LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

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-BG080416.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	3.169	53	56	62	rVB6	18301	28396	0.32%	0.062%
2	3.257	68	71	74	rBV5	14522	20606	0.23%	0.045%
3	3.392	90	94	96	rVB5	9703	11724	0.13%	0.025%
4	3.610	130	131	136	rVB5	8216	10794	0.12%	0.023%
5	4.009	196	199	205	rBV4	17426	33170	0.38%	0.072%
6	4.385	260	263	267	rVB5	10231	13914	0.16%	0.030%
7	4.467	272	277	278	rBV4	9683	9727	0.11%	0.021%
8	5.255	407	411	414	rBV6	7866	10896	0.12%	0.024%
9	5.307	416	420	423	rVB5	6354	10480	0.12%	0.023%
10	5.660	475	480	482	rBV5	7253	8900	0.10%	0.019%
11	6.259	581	582	586	rBV4	6861	9746	0.11%	0.021%
12	7.269	747	754	763	rVB	86921	174955	1.99%	0.380%
13	7.422	774	780	788	rBV	66116	122106	1.39%	0.265%
14	7.640	811	817	825	rBV	84986	161150	1.83%	0.350%
15	8.004	876	879	883	rBV6	9723	13303	0.15%	0.029%
16	8.110	888	897	904	rBV	588993	1018111	11.58%	2.211%
17	8.832	1011	1020	1027	rBV2	95280	175098	1.99%	0.380%
18	8.950	1037	1040	1044	rBV5	8027	9803	0.11%	0.021%
19	9.290	1092	1098	1105	rBV	83017	164214	1.87%	0.357%
20	9.472	1125	1129	1134	rVB7	5668	10563	0.12%	0.023%
21	9.972	1211	1214	1216	rBV4	8266	10035	0.11%	0.022%
22	10.019	1216	1222	1229	rVV3	87377	164100	1.87%	0.356%
23	10.095	1232	1235	1240	rBV7	7001	14182	0.16%	0.031%
24	10.577	1311	1317	1326	rBV2	106835	227817	2.59%	0.495%
25	10.947	1371	1380	1386	rBV	917781	1594394	18.13%	3.462%
26	11.000	1386	1389	1395	rVB2	44264	66760	0.76%	0.145%
27	11.088	1399	1404	1414	rBV2	47721	103569	1.18%	0.225%
28	11.998	1553	1559	1561	rBV7	7957	11898	0.14%	0.026%
29	12.427	1629	1632	1635	rVB4	8699	9799	0.11%	0.021%
30	12.609	1657	1663	1670	rVB4	26407	47938	0.55%	0.104%
31	12.827	1694	1700	1704	rBV2	24307	36863	0.42%	0.080%
32	13.320	1782	1784	1790	rVB7	7932	10709	0.12%	0.023%
33	13.937	1886	1889	1890	rVV3	11959	13078	0.15%	0.028%
34	13.949	1890	1891	1896	rVB5	15946	19024	0.22%	0.041%

LSC Area Percent Report

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Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

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-BG080416.M
Title : SVOA CALIBRATION

35	14.101	1911	1917	1922	rBV7	22039	47604	0.54%	0.103%
36	14.178	1922	1930	1935	rVV	225605	376916	4.29%	0.818%
37	14.225	1935	1938	1944	rVB	113920	156278	1.78%	0.339%
38	14.325	1951	1955	1956	rBV4	7986	11984	0.14%	0.026%
39	14.378	1962	1964	1968	rVB5	9480	10130	0.12%	0.022%
40	14.477	1975	1981	1993	rVB	240980	411104	4.67%	0.893%
41	14.648	2007	2010	2015	rBV7	10660	16620	0.19%	0.036%
42	14.783	2022	2033	2040	rVV2	1479318	2403604	27.33%	5.219%
43	14.847	2040	2044	2051	rVB2	72213	122722	1.40%	0.266%
44	14.924	2053	2057	2058	rBV4	9479	10559	0.12%	0.023%
45	15.024	2064	2074	2082	rBV4	66419	154438	1.76%	0.335%
46	15.094	2084	2086	2088	rVV3	16917	14409	0.16%	0.031%
47	15.182	2090	2101	2110	rBV	66304	146218	1.66%	0.317%
48	15.259	2110	2114	2116	rBV4	12717	16425	0.19%	0.036%
49	15.382	2133	2135	2136	rBV2	11925	8815	0.10%	0.019%
50	15.453	2143	2147	2151	rVB7	16303	27046	0.31%	0.059%
51	15.552	2158	2164	2166	rBV7	14924	28320	0.32%	0.061%
52	15.605	2170	2173	2177	rVV5	19229	26810	0.30%	0.058%
53	15.658	2179	2182	2185	rBV5	11041	14770	0.17%	0.032%
54	15.723	2191	2193	2198	rBV6	9659	14662	0.17%	0.032%
55	15.781	2198	2203	2208	rVV	337023	503860	5.73%	1.094%
56	15.834	2208	2212	2217	rVV2	97386	149704	1.70%	0.325%
57	15.917	2221	2226	2230	rBV7	35081	60615	0.69%	0.132%
58	15.999	2238	2240	2246	rVB7	20267	25472	0.29%	0.055%
59	16.134	2259	2263	2270	rVB10	29640	58427	0.66%	0.127%
60	16.275	2285	2287	2290	rBV4	13623	13959	0.16%	0.030%
61	16.469	2318	2320	2323	rBV4	29264	40495	0.46%	0.088%
62	16.727	2362	2364	2365	rBV2	10333	9402	0.11%	0.020%
63	16.821	2377	2380	2381	rBV3	17486	16812	0.19%	0.037%
64	16.886	2389	2391	2395	rBV5	18011	23807	0.27%	0.052%
65	17.039	2415	2417	2418	rBV2	18022	15071	0.17%	0.033%
66	17.268	2453	2456	2461	rBV7	38985	73083	0.83%	0.159%
67	17.326	2464	2466	2467	rBV2	28812	18635	0.21%	0.040%
68	17.362	2470	2472	2473	rBV2	38074	25329	0.29%	0.055%
69	17.550	2498	2504	2508	rBV	1789799	2836256	32.25%	6.158%
70	17.591	2508	2511	2516	rVB	812901	1106190	12.58%	2.402%
71	17.644	2516	2520	2524	rBV2	326516	525626	5.98%	1.141%

LSC Area Percent Report

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Misc :
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Instrument : BNA_G
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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Title : SVOA CALIBRATION

72	17.679	2524	2526	2531	rVB	162662	221602	2.52%	0.481%
73	18.354	2639	2641	2643	rBV3	31682	36774	0.42%	0.080%
74	18.425	2647	2653	2658	rVV3	295737	589484	6.70%	1.280%
75	18.478	2658	2662	2666	rVB	184971	292237	3.32%	0.635%
76	18.554	2672	2675	2681	rBV2	45377	85505	0.97%	0.186%
77	18.619	2681	2686	2690	rBV2	176605	338515	3.85%	0.735%
78	18.924	2734	2738	2742	rBV	99377	144339	1.64%	0.313%
79	18.971	2742	2746	2747	rBV4	77365	105099	1.20%	0.228%
80	19.336	2805	2808	2813	rBV3	130588	185388	2.11%	0.403%
81	19.606	2849	2854	2859	rVB	1111525	1544181	17.56%	3.353%
82	19.688	2864	2868	2877	rVB3	548330	924058	10.51%	2.006%
83	19.941	2906	2911	2913	rBV2	561421	858392	9.76%	1.864%
84	19.970	2913	2916	2924	rVB	1002626	1392532	15.83%	3.024%
85	20.035	2924	2927	2933	rBV4	82764	114453	1.30%	0.249%
86	20.458	2996	2999	3004	rVB2	189152	294380	3.35%	0.639%
87	20.557	3013	3016	3019	rBV3	139349	177758	2.02%	0.386%
88	20.763	3048	3051	3054	rVB2	113925	133061	1.51%	0.289%
89	20.927	3073	3079	3081	rBV2	493822	892573	10.15%	1.938%
90	20.992	3084	3090	3098	rVB3	683448	2224206	25.29%	4.829%
91	21.433	3159	3165	3170	rBV2	625065	1271857	14.46%	2.761%
92	21.509	3173	3178	3183	rVV	818683	1462799	16.63%	3.176%
93	21.574	3183	3189	3199	rVB	5442071	8794218	100.00%	19.094%
94	21.867	3231	3239	3245	rBV3	1977265	4271675	48.57%	9.275%
95	21.920	3245	3248	3259	rVB	484598	903797	10.28%	1.962%
96	22.672	3374	3376	3380	rVB	131109	155466	1.77%	0.338%
97	24.188	3626	3634	3637	rBV3	600670	1566375	17.81%	3.401%
98	25.263	3809	3817	3827	rBV	1114064	3206136	36.46%	6.961%

Sum of corrected areas: 46056859

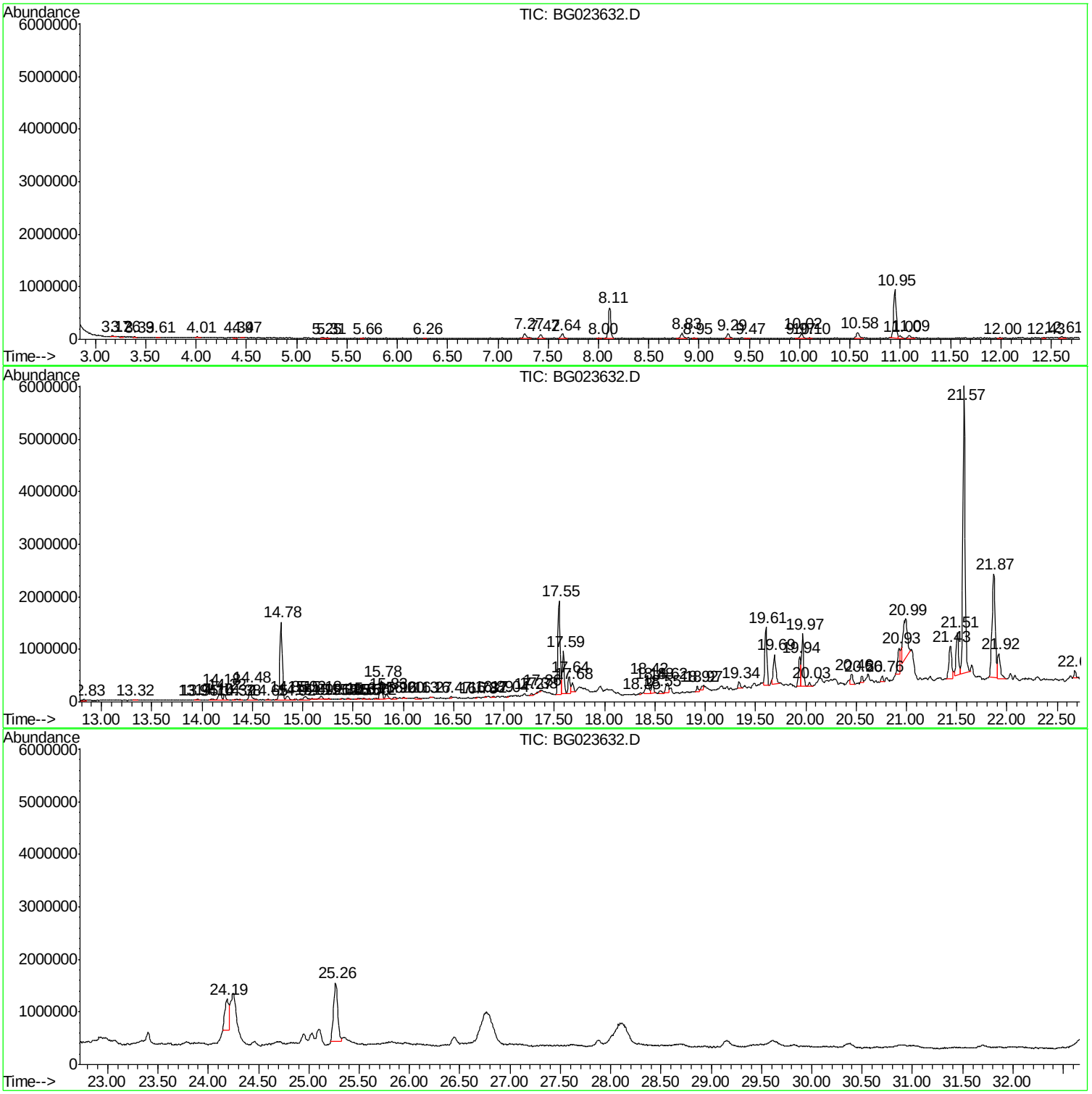
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
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Operator : UM/SJ
Sample : H4360-17 4X
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Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

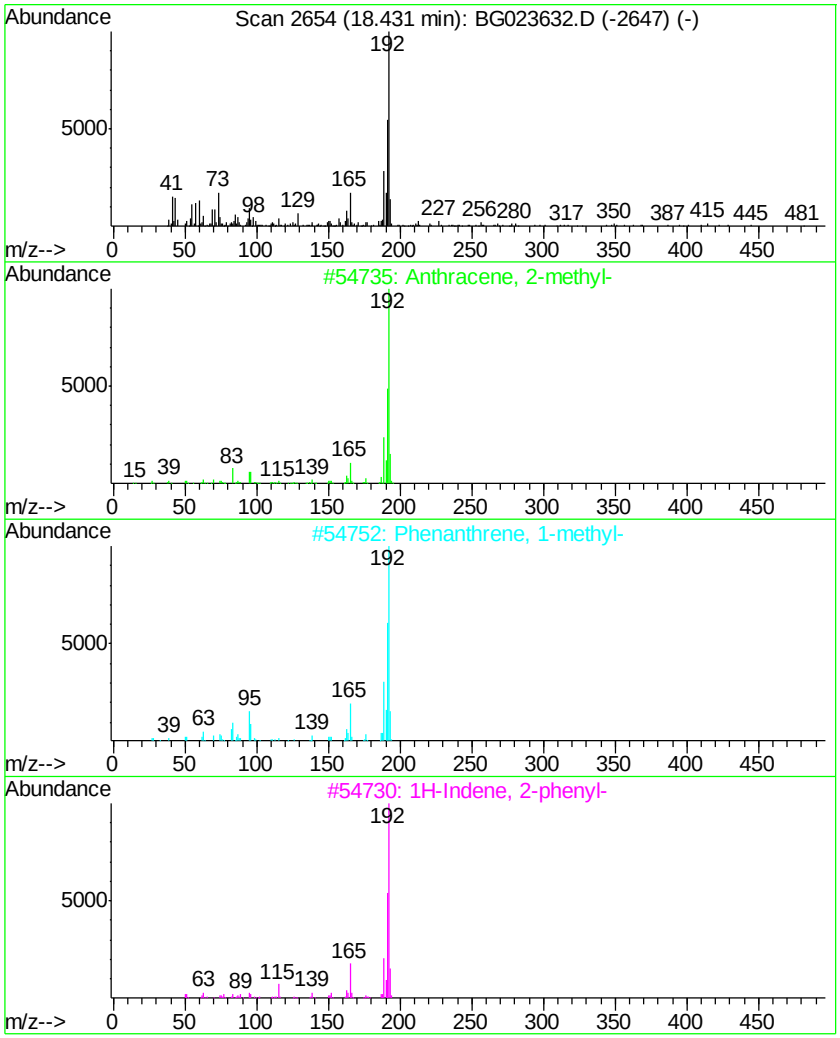
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

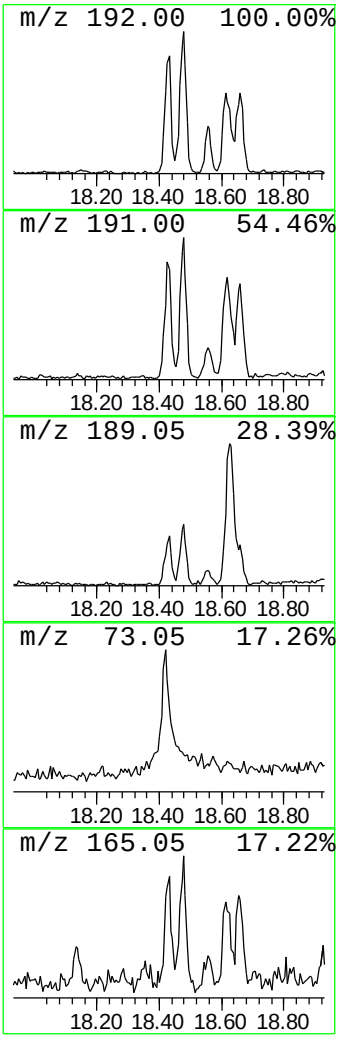
****

Peak Number 1 Anthracene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.43	4.16 ng/ul	589484	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	83
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

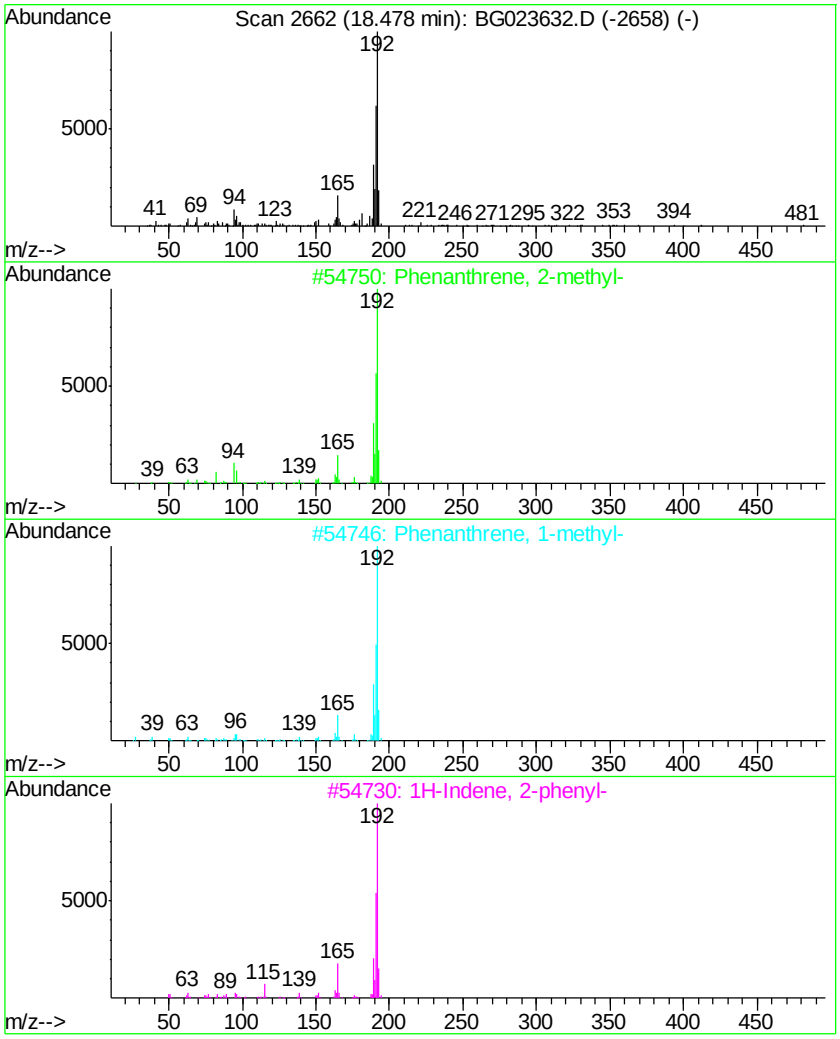
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

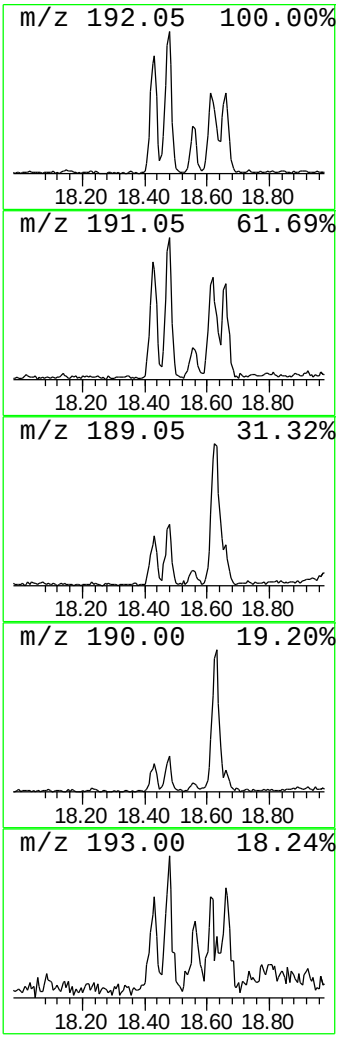
****

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	2.06 ng/ul	292237	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	90





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Sample : H4360-17 4X
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Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

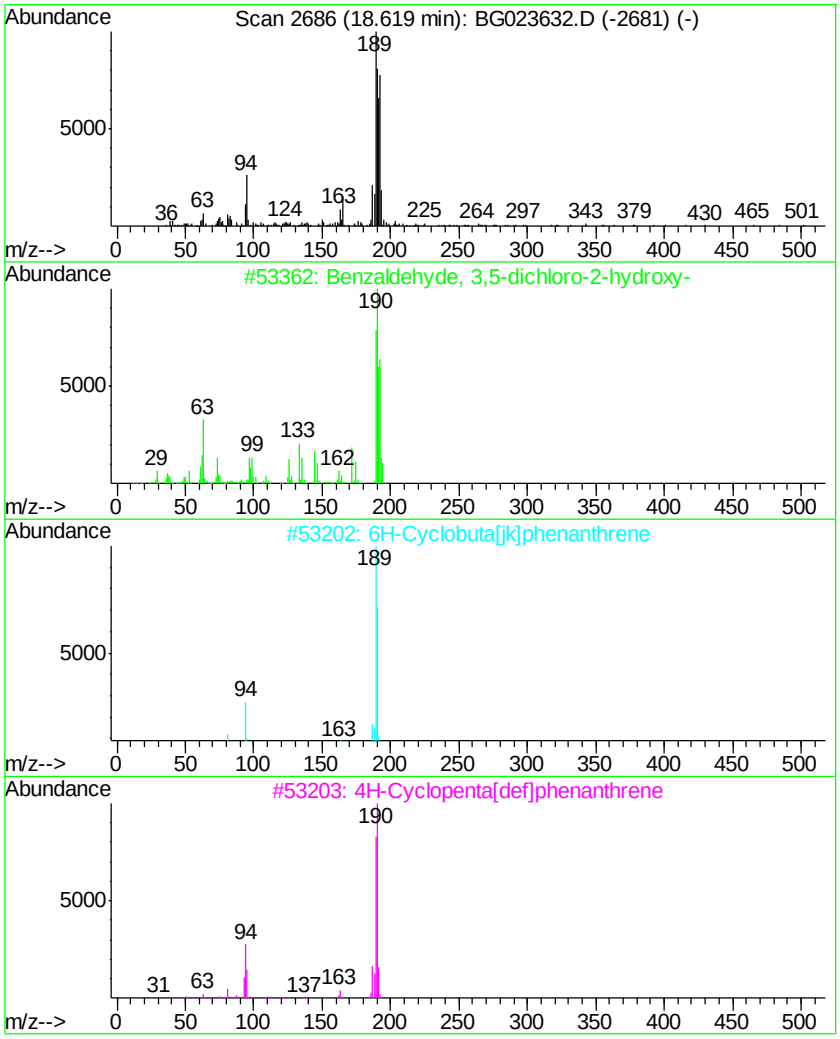
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

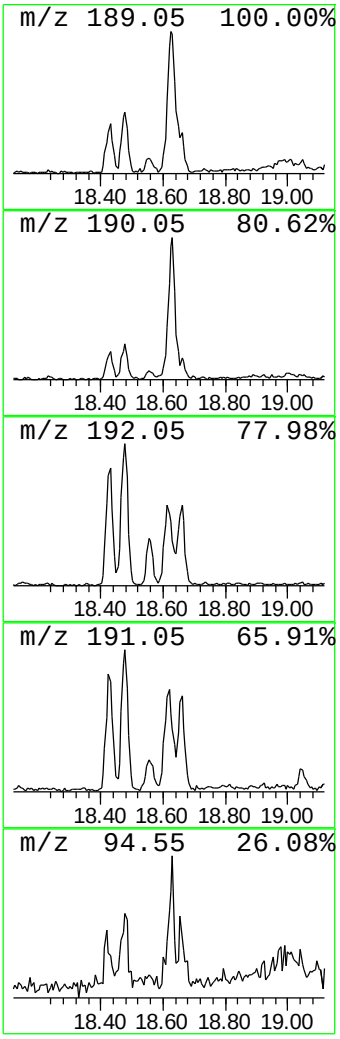
****

Peak Number 3 Benzaldehyde, 3,5-dichloro-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.62	2.39 ng/ul	338515	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	53
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	35
5	Dodecanoic acid, 2,2,2-trichloro...	330	C14H25Cl3O2	1000360-54-3	35





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Sample : H4360-17 4X
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ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

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

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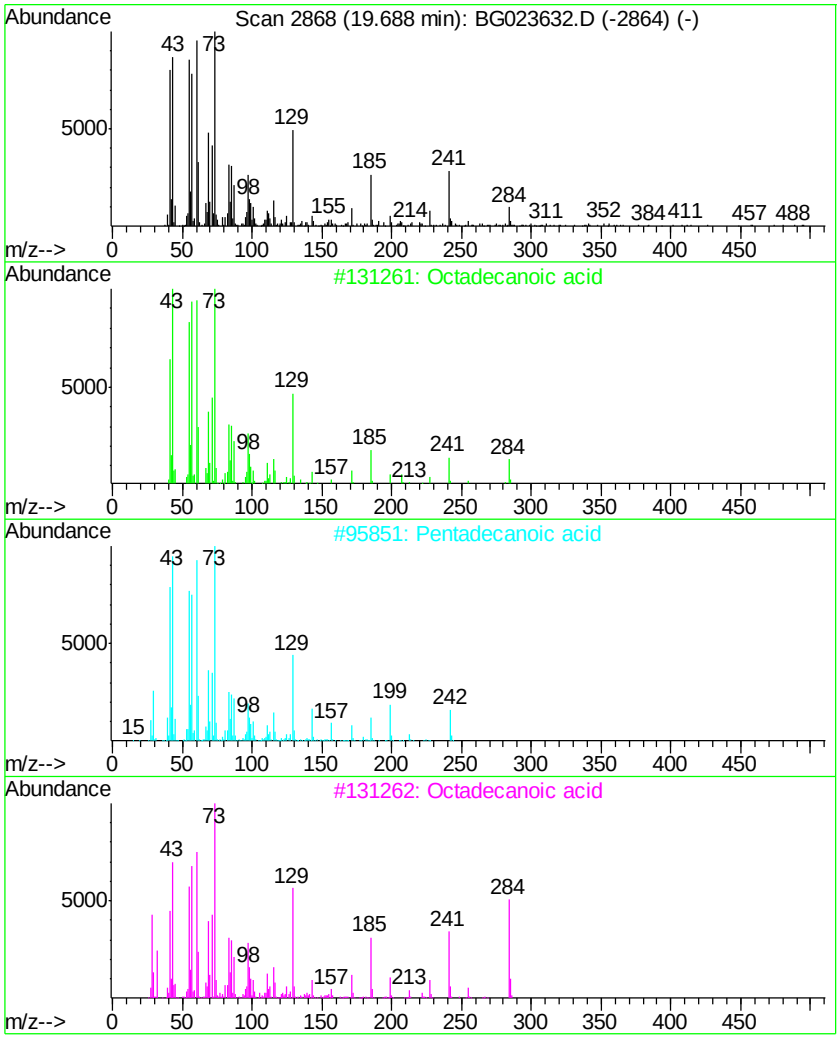
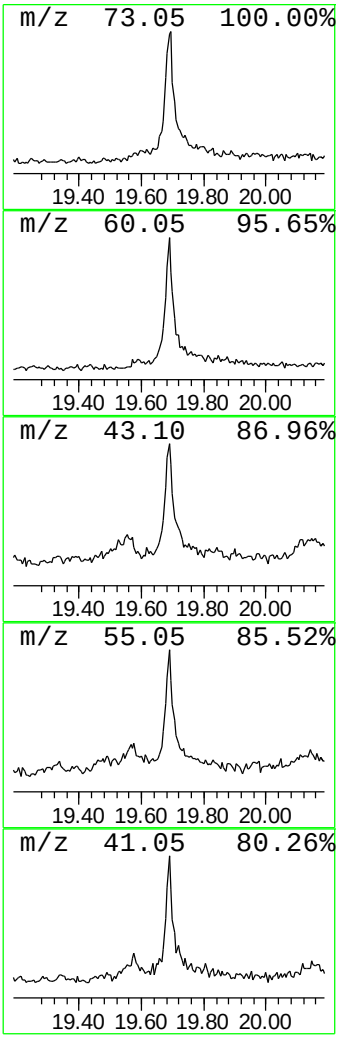
Peak Number 4 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.69	6.52 ng/ul	924058	Phenanthrene-d10	17.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	92
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Octadecanoic acid	284	C18H36O2	000057-11-4	91
4	Octadecanoic acid	284	C18H36O2	000057-11-4	87
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87





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Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

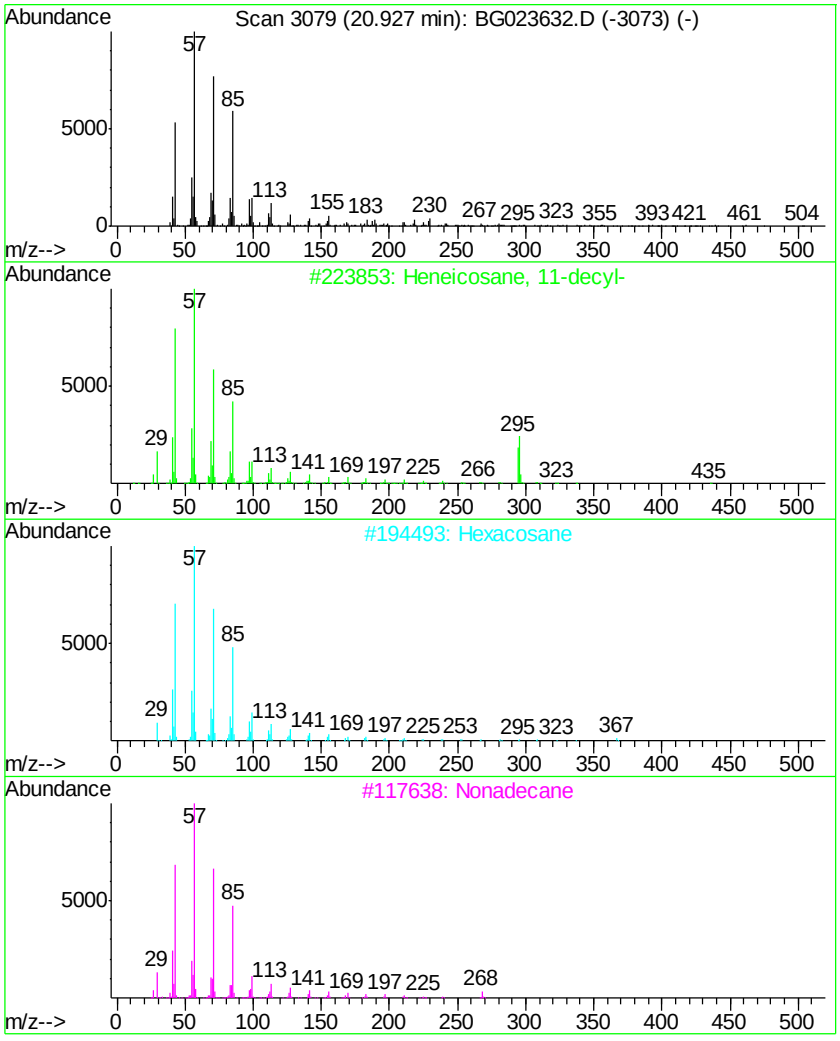
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

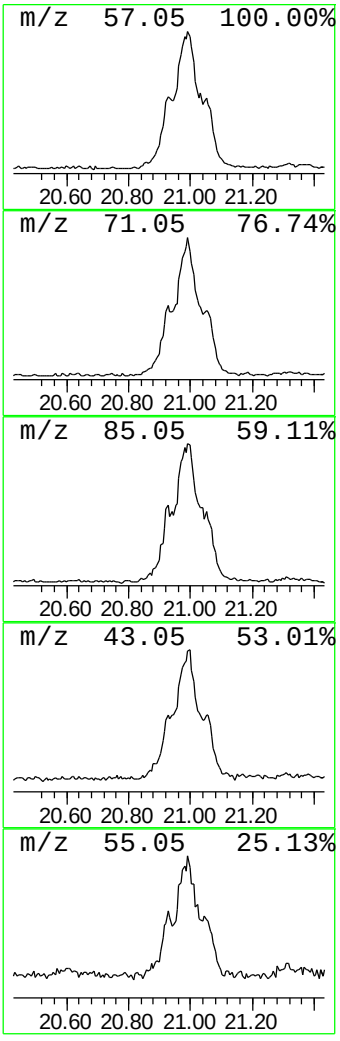
****

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.93	4.18 ng/ul	892573	Chrysene-d12	21.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	93
2	Hexacosane	366	C26H54	000630-01-3	90
3	Nonadecane	268	C19H40	000629-92-5	87
4	Heneicosane	296	C21H44	000629-94-7	87
5	Heptadecane	240	C17H36	000629-78-7	87





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Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

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

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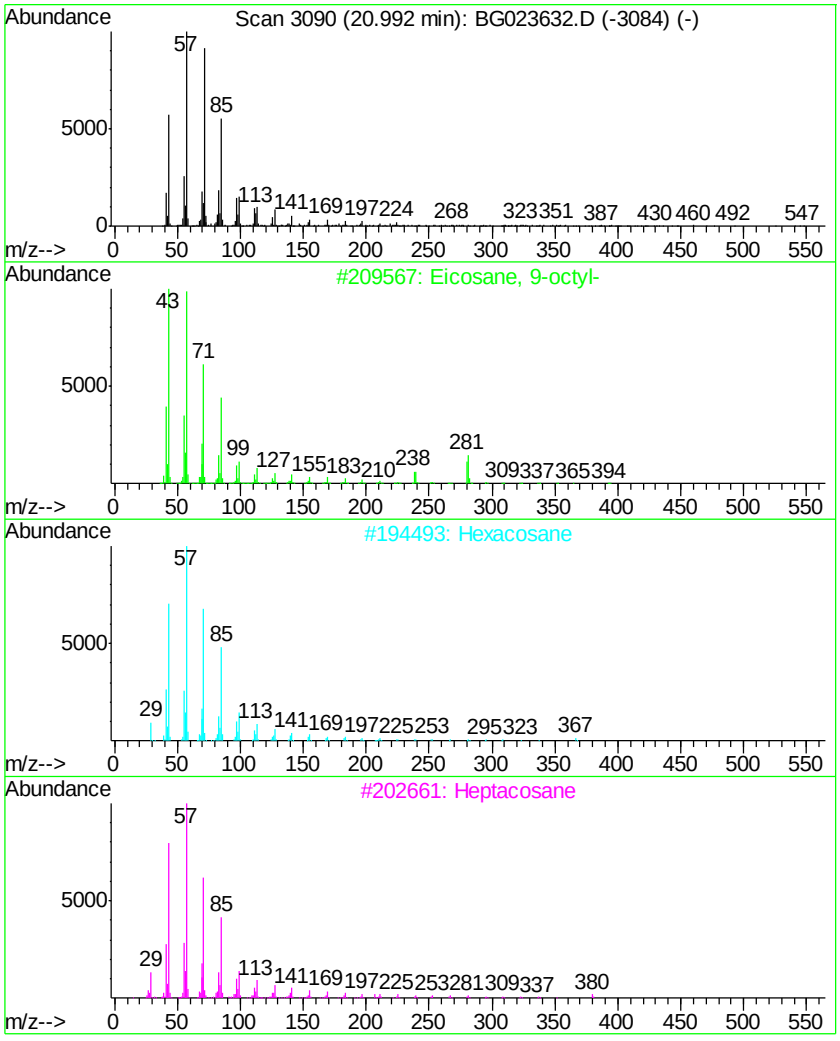
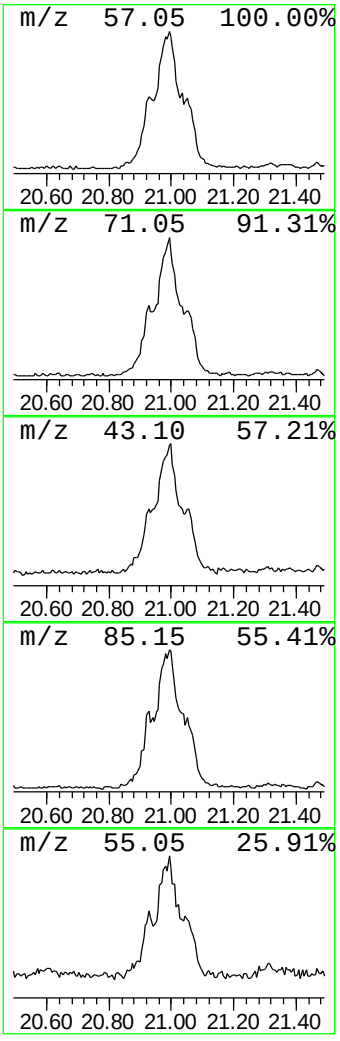
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.99	10.41 ng/ul	2224210	Chrysene-d12	21.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 9-octyl-	394	C28H58	013475-77-9	91
2	Hexacosane	366	C26H54	000630-01-3	91
3	Heptacosane	380	C27H56	000593-49-7	91
4	Docosane, 9-octyl-	422	C30H62	055319-83-0	83
5	Heneicosane	296	C21H44	000629-94-7	80





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

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

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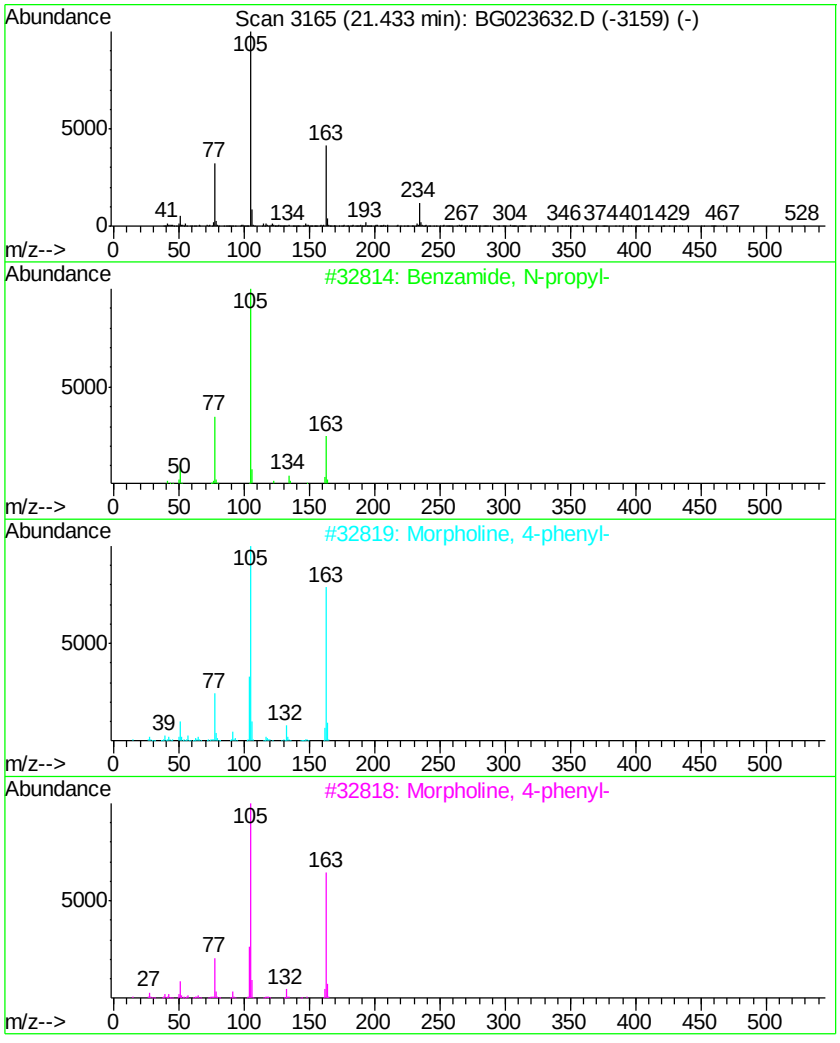
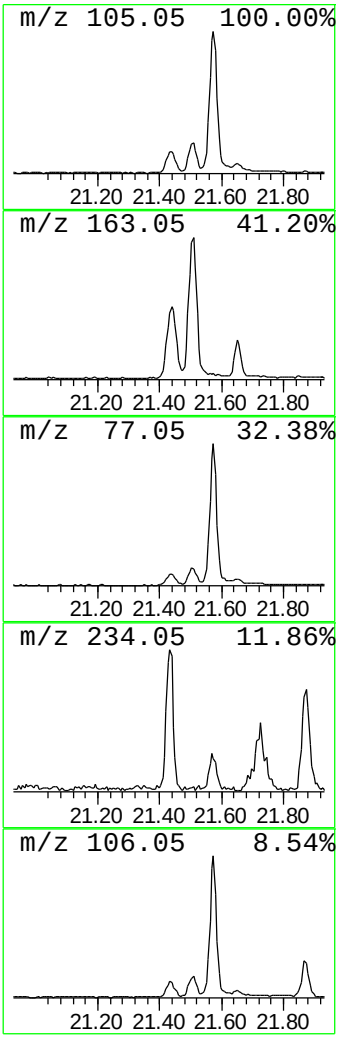
Peak Number 7 Benzamide, N-propyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.43	5.95 ng/ul	1271860	Chrysene-d12	21.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzamide, N-propyl-	163	C10H13NO	010546-70-0	50
2	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	50
3	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	38
4	(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	28
5	1,3-Oxazolidine-4-carboxylic aci...	291	C16H21NO4	1000196-90-4	27





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Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

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

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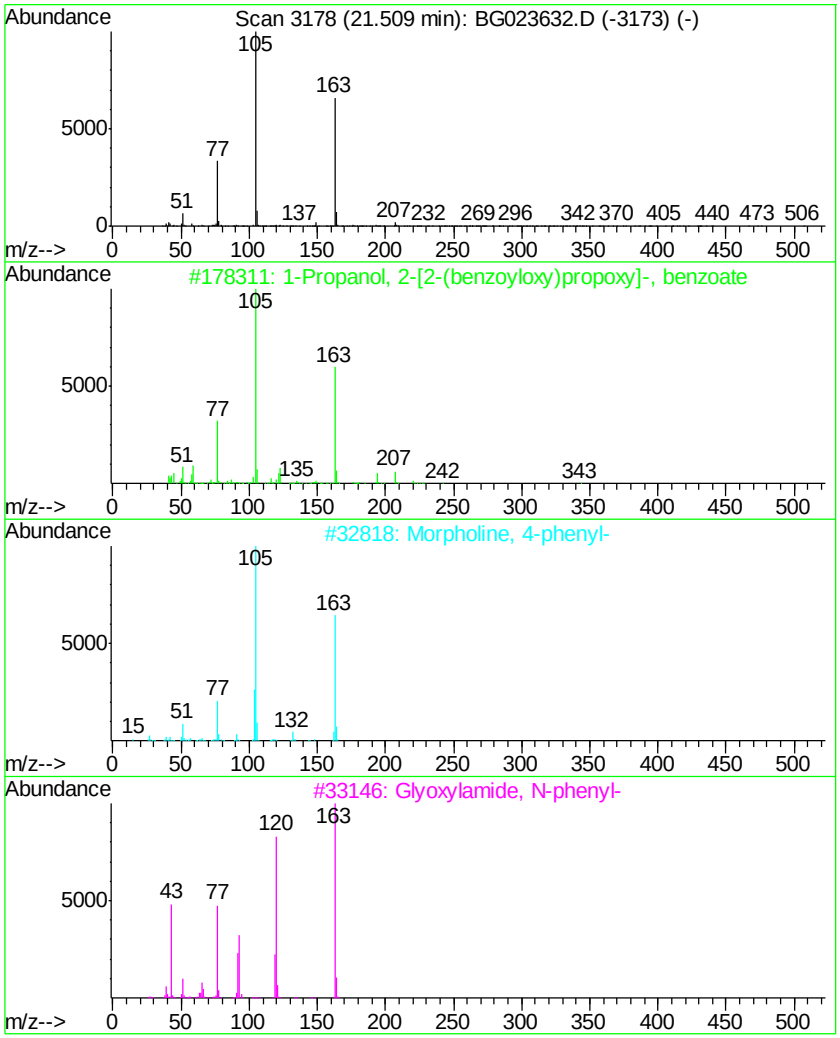
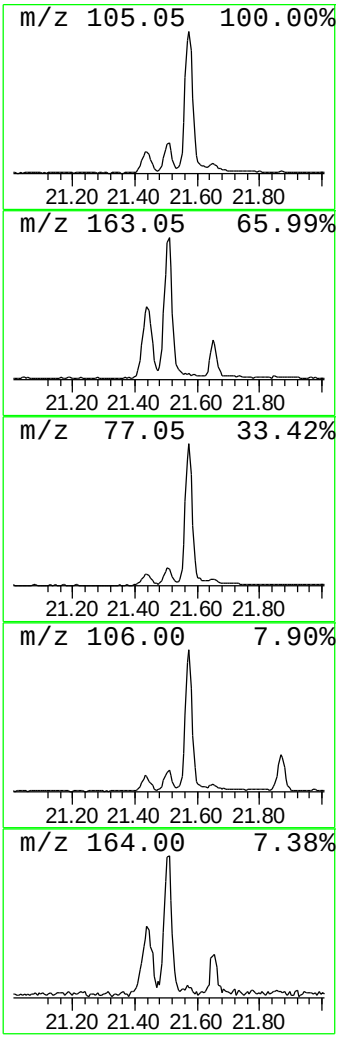
Peak Number 8 1-Propanol, 2-[2-(benzoylox... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.51	6.85 ng/ul	1462800	Chrysene-d12	21.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanol, 2-[2-(benzoyloxy)pro...	342	C20H22O5	020109-39-1	64
2	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	64
3	Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	53
4	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	50
5	Benzoic acid, 2-benzyloxy-3-brom...	348	C17H17BrO3	1000191-33-3	42





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
Quant Title : SVOA CALIBRATION

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

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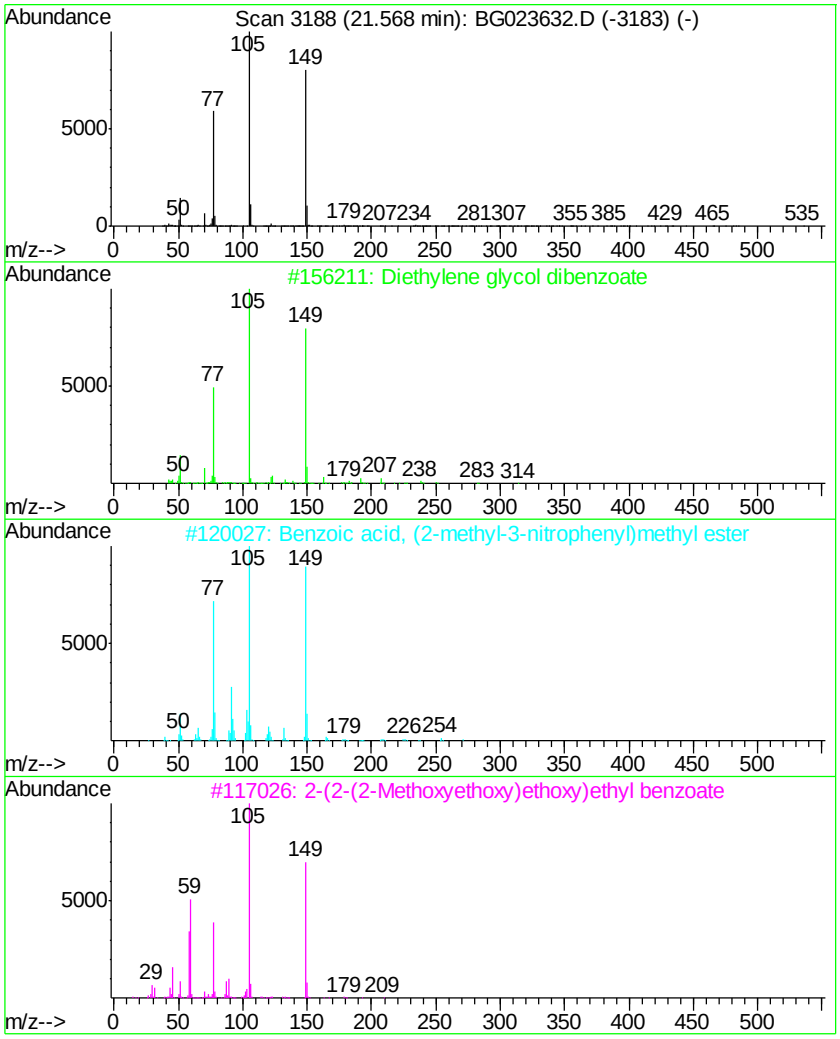
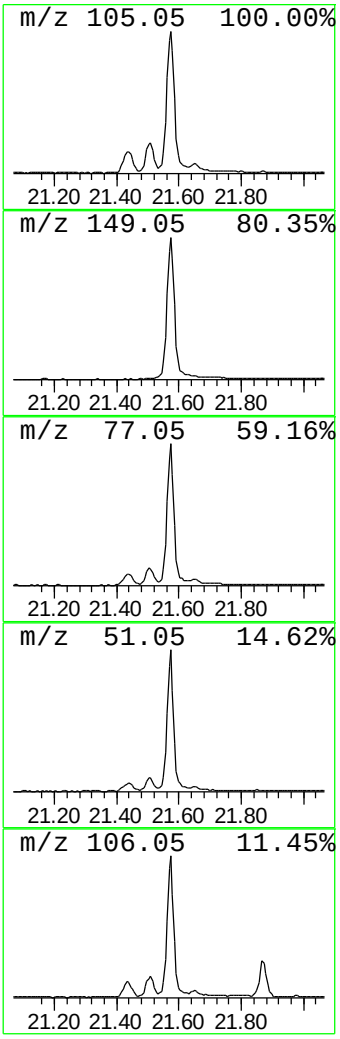
Peak Number 9 Diethylene glycol dibenzoate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.57	41.17 ng/ul	8794220	Chrysene-d12	21.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	78
2	Benzoic acid, (2-methyl-3-nitrop...	271	C15H13NO4	1000367-95-0	74
3	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74
4	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	74
5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081116\
Data File : BG023632.D
Acq On : 11 Aug 2016 21:34
Operator : UM/SJ
Sample : H4360-17 4X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : PR001-SS003-2430-01

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.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
Anthracene, 2-met...	18.43	4.2	ng/ul	589484	4	17.55	2836260	20.0
Phenanthrene, 2-m...	18.48	2.1	ng/ul	292237	4	17.55	2836260	20.0
Benzaldehyde, 3,5...	18.62	2.4	ng/ul	338515	4	17.55	2836260	20.0
Octadecanoic acid	19.69	6.5	ng/ul	924058	4	17.55	2836260	20.0
(DEL) Alkane: Str...	20.93	4.2	ng/ul	892573	5	21.87	4271680	20.0
(DEL) Alkane: Str...	20.99	10.4	ng/ul	2224210	5	21.87	4271680	20.0
Benzamide, N-propyl-	21.43	6.0	ng/ul	1271860	5	21.87	4271680	20.0
1-Propanol, 2-[2-...	21.51	6.8	ng/ul	1462800	5	21.87	4271680	20.0
Diethylene glycol...	21.57	41.2	ng/ul	8794220	5	21.87	4271680	20.0