LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_F\METHODS\8270-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.122	3	6	16	rVB	199506	260602	6.79%	0.753%
2	5.116	512	515	524	rBV	75734	110699	2.88%	0.320%
3	5.504	577	581	596	rVB	3252029	3495525	91.09%	10.096%
4	6.522	749	754	770	rBV	3055858	3837562	100.00%	11.083%
5	6.651	772	776	780	rVV	3345898	3365416	87.70%	9.720%
6	6.698	782	784	791	rVB2	48388	51524	1.34%	0.149%
7	6.875	811	814	821	rBV	1130385	1053230	27.45%	3.042%
8	7.033	837	841	848	rBV	2795666	2609122	67.99%	7.536%
9	7.445	907	911	921	rBV	2250605	2221991	57.90%	6.417%
10	8.163	1029	1033	1036	rBV	1467425	1326611	34.57%	3.831%
11	8.880	1150	1155	1161	rBV	57133	66271	1.73%	0.191%
12	8.975	1168	1171	1177	rVB	37266	38732	1.01%	0.112%
13	9.233	1211	1215	1224	rBV	3026552	2947442	76.81%	8.513%
14	9.304	1224	1227	1230	rVV	82088	75547	1.97%	0.218%
15	9.498	1257	1260	1266	rBV3	28197	49570	1.29%	0.143%
16	9.627	1276	1282	1289	rVB	190416	228917	5.97%	0.661%
17	9.780	1305	1308	1313	rBV	79644	80277	2.09%	0.232%
18	9.833	1313	1317	1320	rVB	163265	126017	3.28%	0.364%
19	9.922	1328	1332	1335	rBV	1190733	1101907	28.71%	3.182%
20	9.951	1335	1337	1345	rVB	76074	86856	2.26%	0.251%
21	10.127	1363	1367	1369	rBV2	50627	57651	1.50%	0.167%
22	10.157	1369	1372	1375	rVB4	42710	46780	1.22%	0.135%
23	10.333	1399	1402	1405	rVB	199203	154453	4.02%	0.446%
24	10.469	1422	1425	1432	rVB	51858	65584	1.71%	0.189%
25	10.551	1435	1439	1442	rVV2	65458	72990	1.90%	0.211%
26	10.716	1463	1467	1475	rBV	1364211	1427145	37.19%	4.122%
27	10.804	1480	1482	1492	rVB	151140	220911	5.76%	0.638%
28	11.221	1550	1553	1556	rBV	37568	42234	1.10%	0.122%
29	11.257	1556	1559	1562	rVV	140727	135710	3.54%	0.392%
30	11.416	1582	1586	1588	rBV	929152	915975	23.87%	2.645%
31	11.439	1588	1590	1594	rVV	548479	464984	12.12%	1.343%
32	11.492	1595	1599	1603	rVB	133338	142288	3.71%	0.411%
33	11.651	1621	1626	1629	rBV2	45089	64159	1.67%	0.185%
34	11.680	1629	1631	1634	rVB	46030	38554	1.00%	0.111%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_F\METHODS\8270-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.921	1668	1672	1674	rBV2	122610	152264	3.97%	0.440%
36	11.945	1674	1676	1681	rVB2	122197	124211	3.24%	0.359%
37	11.992	1681	1684	1687	rBV	50561	50024	1.30%	0.144%
38	12.027	1687	1690	1693	rBV2	133102	156651	4.08%	0.452%
39	12.210	1718	1721	1724	rBV	62349	58559	1.53%	0.169%
40	12.245	1725	1727	1732	rVB2	50345	50467	1.32%	0.146%
41	12.463	1762	1764	1769	rBV2	67720	106867	2.78%	0.309%
42	12.563	1777	1781	1784	rBV3	55684	77726	2.03%	0.224%
43	12.633	1789	1793	1796	rBV	988181	864222	22.52%	2.496%
44	12.798	1820	1821	1826	rVB2	75091	67590	1.76%	0.195%
45	12.862	1826	1832	1837	rBV	845544	983244	25.62%	2.840%
46	12.910	1838	1840	1844	rVB2	47327	45202	1.18%	0.131%
47	12.998	1851	1855	1858	rBV	1996009	1810591	47.18%	5.229%
48	13.027	1858	1860	1864	rVB	52923	55659	1.45%	0.161%
49	13.080	1865	1869	1873	rBV2	64859	128760	3.36%	0.372%
50	13.192	1886	1888	1894	rVB2	135705	115283	3.00%	0.333%
51	13.257	1896	1899	1901	rBV2	82362	68791	1.79%	0.199%
52	13.298	1904	1906	1908	rVB	68017	49430	1.29%	0.143%
53	13.421	1925	1927	1932	rBV2	59847	73118	1.91%	0.211%
54	13.709	1973	1976	1979	rBV	76787	77444	2.02%	0.224%
55	13.880	2001	2005	2009	rBV3	184500	239216	6.23%	0.691%
56	14.015	2026	2028	2031	rBV	261919	208739	5.44%	0.603%
57	14.068	2032	2037	2039	rVV2	783310	1081366	28.18%	3.123%
58	14.092	2039	2041	2045	rVB	407252	343090	8.94%	0.991%
59	15.133	2215	2218	2221	rBV	250509	340050	8.86%	0.982%
60	15.574	2291	2293	2297	rVB	275728	312421	8.14%	0.902%

Sum of corrected areas: 34624221

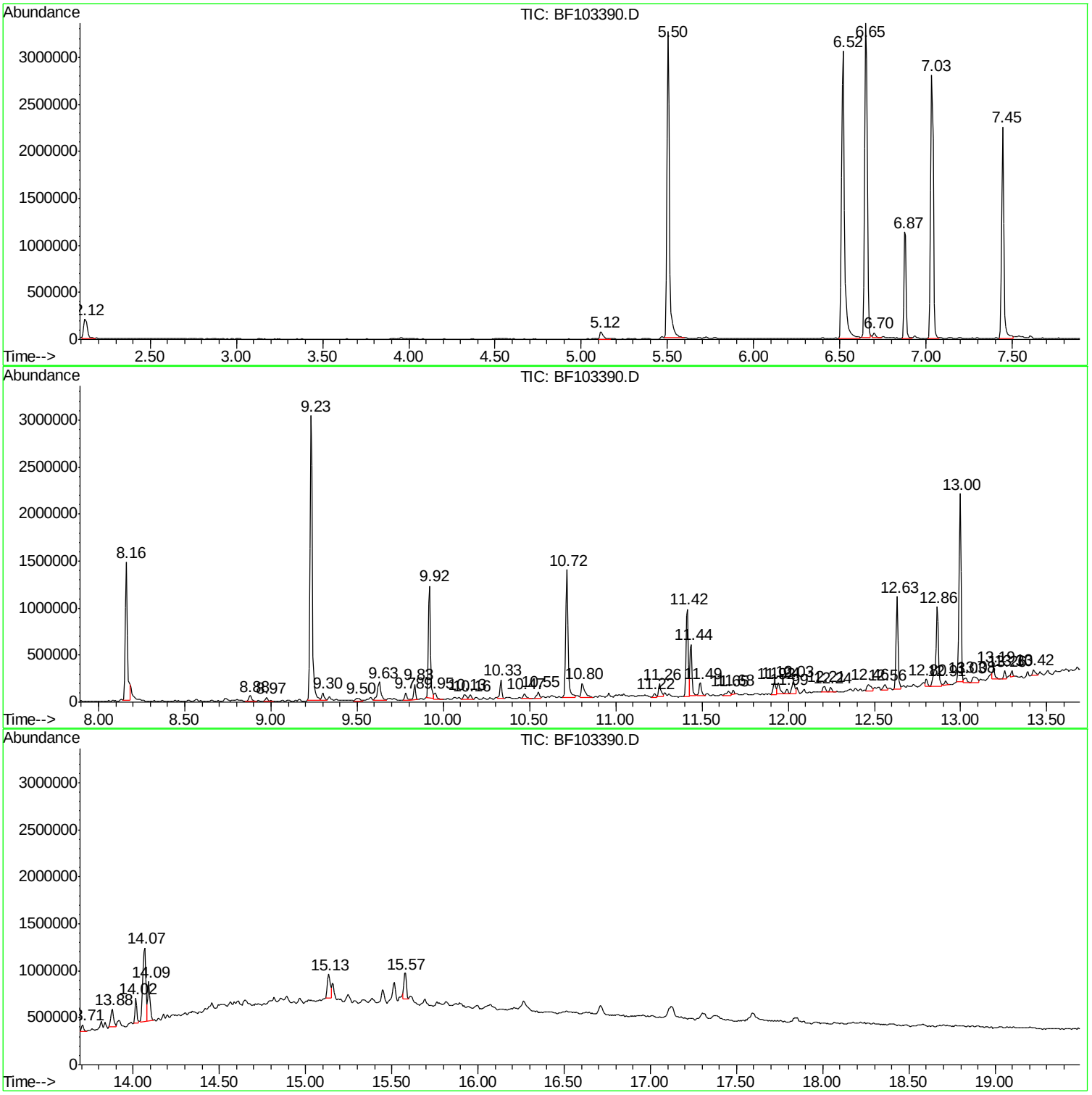
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

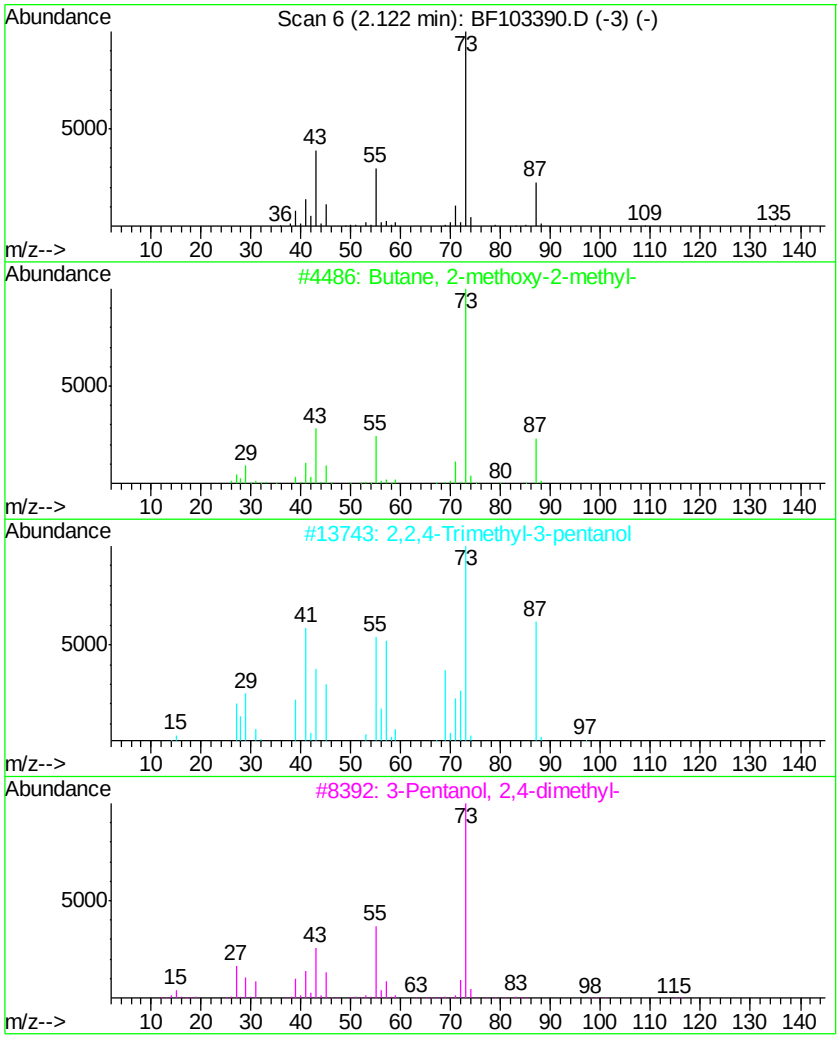
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

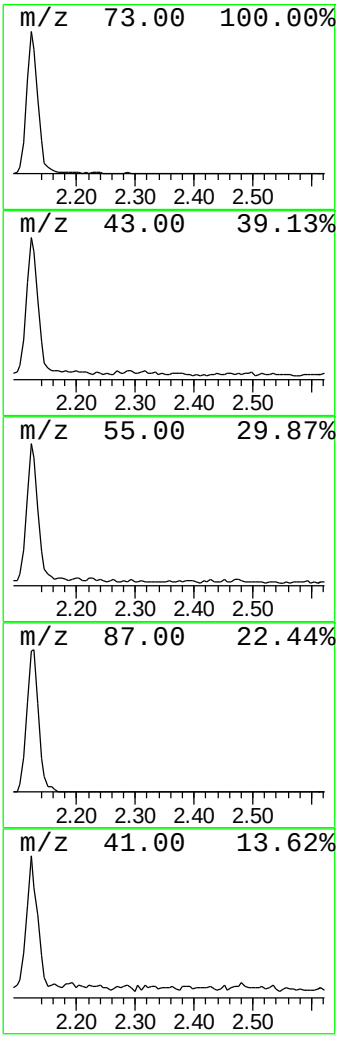
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.12	4.95 ng	260602	1,4-Dichlorobenzene-d4	6.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
5	2,5-Dimethyl-4-hydroxy-3-hexanone	144	C8H16O2	000815-77-0	28





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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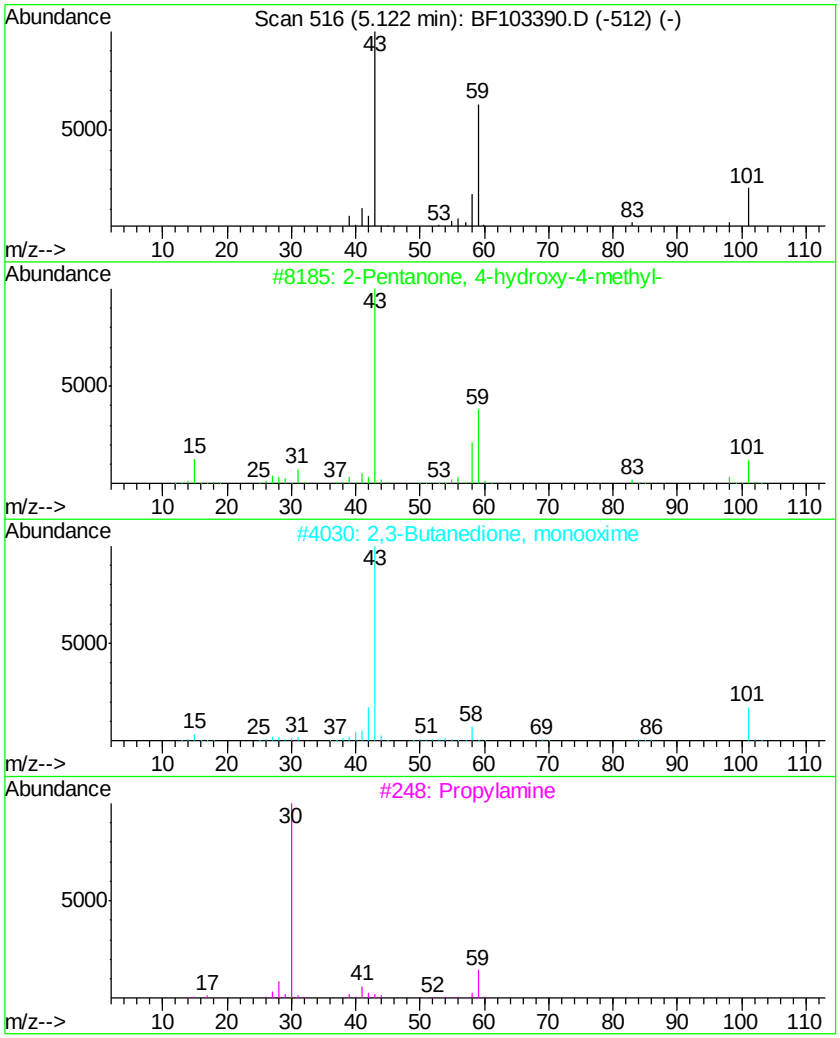
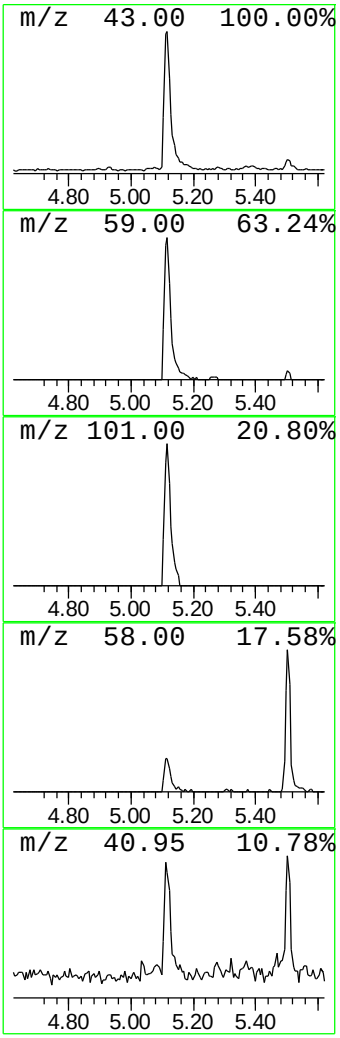
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	2.10 ng	110699	1,4-Dichlorobenzene-d4	6.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3	Propylamine	59	C3H9N	000107-10-8	9
4	Diacetamide	101	C4H7NO2	000625-77-4	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

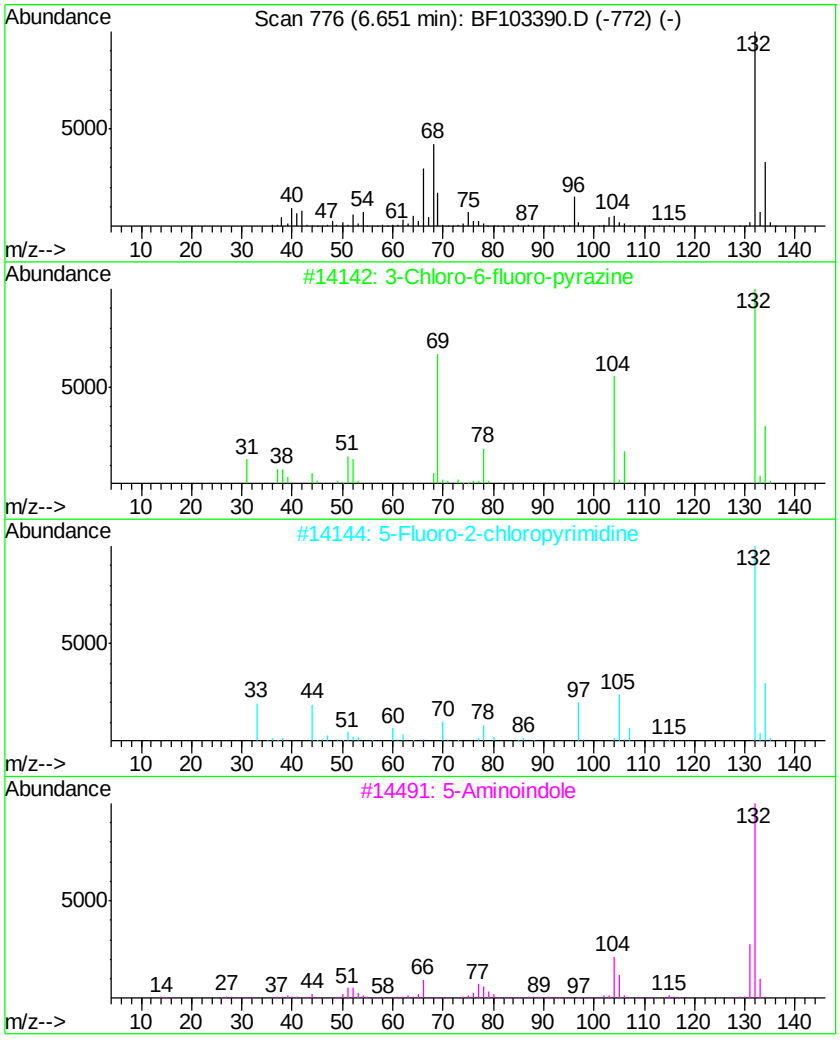
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

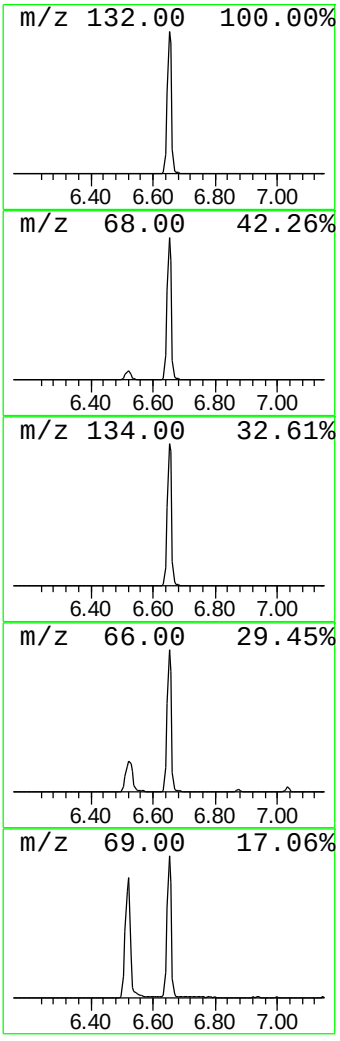
****

Peak Number 3 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	63.91 ng	3365420	1,4-Dichlorobenzene-d4	6.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3	5-Aminoindole	132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

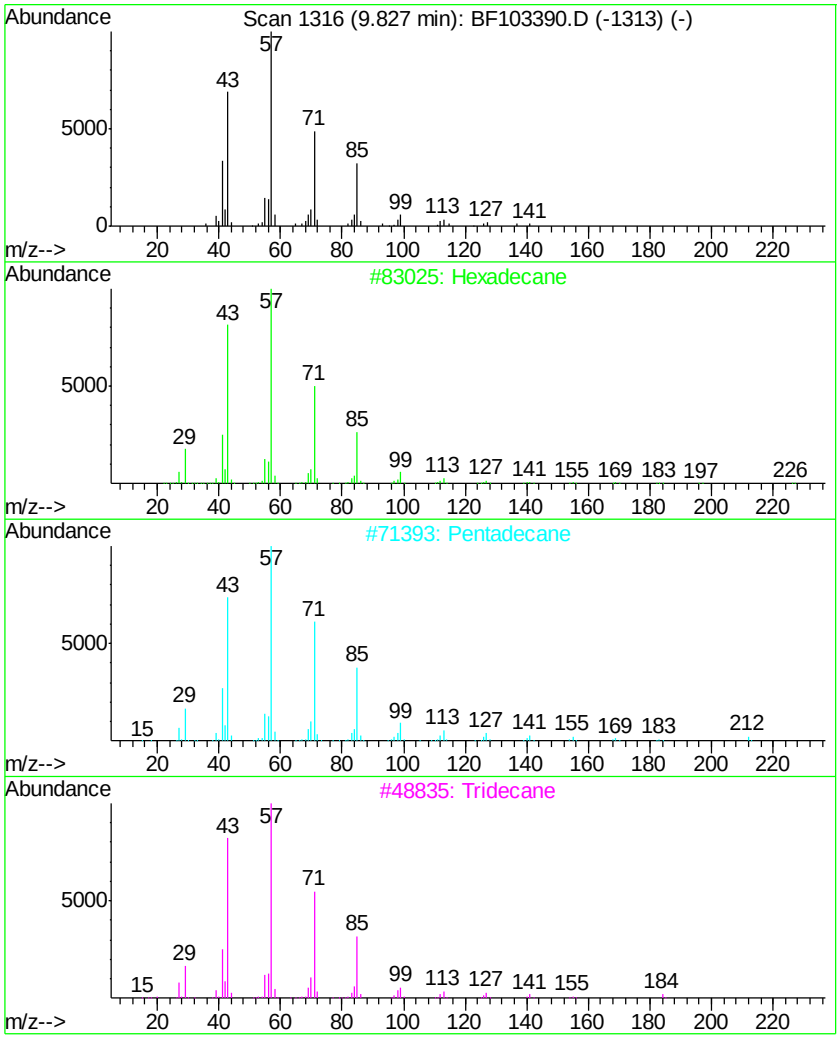
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

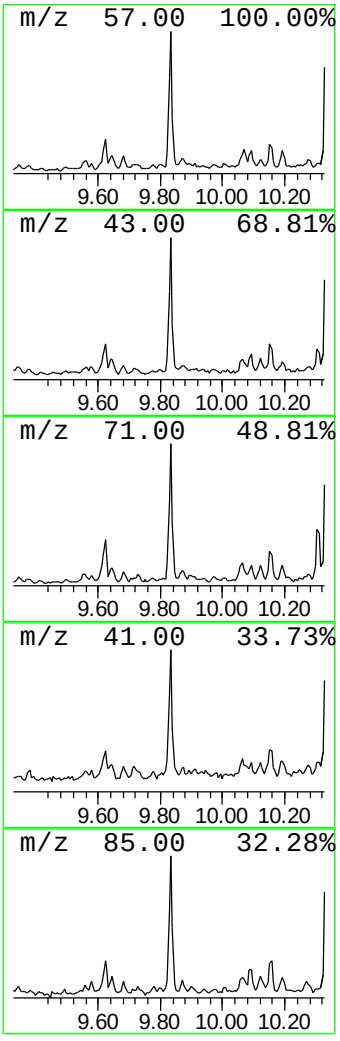
****

Peak Number 5 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	2.29 ng	126017	Acenaphthene-d10	9.92

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Pentadecane	212	C15H32	000629-62-9	90
3	Tridecane	184	C13H28	000629-50-5	90
4	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86
5	Tridecane, 6-methyl-	198	C14H30	013287-21-3	78





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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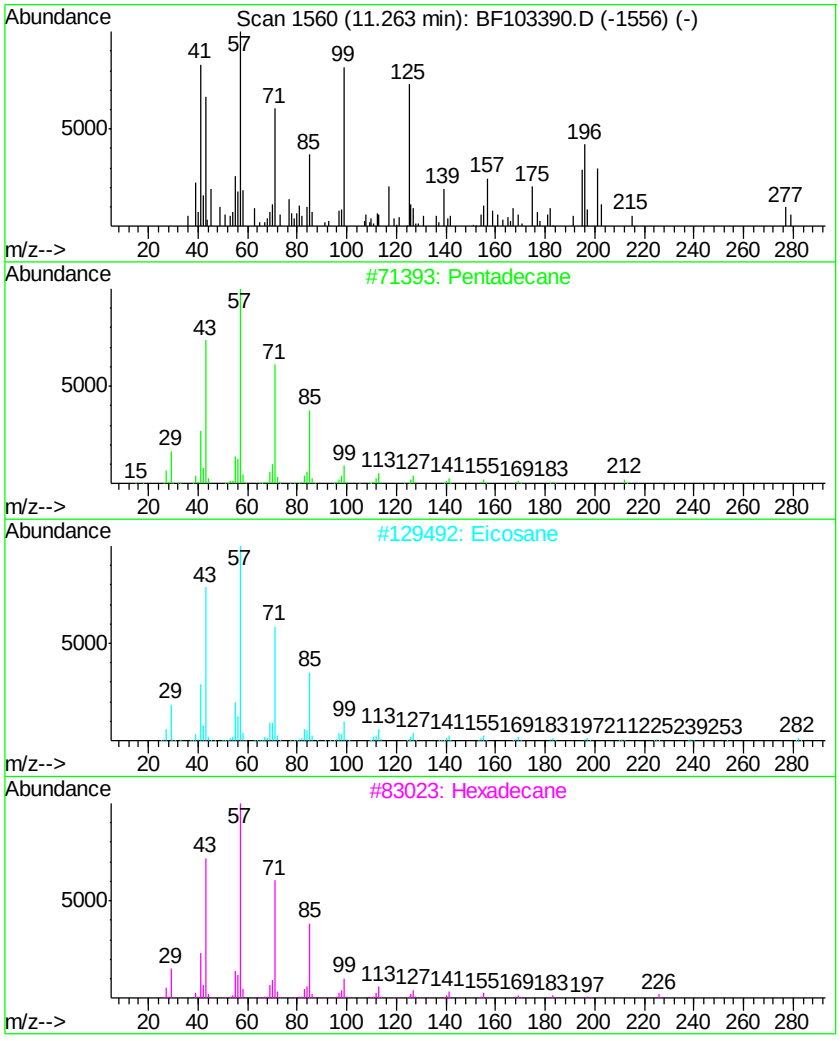
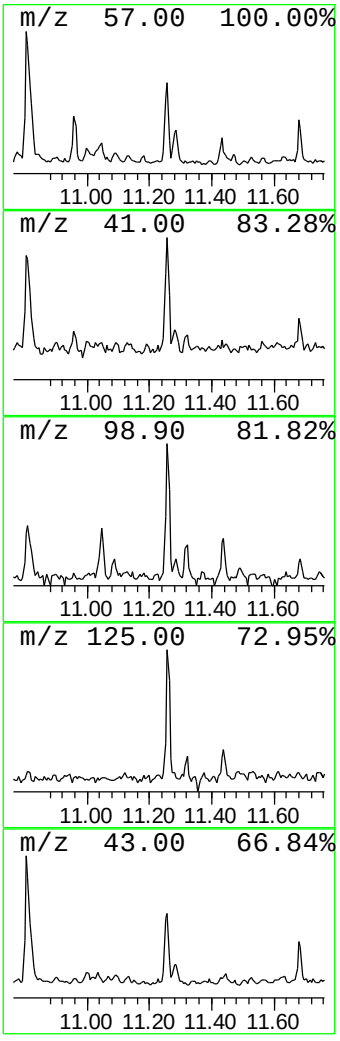
Peak Number 8 unknown11.26 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	2.96 ng	135710	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	49
2	Eicosane	282	C20H42	000112-95-8	46
3	Hexadecane	226	C16H34	000544-76-3	46
4	Heptadecane	240	C17H36	000629-78-7	46
5	Tetradecane	198	C14H30	000629-59-4	43





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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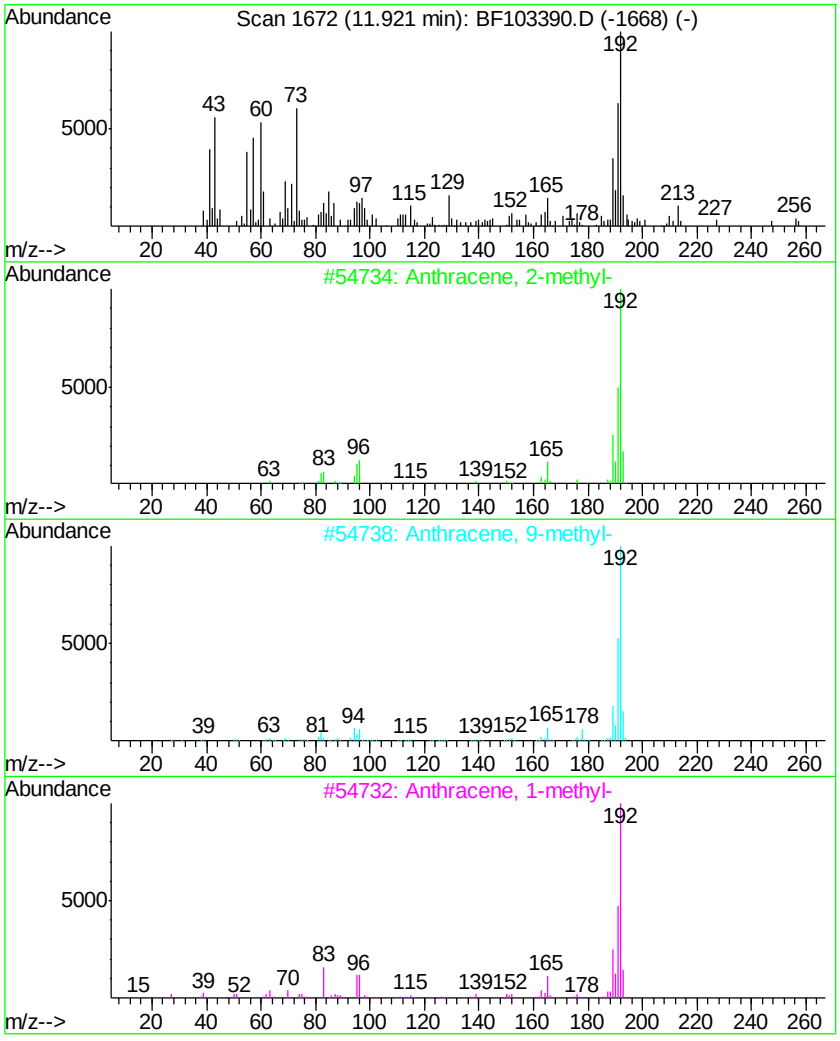
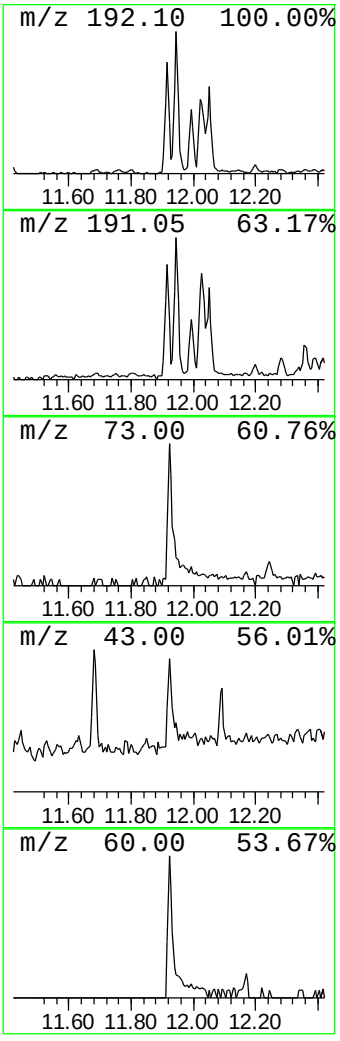
Peak Number 9 Anthracene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	3.32 ng	152264	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	78
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	78
4	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	60
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	60





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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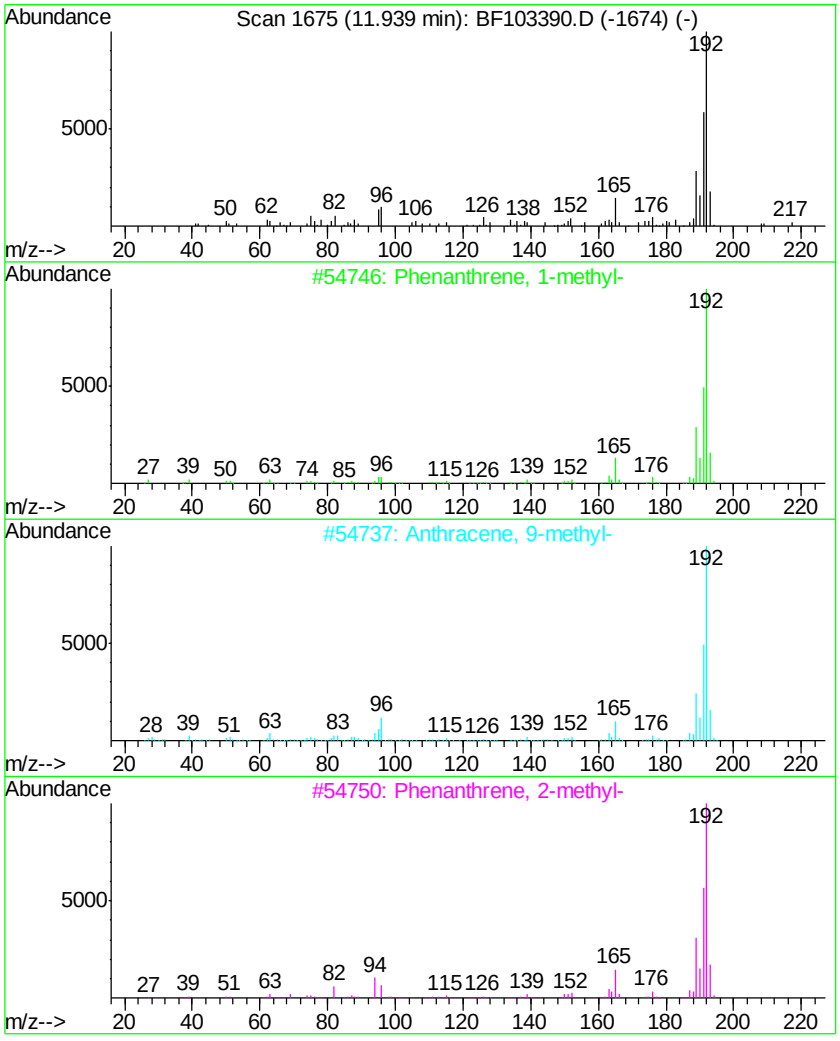
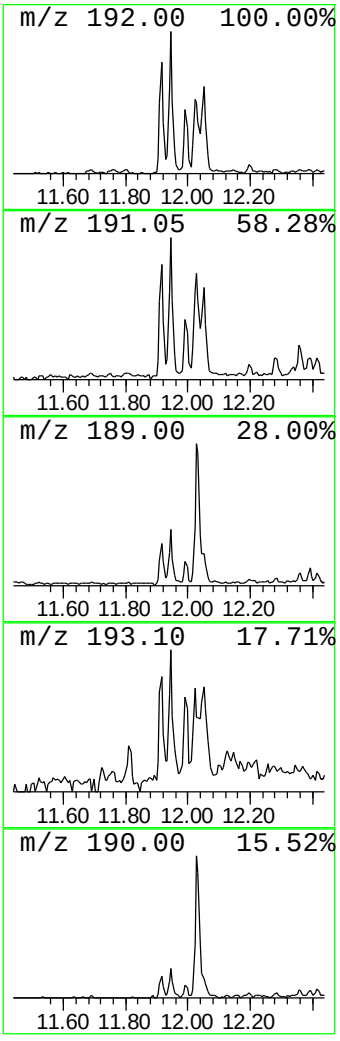
Peak Number 10 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	2.71 ng	124211	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91





Data Path : Z:\HPCHEM1\BNA F\Data\BF030218\
Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

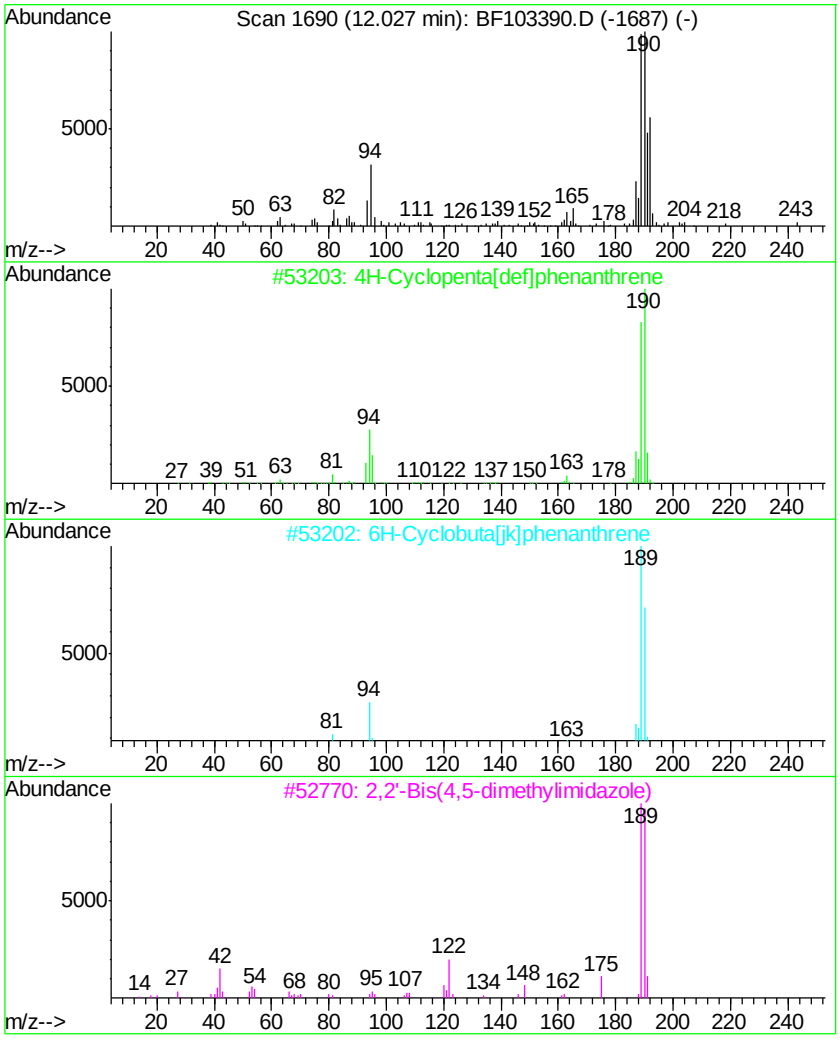
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

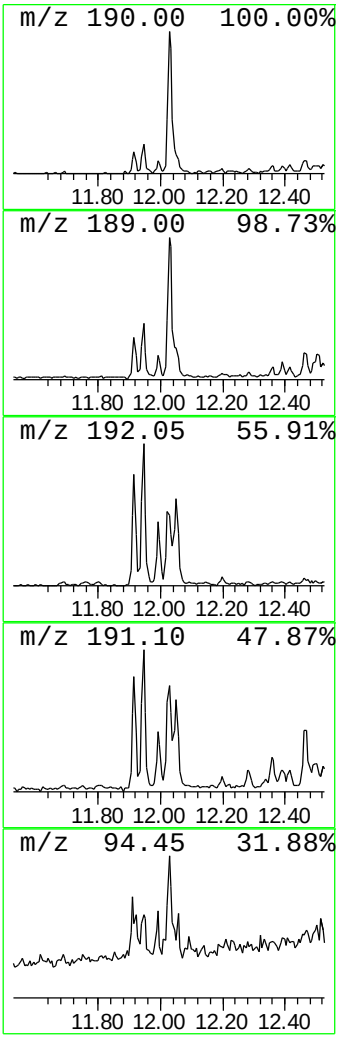
****

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	3.42 ng	156651	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	53
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
4	Methyl diselenide	190	C2H6Se2	007101-31-7	35
5	Megastigmatrienone	190	C13H18O	038818-55-2	18





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Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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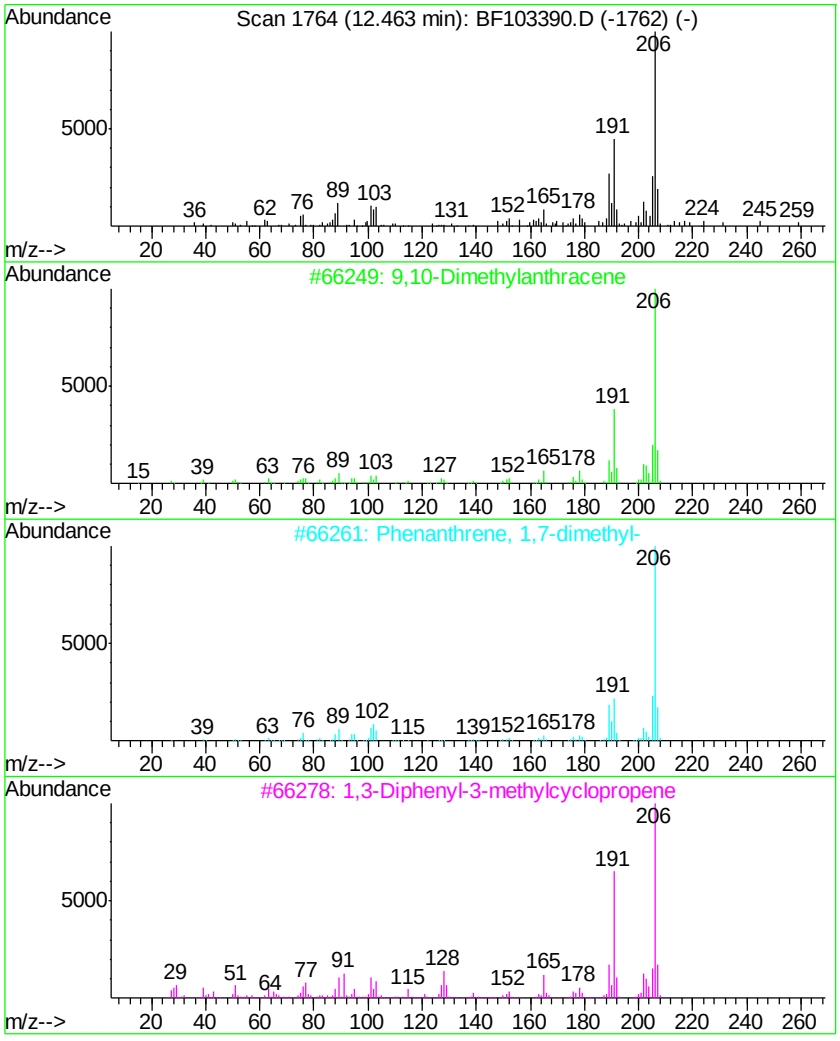
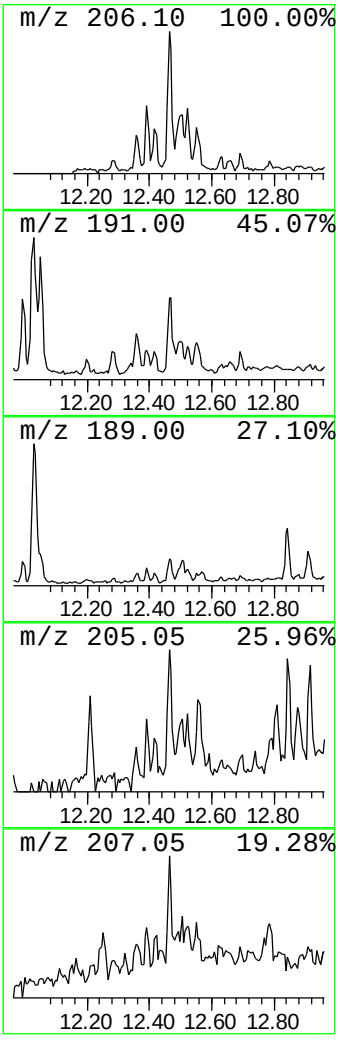
Peak Number 12 9,10-Dimethylanthracene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	2.33 ng	106867	Phenanthrene-d10	11.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylanthracene	206	C16H14	000781-43-1	93
2	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	92
4	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91
5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	90





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Misc :
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Instrument : BNA_F
ClientSampleId : SP-22

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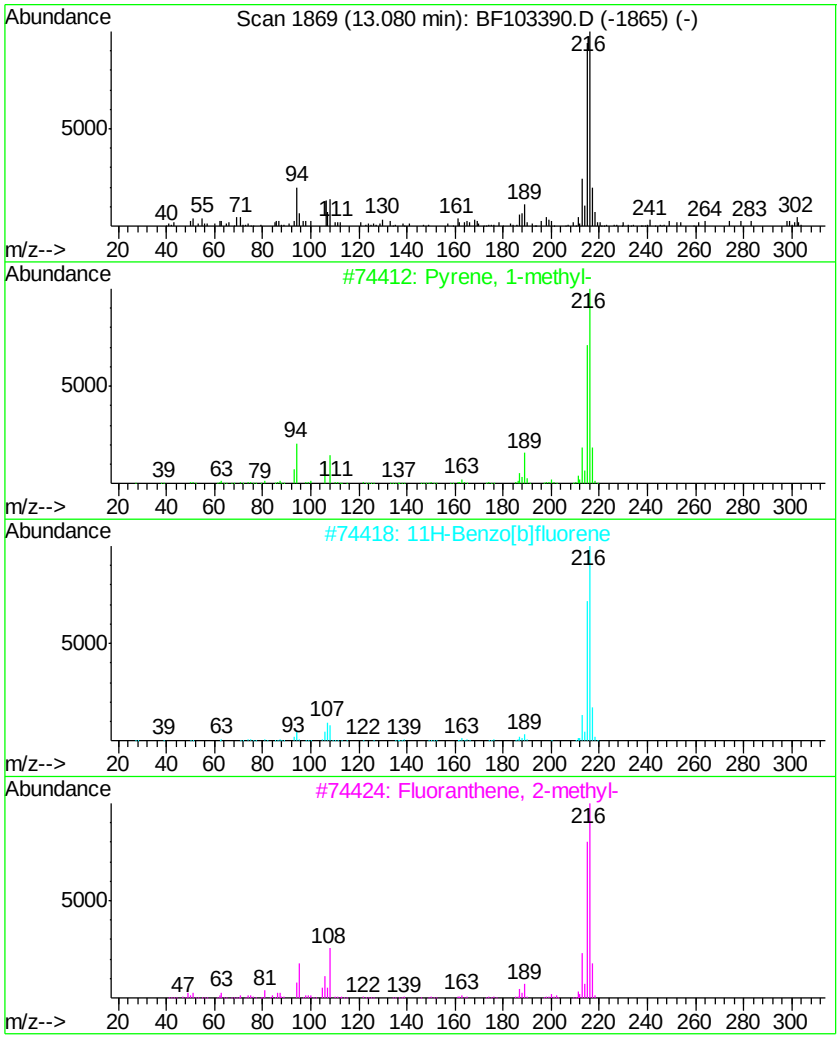
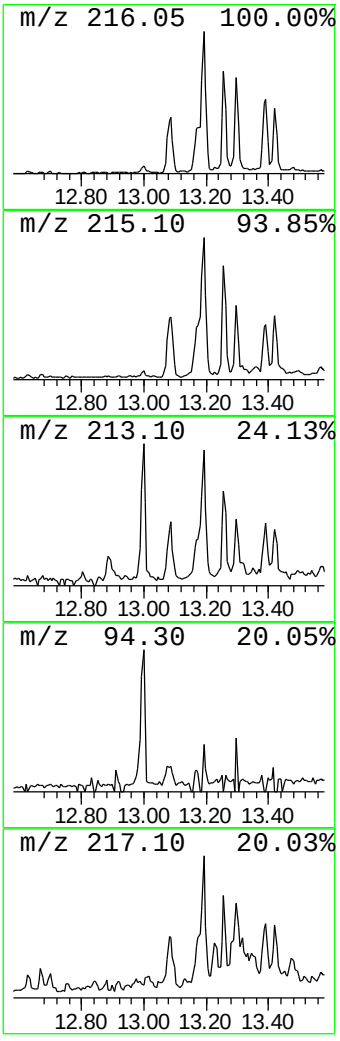
Peak Number 13 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	2.38 ng	128760	Chrysene-d12	14.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	64





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Data File : BF103390.D
Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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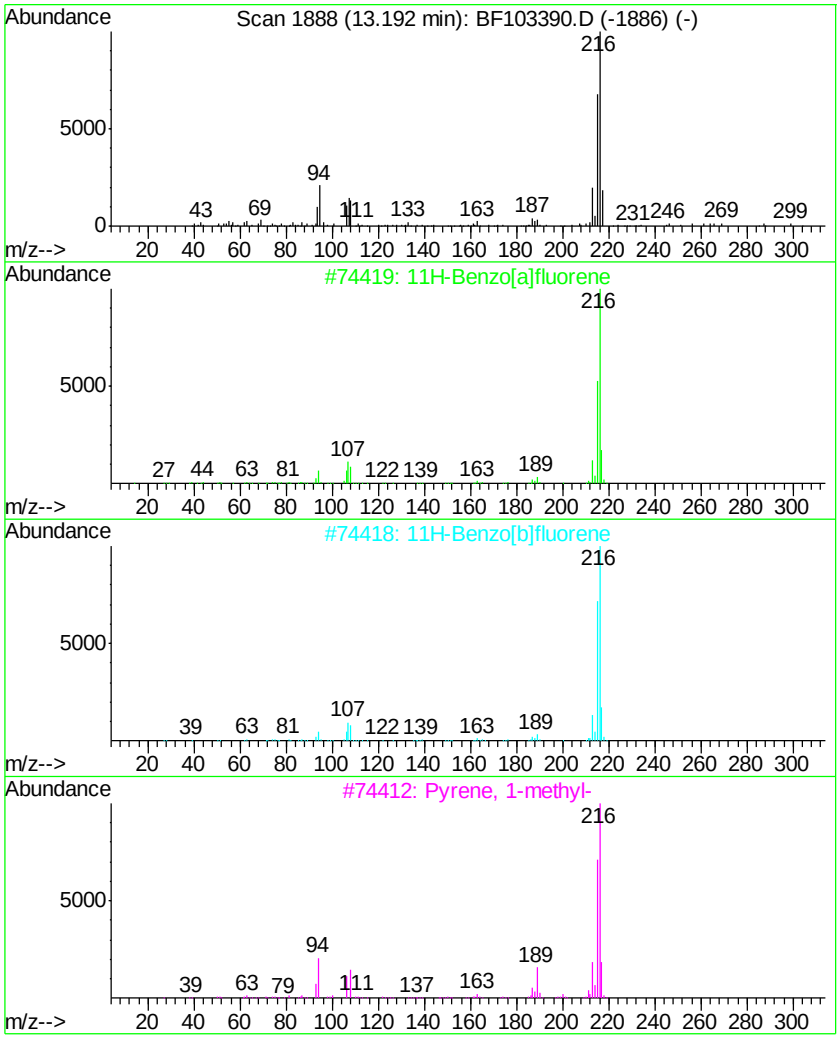
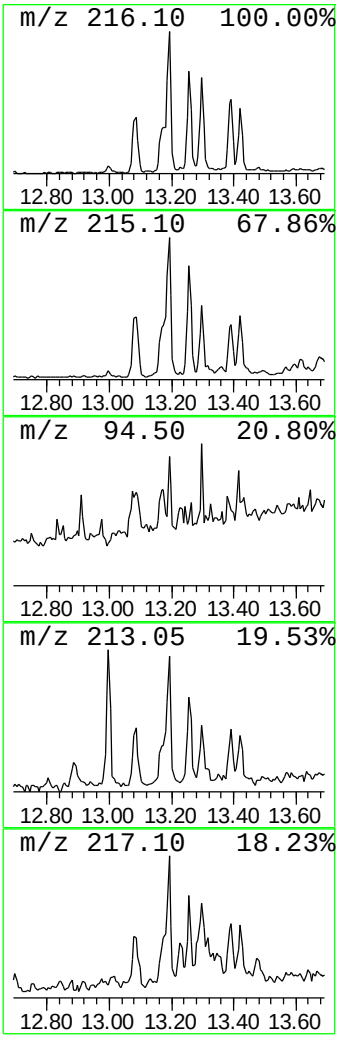
Peak Number 14 11H-Benzo[a]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	2.13 ng	115283	Chrysene-d12	14.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	80
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	70





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Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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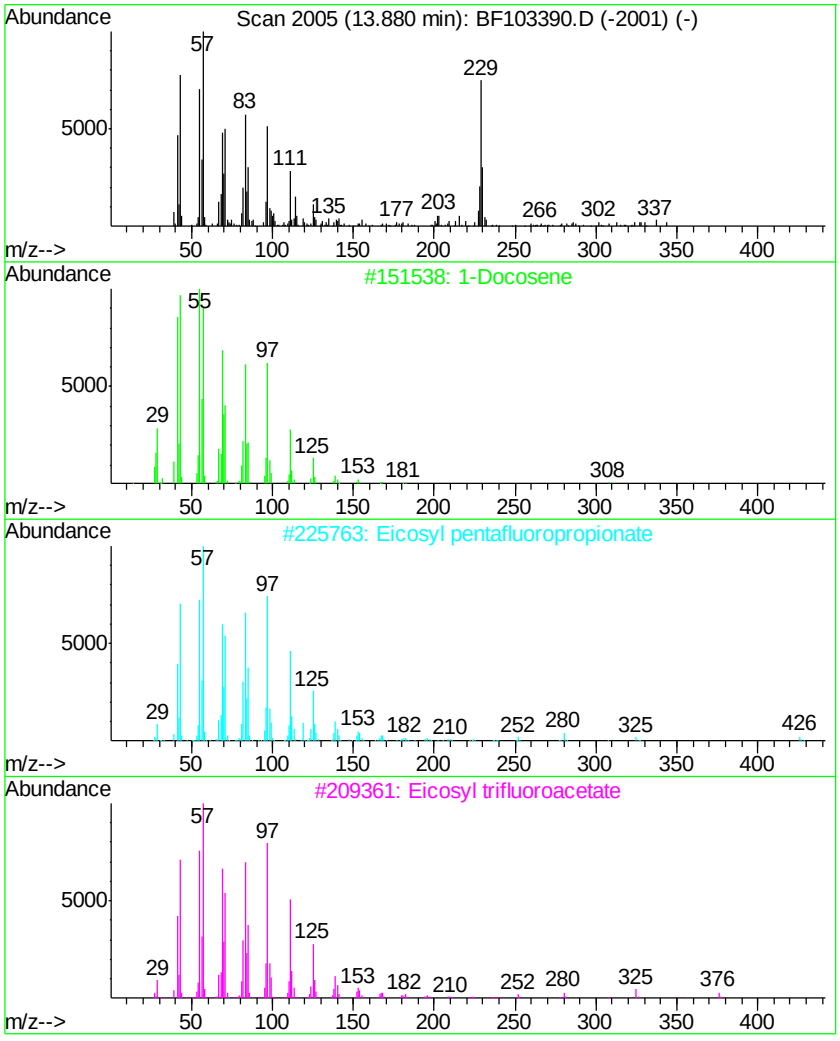
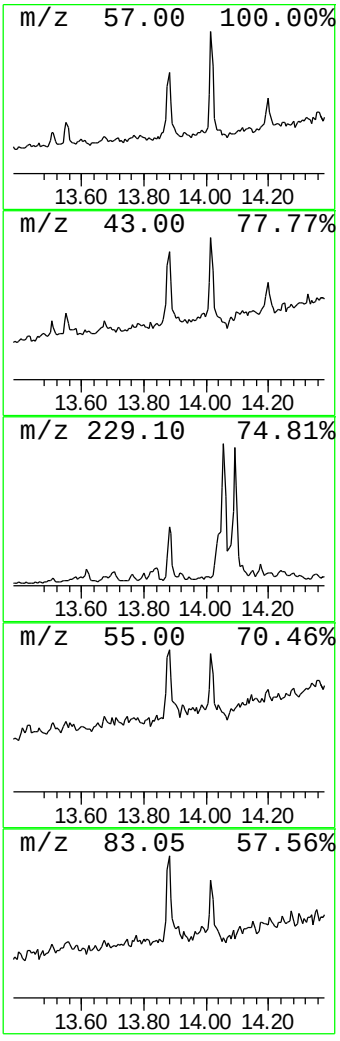
Peak Number 15 1-Docosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	4.42 ng	239216	Chrysene-d12	14.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	62
2	Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	55
3	Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	55
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	49
5	Eicosyl heptafluorobutyrate	494	C24H41F7O2	1000351-83-5	49





Tentatively Identified Compound (LSC) summary

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Acq On : 3 Mar 2018 2:19
Operator : SJ/JU
Sample : J1724-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-22

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butane, 2-methoxy...	2.12	5.0	ng	260602	1	6.88	1053230	20.0
2-Pentanone, 4-hy...	5.12	2.1	ng	110699	1	6.88	1053230	20.0
unknown6.65	6.65	63.9	ng	3365420	1	6.88	1053230	20.0
Hexadecane	9.83	2.3	ng	126017	3	9.92	1101910	20.0
unknown11.26	11.26	3.0	ng	135710	4	11.42	915975	20.0
Anthracene, 2-met...	11.92	3.3	ng	152264	4	11.42	915975	20.0
Phenanthrene, 1-m...	11.94	2.7	ng	124211	4	11.42	915975	20.0
4H-Cyclopenta[def...	12.03	3.4	ng	156651	4	11.42	915975	20.0
9,10-Dimethylanth...	12.46	2.3	ng	106867	4	11.42	915975	20.0
Pyrene, 1-methyl-	13.08	2.4	ng	128760	5	14.07	1081370	20.0
11H-Benzo[a]fluorene	13.19	2.1	ng	115283	5	14.07	1081370	20.0
1-Docosene	13.88	4.4	ng	239216	5	14.07	1081370	20.0