LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098176.D
Acq On : 31 Aug 2017 00:46
Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

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-BF081517.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	3.681	265	271	277	rBV	36372	54203	5.62%	0.361%
2	4.898	476	478	490	rBV	47745	54002	5.60%	0.359%
3	5.322	546	550	562	rBV	582636	542690	56.24%	3.611%
4	6.357	722	726	742	rBV	588050	588976	61.04%	3.919%
5	6.492	745	749	756	rVB	599297	559209	57.96%	3.721%
6	6.722	785	788	792	rBV	825315	725867	75.23%	4.830%
7	6.881	811	815	818	rBV	508667	445024	46.12%	2.961%
8	6.987	830	833	838	rBV4	12266	16243	1.68%	0.108%
9	7.287	880	884	894	rVB	347029	325668	33.75%	2.167%
10	7.645	943	945	949	rVB3	16947	17110	1.77%	0.114%
11	8.010	1001	1007	1009	rBV	966192	867540	89.91%	5.773%
12	8.075	1016	1018	1021	rVB3	17735	14515	1.50%	0.097%
13	8.216	1039	1042	1044	rBV	14121	12195	1.26%	0.081%
14	8.286	1049	1054	1055	rBV2	17669	20217	2.10%	0.135%
15	8.357	1063	1066	1069	rBV3	14388	15223	1.58%	0.101%
16	8.434	1075	1079	1081	rBV	19187	22917	2.38%	0.153%
17	8.510	1088	1092	1095	rBV4	12965	16422	1.70%	0.109%
18	8.551	1095	1099	1101	rBV2	20967	24734	2.56%	0.165%
19	8.604	1104	1108	1110	rBV4	13471	15452	1.60%	0.103%
20	8.722	1126	1128	1131	rBV	41617	31227	3.24%	0.208%
21	8.816	1141	1144	1153	rVB	45976	54698	5.67%	0.364%
22	8.886	1153	1156	1159	rBV5	18294	20994	2.18%	0.140%
23	8.916	1159	1161	1165	rVB3	14000	14964	1.55%	0.100%
24	8.957	1165	1168	1172	rBV6	8126	14252	1.48%	0.095%
25	8.998	1172	1175	1178	rBV4	14533	12516	1.30%	0.083%
26	9.033	1178	1181	1184	rVB2	28959	23786	2.47%	0.158%
27	9.081	1185	1189	1194	rBV	700173	564972	58.55%	3.760%
28	9.169	1199	1204	1208	rBV5	28227	50021	5.18%	0.333%
29	9.281	1220	1223	1227	rVB5	11381	14848	1.54%	0.099%
30	9.351	1231	1235	1238	rVV3	36268	51095	5.30%	0.340%
31	9.392	1238	1242	1244	rVV5	10932	18867	1.96%	0.126%
32	9.422	1244	1247	1249	rVV	53746	57994	6.01%	0.386%
33	9.445	1249	1251	1253	rVV	31431	29046	3.01%	0.193%
34	9.475	1253	1256	1260	rVV2	85141	119339	12.37%	0.794%

LSC Area Percent Report

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Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

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-BF081517.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.533	1260	1266	1271	rVB3	22613	37954	3.93%	0.253%
36	9.622	1278	1281	1287	rBV	41425	45256	4.69%	0.301%
37	9.669	1287	1289	1292	rVB	39753	28620	2.97%	0.190%
38	9.698	1292	1294	1298	rVB4	15326	19132	1.98%	0.127%
39	9.757	1298	1304	1307	rBV	845894	777905	80.62%	5.177%
40	9.792	1307	1310	1313	rVB	104936	98480	10.21%	0.655%
41	9.869	1320	1323	1327	rVB4	17535	22121	2.29%	0.147%
42	9.916	1327	1331	1335	rBV6	13182	19676	2.04%	0.131%
43	9.963	1335	1339	1344	rVV	80372	75233	7.80%	0.501%
44	10.075	1356	1358	1360	rBV3	20061	17718	1.84%	0.118%
45	10.163	1370	1373	1375	rBV3	18155	15403	1.60%	0.103%
46	10.192	1375	1378	1380	rVB2	11921	12202	1.26%	0.081%
47	10.304	1394	1397	1403	rVB	120062	110306	11.43%	0.734%
48	10.416	1413	1416	1418	rVB2	18095	15269	1.58%	0.102%
49	10.463	1421	1424	1428	rVB2	23183	23218	2.41%	0.155%
50	10.545	1434	1438	1442	rBV	234161	225703	23.39%	1.502%
51	10.675	1456	1460	1463	rBV3	28724	42011	4.35%	0.280%
52	10.851	1486	1490	1492	rBV2	17878	19408	2.01%	0.129%
53	11.092	1527	1531	1533	rBV4	10243	14468	1.50%	0.096%
54	11.139	1536	1539	1544	rVB2	48161	49846	5.17%	0.332%
55	11.198	1545	1549	1552	rBV	16299	15050	1.56%	0.100%
56	11.239	1552	1556	1558	rBV	754395	641785	66.52%	4.271%
57	11.263	1558	1560	1564	rVV	499669	392315	40.66%	2.611%
58	11.316	1564	1569	1572	rVV	128172	109230	11.32%	0.727%
59	11.351	1572	1575	1578	rVB4	14668	13976	1.45%	0.093%
60	11.469	1590	1595	1602	rBV2	56190	86172	8.93%	0.573%
61	11.539	1604	1607	1609	rVV2	22914	23005	2.38%	0.153%
62	11.569	1609	1612	1619	rVV3	25551	33784	3.50%	0.225%
63	11.639	1621	1624	1630	rVB	288450	232171	24.06%	1.545%
64	11.739	1637	1641	1644	rBV	67603	65834	6.82%	0.438%
65	11.769	1644	1646	1650	rVV	82851	67040	6.95%	0.446%
66	11.816	1650	1654	1657	rVV2	27684	27754	2.88%	0.185%
67	11.851	1657	1660	1663	rVV	134460	136330	14.13%	0.907%
68	11.874	1663	1664	1669	rVB	44499	27909	2.89%	0.186%
69	12.039	1689	1692	1694	rBV	41064	40989	4.25%	0.273%
70	12.063	1694	1696	1700	rVB5	26160	28108	2.91%	0.187%
71	12.216	1719	1722	1724	rBV	22365	18178	1.88%	0.121%

LSC Area Percent Report

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Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

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

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	12.239	1724	1726	1729	rVB	15667	13569	1.41%	0.090%
73	12.286	1731	1734	1737	rBV	48109	49630	5.14%	0.330%
74	12.322	1737	1740	1742	rBV	638310	577440	59.85%	3.843%
75	12.345	1742	1744	1747	rVB	889138	642809	66.62%	4.278%
76	12.427	1755	1758	1759	rBV	166778	130615	13.54%	0.869%
77	12.451	1759	1762	1765	rVV	704584	590774	61.23%	3.931%
78	12.516	1767	1773	1775	rBV7	12312	20434	2.12%	0.136%
79	12.598	1785	1787	1791	rBV2	17238	18357	1.90%	0.122%
80	12.674	1795	1800	1803	rBV	548118	604784	62.68%	4.025%
81	12.733	1807	1810	1815	rVB2	23912	29386	3.05%	0.196%
82	12.798	1818	1821	1822	rBV2	34833	31237	3.24%	0.208%
83	12.821	1822	1825	1831	rVB	512528	458250	47.49%	3.049%
84	12.898	1833	1838	1841	rBV2	39510	68467	7.10%	0.456%
85	12.980	1849	1852	1853	rBV3	33427	32845	3.40%	0.219%
86	13.004	1853	1856	1860	rVV	122935	116796	12.10%	0.777%
87	13.069	1864	1867	1870	rVB	112821	109313	11.33%	0.727%
88	13.110	1870	1874	1876	rBV2	60956	63596	6.59%	0.423%
89	13.198	1887	1889	1892	rVB	34208	26320	2.73%	0.175%
90	13.227	1892	1894	1898	rBV	36187	41452	4.30%	0.276%
91	13.263	1898	1900	1904	rVB5	21774	22526	2.33%	0.150%
92	13.621	1959	1961	1963	rBV	30781	25428	2.64%	0.169%
93	13.651	1964	1966	1968	rBV	28375	19925	2.07%	0.133%
94	13.727	1976	1979	1981	rBV3	33623	38206	3.96%	0.254%
95	13.874	1999	2004	2006	rBV2	822671	964869	100.00%	6.421%
96	13.898	2006	2008	2011	rVB	247348	193498	20.05%	1.288%
97	13.974	2019	2021	2025	rBV	59831	51228	5.31%	0.341%
98	14.886	2172	2176	2183	rBV	191099	361883	37.51%	2.408%
99	15.227	2231	2234	2239	rBV	158168	195971	20.31%	1.304%
100	15.286	2240	2244	2248	rBV	387026	481066	49.86%	3.201%

Sum of corrected areas: 15027059

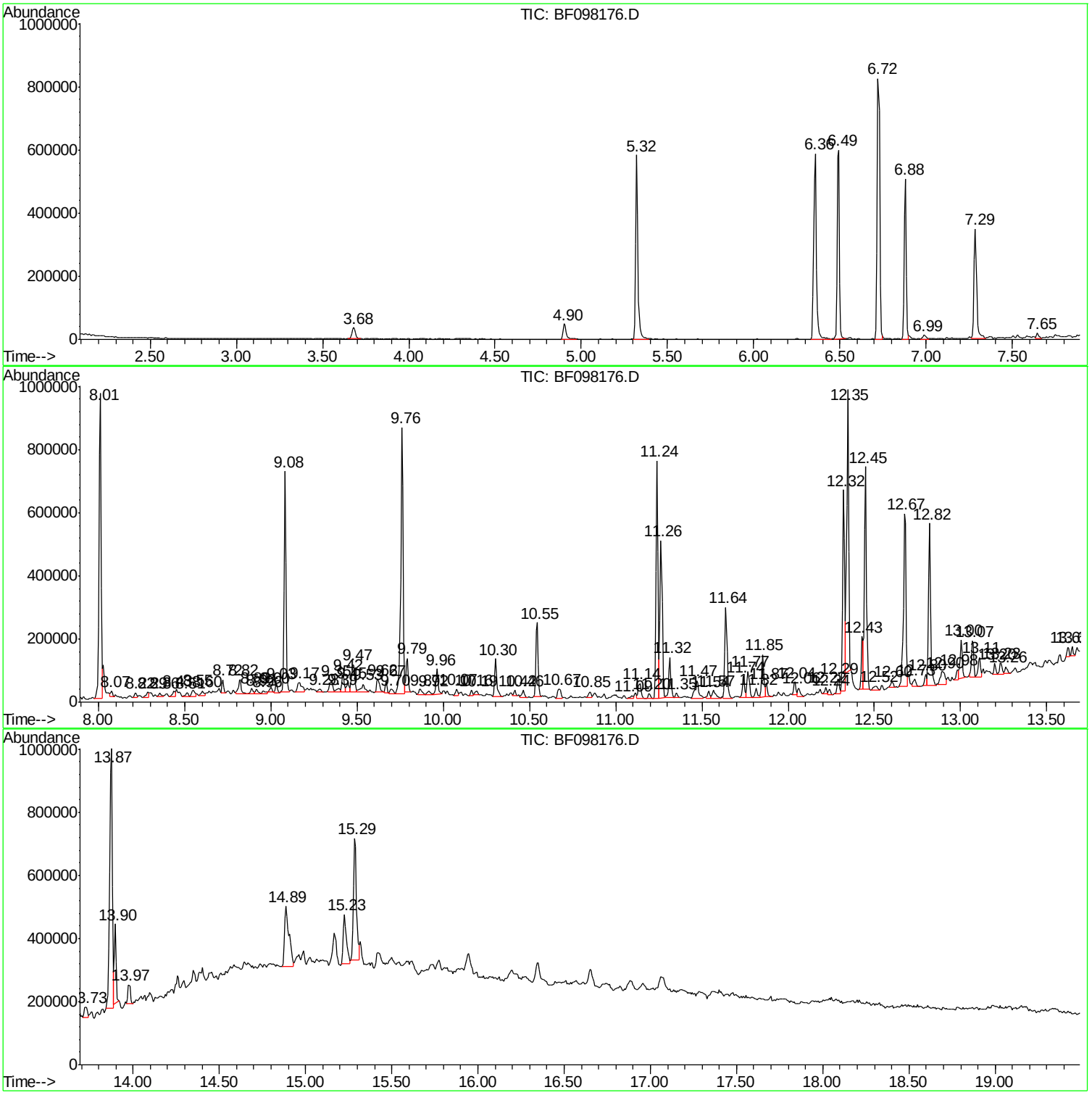
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098176.D
Acq On : 31 Aug 2017 00:46
Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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\BF083017\
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Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

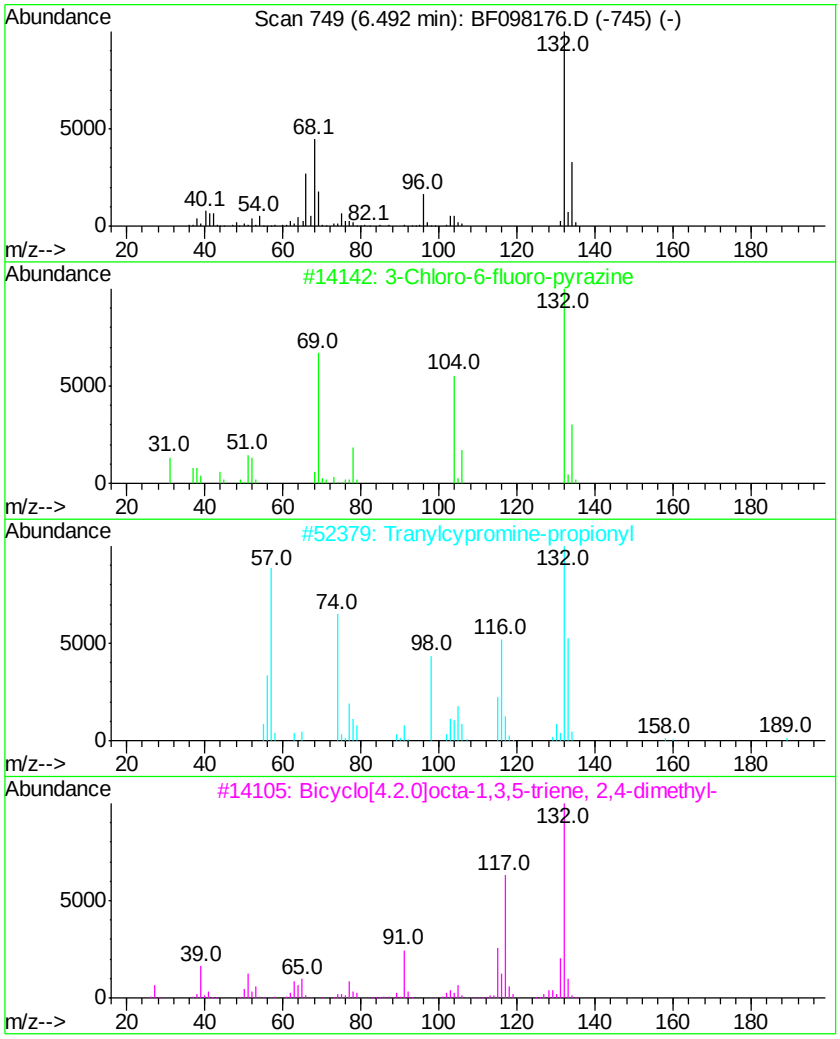
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

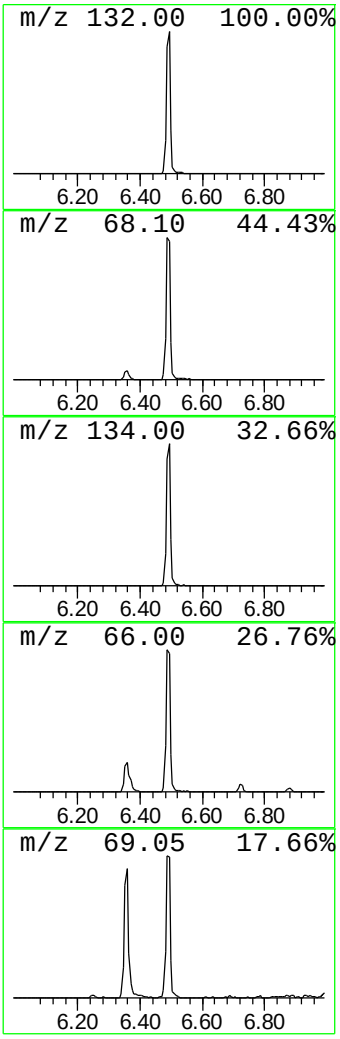
****

Peak Number 1 unknown6.49 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	15.41 ng	559209	1,4-Dichlorobenzene-d4	6.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	Xenon	132	Xe	007440-63-3	9
5	1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
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Acq On : 31 Aug 2017 00:46
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Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

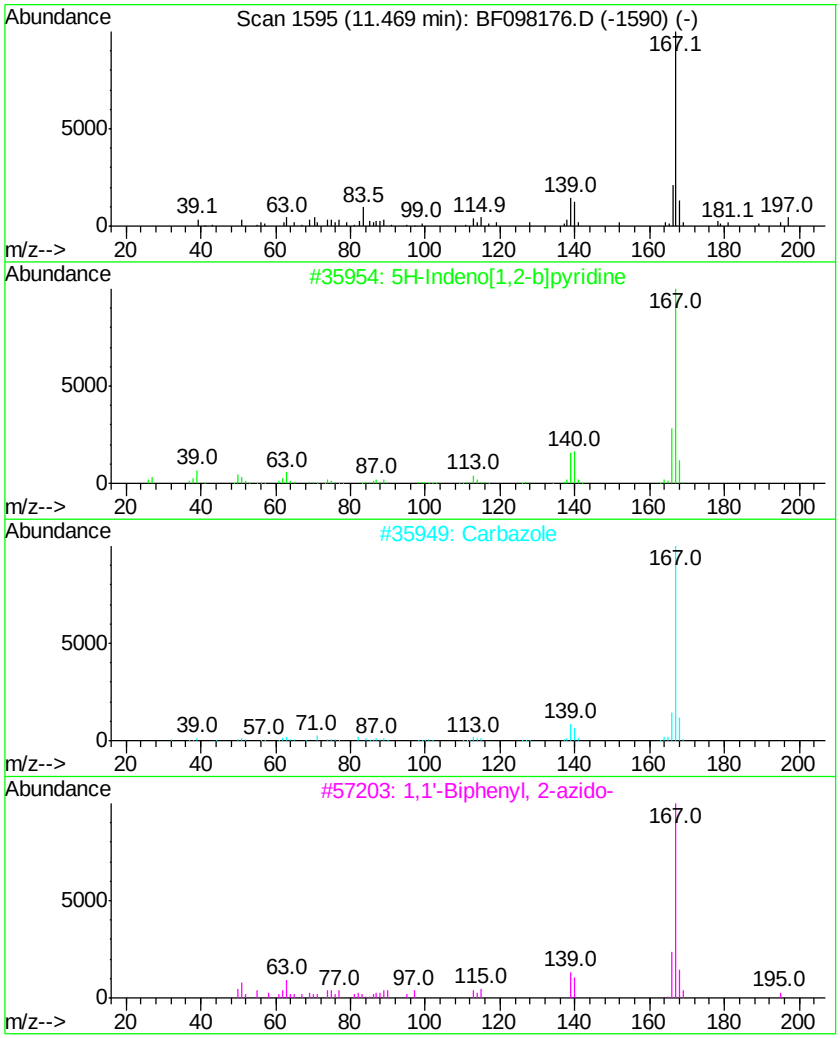
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

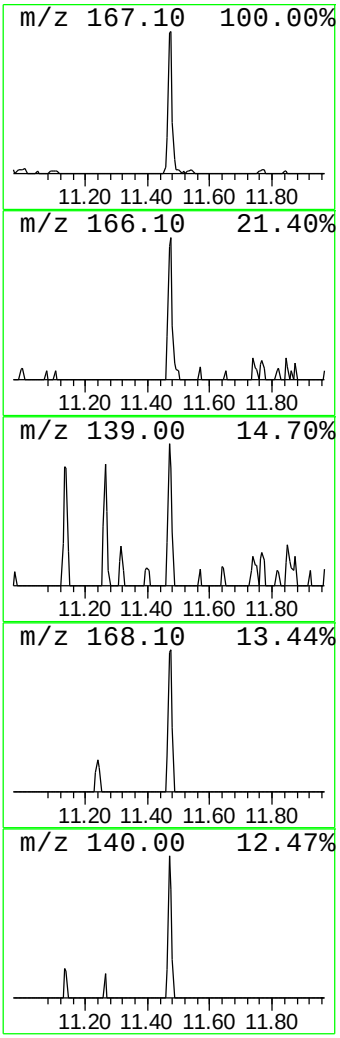
****

Peak Number 3 5H-Indeno[1,2-b]pyridine Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	2.69 ng	86172	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	94
2	Carbazole	167	C12H9N	000086-74-8	87
3	1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	83
4	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	83
5	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	72





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Sample : I5007-01 5X
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Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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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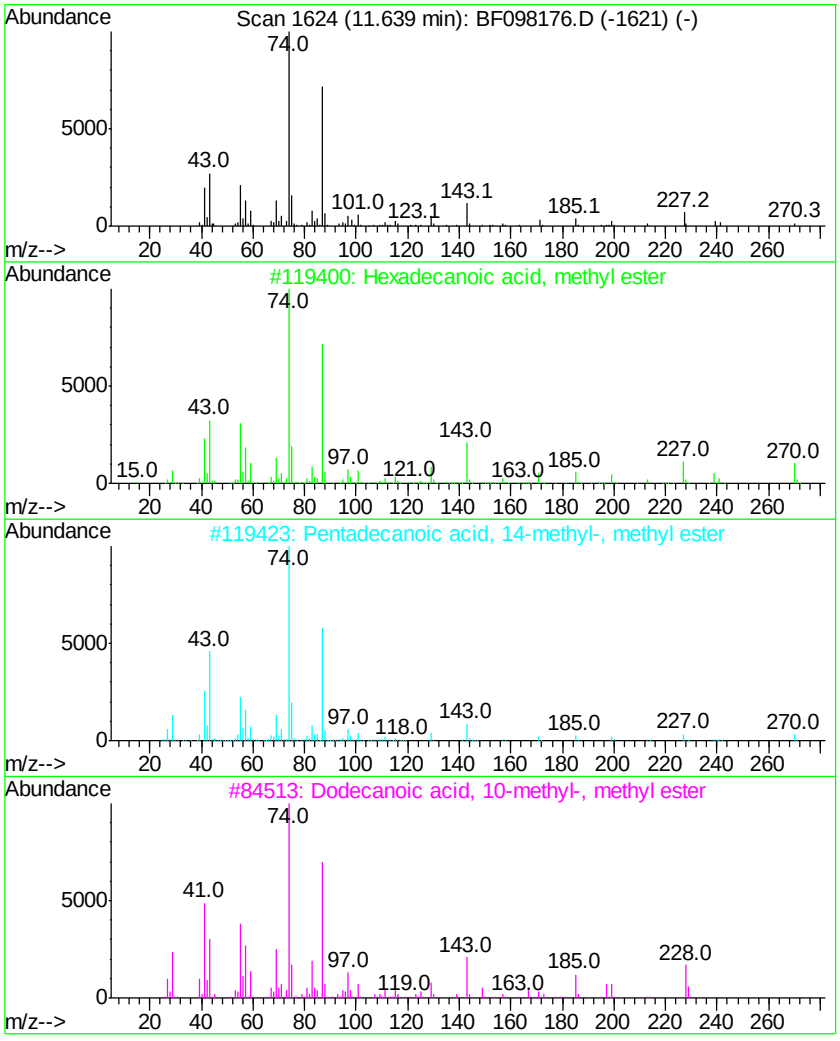
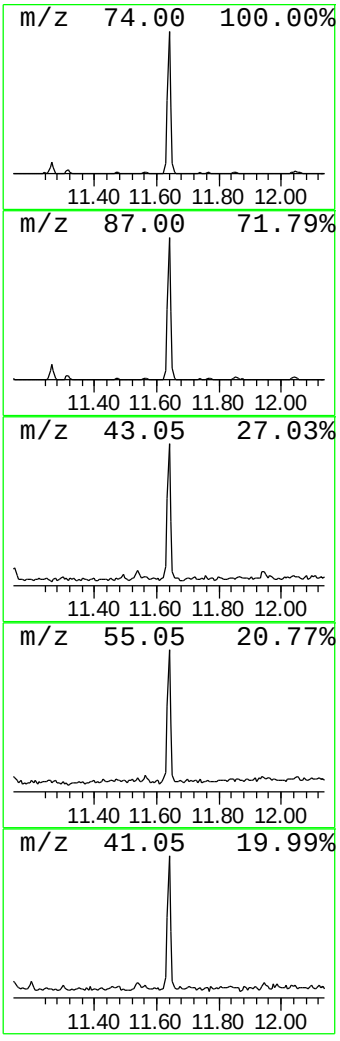
Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	7.24 ng	232171	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
3	Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	87
4	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	87
5	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	87





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Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

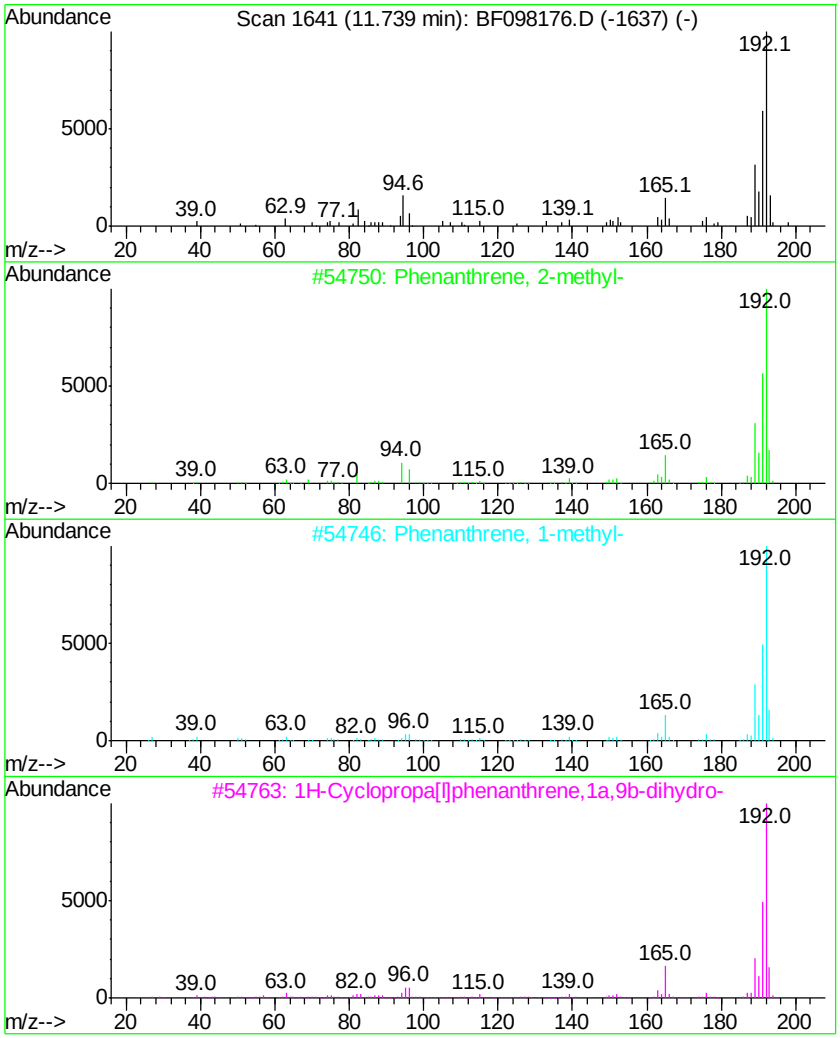
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

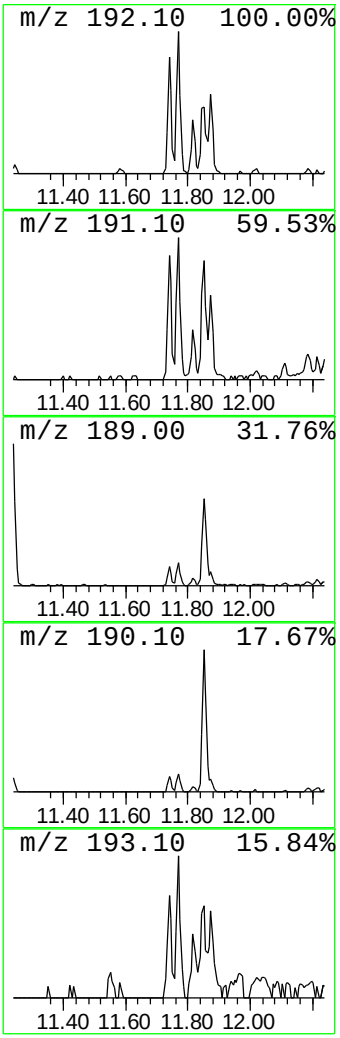
****

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	2.05 ng	65834	Phenanthrene-d10	11.24

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	96
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098176.D
Acq On : 31 Aug 2017 00:46
Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

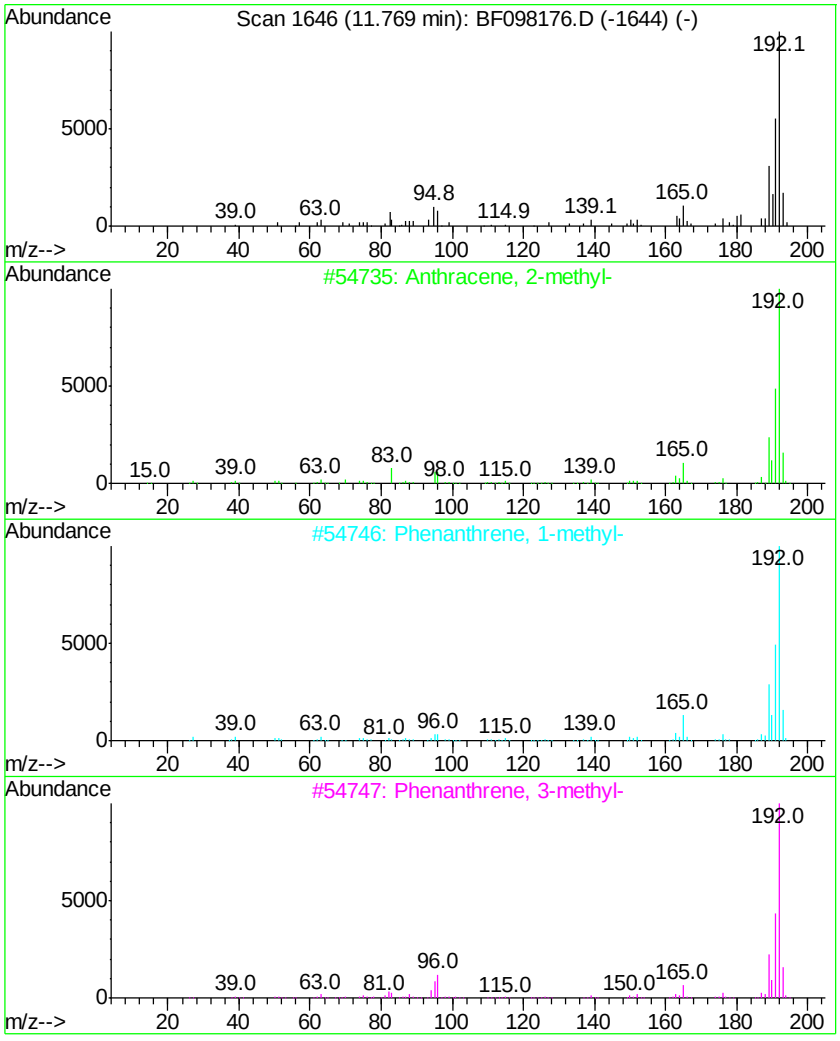
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

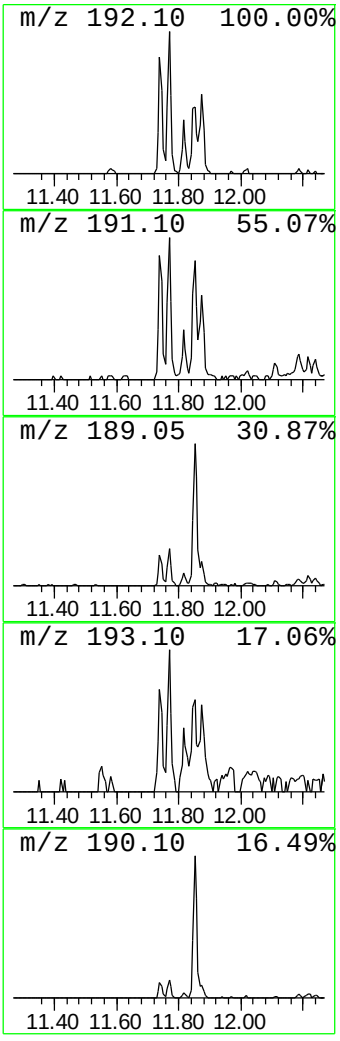
****

Peak Number 6 Anthracene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	2.09 ng	67040	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	94
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098176.D
Acq On : 31 Aug 2017 00:46
Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

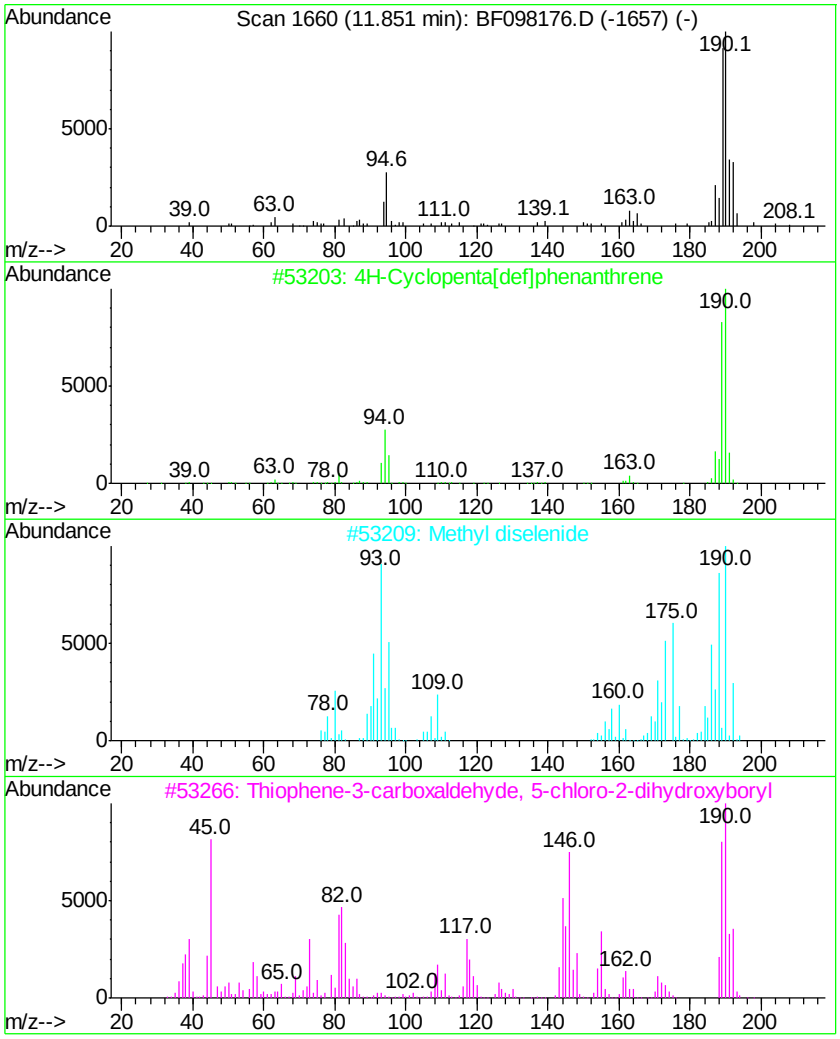
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

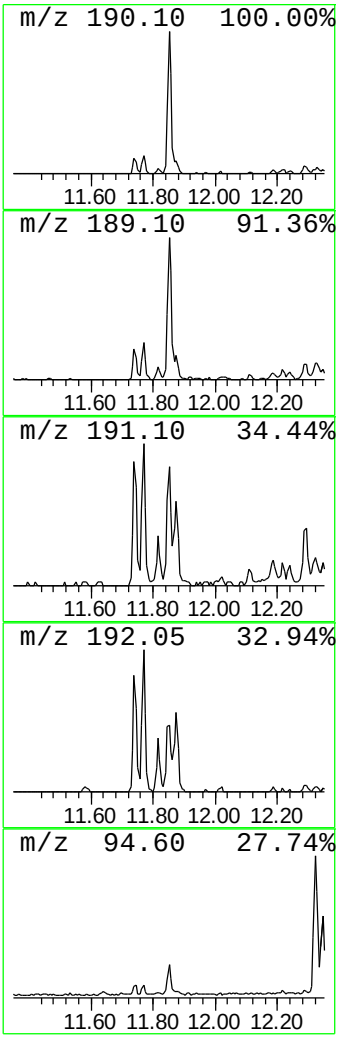
****

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	4.25 ng	136330	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	Methyl diselenide	190	C2H6Se2	007101-31-7	49
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	47
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	28
5	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	12





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Operator : SJ/JU
Sample : I5007-01 5X
Misc :
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Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
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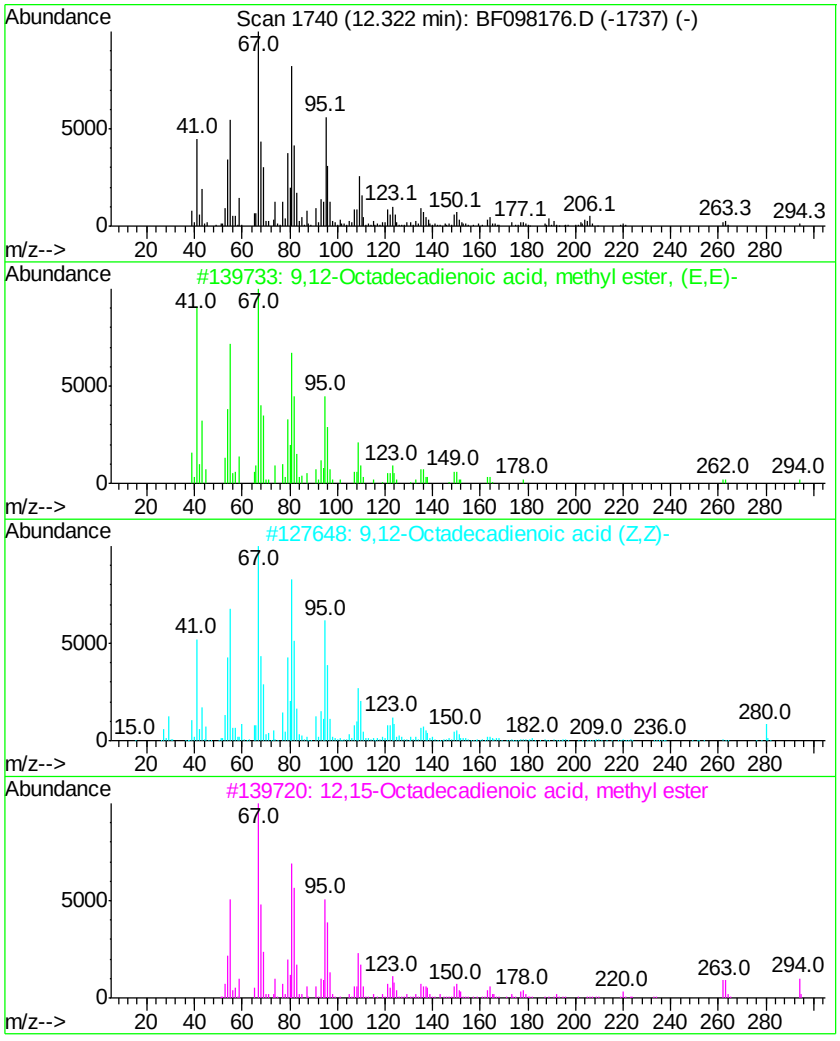
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

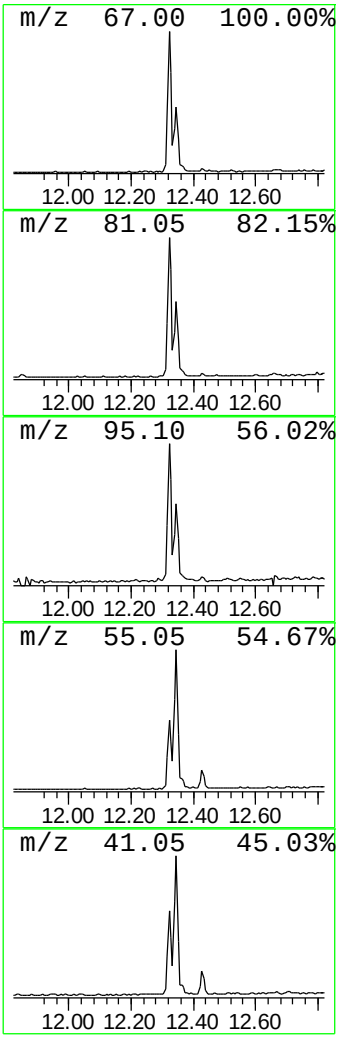
****

Peak Number 8 9,12-Octadecadienoic acid, ... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	17.99 ng	577440	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	95
3	12,15-Octadecadienoic acid, meth...	294	C19H34O2	057156-97-5	91
4	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	91
5	2-Chloroethyl linoleate	342	C20H35ClO2	025525-76-2	87





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098176.D
Acq On : 31 Aug 2017 00:46
Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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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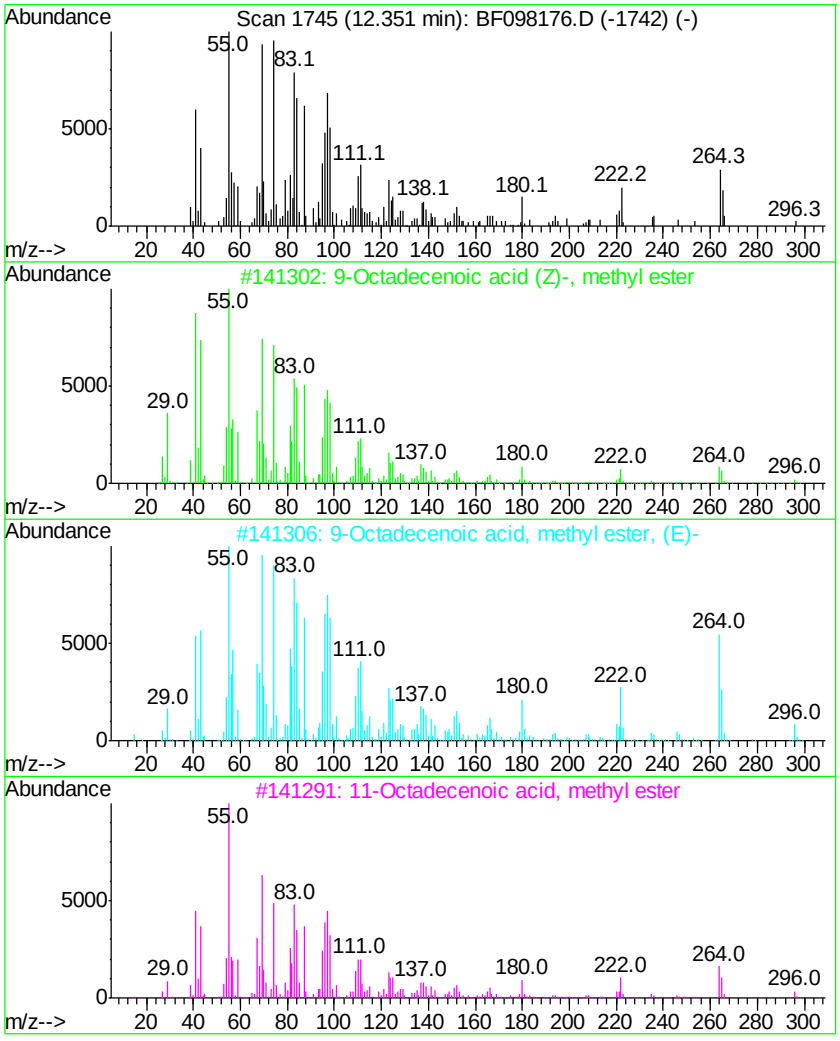
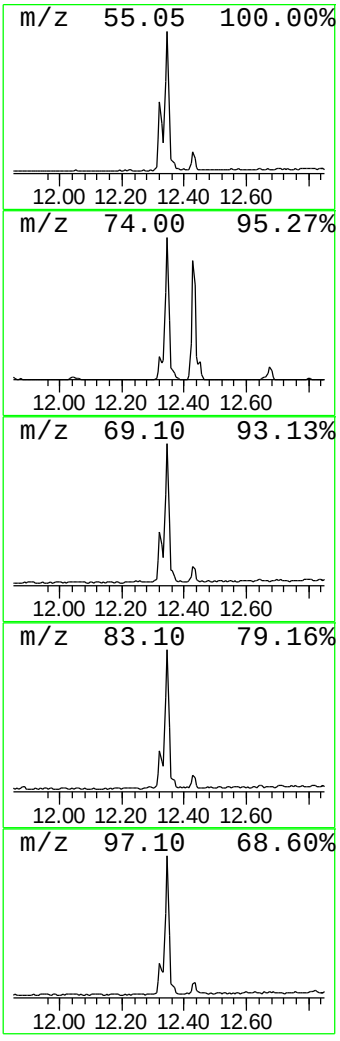
Peak Number 9 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	20.03 ng	642809	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	94
2	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	91
3	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	86
4	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	86
5	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	76





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ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

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TIC Integration Parameters: LSCINT.P

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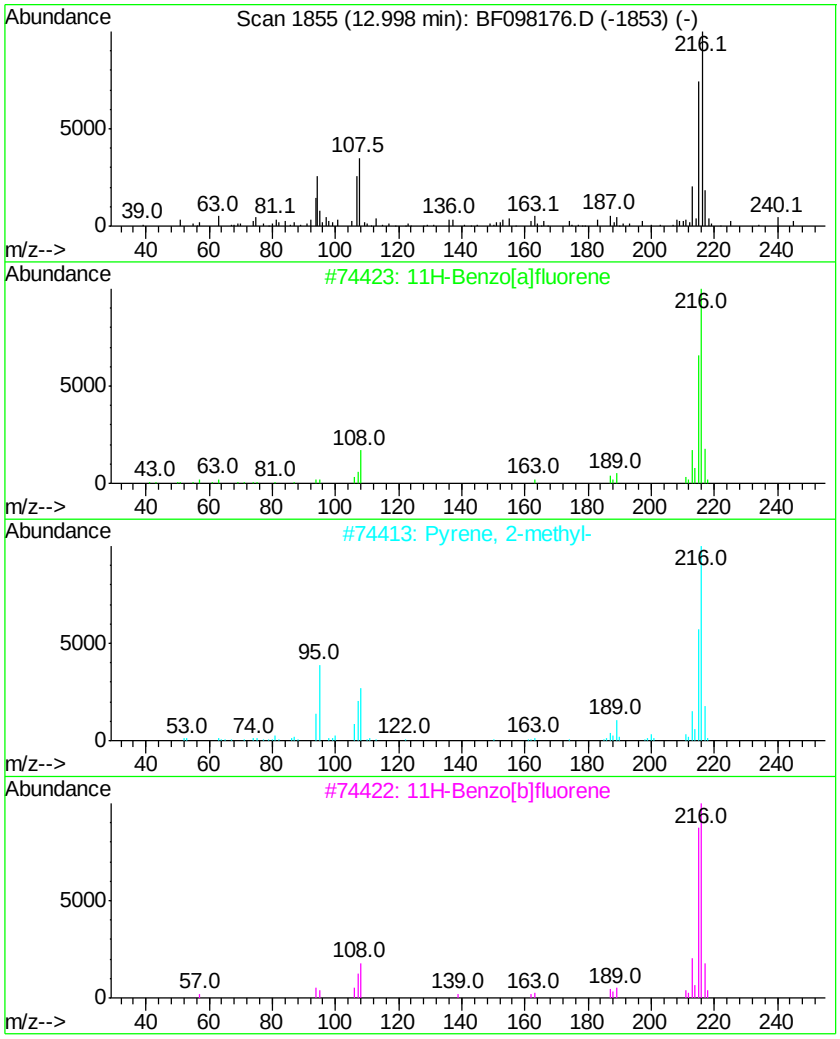
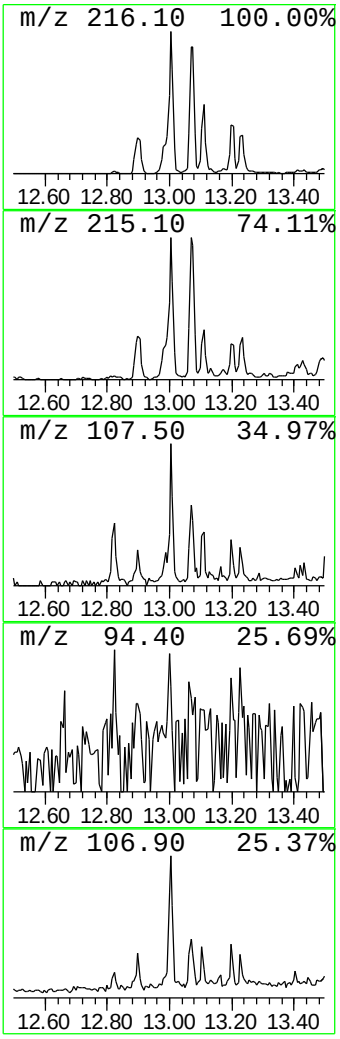
Peak Number 10 11H-Benzo[a]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.00	2.42 ng	116796	Chrysene-d12	13.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90





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Acq On : 31 Aug 2017 00:46
Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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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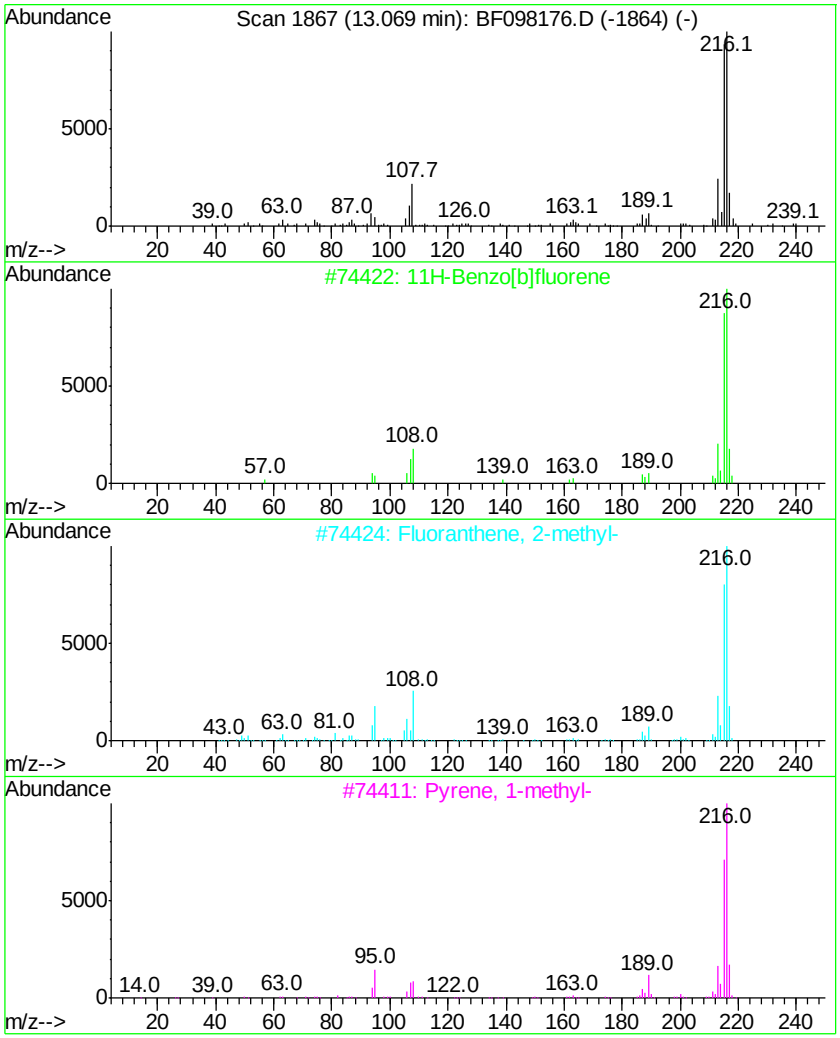
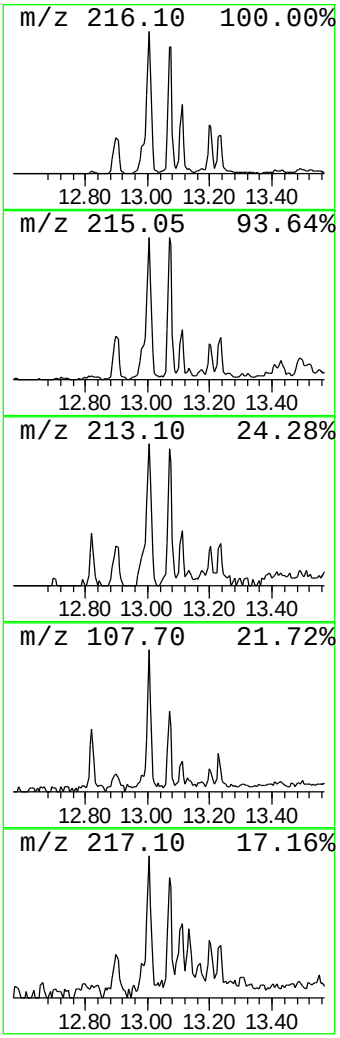
Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.07	2.27 ng	109313	Chrysene-d12	13.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	98
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF083017\
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Operator : SJ/JU
Sample : I5007-01 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-082817

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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
unknown6.49	6.49	15.4	ng	559209	1	6.72	725867	20.0
5H-Indeno[1,2-b]p...	11.47	2.7	ng	86172	4	11.24	641785	20.0
Hexadecanoic acid...	11.64	7.2	ng	232171	4	11.24	641785	20.0
Phenanthrene, 2-m...	11.74	2.0	ng	65834	4	11.24	641785	20.0
Anthracene, 2-met...	11.77	2.1	ng	67040	4	11.24	641785	20.0
4H-Cyclopenta[def...	11.85	4.3	ng	136330	4	11.24	641785	20.0
9,12-Octadecadien...	12.32	18.0	ng	577440	4	11.24	641785	20.0
9-Octadecenoic ac...	12.35	20.0	ng	642809	4	11.24	641785	20.0
11H-Benzo[a]fluorene	13.00	2.4	ng	116796	5	13.87	964869	20.0
11H-Benzo[b]fluorene	13.07	2.3	ng	109313	5	13.87	964869	20.0