LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070918\
Data File : BF107239.D
Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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	5.184	480	484	497	rVB	73299	82713	10.09%	0.968%
2	5.625	555	559	585	rBV	313938	650137	79.30%	7.609%
3	6.619	725	728	745	rBV	330359	610789	74.50%	7.148%
4	6.737	745	748	754	rVV	525295	630468	76.90%	7.379%
5	6.784	754	756	761	rVB	10208	11974	1.46%	0.140%
6	6.948	780	784	787	rBV	216747	182658	22.28%	2.138%
7	7.101	806	810	814	rBV	541772	479561	58.49%	5.612%
8	7.513	876	880	890	rBV	374251	348660	42.53%	4.080%
9	8.231	998	1002	1005	rBV	213394	190709	23.26%	2.232%
10	8.795	1095	1098	1102	rVB2	8587	8824	1.08%	0.103%
11	8.948	1121	1124	1129	rBB	13836	12242	1.49%	0.143%
12	9.301	1180	1184	1188	rBV	710872	616594	75.21%	7.216%
13	9.695	1245	1251	1257	rVB	144239	128741	15.70%	1.507%
14	9.783	1261	1266	1270	rBV2	10342	13205	1.61%	0.155%
15	9.854	1275	1278	1281	rVV	17862	16316	1.99%	0.191%
16	9.895	1281	1285	1289	rVB	10519	9648	1.18%	0.113%
17	9.989	1298	1301	1305	rVV	211572	195041	23.79%	2.283%
18	10.025	1305	1307	1311	rVB	22940	16767	2.05%	0.196%
19	10.201	1333	1337	1339	rBV2	12247	13617	1.66%	0.159%
20	10.395	1366	1370	1372	rBV2	16614	12938	1.58%	0.151%
21	10.542	1392	1395	1401	rVB2	16130	18325	2.24%	0.214%
22	10.613	1401	1407	1412	rBV5	8607	12818	1.56%	0.150%
23	10.789	1433	1437	1443	rBV	430780	407821	49.74%	4.773%
24	10.872	1448	1451	1455	rVV2	14540	21288	2.60%	0.249%
25	11.295	1520	1523	1525	rBV2	11876	11150	1.36%	0.130%
26	11.319	1525	1527	1530	rVV3	11360	10344	1.26%	0.121%
27	11.489	1552	1556	1558	rBV	286568	243748	29.73%	2.853%
28	11.513	1558	1560	1564	rVB	167836	134624	16.42%	1.576%
29	11.566	1566	1569	1573	rBV	55857	48044	5.86%	0.562%
30	11.725	1594	1596	1598	rBV	14261	12189	1.49%	0.143%
31	11.748	1598	1600	1603	rVB2	14353	10691	1.30%	0.125%
32	11.824	1611	1613	1619	rVB6	8670	10755	1.31%	0.126%
33	11.989	1638	1641	1644	rBV	39263	42180	5.14%	0.494%
34	12.019	1644	1646	1651	rVV2	38982	41003	5.00%	0.480%

LSC Area Percent Report

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Sample : J3879-11
Misc :
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Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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	12.072	1652	1655	1657	rVV	18668	16308	1.99%	0.191%
36	12.107	1657	1661	1663	rVV2	47431	54197	6.61%	0.634%
37	12.130	1663	1665	1667	rVV	23457	19136	2.33%	0.224%
38	12.154	1667	1669	1672	rVB	13682	11777	1.44%	0.138%
39	12.283	1688	1691	1695	rBV	27715	25091	3.06%	0.294%
40	12.324	1695	1698	1702	rVB2	22132	21066	2.57%	0.247%
41	12.436	1714	1717	1720	rVB	11033	9221	1.12%	0.108%
42	12.466	1720	1722	1724	rVB	11931	10059	1.23%	0.118%
43	12.495	1724	1727	1729	rBV2	13174	12030	1.47%	0.141%
44	12.542	1732	1735	1738	rBV	46672	44254	5.40%	0.518%
45	12.642	1749	1752	1754	rBV2	21210	17888	2.18%	0.209%
46	12.713	1760	1764	1767	rBV	382003	328232	40.04%	3.841%
47	12.807	1778	1780	1783	rBV2	11700	12053	1.47%	0.141%
48	12.924	1796	1800	1801	rBV2	76954	79958	9.75%	0.936%
49	12.948	1801	1804	1808	rVV	325233	322997	39.40%	3.780%
50	12.989	1809	1811	1815	rVB2	27811	28788	3.51%	0.337%
51	13.077	1821	1826	1829	rBV	825155	819856	100.00%	9.595%
52	13.113	1829	1832	1835	rVB	20833	22392	2.73%	0.262%
53	13.166	1836	1841	1845	rBV4	31291	55953	6.82%	0.655%
54	13.271	1852	1859	1862	rBV	67762	99646	12.15%	1.166%
55	13.336	1868	1870	1872	rBV	28738	21825	2.66%	0.255%
56	13.471	1891	1893	1895	rBV	32226	26619	3.25%	0.312%
57	13.501	1896	1898	1901	rVV	25404	23505	2.87%	0.275%
58	13.618	1915	1918	1921	rBV2	34618	34054	4.15%	0.399%
59	13.789	1944	1947	1950	rBV	34616	32243	3.93%	0.377%
60	13.895	1963	1965	1967	rBV	31568	24655	3.01%	0.289%
61	13.918	1967	1969	1972	rVV	30862	28574	3.49%	0.334%
62	13.948	1972	1974	1980	rVB3	69505	82708	10.09%	0.968%
63	14.148	2003	2008	2011	rBV2	273849	395587	48.25%	4.630%
64	14.177	2011	2013	2018	rVB	138820	118965	14.51%	1.392%
65	14.260	2025	2027	2032	rVB2	41786	42262	5.15%	0.495%
66	14.571	2078	2080	2084	rBV	45543	48780	5.95%	0.571%
67	15.236	2190	2193	2201	rVB	110283	220782	26.93%	2.584%
68	15.630	2257	2260	2264	rBV	61814	75954	9.26%	0.889%
69	15.695	2268	2271	2274	rBV	93410	100629	12.27%	1.178%

Sum of corrected areas: 8544626

LSC Report - Integrated Chromatogram

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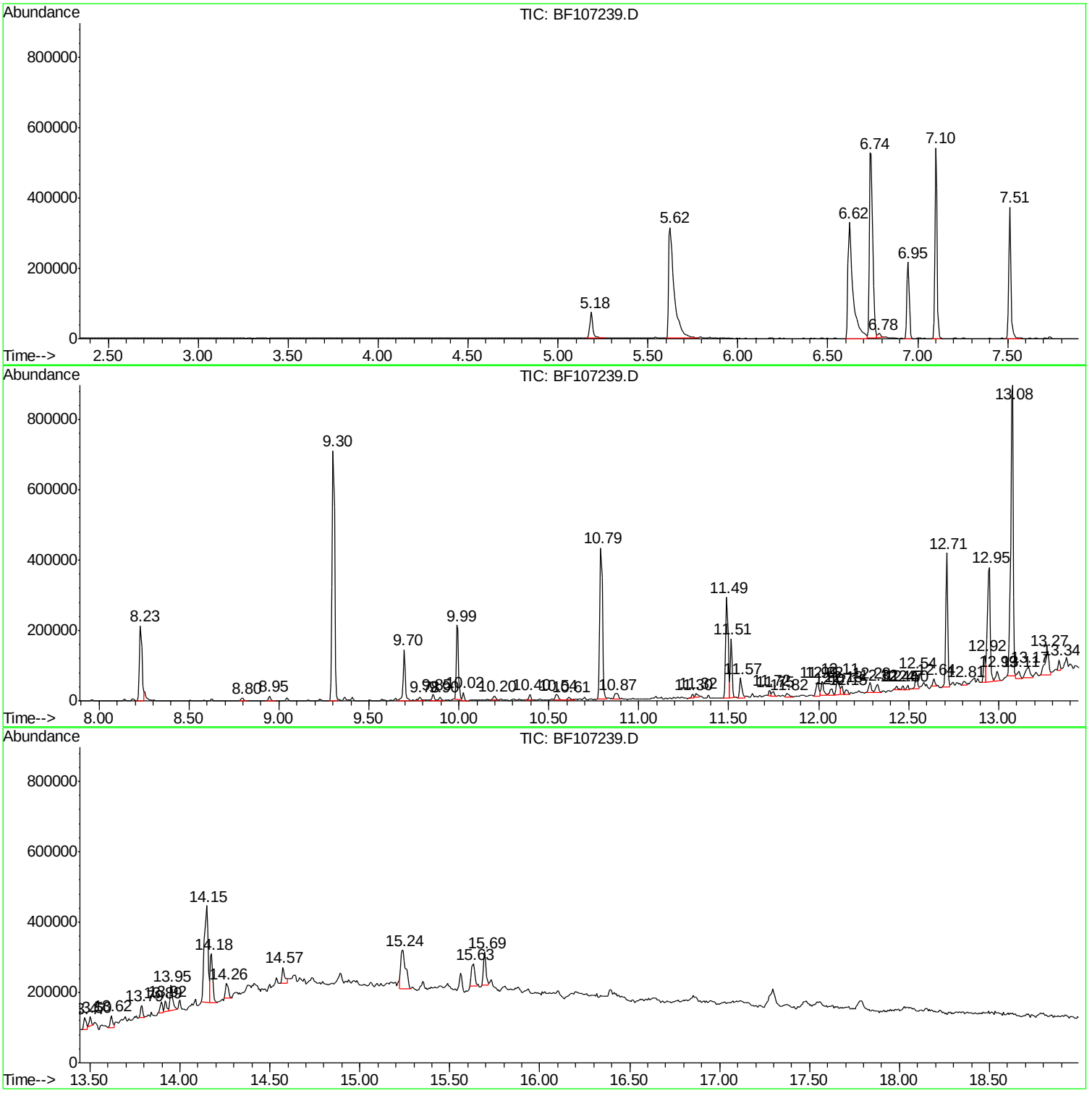
Data File : BF107239.D
Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SURCHARGE-SP-6

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



Library Search Compound Report

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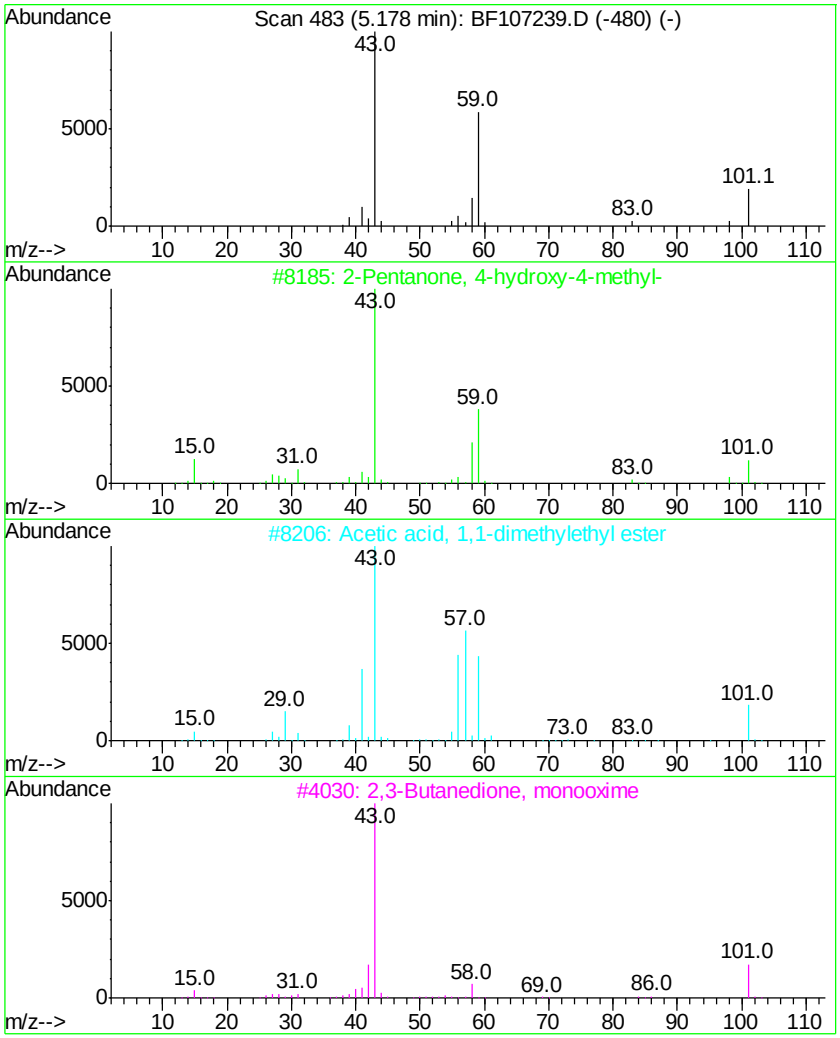
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

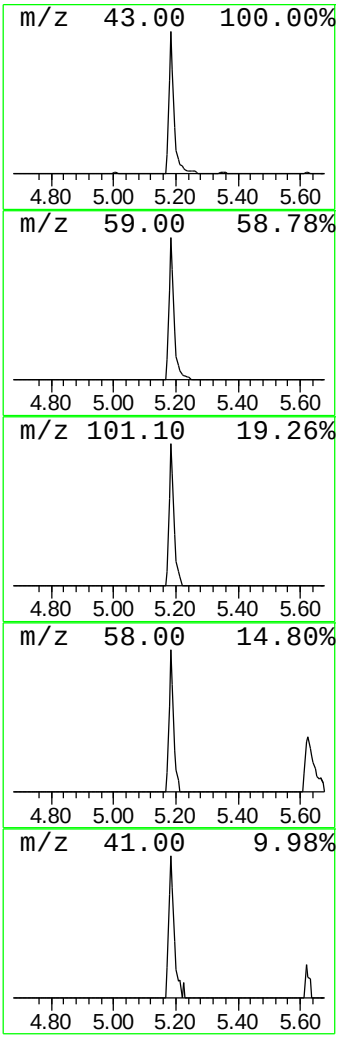
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	9.06 ng	82713	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	56
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Library Search Compound Report

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Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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

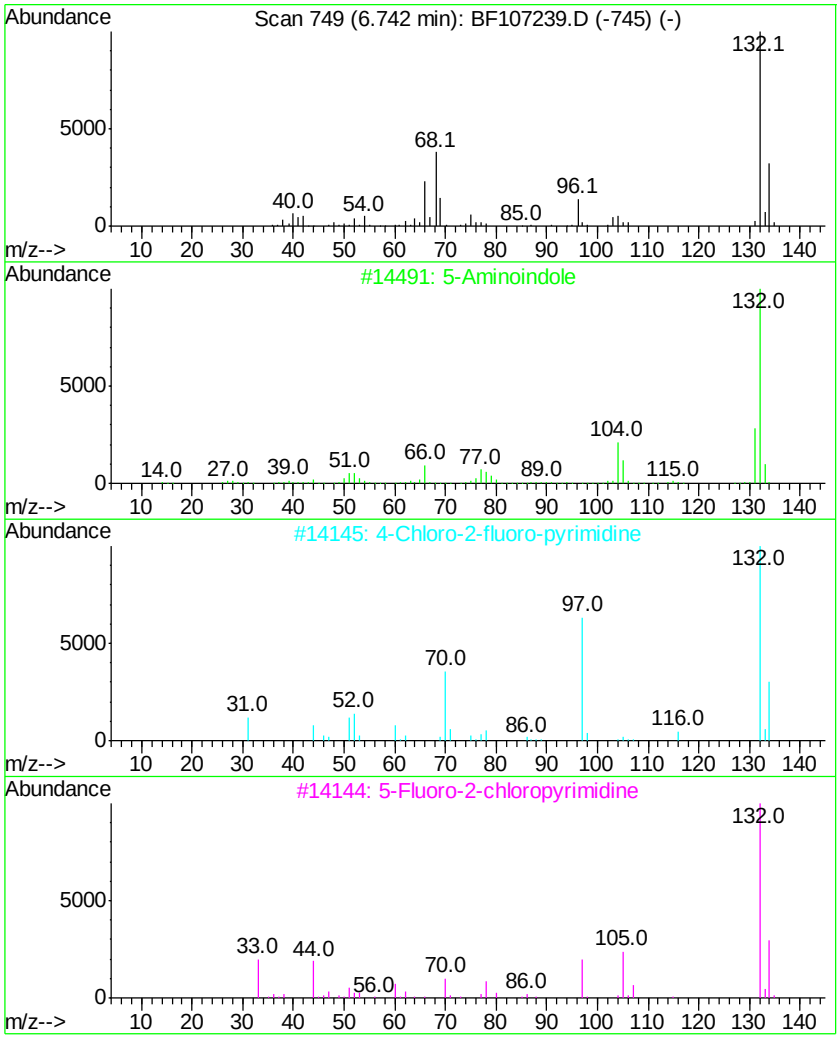
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

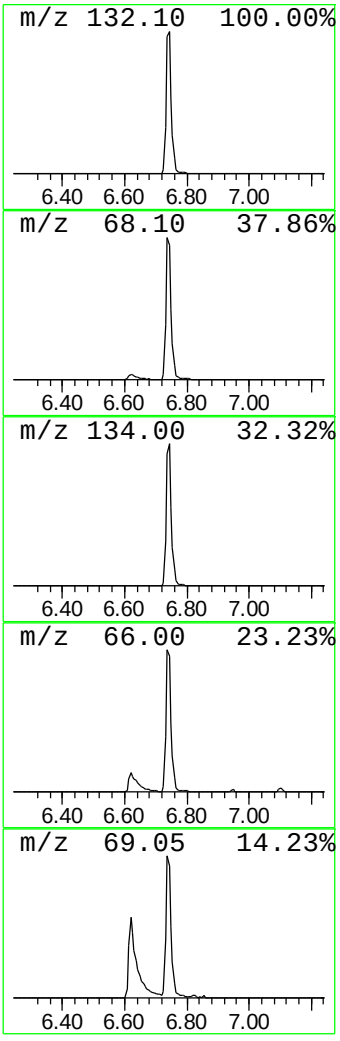
****

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	69.03 ng	630468	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	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9





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Sample : J3879-11
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Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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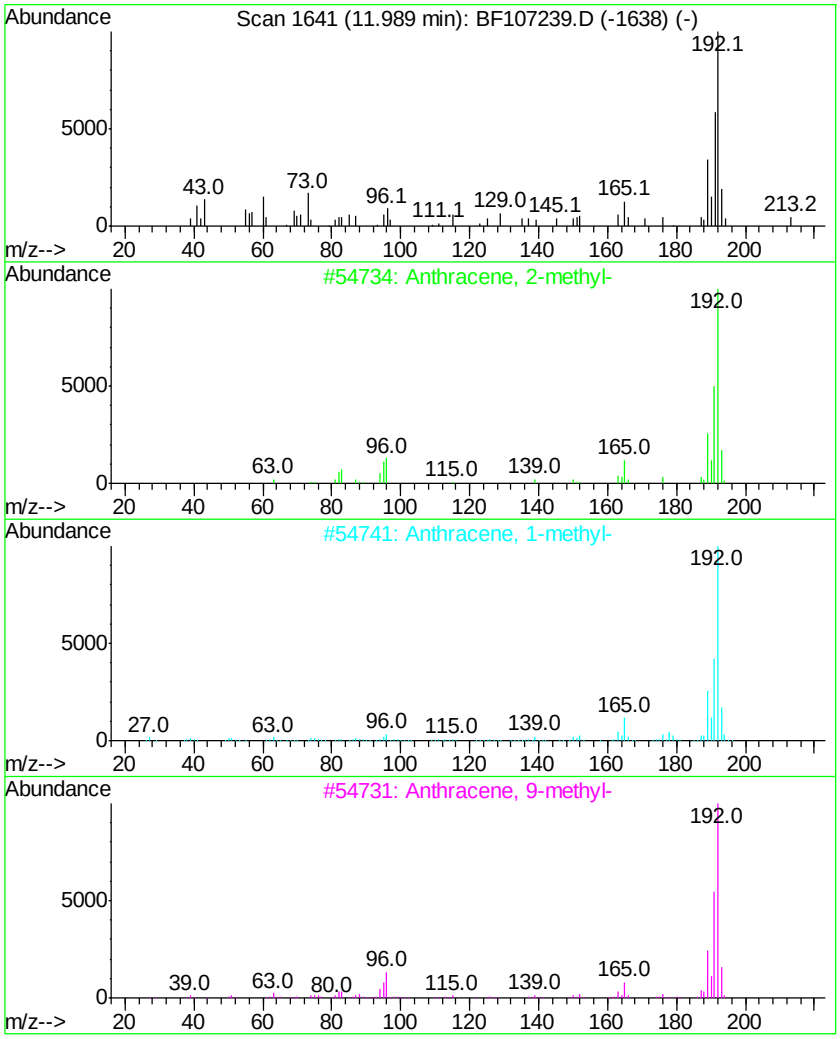
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

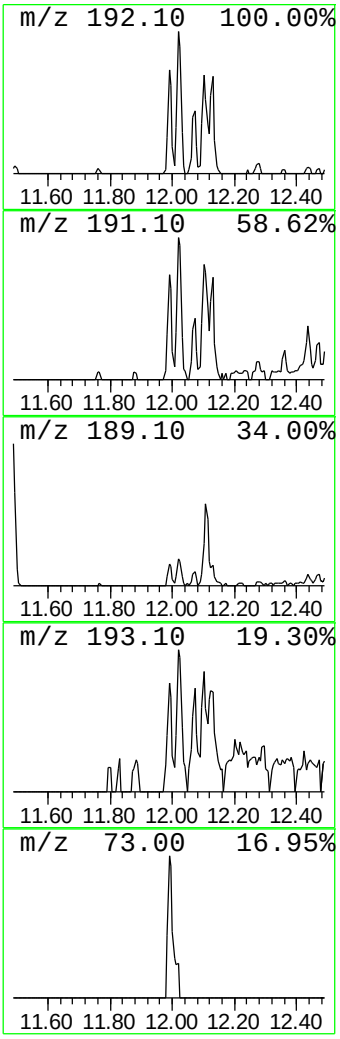
****

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	3.46 ng	42180	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	89
4	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	86
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81





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Instrument : BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
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TIC Library : C:\DATABASE\NIST11.L
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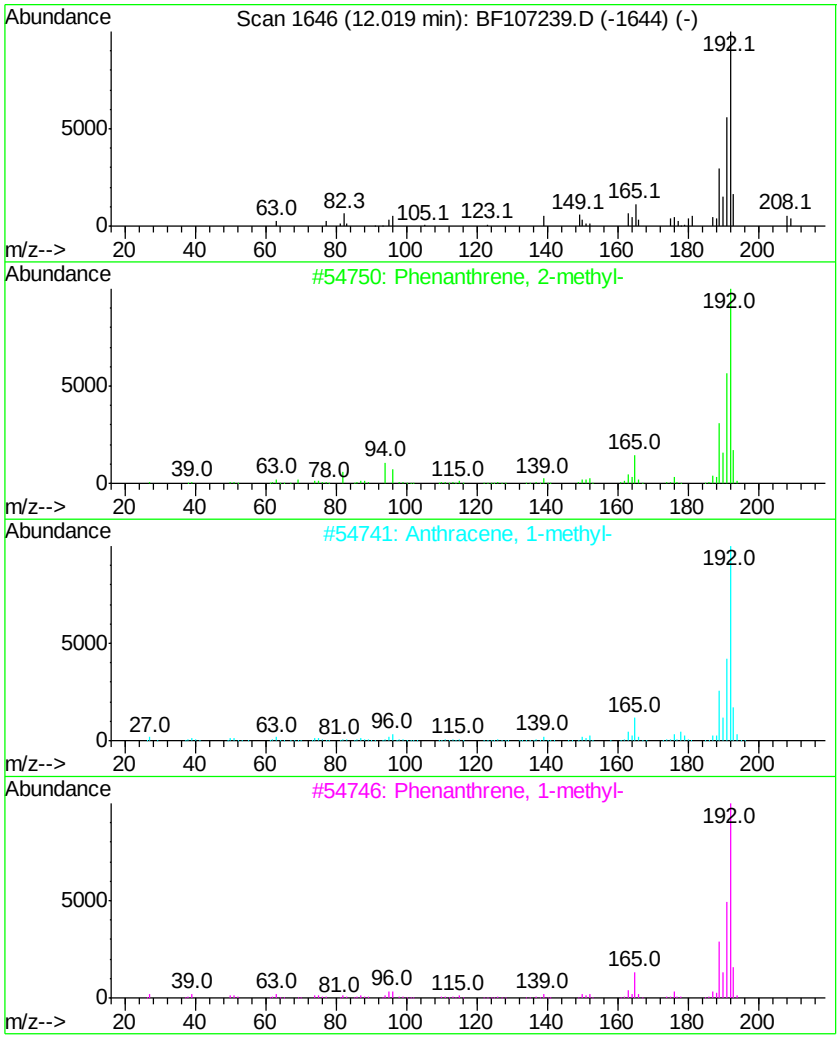
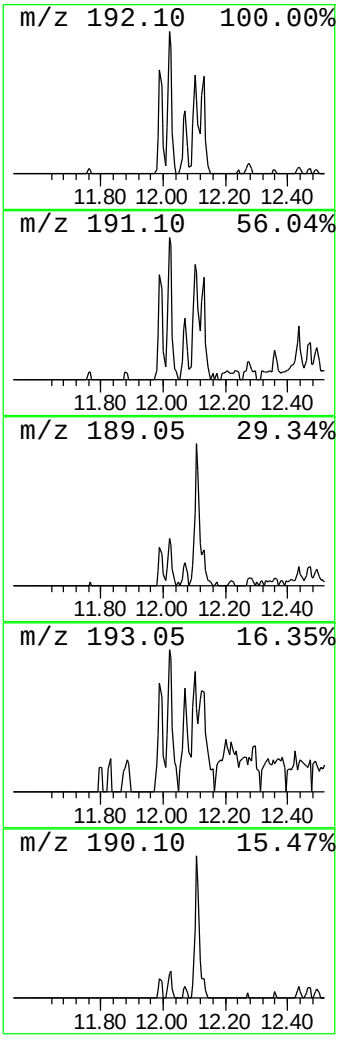
****

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	3.36 ng	41003	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	89





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Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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

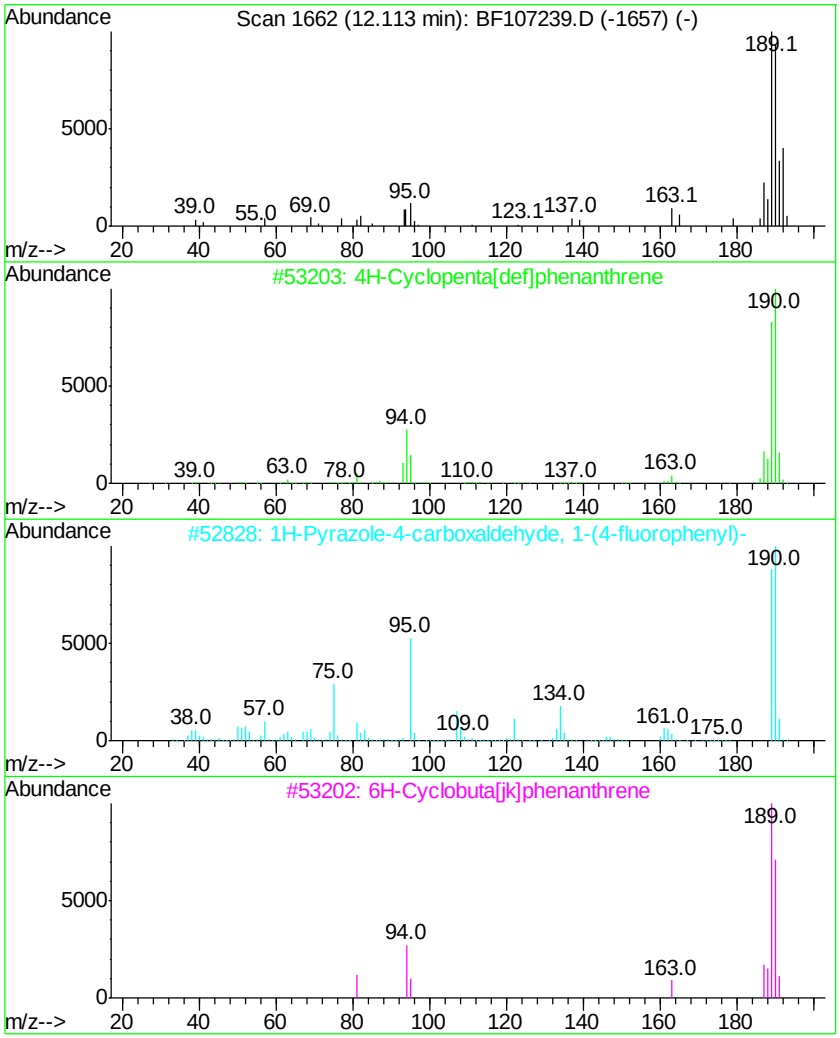
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

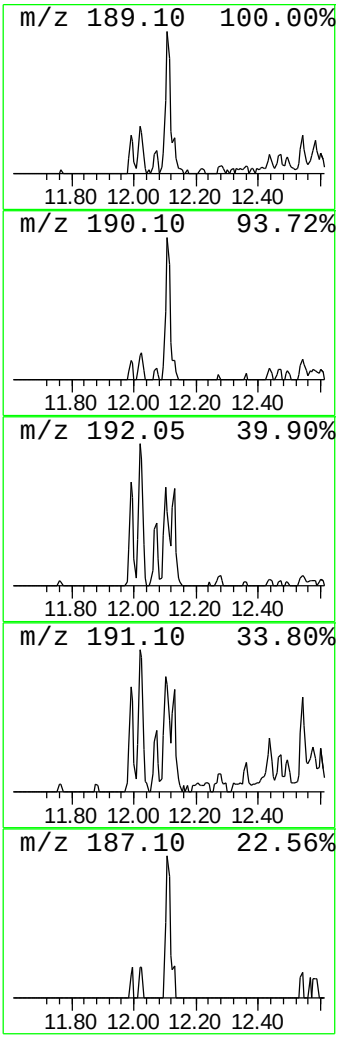
****

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	4.45 ng	54197	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	1H-Pyrazole-4-carboxaldehyde, 1-...	190	C10H7FN2O	1000338-05-6	43
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42
4	5-Thiazolecarboxamide, 4-amino-2...	189	C5H7N3OS2	1000338-08-9	22
5	Benzene, 1,1'-(1,2-propadienylid...	192	C15H12	014251-57-1	22





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Misc :
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Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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

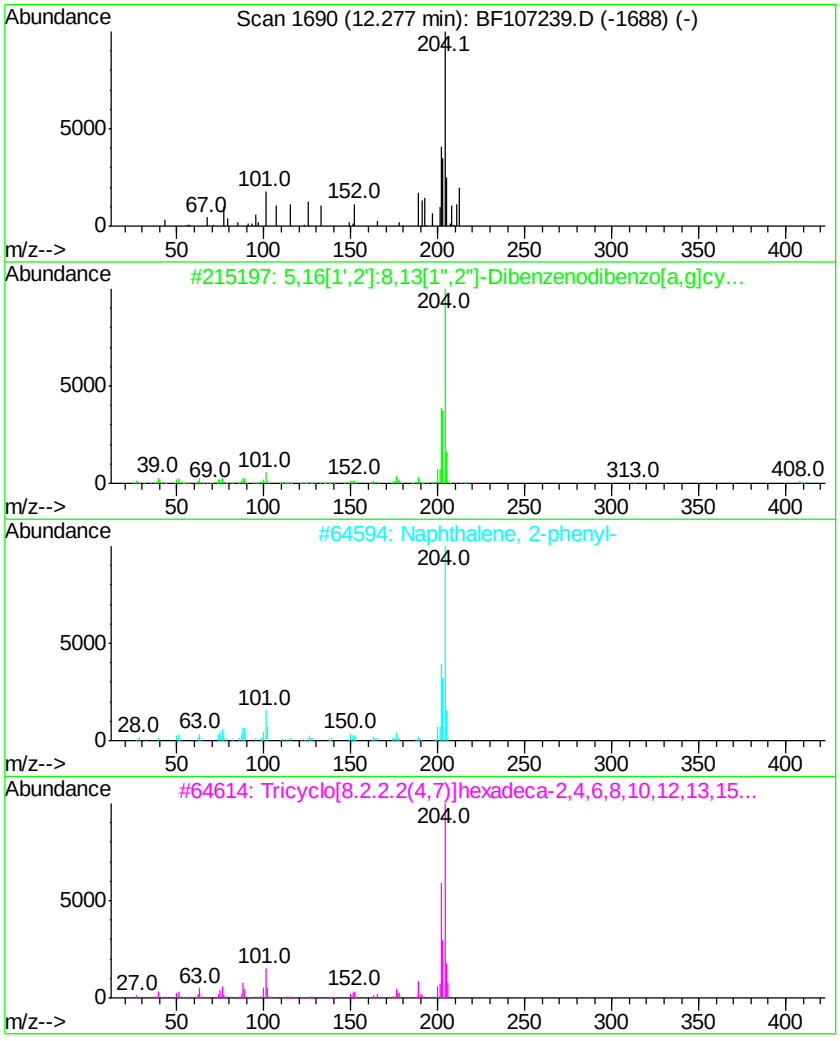
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

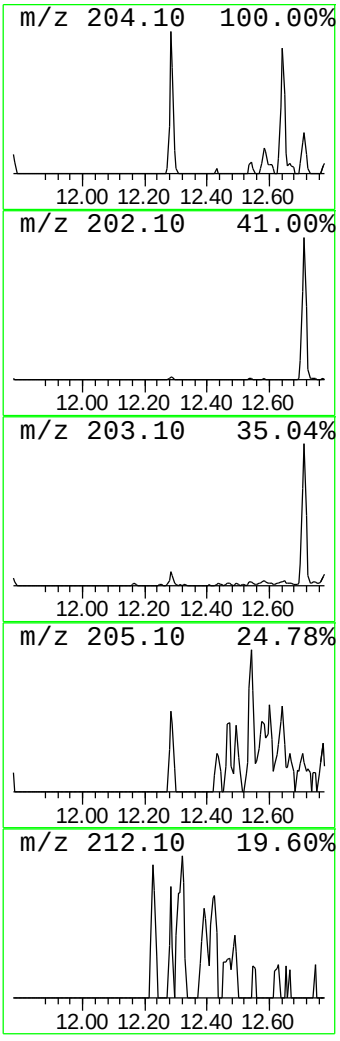
****

Peak Number 7 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	2.06 ng	25091	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	72
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
4	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	52
5	3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	49





Library Search Compound Report

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Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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

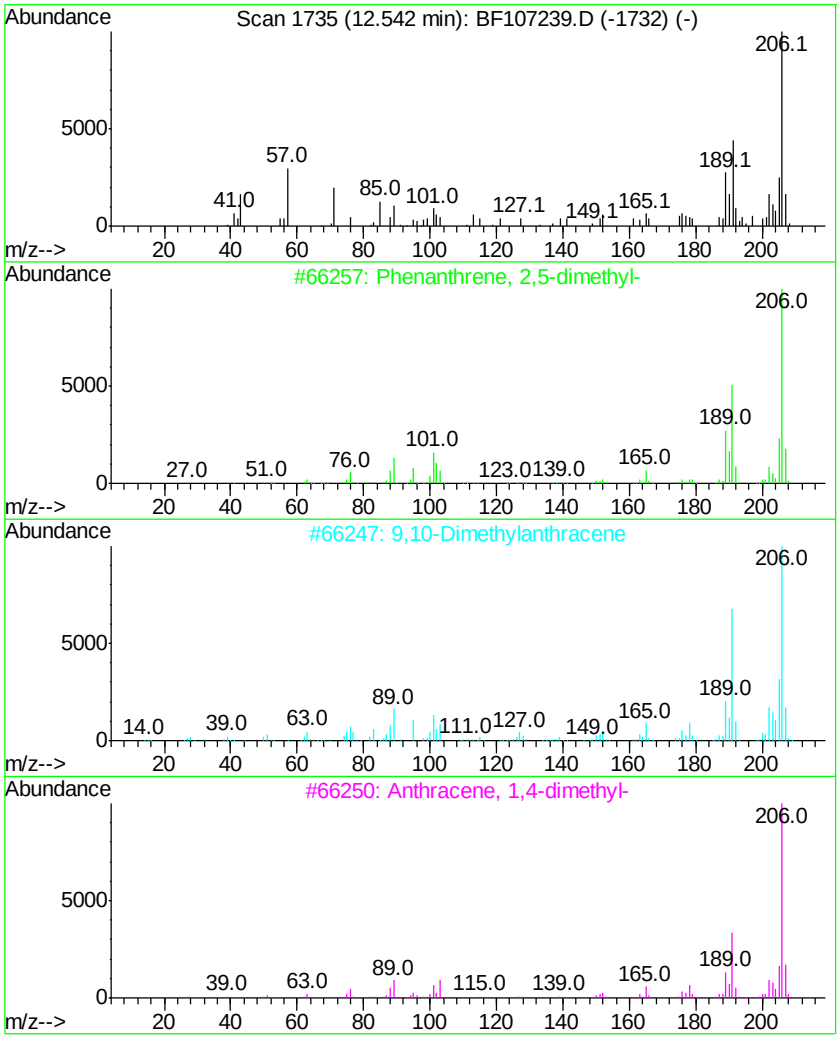
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

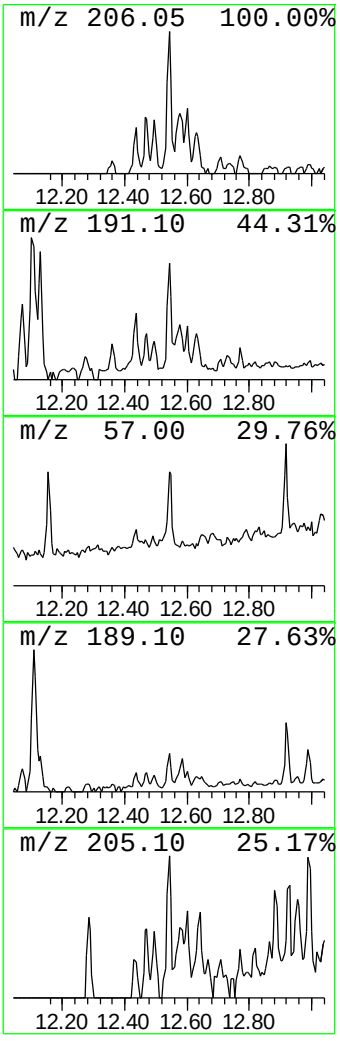
****

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	3.63 ng	44254	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
3	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	90
4	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90





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Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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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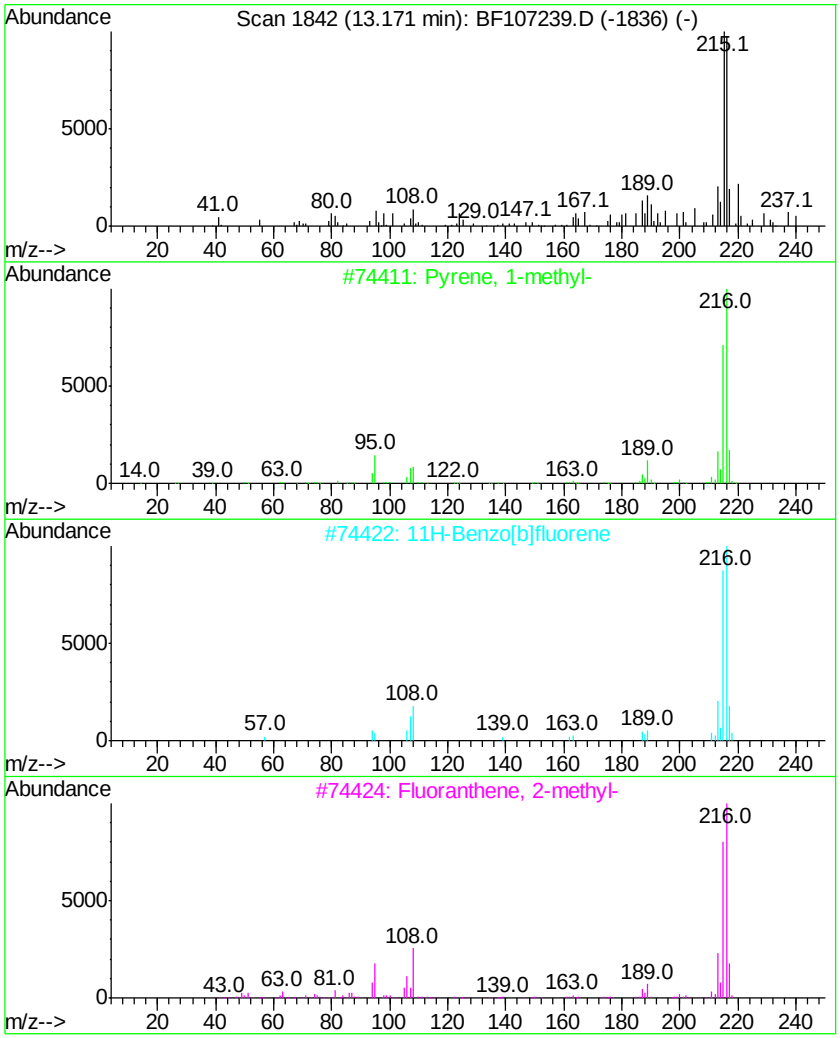
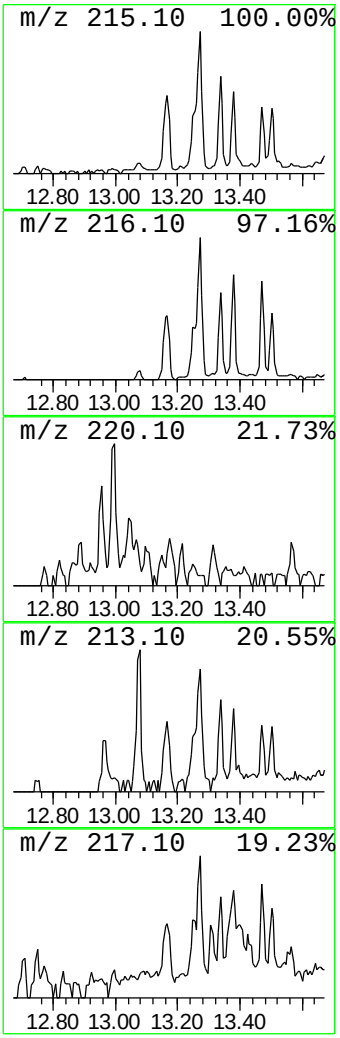
Peak Number 9 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.17	2.83 ng	55953	Chrysene-d12	14.15

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	89
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	76
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5	Benzoic acid, 3-(3,5-dimethyl-1-...	216	C12H12N2O2	312531-88-7	58





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Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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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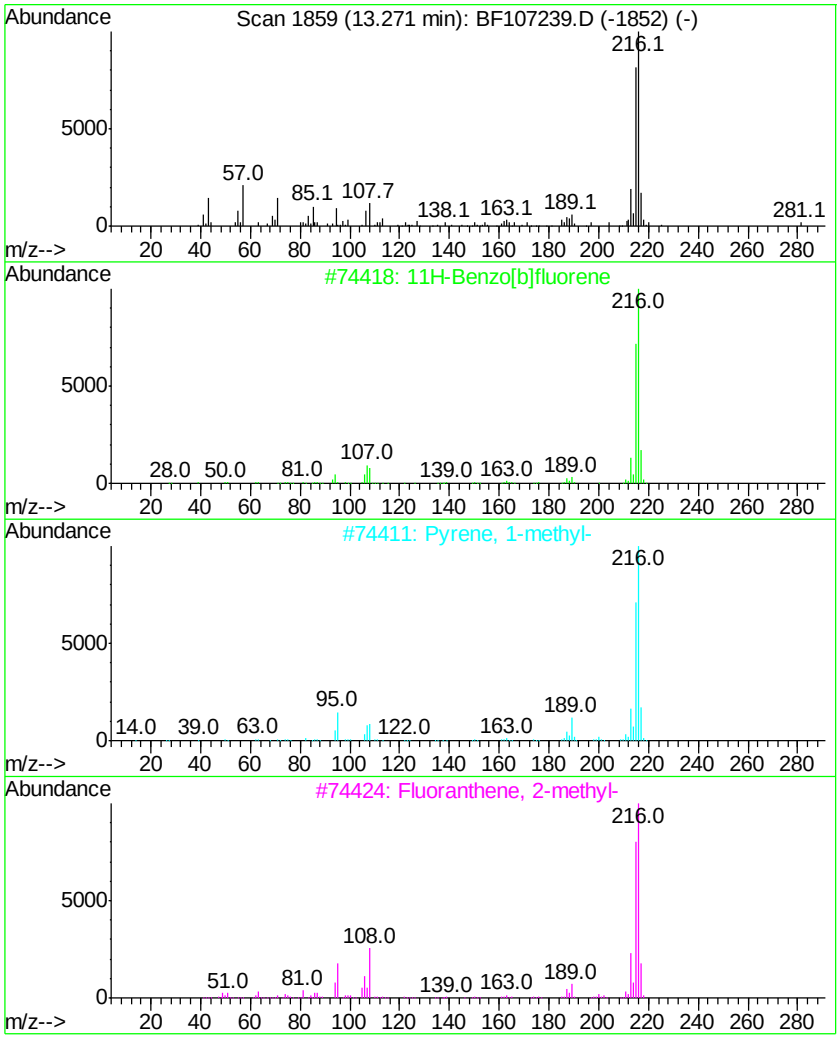
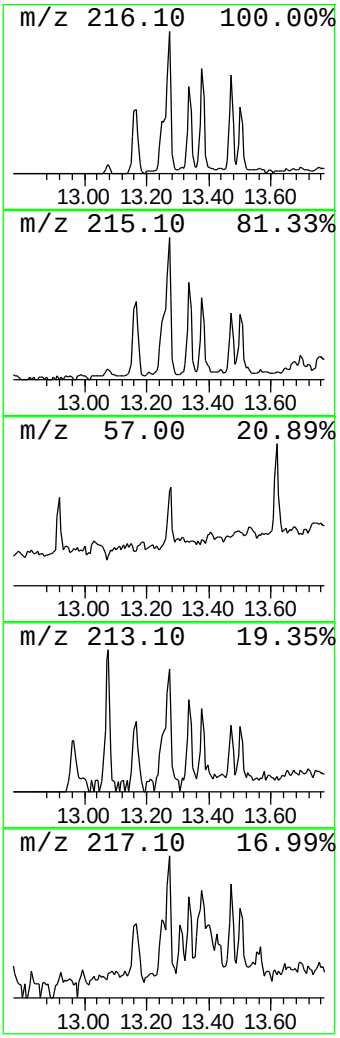
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	5.04 ng	99646	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	64





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Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
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TIC Integration Parameters: LSCINT.P

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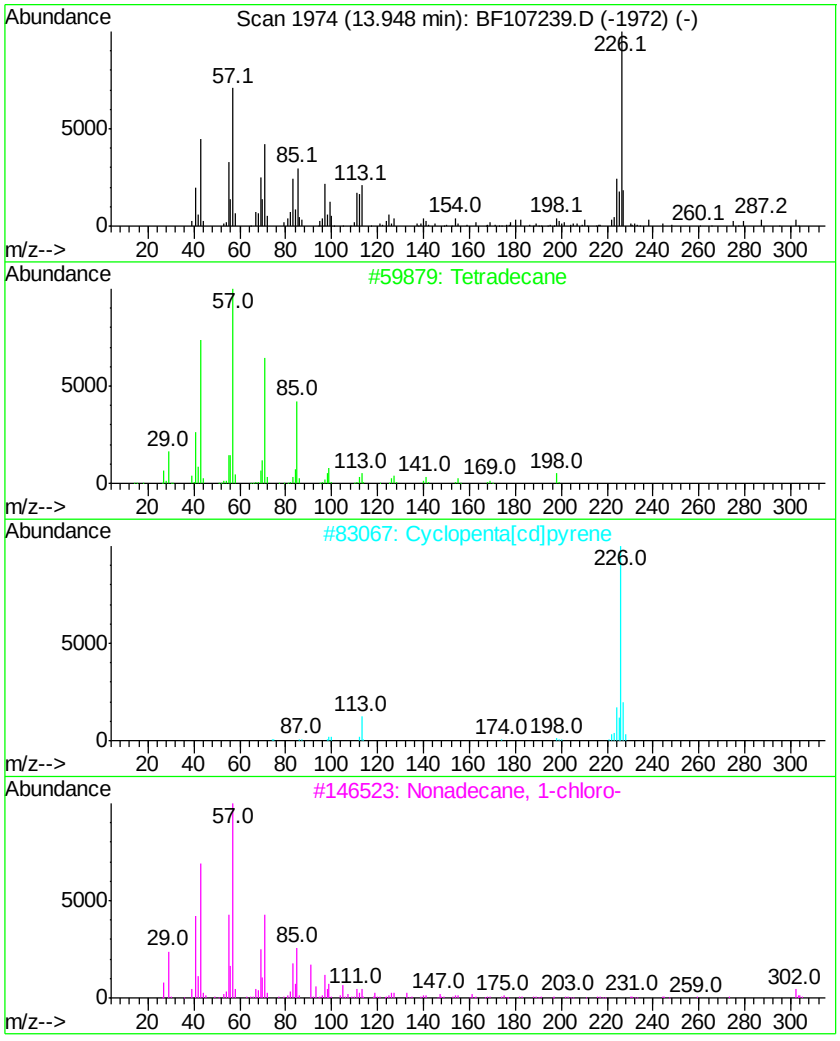
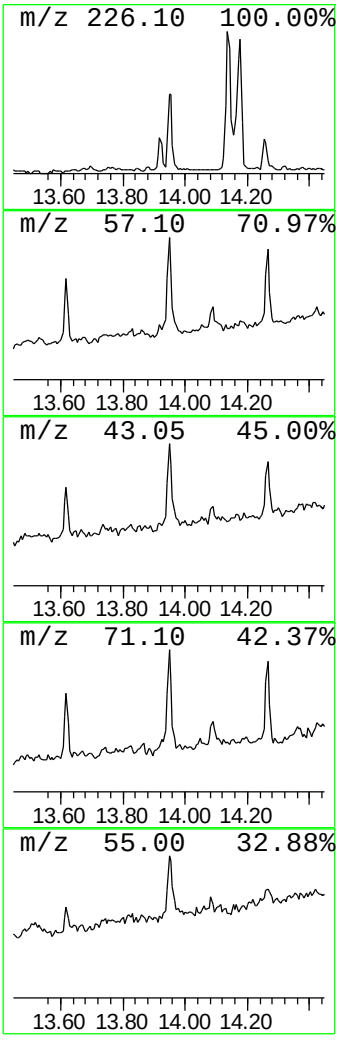
Peak Number 11 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	4.18 ng	82708	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	64
2	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	47
3	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	43
4	Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	35
5	Oxalic acid, isobutyl tridecyl e...	328	C19H36O4	1000309-37-8	30





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Data File : BF107239.D
Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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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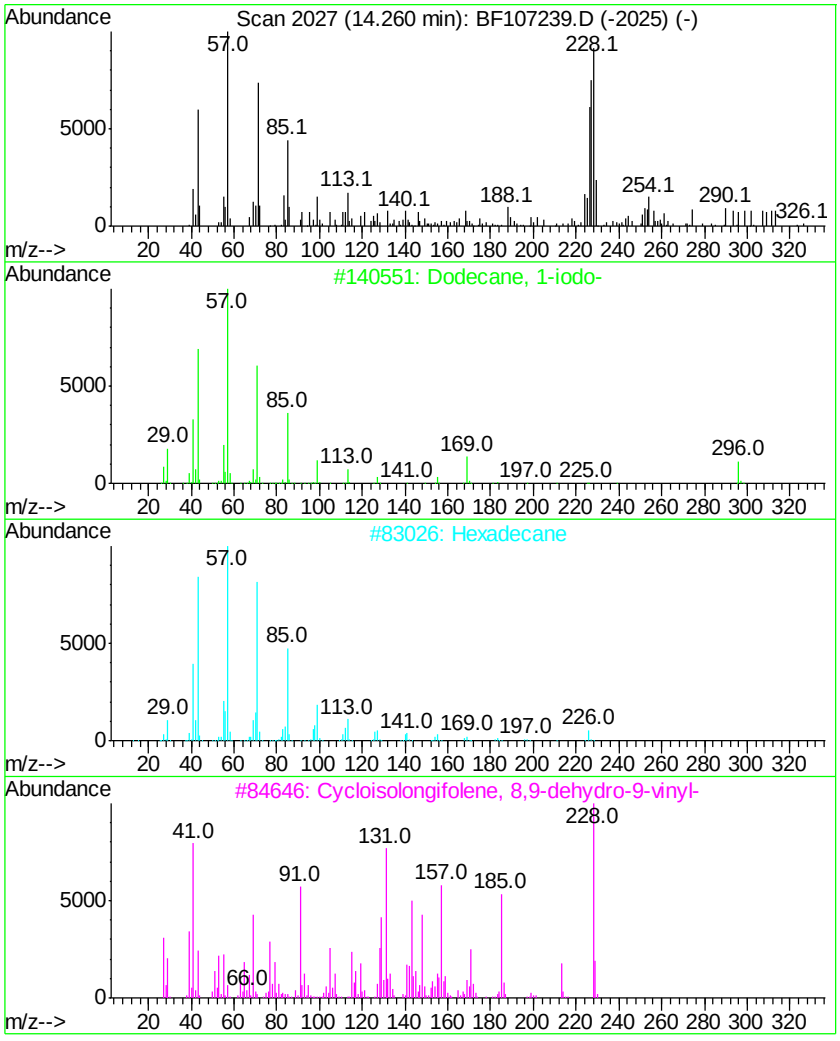
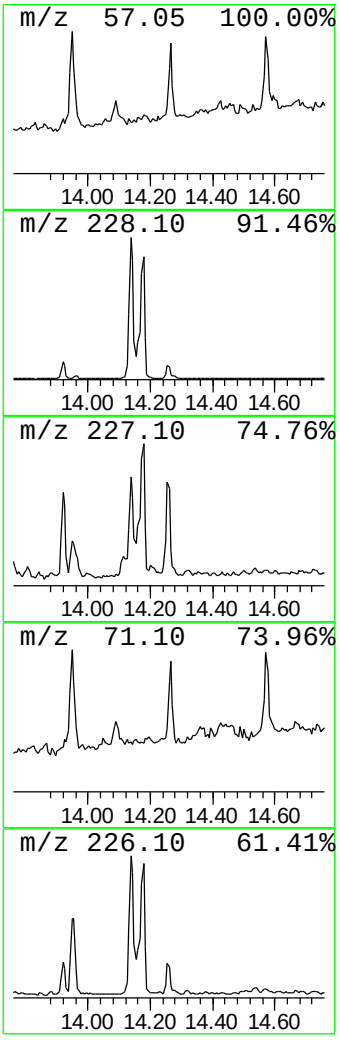
Peak Number 12 unknown14.26 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	2.14 ng	42262	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	43
2	Hexadecane	226	C16H34	000544-76-3	27
3	Cycloisolongifolene, 8,9-dehydro...	228	C17H24	1000151-80-0	25
4	Methoxyacetic acid, 4-hexadecyl ...	314	C19H38O3	1000282-99-0	25
5	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	22





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Data File : BF107239.D
Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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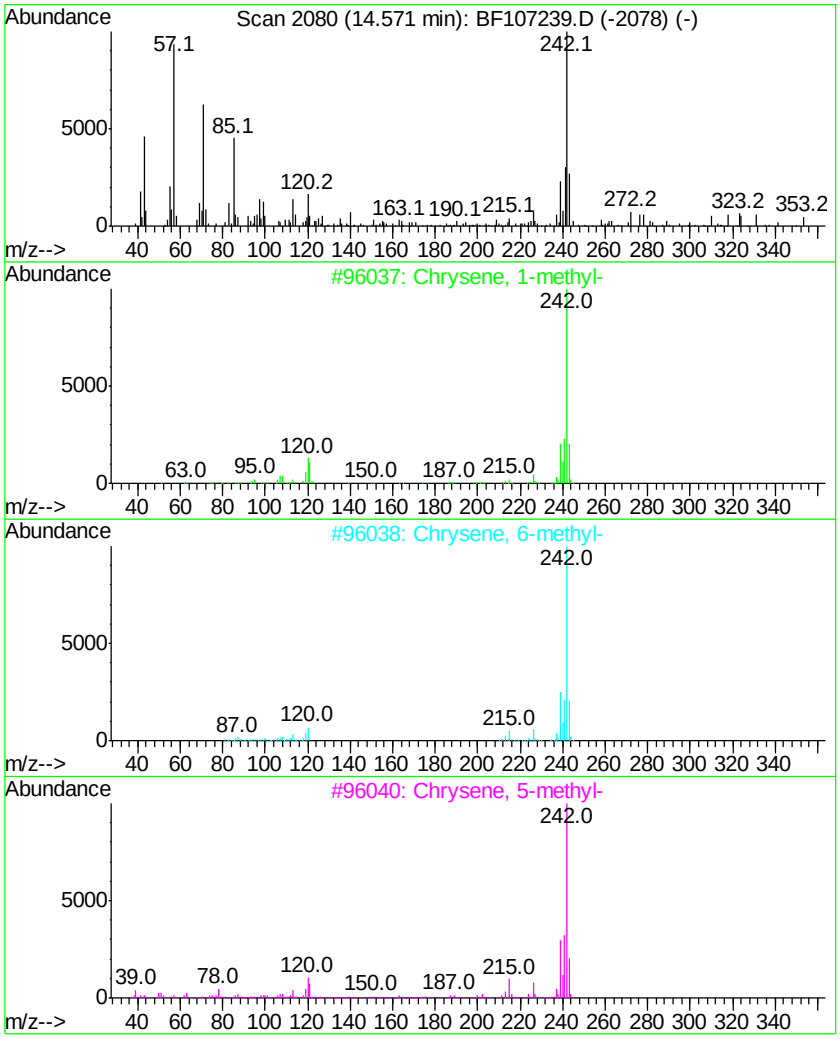
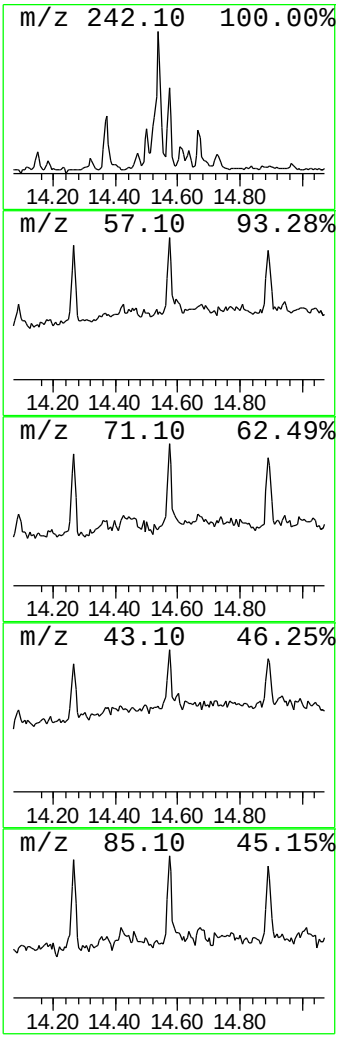
Peak Number 13 Chrysene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	2.47 ng	48780	Chrysene-d12	14.15

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Chrysene, 1-methyl-	242	C19H14	003351-28-8	68
2	Chrysene, 6-methyl-	242	C19H14	001705-85-7	49
3	Chrysene, 5-methyl-	242	C19H14	003697-24-3	49
4	Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	45
5	Benzo[c]phenanthrene, 5-methyl-	242	C19H14	000652-04-0	41





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF070918\
Data File : BF107239.D
Acq On : 9 Jul 2018 19:55
Operator : JU/SJ
Sample : J3879-11
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SURCHARGE-SP-6

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	9.1	ng	82713	1	6.95	182658	20.0
unknown6.74	6.74	69.0	ng	630468	1	6.95	182658	20.0
Anthracene, 2-met...	11.99	3.5	ng	42180	4	11.49	243748	20.0
Phenanthrene, 2-m...	12.02	3.4	ng	41003	4	11.49	243748	20.0
4H-Cyclopenta[def...	12.11	4.5	ng	54197	4	11.49	243748	20.0
5,16[1',2']:8,13[...	12.28	2.1	ng	25091	4	11.49	243748	20.0
Phenanthrene, 2,5...	12.54	3.6	ng	44254	4	11.49	243748	20.0
Pyrene, 1-methyl-	13.17	2.8	ng	55953	5	14.15	395587	20.0
11H-Benzo[b]fluorene	13.27	5.0	ng	99646	5	14.15	395587	20.0
Tetradecane	13.95	4.2	ng	82708	5	14.15	395587	20.0
unknown14.26	14.26	2.1	ng	42262	5	14.15	395587	20.0
Chrysene, 1-methyl-	14.57	2.5	ng	48780	5	14.15	395587	20.0