LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

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-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.275	28	32	41	rVB	112566	150172	8.38%	0.490%
2	4.099	339	342	348	rVB	18166	22024	1.23%	0.072%
3	5.193	525	528	535	rVB	70977	81498	4.55%	0.266%
4	5.587	589	595	598	rBV	1838170	1684348	93.98%	5.499%
5	5.816	631	634	639	rVB3	18057	23974	1.34%	0.078%
6	5.916	647	651	657	rBV	28856	28247	1.58%	0.092%
7	6.587	760	765	775	rBV	1900125	1706954	95.24%	5.573%
8	6.740	787	791	794	rBV	1962243	1724037	96.20%	5.628%
9	6.793	798	800	804	rVB2	25626	23870	1.33%	0.078%
10	6.969	827	830	834	rBV	1041422	896426	50.02%	2.927%
11	7.016	834	838	842	rVB2	24012	23140	1.29%	0.076%
12	7.128	853	857	860	rBV	1669262	1321708	73.75%	4.315%
13	7.228	871	874	878	rVB	134954	107093	5.98%	0.350%
14	7.528	921	925	928	rBV	1167651	1007706	56.23%	3.290%
15	8.257	1045	1049	1050	rBV	1274283	1208819	67.45%	3.946%
16	8.275	1050	1052	1055	rVV	1270586	969225	54.08%	3.164%
17	8.328	1058	1061	1065	rVB	58401	47274	2.64%	0.154%
18	8.828	1143	1146	1151	rVB2	22728	24100	1.34%	0.079%
19	8.963	1166	1169	1173	rBV	102232	83035	4.63%	0.271%
20	9.063	1182	1186	1189	rBV	98168	79770	4.45%	0.260%
21	9.328	1227	1231	1234	rBV	2135594	1792228	100.00%	5.851%
22	9.398	1240	1243	1245	rBV	48146	42822	2.39%	0.140%
23	9.428	1245	1248	1252	rVB	55987	44473	2.48%	0.145%
24	9.592	1272	1276	1279	rBV3	47930	54793	3.06%	0.179%
25	9.663	1285	1288	1291	rBV	43308	44790	2.50%	0.146%
26	9.716	1294	1297	1304	rVB	240879	206975	11.55%	0.676%
27	9.781	1304	1308	1310	rBV3	23860	31350	1.75%	0.102%
28	9.875	1317	1324	1327	rBV	212528	206756	11.54%	0.675%
29	9.928	1329	1333	1336	rVB	86343	73666	4.11%	0.240%
30	10.010	1344	1347	1351	rVV	1319423	1212847	67.67%	3.960%
31	10.045	1351	1353	1355	rVB	73538	52094	2.91%	0.170%
32	10.104	1360	1363	1368	rBV6	12395	20817	1.16%	0.068%
33	10.157	1368	1372	1375	rBV5	14669	23043	1.29%	0.075%
34	10.216	1379	1382	1385	rVB	140111	114528	6.39%	0.374%

LSC Area Percent Report

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Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

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-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	10.251	1385	1388	1392	rBV2	33985	34265	1.91%	0.112%
36	10.322	1397	1400	1403	rBV4	23179	26685	1.49%	0.087%
37	10.404	1411	1414	1415	rBV2	27284	22101	1.23%	0.072%
38	10.428	1415	1418	1421	rVB	101684	93375	5.21%	0.305%
39	10.557	1437	1440	1447	rVB	126583	130465	7.28%	0.426%
40	10.645	1450	1455	1458	rVV	41896	53703	3.00%	0.175%
41	10.716	1462	1467	1471	rBV2	35893	44711	2.49%	0.146%
42	10.798	1478	1481	1485	rVB	1068716	1005368	56.10%	3.282%
43	10.904	1496	1499	1508	rVB3	88378	156535	8.73%	0.511%
44	11.110	1529	1534	1536	rBV4	29765	44305	2.47%	0.145%
45	11.257	1557	1559	1565	rVB6	19186	25762	1.44%	0.084%
46	11.310	1565	1568	1571	rVV2	36807	36268	2.02%	0.118%
47	11.351	1572	1575	1577	rVV	83646	68564	3.83%	0.224%
48	11.380	1577	1580	1582	rVV	74595	79462	4.43%	0.259%
49	11.398	1582	1583	1587	rVB	65959	55889	3.12%	0.182%
50	11.504	1597	1601	1603	rBV	1194321	1108142	61.83%	3.618%
51	11.527	1603	1605	1609	rVV	993803	822501	45.89%	2.685%
52	11.580	1609	1614	1617	rVV	376667	309139	17.25%	1.009%
53	11.616	1617	1620	1625	rVB2	45951	65579	3.66%	0.214%
54	11.663	1625	1628	1633	rBV7	22182	38951	2.17%	0.127%
55	11.733	1637	1640	1644	rVB	117235	96400	5.38%	0.315%
56	11.780	1645	1648	1650	rBV	55584	45366	2.53%	0.148%
57	11.839	1655	1658	1662	rVV2	32101	50585	2.82%	0.165%
58	11.880	1663	1665	1669	rVB5	27781	23924	1.33%	0.078%
59	12.010	1684	1687	1689	rBV2	211188	228921	12.77%	0.747%
60	12.033	1689	1691	1697	rVB2	157619	148890	8.31%	0.486%
61	12.080	1697	1699	1702	rVB	59181	49108	2.74%	0.160%
62	12.122	1702	1706	1708	rBV2	228301	224918	12.55%	0.734%
63	12.186	1715	1717	1719	rBV2	52980	43947	2.45%	0.143%
64	12.298	1734	1736	1739	rBV	79475	71193	3.97%	0.232%
65	12.327	1739	1741	1745	rVV3	58630	58010	3.24%	0.189%
66	12.557	1777	1780	1782	rBV	62117	64831	3.62%	0.212%
67	12.645	1792	1795	1798	rBV	64975	63124	3.52%	0.206%
68	12.722	1805	1808	1812	rBV	1289524	1221321	68.15%	3.987%
69	12.816	1822	1824	1827	rVB	138408	119739	6.68%	0.391%
70	12.933	1841	1844	1845	rBV	207793	177108	9.88%	0.578%
71	12.957	1845	1848	1853	rVV	1216650	1101109	61.44%	3.595%

LSC Area Percent Report

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Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

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-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	13.004	1853	1856	1859	rVB2	68506	70044	3.91%	0.229%
73	13.092	1864	1871	1874	rBV	1647563	1560562	87.07%	5.095%
74	13.169	1880	1884	1888	rBV2	82375	150794	8.41%	0.492%
75	13.257	1896	1899	1900	rBV3	65667	69197	3.86%	0.226%
76	13.280	1900	1903	1906	rVB2	200978	179607	10.02%	0.586%
77	13.310	1906	1908	1910	rBV2	67308	54179	3.02%	0.177%
78	13.345	1912	1914	1917	rVB	172493	155321	8.67%	0.507%
79	13.386	1918	1921	1924	rVB2	79266	81085	4.52%	0.265%
80	13.480	1935	1937	1939	rBV	61167	42361	2.36%	0.138%
81	13.510	1940	1942	1945	rBV2	59836	70960	3.96%	0.232%
82	13.933	2012	2014	2016	rVB	82712	58197	3.25%	0.190%
83	13.974	2016	2021	2024	rBV4	228339	315106	17.58%	1.029%
84	14.116	2043	2045	2047	rVB2	112593	83411	4.65%	0.272%
85	14.157	2047	2052	2055	rVV2	1157659	1575539	87.91%	5.144%
86	14.186	2055	2057	2063	rVB	538853	473448	26.42%	1.546%
87	14.780	2155	2158	2162	rBV	421583	416575	23.24%	1.360%
88	15.239	2232	2236	2239	rBV	404118	615815	34.36%	2.010%
89	15.563	2288	2291	2298	rVB	232857	332006	18.52%	1.084%
90	15.633	2299	2303	2309	rVB	351717	483920	27.00%	1.580%
91	15.698	2310	2314	2318	rBV	519444	697633	38.93%	2.278%

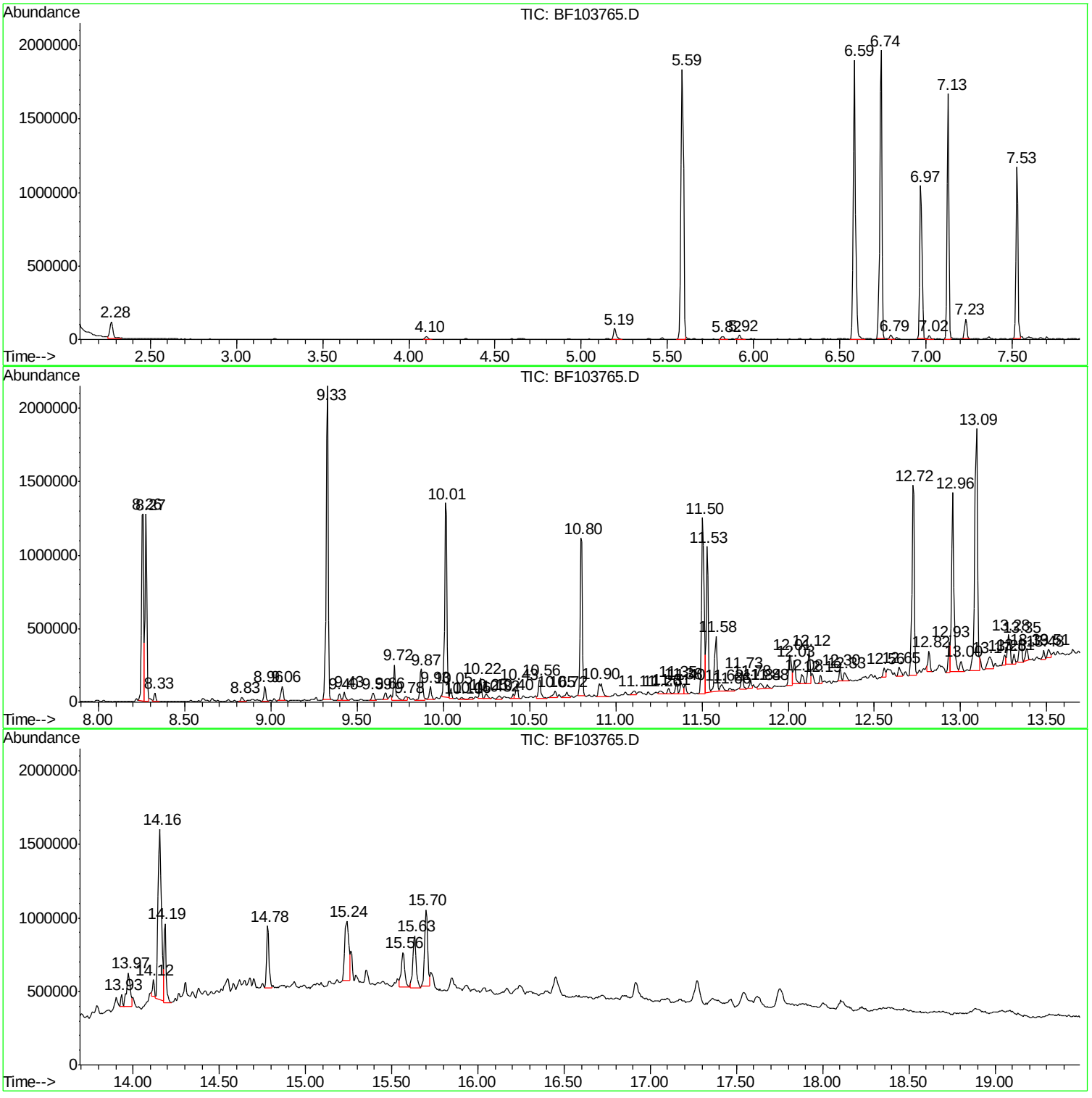
Sum of corrected areas: 30630626

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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\BF031918\
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Acq On : 19 Mar 2018 19:36
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

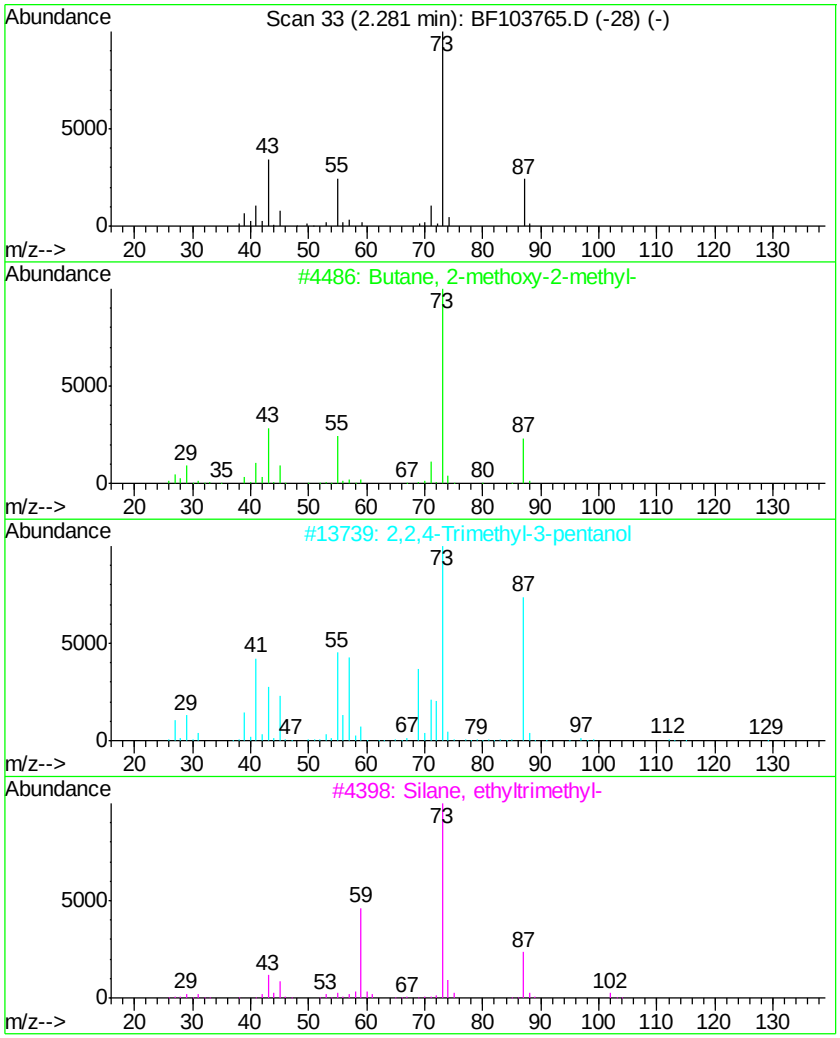
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

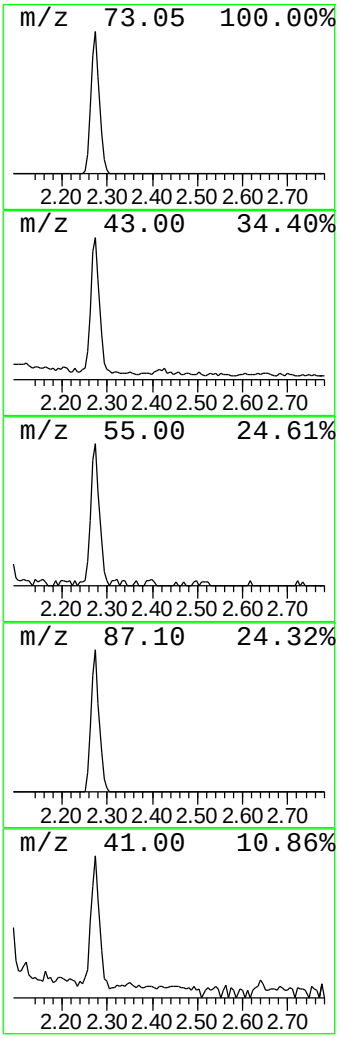
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.28	3.35 ng	150172	1,4-Dichlorobenzene-d4	6.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4	1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	25
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23





Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

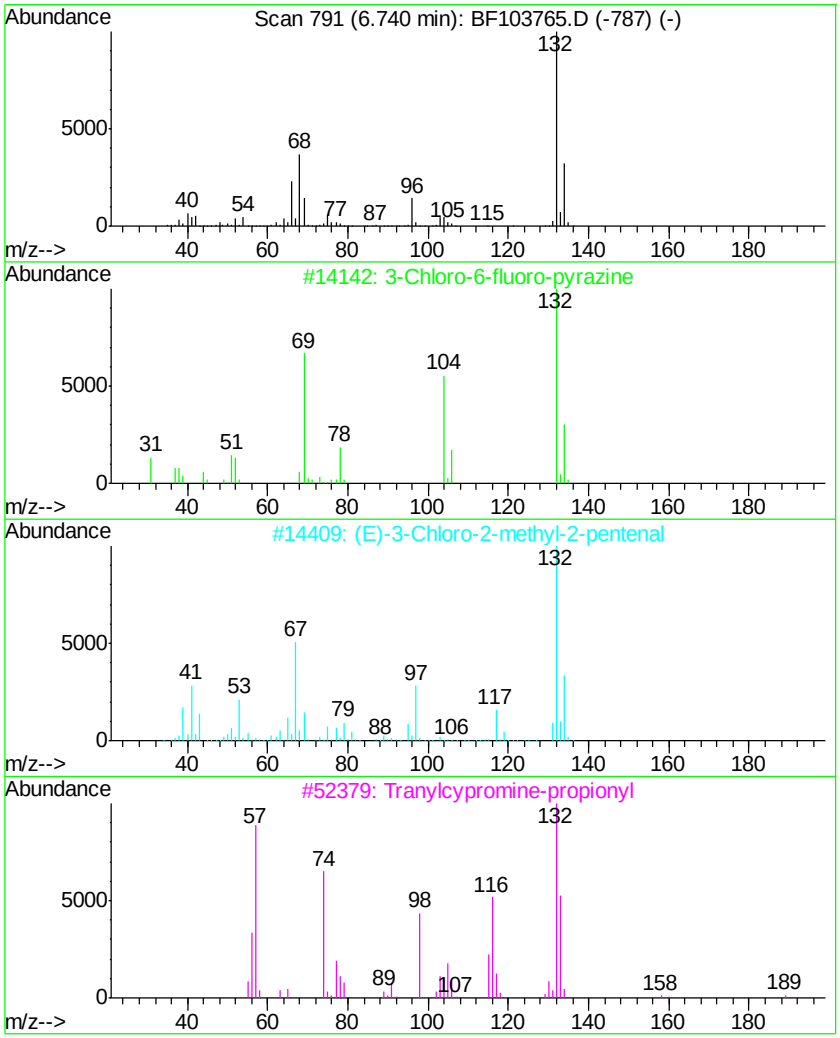
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

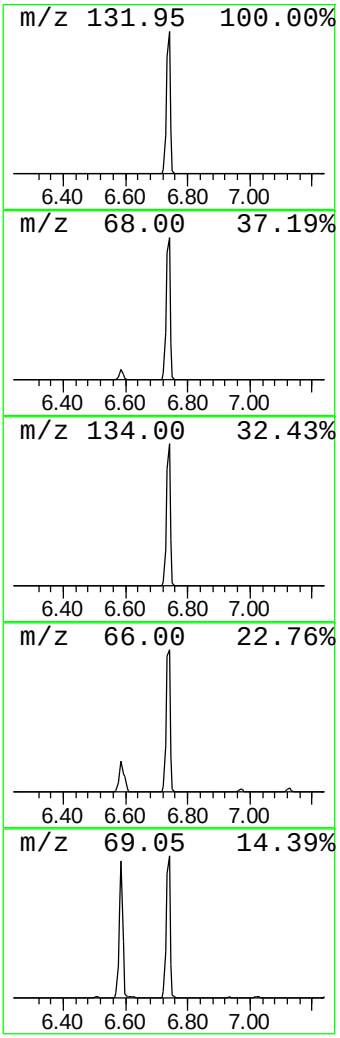
****

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	38.46 ng	1724040	1,4-Dichlorobenzene-d4	6.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	5-Aminoindole	132	C8H8N2	005192-03-0	12
5	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
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Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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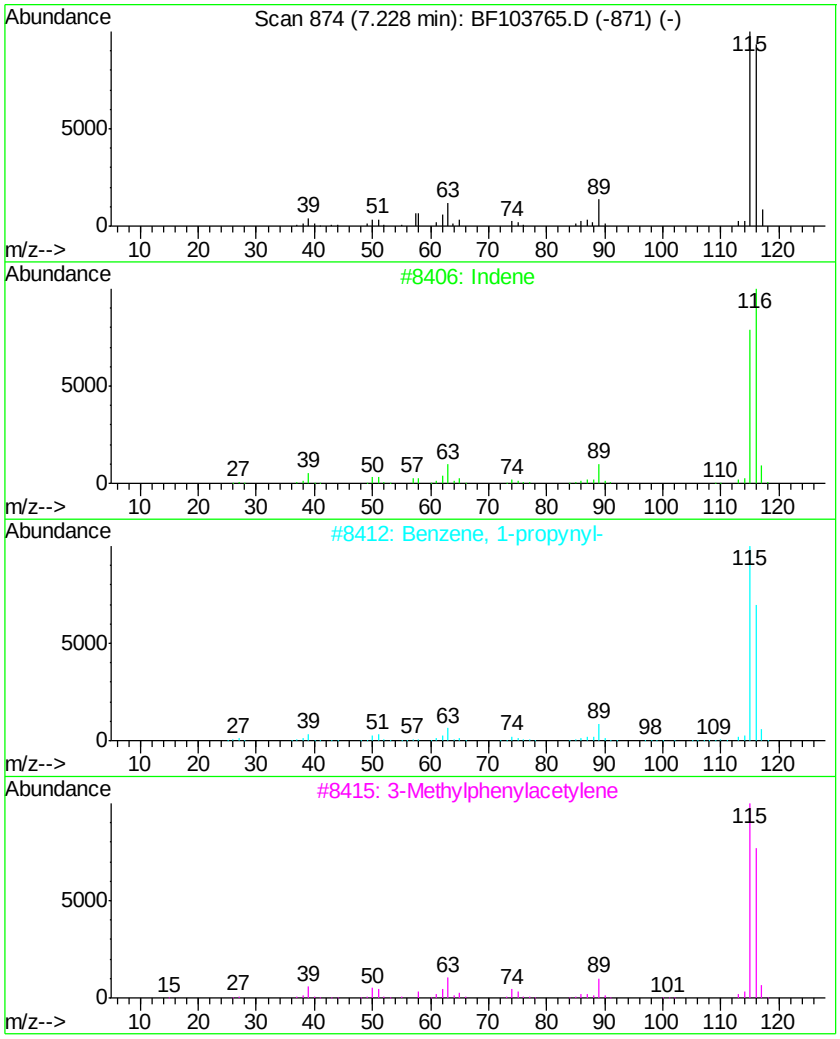
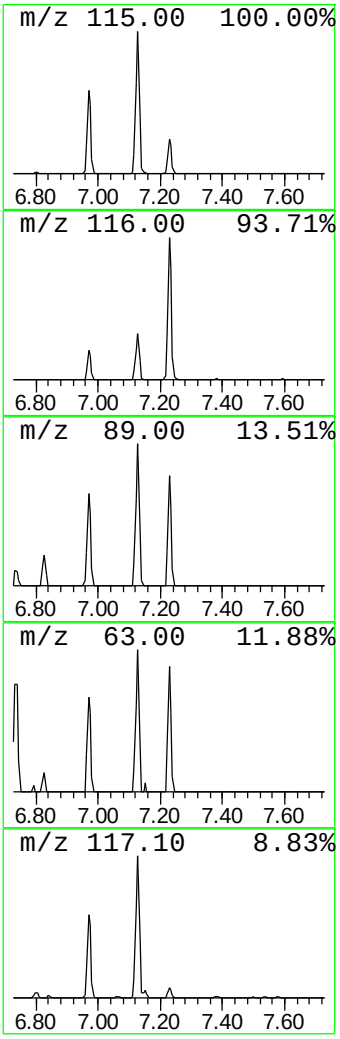
Peak Number 4 Indene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.23	2.39 ng	107093	1,4-Dichlorobenzene-d4	6.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
3	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
4	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
5	2-Methylphenylacetylene	116	C9H8	000766-47-2	91





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Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

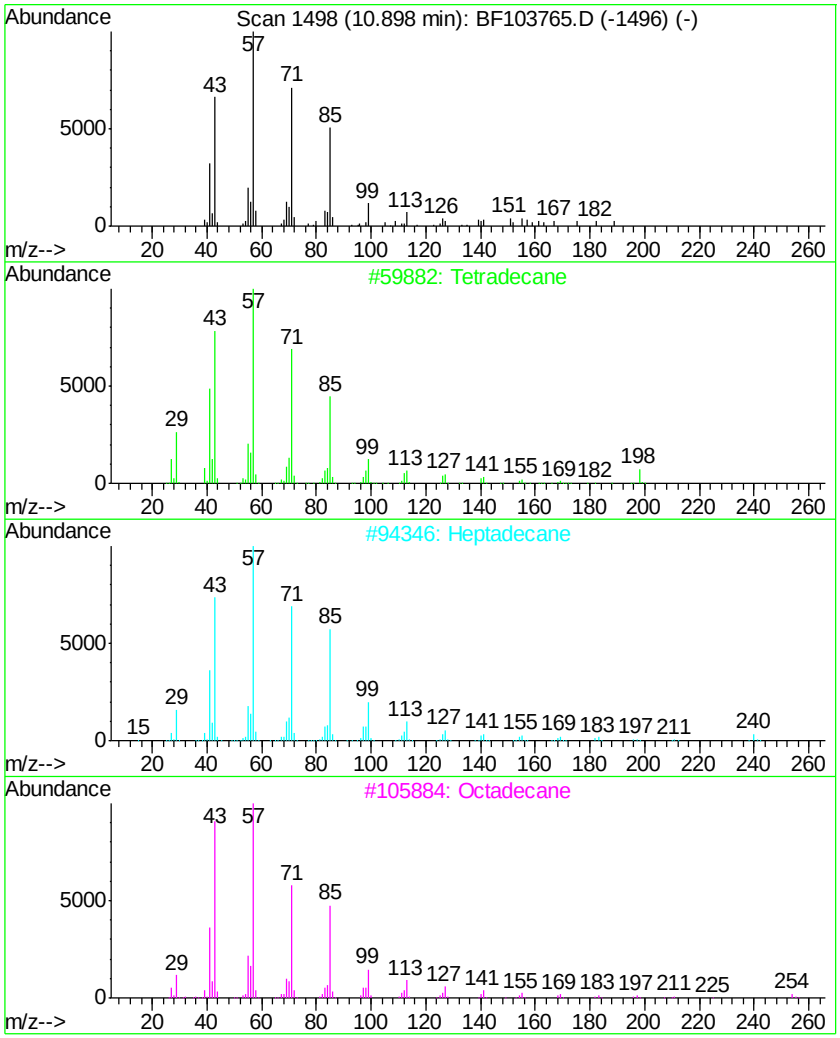
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

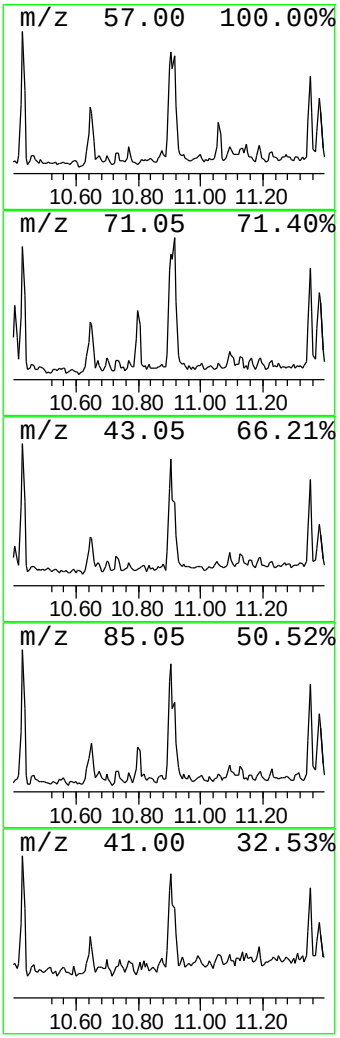
****

Peak Number 6 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	2.83 ng	156535	Phenanthrene-d10	11.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Heptadecane	240	C17H36	000629-78-7	90
3	Octadecane	254	C18H38	000593-45-3	86
4	Pentadecane	212	C15H32	000629-62-9	86
5	Octacosane	394	C28H58	000630-02-4	86





Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

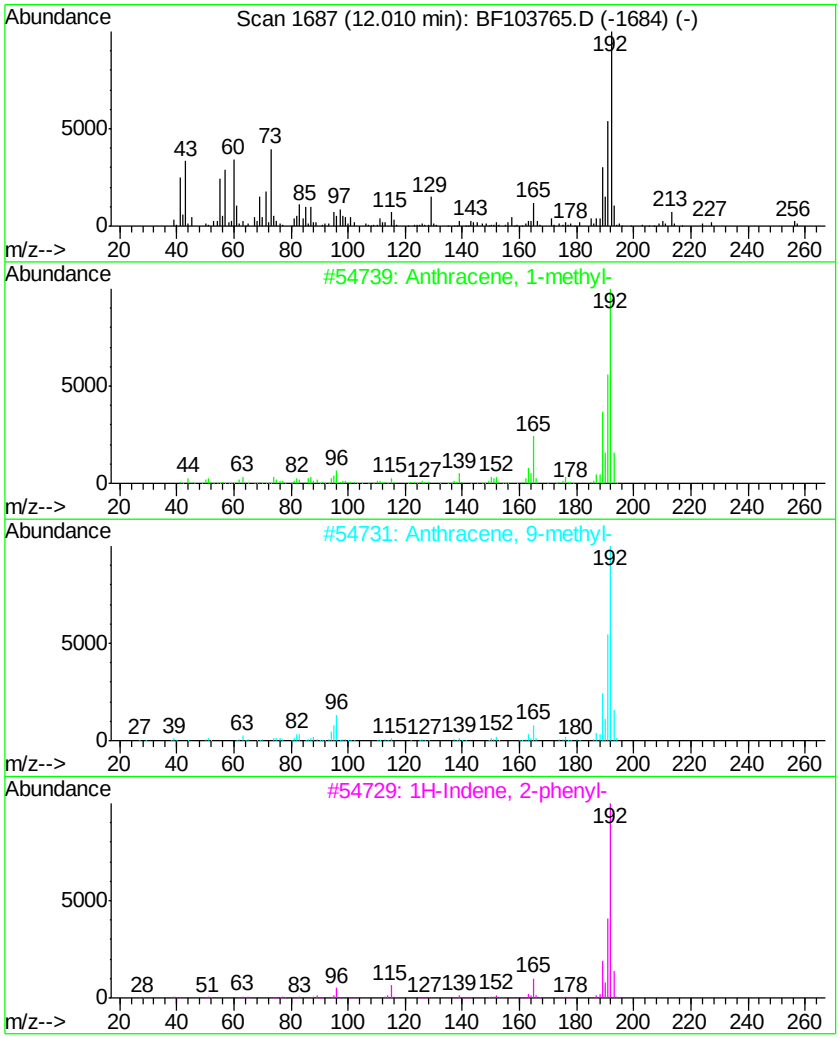
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

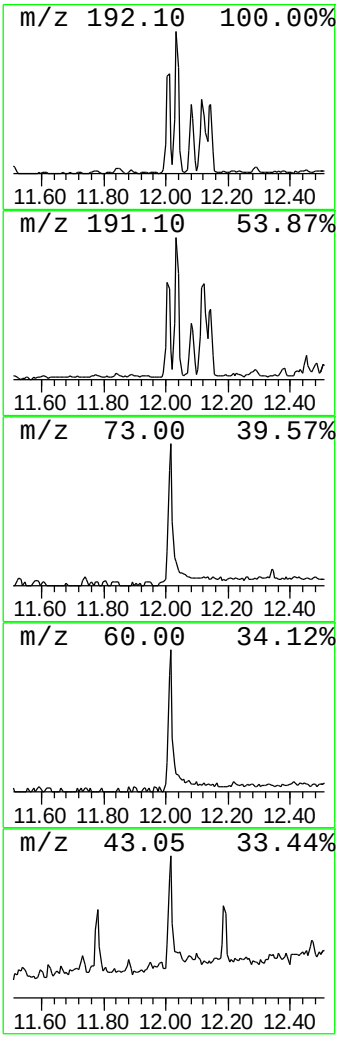
****

Peak Number 7 Anthracene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	4.13 ng	228921	Phenanthrene-d10	11.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	86
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	64
4	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	64
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	60





Data Path : Z:\HPCHEM1\BNA F\DATA\BF031918\
Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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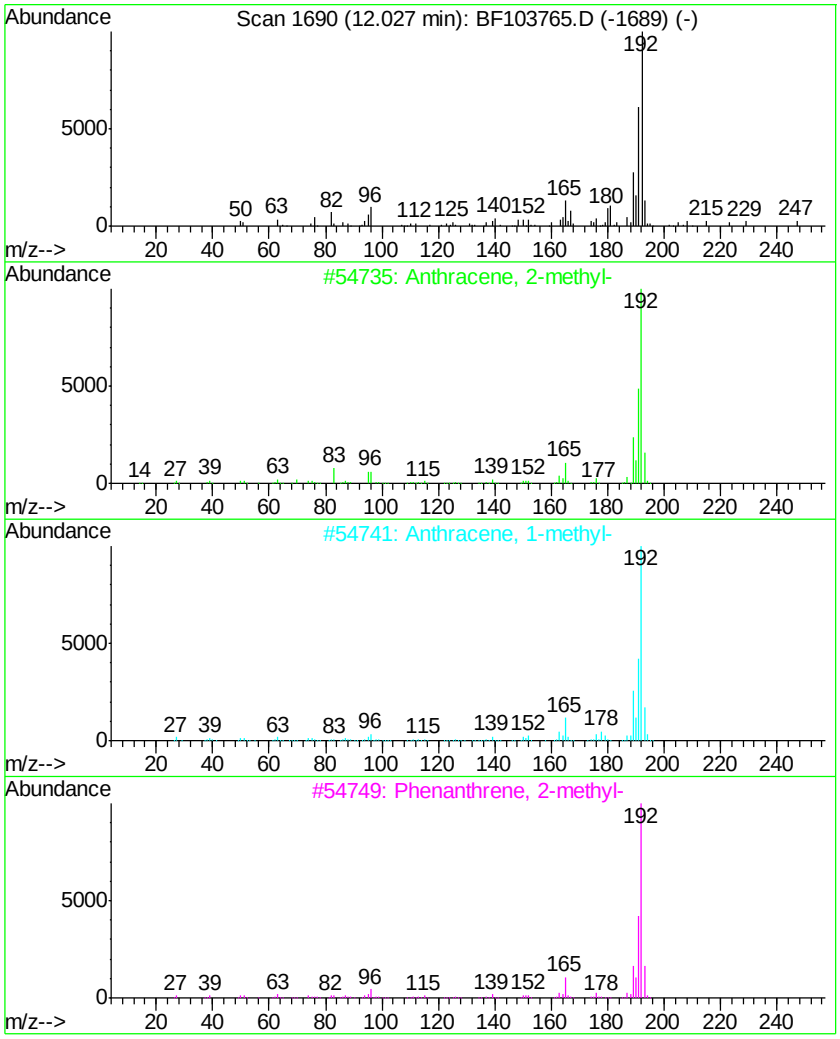
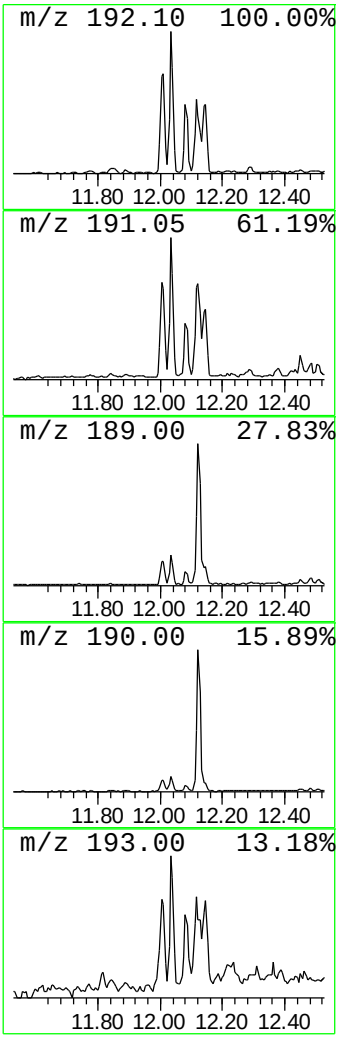
Peak Number 8 Anthracene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	2.69 ng	148890	Phenanthrene-d10	11.50

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	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	87
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	87





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Data File : BF103765.D
Acq On : 19 Mar 2018 19:36
Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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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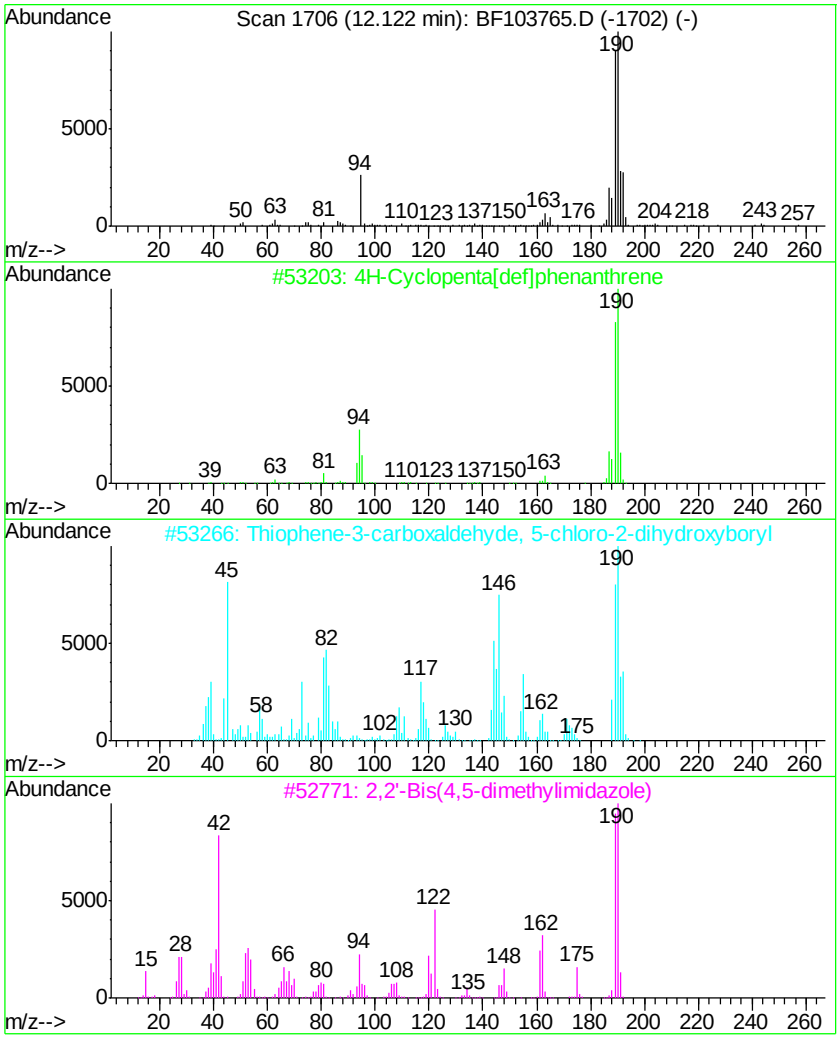
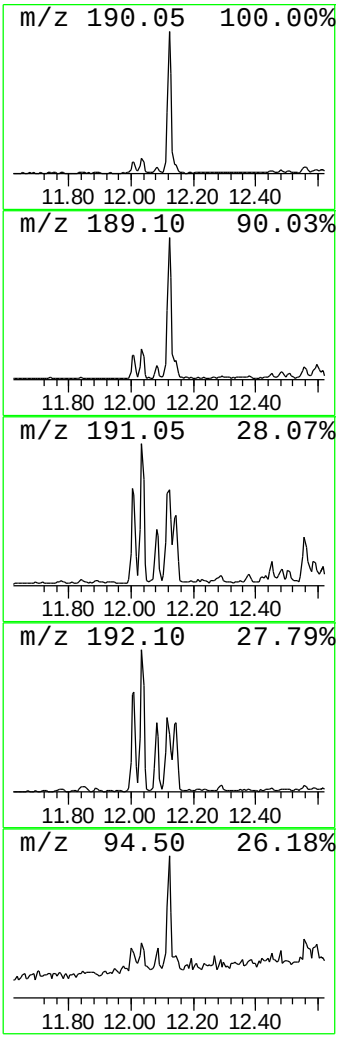
Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	4.06 ng	224918	Phenanthrene-d10	11.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	32
5	2,6-Dichlorobenzaldoxime	189	C7H5Cl2NO	025185-95-9	18





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Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
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TIC Library : C:\DATABASE\NIST11.L
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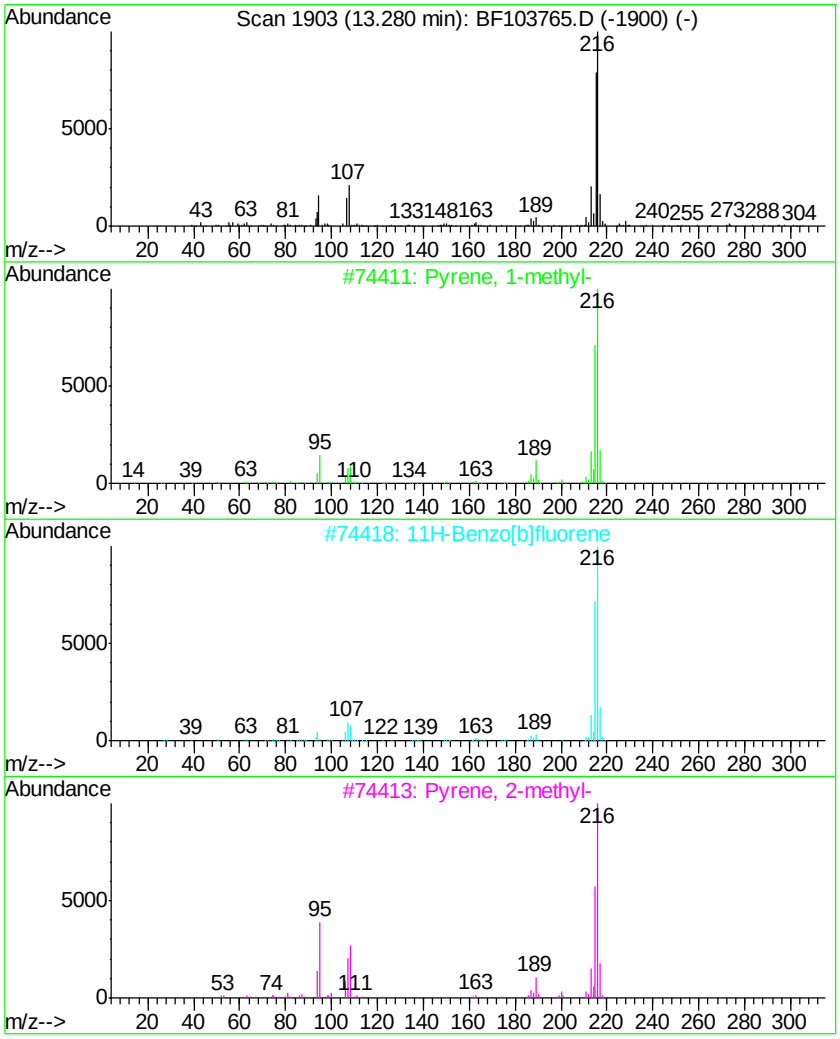
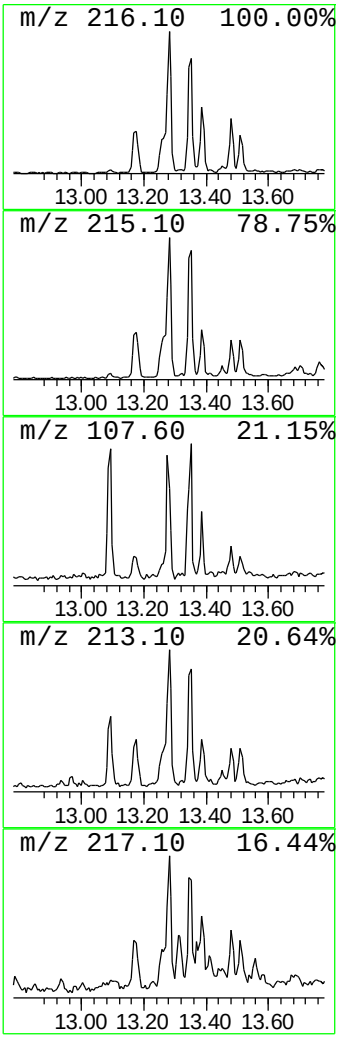
Peak Number 11 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.28	2.28 ng	179607	Chrysene-d12	14.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	76
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	68
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	68





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Operator : JU/SJ
Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
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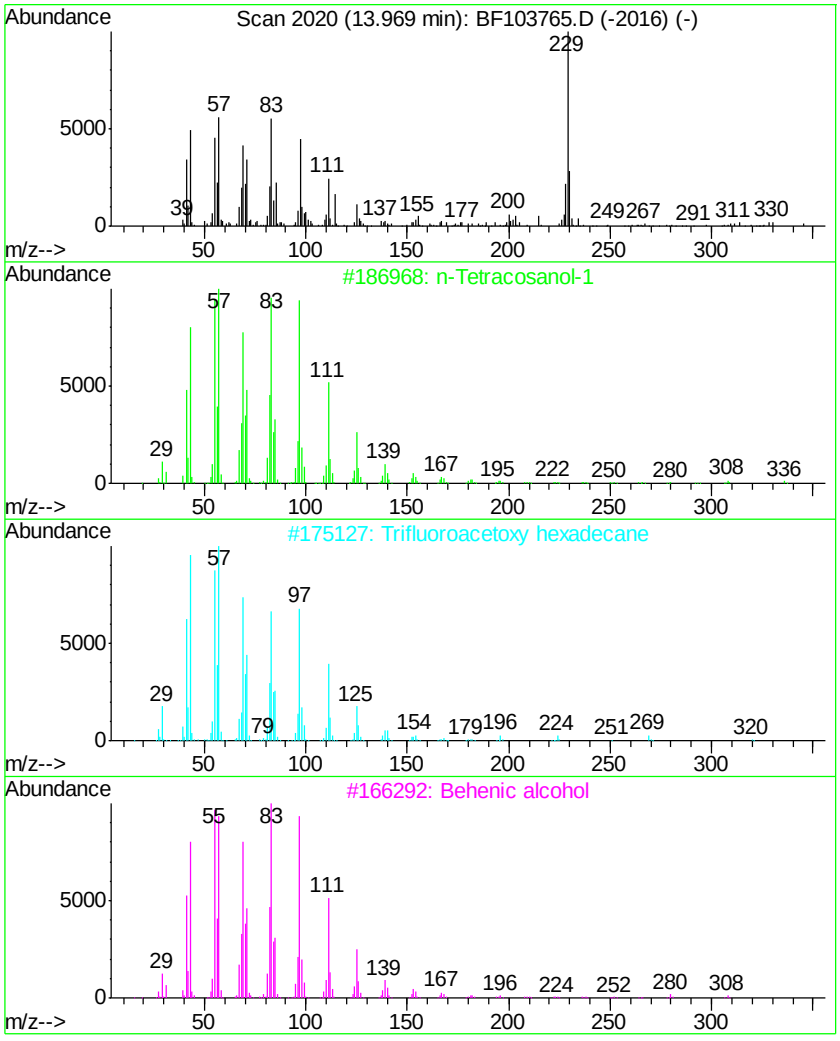
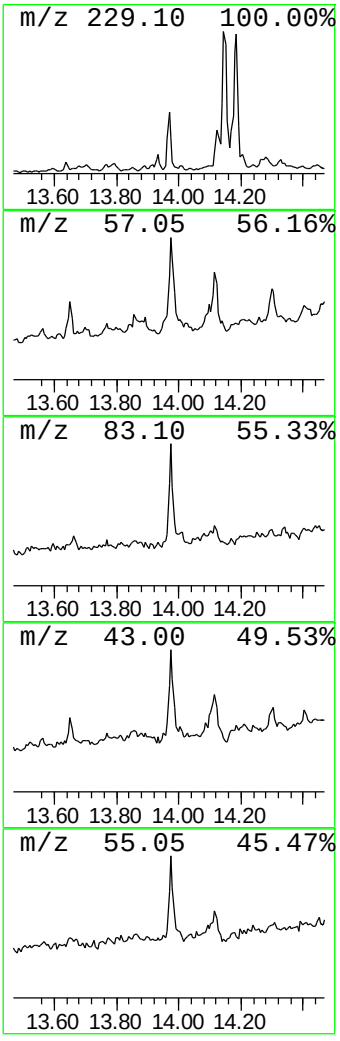
Peak Number 12 n-Tetracosanol-1 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	4.00 ng	315106	Chrysene-d12	14.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	90
2	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	81
3	Behenic alcohol	326	C22H46O	000661-19-8	81
4	1-Heneicosanol	312	C21H44O	015594-90-8	81
5	1-Heptacosanol	396	C27H56O	002004-39-9	76





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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
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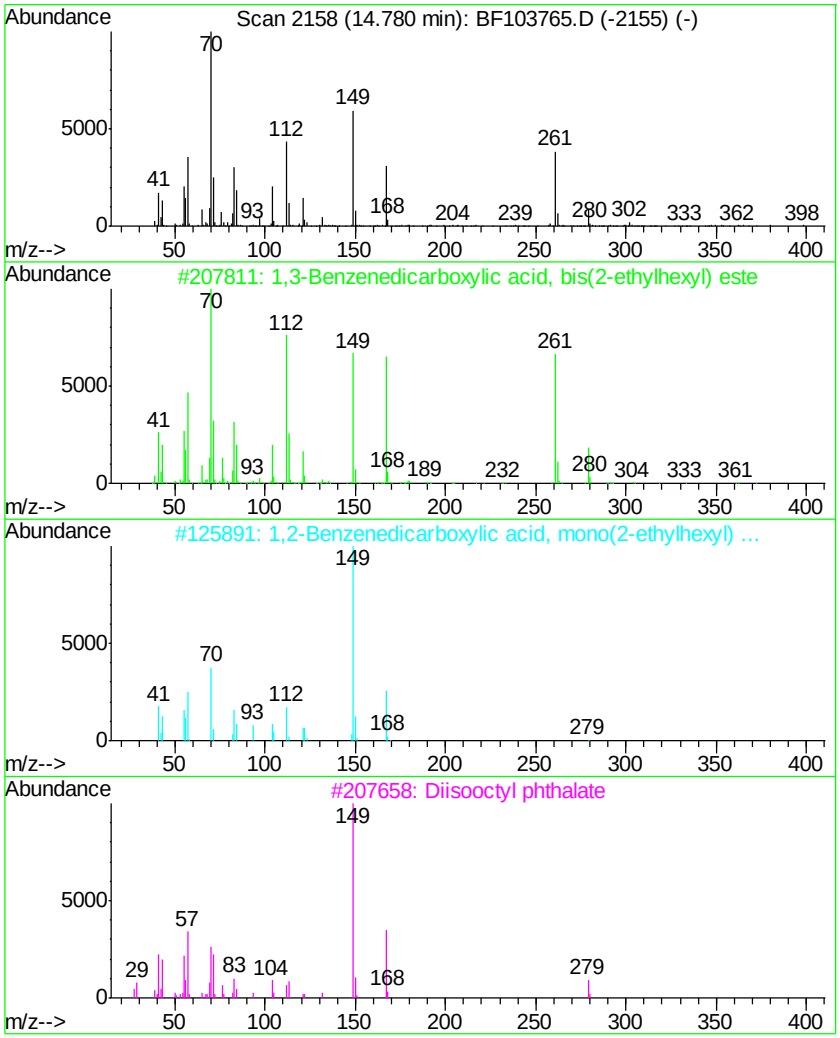
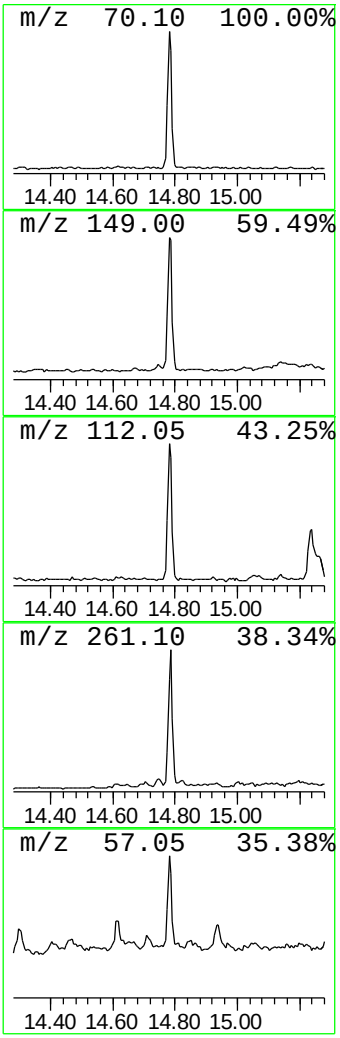
Peak Number 13 1,3-Benzenedicarboxylic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	5.29 ng	416575	Chrysene-d12	14.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	58
2	1,2-Benzenedicarboxylic acid, mo...	278	C16H22O4	004376-20-9	38
3	Diisooctyl phthalate	390	C24H38O4	000131-20-4	22
4	Bromoacetic acid, 2-ethylhexyl e...	250	C10H19BrO2	068144-73-0	16
5	Carbonic acid, propargyl 2-ethyl...	212	C12H20O3	1000357-90-9	14





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Sample : J1965-02 2X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-112(6-7)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
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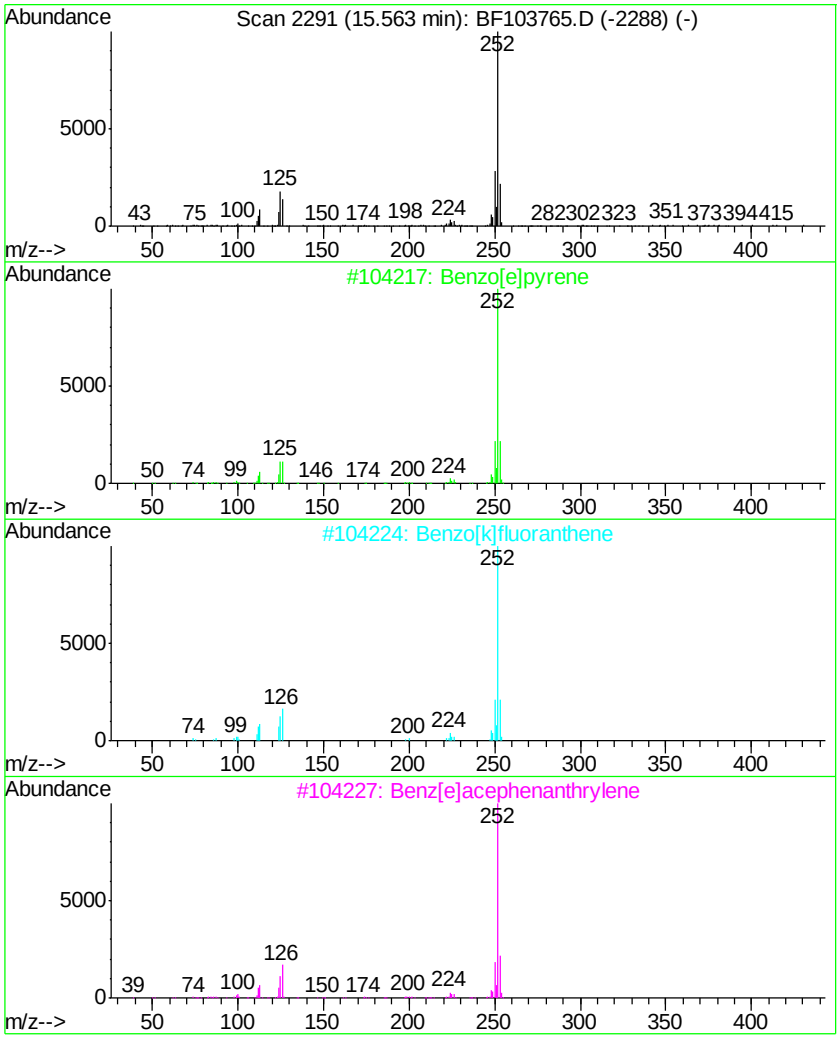
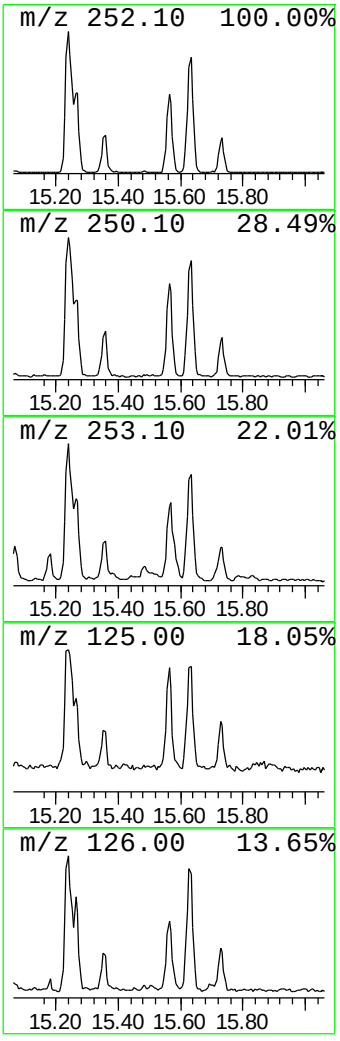
Peak Number 14 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.56	9.52 ng	332006	Perylene-d12	15.70

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	96
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
4	Perylene	252	C20H12	000198-55-0	93
5	Benzo[a]pyrene	252	C20H12	000050-32-8	90





Tentatively Identified Compound (LSC) summary

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Misc :
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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
Butane, 2-methoxy...	2.28	3.4	ng	150172	1	6.97	896426	20.0
unknown6.74	6.74	38.5	ng	1724040	1	6.97	896426	20.0
Indene	7.23	2.4	ng	107093	1	6.97	896426	20.0
Tetradecane	10.90	2.8	ng	156535	4	11.50	1108140	20.0
Anthracene, 1-met...	12.01	4.1	ng	228921	4	11.50	1108140	20.0
Anthracene, 2-met...	12.03	2.7	ng	148890	4	11.50	1108140	20.0
4H-Cyclopenta[def...	12.12	4.1	ng	224918	4	11.50	1108140	20.0
Pyrene, 1-methyl-	13.28	2.3	ng	179607	5	14.16	1575540	20.0
n-Tetracosanol-1	13.97	4.0	ng	315106	5	14.16	1575540	20.0
1,3-Benzenedicarb...	14.78	5.3	ng	416575	5	14.16	1575540	20.0
Benzo[e]pyrene	15.56	9.5	ng	332006	6	15.70	697633	20.0