LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093222.D
Acq On : 27 Feb 2017 23:24
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.979	251	253	262	rBV	1067769	1714103	41.52%	3.249%
2	5.424	289	292	294	rBV	3830830	3990272	96.65%	7.563%
3	6.465	381	383	386	rBV	2748906	4126225	99.94%	7.820%
4	6.590	391	394	396	rBV	4026222	3910698	94.72%	7.412%
5	6.819	412	414	416	rBV	977612	1003970	24.32%	1.903%
6	6.967	425	427	430	rBV	2453492	2875897	69.66%	5.451%
7	7.390	461	464	466	rBV	2068660	2652698	64.25%	5.028%
8	8.099	524	526	530	rBV	1549252	1425447	34.53%	2.702%
9	8.819	586	589	591	rBV	56550	79038	1.91%	0.150%
10	8.911	595	597	601	rVB	138934	125828	3.05%	0.238%
11	9.185	617	621	623	rBV	2854125	4128580	100.00%	7.825%
12	9.436	641	643	645	rBV	49694	71887	1.74%	0.136%
13	9.516	648	650	651	rVV	120368	127294	3.08%	0.241%
14	9.573	653	655	658	rVV	1081459	1027895	24.90%	1.948%
15	9.631	658	660	665	rVB	70732	83544	2.02%	0.158%
16	9.711	665	667	670	rBV	52607	56513	1.37%	0.107%
17	9.791	672	674	677	rBV	35954	50845	1.23%	0.096%
18	9.859	677	680	681	rVV	1247014	1451322	35.15%	2.751%
19	9.882	681	682	684	rVB	129861	143039	3.46%	0.271%
20	10.008	691	693	695	rBV3	30766	41971	1.02%	0.080%
21	10.053	695	697	699	rVV	71551	101855	2.47%	0.193%
22	10.099	699	701	703	rVB	49673	49103	1.19%	0.093%
23	10.168	706	707	709	rBV	38002	41742	1.01%	0.079%
24	10.259	713	715	716	rBV	54148	64248	1.56%	0.122%
25	10.282	716	717	719	rVB	98228	75892	1.84%	0.144%
26	10.396	726	727	730	rVB	136443	168282	4.08%	0.319%
27	10.465	730	733	734	rBV	45617	68425	1.66%	0.130%
28	10.488	734	735	740	rVB3	41166	66982	1.62%	0.127%
29	10.556	740	741	743	rVB	57875	53687	1.30%	0.102%
30	10.648	746	749	751	rBV	2094734	2045310	49.54%	3.876%
31	10.762	757	759	763	rVB2	101515	155906	3.78%	0.295%
32	10.854	763	767	770	rBV2	26680	52884	1.28%	0.100%
33	10.945	772	775	777	rBV2	61757	116262	2.82%	0.220%
34	11.036	779	783	784	rVB3	26790	47545	1.15%	0.090%

LSC Area Percent Report

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Acq On : 27 Feb 2017 23:24
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.082	784	787	791	rBV4	31979	96590	2.34%	0.183%
36	11.151	791	793	795	rVB	75038	65160	1.58%	0.123%
37	11.208	795	798	799	rBV2	92258	148319	3.59%	0.281%
38	11.231	799	800	803	rVB	96035	126376	3.06%	0.240%
39	11.345	807	810	811	rBV	1165440	1646282	39.88%	3.120%
40	11.368	811	812	814	rVV	1475740	1057168	25.61%	2.004%
41	11.414	814	816	818	rVB	317567	294281	7.13%	0.558%
42	11.574	827	830	834	rBV2	131263	186582	4.52%	0.354%
43	11.631	834	835	837	rVV	96319	86136	2.09%	0.163%
44	11.676	837	839	841	rVB	51624	73726	1.79%	0.140%
45	11.756	841	846	848	rVB2	35723	74092	1.79%	0.140%
46	11.836	851	853	855	rBV	161185	246063	5.96%	0.466%
47	11.871	855	856	858	rVB	331126	249790	6.05%	0.473%
48	11.916	858	860	861	rBV2	118413	112305	2.72%	0.213%
49	11.951	861	863	868	rVB	348068	536049	12.98%	1.016%
50	12.042	868	871	873	rBV2	37961	82696	2.00%	0.157%
51	12.134	875	879	881	rBV	124684	174327	4.22%	0.330%
52	12.168	881	882	885	rVB	85734	69358	1.68%	0.131%
53	12.282	889	892	894	rBV2	59066	85736	2.08%	0.162%
54	12.397	899	902	903	rBV	108122	171198	4.15%	0.324%
55	12.431	903	905	908	rVB	98164	135026	3.27%	0.256%
56	12.477	908	909	913	rVB	93933	130810	3.17%	0.248%
57	12.557	913	916	918	rBV	1430420	1409419	34.14%	2.671%
58	12.614	918	921	923	rVB3	39789	86475	2.09%	0.164%
59	12.705	925	929	931	rBV	68570	104730	2.54%	0.198%
60	12.785	931	936	938	rBV	1353890	1572970	38.10%	2.981%
61	12.831	938	940	943	rVB2	66802	100743	2.44%	0.191%
62	12.922	946	948	951	rVB	2084048	2918373	70.69%	5.531%
63	13.002	952	955	959	rBV3	106110	200792	4.86%	0.381%
64	13.105	961	964	966	rVB	129088	265130	6.42%	0.502%
65	13.174	969	970	972	rBV	76297	78747	1.91%	0.149%
66	13.220	972	974	977	rVB2	116244	170884	4.14%	0.324%
67	13.311	980	982	983	rBV	77856	83373	2.02%	0.158%
68	13.334	983	984	986	rBV	89895	104420	2.53%	0.198%
69	13.494	996	998	1000	rBV3	37591	86490	2.09%	0.164%
70	13.620	1005	1009	1011	rBV	106415	139439	3.38%	0.264%
71	13.734	1016	1019	1020	rBV	99233	150510	3.65%	0.285%

LSC Area Percent Report

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Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	13.757	1020	1021	1022	rVV	100789	73316	1.78%	0.139%
73	13.780	1022	1023	1025	rVV	69298	106181	2.57%	0.201%
74	13.825	1025	1027	1031	rVB2	231781	370866	8.98%	0.703%
75	13.894	1031	1033	1036	rBV	31188	51002	1.24%	0.097%
76	13.985	1037	1041	1045	rBV3	1044365	2248503	54.46%	4.262%
77	14.088	1047	1050	1051	rBV2	87540	111705	2.71%	0.212%
78	14.145	1053	1055	1057	rBV2	35392	66341	1.61%	0.126%
79	14.362	1072	1074	1076	rVV	92246	155396	3.76%	0.295%
80	14.442	1079	1081	1084	rBV3	57360	123707	3.00%	0.234%
81	14.682	1100	1102	1106	rBV4	36823	84281	2.04%	0.160%
82	14.854	1115	1117	1120	rBV2	48636	80680	1.95%	0.153%
83	15.003	1127	1130	1134	rBV	486634	899074	21.78%	1.704%
84	15.105	1137	1139	1141	rVB	78508	83712	2.03%	0.159%
85	15.197	1144	1147	1149	rBV2	34938	79270	1.92%	0.150%
86	15.288	1153	1155	1158	rVB	280675	354458	8.59%	0.672%
87	15.345	1158	1160	1163	rBV	388557	494800	11.98%	0.938%
88	15.414	1163	1166	1167	rBV	579034	909675	22.03%	1.724%
89	15.825	1199	1202	1204	rBV	45724	76214	1.85%	0.144%
90	15.871	1204	1206	1210	rBV4	30087	85565	2.07%	0.162%
91	15.940	1210	1212	1214	rVB2	33007	45777	1.11%	0.087%
92	16.797	1284	1287	1297	rVB	165344	395441	9.58%	0.749%
93	16.968	1299	1302	1304	rBV	35295	69686	1.69%	0.132%
94	17.220	1321	1324	1329	rVB	137836	281550	6.82%	0.534%

Sum of corrected areas: 52762954

LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093222.D
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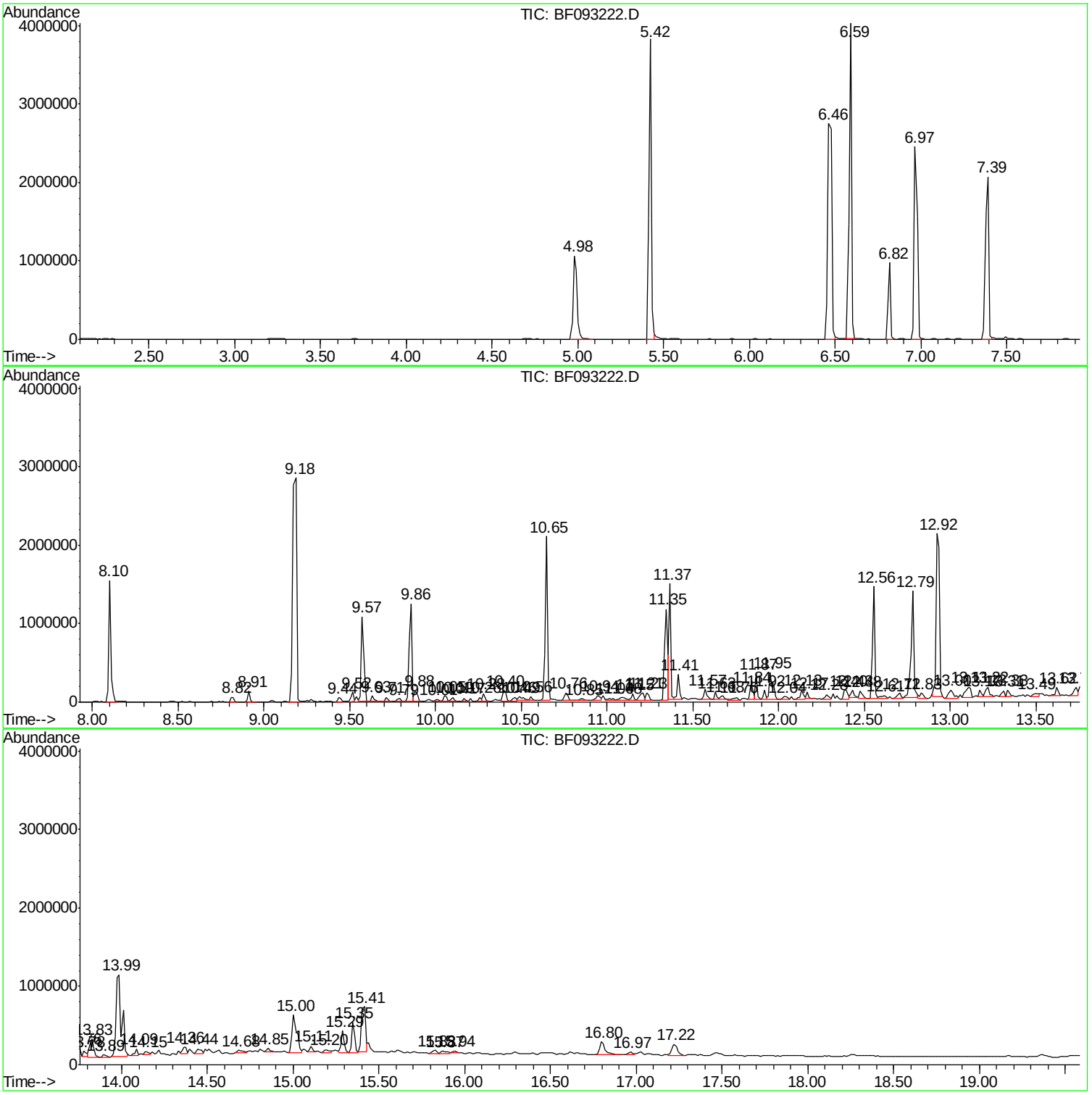
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093222.D
Acq On : 27 Feb 2017 23:24
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

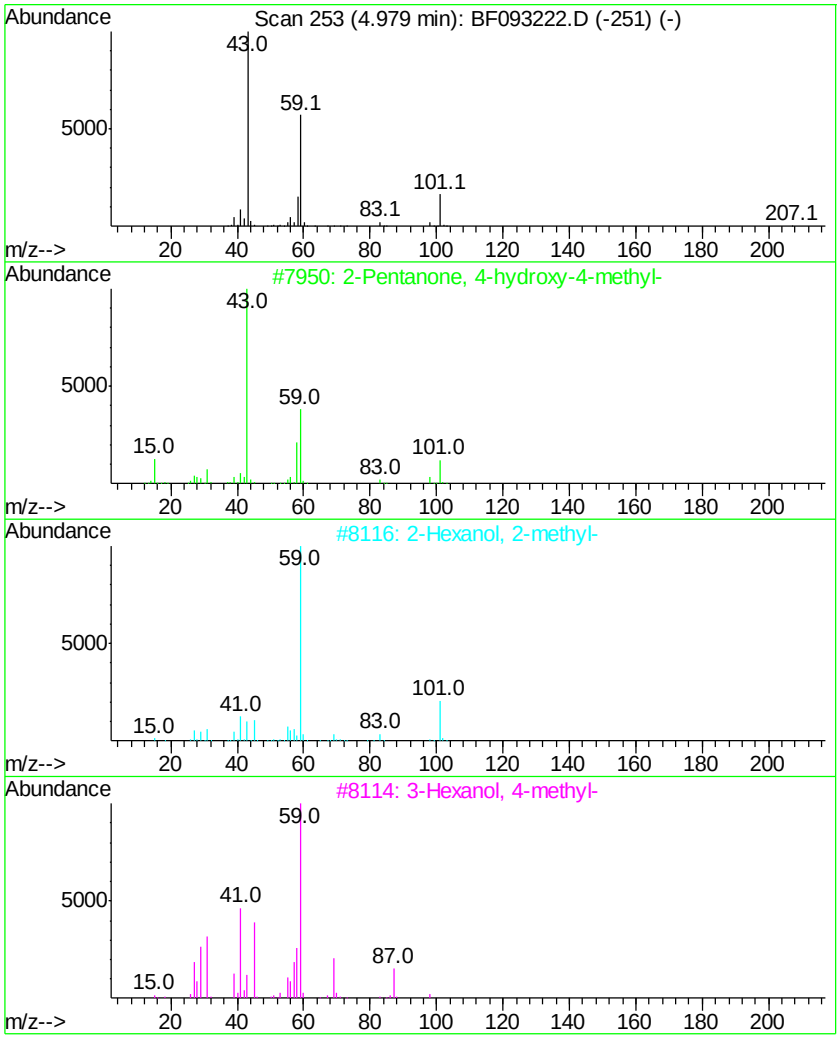
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

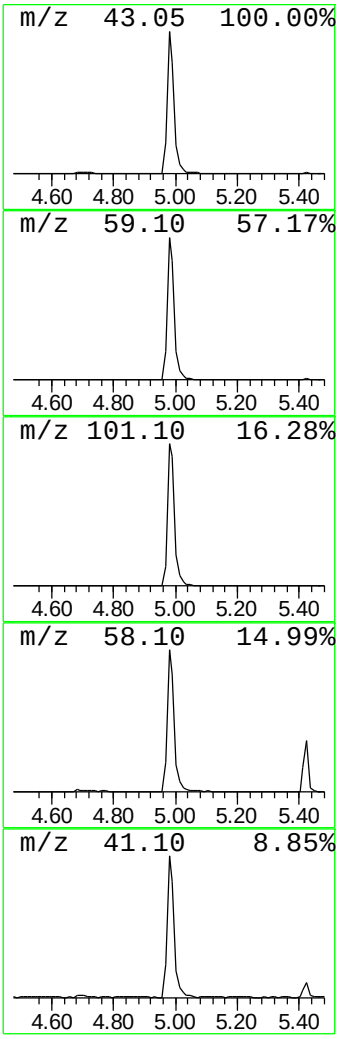
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	34.15 ng	1714100	1,4-Dichlorobenzene-d4	6.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093222.D
Acq On : 27 Feb 2017 23:24
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

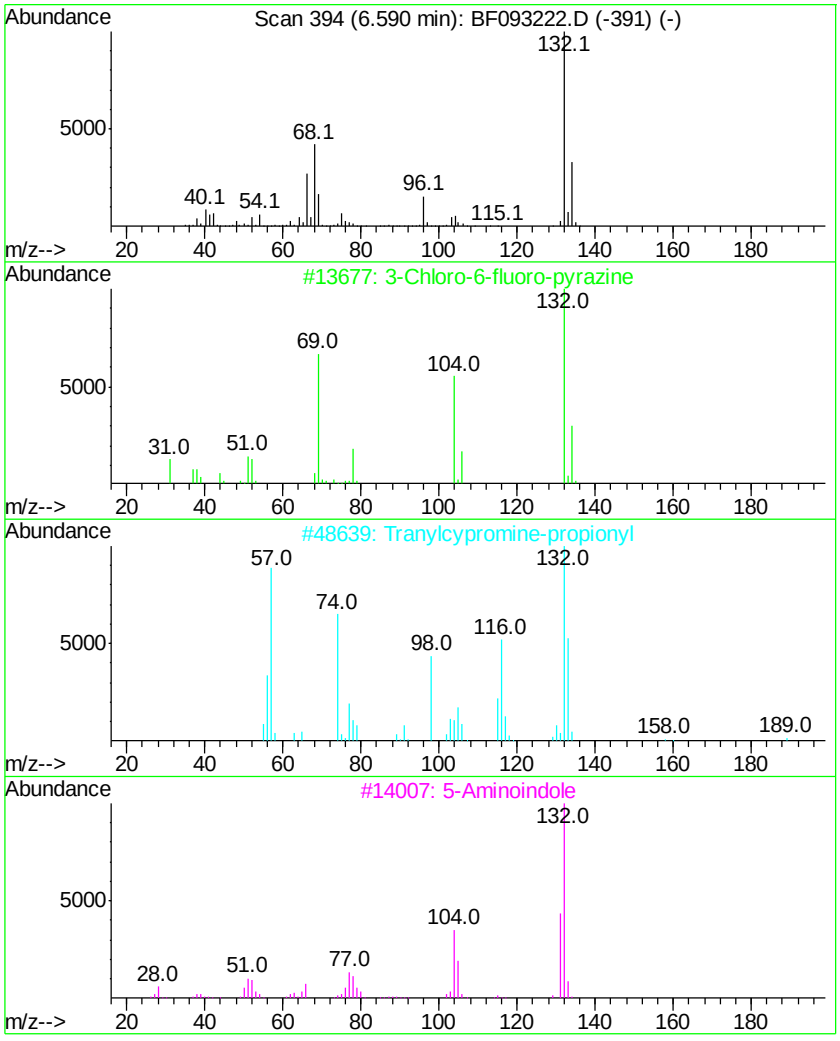
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

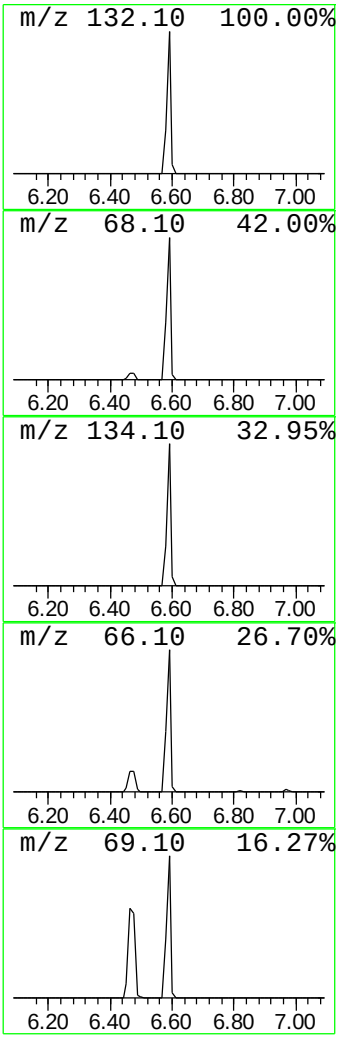
****

Peak Number 2 unknown6.59 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	77.90 ng	3910700	1,4-Dichlorobenzene-d4	6.82

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	5-Aminoindole	132	C8H8N2	005192-03-0	9
4	Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5	Xenon	132	Xe	007440-63-3	9





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Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

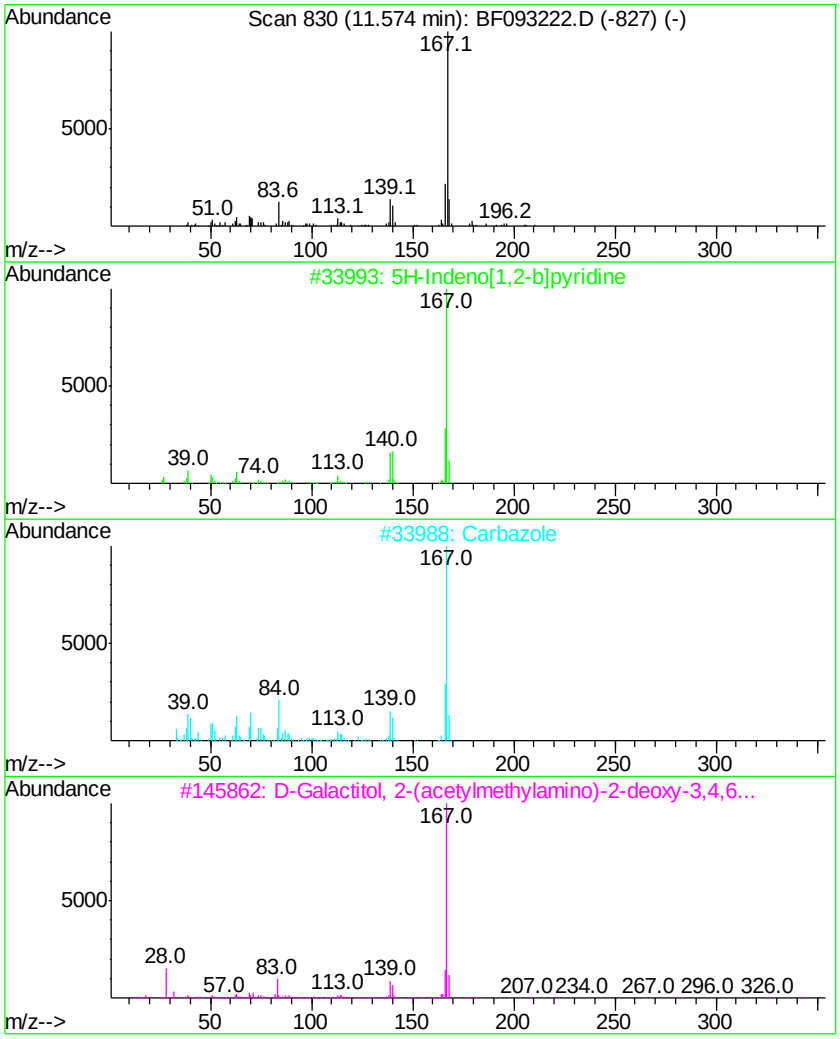
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

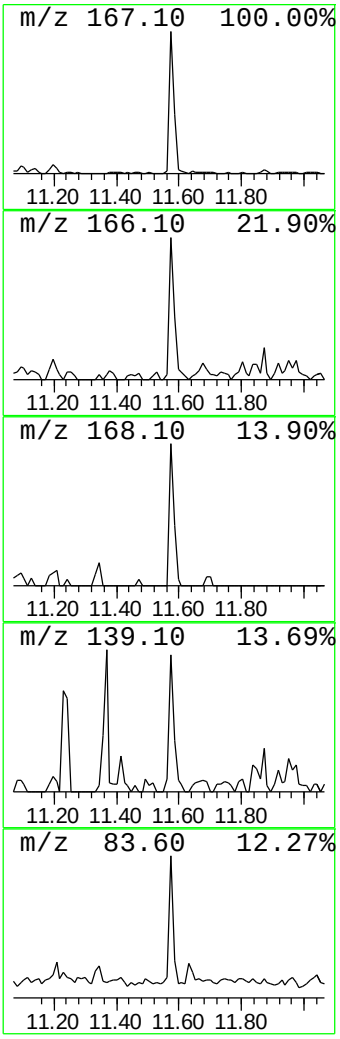
****

Peak Number 4 5H-Indeno[1,2-b]pyridine Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	2.27 ng	186582	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	93
2	Carbazole	167	C12H9N	000086-74-8	90
3	D-Galactitol, 2-(acetylmethylami...	363	C16H29NO8	074410-42-7	90
4	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	87
5	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	87





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
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Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

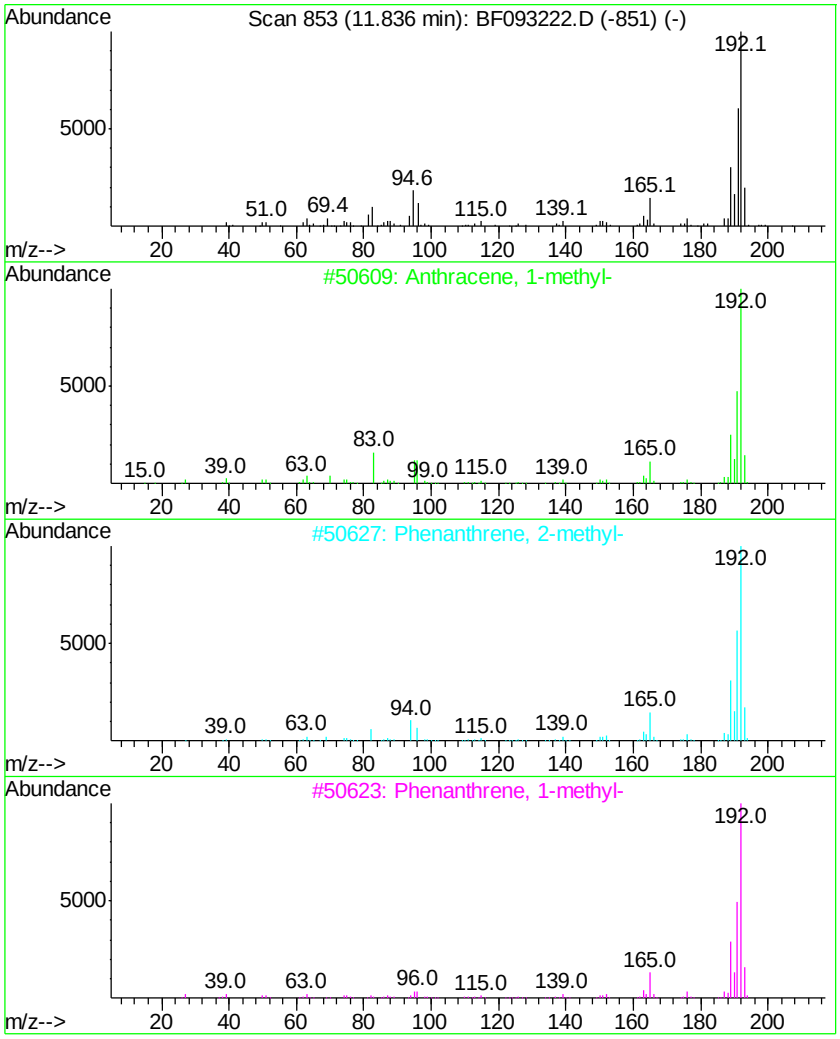
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

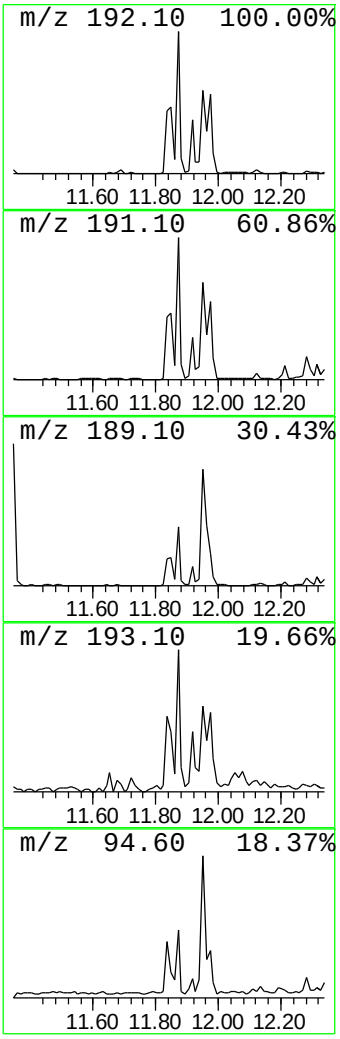
****

Peak Number 5 Anthracene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	2.99 ng	246063	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91





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Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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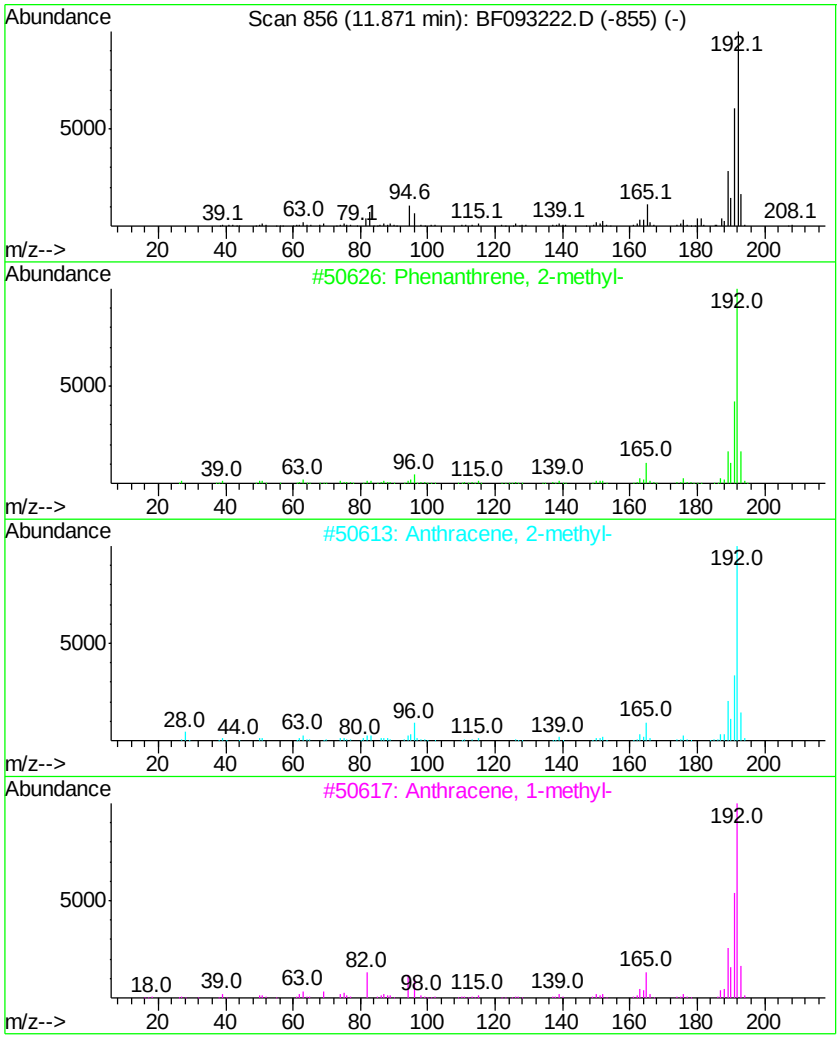
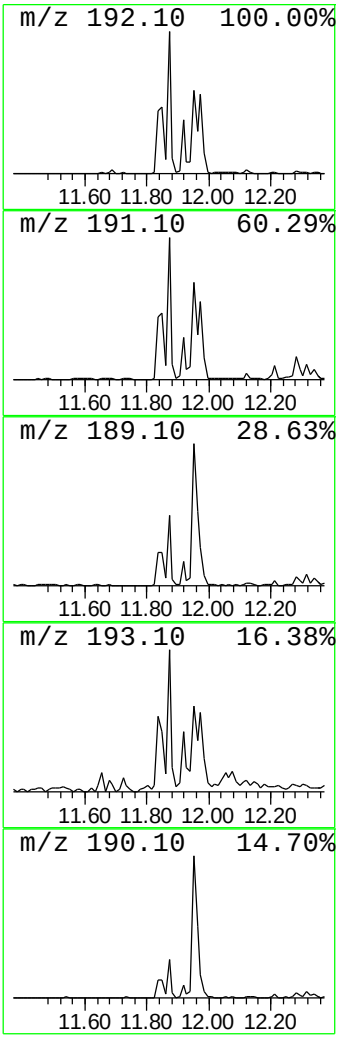
Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	3.03 ng	249790	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	89





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022717\
Data File : BF093222.D
Acq On : 27 Feb 2017 23:24
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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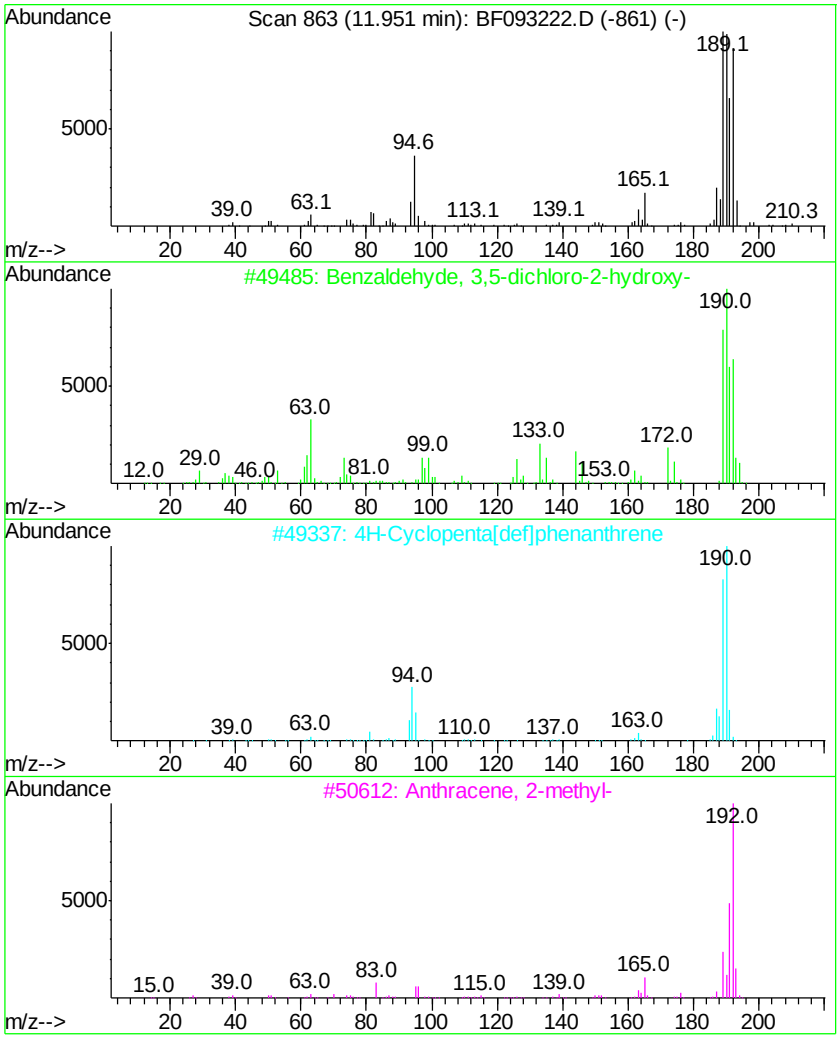
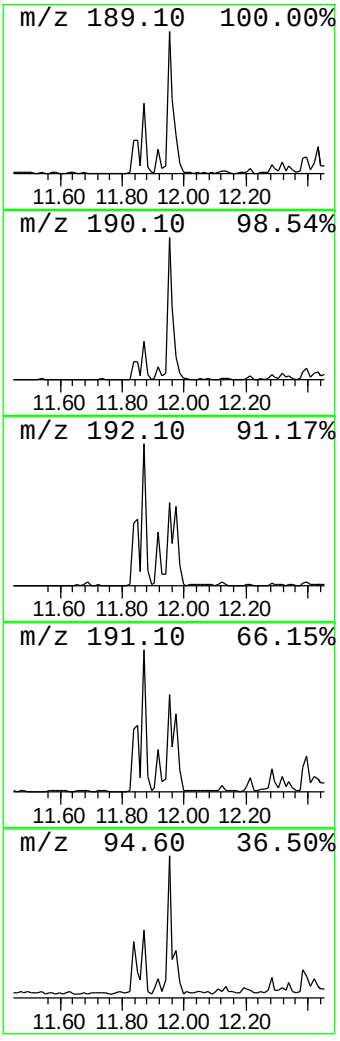
Peak Number 7 Benzaldehyde, 3,5-dichloro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	6.51 ng	536049	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	42
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	38
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38





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Data File : BF093222.D
Acq On : 27 Feb 2017 23:24
Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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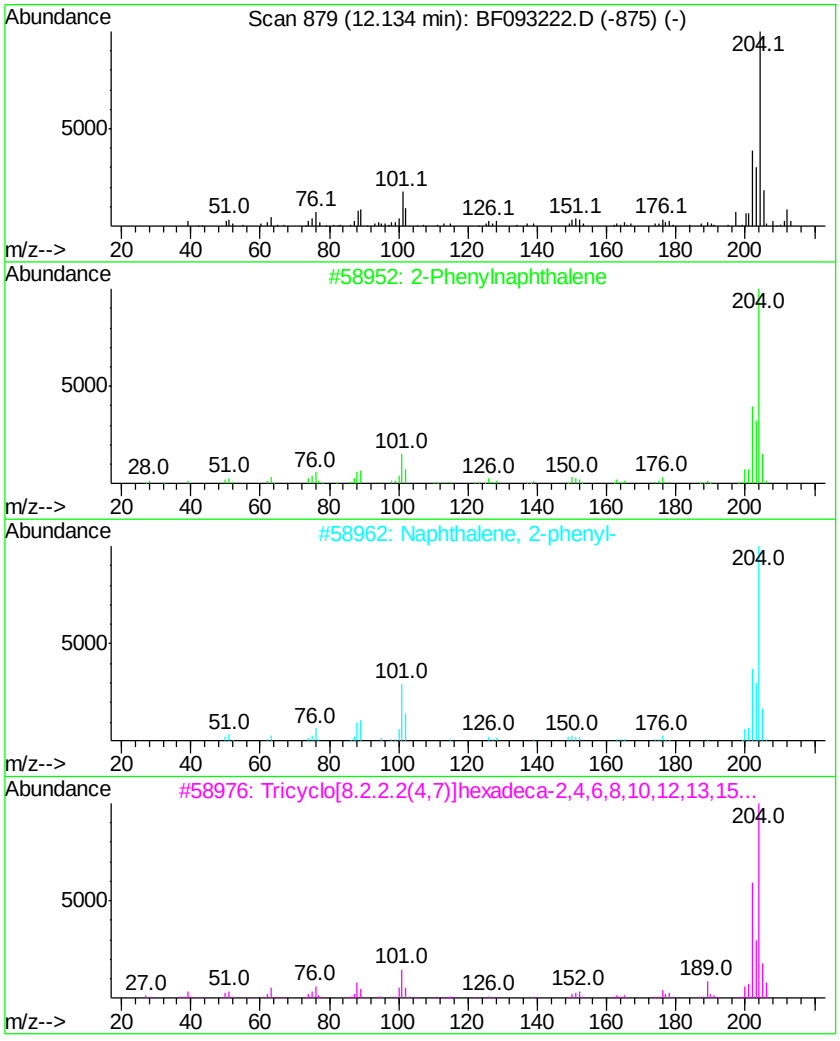
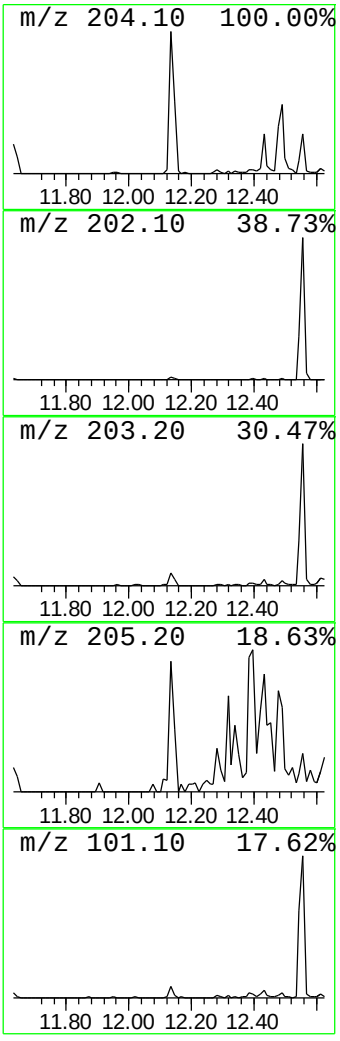
Peak Number 8 2-Phenylnaphthalene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	2.12 ng	174327	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenylnaphthalene	204	C16H12	035465-71-5	96
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	86
4	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	78
5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	78





Library Search Compound Report

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Data File : BF093222.D
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Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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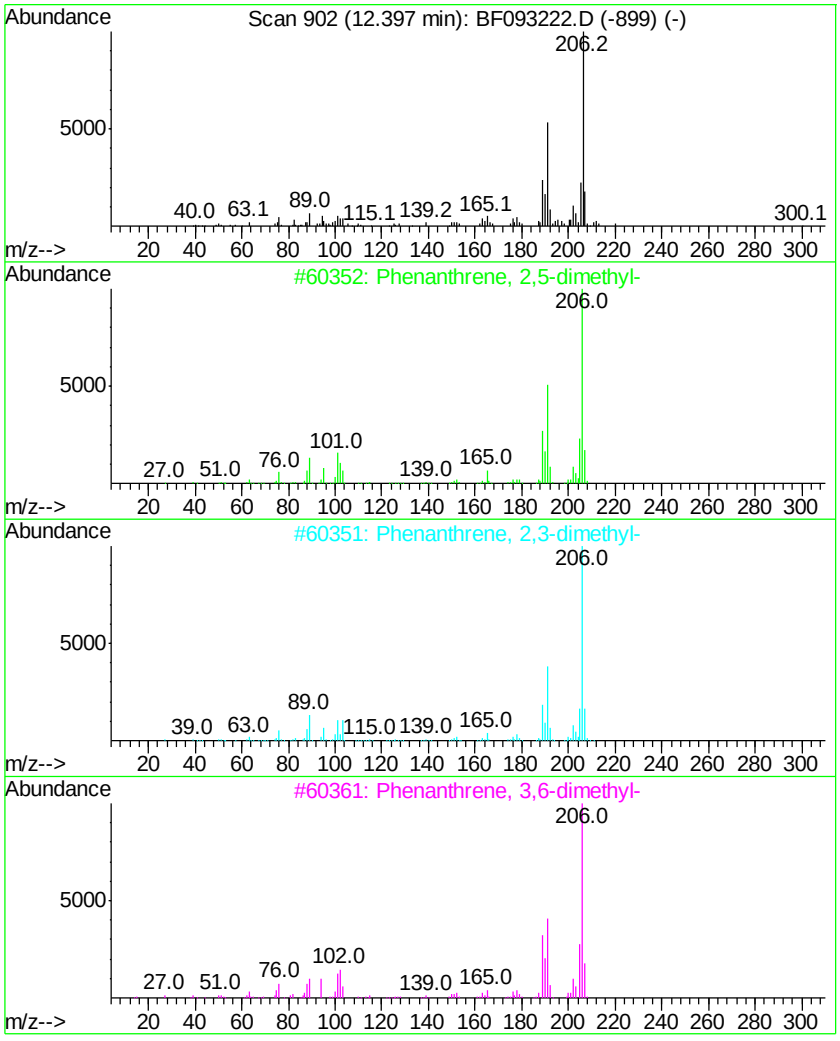
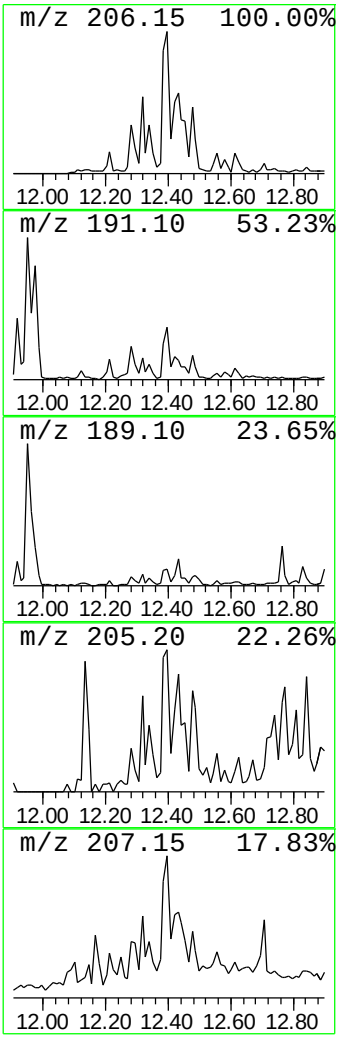
Peak Number 9 Phenanthrene, 2,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	2.08 ng	171198	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	86
4	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	86
5	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	83





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Operator : SJ/MA
Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
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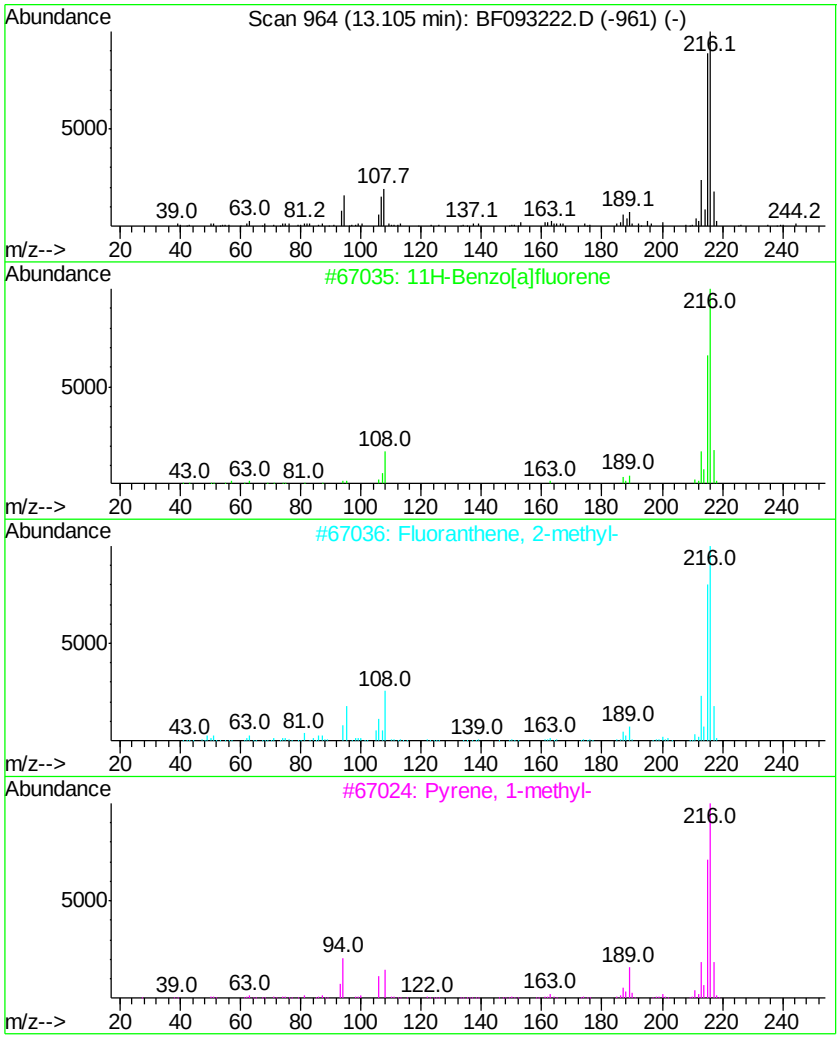
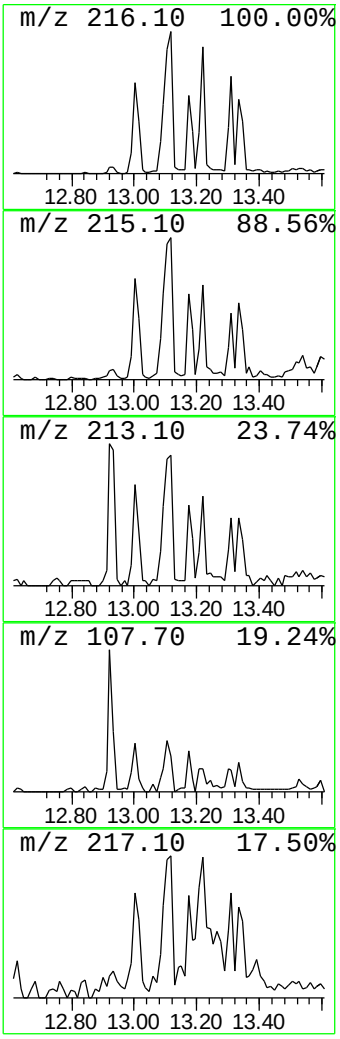
Peak Number 10 11H-Benzo[a]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	2.36 ng	265130	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	95
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	68





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ClientSampleId : E2-B1E-WSW-1

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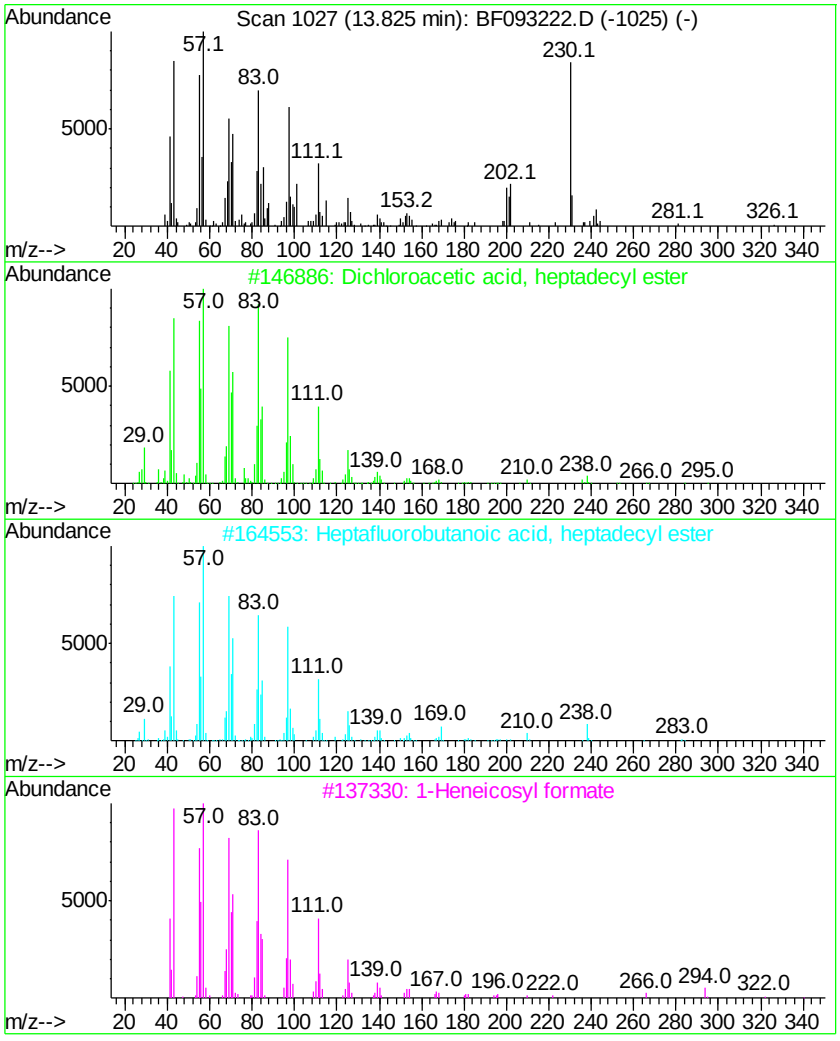
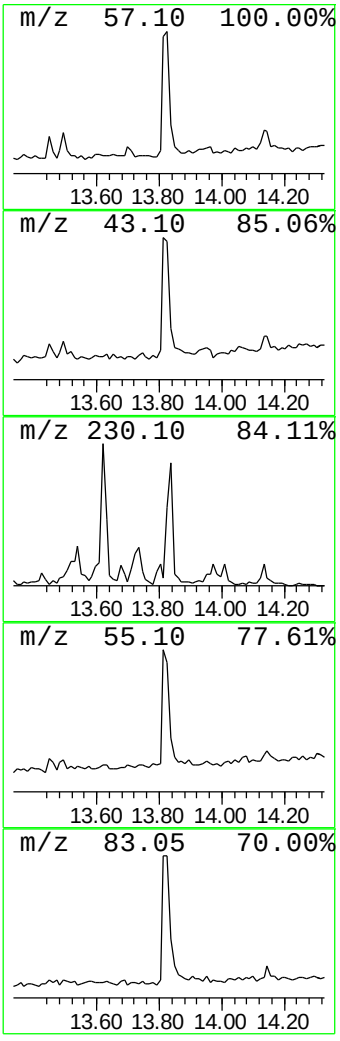
Peak Number 11 Dichloroacetic acid, heptad... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	3.30 ng	370866	Chrysene-d12	13.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	87
2	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	70
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	55
4	2- Chloropropionic acid, pentade...	318	C18H35ClO2	1000292-44-1	55
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	55





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Sample : I1997-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : E2-B1E-WSW-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
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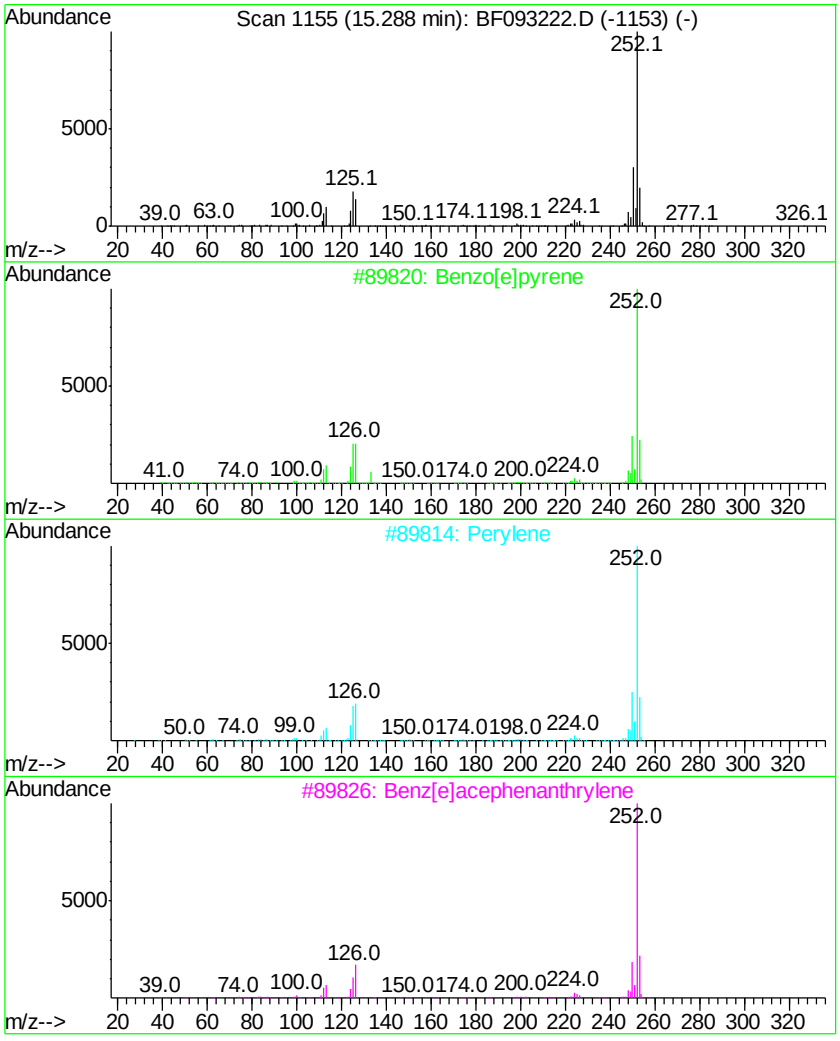
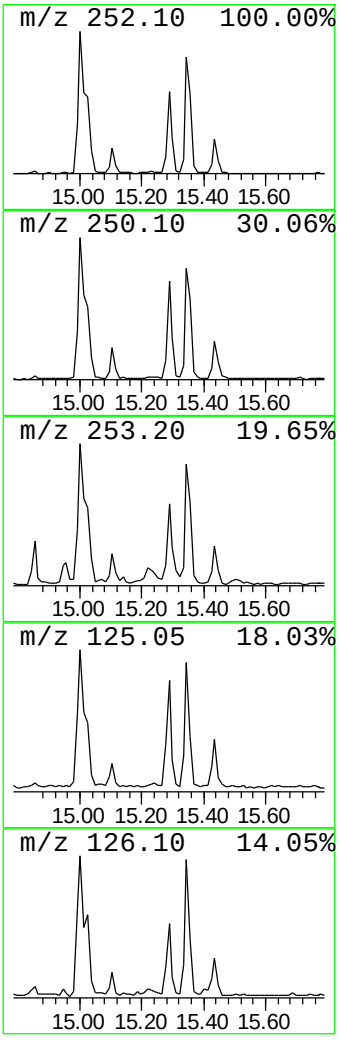
Peak Number 12 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	7.79 ng	354458	Perylene-d12	15.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2	Perylene	252	C20H12	000198-55-0	98
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
5	Benzo[a]pyrene	252	C20H12	000050-32-8	89





Tentatively Identified Compound (LSC) summary

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.98	34.1	ng	1714100	1	6.82	1003970	20.0
unknown6.59	6.59	77.9	ng	3910700	1	6.82	1003970	20.0
5H-Indeno[1,2-b]p...	11.57	2.3	ng	186582	4	11.35	1646280	20.0
Anthracene, 1-met...	11.84	3.0	ng	246063	4	11.35	1646280	20.0
Phenanthrene, 2-m...	11.87	3.0	ng	249790	4	11.35	1646280	20.0
Benzaldehyde, 3,5...	11.95	6.5	ng	536049	4	11.35	1646280	20.0
2-Phenylnaphthalene	12.13	2.1	ng	174327	4	11.35	1646280	20.0
Phenanthrene, 2,5...	12.40	2.1	ng	171198	4	11.35	1646280	20.0
11H-Benzo[a]fluorene	13.11	2.4	ng	265130	5	13.99	2248500	20.0
Dichloroacetic ac...	13.83	3.3	ng	370866	5	13.99	2248500	20.0
Benzo[e]pyrene	15.29	7.8	ng	354458	6	15.41	909675	20.0