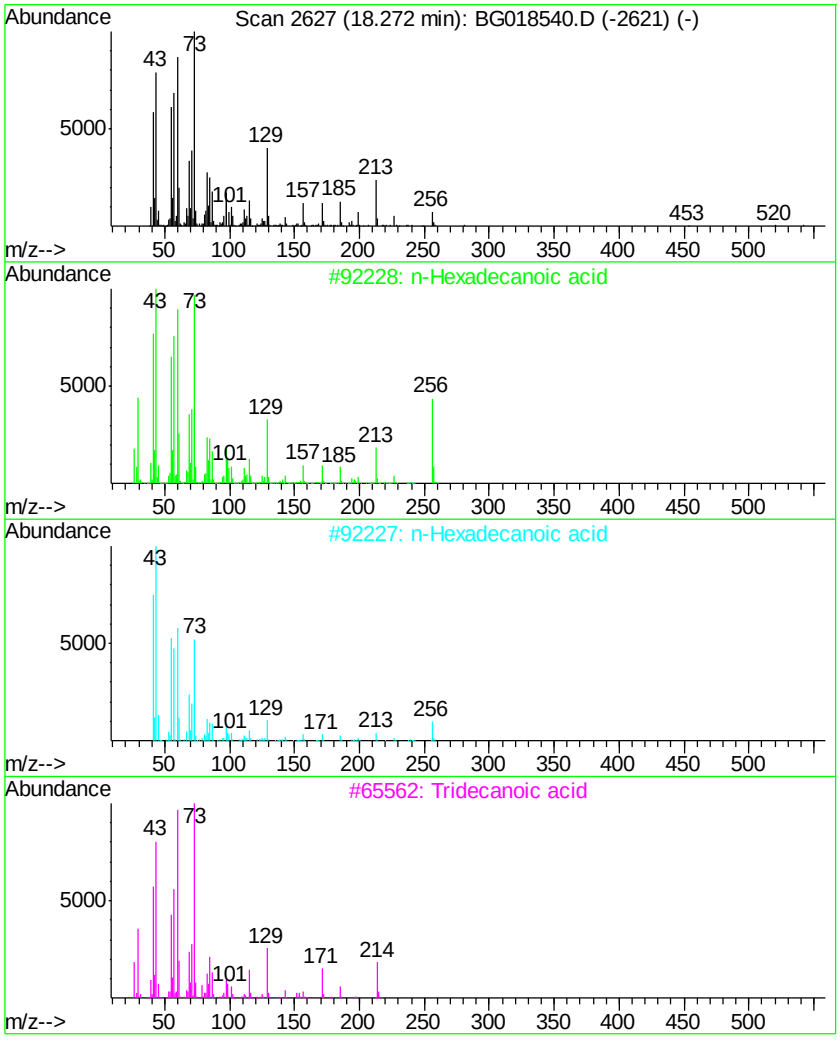
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

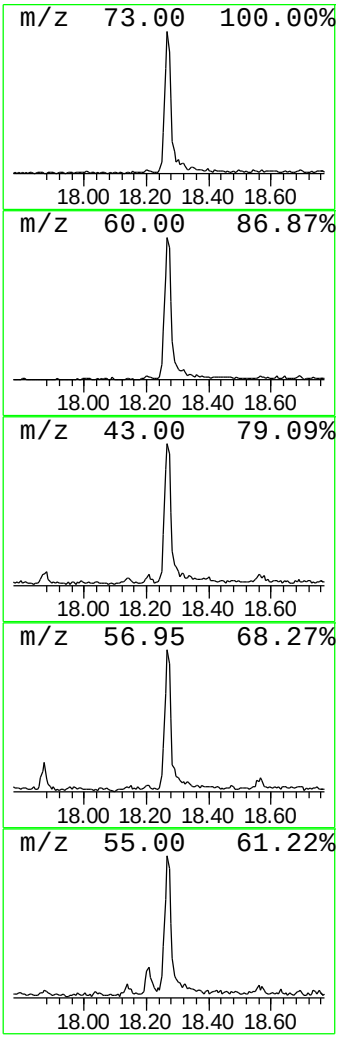
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.27	8.63 ng/ul	763601	Phenanthrene-d10	17.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	90
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	78





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

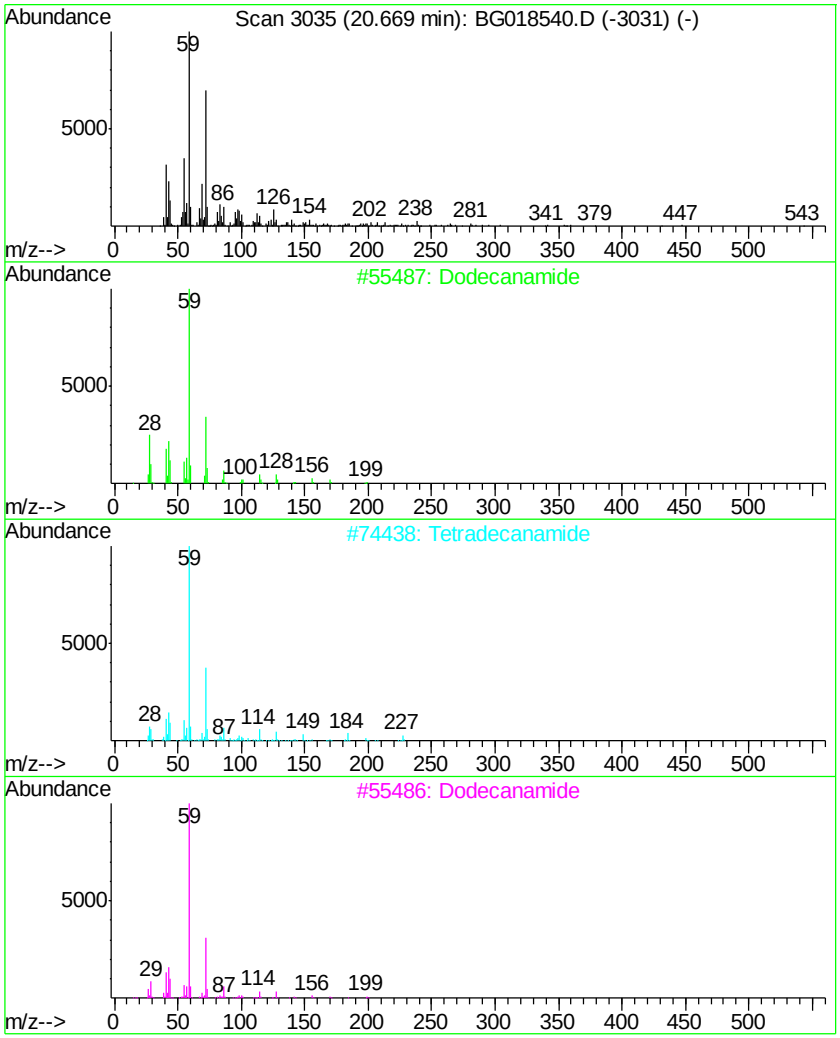
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

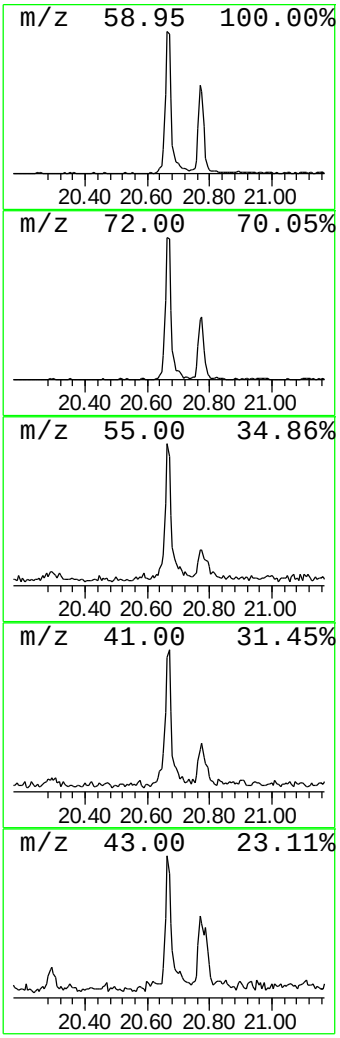
****

Peak Number 4 Dodecanamide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.67	4.76 ng/ul	494113	Chrysene-d12	21.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanamide	199	C12H25NO	001120-16-7	81
2	Tetradecanamide	227	C14H29NO	000638-58-4	78
3	Dodecanamide	199	C12H25NO	001120-16-7	74
4	Octadecanamide	283	C18H37NO	000124-26-5	64
5	Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	64





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

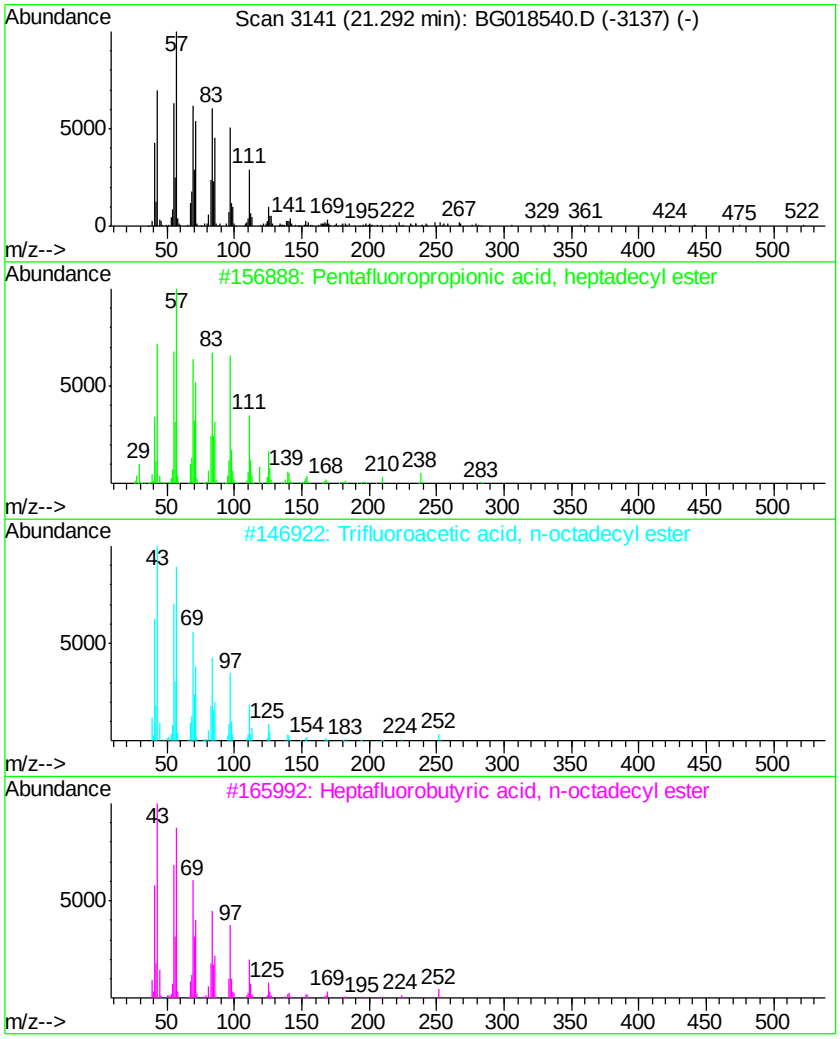
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

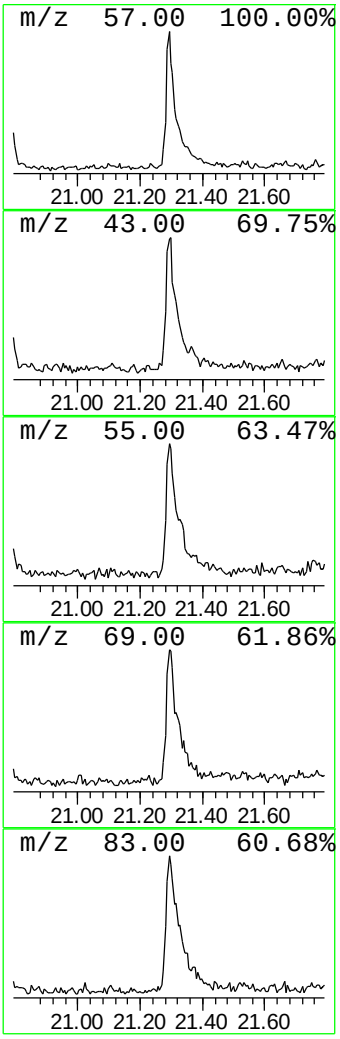
****

Peak Number 5 Pentafluoropropionic acid, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.29	4.45 ng/ul	461897	Chrysene-d12	21.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87
2	Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	83
3	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	80
4	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	80
5	17-Pentatriacontene	491	C35H70	006971-40-0	72





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

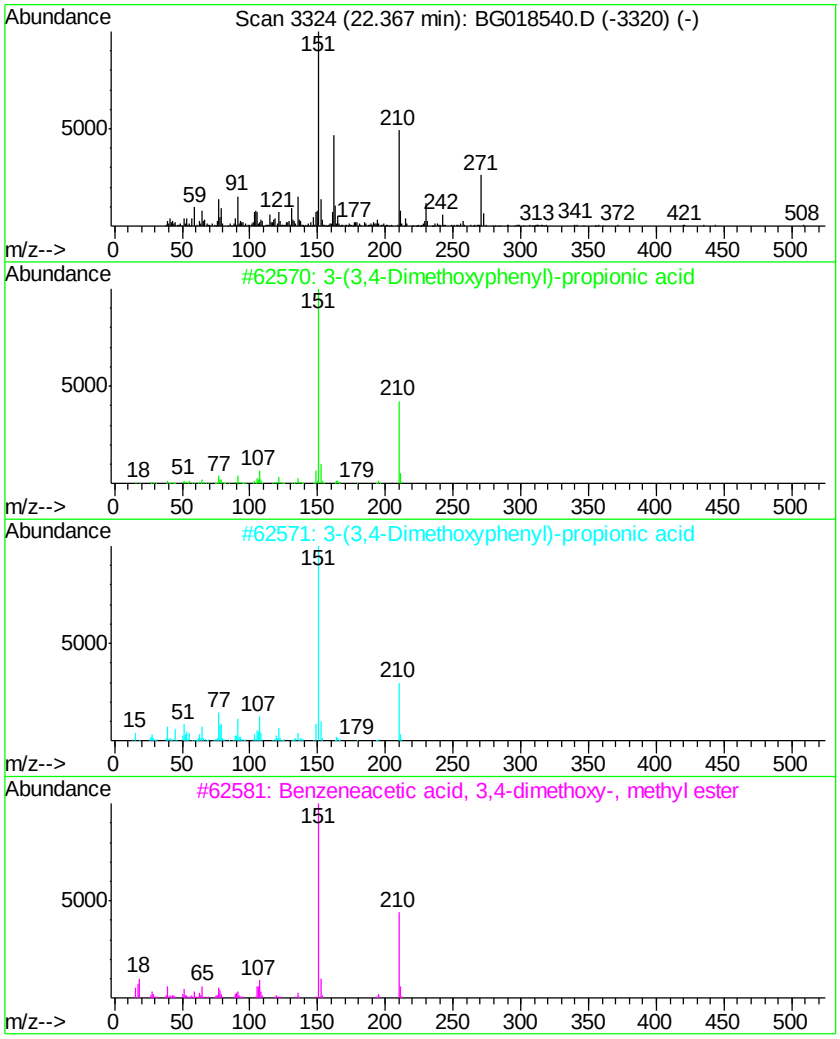
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

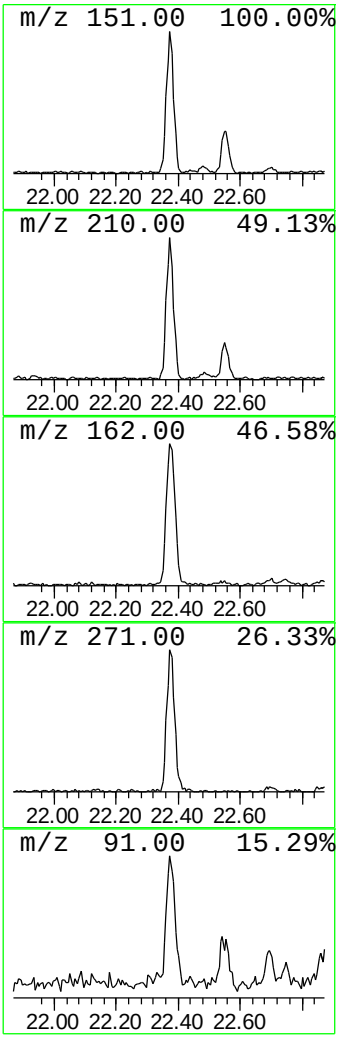
****

Peak Number 6 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.37	7.00 ng/ul	726337	Chrysene-d12	21.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(3,4-Dimethoxyphenyl)-propioni...	210	C11H14O4	002107-70-2	49
2	3-(3,4-Dimethoxyphenyl)-propioni...	210	C11H14O4	002107-70-2	47
3	Benzeneacetic acid, 3,4-dimethox...	210	C11H14O4	015964-79-1	46
4	Benzeneacetic acid, 3,4-dimethox...	210	C11H14O4	015964-79-1	43
5	Acetic acid, 2-(3,5-dimethoxyphe...	210	C11H14O4	006512-32-9	38





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

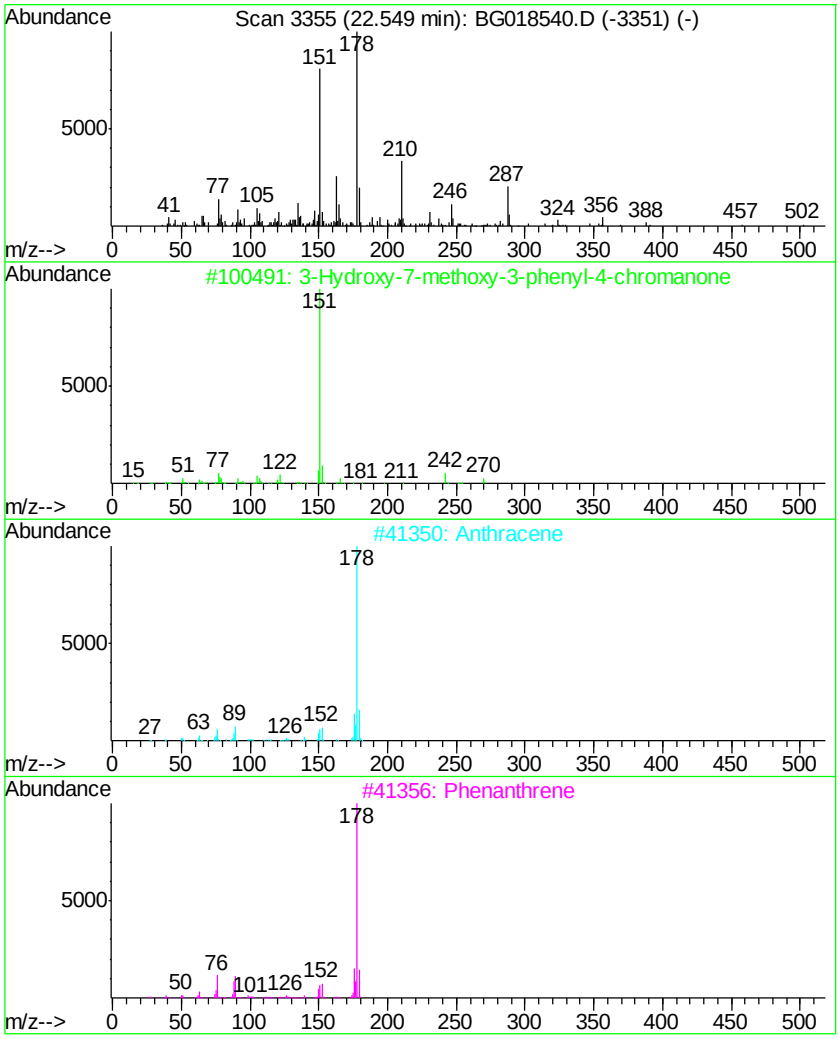
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

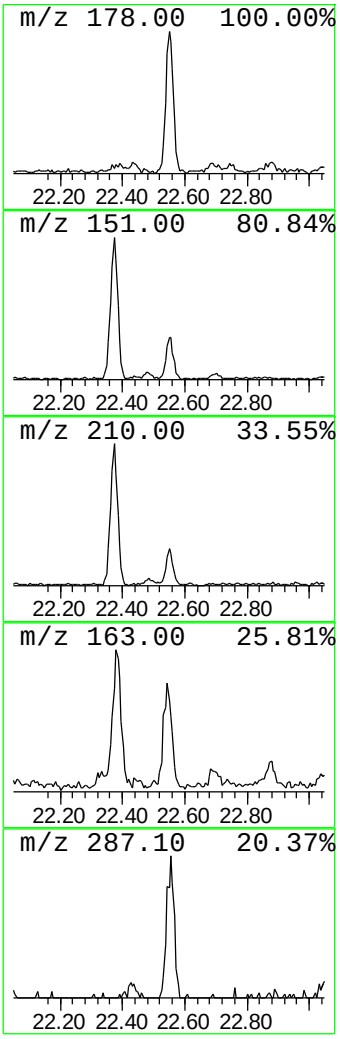
****

Peak Number 7 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.55	3.58 ng/ul	371308	Chrysene-d12	21.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hydroxy-7-methoxy-3-phenyl-4-c...	270	C16H14O4	018380-57-9	35
2	Anthracene	178	C14H10	000120-12-7	30
3	Phenanthrene	178	C14H10	000085-01-8	30
4	Benzene, 1,2-dimethoxy-4-(2-prop...	178	C11H14O2	000093-15-2	27
5	Phenanthrene	178	C14H10	000085-01-8	27





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Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

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Quant Title : SVOA CALIBRATION

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

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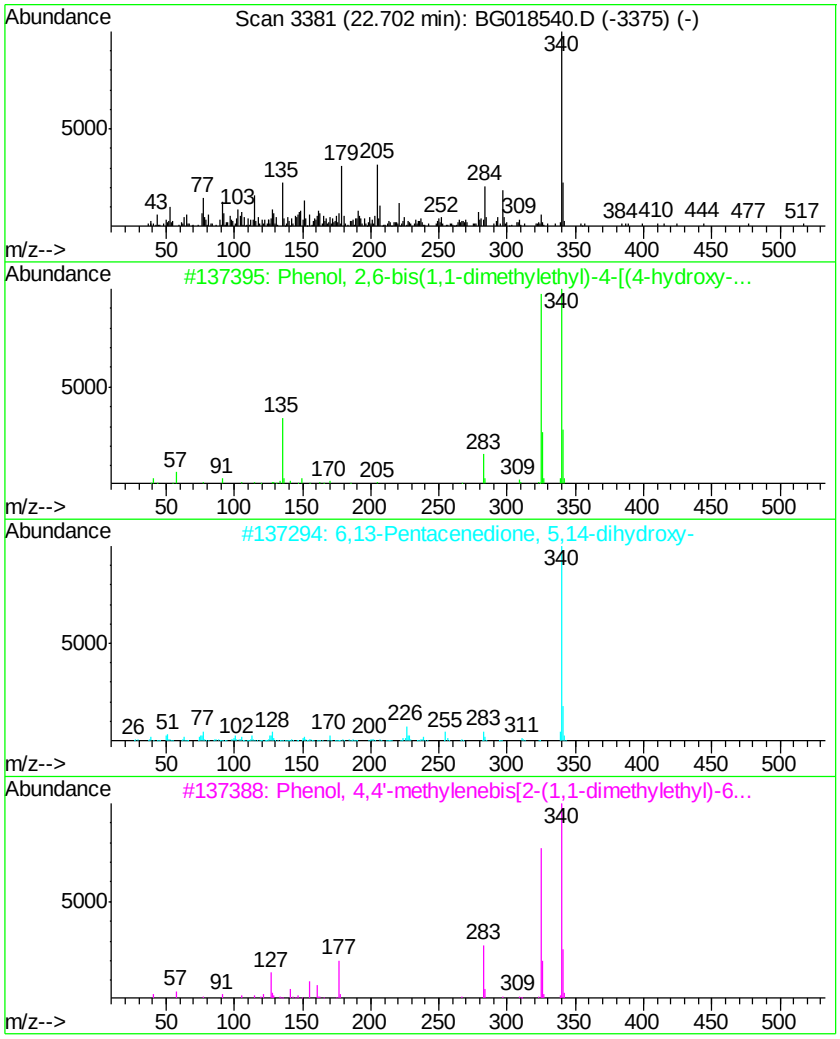
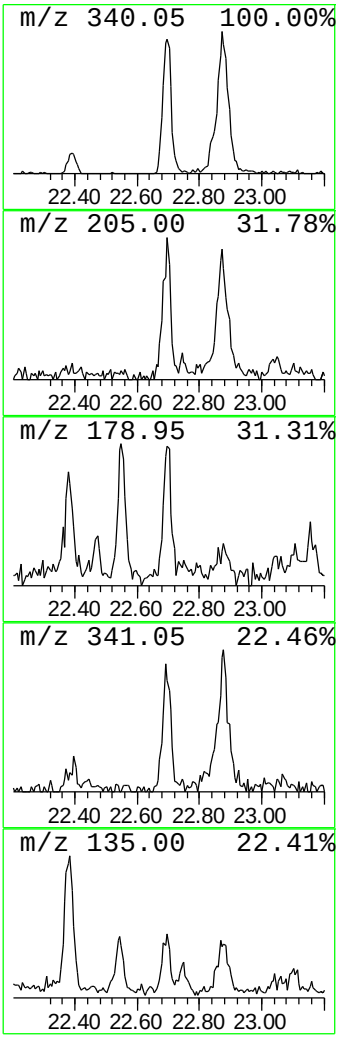
Peak Number 8 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.70	2.59 ng/ul	268991	Chrysene-d12	21.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-bis(1,1-dimethylethy...	340	C23H32O2	020690-84-0	58
2	6,13-Pentacenedione, 5,14-dihydr...	340	C22H12O4	068970-90-1	50
3	Phenol, 4,4'-methylenebis[2-(1,1...	340	C23H32O2	000096-65-1	46
4	Colchicine, 7-deacetoamino-5,6-d...	340	C20H20O5	076265-62-8	43
5	Estra-1,3,5(10),7-tetraen-17-one...	340	C21H28O2Si	069688-22-8	40





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG082915\
Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

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

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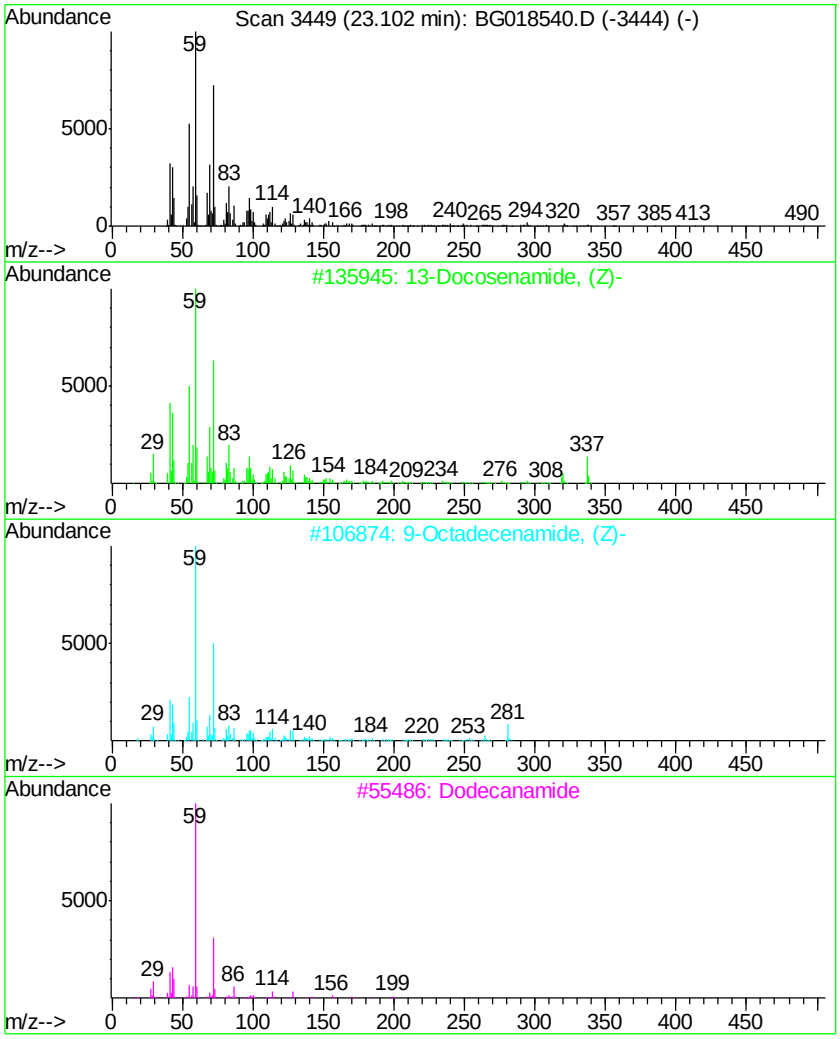
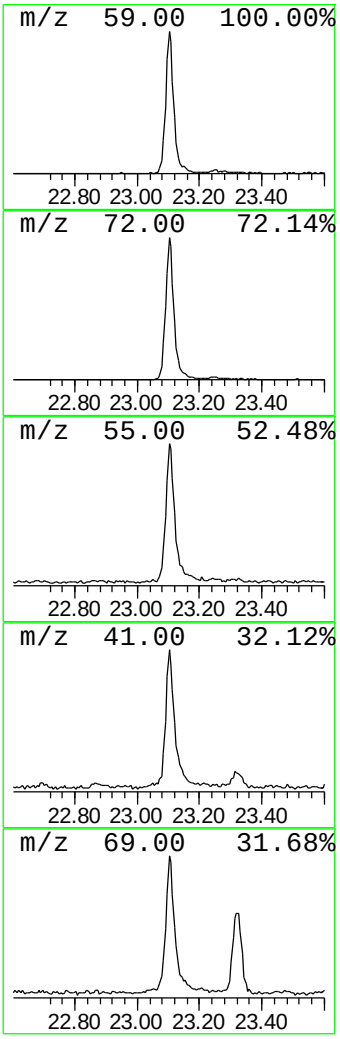
Peak Number 9 13-Docosenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.10	19.68 ng/ul	2042610	Chrysene-d12	21.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	87
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	80
3	Dodecanamide	199	C12H25NO	001120-16-7	59
4	Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	50
5	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	38





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Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
Quant Title : SVOA CALIBRATION

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

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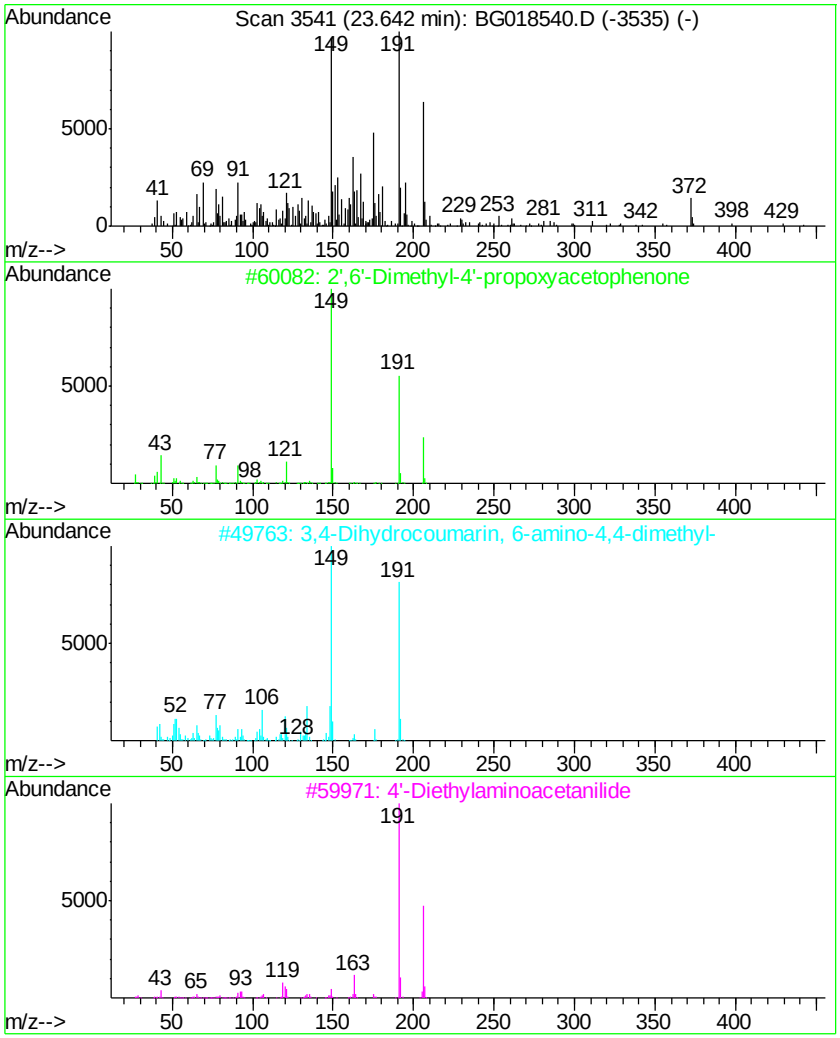
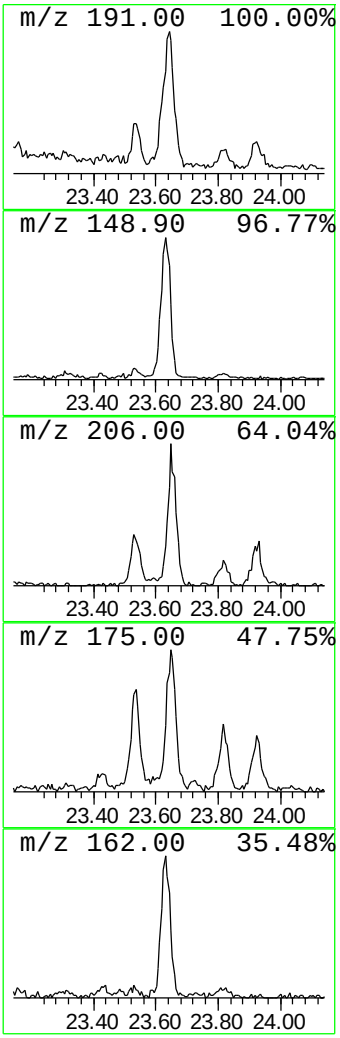
Peak Number 10 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.64	8.60 ng/ul	881882	Perylene-d12	24.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2',6'-Dimethyl-4'-propoxyacetoph...	206	C13H18O2	1000195-98-1	49
2	3,4-Dihydrocoumarin, 6-amino-4,4...	191	C11H13NO2	1000128-49-9	38
3	4'-Diethylaminoacetanilide	206	C12H18N2O	005326-57-8	25
4	2-Cyano-3-(5-methyl-furan-2-yl)-...	191	C10H9NO3	073403-31-3	25
5	4-Diethylaminophenyl isothiocyanate	206	C11H14N2S	084381-54-4	22





Tentatively Identified Compound (LSC) summary

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Data File : BG018540.D
Acq On : 29 Aug 2015 17:30
Operator : UM/MA
Sample : G3476-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BCTJ4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.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
Butane, 2-methoxy...	3.18	43.1	ng/ul	924627	1	8.01	429188	20.0
n-Hexadecanoic acid	18.27	8.6	ng/ul	763601	4	17.38	1769200	20.0
Dodecanamide	20.67	4.8	ng/ul	494113	5	21.64	2075470	20.0
Pentafluoropropio...	21.29	4.5	ng/ul	461897	5	21.64	2075470	20.0
unknown-01	22.37	7.0	ng/ul	726337	5	21.64	2075470	20.0
unknown-02	22.55	3.6	ng/ul	371308	5	21.64	2075470	20.0
Phenol, 2,6-bis(1...	22.70	2.6	ng/ul	268991	5	21.64	2075470	20.0
13-Docosenamide, ...	23.10	19.7	ng/ul	2042610	5	21.64	2075470	20.0
unknown-03	23.64	8.6	ng/ul	881882	6	24.83	2051570	20.0