LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101316\
Data File : BG024318.D
Acq On : 13 Oct 2016 13:26
Operator : UM/SJ
Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

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-BG0100716.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	3.646	131	137	143	rBV3	45314	82934	1.67%	0.139%
2	3.793	159	162	166	rVB5	12063	15858	0.32%	0.027%
3	3.840	166	170	172	rBV4	14578	18379	0.37%	0.031%
4	3.916	180	183	190	rBV	64552	106287	2.15%	0.179%
5	4.081	207	211	216	rBV8	12635	27563	0.56%	0.046%
6	4.410	261	267	274	rBV8	22201	49499	1.00%	0.083%
7	4.745	318	324	328	rBV5	43295	73159	1.48%	0.123%
8	4.786	328	331	336	rVB5	27824	42057	0.85%	0.071%
9	5.027	367	372	375	rBV7	13805	19476	0.39%	0.033%
10	5.356	420	428	438	rVV	119536	207942	4.20%	0.350%
11	5.573	459	465	472	rVB3	62601	117604	2.37%	0.198%
12	5.644	472	477	483	rBV10	17400	39299	0.79%	0.066%
13	5.973	530	533	538	rVB6	11618	21930	0.44%	0.037%
14	6.255	571	581	587	rBV	97691	184655	3.73%	0.310%
15	6.760	659	667	672	rBV8	9845	20398	0.41%	0.034%
16	6.907	686	692	698	rVV9	15065	26759	0.54%	0.045%
17	7.218	738	745	754	rBV7	21259	64400	1.30%	0.108%
18	7.418	771	779	792	rBV	832997	1580253	31.90%	2.657%
19	7.606	802	811	822	rBV	569785	984448	19.87%	1.655%
20	7.700	822	827	832	rVV8	13832	17492	0.35%	0.029%
21	7.812	836	846	858	rVB	774895	1384648	27.95%	2.328%
22	8.106	894	896	907	rVB	10987	20676	0.42%	0.035%
23	8.288	919	927	942	rBV2	592079	1103509	22.28%	1.855%
24	8.975	1035	1044	1053	rVB2	1001364	1792419	36.18%	3.014%
25	9.122	1059	1069	1077	rBV3	36140	75979	1.53%	0.128%
26	9.386	1105	1114	1120	rBV2	80139	158838	3.21%	0.267%
27	9.463	1120	1127	1134	rVV	799453	1493751	30.15%	2.511%
28	9.527	1134	1138	1144	rVB4	86223	143799	2.90%	0.242%
29	9.851	1189	1193	1200	rVB7	23848	35708	0.72%	0.060%
30	10.185	1242	1250	1258	rVV	698143	1302917	26.30%	2.191%
31	10.720	1333	1341	1353	rVB	1029935	1908103	38.52%	3.208%
32	10.873	1361	1367	1375	rVB10	9926	19740	0.40%	0.033%
33	11.120	1400	1409	1417	rVB	863893	1590868	32.11%	2.675%
34	11.243	1422	1430	1447	rBV2	923394	1808624	36.51%	3.041%

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Operator : UM/SJ
Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

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-BG0100716.M
Title : SVOA CALIBRATION

35	11.390	1452	1455	1463	rVB8	9675	19394	0.39%	0.033%
36	11.460	1463	1467	1471	rBV6	9993	16925	0.34%	0.028%
37	11.754	1509	1517	1519	rBV8	13241	26632	0.54%	0.045%
38	11.854	1526	1534	1537	rBV8	15811	23896	0.48%	0.040%
39	12.677	1669	1674	1681	rVB8	29007	51781	1.05%	0.087%
40	12.806	1691	1696	1701	rVB8	12952	26445	0.53%	0.044%
41	12.871	1701	1707	1712	rBV6	16093	40329	0.81%	0.068%
42	13.147	1751	1754	1759	rVB6	12016	22948	0.46%	0.039%
43	13.258	1767	1773	1777	rVB6	24129	40887	0.83%	0.069%
44	13.341	1781	1787	1792	rVB9	15529	20328	0.41%	0.034%
45	13.499	1809	1814	1818	rVV6	28703	50058	1.01%	0.084%
46	13.535	1818	1820	1828	rVV9	18646	27942	0.56%	0.047%
47	13.770	1854	1860	1865	rBV5	38866	61987	1.25%	0.104%
48	14.298	1940	1950	1954	rBV	1738836	2686299	54.23%	4.516%
49	14.345	1954	1958	1964	rVB	1111874	1600823	32.31%	2.691%
50	14.510	1981	1986	1989	rBV7	11185	19632	0.40%	0.033%
51	14.604	1994	2002	2011	rBV	1821786	2992336	60.40%	5.031%
52	14.686	2012	2016	2019	rVB6	12130	14862	0.30%	0.025%
53	14.757	2025	2028	2033	rBV7	18425	25324	0.51%	0.043%
54	14.903	2043	2053	2061	rVB	1372440	2213481	44.68%	3.721%
55	15.062	2072	2080	2089	rBV	877510	1398347	28.23%	2.351%
56	15.174	2095	2099	2102	rBV5	13719	18802	0.38%	0.032%
57	15.262	2110	2114	2123	rBV10	27240	50074	1.01%	0.084%
58	15.444	2142	2145	2152	rVB9	15785	23515	0.47%	0.040%
59	15.562	2163	2165	2170	rVB5	9427	15045	0.30%	0.025%
60	15.667	2178	2183	2185	rBV3	24744	37092	0.75%	0.062%
61	15.703	2185	2189	2196	rVB5	42877	66218	1.34%	0.111%
62	15.891	2213	2221	2230	rBV	2468550	3961399	79.97%	6.660%
63	15.990	2230	2238	2249	rVB	832146	1271970	25.68%	2.139%
64	16.120	2255	2260	2274	rVB10	43016	89039	1.80%	0.150%
65	16.249	2281	2282	2286	rVB3	13682	15118	0.31%	0.025%
66	16.443	2311	2315	2320	rBV8	8649	15235	0.31%	0.026%
67	16.560	2328	2335	2345	rBV3	78666	191130	3.86%	0.321%
68	16.725	2355	2363	2366	rBV6	54797	98643	1.99%	0.166%
69	16.842	2377	2383	2385	rBV7	18830	30404	0.61%	0.051%
70	16.901	2390	2393	2398	rBV7	13604	22343	0.45%	0.038%
71	16.995	2402	2409	2416	rBV10	40390	76655	1.55%	0.129%

LSC Area Percent Report

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Operator : UM/SJ
Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

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-BG0100716.M
Title : SVOA CALIBRATION

72	17.077	2420	2423	2429	rVB7	15777	26402	0.53%	0.044%
73	17.136	2429	2433	2438	rBV7	15975	28808	0.58%	0.048%
74	17.354	2466	2470	2474	rBV	88232	108370	2.19%	0.182%
75	17.406	2475	2479	2488	rVB7	25342	55359	1.12%	0.093%
76	17.512	2494	2497	2501	rVB5	20635	34117	0.69%	0.057%
77	17.641	2510	2519	2528	rBV2	1595808	2510334	50.67%	4.221%
78	17.741	2530	2536	2548	rVB2	2461333	3858458	77.89%	6.487%
79	17.900	2560	2563	2567	rVB5	24670	33916	0.68%	0.057%
80	18.094	2592	2596	2600	rVB4	55095	69619	1.41%	0.117%
81	18.182	2610	2611	2621	rBV10	12919	33245	0.67%	0.056%
82	18.270	2622	2626	2631	rVB7	29211	41469	0.84%	0.070%
83	18.364	2637	2642	2650	rBV7	39287	100755	2.03%	0.169%
84	18.435	2651	2654	2658	rBV6	29306	42428	0.86%	0.071%
85	18.488	2658	2663	2674	rBV	723115	988612	19.96%	1.662%
86	18.781	2710	2713	2715	rVB4	18685	18665	0.38%	0.031%
87	19.140	2772	2774	2783	rBV9	17474	27570	0.56%	0.046%
88	19.339	2805	2808	2815	rBV9	21066	33643	0.68%	0.057%
89	19.404	2815	2819	2821	rBV5	24229	30107	0.61%	0.051%
90	19.633	2850	2858	2867	rBV7	119104	350595	7.08%	0.589%
91	19.745	2872	2877	2887	rBV2	470518	647996	13.08%	1.089%
92	19.892	2899	2902	2909	rVB4	59613	105322	2.13%	0.177%
93	20.009	2915	2922	2936	rBV	2762047	4953842	100.00%	8.329%
94	21.225	3126	3129	3138	rVB9	105690	207343	4.19%	0.349%
95	21.531	3175	3181	3187	rBV2	326942	556387	11.23%	0.935%
96	21.942	3243	3251	3259	rVB	1666372	2886809	58.27%	4.853%
97	23.699	3544	3550	3559	rVB2	204895	449268	9.07%	0.755%
98	25.150	3786	3797	3811	rVB2	1364544	4139065	83.55%	6.959%
99	25.385	3827	3837	3849	rVB3	971038	2927729	59.10%	4.922%
100	26.590	4037	4042	4054	rVB9	95438	309561	6.25%	0.520%

Sum of corrected areas: 59479103

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101316\
Data File : BG024318.D
Acq On : 13 Oct 2016 13:26
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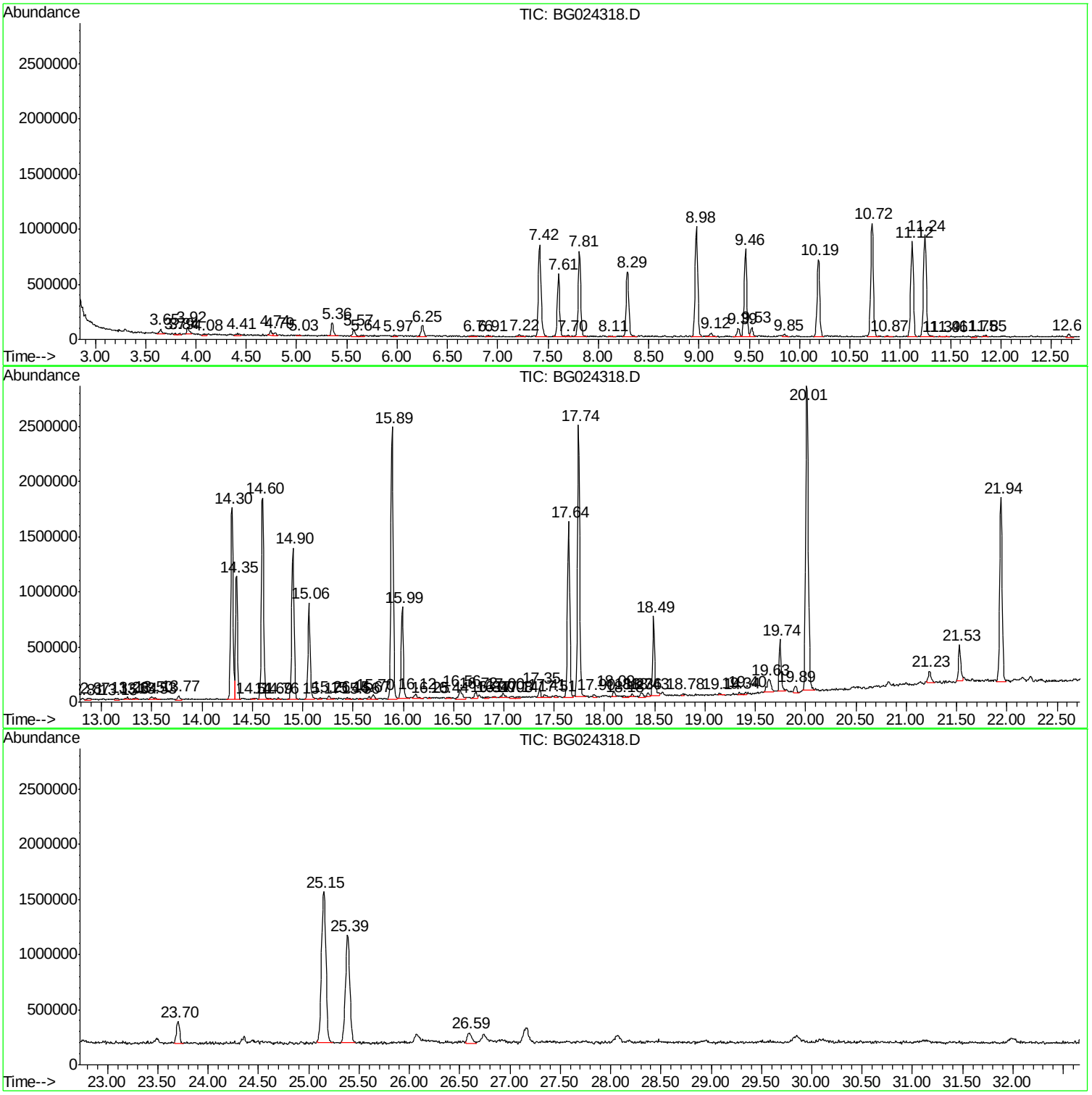
Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BDNH8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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



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Instrument : BNA_G
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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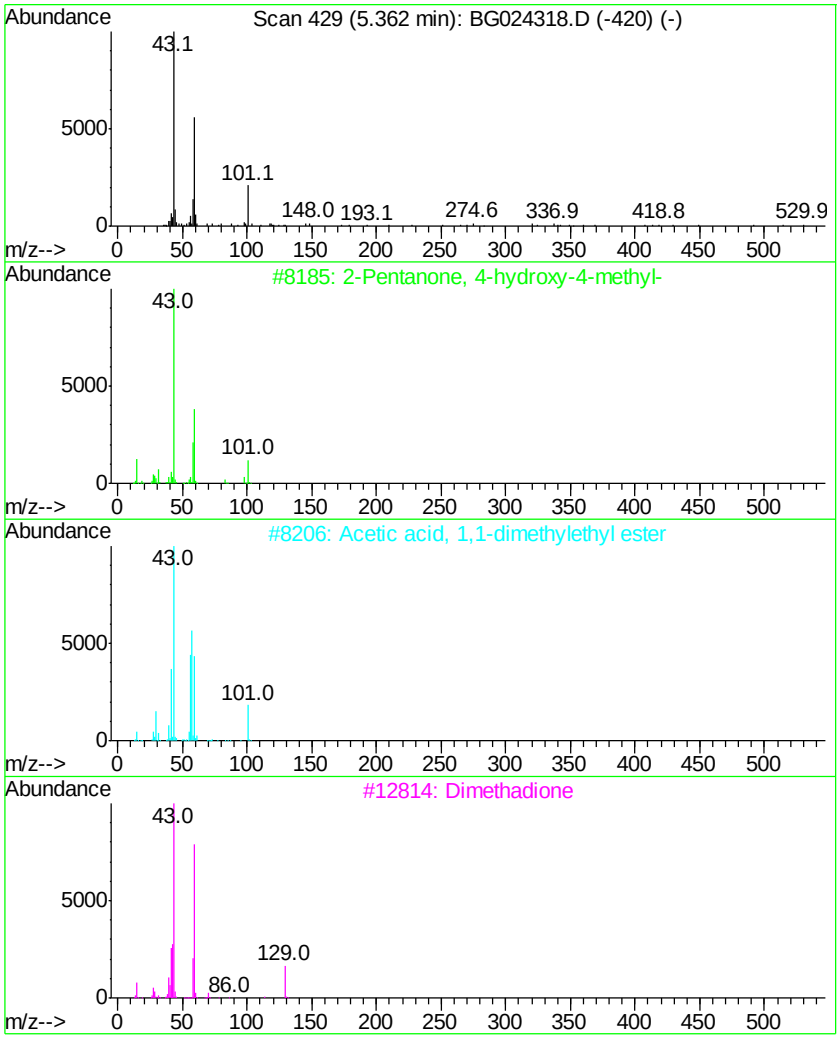
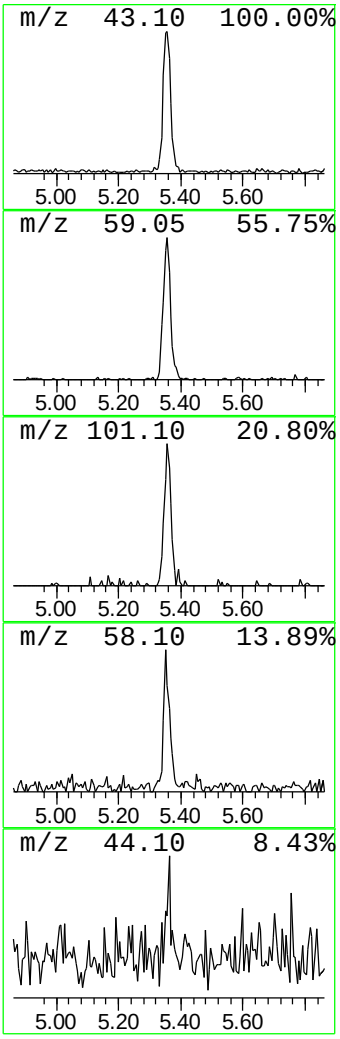
Peak Number 1 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.36	3.77 ng/ul	207942	1,4-Dichlorobenzene-d4	8.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3	Dimethadione	129	C5H7NO3	000695-53-4	32
4	1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	32
5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	25





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101316\
Data File : BG024318.D
Acq On : 13 Oct 2016 13:26
Operator : UM/SJ
Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

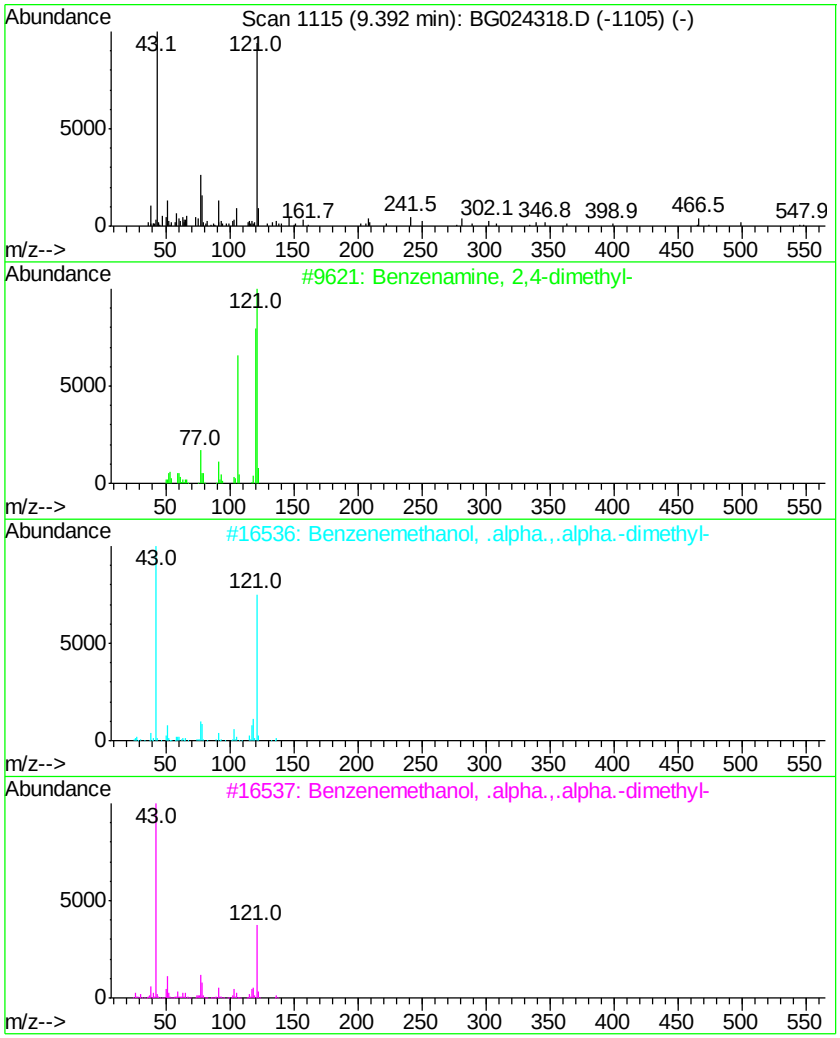
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

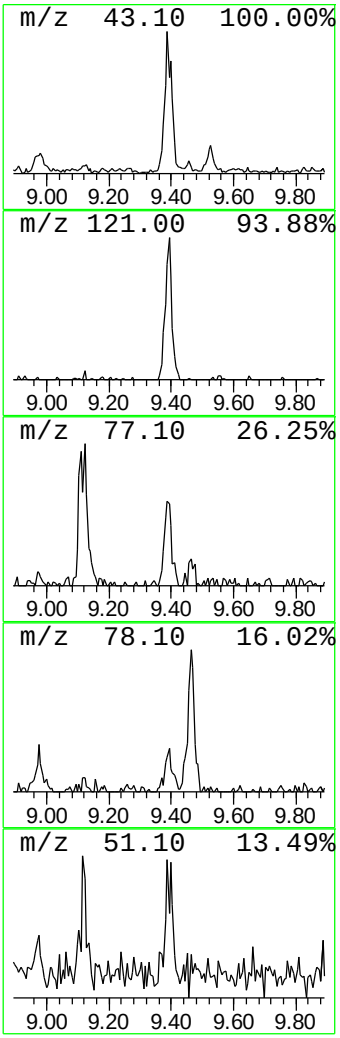
****

Peak Number 4 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.39	2.88 ng/ul	158838	1,4-Dichlorobenzene-d4	8.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,4-dimethyl-	121	C8H11N	000095-68-1	22
2	Benzenemethanol, .alpha.,.alpha....	136	C9H12O	000617-94-7	18
3	Benzenemethanol, .alpha.,.alpha....	136	C9H12O	000617-94-7	18
4	3-Hydroxy-2-methyl-3-phenylbutyr...	208	C12H16O3	057956-39-5	14
5	3-Methoxy-3-phenylpropyl chloride	184	C10H13ClO	1000370-35-5	14





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Acq On : 13 Oct 2016 13:26
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Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

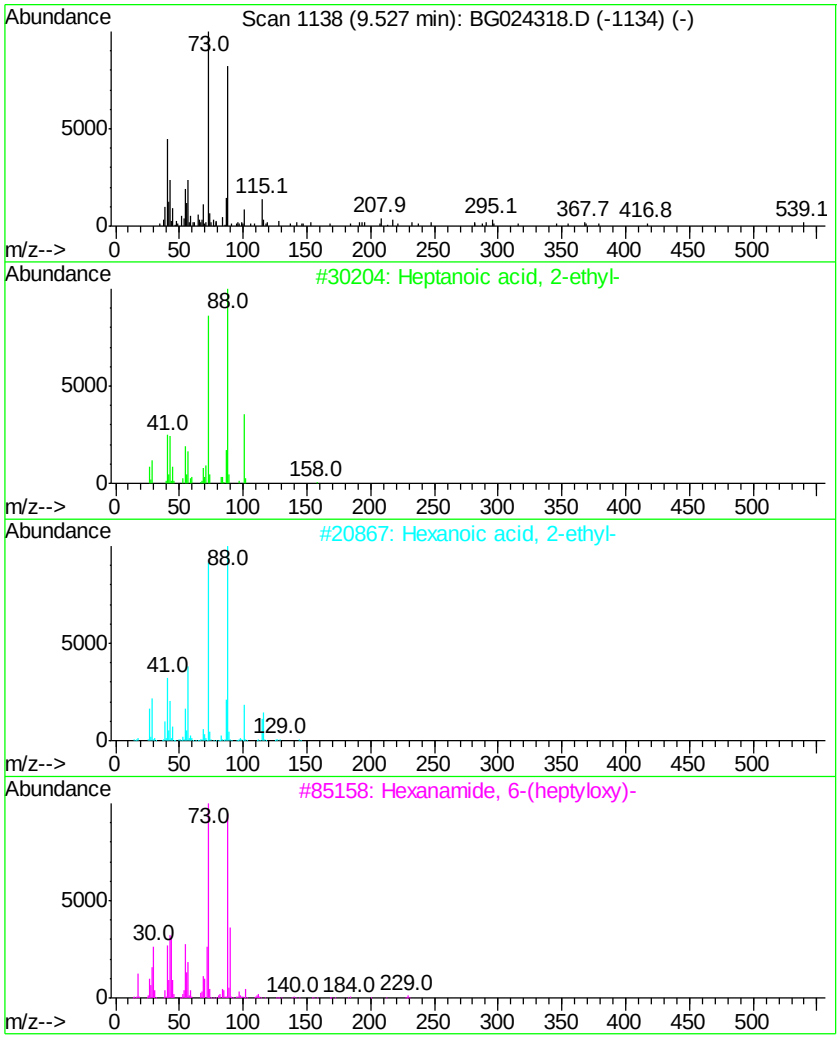
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

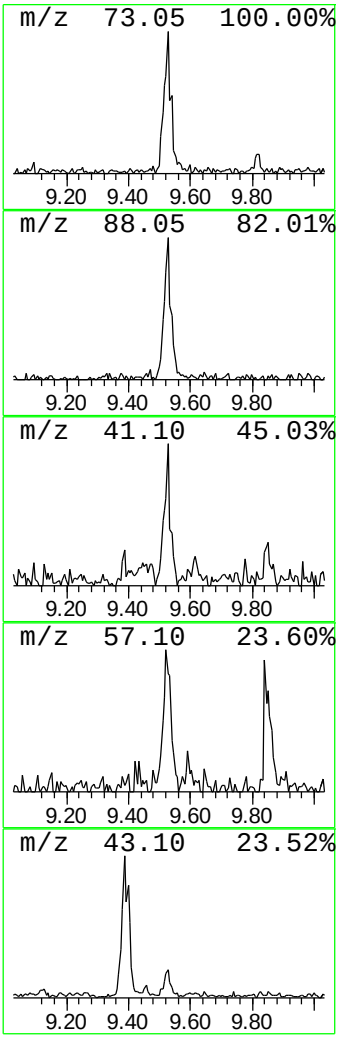
****

Peak Number 5 Heptanoic acid, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	2.61 ng/ul	143799	1,4-Dichlorobenzene-d4	8.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	64
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	59
3	Hexanamide, 6-(heptyloxy)-	229	C13H27NO2	055191-08-7	45
4	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	45
5	1,4-Dioxane	88	C4H8O2	000123-91-1	40





Library Search Compound Report

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Data File : BG024318.D
Acq On : 13 Oct 2016 13:26
Operator : UM/SJ
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Instrument : BNA_G
ClientSampleId : BDNH8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

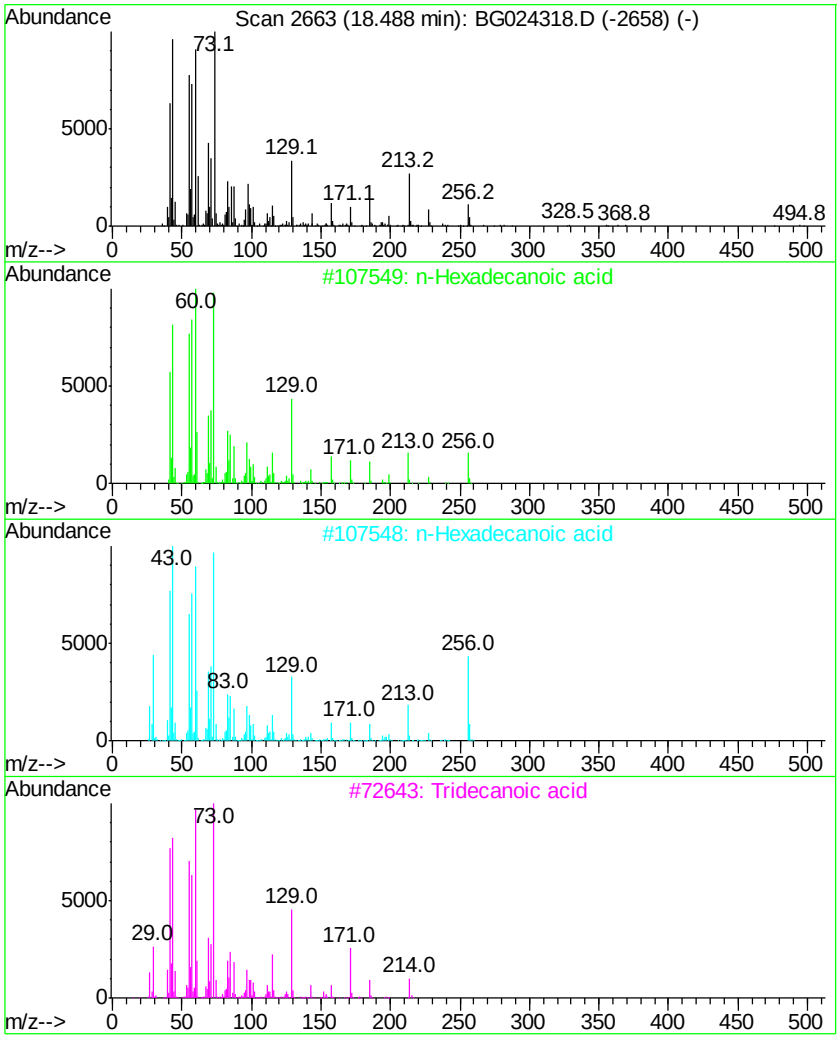
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

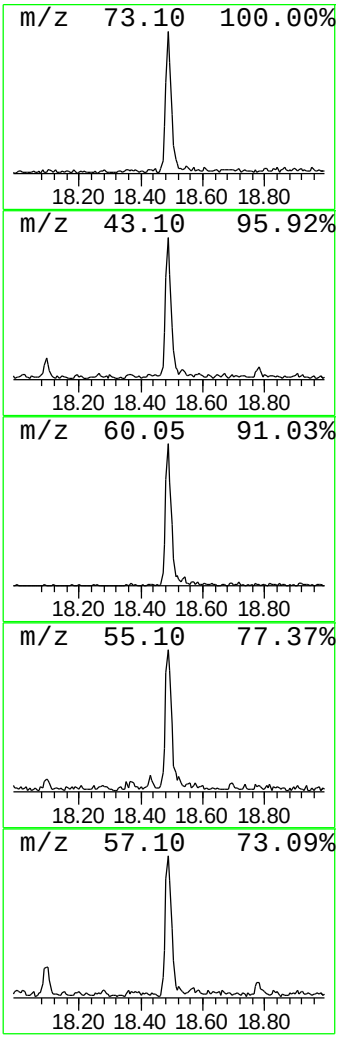
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.49	7.88 ng/ul	988612	Phenanthrene-d10	17.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	Tridecanoic acid	214	C13H26O2	000638-53-9	90
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	80





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Data File : BG024318.D
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Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

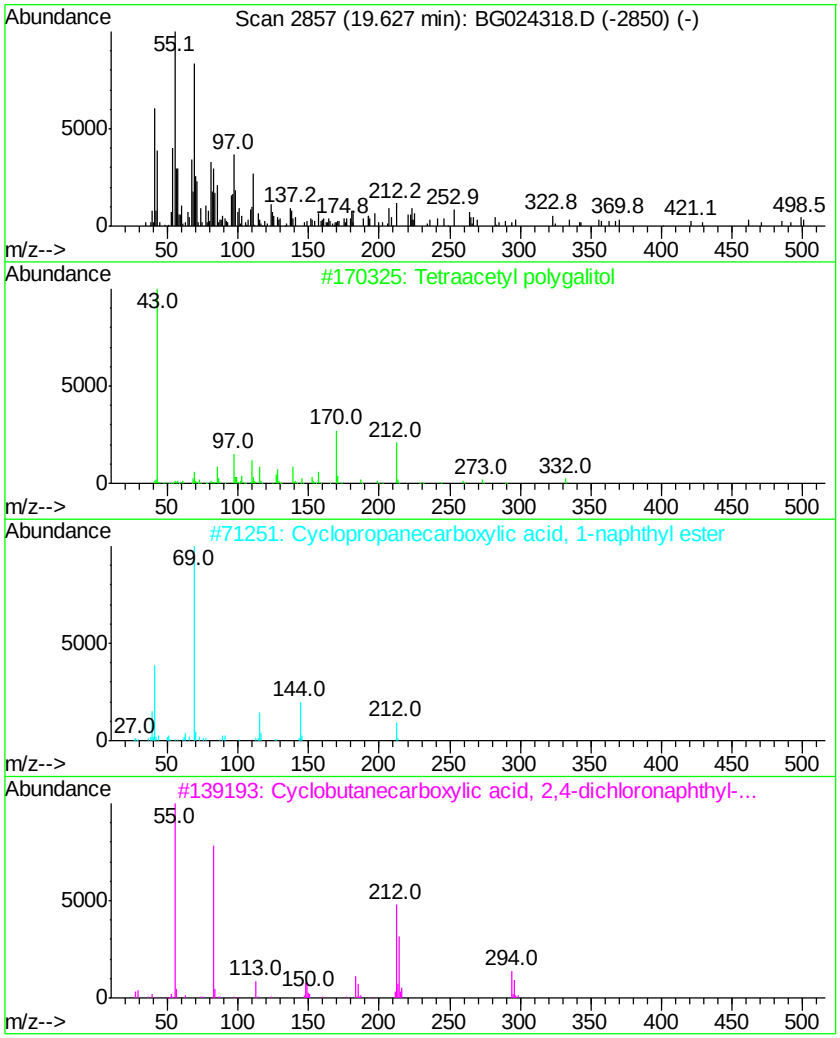
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

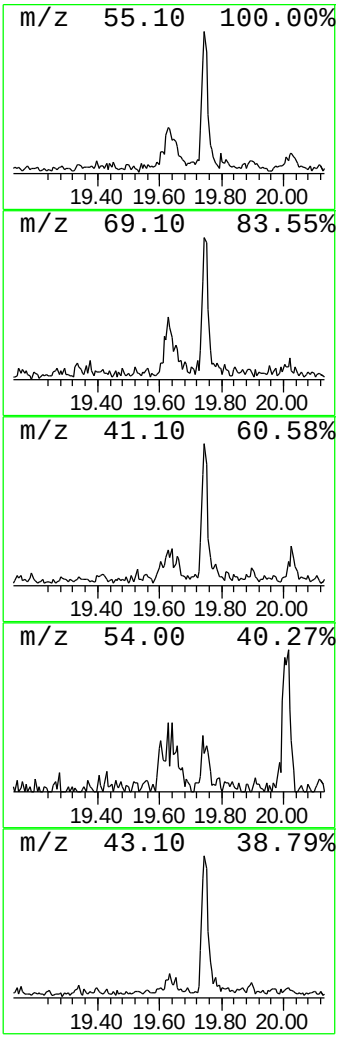
****

Peak Number 7 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.63	2.79 ng/ul	350595	Phenanthrene-d10	17.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetraacetyl polygalitol	332	C14H20O9	1000130-06-9	12
2	Cyclopropanecarboxylic acid, 1-n...	212	C14H12O2	1000279-96-1	12
3	Cyclobutanecarboxylic acid, 2,4-...	294	C15H12Cl2O2	1000331-09-6	10
4	Ethyl-1,1,2,3,3,3-hexafluoroprop...	212	C5H6F6S	000380-35-8	10
5	4,7-Ethanoisobenzofuran-1,3-dion...	264	C14H16O5	052918-82-8	10





Data Path : Z:\HPCHEM1\BNA G\DATA\BG101316\
Data File : BG024318.D
Acq On : 13 Oct 2016 13:26
Operator : UM/SJ
Sample : H5166-08
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDNH8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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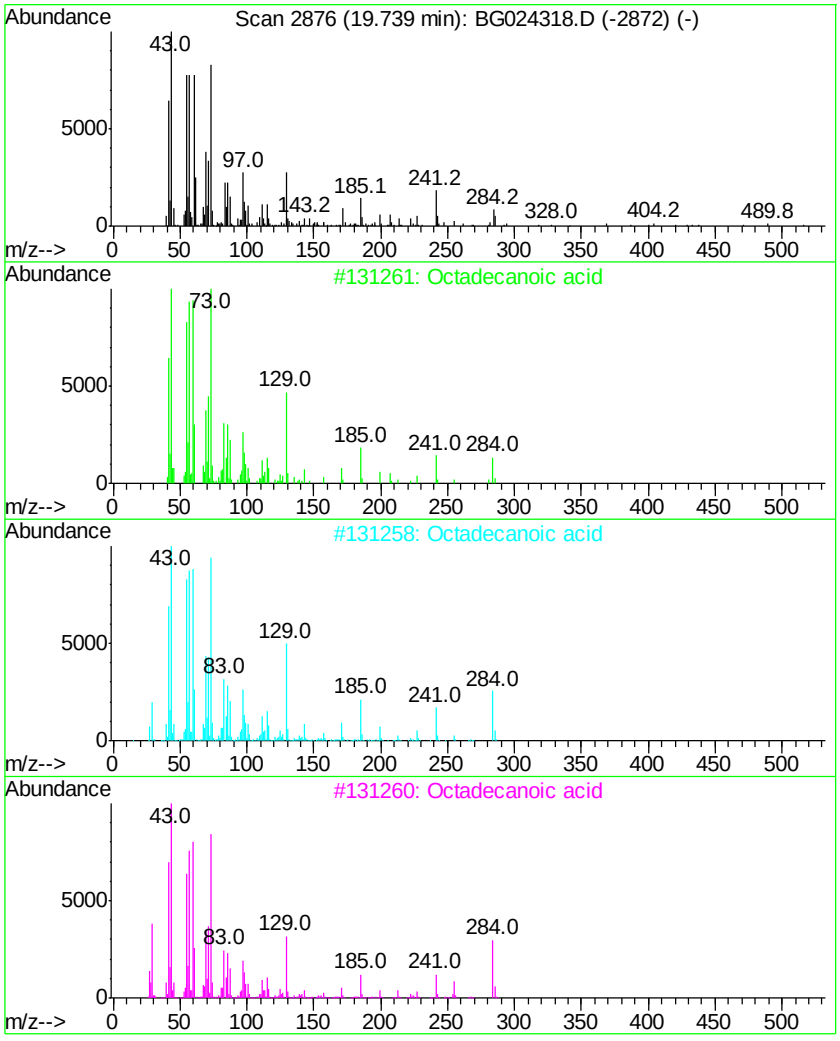
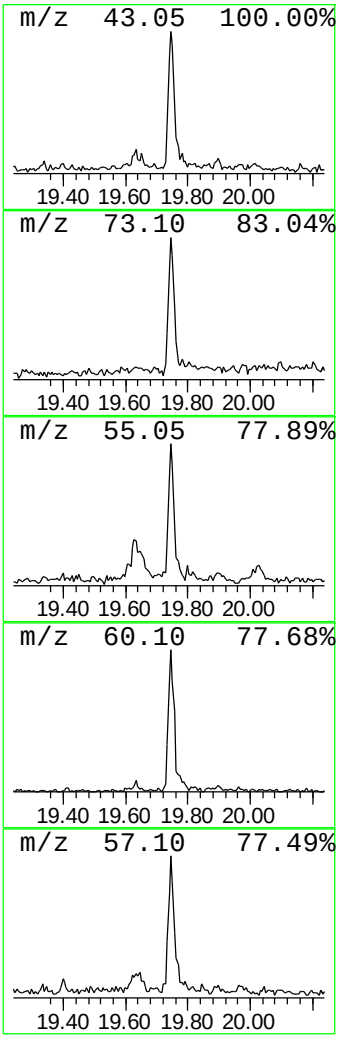
Peak Number 8 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.74	5.16 ng/ul	647996	Phenanthrene-d10	17.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2	Octadecanoic acid	284	C18H36O2	000057-11-4	97
3	Octadecanoic acid	284	C18H36O2	000057-11-4	96
4	Octadecanoic acid	284	C18H36O2	000057-11-4	94
5	Octadecanoic acid	284	C18H36O2	000057-11-4	78





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Data File : BG024318.D
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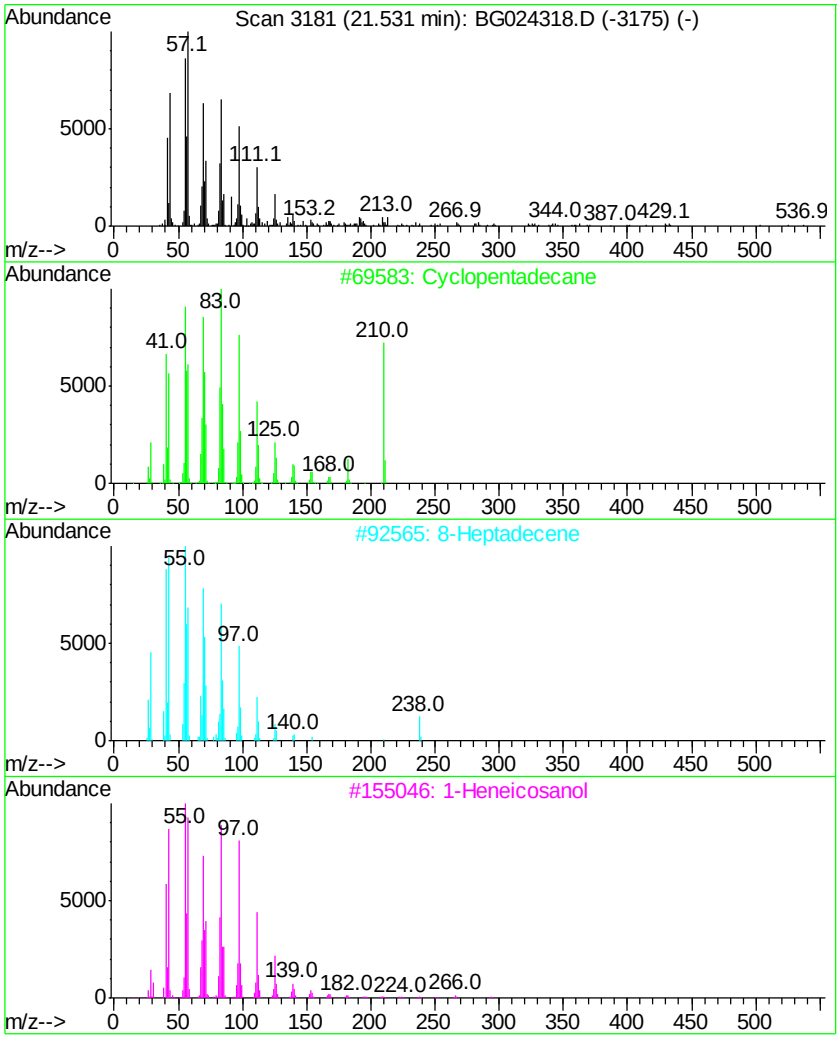
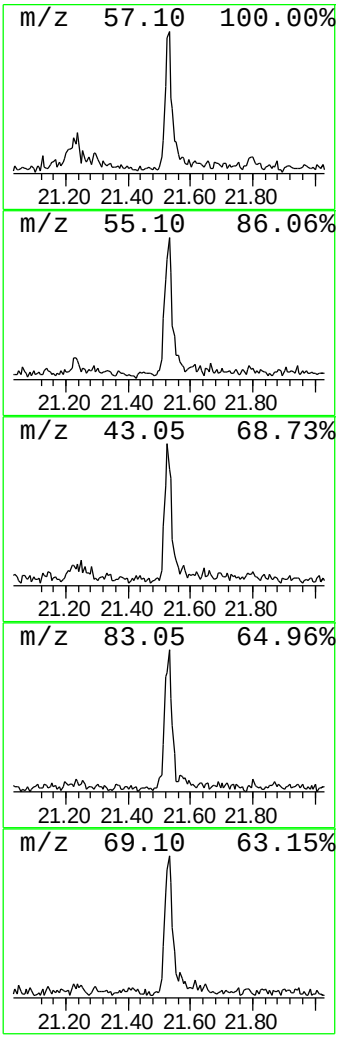
Peak Number 9 (DEL) Alkane: Cyclic21.53 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.53	3.85 ng/ul	556387	Chrysene-d12	21.94

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	83
2	8-Heptadecene	238	C17H34	002579-04-6	83
3	1-Heneicosanol	312	C21H44O	015594-90-8	76
4	1-Pentadecene	210	C15H30	013360-61-7	74
5	Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	74





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Instrument : BNA_G
ClientSampleId : BDNH8

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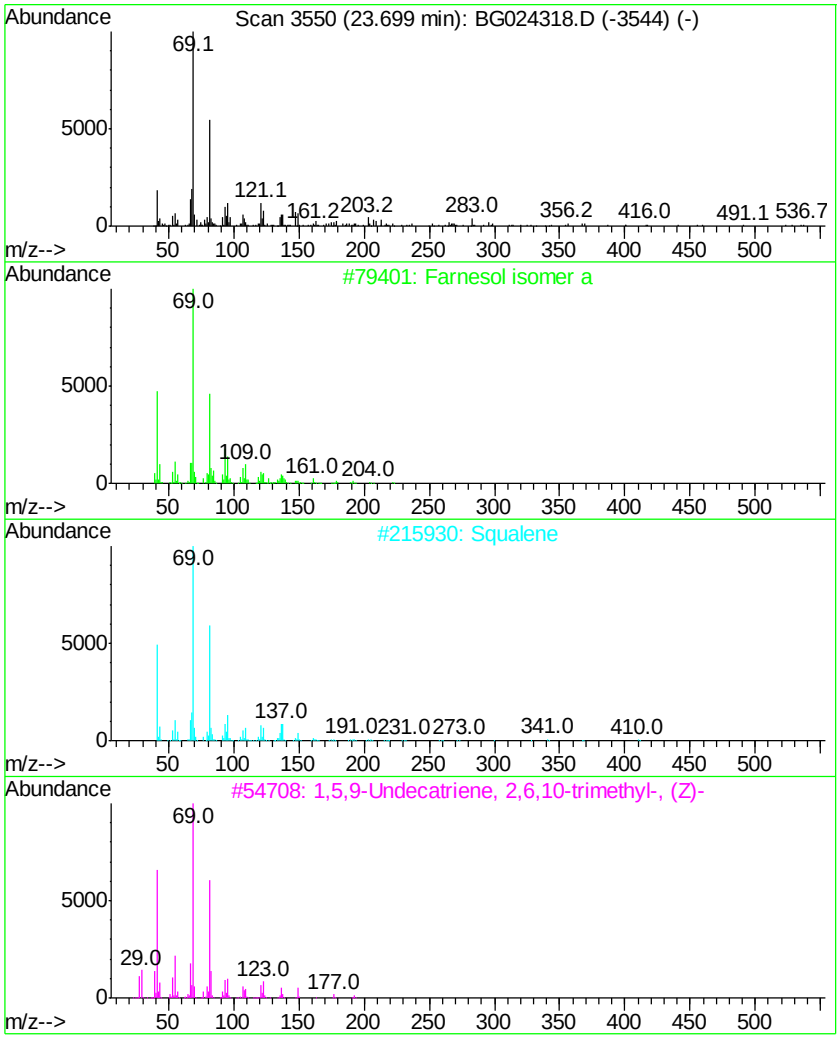
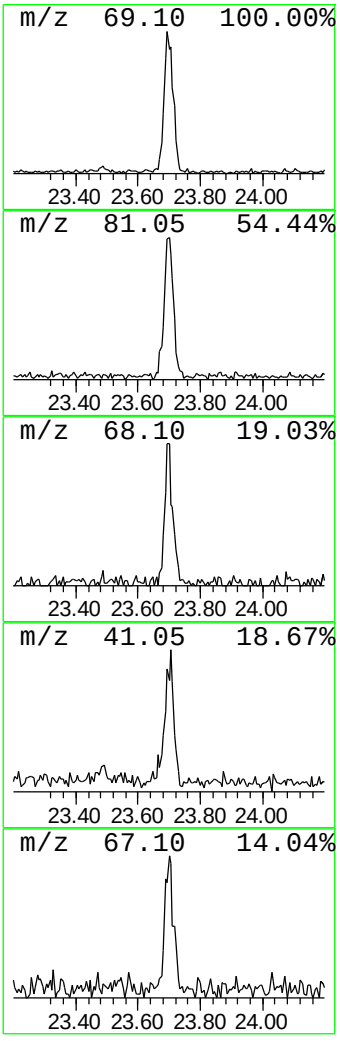
Peak Number 10 Farnesol isomer a Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.70	3.07 ng/ul	449268	Perylene-d12	25.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Farnesol isomer a	222	C15H26O	1000108-92-4	81
2	Squalene	410	C30H50	000111-02-4	72
3	1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	70
4	Squalene	410	C30H50	000111-02-4	64
5	Squalene	410	C30H50	000111-02-4	64





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Instrument : BNA_G
ClientSampleId : BDNH8

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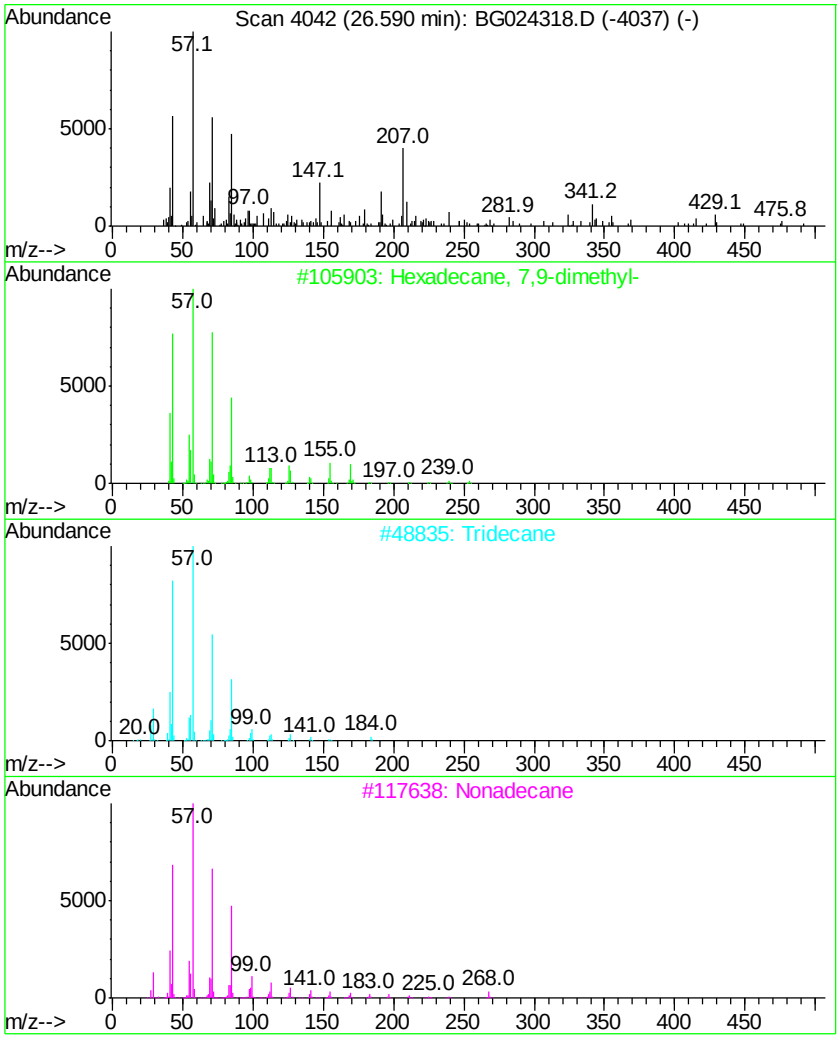
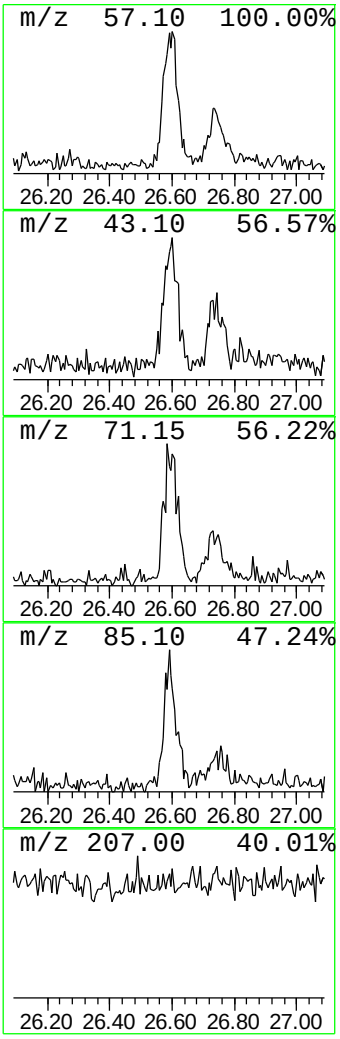
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.59	2.11 ng/ul	309561	Perylene-d12	25.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	46
2	Tridecane	184	C13H28	000629-50-5	41
3	Nonadecane	268	C19H40	000629-92-5	41
4	Tricosane	324	C23H48	000638-67-5	41
5	Hexadecane	226	C16H34	000544-76-3	41





Tentatively Identified Compound (LSC) summary

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Misc :
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Instrument : BNA_G
ClientSampleId : BDNH8

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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
unknown-01	5.36	3.8	ng/ul	207942	1	8.29	1103510	20.0
unknown-02	9.39	2.9	ng/ul	158838	1	8.29	1103510	20.0
Heptanoic acid, 2...	9.53	2.6	ng/ul	143799	1	8.29	1103510	20.0
n-Hexadecanoic acid	18.49	7.9	ng/ul	988612	4	17.64	2510330	20.0
unknown-03	19.63	2.8	ng/ul	350595	4	17.64	2510330	20.0
Octadecanoic acid	19.74	5.2	ng/ul	647996	4	17.64	2510330	20.0
(DEL) Alkane: Cyc...	21.53	3.9	ng/ul	556387	5	21.94	2886810	20.0
Farnesol isomer a	23.70	3.1	ng/ul	449268	6	25.39	2927730	20.0
(DEL) Alkane: Str...	26.59	2.1	ng/ul	309561	6	25.39	2927730	20.0