LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

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-BF043015.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	1.347	12	14	17	rBV	1485665	1919178	37.05%	5.414%
2	1.484	24	26	39	rVB2	118672	356556	6.88%	1.006%
3	1.655	39	41	43	rVV	77444	116750	2.25%	0.329%
4	1.690	43	44	53	rVV	173170	365286	7.05%	1.030%
5	1.815	53	55	96	rVB	2611737	5180500	100.00%	14.614%
6	4.993	330	333	339	rVB	89650	114783	2.22%	0.324%
7	5.507	375	378	381	rBV	464747	461501	8.91%	1.302%
8	6.787	487	490	492	rBV	339682	474897	9.17%	1.340%
9	6.925	500	502	505	rBV	505920	506487	9.78%	1.429%
10	7.210	525	527	530	rBV	193258	262884	5.07%	0.742%
11	7.405	541	544	546	rBV	390623	387522	7.48%	1.093%
12	7.907	586	588	590	rBV	347637	299683	5.78%	0.845%
13	8.788	663	665	667	rBV	285552	371142	7.16%	1.047%
14	10.136	781	783	786	rBV	670991	612863	11.83%	1.729%
15	10.788	838	840	843	rVB	110847	121671	2.35%	0.343%
16	10.971	854	856	858	rBV	374196	407336	7.86%	1.149%
17	11.005	858	859	862	rVB	72121	59754	1.15%	0.169%
18	11.645	913	915	918	rBV	65478	64580	1.25%	0.182%
19	11.942	939	941	944	rVB2	330614	398179	7.69%	1.123%
20	12.685	1004	1006	1010	rVB	32227	56860	1.10%	0.160%
21	12.811	1015	1017	1018	rBV	462615	446396	8.62%	1.259%
22	12.845	1018	1020	1022	rVV	719618	966958	18.67%	2.728%
23	12.902	1022	1025	1027	rVV	289966	278116	5.37%	0.785%
24	13.108	1041	1043	1045	rBV	74025	63185	1.22%	0.178%
25	13.439	1069	1072	1073	rBV	129068	122706	2.37%	0.346%
26	13.474	1073	1075	1078	rVB	183151	180981	3.49%	0.511%
27	13.531	1078	1080	1082	rBV	84683	89132	1.72%	0.251%
28	13.577	1082	1084	1085	rVV	193759	261278	5.04%	0.737%
29	13.599	1085	1086	1089	rVB	83766	92829	1.79%	0.262%
30	13.817	1103	1105	1109	rVB2	104752	200468	3.87%	0.566%
31	14.125	1128	1132	1134	rBV	97106	127447	2.46%	0.360%
32	14.171	1134	1136	1139	rVV2	45518	71277	1.38%	0.201%
33	14.228	1139	1141	1146	rVB2	88218	134161	2.59%	0.378%
34	14.320	1146	1149	1152	rBV	1911455	1949409	37.63%	5.499%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

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-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.400	1153	1156	1157	rBV	43576	58622	1.13%	0.165%
36	14.434	1157	1159	1161	rVB	89023	93584	1.81%	0.264%
37	14.502	1161	1165	1166	rBV2	87808	123611	2.39%	0.349%
38	14.594	1170	1173	1176	rBV	1307175	2032848	39.24%	5.735%
39	14.662	1178	1179	1184	rVB3	60802	103189	1.99%	0.291%
40	14.754	1185	1187	1189	rBV	95789	114312	2.21%	0.322%
41	14.811	1189	1192	1194	rVB2	575198	867735	16.75%	2.448%
42	14.880	1194	1198	1200	rBV2	86412	164705	3.18%	0.465%
43	14.982	1203	1207	1208	rBV3	101370	172171	3.32%	0.486%
44	15.017	1208	1210	1212	rVB	204823	235735	4.55%	0.665%
45	15.097	1215	1217	1219	rVB	111354	119183	2.30%	0.336%
46	15.142	1219	1221	1223	rBV	215787	247241	4.77%	0.697%
47	15.257	1229	1231	1233	rBV	103849	106487	2.06%	0.300%
48	15.291	1233	1234	1236	rBV	87977	96457	1.86%	0.272%
49	15.394	1241	1243	1246	rBV3	38691	82384	1.59%	0.232%
50	15.508	1249	1253	1254	rBV4	34733	72291	1.40%	0.204%
51	15.645	1263	1265	1269	rVB	149963	201420	3.89%	0.568%
52	15.783	1275	1277	1279	rBV2	160256	251115	4.85%	0.708%
53	15.828	1279	1281	1284	rVV2	158579	346493	6.69%	0.977%
54	15.874	1284	1285	1287	rVV	100279	133735	2.58%	0.377%
55	15.920	1287	1289	1291	rVB	145897	196723	3.80%	0.555%
56	15.988	1293	1295	1299	rVV3	79564	177370	3.42%	0.500%
57	16.103	1301	1305	1307	rVV2	1164571	2049743	39.57%	5.782%
58	16.148	1307	1309	1311	rVV	956417	1125799	21.73%	3.176%
59	16.240	1314	1317	1319	rBV2	163489	216240	4.17%	0.610%
60	16.297	1319	1322	1323	rVV2	54980	80184	1.55%	0.226%
61	16.365	1326	1328	1331	rVV3	64900	109007	2.10%	0.308%
62	16.411	1331	1332	1334	rVV	81867	94039	1.82%	0.265%
63	16.537	1340	1343	1345	rBV3	33988	83991	1.62%	0.237%
64	16.571	1345	1346	1348	rVV	44126	54358	1.05%	0.153%
65	16.617	1348	1350	1352	rVB	133526	189591	3.66%	0.535%
66	16.663	1352	1354	1357	rBV	81724	101658	1.96%	0.287%
67	16.743	1359	1361	1363	rBV	68084	72489	1.40%	0.204%
68	16.777	1363	1364	1365	rBV	72732	68706	1.33%	0.194%
69	16.891	1372	1374	1376	rVB	51542	63704	1.23%	0.180%
70	17.028	1383	1386	1388	rBV	59354	118459	2.29%	0.334%
71	17.234	1402	1404	1407	rBV2	52221	123154	2.38%	0.347%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

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-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	17.440	1419	1422	1428	rBV	996082	2194378	42.36%	6.190%
73	17.566	1431	1433	1435	rVB	182162	246042	4.75%	0.694%
74	17.771	1448	1451	1454	rVB	663176	831736	16.06%	2.346%
75	17.840	1454	1457	1459	rBV	837853	1131580	21.84%	3.192%
76	17.897	1460	1462	1464	rBV	405627	649823	12.54%	1.833%
77	19.177	1572	1574	1577	rVB2	88713	181108	3.50%	0.511%
78	19.337	1585	1588	1593	rVB2	419598	860996	16.62%	2.429%
79	19.520	1601	1604	1606	rBV	89257	193646	3.74%	0.546%
80	19.566	1606	1608	1612	rVV2	63421	141544	2.73%	0.399%
81	19.760	1622	1625	1631	rVB	352207	699719	13.51%	1.974%
82	19.989	1642	1645	1652	rVB	102523	246619	4.76%	0.696%

Sum of corrected areas: 35448479

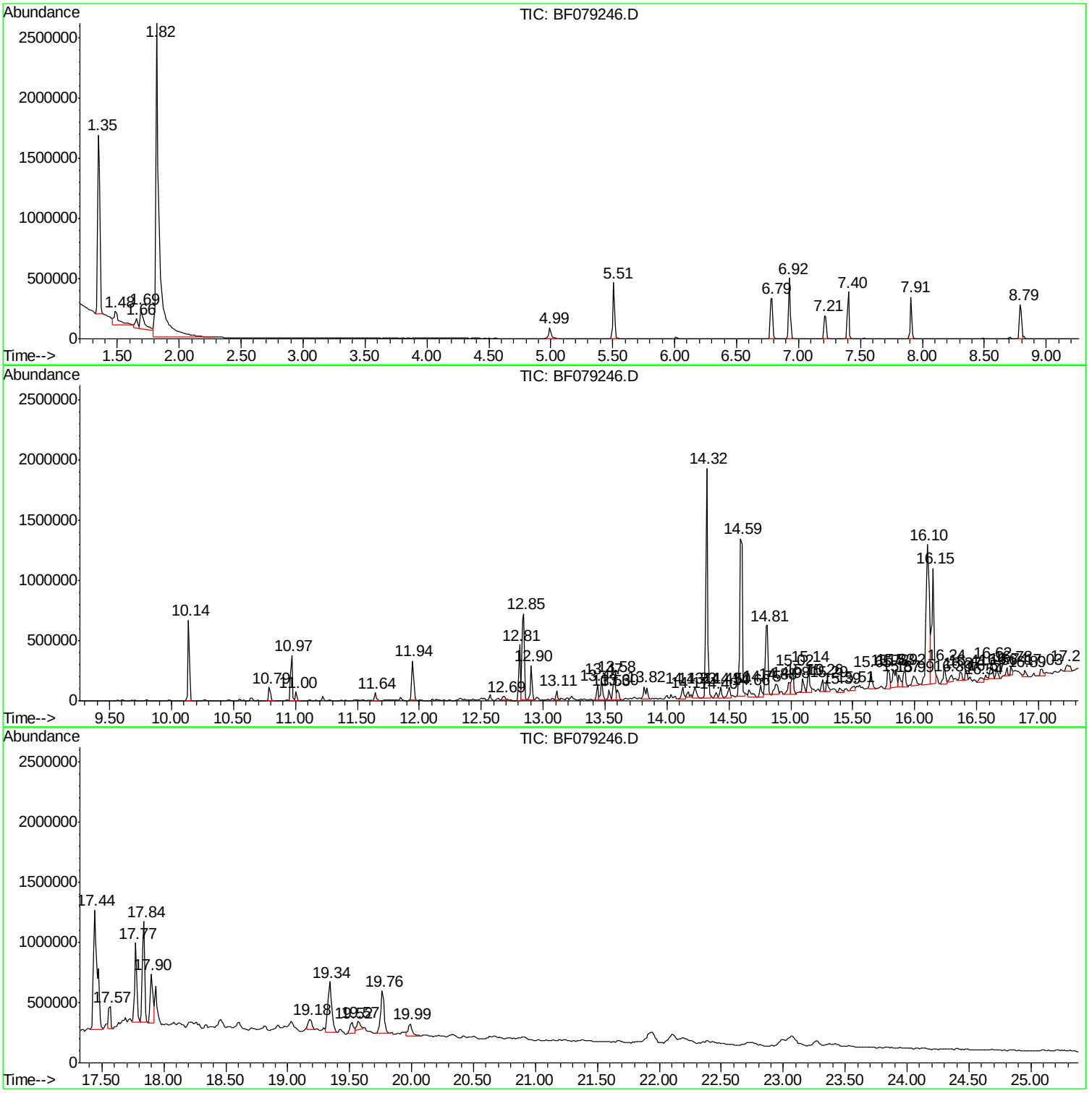
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

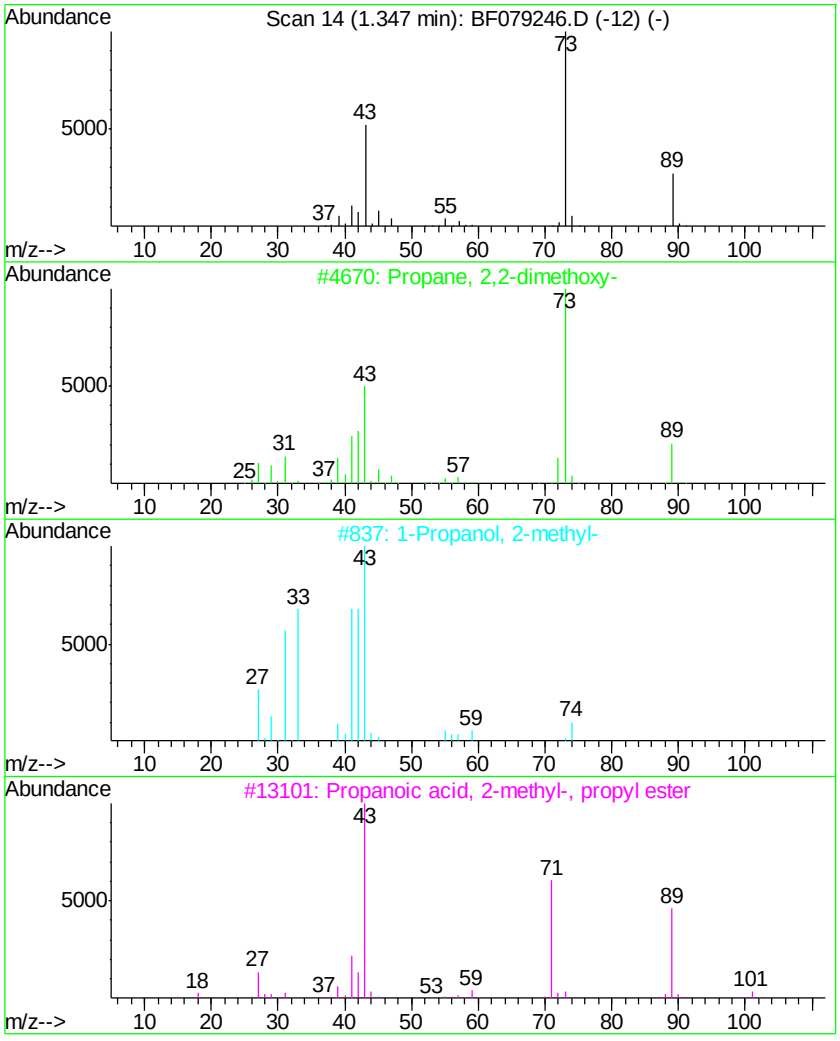
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

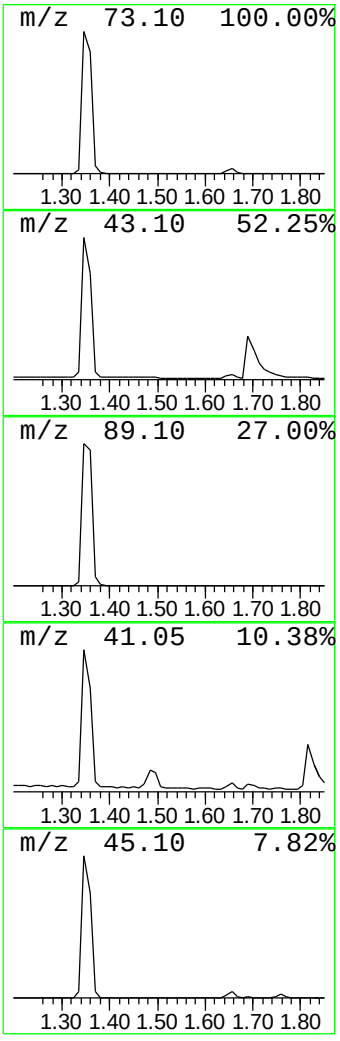
****

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.35	146.01 ng	1919180	1,4-Dichlorobenzene-d4	7.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2	1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
4	Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5	2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

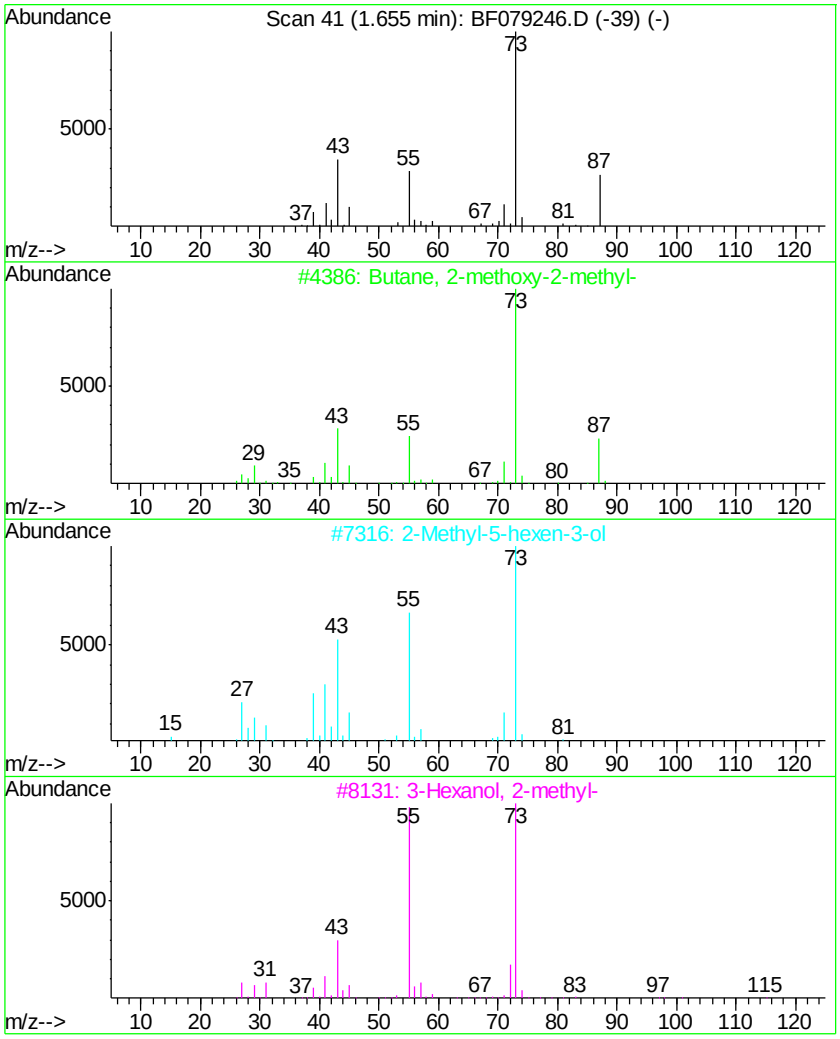
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

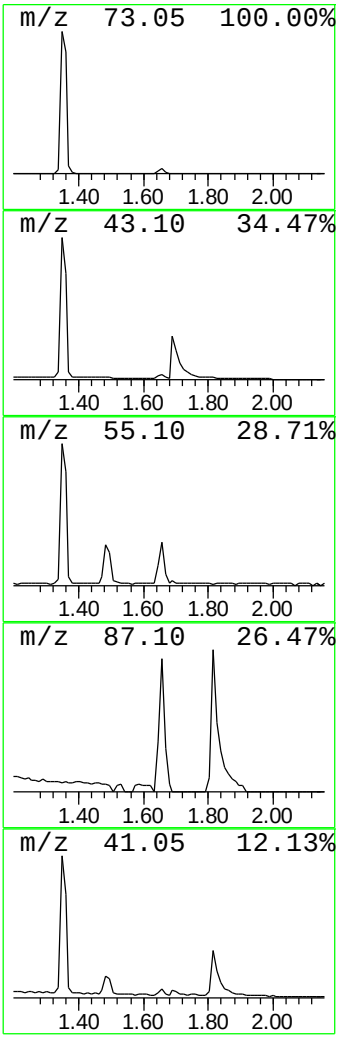
****

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.66	8.88 ng	116750	1,4-Dichlorobenzene-d4	7.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
3	3-Hexanol, 2-methyl-	116	C7H16O	000617-29-8	17
4	Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

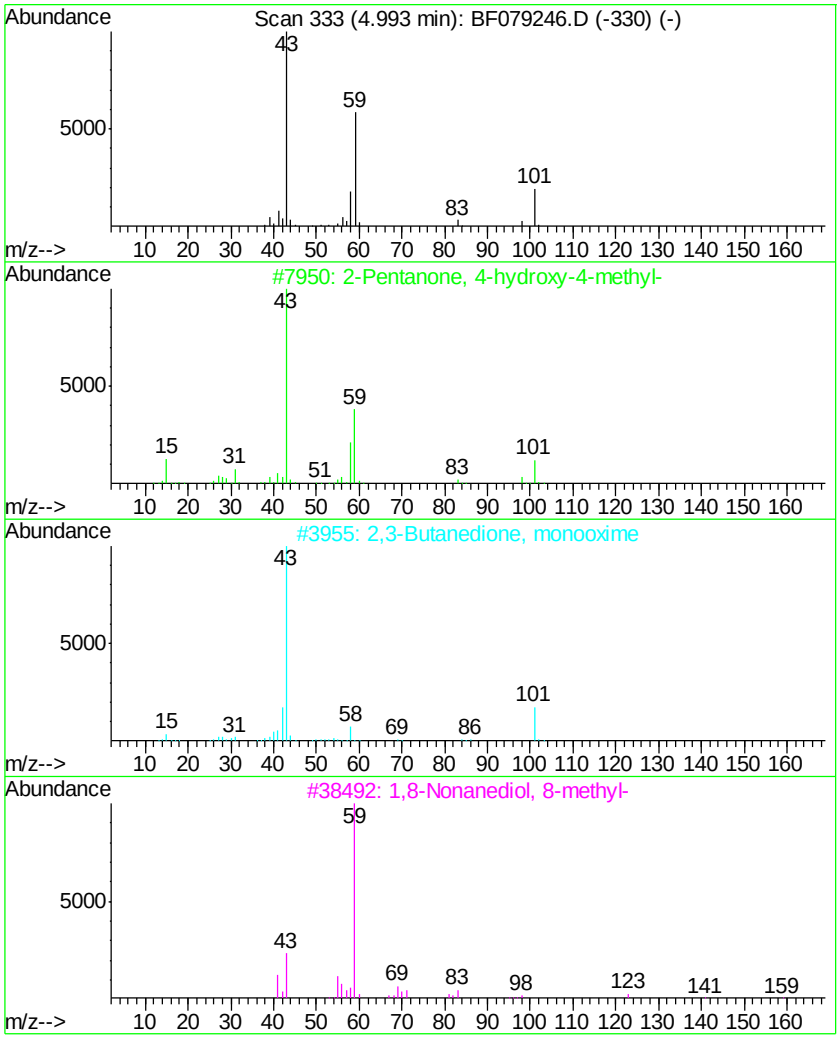
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

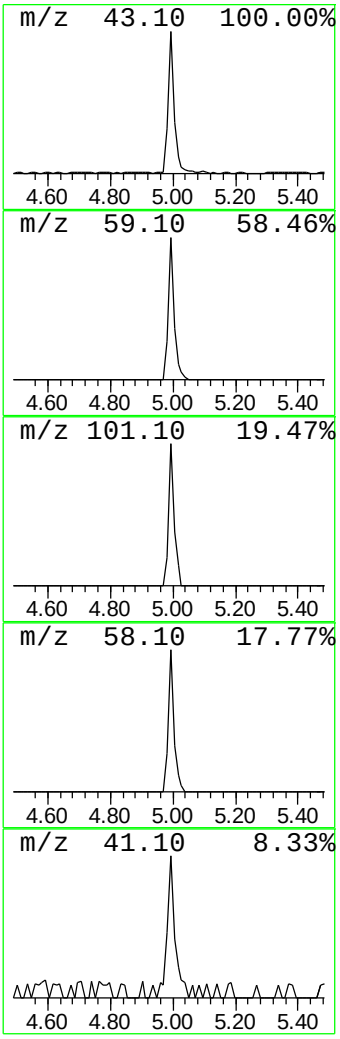
****

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	8.73 ng	114783	1,4-Dichlorobenzene-d4	7.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32
3	1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	23
4	5-Hexen-2-one	98	C6H10O	000109-49-9	10
5	Acetone	58	C3H6O	000067-64-1	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
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Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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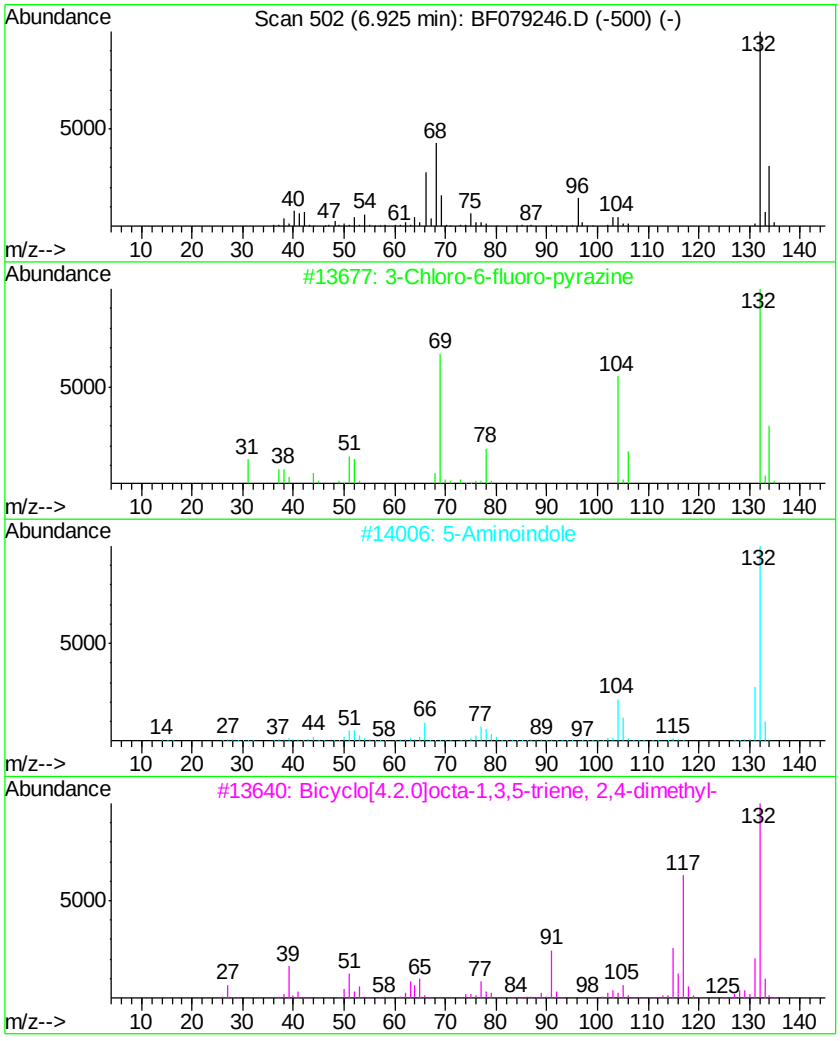
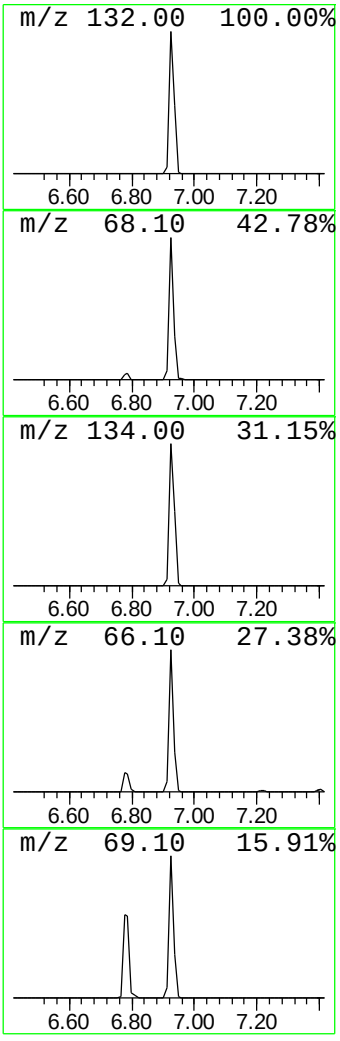
Peak Number 7 unknown6.92 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.92	38.53 ng	506487	1,4-Dichlorobenzene-d4	7.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

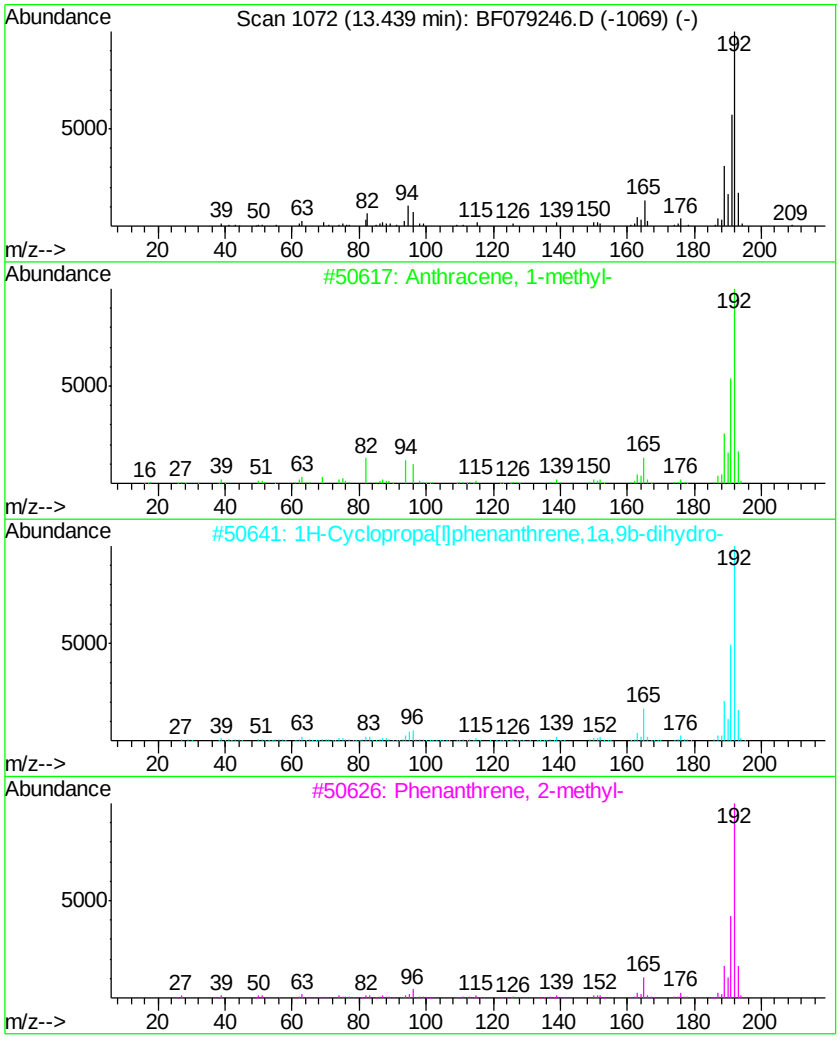
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

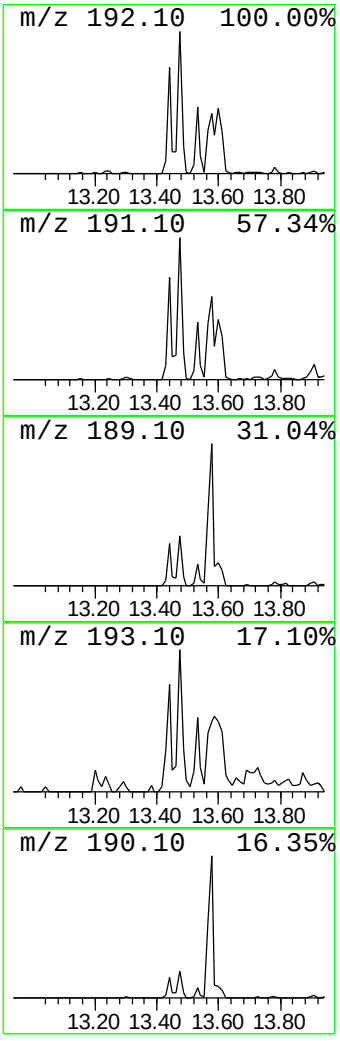
****

Peak Number 9 Anthracene, 1-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	5.50 ng	122706	Phenanthrene-d10	12.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	97
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96





Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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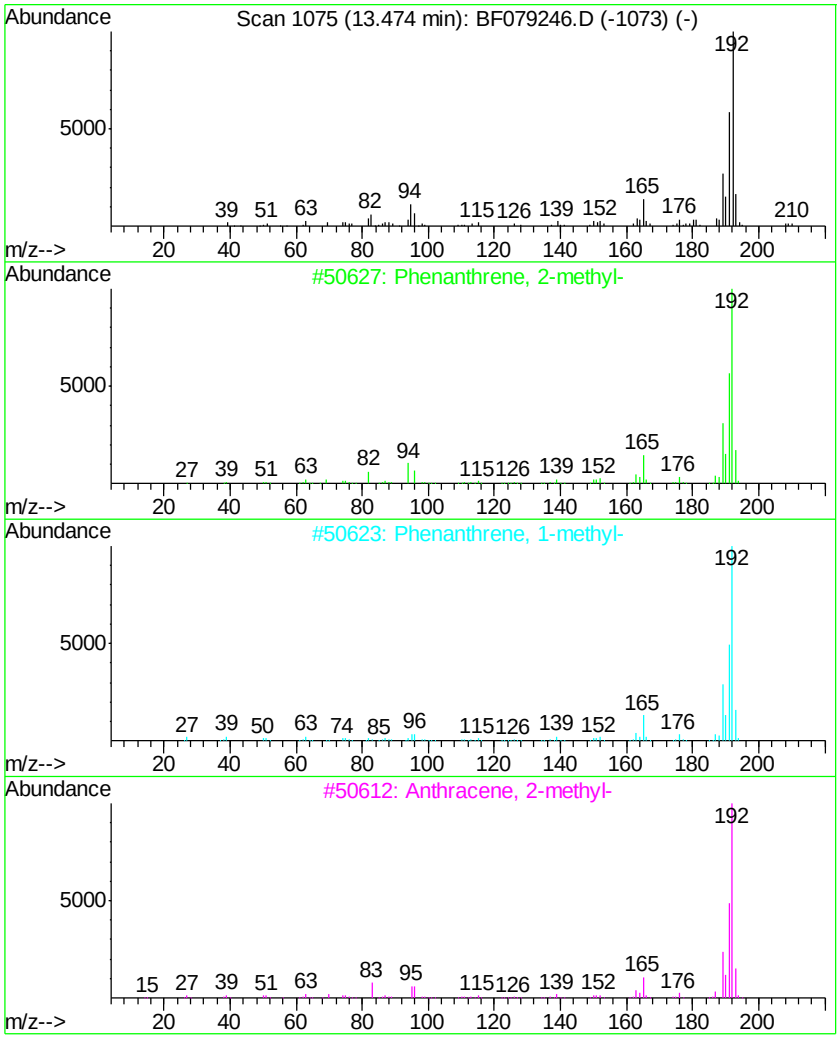
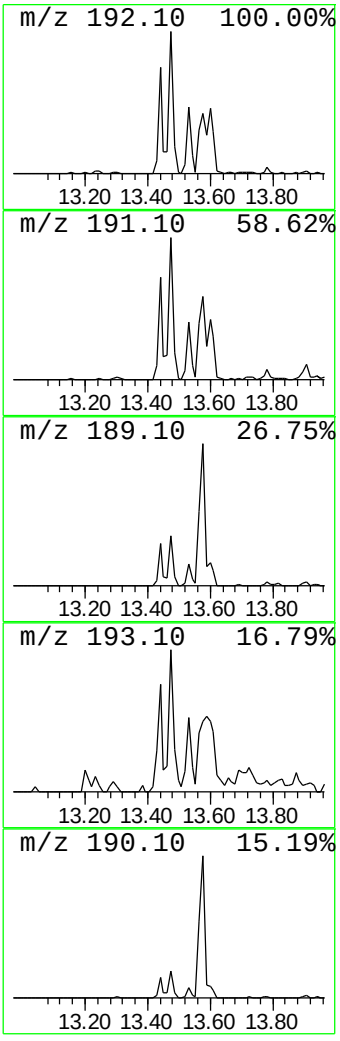
Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	8.11 ng	180981	Phenanthrene-d10	12.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95





Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
Data File : BF079246.D
Acq On : 14 May 2015 22:07
Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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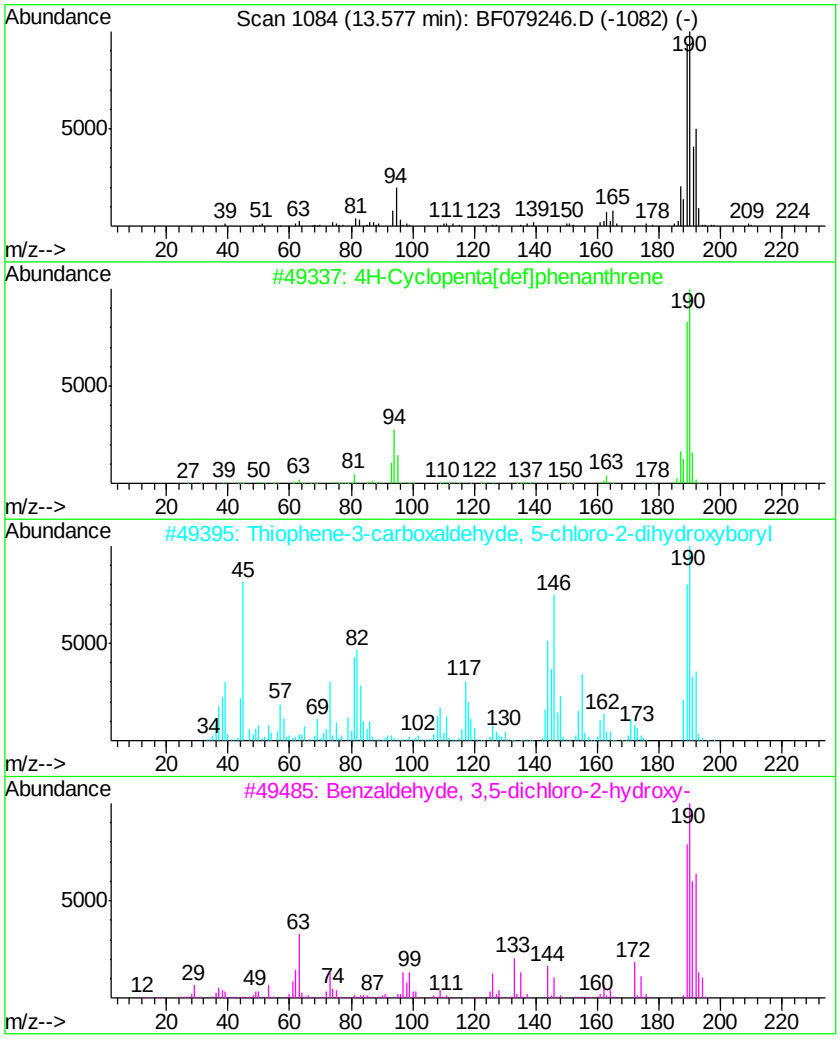
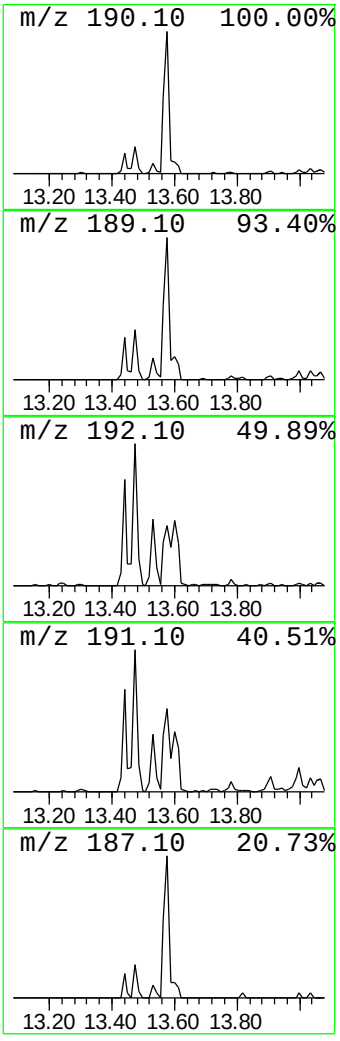
Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.58	11.71 ng	261278	Phenanthrene-d10	12.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
4	Methyl diselenide	190	C2H6Se2	007101-31-7	45
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38





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Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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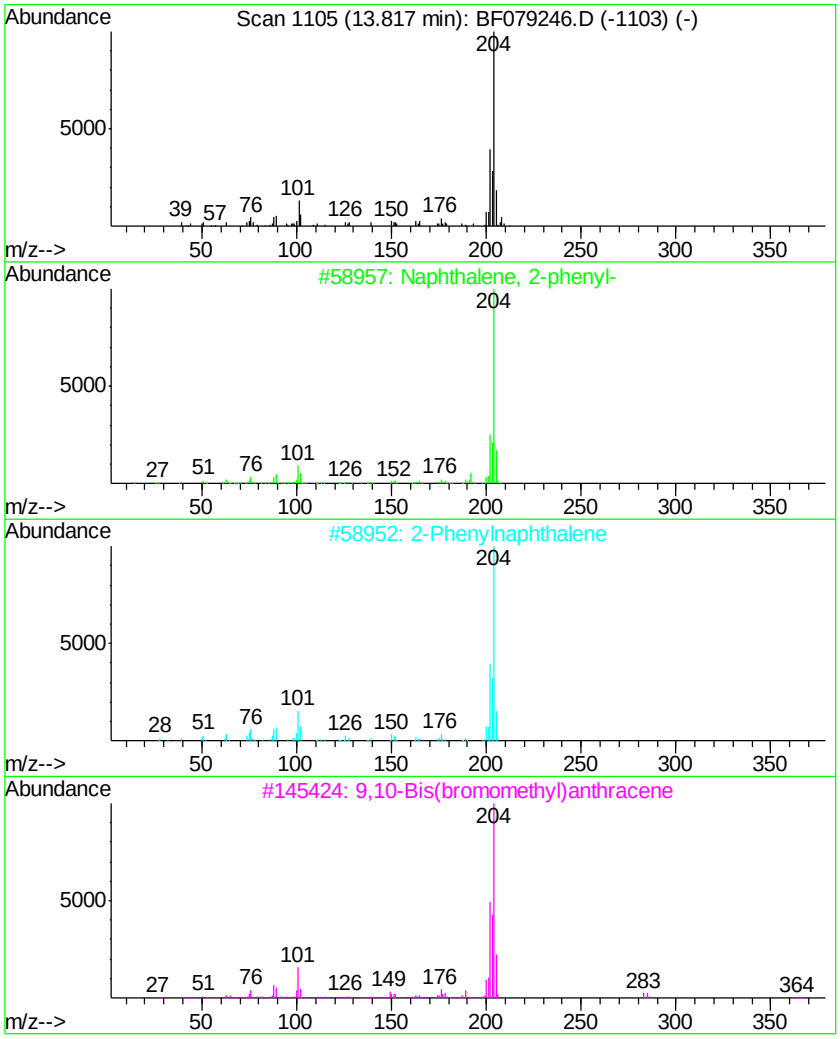
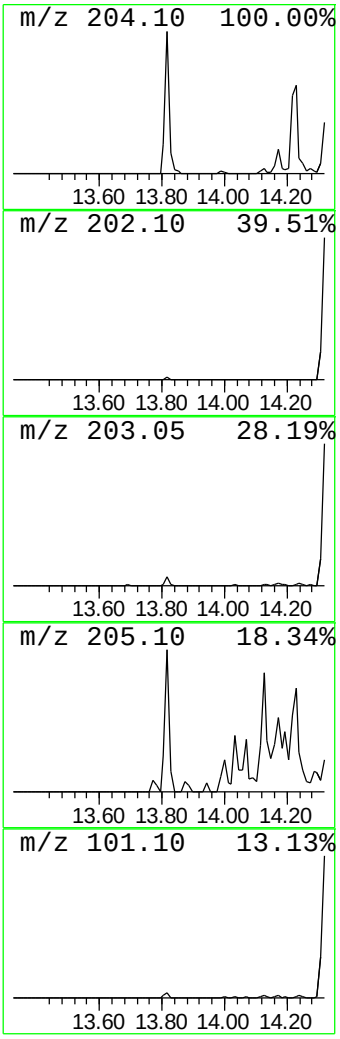
Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	8.98 ng	200468	Phenanthrene-d10	12.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	89
2	2-Phenylnaphthalene	204	C16H12	035465-71-5	86
3	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	76
4	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	72
5	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64





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Data File : BF079246.D
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Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

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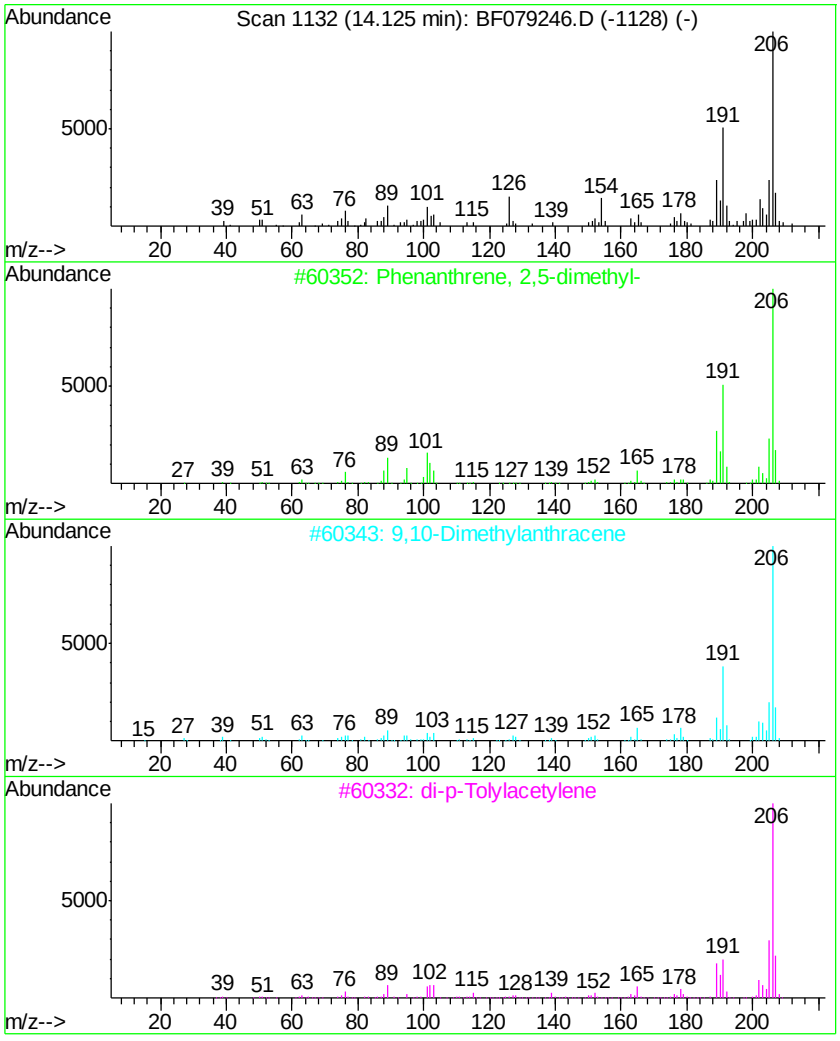
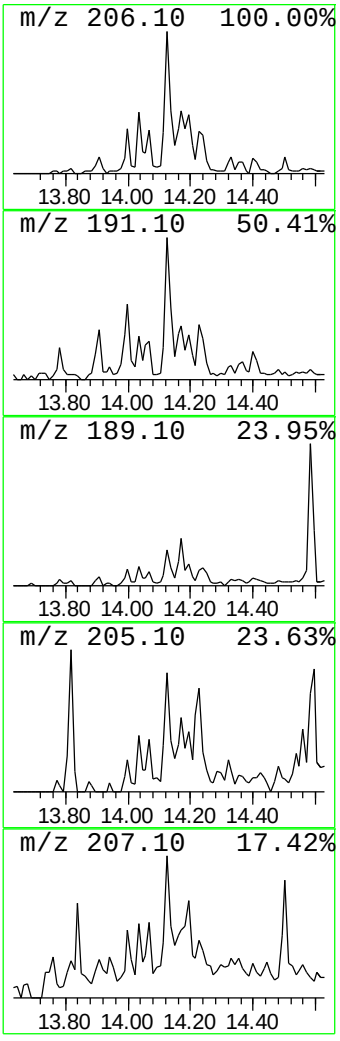
Peak Number 13 Phenanthrene, 2,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	5.71 ng	127447	Phenanthrene-d10	12.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2	9,10-Dimethylanthracene	206	C16H14	000781-43-1	94
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	91
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91





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Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

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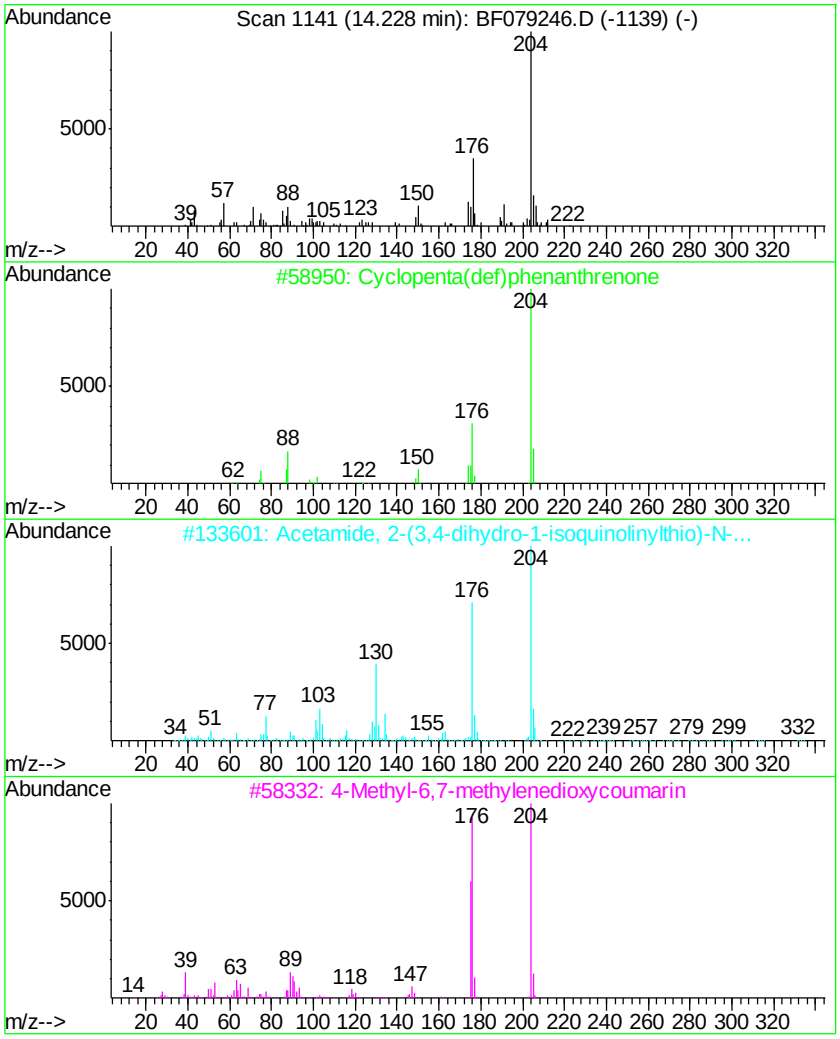
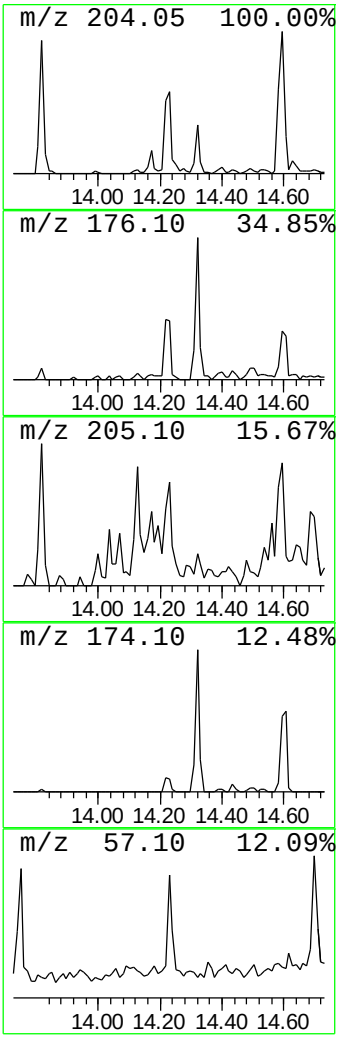
Peak Number 14 Cyclopenta(def)phenanthrenone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	6.01 ng	134161	Phenanthrene-d10	12.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	87
2	Acetamide, 2-(3,4-dihydro-1-isoq...	332	C17H14F2N2OS	1000270-76-1	53
3	4-Methyl-6,7-methylenedioxycoumarin	204	C11H8O4	015071-04-2	52
4	3'-Carboxybenzo[1',2',-b]-1,4-di...	204	C11H12N2O2	138023-41-3	52
5	9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	47





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Operator : TP/IZ
Sample : G2215-09 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-38S

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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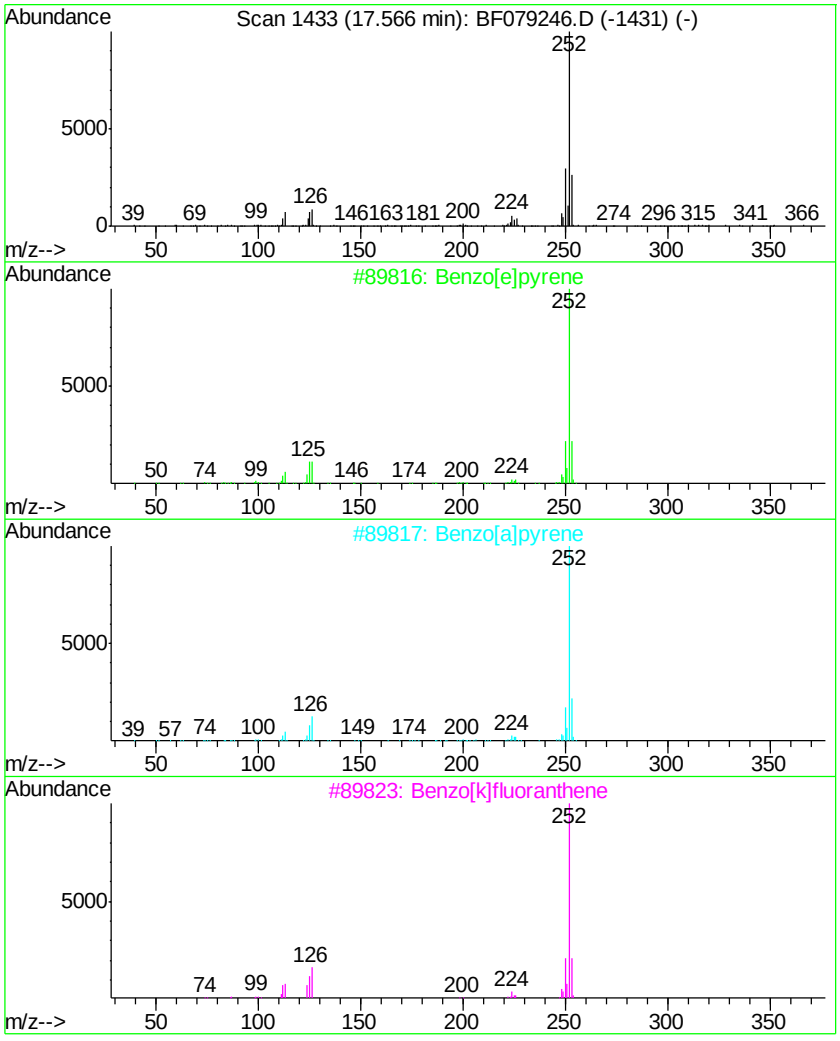
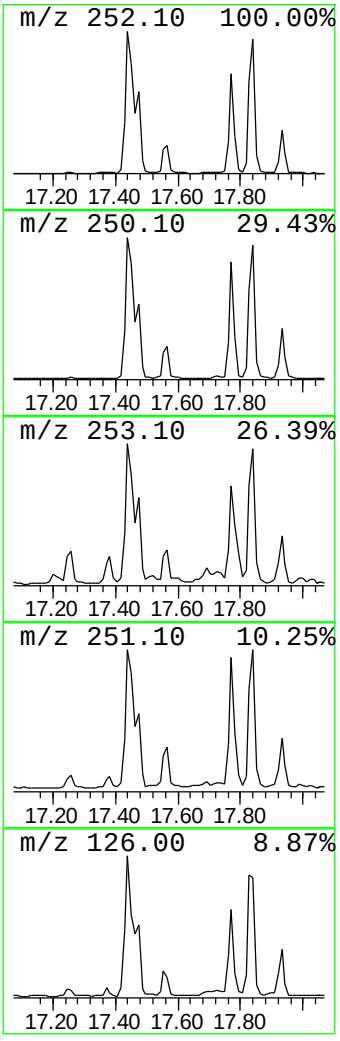
Peak Number 15 Benzo[e]pyrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.57	7.57 ng	246042	Perylene-d12	17.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	91
2	Benzo[a]pyrene	252	C20H12	000050-32-8	90
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
4	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	81
5	Perylene	252	C20H12	000198-55-0	74





Tentatively Identified Compound (LSC) summary

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Propane, 2,2-dime...	1.35	146.0	ng	1919180	1	7.22	262884	20.0
Butane, 2-methoxy...	1.66	8.9	ng	116750	1	7.22	262884	20.0
2-Pentanone, 4-hy...	4.99	8.7	ng	114783	1	7.22	262884	20.0
unknown6.92	6.92	38.5	ng	506487	1	7.22	262884	20.0
Anthracene, 1-met...	13.44	5.5	ng	122706	4	12.81	446396	20.0
Phenanthrene, 2-m...	13.47	8.1	ng	180981	4	12.81	446396	20.0
4H-Cyclopenta[def...	13.58	11.7	ng	261278	4	12.81	446396	20.0
Naphthalene, 2-ph...	13.82	9.0	ng	200468	4	12.81	446396	20.0
Phenanthrene, 2,5...	14.13	5.7	ng	127447	4	12.81	446396	20.0
Cyclopenta(def)ph...	14.23	6.0	ng	134161	4	12.81	446396	20.0
Benzo[e]pyrene	17.57	7.6	ng	246042	6	17.90	649823	20.0