LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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	5.189	481	485	496	rBV	144764	166454	11.23%	1.037%
2	5.595	551	554	567	rBV	1141183	1185043	79.96%	7.380%
3	6.601	721	725	741	rBV	1240510	1224142	82.60%	7.624%
4	6.731	743	747	751	rBV	1283876	1252167	84.49%	7.798%
5	6.954	782	785	789	rBV	464414	361692	24.40%	2.253%
6	7.113	808	812	815	rBV	1168904	1024600	69.13%	6.381%
7	7.519	877	881	892	rBV	814392	794185	53.59%	4.946%
8	8.236	1000	1003	1022	rVB	473112	457251	30.85%	2.848%
9	9.313	1181	1186	1189	rBV	1655273	1482049	100.00%	9.230%
10	9.701	1249	1252	1262	rVB	103714	100590	6.79%	0.626%
11	9.995	1299	1302	1306	rBV	486151	474419	32.01%	2.955%
12	10.030	1306	1308	1314	rVB	46951	36904	2.49%	0.230%
13	10.401	1369	1371	1375	rVB	22861	16894	1.14%	0.105%
14	10.613	1400	1407	1413	rBV3	26779	42897	2.89%	0.267%
15	10.795	1434	1438	1449	rBV	949320	881299	59.46%	5.489%
16	10.877	1449	1452	1456	rVB2	24453	29765	2.01%	0.185%
17	11.495	1553	1557	1559	rBV	516424	482066	32.53%	3.002%
18	11.513	1559	1560	1564	rVV	124780	107667	7.26%	0.671%
19	11.571	1567	1570	1573	rVB	30566	26930	1.82%	0.168%
20	11.695	1585	1591	1596	rBV4	24995	37178	2.51%	0.232%
21	11.995	1639	1642	1644	rBV	33015	27582	1.86%	0.172%
22	12.024	1644	1647	1650	rVB2	37986	29432	1.99%	0.183%
23	12.071	1653	1655	1658	rVB	26260	20147	1.36%	0.125%
24	12.107	1658	1661	1664	rBV2	41626	52161	3.52%	0.325%
25	12.242	1679	1684	1686	rBV2	26196	21078	1.42%	0.131%
26	12.436	1712	1717	1721	rBV3	22339	24622	1.66%	0.153%
27	12.542	1732	1735	1738	rBV2	30834	32891	2.22%	0.205%
28	12.642	1747	1752	1756	rBV3	11358	19369	1.31%	0.121%
29	12.713	1760	1764	1768	rBV	172549	144841	9.77%	0.902%
30	12.924	1796	1800	1801	rBV2	119611	99076	6.69%	0.617%
31	12.942	1801	1803	1809	rVV	222631	263617	17.79%	1.642%
32	12.989	1809	1811	1815	rVB3	18070	17561	1.18%	0.109%
33	13.077	1821	1826	1829	rBV	1600575	1437606	97.00%	8.953%
34	13.165	1837	1841	1844	rBV2	25233	35030	2.36%	0.218%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.271	1852	1859	1862	rBV	52499	82233	5.55%	0.512%
36	13.336	1864	1870	1873	rBV	62159	66654	4.50%	0.415%
37	13.377	1873	1877	1879	rBV2	49251	58932	3.98%	0.367%
38	13.401	1879	1881	1884	rVB	32778	32696	2.21%	0.204%
39	13.471	1890	1893	1895	rBV2	30086	24142	1.63%	0.150%
40	13.501	1895	1898	1901	rVB	37326	32049	2.16%	0.200%
41	13.624	1915	1919	1921	rBV2	32329	32388	2.19%	0.202%
42	13.759	1939	1942	1944	rBV3	16465	20171	1.36%	0.126%
43	13.783	1944	1946	1950	rVB	20642	18661	1.26%	0.116%
44	13.895	1960	1965	1968	rBV4	38379	58236	3.93%	0.363%
45	13.924	1968	1970	1972	rVB	37461	30088	2.03%	0.187%
46	13.954	1972	1975	1981	rBV3	114321	136810	9.23%	0.852%
47	14.154	2002	2009	2011	rBV2	568237	705009	47.57%	4.391%
48	14.177	2011	2013	2020	rVB	190235	170183	11.48%	1.060%
49	14.259	2024	2027	2033	rVV	156495	172148	11.62%	1.072%
50	14.307	2033	2035	2041	rVB3	41678	49824	3.36%	0.310%
51	14.477	2061	2064	2067	rBV3	16931	19114	1.29%	0.119%
52	14.512	2067	2070	2072	rBV	44636	44290	2.99%	0.276%
53	14.548	2072	2076	2080	rVV	80725	115451	7.79%	0.719%
54	14.583	2080	2082	2087	rVV2	40578	65918	4.45%	0.411%
55	14.654	2091	2094	2096	rVV	52605	61946	4.18%	0.386%
56	14.683	2096	2099	2100	rVV2	53091	50726	3.42%	0.316%
57	14.707	2100	2103	2106	rVV	71214	82456	5.56%	0.514%
58	14.742	2106	2109	2117	rVB3	55516	101326	6.84%	0.631%
59	14.995	2149	2152	2159	rVB	57862	70484	4.76%	0.439%
60	15.071	2163	2165	2169	rVV4	21084	21497	1.45%	0.134%
61	15.248	2191	2195	2198	rBV	121776	197397	13.32%	1.229%
62	15.359	2211	2214	2218	rVB	49179	57816	3.90%	0.360%
63	15.571	2246	2250	2255	rVB	116931	141181	9.53%	0.879%
64	15.636	2257	2261	2266	rBV	190623	248519	16.77%	1.548%
65	15.706	2268	2273	2277	rBV	259075	340786	22.99%	2.122%
66	17.289	2537	2542	2550	rVB2	65855	136014	9.18%	0.847%
67	17.471	2570	2573	2577	rBV2	22422	41966	2.83%	0.261%
68	17.777	2619	2625	2635	rVB2	63304	160319	10.82%	0.998%

Sum of corrected areas: 16056635

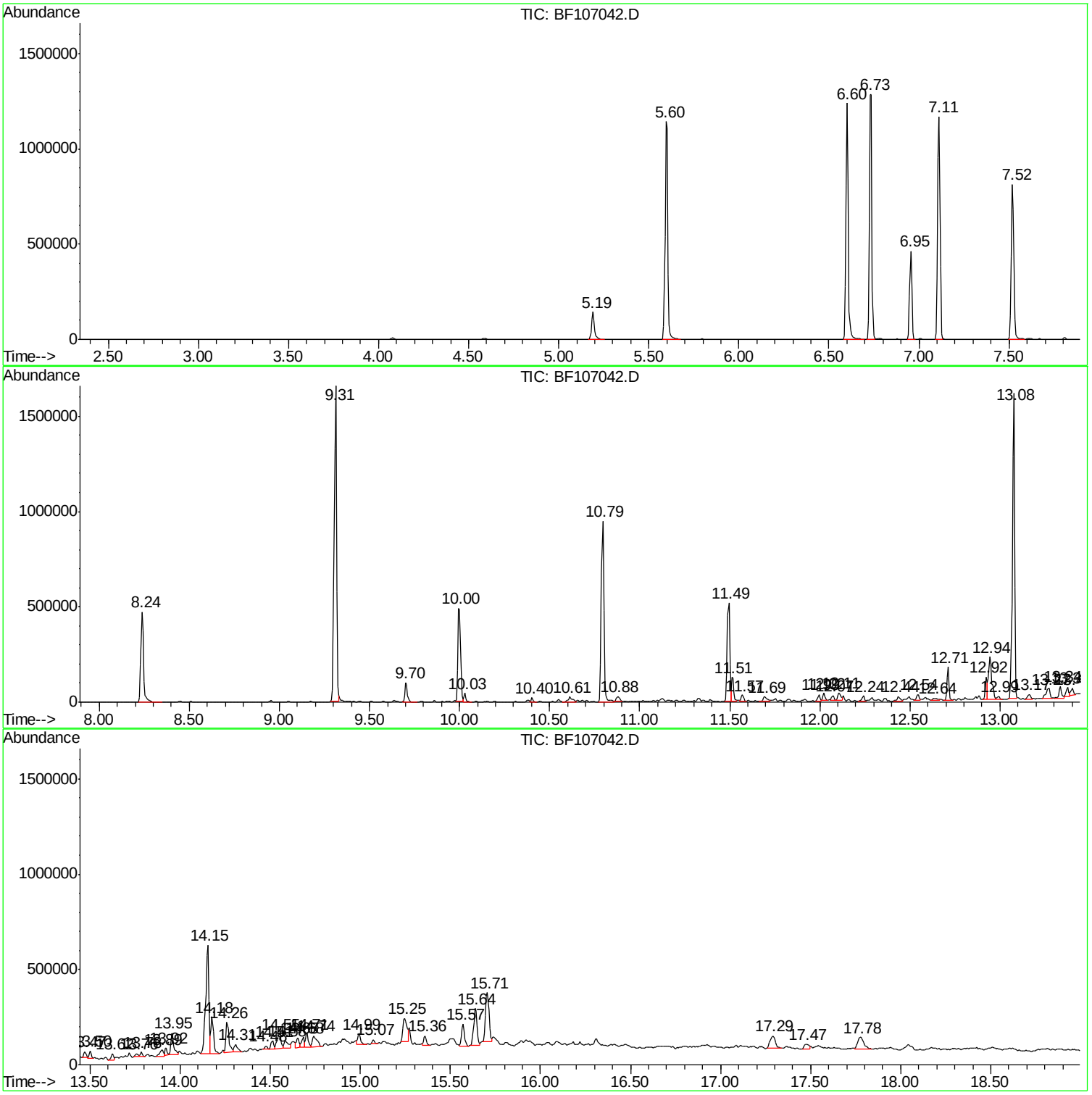
LSC Report - Integrated Chromatogram

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

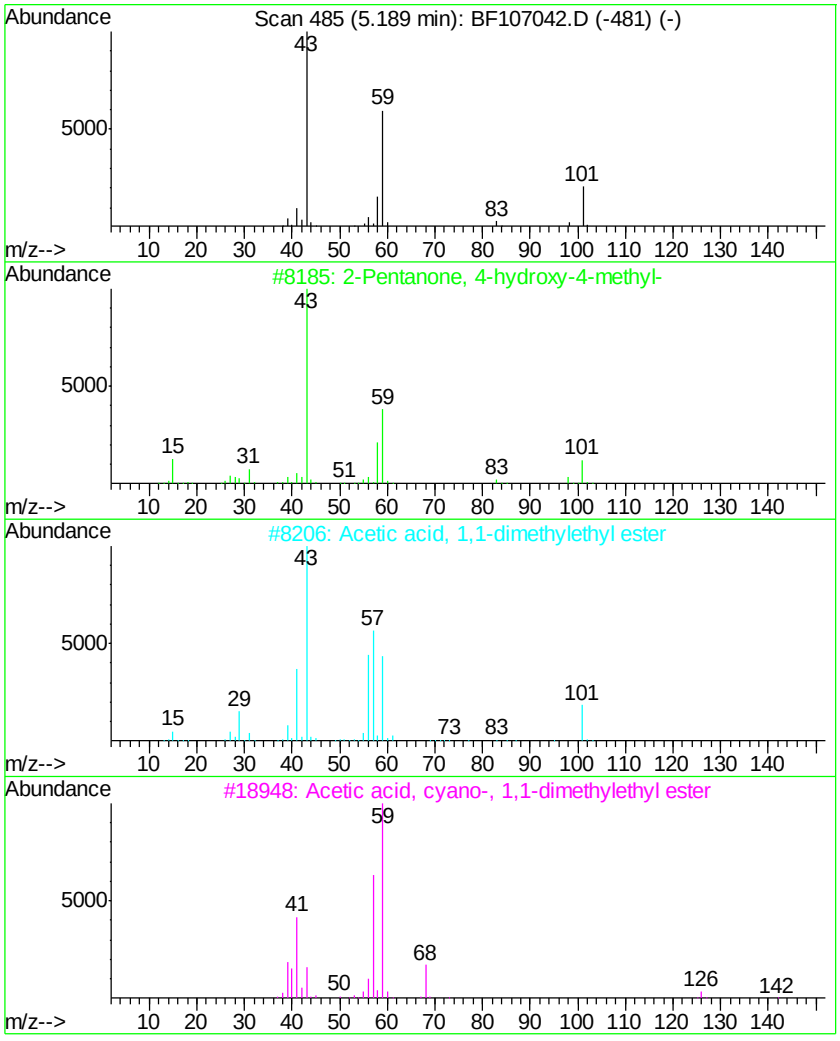
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

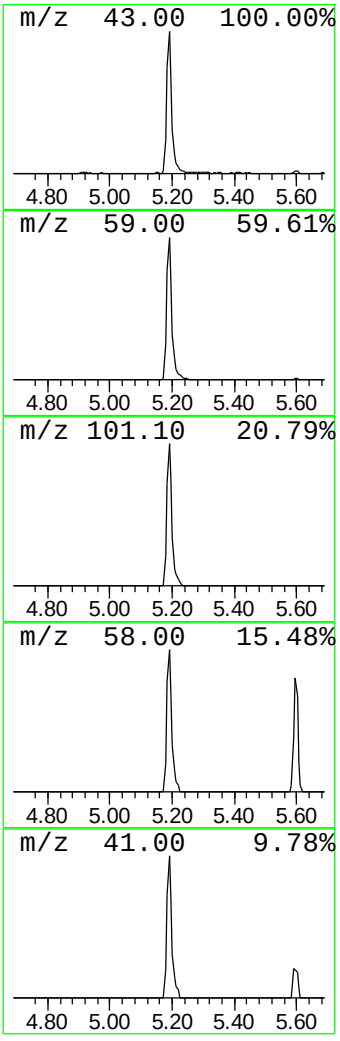
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	9.20 ng	166454	1,4-Dichlorobenzene-d4	6.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

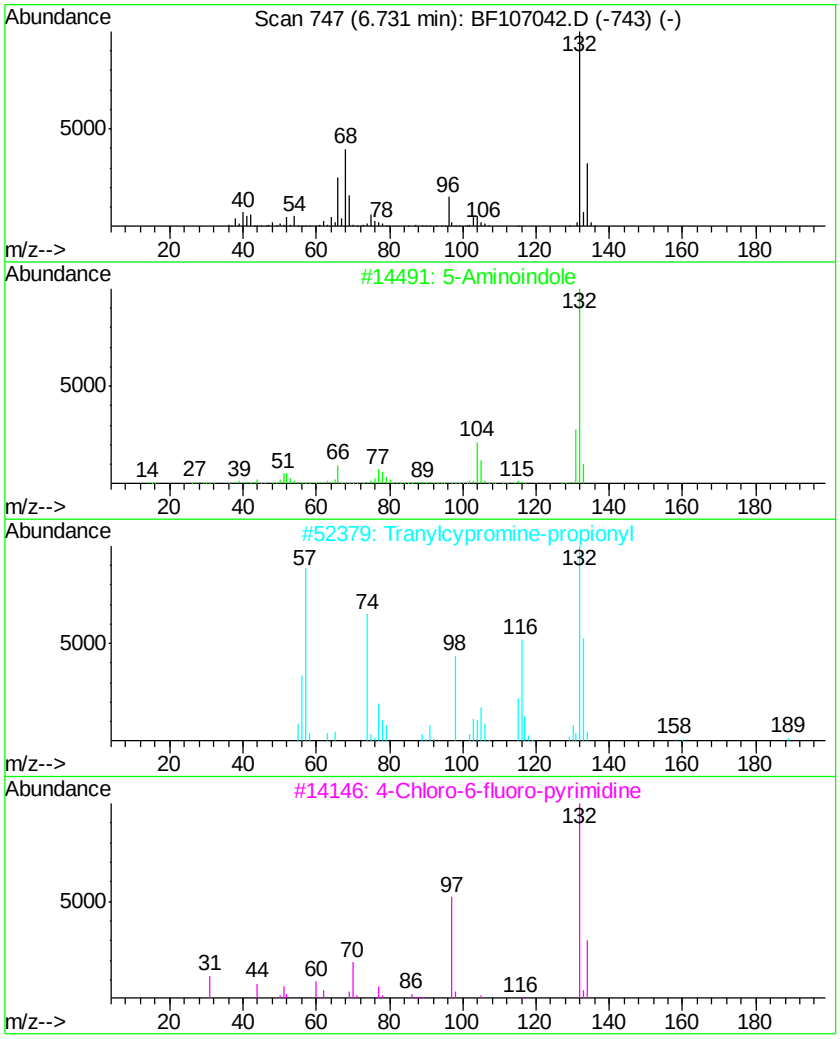
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

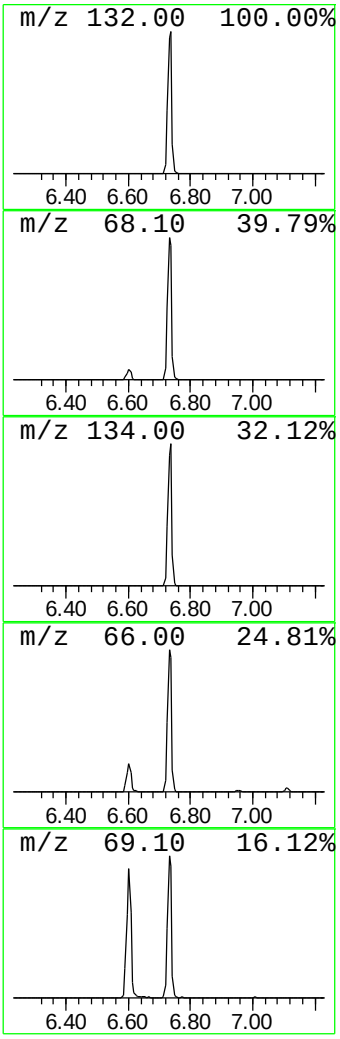
****

Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	69.24 ng	1252170	1,4-Dichlorobenzene-d4	6.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Aminoindole	132	C8H8N2	005192-03-0	12
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4	Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

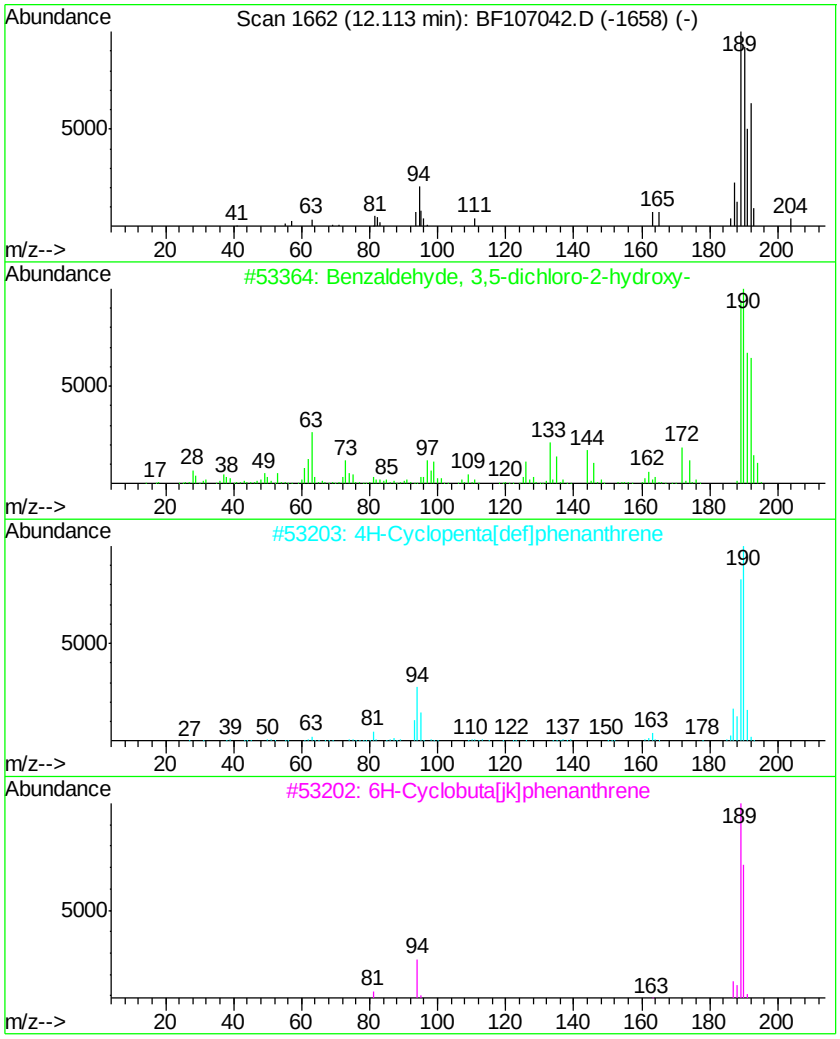
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

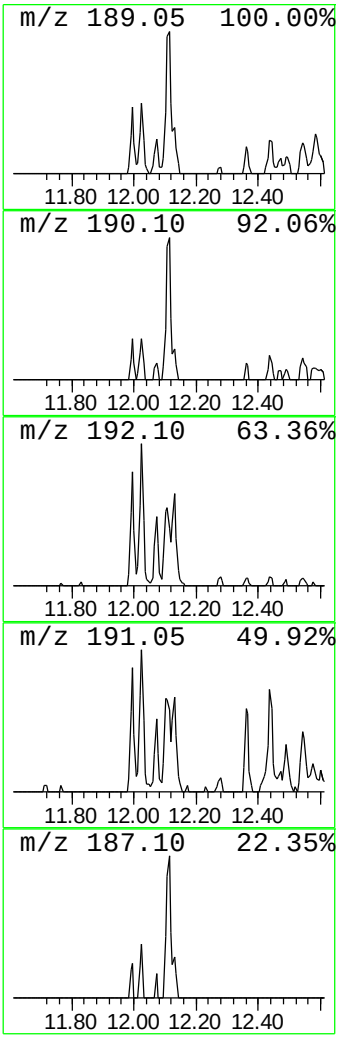
****

Peak Number 4 Benzaldehyde, 3,5-dichloro-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	2.16 ng	52161	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	58
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	41
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	38





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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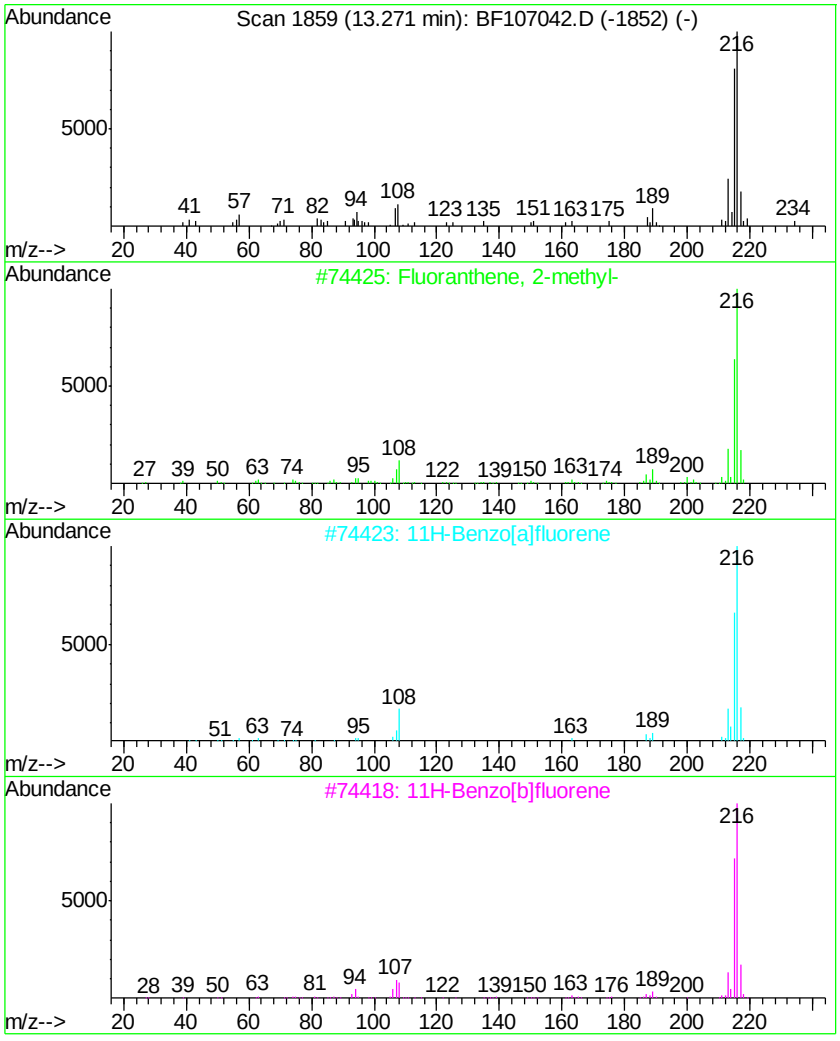
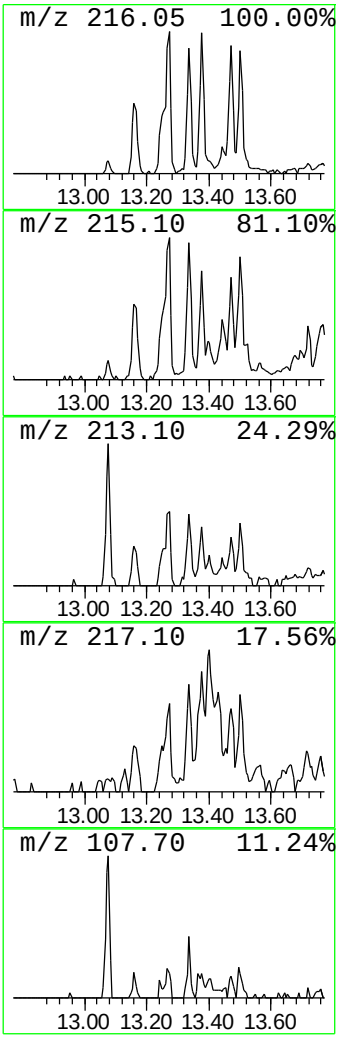
Peak Number 5 Fluoranthene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	2.33 ng	82233	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
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Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

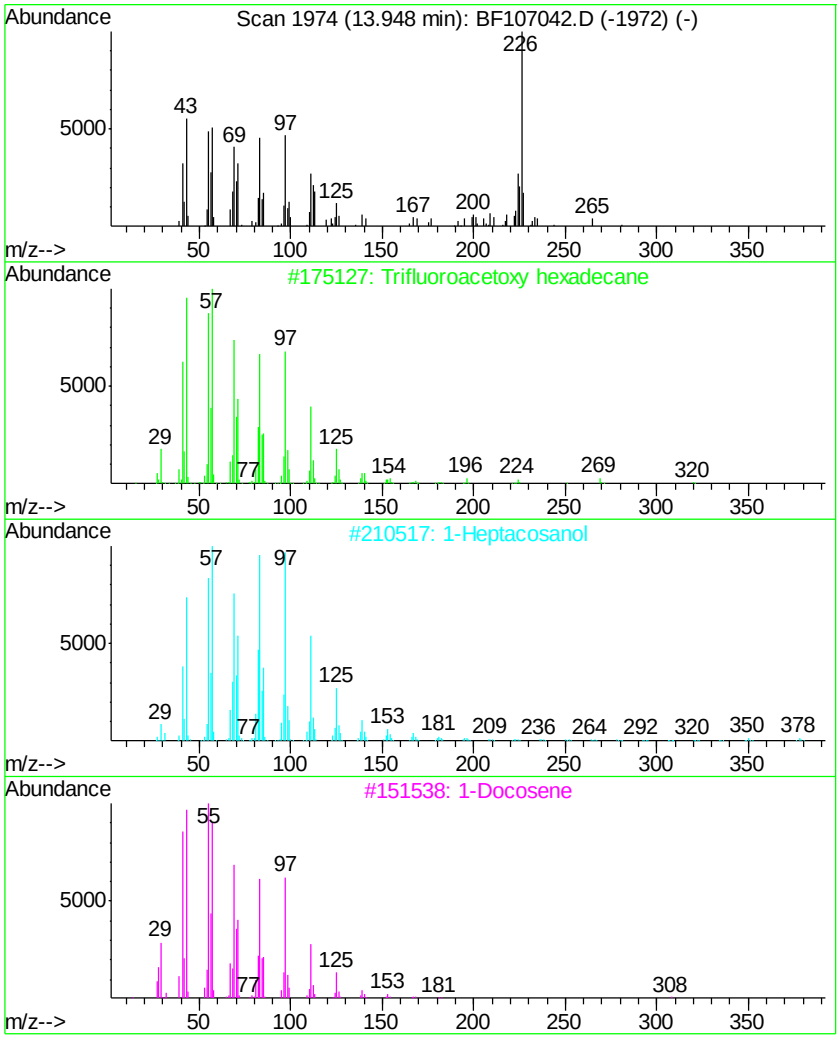
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

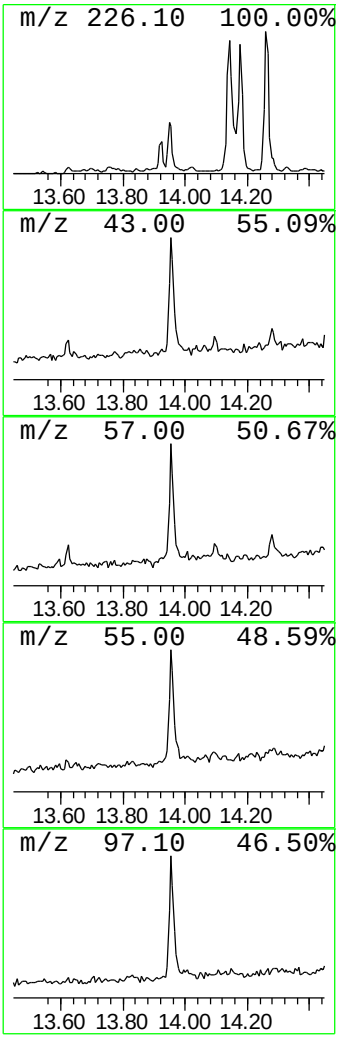
****

Peak Number 6 Trifluoroacetoxy hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	3.88 ng	136810	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	76
2	1-Heptacosanol	396	C27H56O	002004-39-9	76
3	1-Docosene	308	C22H44	001599-67-3	60
4	Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	60
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
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Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

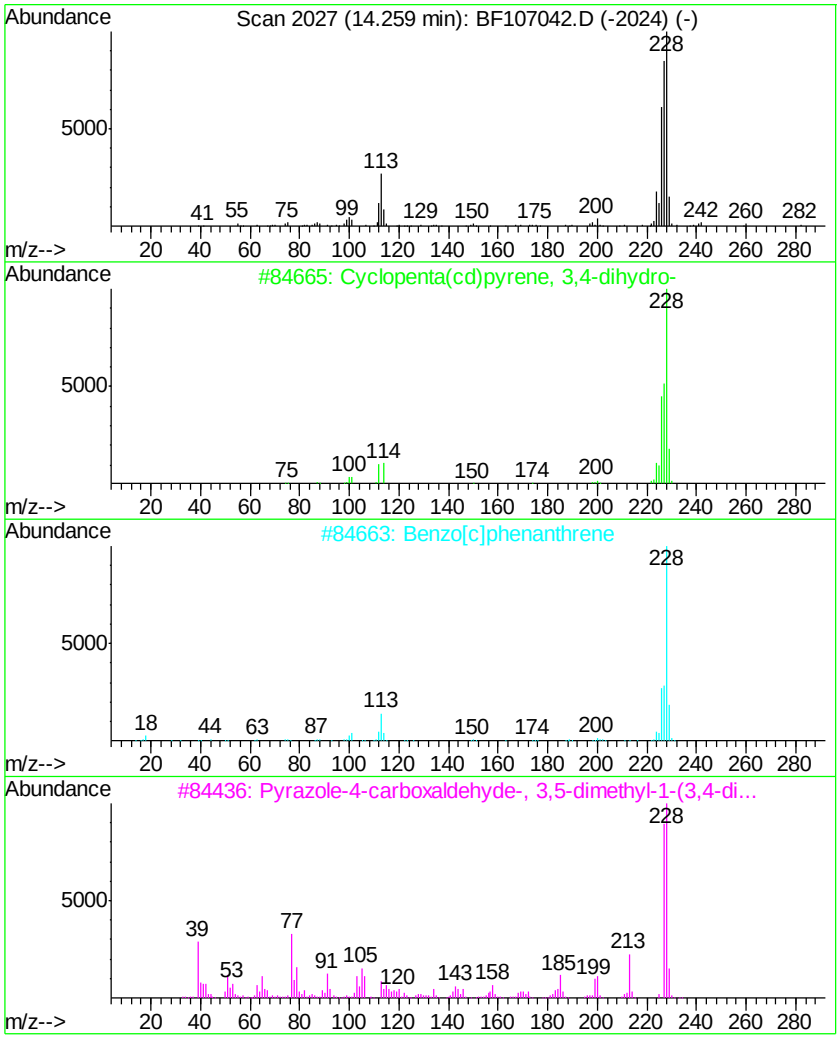
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

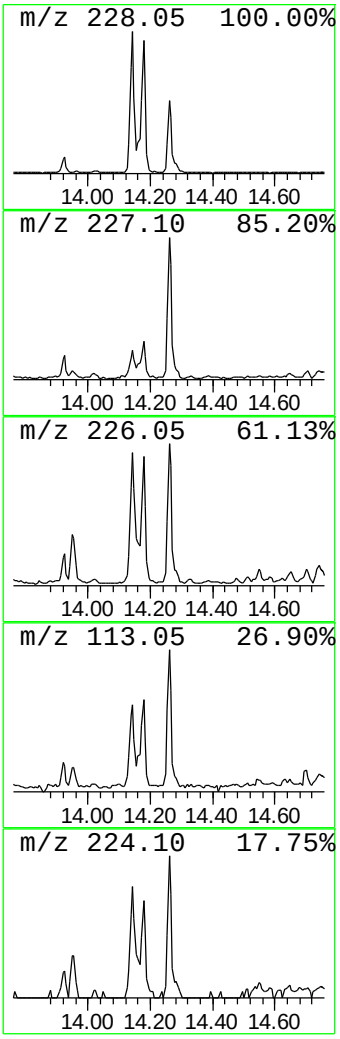
****

Peak Number 7 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	4.88 ng	172148	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	87
2	Benzo[c]phenanthrene	228	C18H12	000195-19-7	53
3	Pyrazole-4-carboxaldehyde-, 3,5-...	228	C14H16N2O	1000272-54-8	50
4	Isothiazol-5-amine, 4-bromo-3-ch...	226	C4H4BrClN2S	1000272-59-9	49
5	Diphenylvinylphosphine oxide	228	C14H13OP	002096-78-8	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

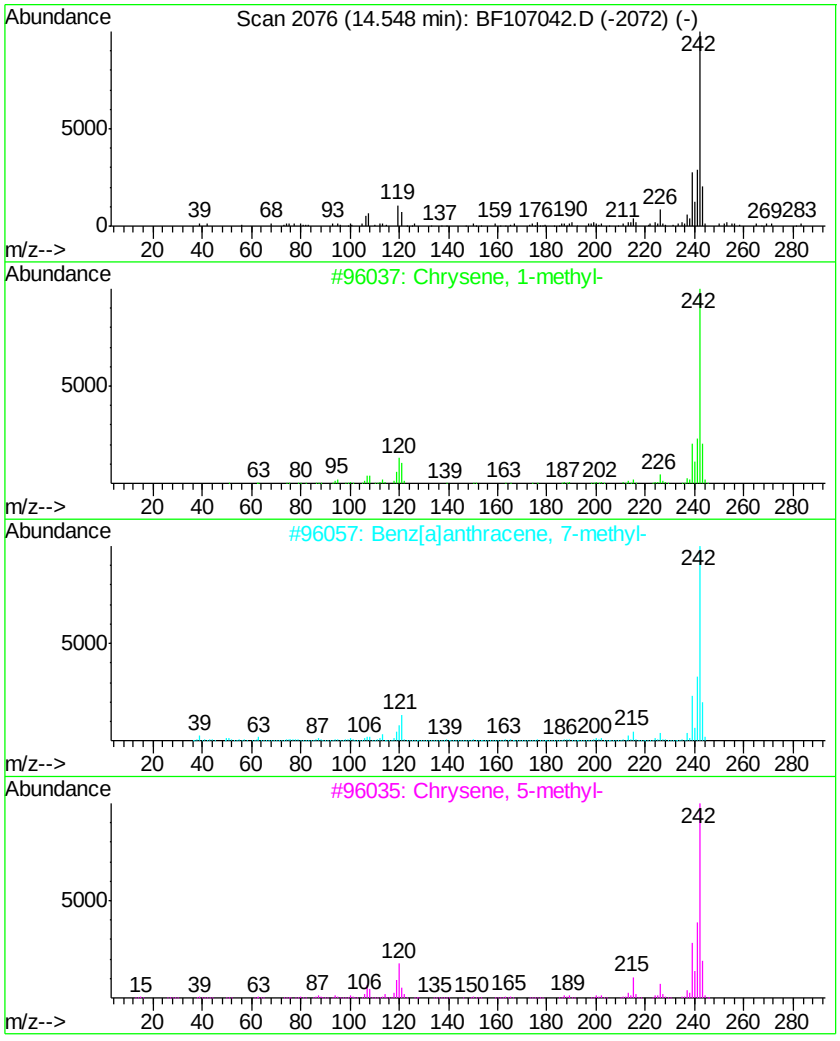
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

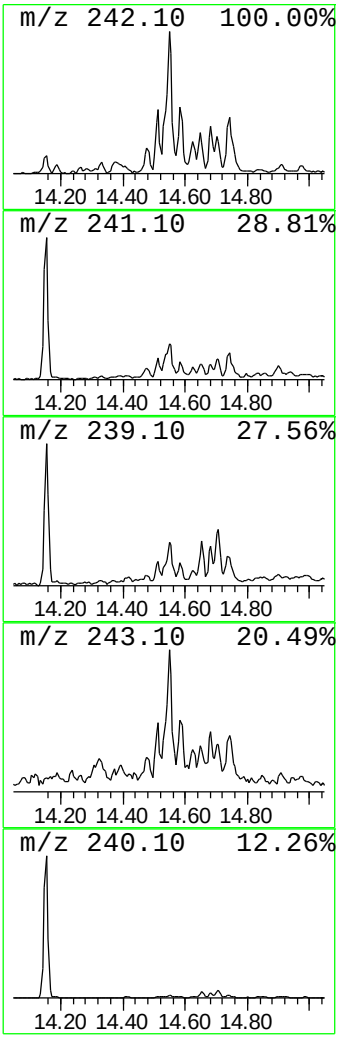
****

Peak Number 8 Chrysene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	3.28 ng	115451	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chrysene, 1-methyl-	242	C19H14	003351-28-8	91
2	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	90
3	Chrysene, 5-methyl-	242	C19H14	003697-24-3	90
4	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	86
5	Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	76





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Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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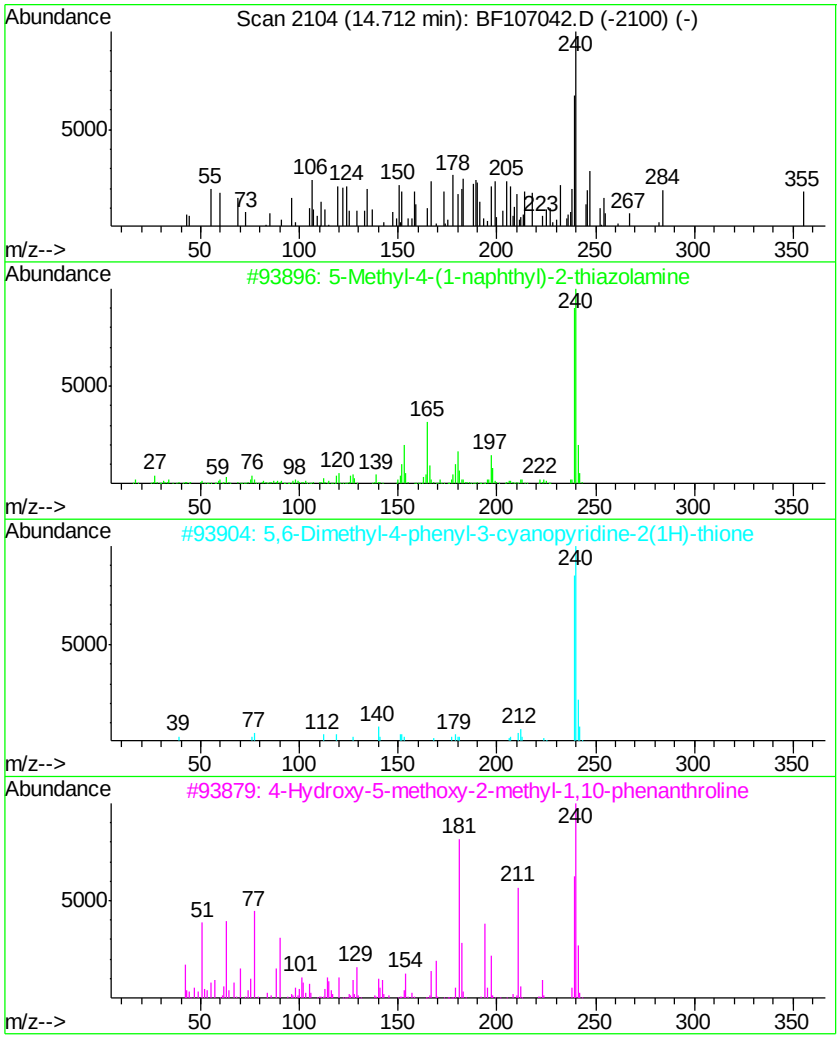
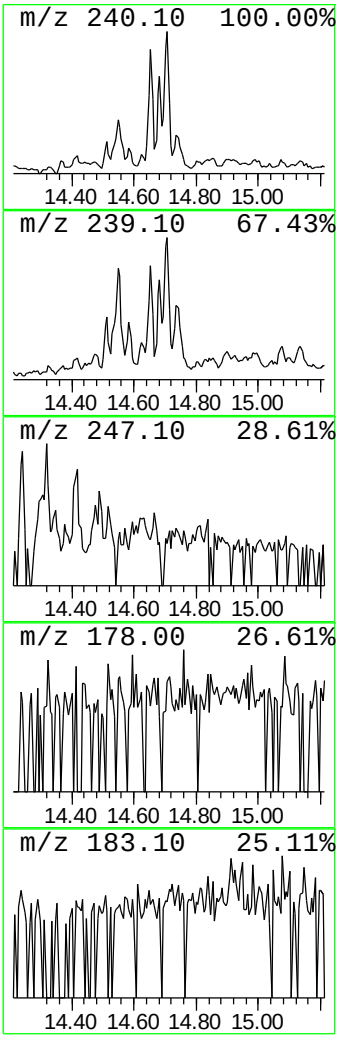
Peak Number 9 unknown14.71 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	2.34 ng	82456	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methyl-4-(1-naphthyl)-2-thiazo...	240	C14H12N2S	107411-05-2	43
2	5,6-Dimethyl-4-phenyl-3-cyanopyr...	240	C14H12N2S	094639-18-6	43
3	4-Hydroxy-5-methoxy-2-methyl-1,1...	240	C14H12N2O2	093533-14-3	38
4	9H-Cyclopenta[a]pyrene	240	C19H12	050861-05-7	38
5	2-Propen-1-one, 1-(2-hydroxyphen...	240	C15H12O3	013323-66-5	37





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Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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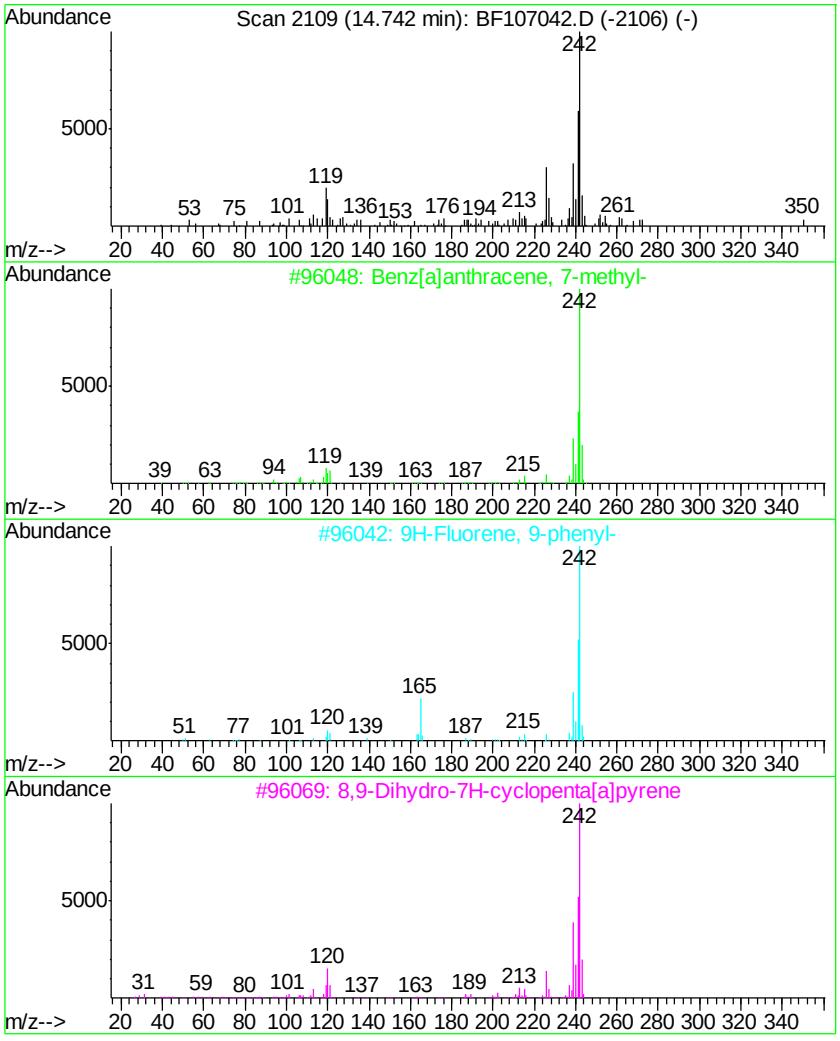
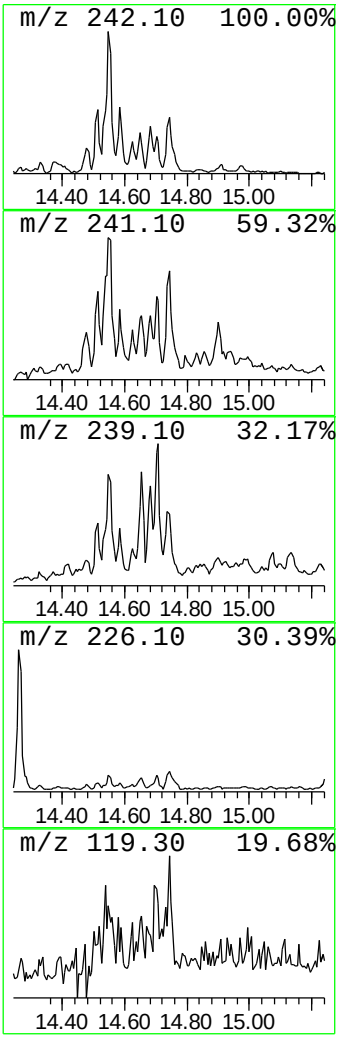
Peak Number 10 Benz[a]anthracene, 7-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.74	2.87 ng	101326	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	81
2	9H-Fluorene, 9-phenyl-	242	C19H14	000789-24-2	64
3	8,9-Dihydro-7H-cyclopenta[a]pyrene	242	C19H14	082979-72-4	53
4	Chrysene, 4-methyl-	242	C19H14	003351-30-2	49
5	Triphenylene, 1-methyl-	242	C19H14	002871-91-2	49





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Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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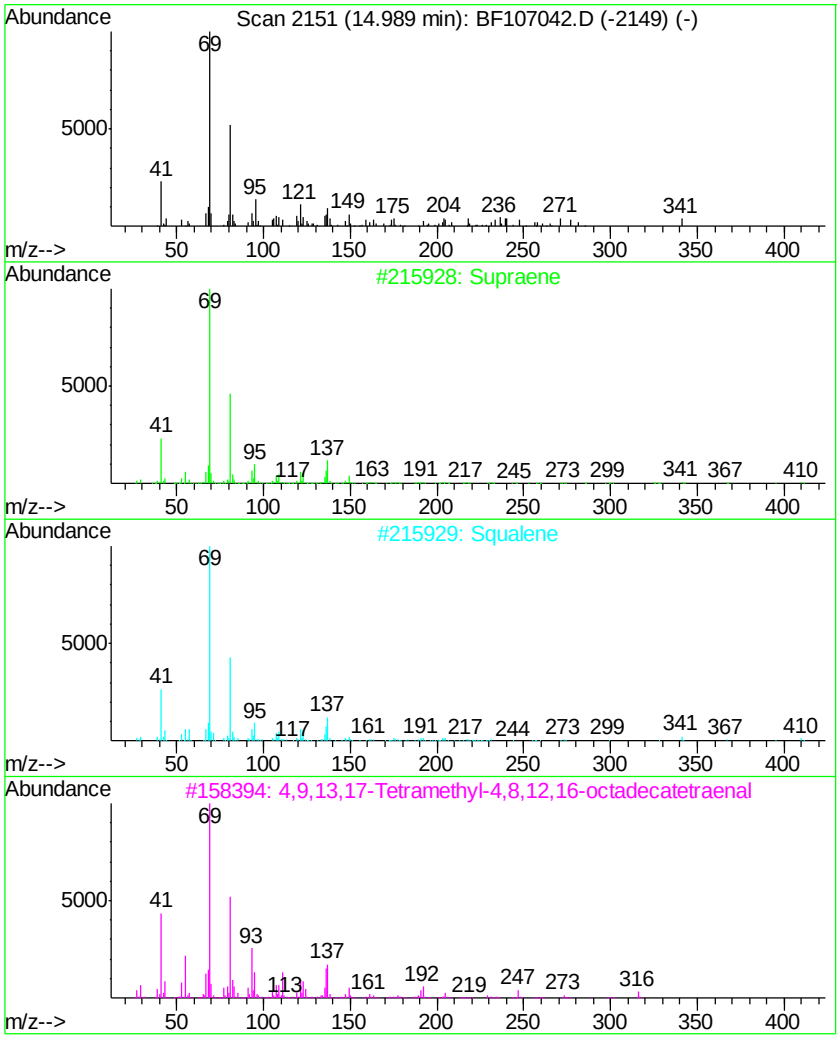
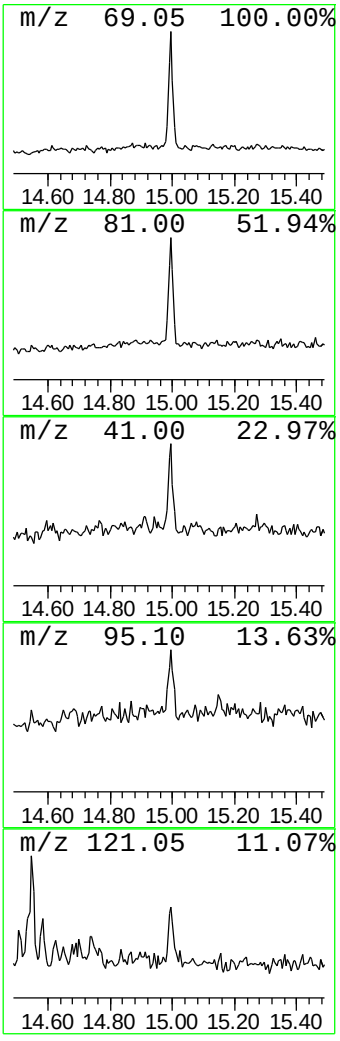
Peak Number 11 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	4.14 ng	70484	Perylene-d12	15.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Supraene	410	C30H50	007683-64-9	74
2	Squalene	410	C30H50	000111-02-4	72
3	4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	64
4	4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64
5	5,9,13-Pentadecatrien-2-one, 6,1...	262	C18H30O	000762-29-8	59





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
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Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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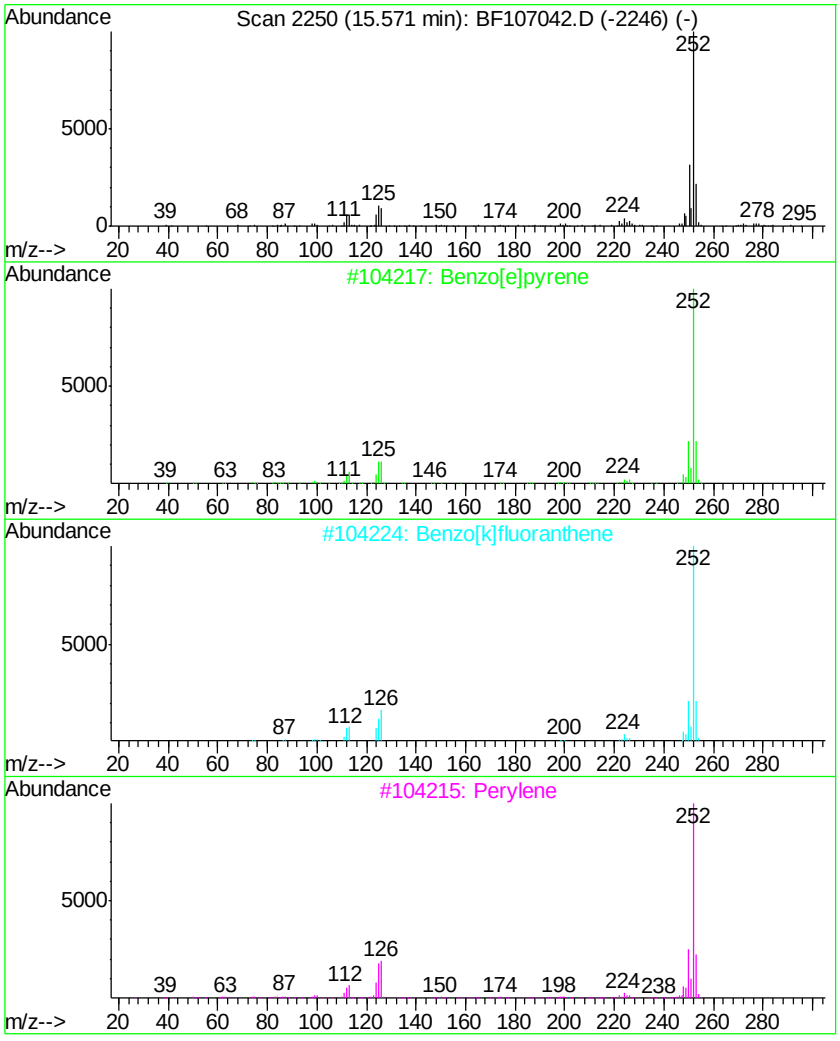
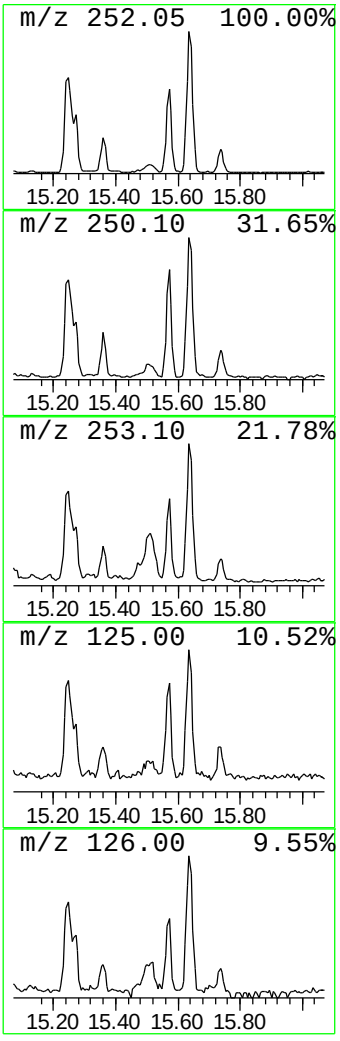
Peak Number 13 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	8.29 ng	141181	Perylene-d12	15.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3	Perylene	252	C20H12	000198-55-0	94
4	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
5	Benzo[a]pyrene	252	C20H12	000050-32-8	81





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Acq On : 2 Jul 2018 14:13
Operator : JU/SJ
Sample : J3774-06
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-6(2-4)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.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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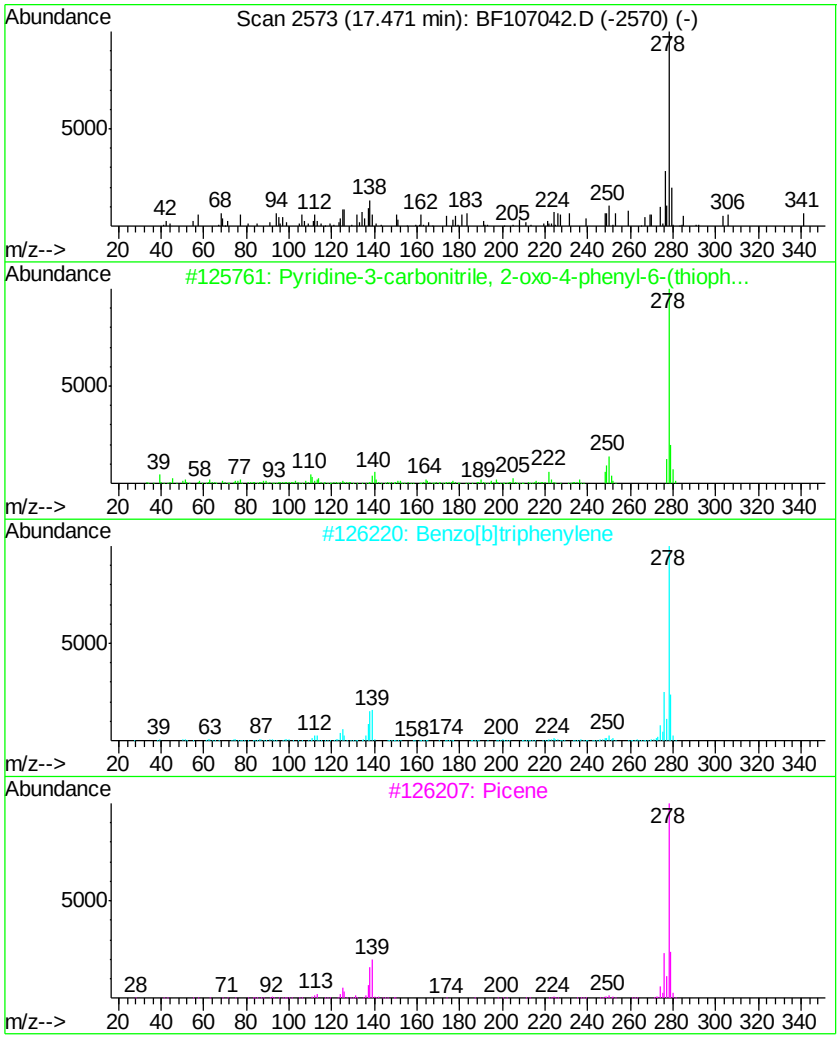
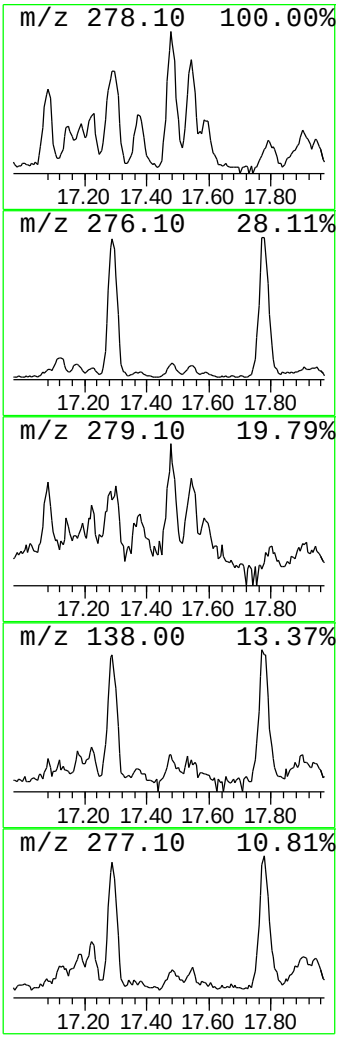
Peak Number 14 Pyridine-3-carbonitrile, 2-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	2.46 ng	41966	Perylene-d12	15.71

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine-3-carbonitrile, 2-oxo-4...	278	C16H10N2OS	1000304-72-3	70
2	Benzo[b]triphenylene	278	C22H14	000215-58-7	64
3	Picene	278	C22H14	000213-46-7	64
4	Estra-1,3,5,7,9,15-hexaen-17-one...	278	C19H18O2	056588-53-5	58
5	Benz[a]anthracene-7,12-dicarboni...	278	C20H10N2	035215-32-8	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070218\
Data File : BF107042.D
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Operator : JU/SJ
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Instrument : BNA_F
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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
2-Pentanone, 4-hy...	5.19	9.2	ng	166454	1	6.95	361692	20.0
unknown6.73	6.73	69.2	ng	1252170	1	6.95	361692	20.0
Benzaldehyde, 3,5...	12.11	2.2	ng	52161	4	11.49	482066	20.0
Fluoranthene, 2-m...	13.27	2.3	ng	82233	5	14.15	705009	20.0
Trifluoroacetoxy ...	13.95	3.9	ng	136810	5	14.15	705009	20.0
Cyclopenta(cd)pyr...	14.26	4.9	ng	172148	5	14.15	705009	20.0
Chrysene, 1-methyl-	14.55	3.3	ng	115451	5	14.15	705009	20.0
unknown14.71	14.71	2.3	ng	82456	5	14.15	705009	20.0
Benz[a]anthracene...	14.74	2.9	ng	101326	5	14.15	705009	20.0
Supraene	14.99	4.1	ng	70484	6	15.71	340786	20.0
Benzo[e]pyrene	15.57	8.3	ng	141181	6	15.71	340786	20.0
Pyridine-3-carbon...	17.47	2.5	ng	41966	6	15.71	340786	20.0