LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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	3.087	168	170	181	rBV	42482	87179	2.48%	0.279%
2	4.928	479	483	494	rVB	206062	302830	8.62%	0.968%
3	5.339	547	553	556	rBV	2955725	3302439	94.03%	10.555%
4	6.369	722	728	731	rBV	3001343	3511999	100.00%	11.225%
5	6.498	745	750	753	rBV	3198268	3278101	93.34%	10.477%
6	6.722	785	788	792	rBV	881148	717027	20.42%	2.292%
7	6.881	811	815	818	rBV	2716519	2518692	71.72%	8.050%
8	7.292	879	885	888	rBV	2152588	2160833	61.53%	6.906%
9	8.004	1002	1006	1022	rVB	991241	980875	27.93%	3.135%
10	8.716	1122	1127	1131	rBV	46648	44626	1.27%	0.143%
11	8.816	1140	1144	1149	rBV	74296	61695	1.76%	0.197%
12	9.086	1184	1190	1193	rBV	3637609	3377285	96.16%	10.794%
13	9.345	1230	1234	1238	rBV3	31579	38260	1.09%	0.122%
14	9.416	1242	1246	1249	rBV	69338	67955	1.93%	0.217%
15	9.475	1253	1256	1263	rVB	524797	474239	13.50%	1.516%
16	9.533	1263	1266	1271	rBV	34852	37505	1.07%	0.120%
17	9.616	1277	1280	1284	rBV	163819	138799	3.95%	0.444%
18	9.757	1297	1304	1308	rBV	1101121	948050	26.99%	3.030%
19	10.069	1353	1357	1360	rBV2	35382	41549	1.18%	0.133%
20	10.163	1369	1373	1375	rBV2	33846	46321	1.32%	0.148%
21	10.304	1394	1397	1403	rVB2	45269	64196	1.83%	0.205%
22	10.545	1432	1438	1441	rBV	1708295	1590364	45.28%	5.083%
23	10.669	1455	1459	1466	rBV2	73531	91652	2.61%	0.293%
24	10.851	1485	1490	1492	rBV3	32995	45379	1.29%	0.145%
25	11.133	1536	1538	1544	rVB2	39518	42339	1.21%	0.135%
26	11.239	1552	1556	1558	rBV	936177	814346	23.19%	2.603%
27	11.263	1558	1560	1563	rVV	377514	289448	8.24%	0.925%
28	11.310	1565	1568	1571	rVB	72790	68384	1.95%	0.219%
29	11.568	1609	1612	1616	rVB2	39304	36920	1.05%	0.118%
30	11.739	1637	1641	1643	rBV	157689	140169	3.99%	0.448%
31	11.768	1643	1646	1649	rVV	166414	152339	4.34%	0.487%
32	11.816	1651	1654	1656	rVV	43424	44108	1.26%	0.141%
33	11.845	1656	1659	1662	rVV2	160422	183732	5.23%	0.587%
34	11.868	1662	1663	1670	rVB	105811	97158	2.77%	0.311%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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	12.033	1686	1691	1693	rBV3	53958	59093	1.68%	0.189%
36	12.057	1693	1695	1701	rVB4	33976	40674	1.16%	0.130%
37	12.216	1719	1722	1724	rBV	54832	51491	1.47%	0.165%
38	12.286	1731	1734	1737	rBV	142327	140994	4.01%	0.451%
39	12.321	1737	1740	1742	rVV2	72473	83954	2.39%	0.268%
40	12.368	1746	1748	1753	rBV3	59388	85147	2.42%	0.272%
41	12.445	1758	1761	1765	rBV	253782	221113	6.30%	0.707%
42	12.539	1775	1777	1780	rVB	46791	41168	1.17%	0.132%
43	12.674	1796	1800	1803	rBV	386994	338719	9.64%	1.083%
44	12.733	1807	1810	1814	rVB2	50903	58333	1.66%	0.186%
45	12.827	1821	1826	1829	rBV	2453996	2434372	69.32%	7.781%
46	12.892	1834	1837	1841	rBV2	60573	92450	2.63%	0.295%
47	12.986	1850	1853	1854	rBV3	42115	50733	1.44%	0.162%
48	13.104	1870	1873	1878	rBV	87024	103459	2.95%	0.331%
49	13.198	1887	1889	1891	rVB	79770	54831	1.56%	0.175%
50	13.227	1892	1894	1898	rVV2	74915	64814	1.85%	0.207%
51	13.721	1975	1978	1982	rBV2	122688	127453	3.63%	0.407%
52	13.868	1999	2003	2006	rBV2	779322	856558	24.39%	2.738%
53	15.286	2241	2244	2254	rVB2	408111	585837	16.68%	1.872%

Sum of corrected areas: 31287986

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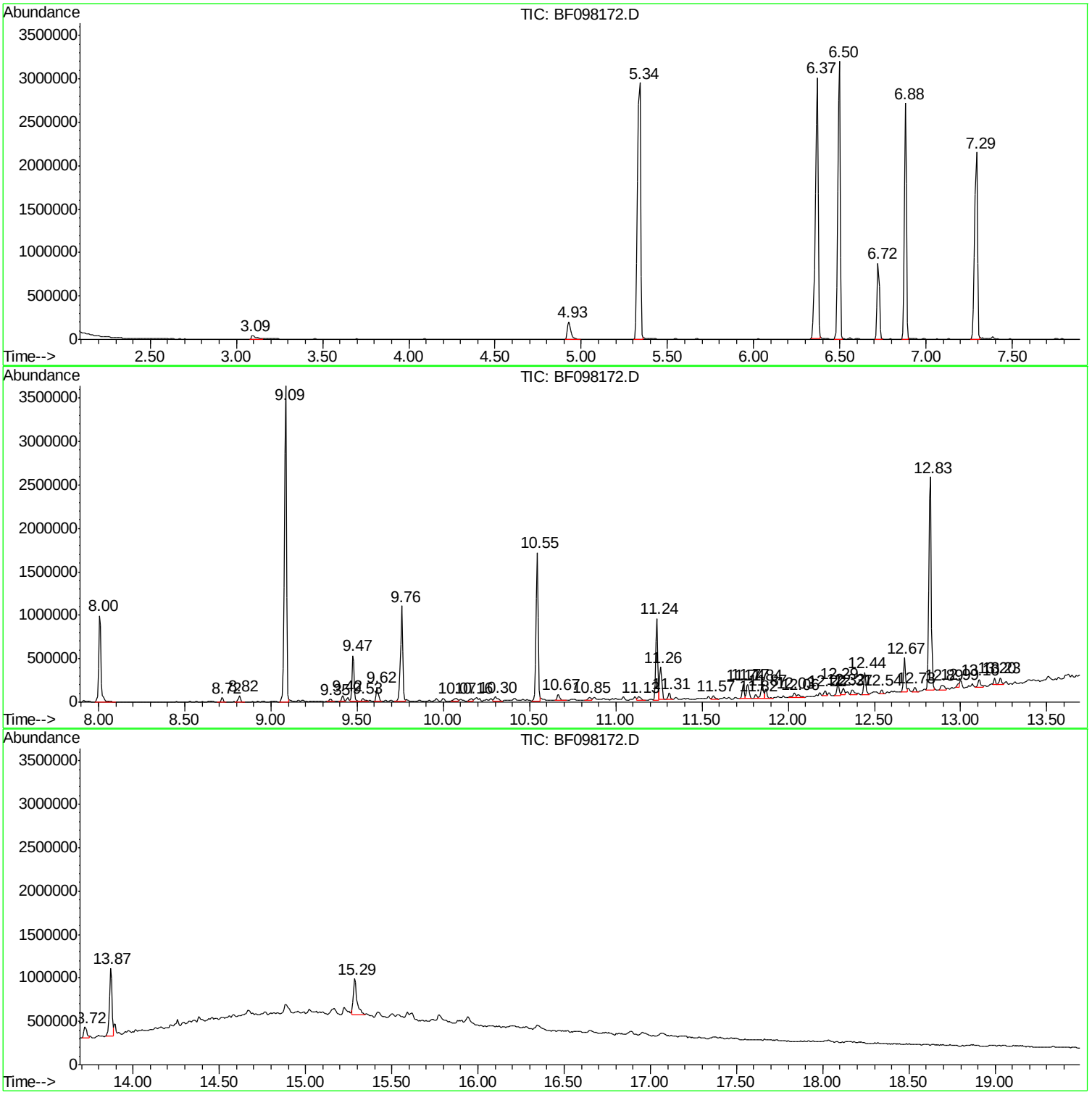
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CL-01-082917-A

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\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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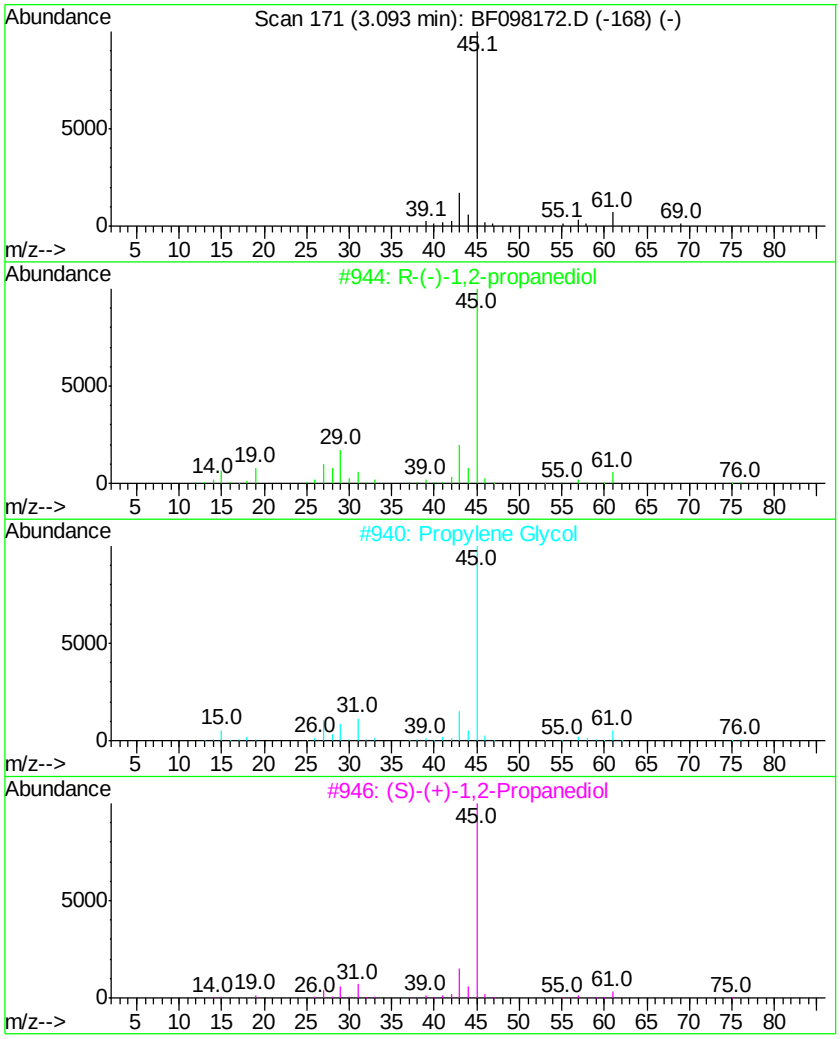
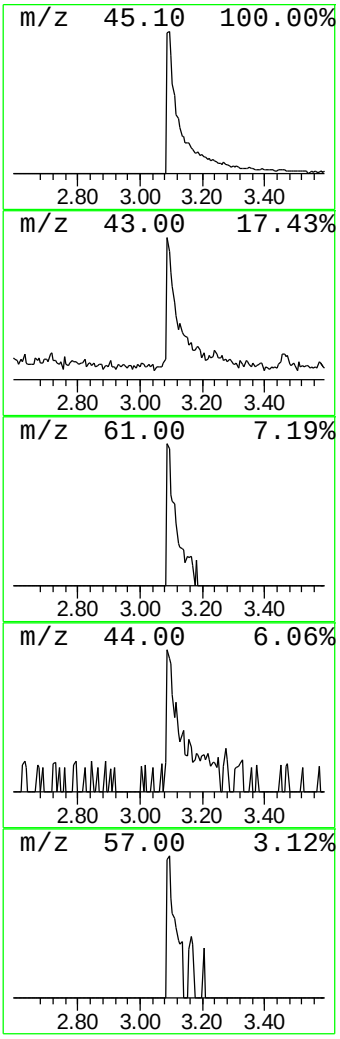
Peak Number 1 unknown3.09 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.09	2.43 ng	87179	1,4-Dichlorobenzene-d4	6.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	43
2	Propylene Glycol	76	C3H8O2	000057-55-6	9
3	(S)-(+)-1,2-Propanediol	76	C3H8O2	004254-15-3	9
4	2-Butanol, (R)-	74	C4H10O	014898-79-4	9
5	Formamide	45	CH3NO	000075-12-7	5





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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

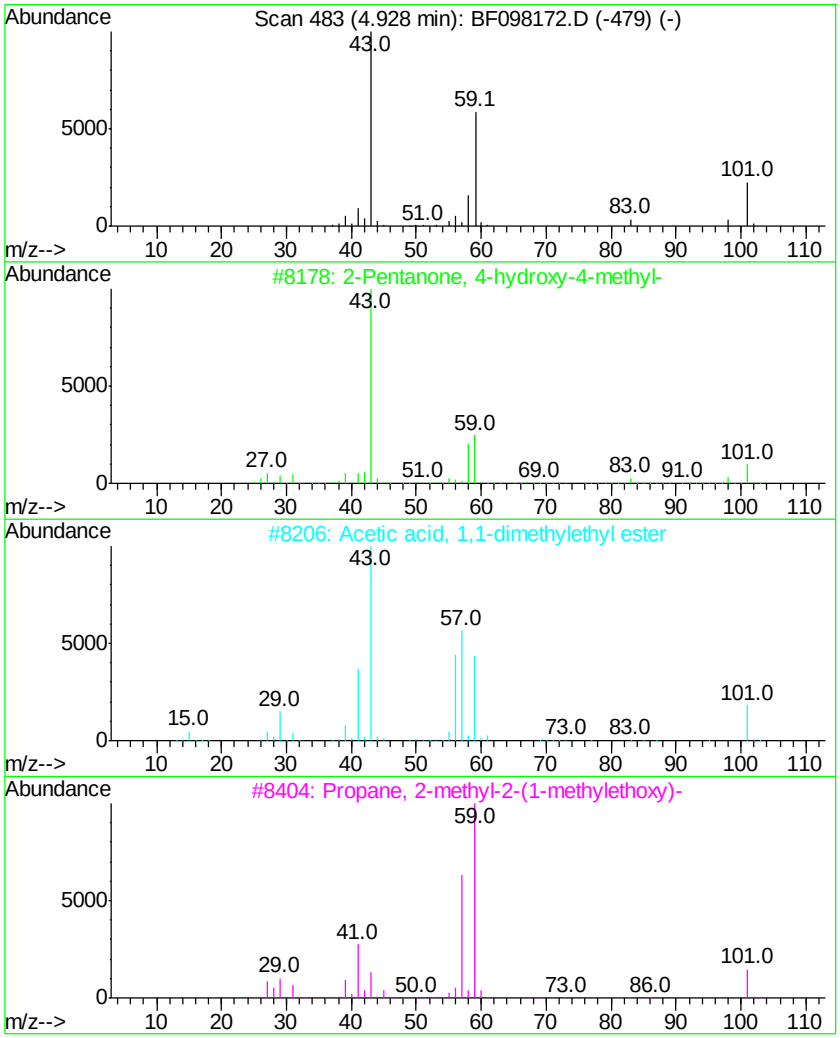
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

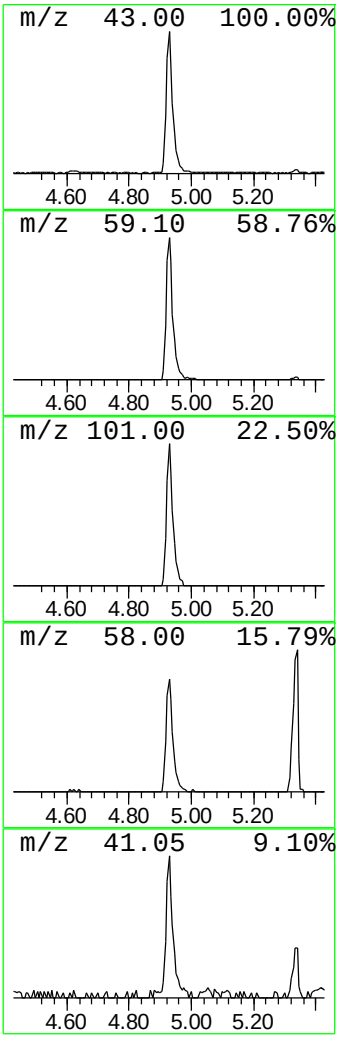
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	8.45 ng	302830	1,4-Dichlorobenzene-d4	6.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	23
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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

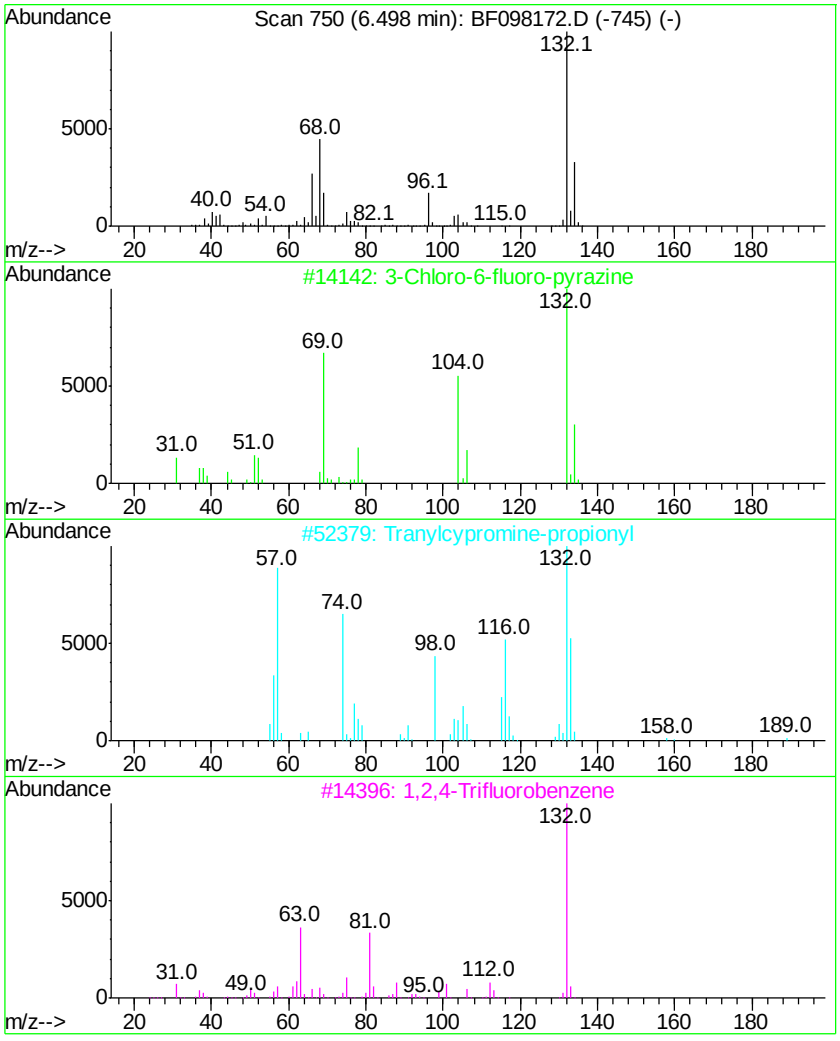
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

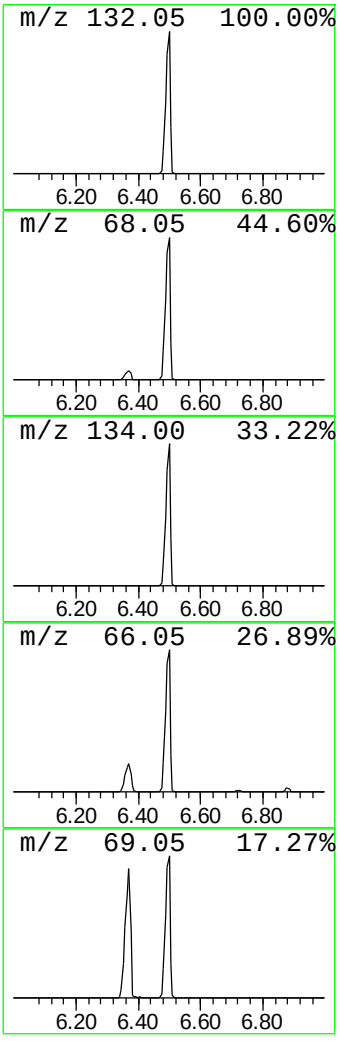
****

Peak Number 3 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	91.44 ng	3278100	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	27
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3	1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9
4	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
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Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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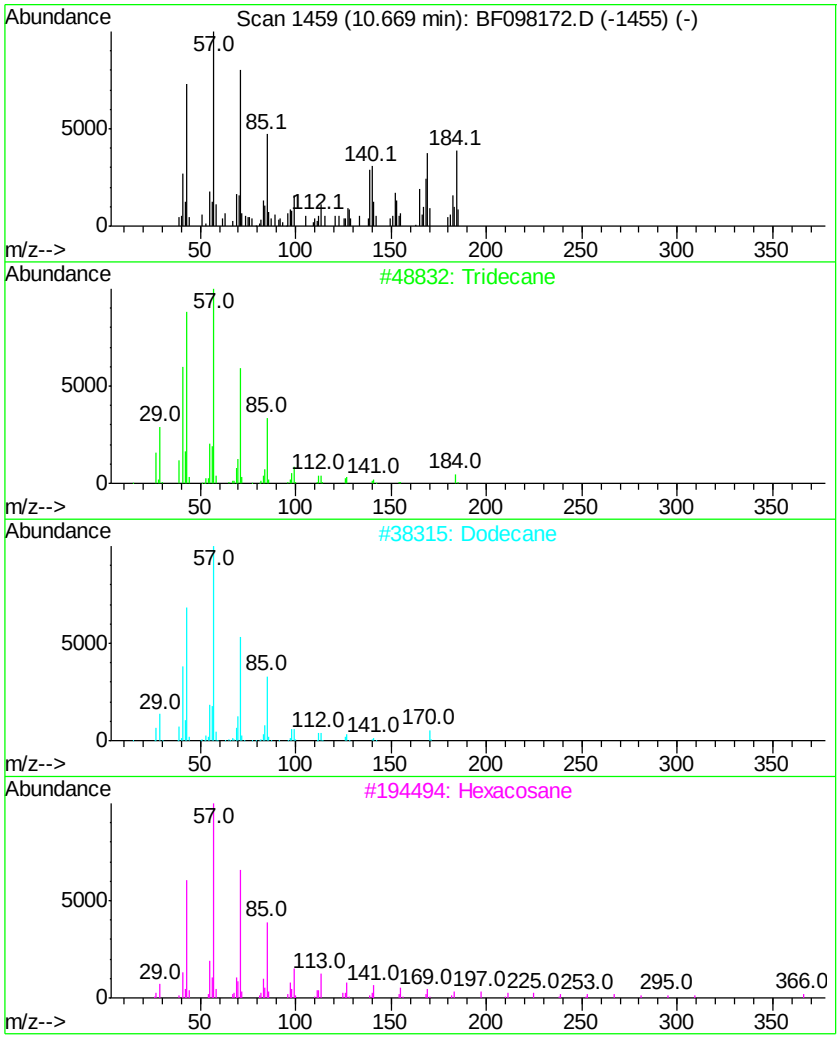
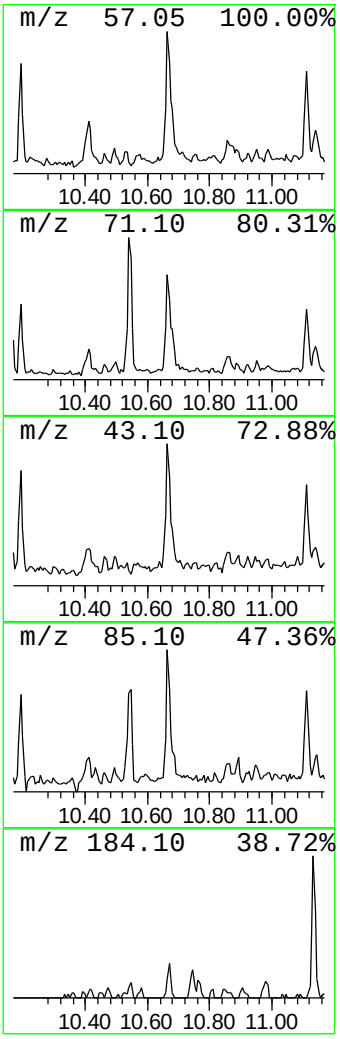
Peak Number 5 unknown10.67 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	2.25 ng	91652	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	49
2	Dodecane	170	C12H26	000112-40-3	46
3	Hexacosane	366	C26H54	000630-01-3	41
4	Hentriacontane	437	C31H64	000630-04-6	38
5	Octacosane	394	C28H58	000630-02-4	38





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Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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
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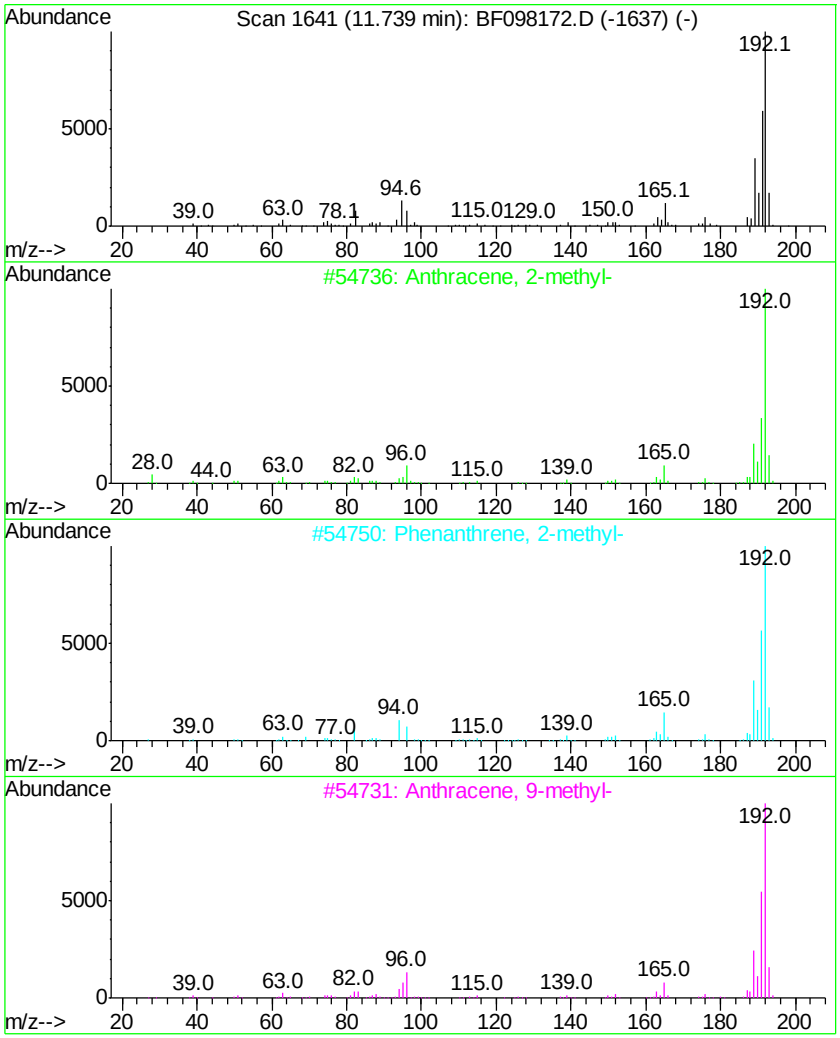
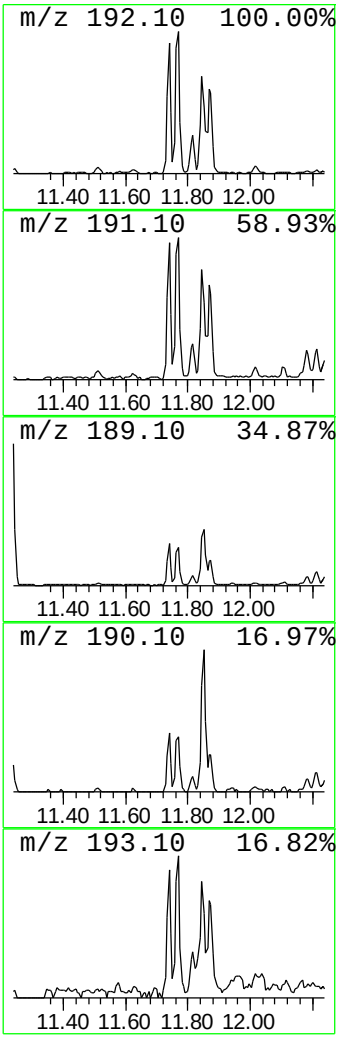
Peak Number 6 Anthracene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	3.44 ng	140169	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	90





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Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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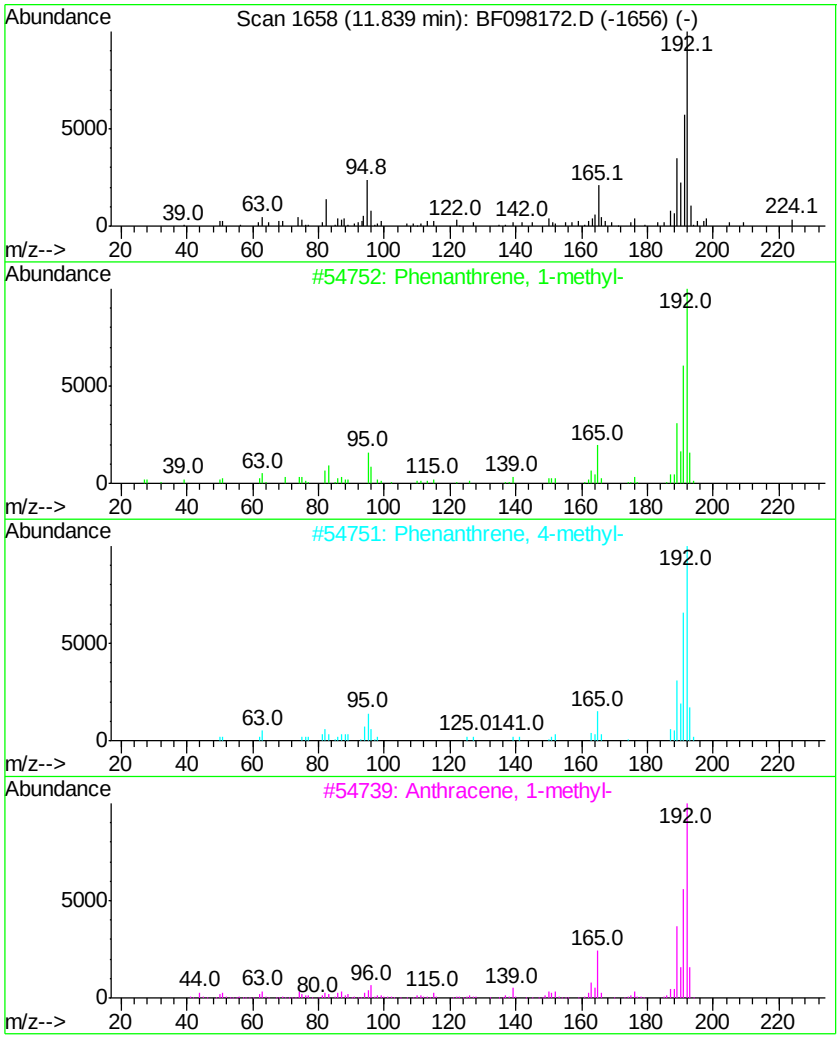
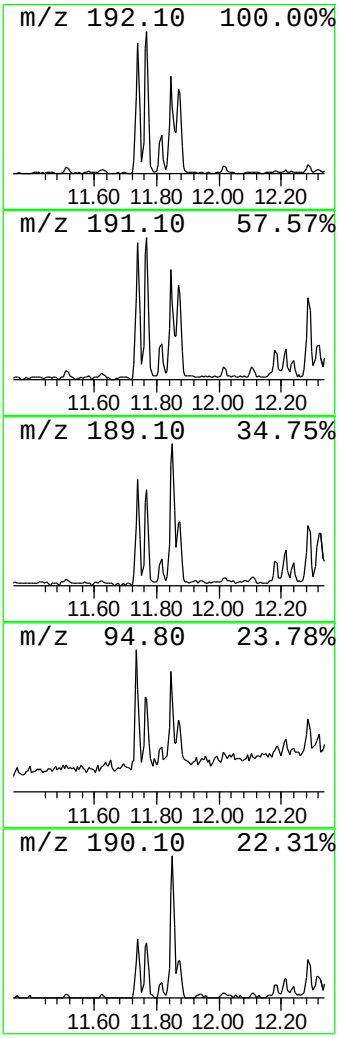
Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	4.51 ng	183732	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	78
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	70
4	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	64
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	64





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ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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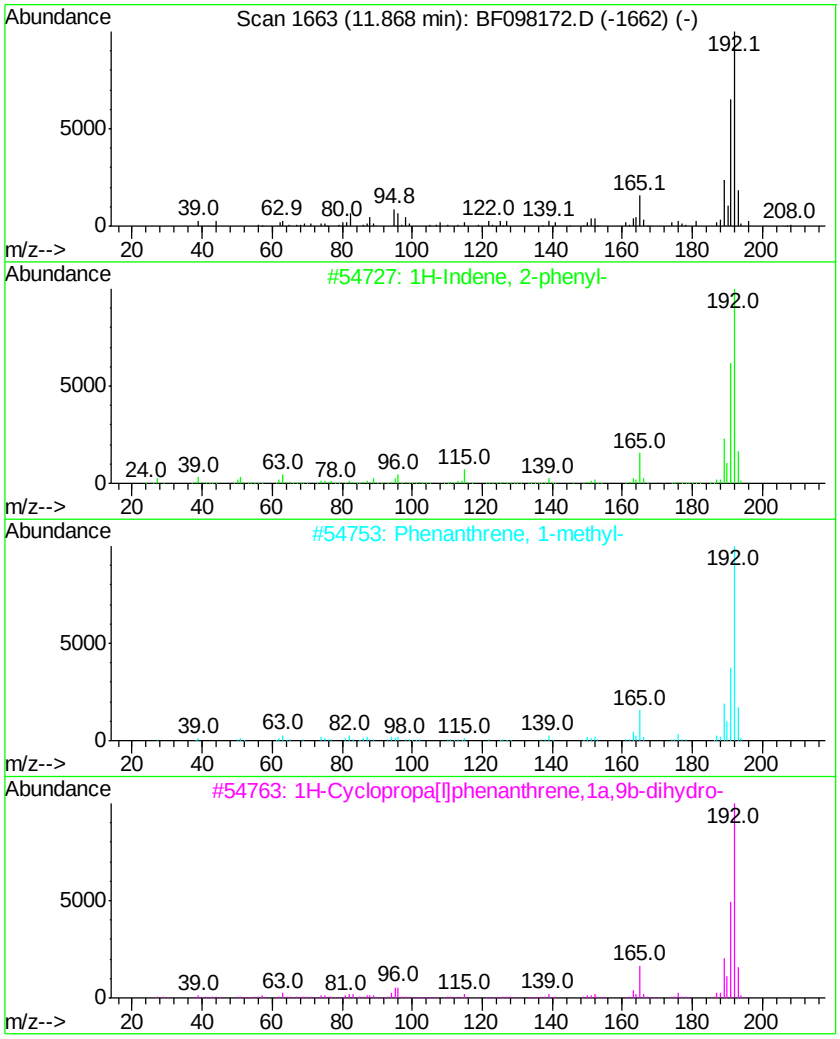
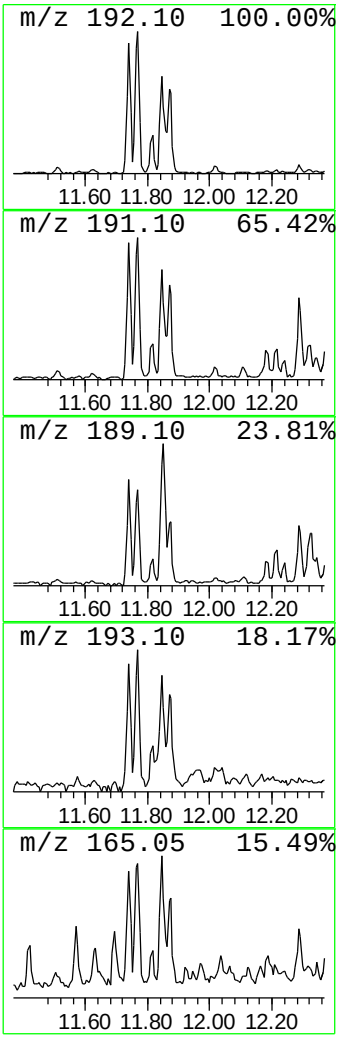
Peak Number 9 1H-Indene, 2-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	2.39 ng	97158	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	87





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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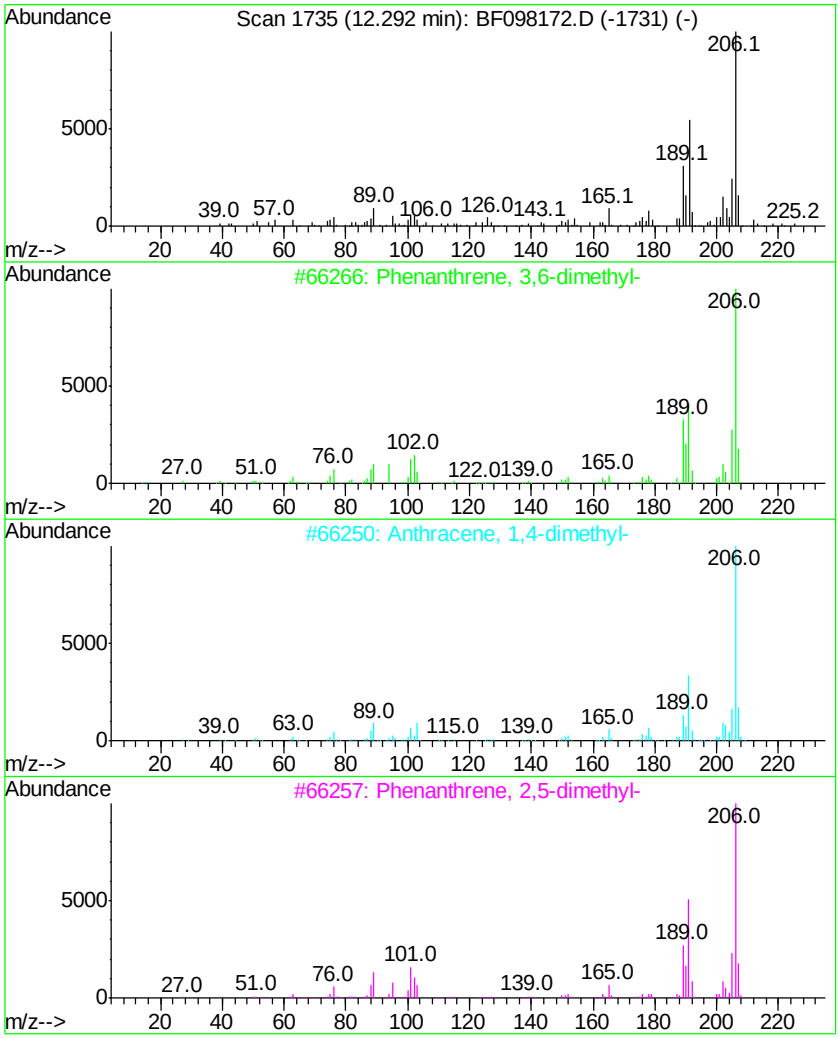
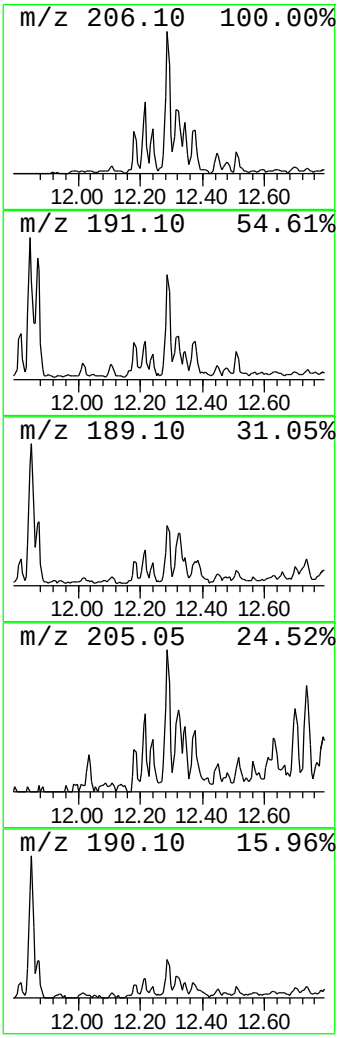
Peak Number 10 Phenanthrene, 3,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	3.46 ng	140994	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
2	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	94
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
4	9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91





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Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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

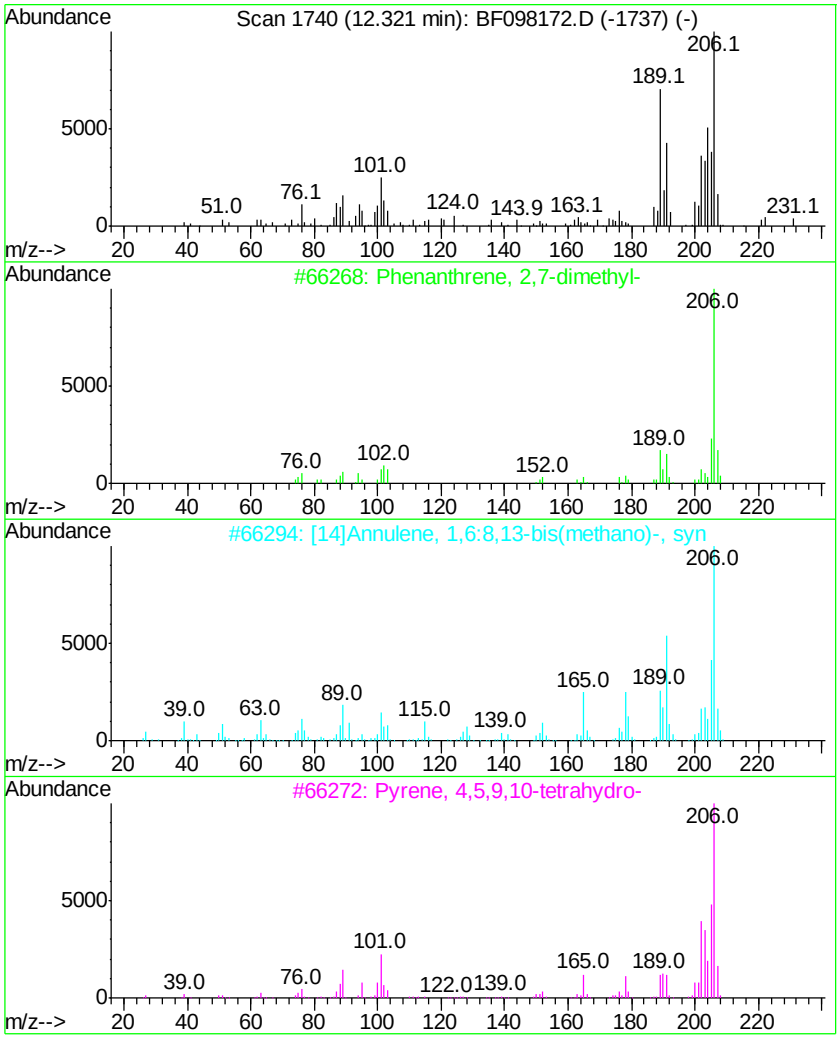
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

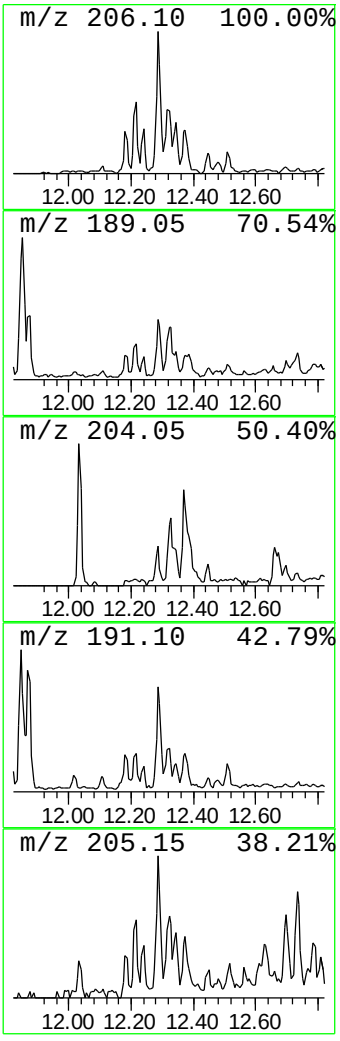
****

Peak Number 11 Phenanthrene, 2,7-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	2.06 ng	83954	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	64
2	[14]Annulene, 1,6:8,13-bis(metha...	206	C16H14	085385-68-8	64
3	Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	64
4	Benzene, 1,1'-(1,3-butadienylide...	206	C16H14	004165-81-5	60
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	55





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Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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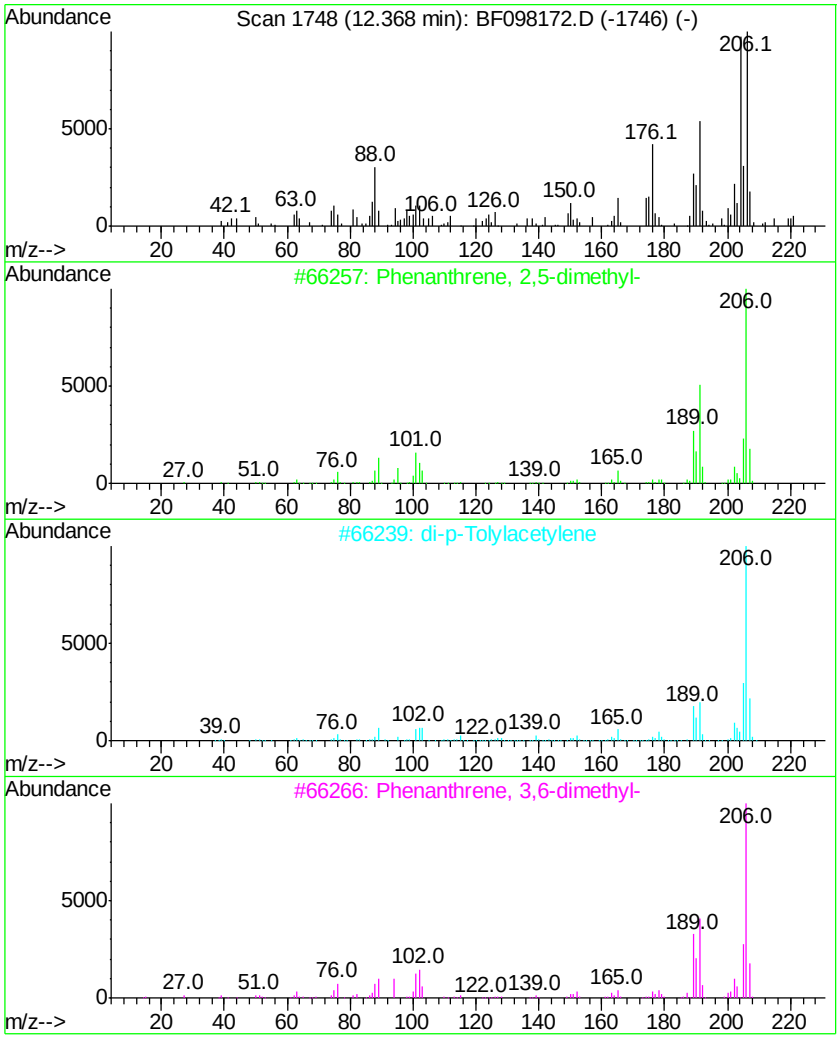
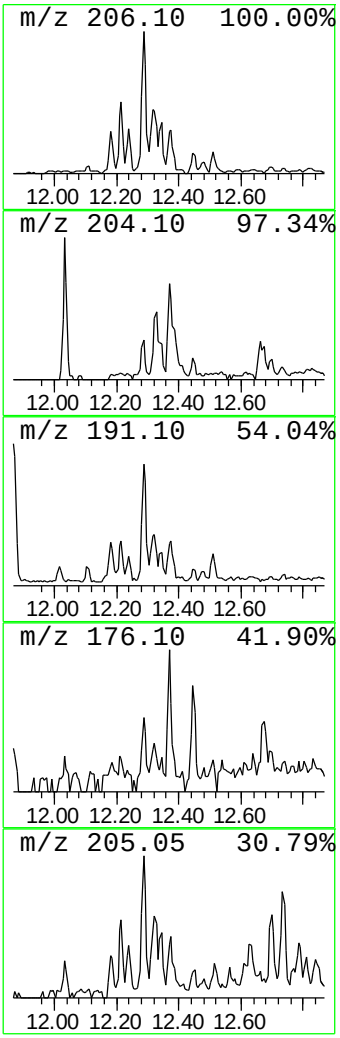
Peak Number 12 Phenanthrene, 2,5-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.37	2.09 ng	85147	Phenanthrene-d10	11.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	64
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	62
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	53
4	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	52
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	50





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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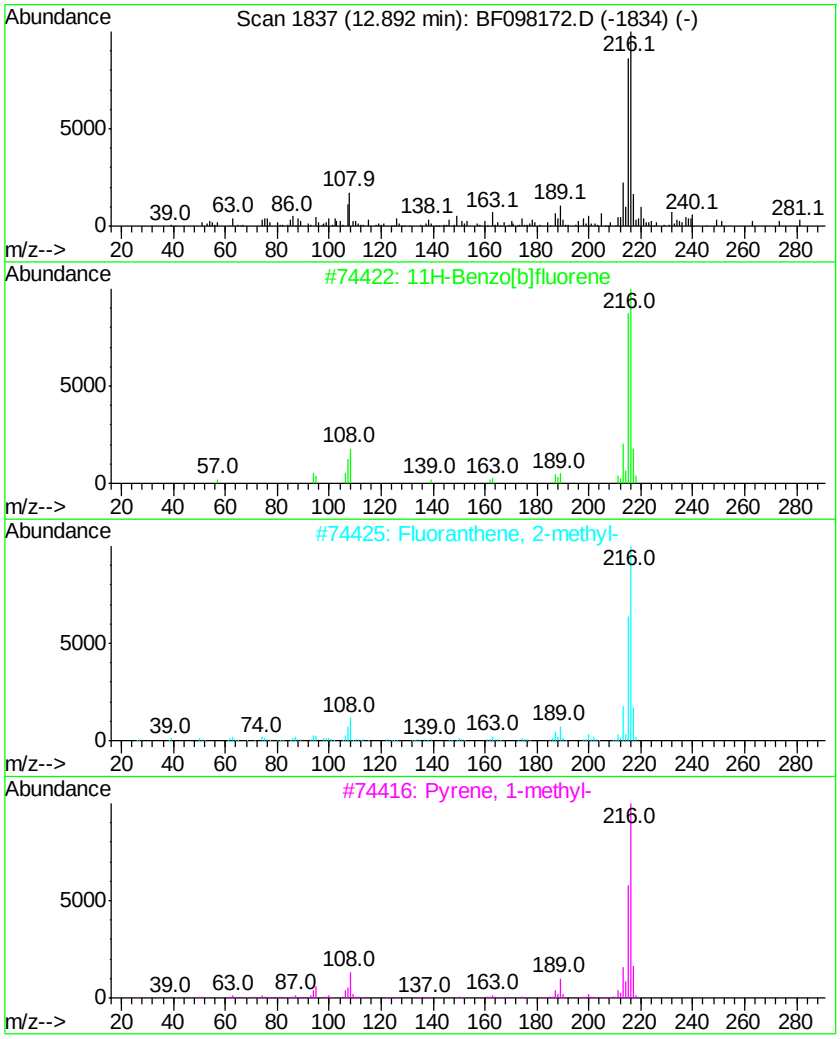
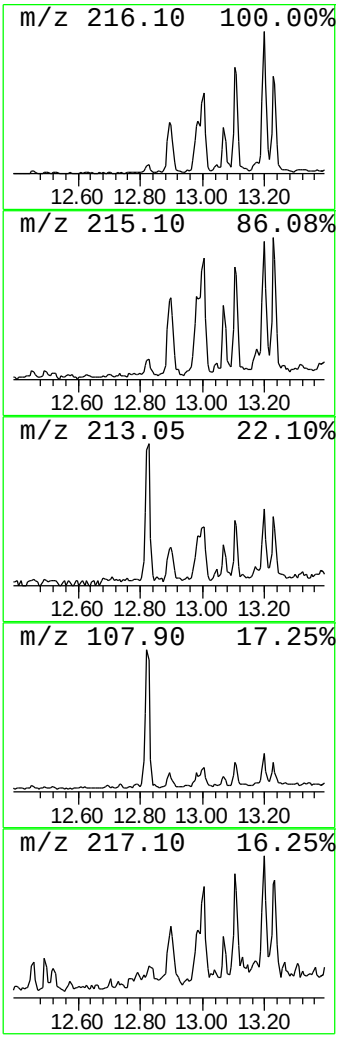
Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	2.16 ng	92450	Chrysene-d12	13.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	89
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	89
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	70





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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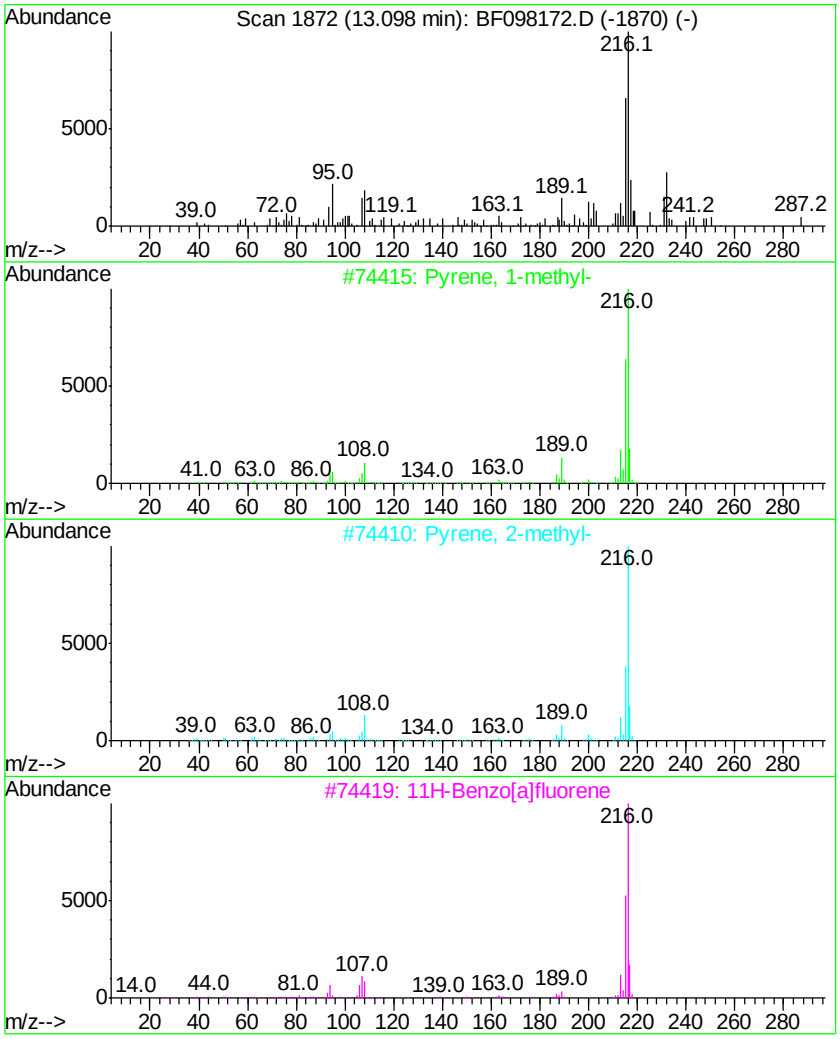
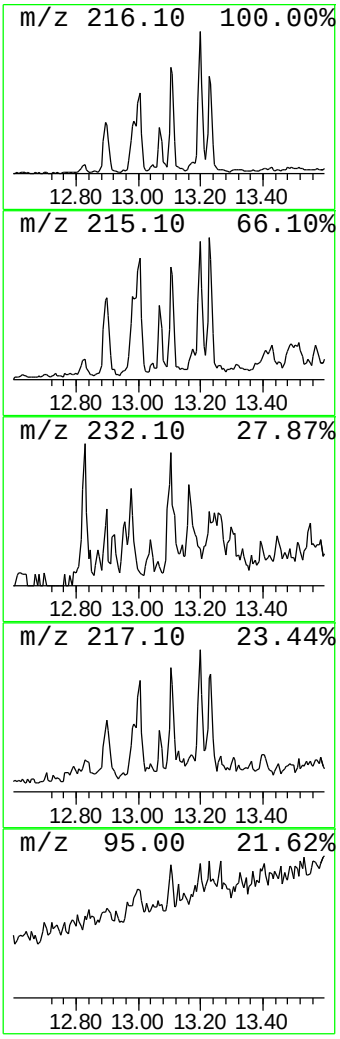
Peak Number 14 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.10	2.42 ng	103459	Chrysene-d12	13.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4	Pyrene, 4-methyl-	216	C17H12	003353-12-6	81
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	78





Data Path : Z:\HPCHEM1\BNA F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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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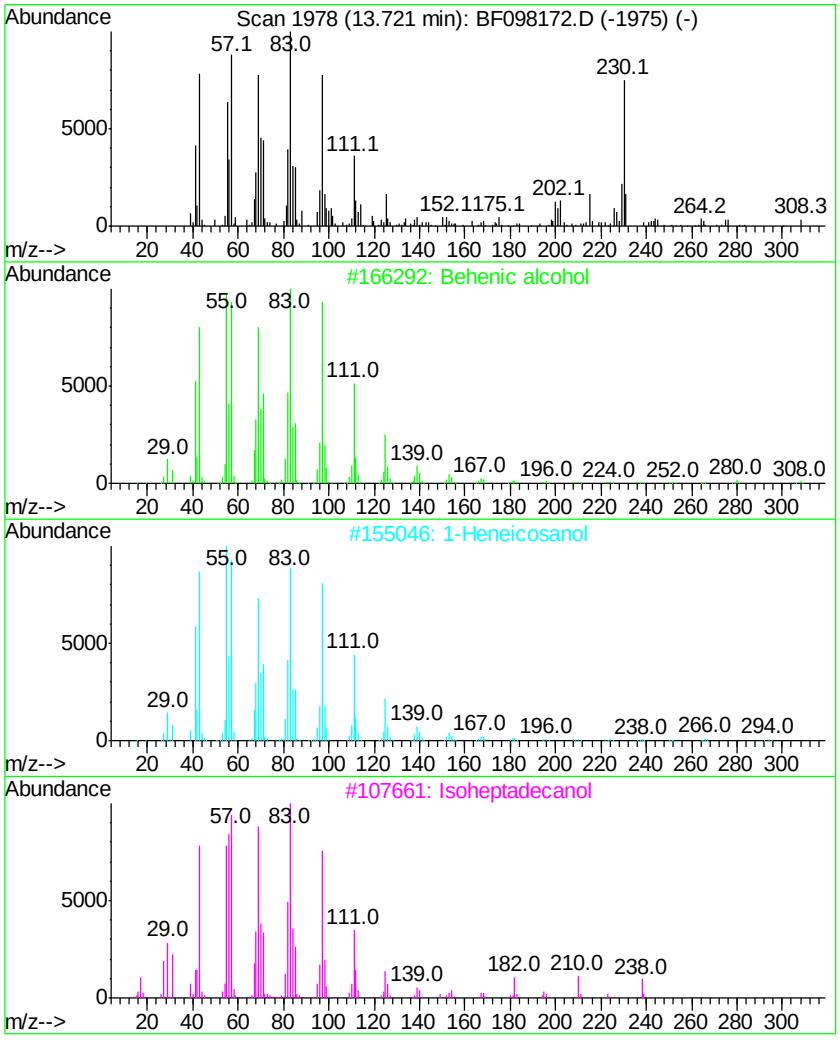
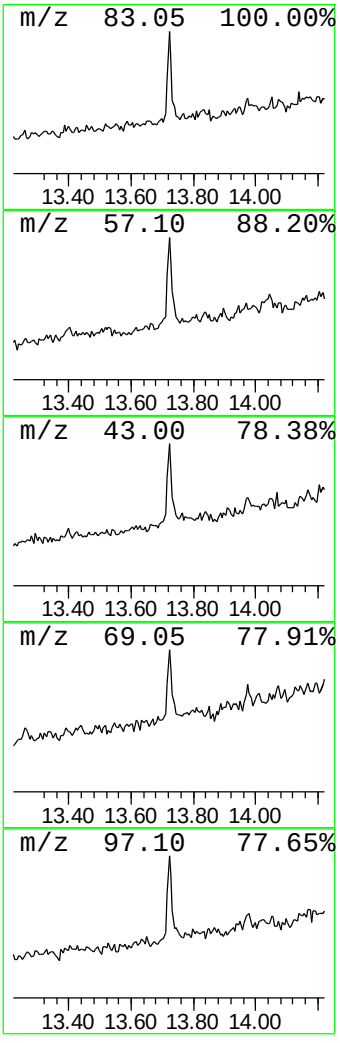
Peak Number 15 Behenic alcohol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	2.98 ng	127453	Chrysene-d12	13.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	76
2	1-Heneicosanol	312	C21H44O	015594-90-8	76
3	Isoheptadecanol	256	C17H36O	057289-07-3	76
4	n-Pentadecanol	228	C15H32O	000629-76-5	70
5	1-Heptacosanol	396	C27H56O	002004-39-9	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF083017\
Data File : BF098172.D
Acq On : 30 Aug 2017 22:54
Operator : SJ/JU
Sample : I5024-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CL-01-082917-A

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown3.09	3.09	2.4	ng	87179	1	6.72	717027	20.0
2-Pentanone, 4-hy...	4.93	8.4	ng	302830	1	6.72	717027	20.0
unknown6.50	6.50	91.4	ng	3278100	1	6.72	717027	20.0
unknown10.67	10.67	2.3	ng	91652	4	11.24	814346	20.0
Anthracene, 2-met...	11.74	3.4	ng	140169	4	11.24	814346	20.0
Phenanthrene, 1-m...	11.84	4.5	ng	183732	4	11.24	814346	20.0
1H-Indene, 2-phenyl-	11.87	2.4	ng	97158	4	11.24	814346	20.0
Phenanthrene, 3,6...	12.29	3.5	ng	140994	4	11.24	814346	20.0
Phenanthrene, 2,7...	12.32	2.1	ng	83954	4	11.24	814346	20.0
Phenanthrene, 2,5...	12.37	2.1	ng	85147	4	11.24	814346	20.0
11H-Benzo[b]fluorene	12.89	2.2	ng	92450	5	13.87	856558	20.0
Pyrene, 1-methyl-	13.10	2.4	ng	103459	5	13.87	856558	20.0
Behenic alcohol	13.72	3.0	ng	127453	5	13.87	856558	20.0