LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

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-BF101819.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	4.281	370	373	378	rBV2	7437	11660	2.24%	0.128%
2	4.922	479	482	494	rBV	17623	24224	4.65%	0.265%
3	5.357	553	556	571	rBV	292718	387162	74.28%	4.243%
4	6.387	728	731	748	rBV	345496	450895	86.51%	4.941%
5	6.510	748	752	758	rBV	484728	472468	90.65%	5.177%
6	6.734	786	790	804	rVB	328627	308603	59.21%	3.382%
7	6.886	813	816	827	rBV	412887	356481	68.39%	3.906%
8	7.298	883	886	905	rBV	248999	274755	52.71%	3.011%
9	8.016	1004	1008	1011	rBV	479781	408357	78.35%	4.475%
10	8.728	1127	1129	1134	rBV	11895	12589	2.42%	0.138%
11	8.828	1143	1146	1151	rBB	14991	12760	2.45%	0.140%
12	9.086	1187	1190	1201	rBV	669789	521213	100.00%	5.712%
13	9.357	1233	1236	1240	rBB2	5067	6816	1.31%	0.075%
14	9.422	1244	1247	1250	rVV2	22129	20913	4.01%	0.229%
15	9.480	1254	1257	1263	rVV	40577	44986	8.63%	0.493%
16	9.545	1265	1268	1274	rBB	5271	6695	1.28%	0.073%
17	9.627	1279	1282	1286	rBV	18349	15721	3.02%	0.172%
18	9.763	1302	1305	1309	rBV	508750	467398	89.68%	5.122%
19	9.798	1309	1311	1318	rVB	51195	43595	8.36%	0.478%
20	9.975	1337	1341	1345	rBV	24268	25842	4.96%	0.283%
21	10.316	1396	1399	1404	rBV	55259	51629	9.91%	0.566%
22	10.475	1422	1426	1429	rVB	8469	8785	1.69%	0.096%
23	10.557	1437	1440	1445	rBV	276638	260185	49.92%	2.851%
24	10.680	1455	1461	1466	rVB6	9625	18983	3.64%	0.208%
25	10.863	1488	1492	1495	rVB2	12267	12962	2.49%	0.142%
26	10.992	1512	1514	1516	rBV3	7517	7163	1.37%	0.078%
27	11.069	1523	1527	1530	rBV2	10419	9385	1.80%	0.103%
28	11.110	1530	1534	1538	rVB2	8141	12931	2.48%	0.142%
29	11.145	1538	1540	1547	rVB	16690	18173	3.49%	0.199%
30	11.204	1547	1550	1554	rBV2	9719	11097	2.13%	0.122%
31	11.251	1554	1558	1560	rBV2	452895	394359	75.66%	4.322%
32	11.274	1560	1562	1566	rVV	450930	374058	71.77%	4.099%
33	11.327	1568	1571	1575	rVB	91193	81846	15.70%	0.897%
34	11.492	1595	1599	1605	rBV2	26473	33509	6.43%	0.367%

LSC Area Percent Report

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Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.539	1605	1607	1610	rVB2	8115	7496	1.44%	0.082%
36	11.710	1634	1636	1640	rBV4	8055	10779	2.07%	0.118%
37	11.751	1640	1643	1646	rVV	43638	43831	8.41%	0.480%
38	11.780	1646	1648	1654	rVV	104705	106184	20.37%	1.164%
39	11.833	1654	1657	1659	rVV	21366	23538	4.52%	0.258%
40	11.863	1659	1662	1665	rVV	67032	76320	14.64%	0.836%
41	11.886	1665	1666	1670	rVV	28152	21806	4.18%	0.239%
42	11.957	1673	1678	1681	rBV7	7457	14917	2.86%	0.163%
43	12.045	1691	1693	1698	rVB	29126	25056	4.81%	0.275%
44	12.092	1698	1701	1705	rBV3	17703	18943	3.63%	0.208%
45	12.198	1714	1719	1721	rBV3	7864	6924	1.33%	0.076%
46	12.227	1721	1724	1727	rVV4	7748	8143	1.56%	0.089%
47	12.257	1727	1729	1734	rVB2	8014	6663	1.28%	0.073%
48	12.304	1734	1737	1740	rBV	26384	26331	5.05%	0.289%
49	12.339	1740	1743	1746	rVV2	23138	28013	5.37%	0.307%
50	12.398	1749	1753	1756	rVB3	14343	18630	3.57%	0.204%
51	12.468	1761	1765	1768	rBV	447871	419568	80.50%	4.598%
52	12.510	1770	1772	1774	rVV2	10658	9223	1.77%	0.101%
53	12.533	1774	1776	1778	rVV3	10138	7946	1.52%	0.087%
54	12.563	1778	1781	1785	rBV3	19112	19123	3.67%	0.210%
55	12.616	1788	1790	1793	rVV2	17664	16556	3.18%	0.181%
56	12.692	1797	1803	1810	rVV	330621	404720	77.65%	4.435%
57	12.745	1810	1812	1817	rVB2	19140	20911	4.01%	0.229%
58	12.833	1821	1827	1830	rBV	418032	379641	72.84%	4.160%
59	12.863	1830	1832	1836	rVB	12817	12125	2.33%	0.133%
60	12.910	1836	1840	1845	rBV4	20881	38441	7.38%	0.421%
61	13.004	1852	1856	1857	rBV3	19314	23548	4.52%	0.258%
62	13.027	1857	1860	1862	rVV2	49498	50747	9.74%	0.556%
63	13.063	1862	1866	1868	rVV	25048	24583	4.72%	0.269%
64	13.092	1868	1871	1873	rVV	39928	36390	6.98%	0.399%
65	13.127	1873	1877	1884	rVV3	35105	52731	10.12%	0.578%
66	13.180	1884	1886	1890	rVB3	9639	8352	1.60%	0.092%
67	13.221	1890	1893	1896	rBV	17199	18072	3.47%	0.198%
68	13.257	1896	1899	1903	rBV2	41187	48607	9.33%	0.533%
69	13.298	1903	1906	1908	rVV4	15292	18475	3.54%	0.202%
70	13.316	1908	1909	1914	rVV5	4950	7386	1.42%	0.081%
71	13.404	1920	1924	1926	rBV4	15298	18667	3.58%	0.205%

LSC Area Percent Report

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Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

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
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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	13.427	1926	1928	1930	rBV2	9668	8879	1.70%	0.097%
73	13.521	1940	1944	1945	rBV4	9585	13784	2.64%	0.151%
74	13.545	1945	1948	1951	rVB	25795	26035	5.00%	0.285%
75	13.645	1962	1965	1967	rBV2	32431	34139	6.55%	0.374%
76	13.668	1967	1969	1971	rVB	27574	20165	3.87%	0.221%
77	13.698	1971	1974	1976	rBV	29767	32101	6.16%	0.352%
78	13.733	1978	1980	1981	rBV2	28776	22991	4.41%	0.252%
79	13.892	2003	2007	2010	rBV2	370537	471223	90.41%	5.164%
80	13.921	2010	2012	2020	rVB	175342	163128	31.30%	1.788%
81	13.998	2023	2025	2028	rVB	26618	23829	4.57%	0.261%
82	14.051	2032	2034	2040	rBV6	19966	44185	8.48%	0.484%
83	14.286	2072	2074	2076	rVB	31445	21648	4.15%	0.237%
84	14.351	2083	2085	2087	rBV3	26955	28286	5.43%	0.310%
85	14.380	2087	2090	2094	rVV4	28820	39034	7.49%	0.428%
86	14.715	2145	2147	2150	rBV	20836	15215	2.92%	0.167%
87	14.921	2178	2182	2185	rBV	142968	202376	38.83%	2.218%
88	15.209	2228	2231	2237	rVB	73385	83385	16.00%	0.914%
89	15.274	2238	2242	2245	rBV	102160	121771	23.36%	1.334%
90	15.333	2247	2252	2255	rBV	205013	265063	50.86%	2.905%
91	15.727	2316	2319	2323	rVB2	39982	50679	9.72%	0.555%
92	16.745	2487	2492	2499	rVB2	52723	98399	18.88%	1.078%
93	16.815	2501	2504	2510	rVB8	24972	41559	7.97%	0.455%
94	17.174	2562	2565	2572	rVB2	27714	50015	9.60%	0.548%

Sum of corrected areas: 9125473

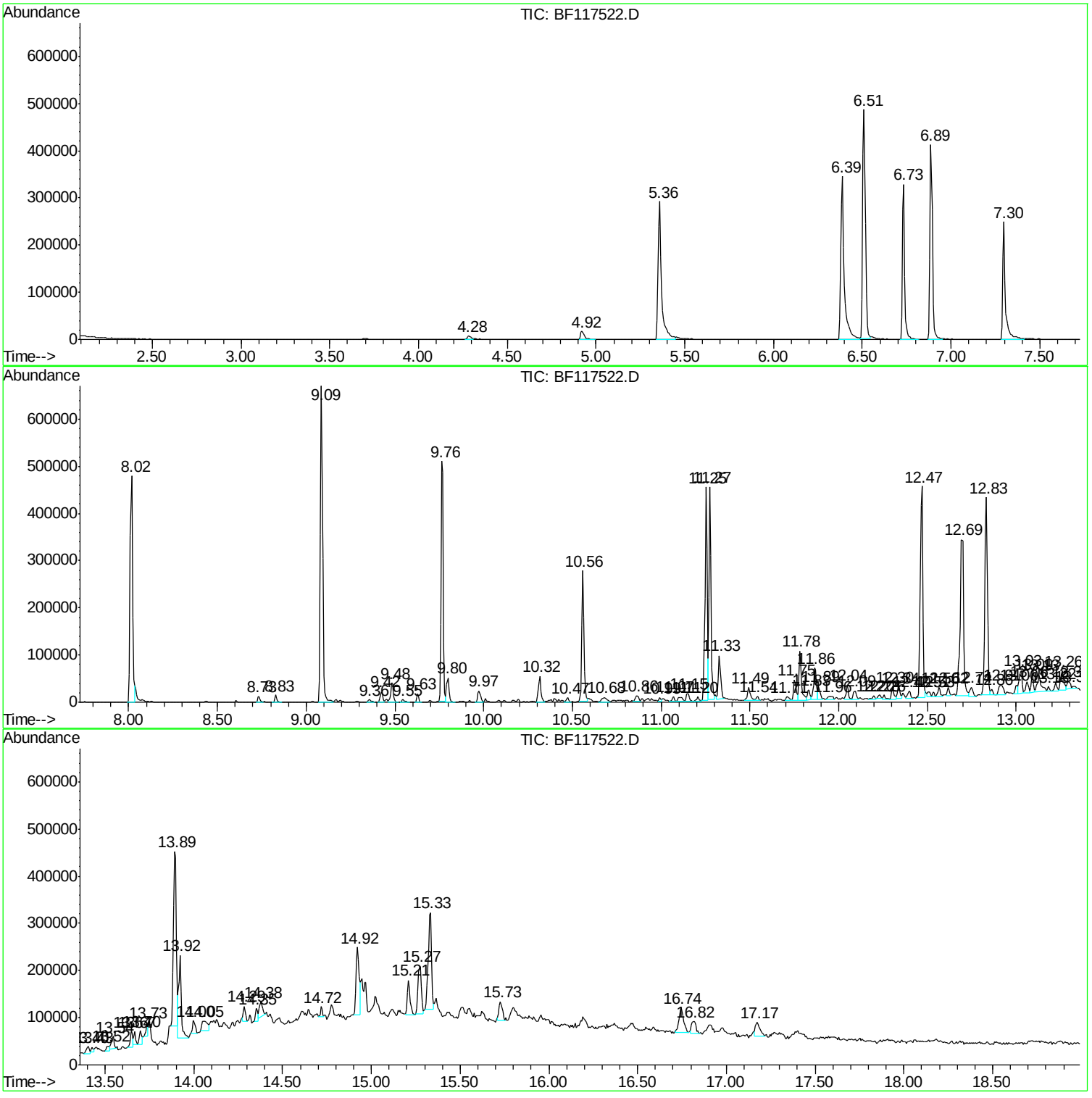
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
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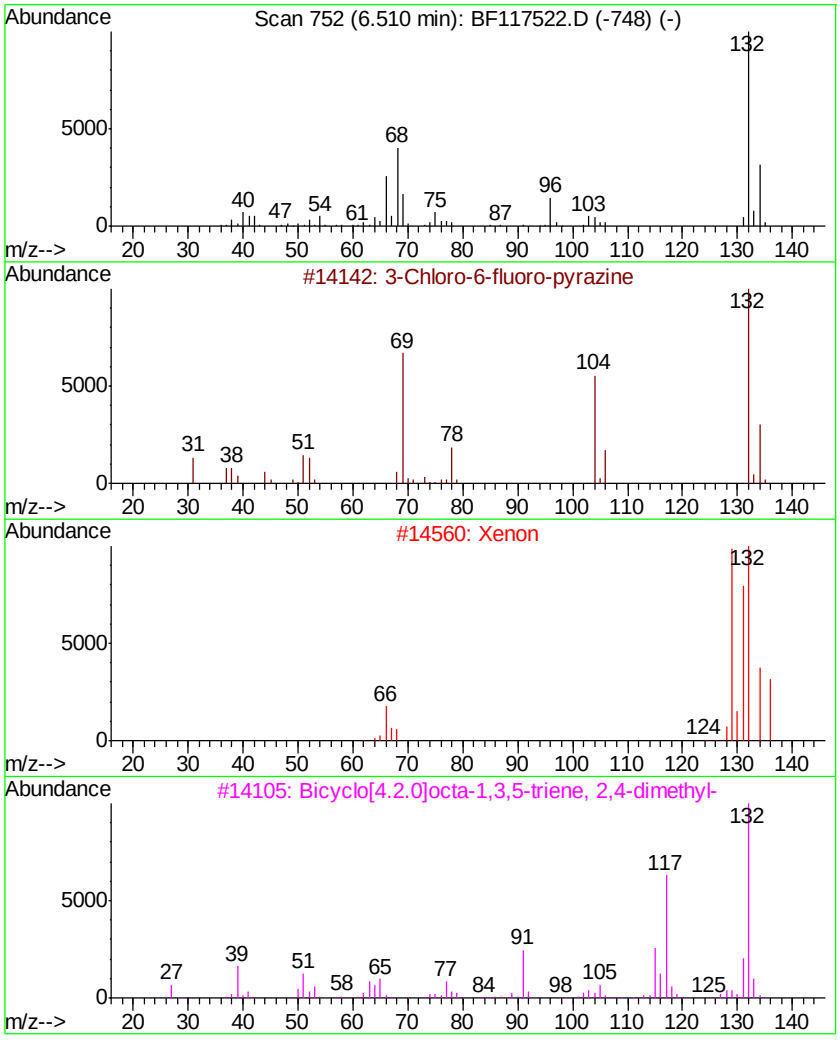
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

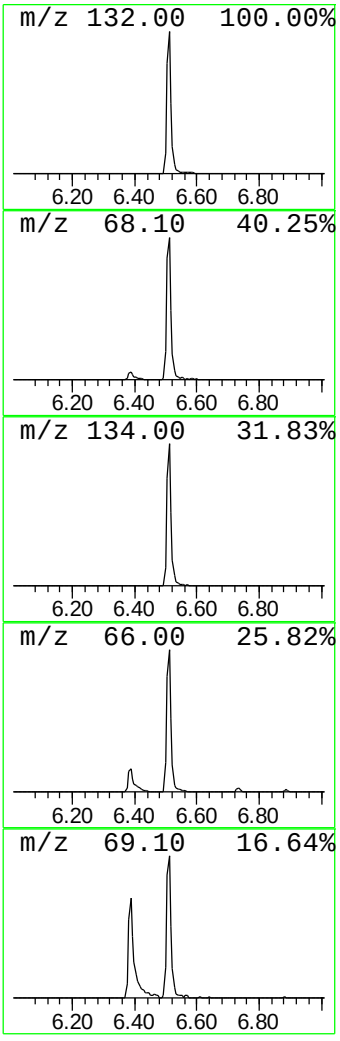
****

Peak Number 1 unknown6.51 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	30.62 ng	472468	1,4-Dichlorobenzene-d4	6.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2	Xenon	132	Xe	007440-63-3	9
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





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Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

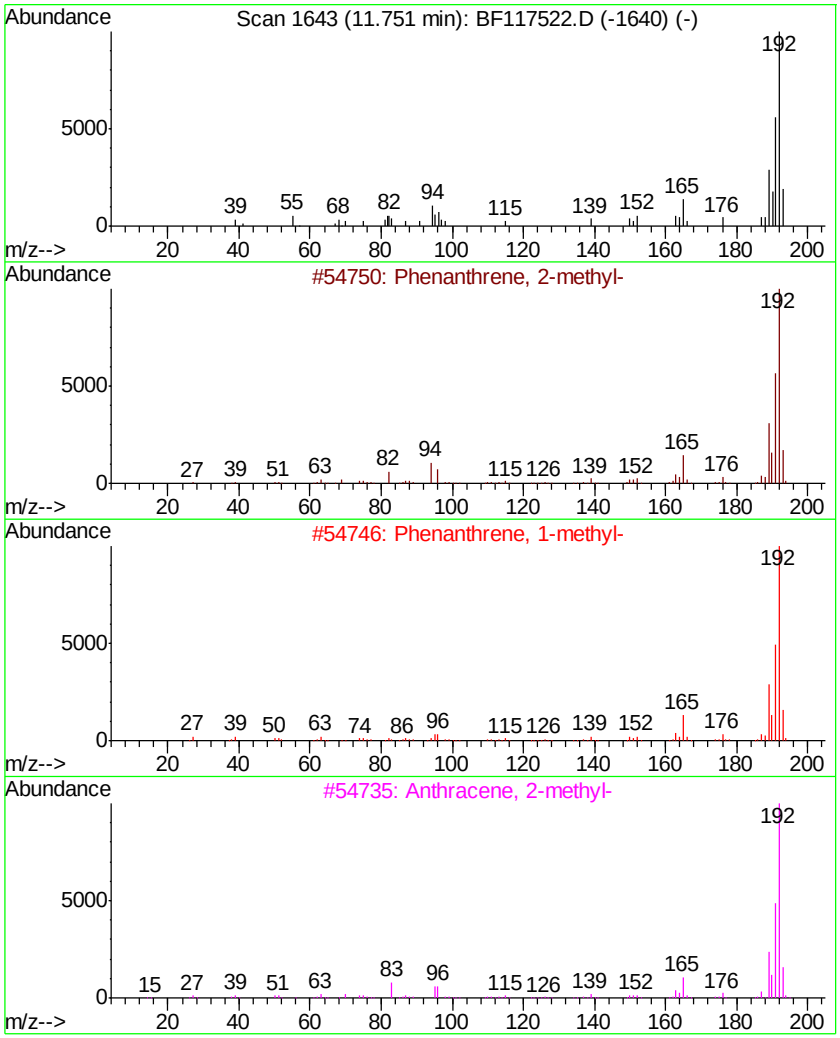
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

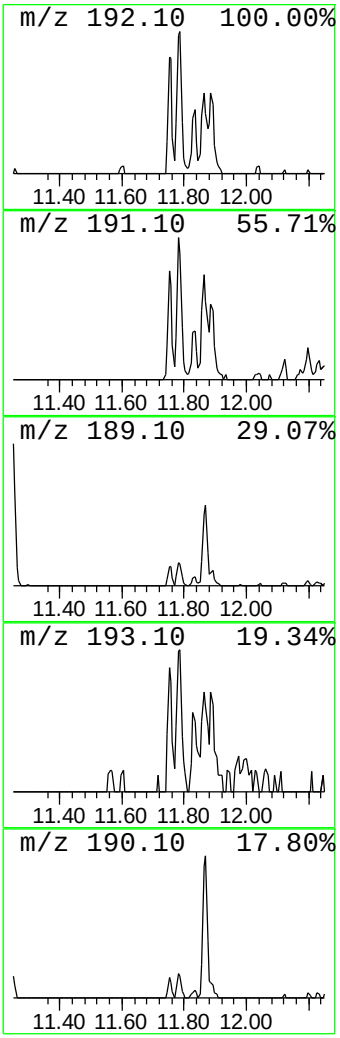
****

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	2.22 ng	43831	Phenanthrene-d10	11.25

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	94
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93





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Misc :
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Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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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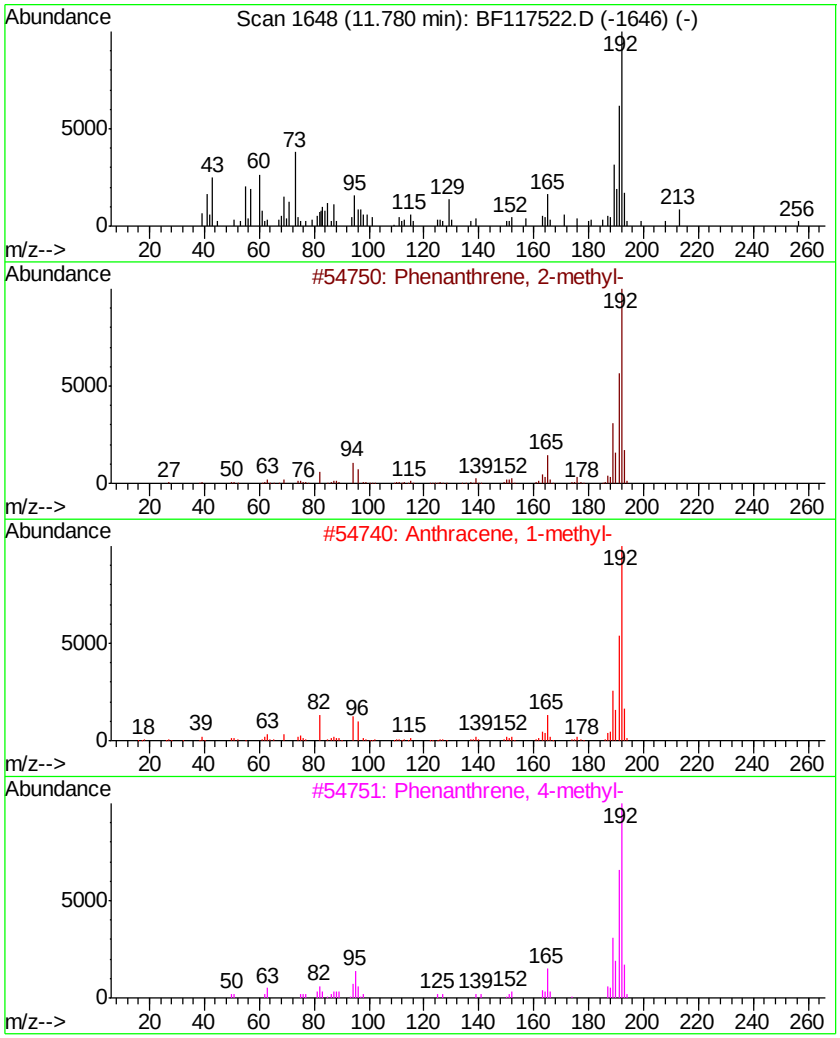
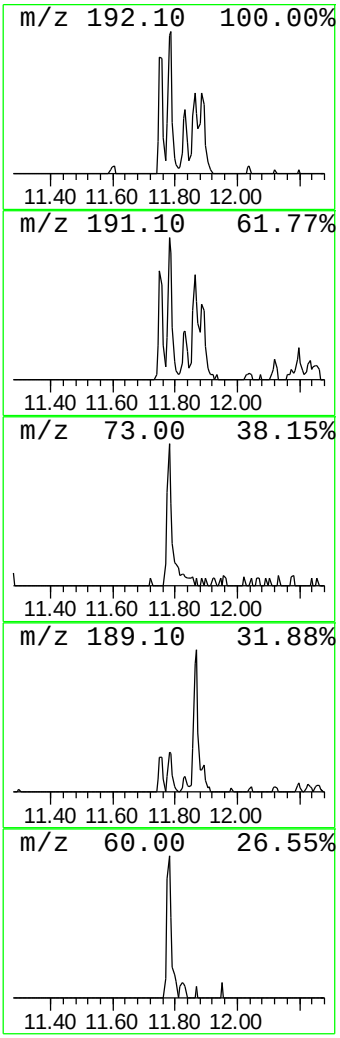
Peak Number 4 Anthracene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	5.39 ng	106184	Phenanthrene-d10	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95





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Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

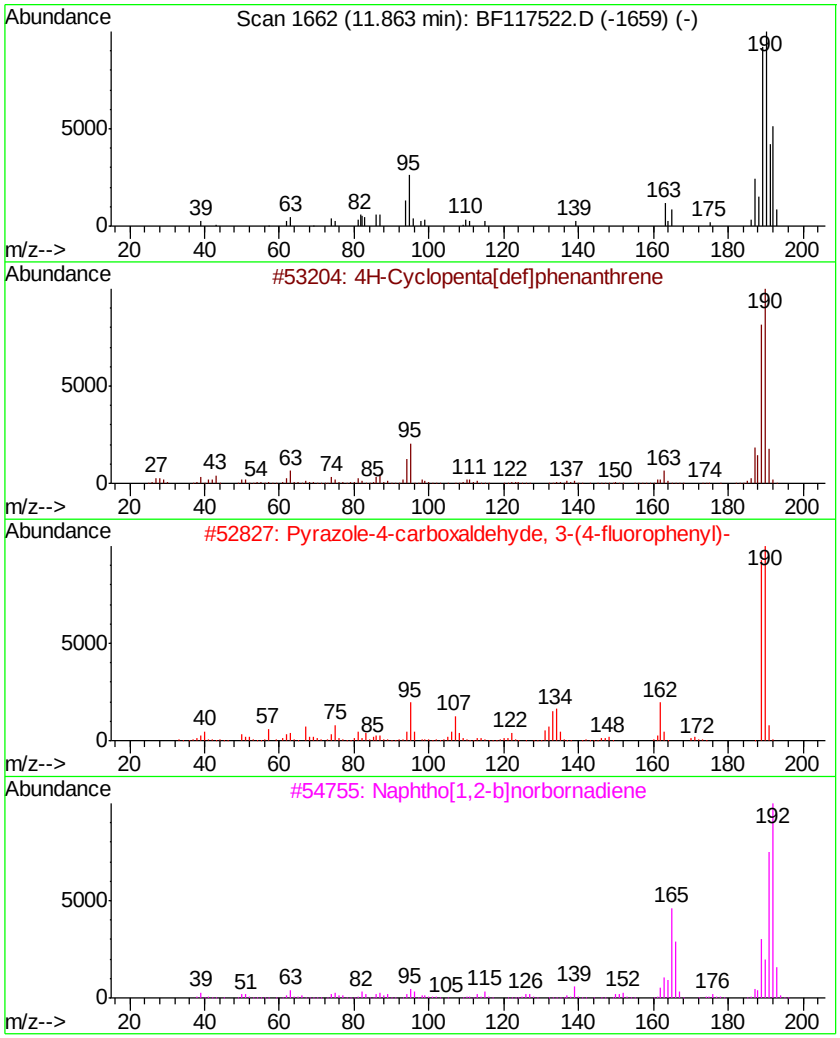
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

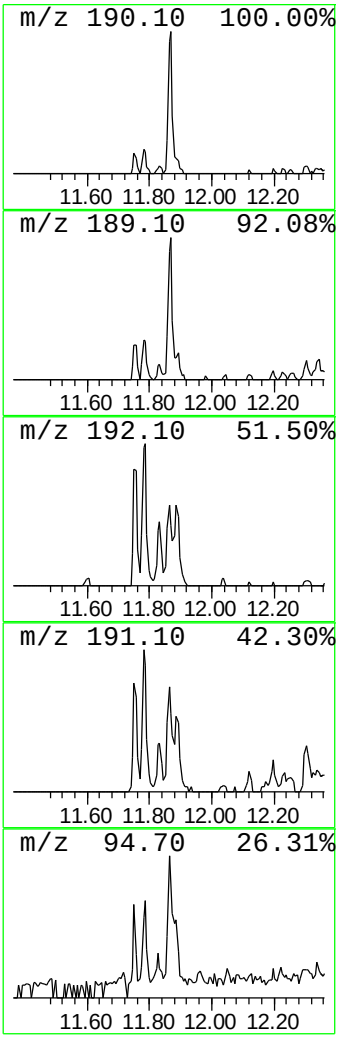
****

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	3.87 ng	76320	Phenanthrene-d10	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	37
3	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	25
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	14
5	Benzene, 1,1'-(2-cyclopropen-1-y...	192	C15H12	022825-21-4	14





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Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

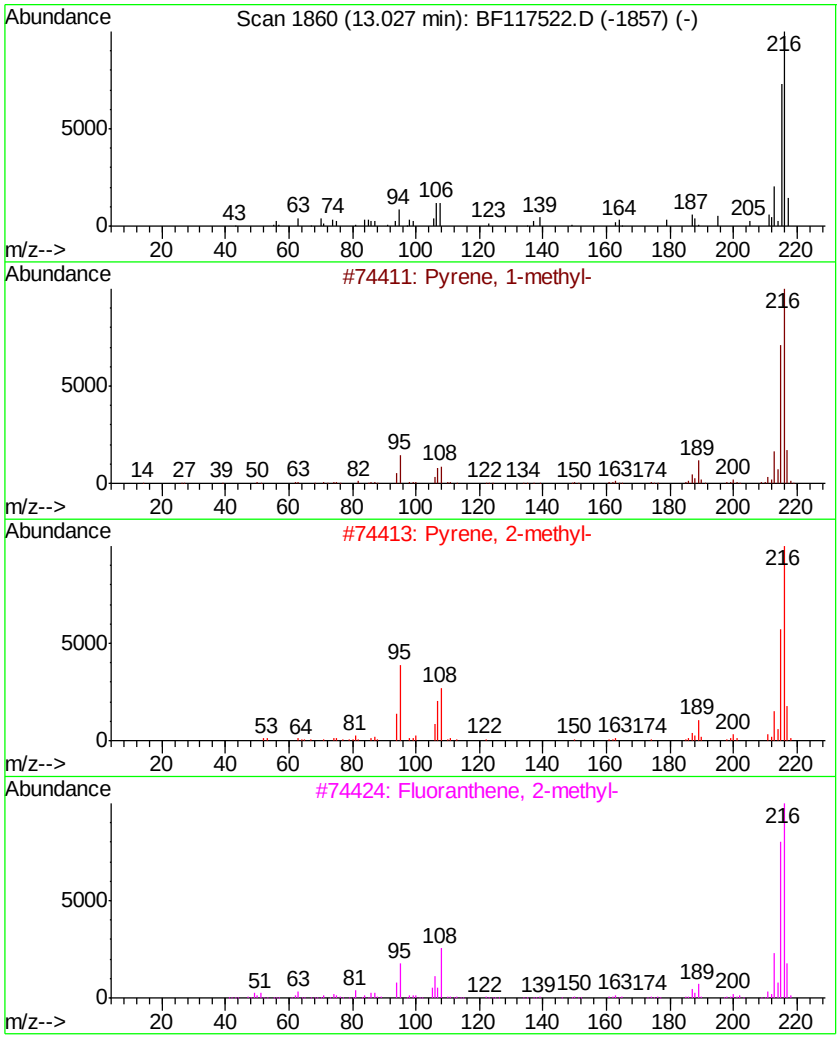
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

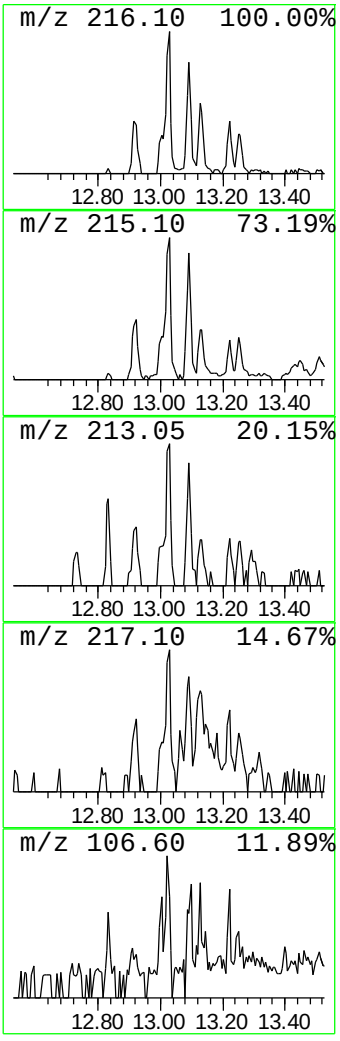
****

Peak Number 6 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	2.15 ng	50747	Chrysene-d12	13.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	86
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	64
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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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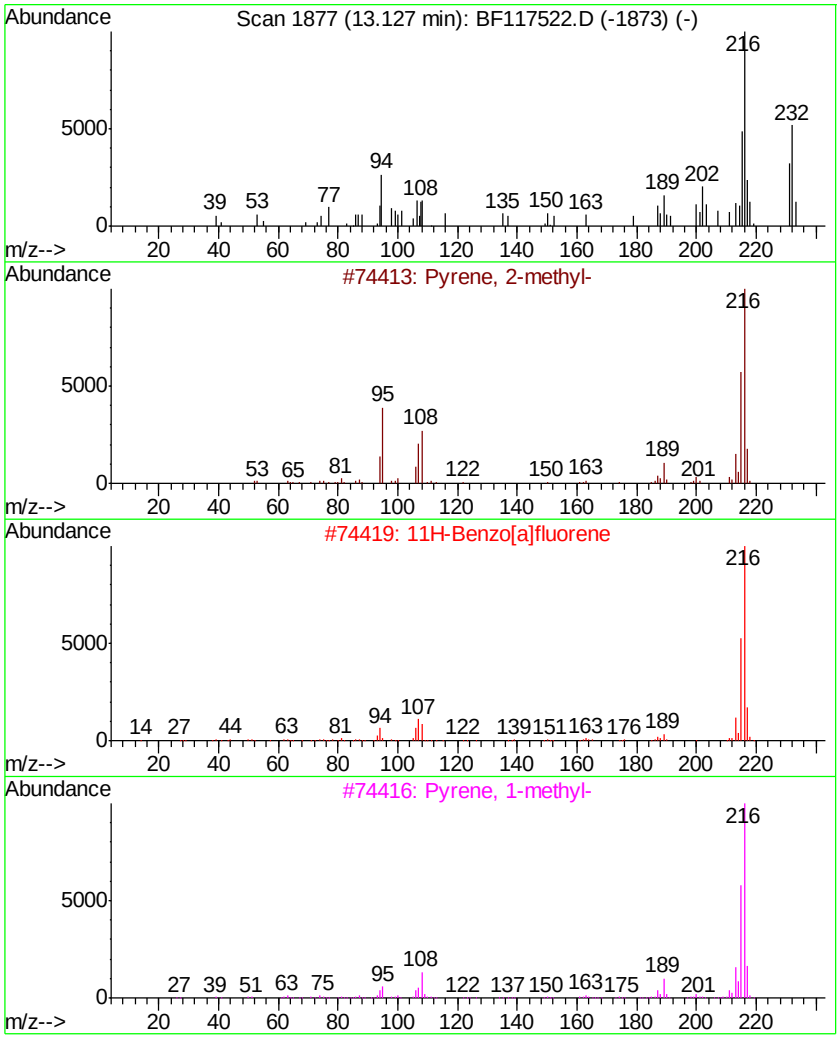
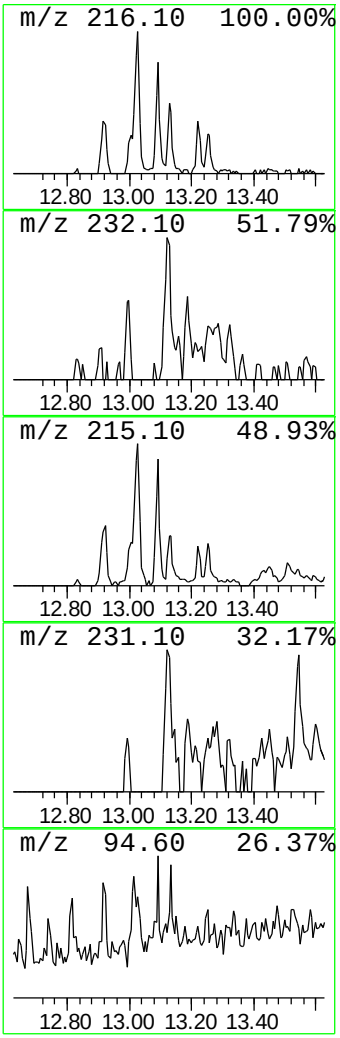
Peak Number 7 unknown13.13 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	2.24 ng	52731	Chrysene-d12	13.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 2-methyl-	216	C17H12	003442-78-2	46
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	46
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	43
4	Pyrene, 4-methyl-	216	C17H12	003353-12-6	38
5	4b,5,6,12-Tetrahydrochrysene	232	C18H16	031570-60-2	38





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Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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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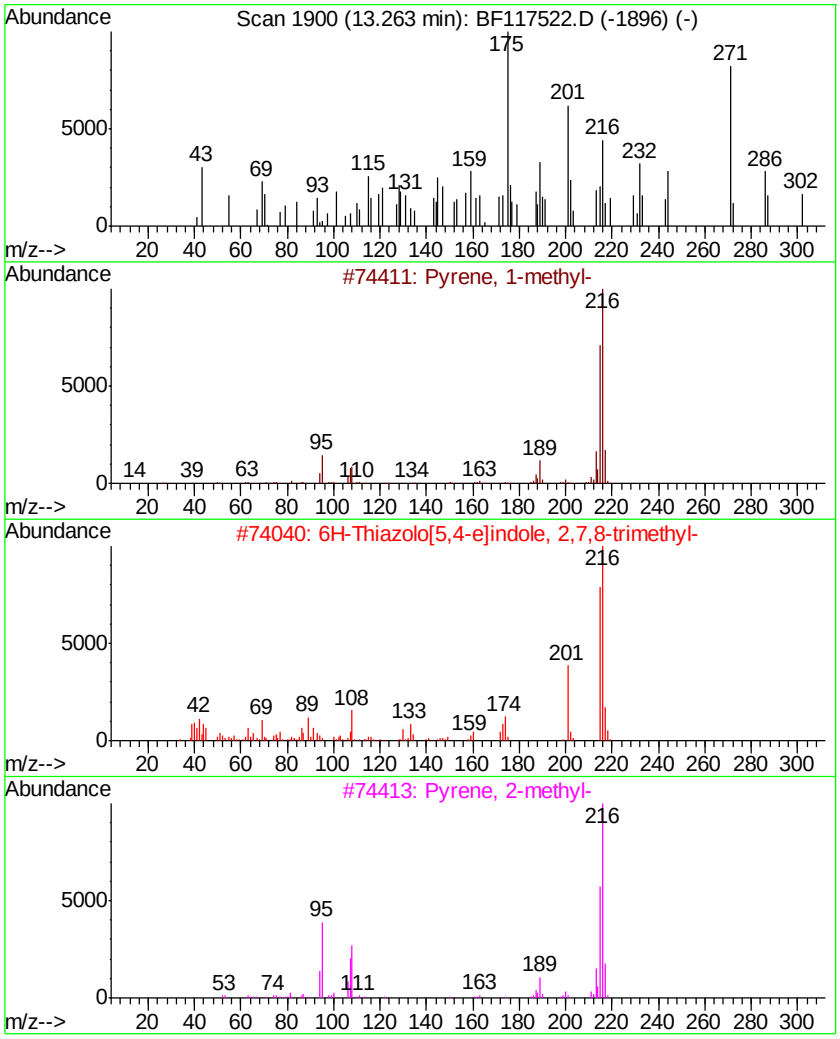
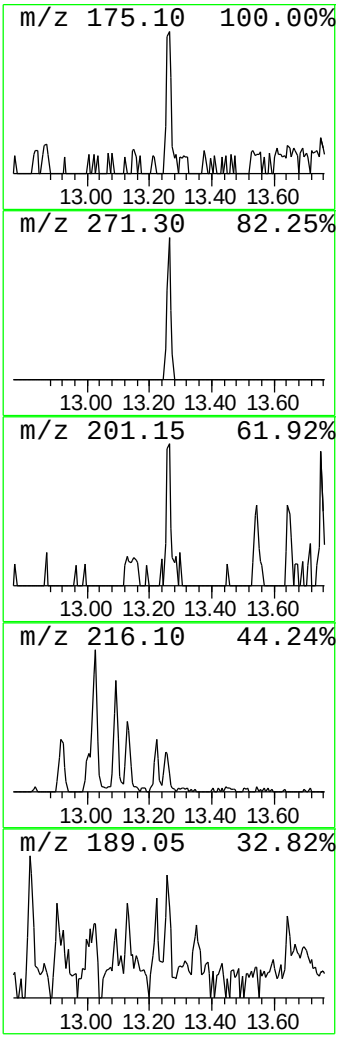
Peak Number 8 unknown13.26 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.26	2.06 ng	48607	Chrysene-d12	13.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	42
2	6H-Thiazolo[5,4-e]indole, 2,7,8-...	216	C12H12N2S	1000338-32-0	38
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	38
4	4-Hydroxy-6-methyl-1-pyridin-3-y...	216	C12H12N2O2	1000318-25-3	30
5	Cyclohexanone, 2-[(4-methoxyphen...	216	C14H16O2	005765-29-7	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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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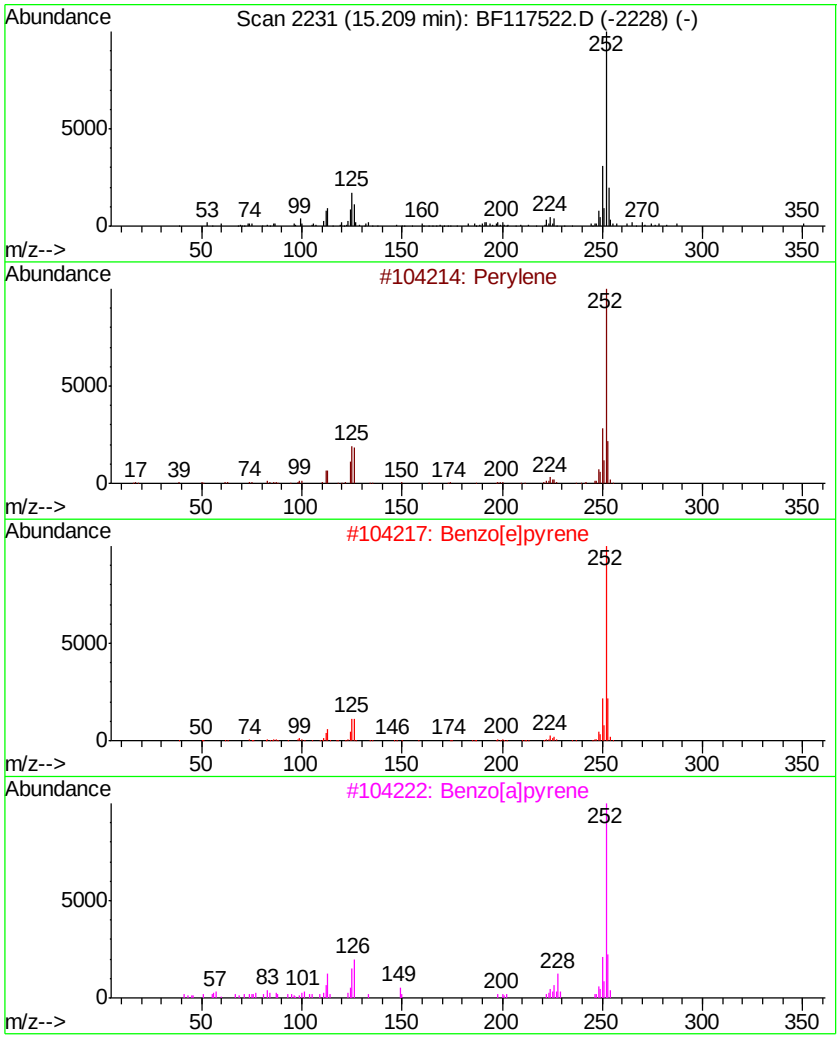
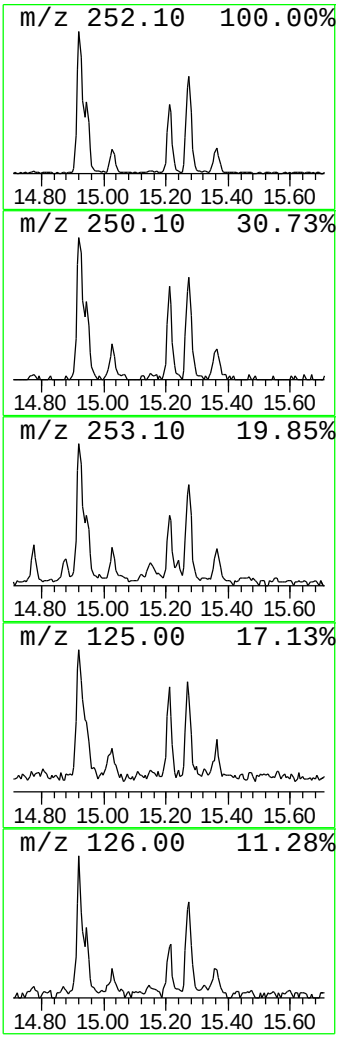
Peak Number 9 Perylene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	6.29 ng	83385	Perylene-d12	15.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perylene	252	C20H12	000198-55-0	94
2	Benzo[e]pyrene	252	C20H12	000192-97-2	93
3	Benzo[a]pyrene	252	C20H12	000050-32-8	91
4	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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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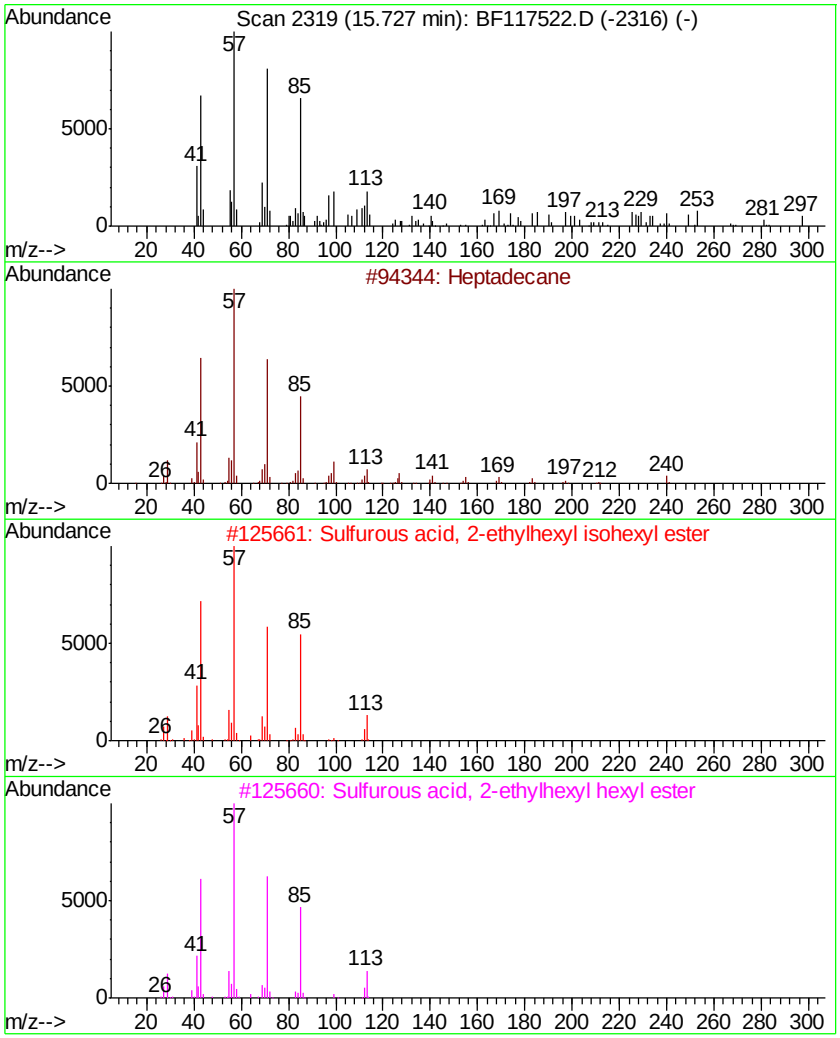
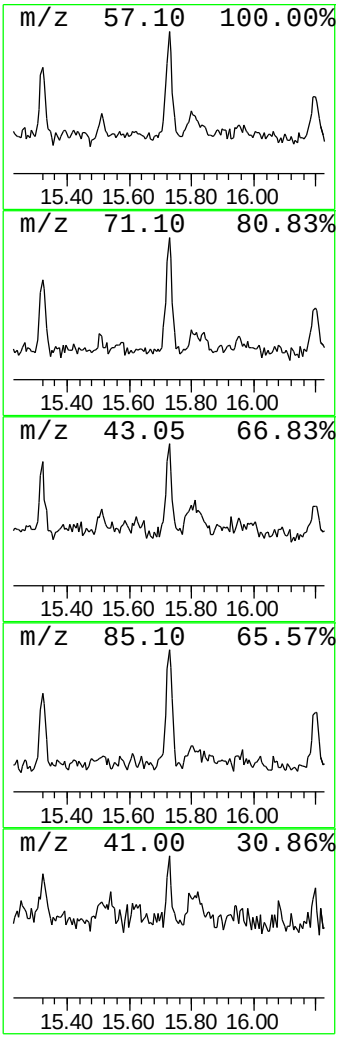
Peak Number 10 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	3.82 ng	50679	Perylene-d12	15.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	81
2	Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	64
3	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
4	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	64
5	Octadecane, 3-methyl-	268	C19H40	006561-44-0	64





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Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101819.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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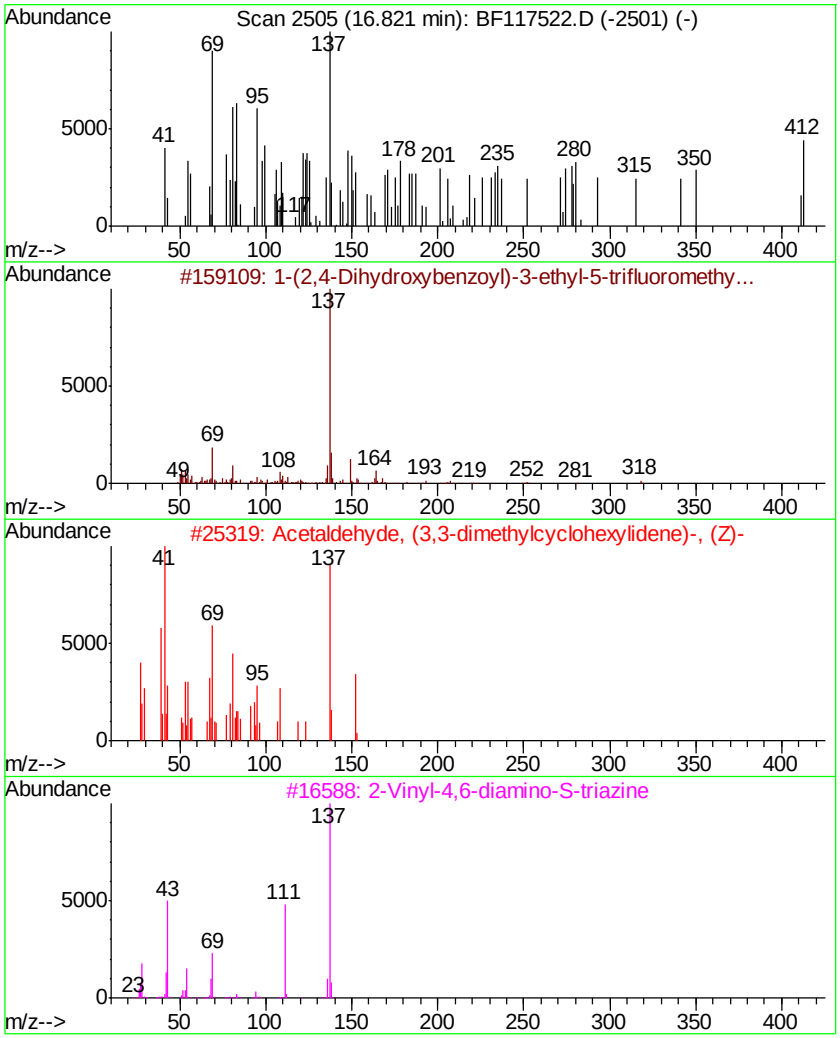
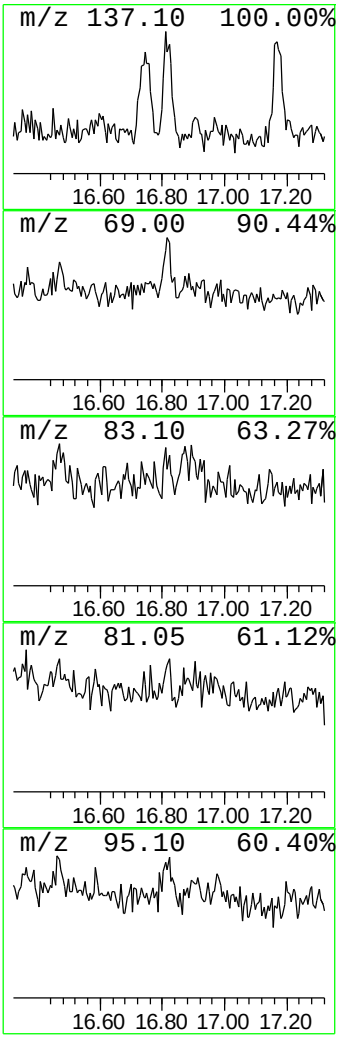
Peak Number 11 unknown16.82 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.82	3.14 ng	41559	Perylene-d12	15.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2,4-Dihydroxybenzoyl)-3-ethyl...	318	C13H13F3N2O4	331835-05-3	27
2	Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-24-1	25
3	2-Vinyl-4,6-diamino-S-triazine	137	C5H7N5	003194-70-5	22
4	Tricyclo[3.2.1.0(2,4)]octane-3,3...	280	C18H20N2O	1000164-56-3	18
5	Pyrazine, 2-methoxy-3-(1-methyle...	152	C8H12N2O	025773-40-4	15





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF102419\
Data File : BF117522.D
Acq On : 24 Oct 2019 20:39
Operator : JU
Sample : K5499-09 2X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : 197-74-(0-1)

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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
unknown6.51	6.51	30.6	ng	472468	1	6.73	308603	20.0
Phenanthrene, 2-m...	11.75	2.2	ng	43831	4	11.25	394359	20.0
Anthracene, 1-met...	11.78	5.4	ng	106184	4	11.25	394359	20.0
4H-Cyclopenta[def...	11.86	3.9	ng	76320	4	11.25	394359	20.0
Pyrene, 1-methyl-	13.03	2.1	ng	50747	5	13.90	471223	20.0
unknown13.13	13.13	2.2	ng	52731	5	13.90	471223	20.0
unknown13.26	13.26	2.1	ng	48607	5	13.90	471223	20.0
Perylene	15.21	6.3	ng	83385	6	15.33	265063	20.0
Heptadecane	15.73	3.8	ng	50679	6	15.33	265063	20.0
unknown16.82	16.82	3.1	ng	41559	6	15.33	265063	20.0