LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
Data File : BF087801.D
Acq On : 7 Jun 2016 22:49
Operator : UM/SJ
Sample : H3379-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

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-BF060716.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF087802.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.690	5	9	11	rVB	3971220	6434123	100.00%	15.140%
2	1.758	14	15	25	rVB	366685	585003	9.09%	1.377%
3	1.896	25	27	42	rVB	1625768	2877535	44.72%	6.771%
4	2.090	42	44	70	rVB2	83451	339973	5.28%	0.800%
5	5.027	298	301	306	rBV	155618	243168	3.78%	0.572%
6	5.439	334	337	339	rBV	2179967	2264578	35.20%	5.329%
7	6.479	424	428	430	rBV	2130884	2660939	41.36%	6.261%
8	6.605	436	439	442	rBV	1983022	2430645	37.78%	5.719%
9	6.845	457	460	462	rBV	412744	566752	8.81%	1.334%
10	6.993	471	473	476	rVB	1729948	1881784	29.25%	4.428%
11	7.405	506	509	511	rBV	1604392	1575751	24.49%	3.708%
12	8.125	569	572	574	rBV	966312	849059	13.20%	1.998%
13	9.199	663	666	669	rBV	2664943	2574090	40.01%	6.057%
14	9.588	699	700	703	rVB	235109	246883	3.84%	0.581%
15	9.873	722	725	727	rBV	1028311	903741	14.05%	2.127%
16	9.908	727	728	730	rVB	94288	67071	1.04%	0.158%
17	10.422	771	773	775	rBV	62878	81314	1.26%	0.191%
18	10.662	791	794	796	rBV	1646940	1733671	26.94%	4.079%
19	10.788	803	805	809	rVB	45497	80472	1.25%	0.189%
20	11.222	839	843	849	rVB3	50447	166555	2.59%	0.392%
21	11.348	852	854	855	rBV	676561	764693	11.88%	1.799%
22	11.428	858	861	862	rVB	222563	205896	3.20%	0.484%
23	11.576	871	874	879	rBV2	61258	82618	1.28%	0.194%
24	11.851	896	898	899	rBV	112818	93555	1.45%	0.220%
25	11.874	899	900	902	rVB	102511	121270	1.88%	0.285%
26	11.919	902	904	906	rBV	50760	65593	1.02%	0.154%
27	11.965	906	908	912	rVB2	192974	289915	4.51%	0.682%
28	12.148	920	924	928	rVB2	103409	135182	2.10%	0.318%
29	12.399	943	946	948	rBV	90498	107636	1.67%	0.253%
30	12.479	952	953	958	rVB	83177	82935	1.29%	0.195%
31	12.559	958	960	962	rBV	1189867	1003471	15.60%	2.361%
32	12.708	969	973	977	rBV4	43313	87687	1.36%	0.206%
33	12.788	977	980	982	rBV	1092137	1256889	19.53%	2.957%
34	12.902	988	990	991	rBV	71919	66394	1.03%	0.156%

LSC Area Percent Report

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Acq On : 7 Jun 2016 22:49
Operator : UM/SJ
Sample : H3379-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

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-BF060716.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.937	991	993	996	rVB	2232635	2098444	32.61%	4.938%
36	13.005	996	999	1003	rBV3	79738	154018	2.39%	0.362%
37	13.108	1003	1008	1011	rBV2	128777	299970	4.66%	0.706%
38	13.177	1012	1014	1016	rBV	151692	150860	2.34%	0.355%
39	13.211	1016	1017	1021	rVB2	102743	138377	2.15%	0.326%
40	13.337	1027	1028	1030	rBV	73098	66783	1.04%	0.157%
41	13.611	1049	1052	1055	rBV2	66311	104487	1.62%	0.246%
42	13.725	1059	1062	1064	rBV	84159	148968	2.32%	0.351%
43	13.851	1072	1073	1080	rVB4	51737	134250	2.09%	0.316%
44	13.977	1081	1084	1085	rBV2	892916	1211456	18.83%	2.851%
45	14.080	1091	1093	1095	rBV	106274	115275	1.79%	0.271%
46	14.365	1116	1118	1119	rVB	90626	88976	1.38%	0.209%
47	14.457	1122	1126	1127	rBV	71355	92436	1.44%	0.218%
48	14.503	1127	1130	1132	rVB4	65512	143141	2.22%	0.337%
49	14.560	1132	1135	1138	rVB4	37611	67147	1.04%	0.158%
50	14.834	1156	1159	1163	rBV3	43619	96365	1.50%	0.227%
51	14.994	1170	1173	1178	rBV	485484	977454	15.19%	2.300%
52	15.097	1179	1182	1183	rVB	84329	134479	2.09%	0.316%
53	15.280	1195	1198	1201	rVB	294042	380615	5.92%	0.896%
54	15.337	1201	1203	1206	rVV	449133	526488	8.18%	1.239%
55	15.394	1206	1208	1214	rVB2	454097	795177	12.36%	1.871%
56	15.554	1219	1222	1224	rBV3	30414	77130	1.20%	0.181%
57	15.725	1235	1237	1241	rVB4	53646	118397	1.84%	0.279%
58	16.068	1265	1267	1273	rVB3	44627	103941	1.62%	0.245%
59	16.606	1310	1314	1319	rBV3	43777	139452	2.17%	0.328%
60	16.766	1325	1328	1333	rVB2	189148	386699	6.01%	0.910%
61	16.937	1340	1343	1345	rBV	43119	79058	1.23%	0.186%
62	17.189	1361	1365	1373	rBV	143541	332742	5.17%	0.783%
63	17.394	1380	1383	1390	rVB2	52289	145738	2.27%	0.343%
64	19.280	1543	1548	1558	rVB	41306	177139	2.75%	0.417%
65	19.452	1558	1563	1567	rBV	23136	86147	1.34%	0.203%

Sum of corrected areas: 42498450

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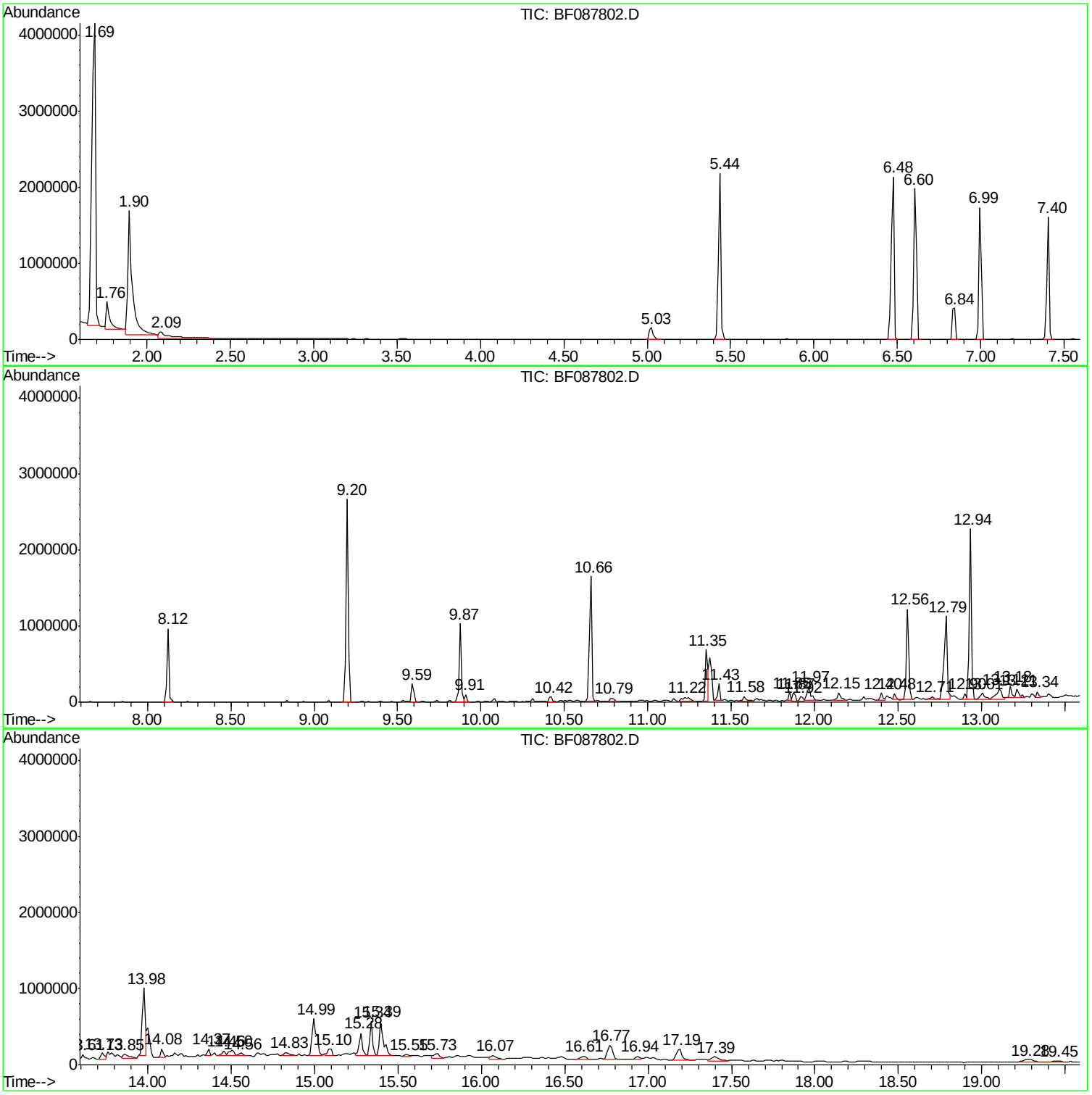
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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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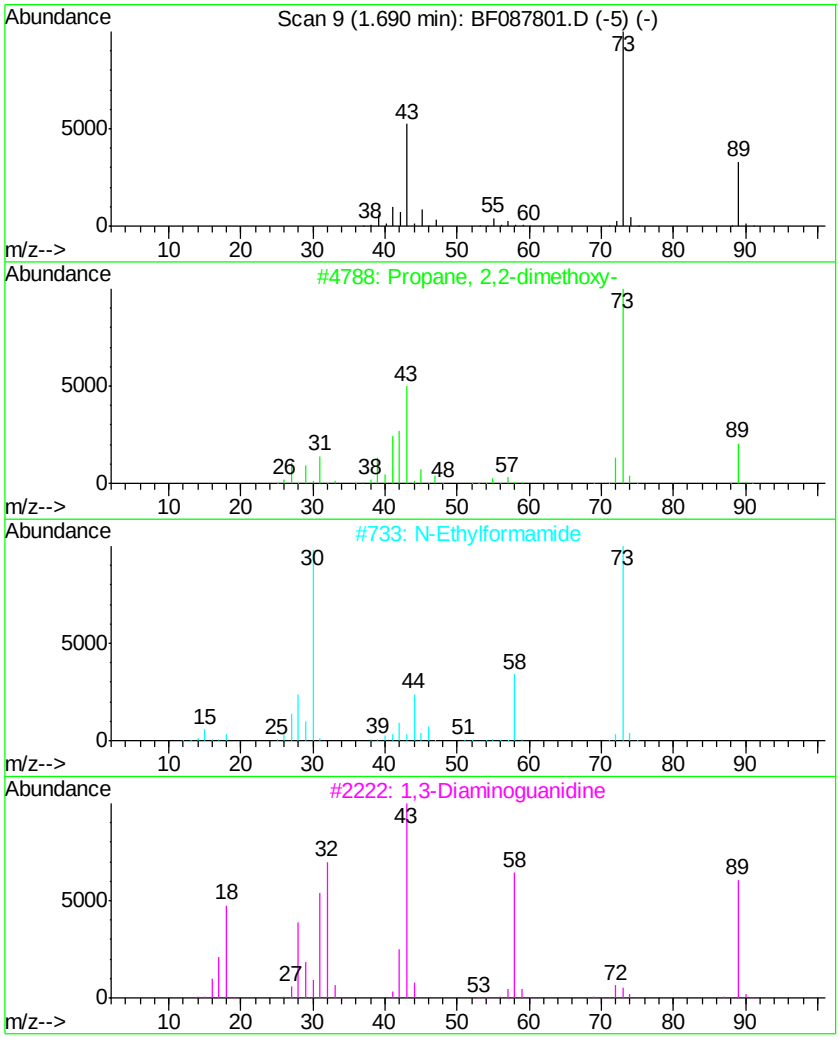
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

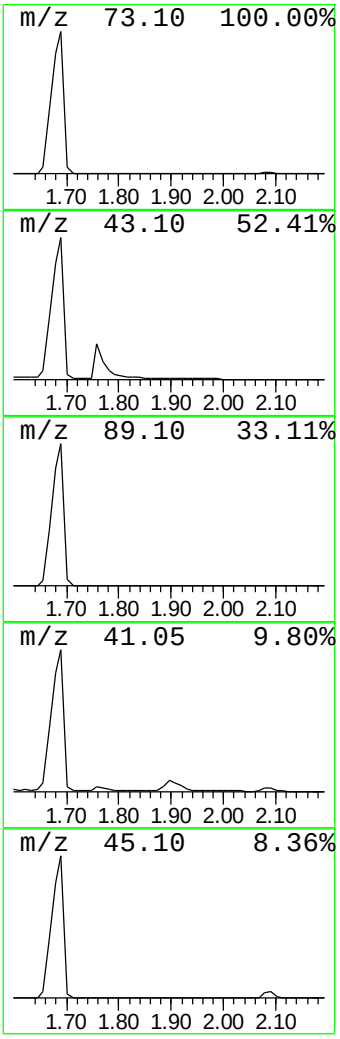
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.69	227.05 ng	6434120	1,4-Dichlorobenzene-d4	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2	N-Ethylformamide	73	C3H7NO	000627-45-2	9
3	1,3-Diaminoguanidine	89	CH7N5	004364-78-7	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





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
Data File : BF087801.D
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Operator : UM/SJ
Sample : H3379-06
Misc :
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Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

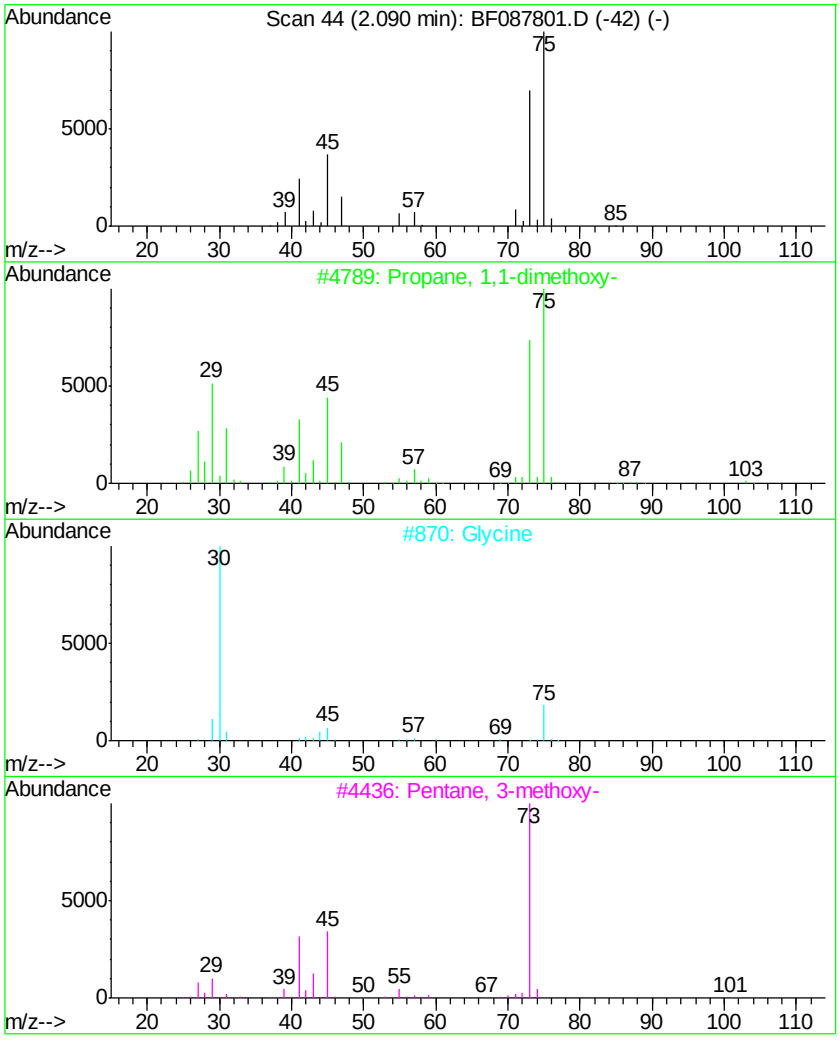
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

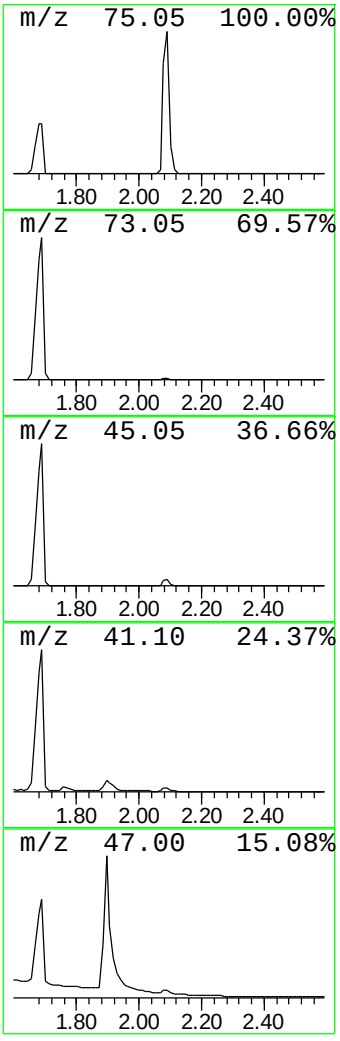
****

Peak Number 4 Propane, 1,1-dimethoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.09	12.00 ng	339973	1,4-Dichlorobenzene-d4	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2	Glycine	75	C2H5NO2	000056-40-6	9
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
4	Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	9
5	Silanol, trimethyl-	90	C3H10OSi	001066-40-6	9





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Operator : UM/SJ
Sample : H3379-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
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TIC Library : C:\Database\NIST11.L
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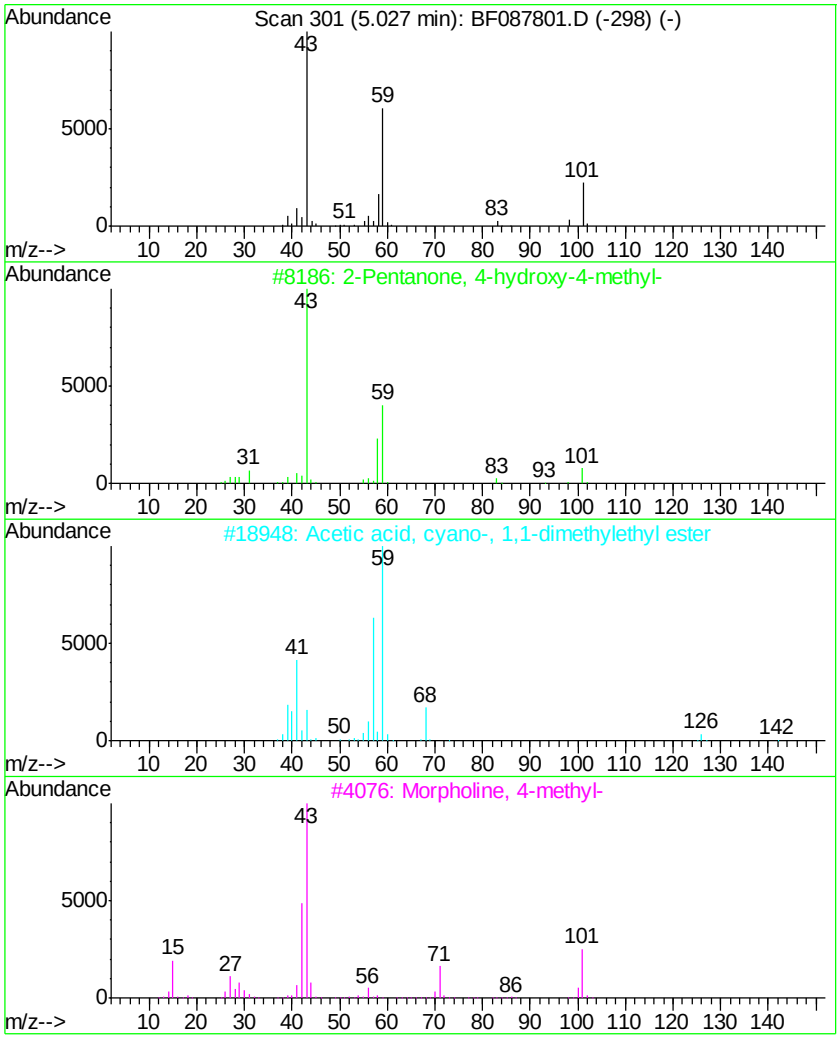
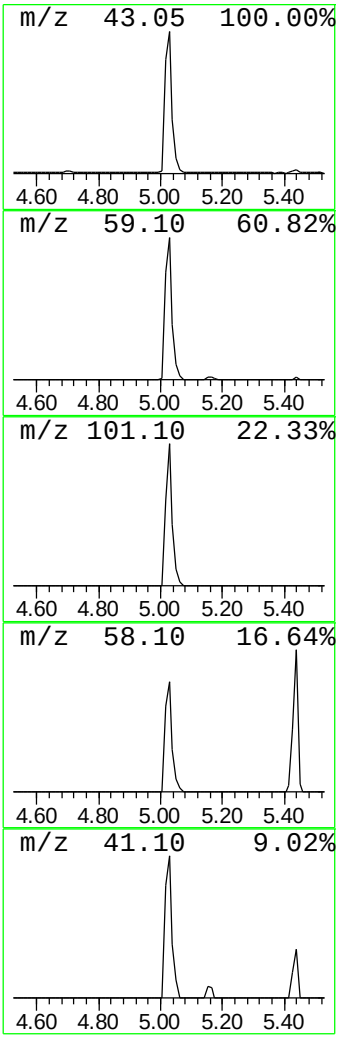
****

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	8.58 ng	243168	1,4-Dichlorobenzene-d4	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9





Library Search Compound Report

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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

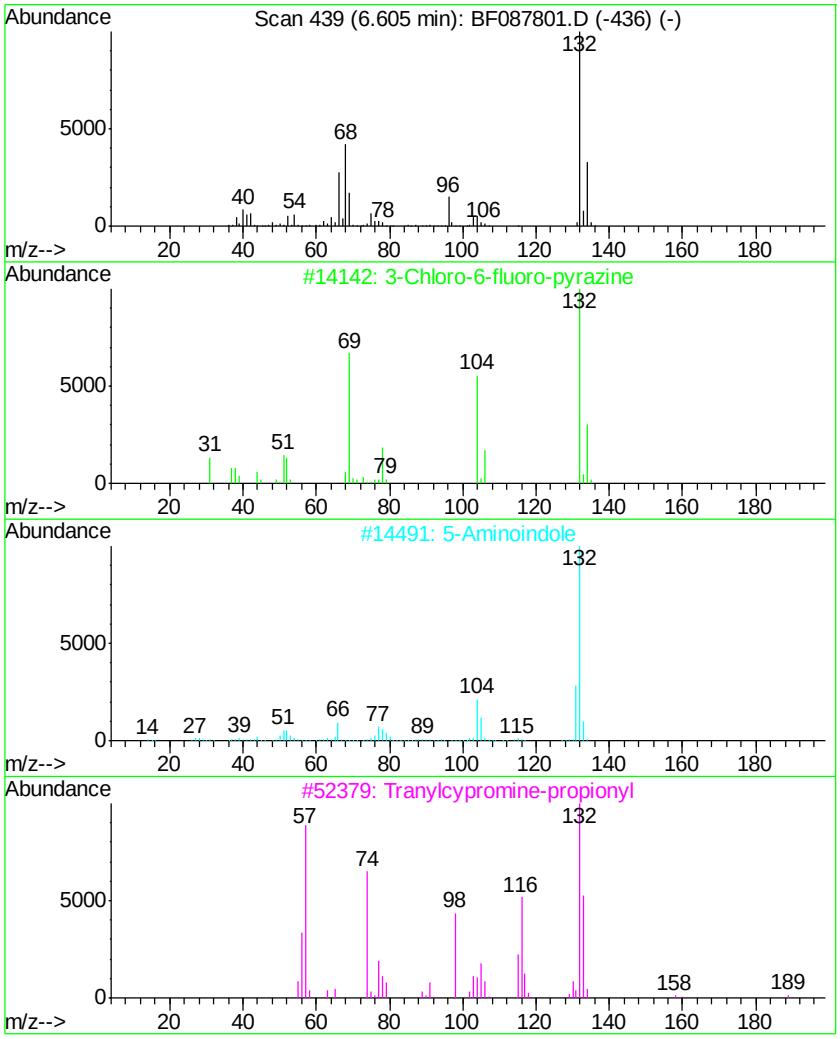
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

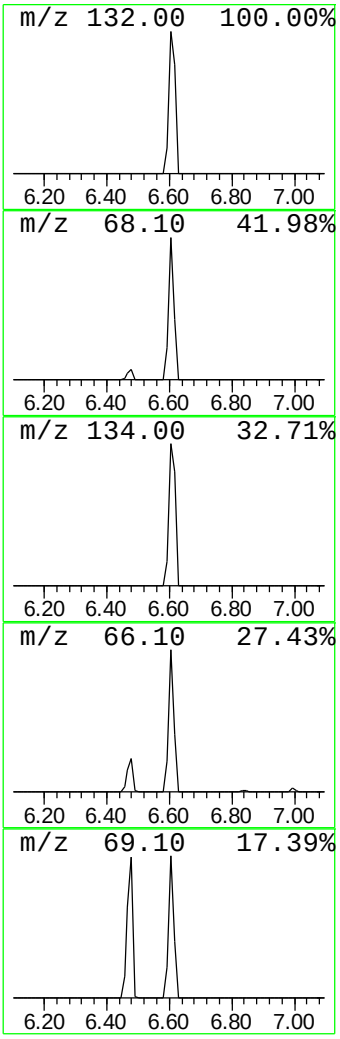
****

Peak Number 6 unknown6.60 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	85.77 ng	2430650	1,4-Dichlorobenzene-d4	6.84

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	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9





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Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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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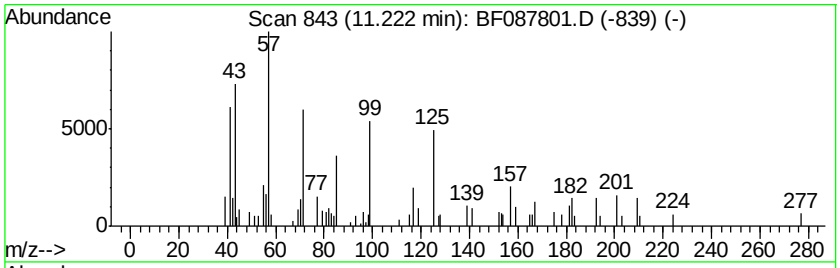
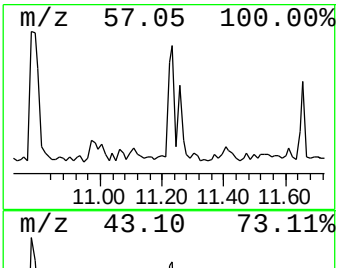
Peak Number 8 unknown11.22 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	4.36 ng	166555	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	22
2	Dodecane, 2-methyl-	184	C13H28	001560-97-0	22
3	Tetradecane	198	C14H30	000629-59-4	14
4	Dodecane	170	C12H26	000112-40-3	14
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	14





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Instrument : BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

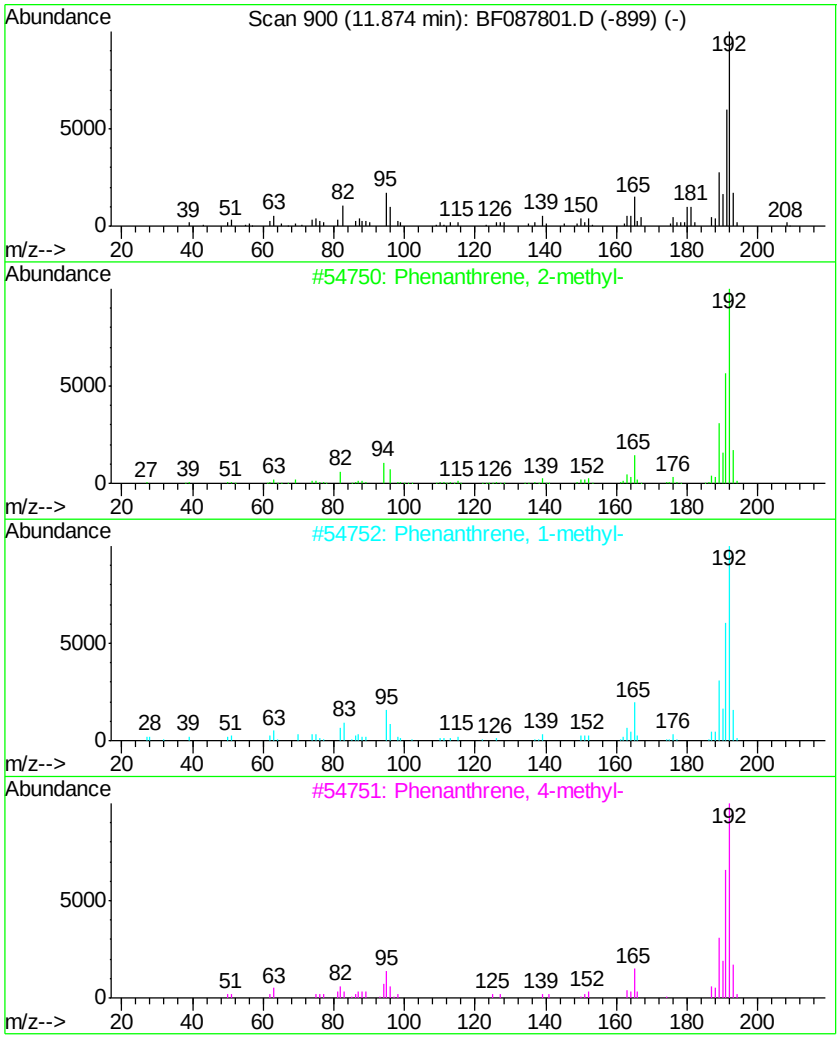
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

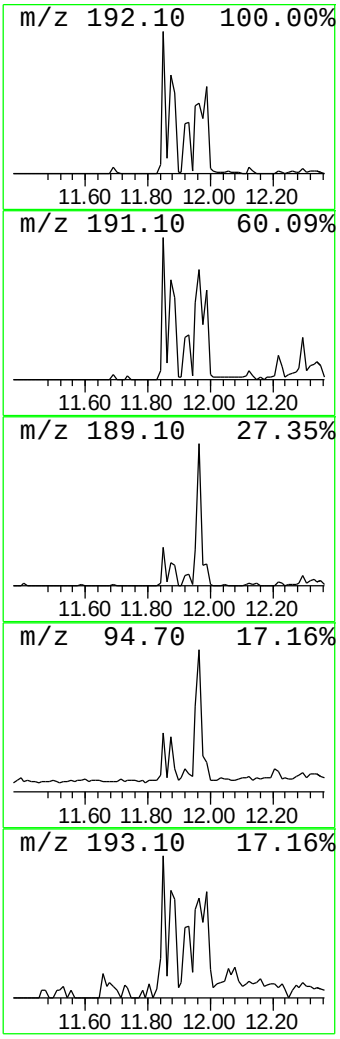
****

Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	3.17 ng	121270	Phenanthrene-d10	11.35

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	96
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94





Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
Data File : BF087801.D
Acq On : 7 Jun 2016 22:49
Operator : UM/SJ
Sample : H3379-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

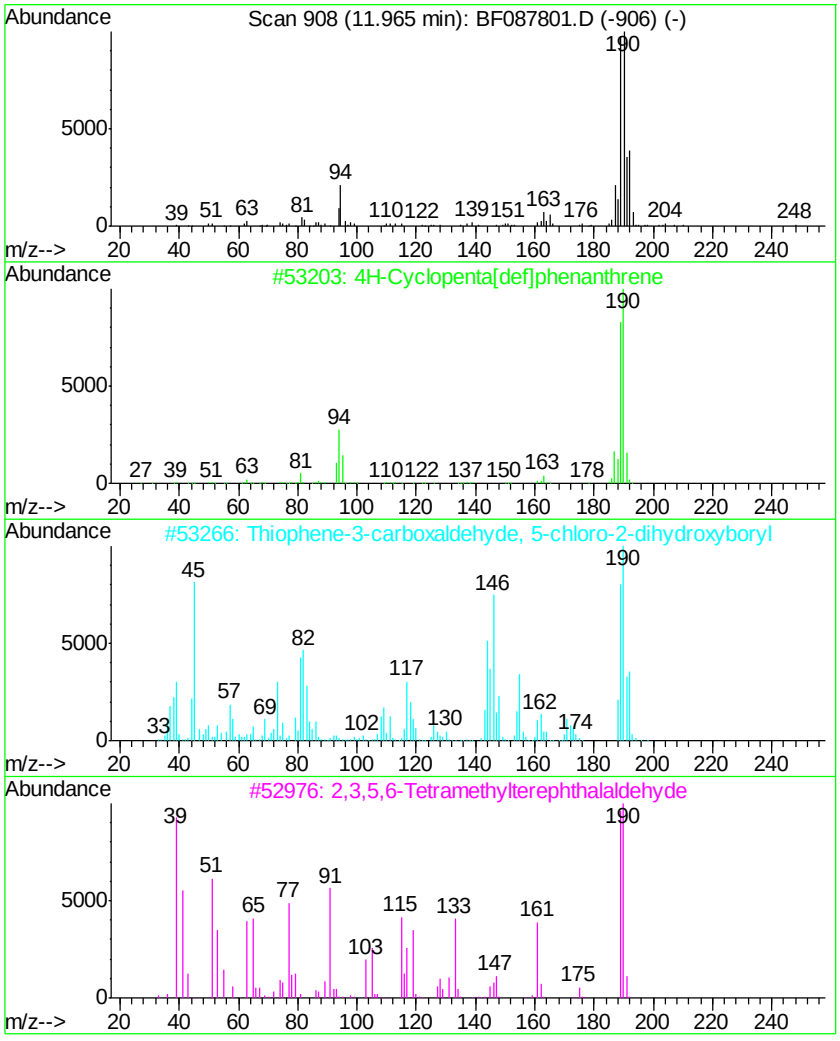
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

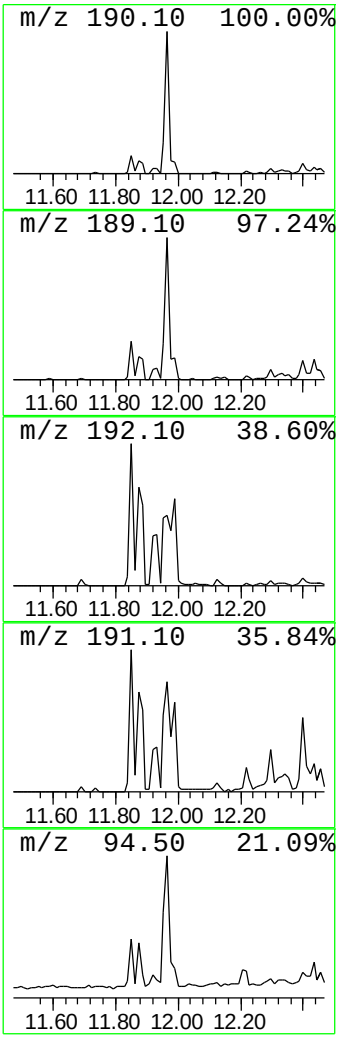
****

Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	7.58 ng	289915	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	45
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42





Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
Data File : BF087801.D
Acq On : 7 Jun 2016 22:49
Operator : UM/SJ
Sample : H3379-06
Misc :
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Instrument : BNA_F
ClientSampleId : SS-3

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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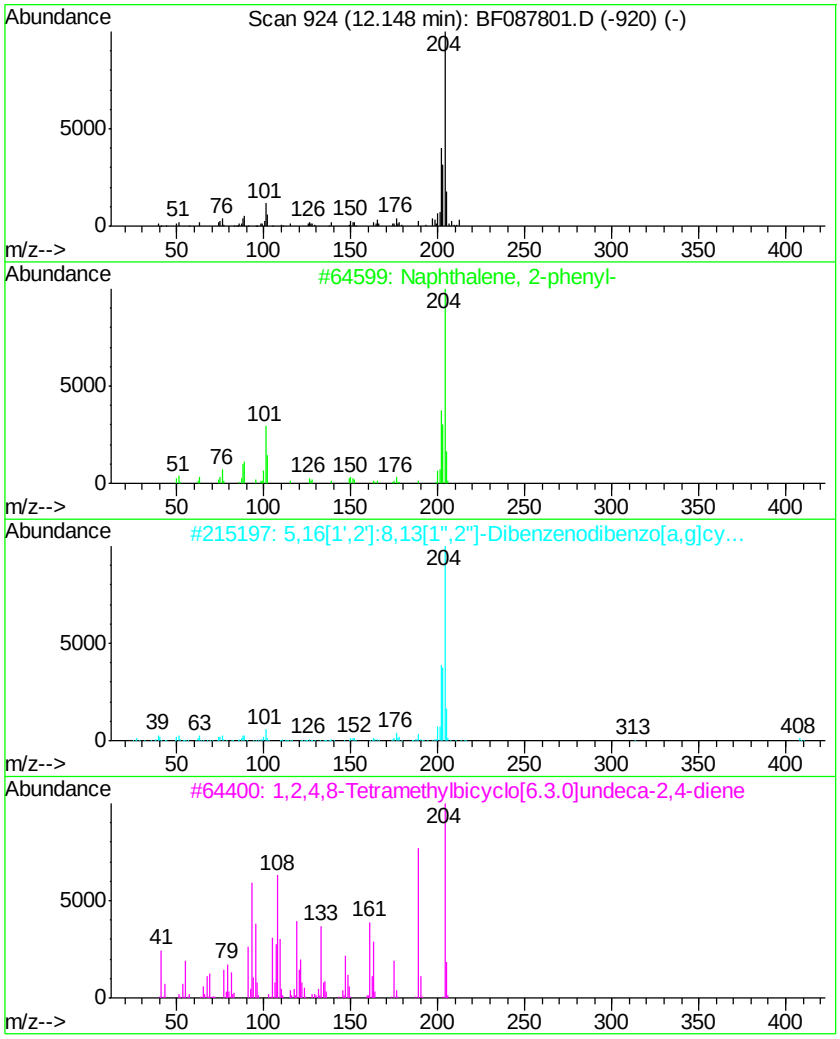
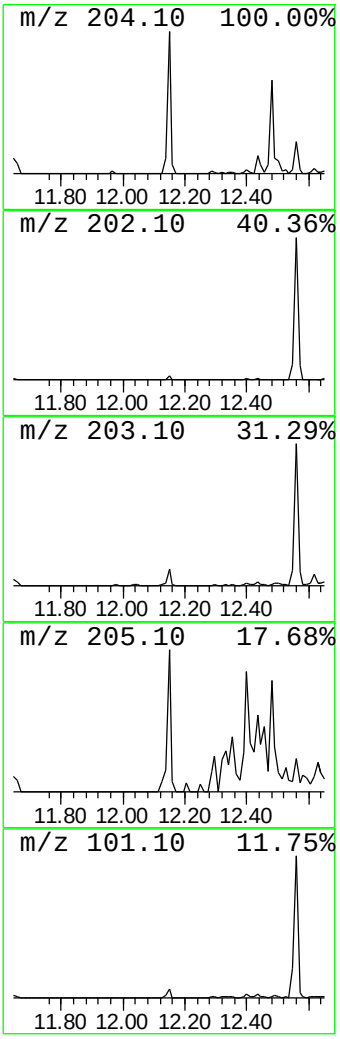
Peak Number 11 Naphthalene, 2-phenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	3.54 ng	135182	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	94
2	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91
3	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	83
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
5	6-Cyanophenanthridine	204	C14H8N2	002176-28-5	55





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Acq On : 7 Jun 2016 22:49
Operator : UM/SJ
Sample : H3379-06
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SS-3

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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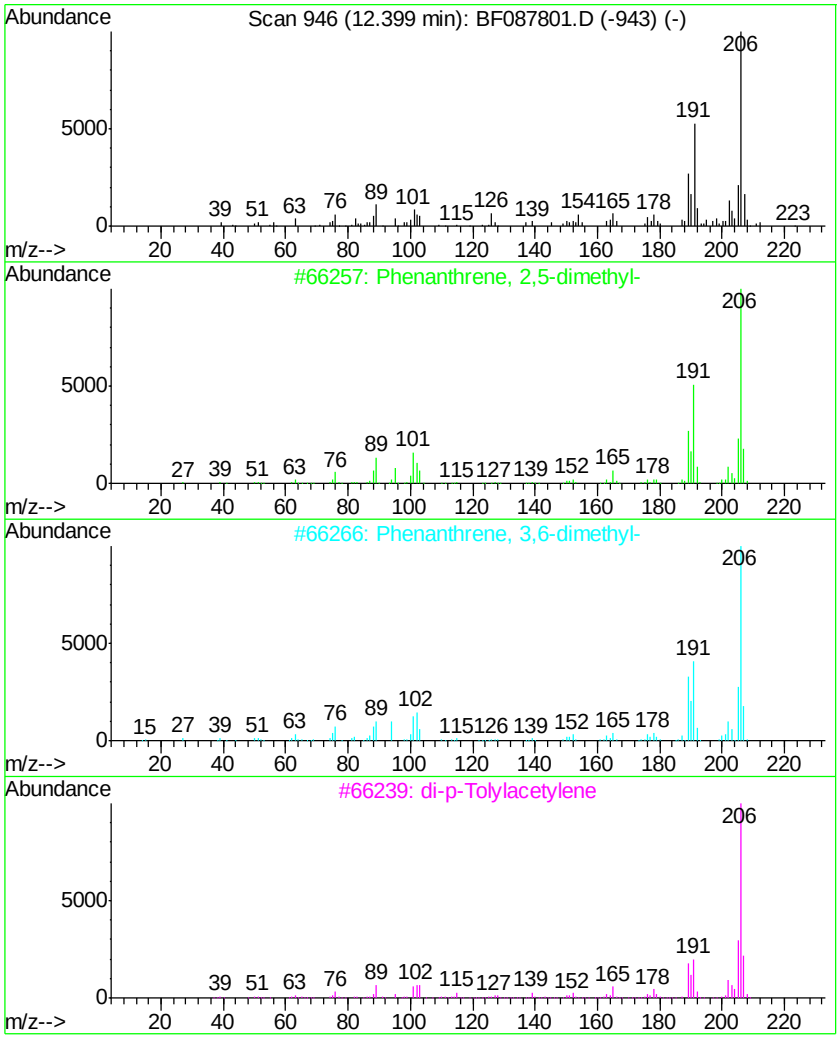
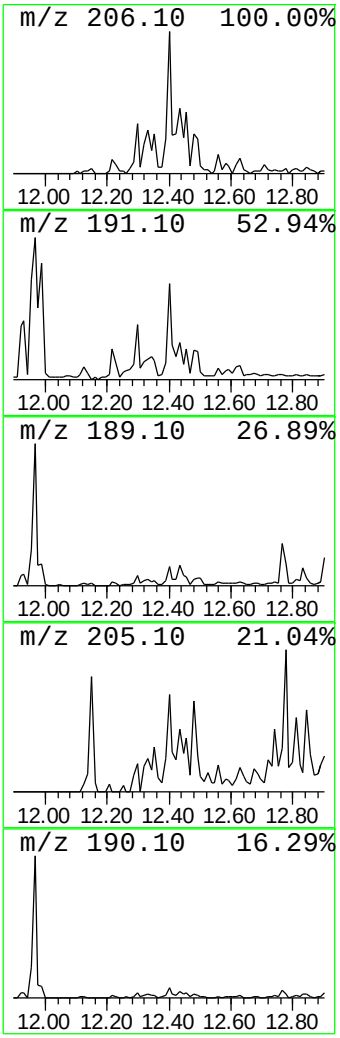
Peak Number 12 Phenanthrene, 2,5-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	2.82 ng	107636	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	92
4	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90
5	1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	89





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Instrument : BNA_F
ClientSampleId : SS-3

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TIC Library : C:\Database\NIST11.L
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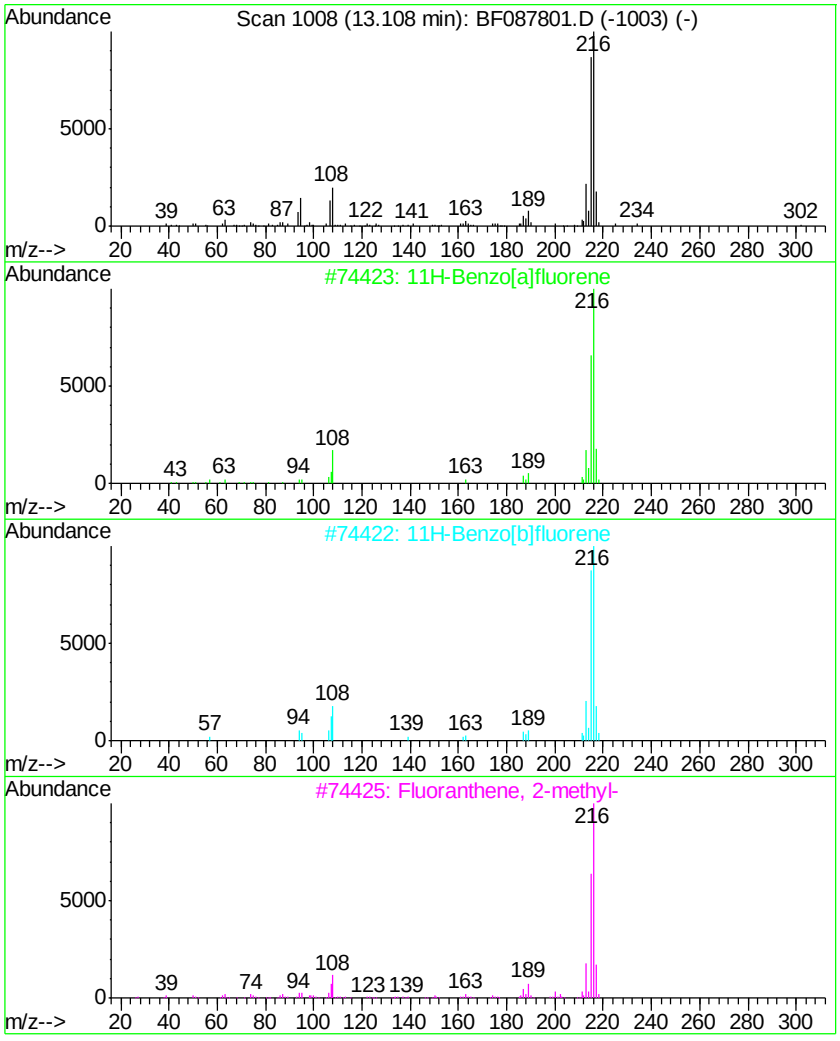
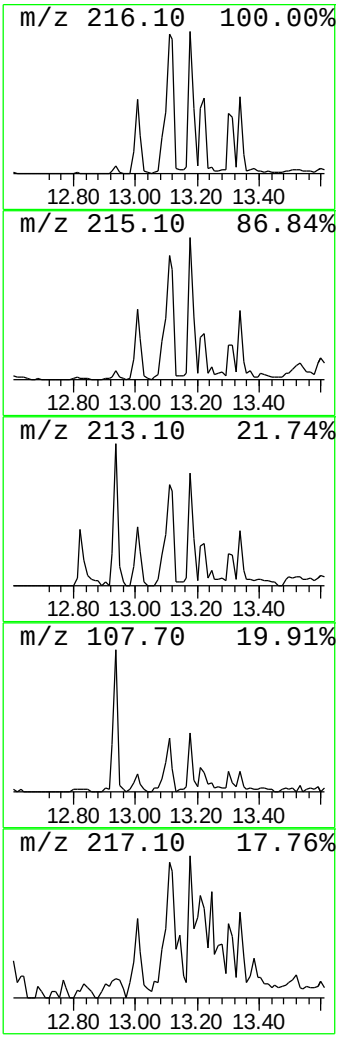
Peak Number 13 11H-Benzo[a]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	4.95 ng	299970	Chrysene-d12	13.98

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	87





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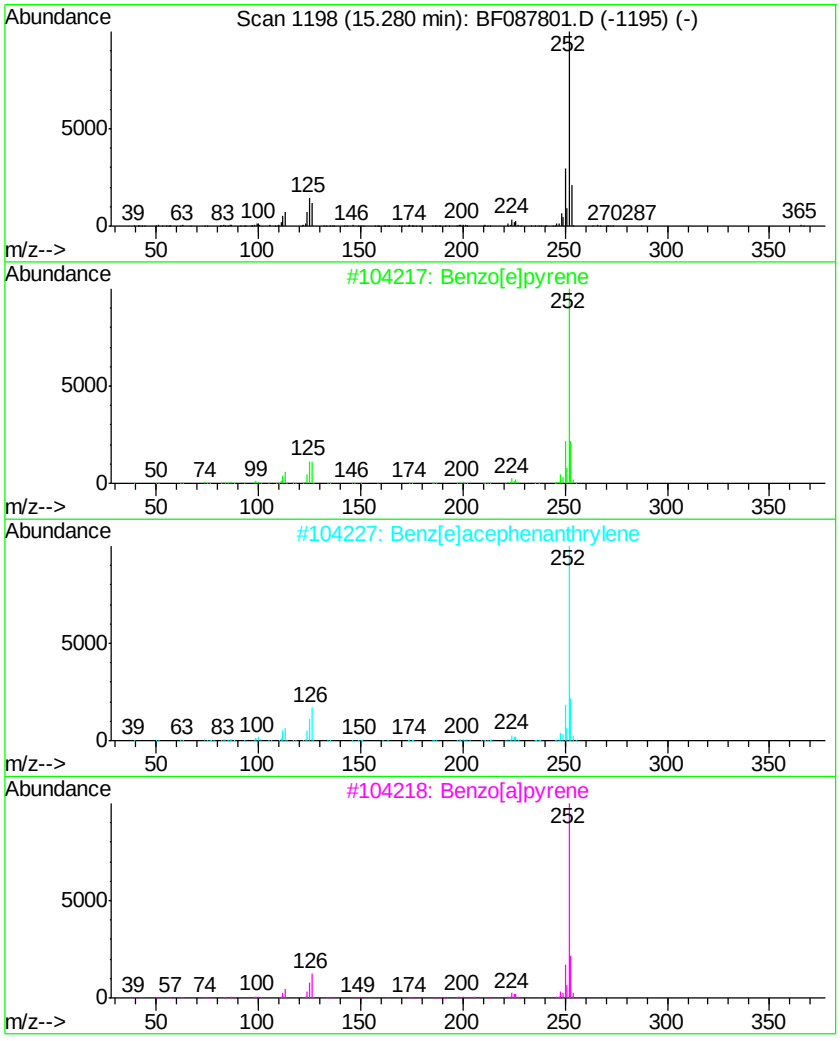
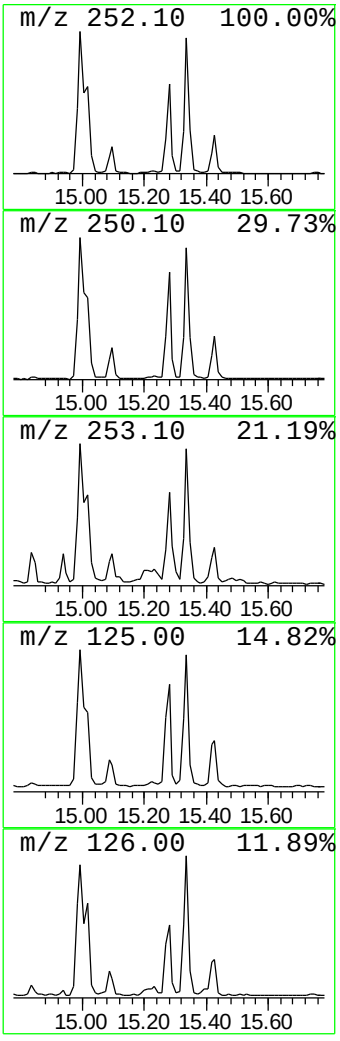
Peak Number 15 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	9.57 ng	380615	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
3	Benzo[a]pyrene	252	C20H12	000050-32-8	96
4	Perylene	252	C20H12	000198-55-0	95
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93





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ALS Vial : 16 Sample Multiplier: 1

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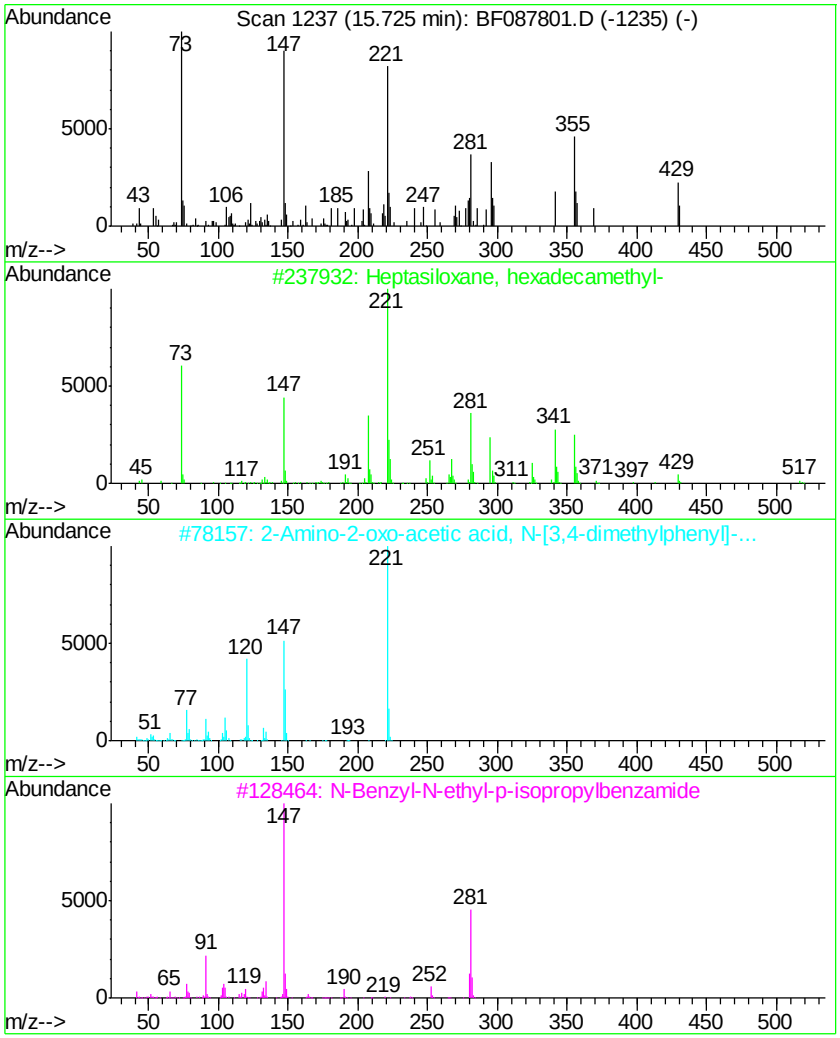
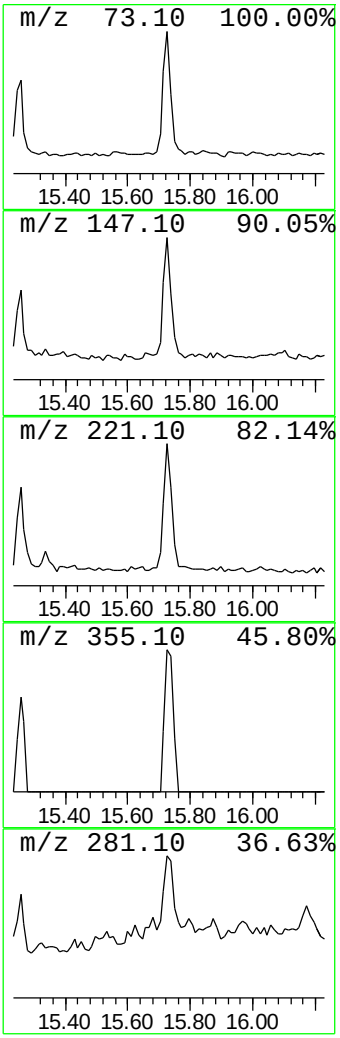
Peak Number 16 unknown15.73 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	2.98 ng	118397	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	38
2	2-Amino-2-oxo-acetic acid, N-[3,...	221	C12H15NO3	024451-17-0	16
3	N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	14
4	3-Trimethylsilyloxystearic acid,...	444	C24H52O3Si2	1000079-42-6	12
5	1,2-Benzenedicarboxylic acid, bi...	310	C14H22O4Si2	002078-22-0	12





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Misc :
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Instrument : BNA_F
ClientSampleId : SS-3

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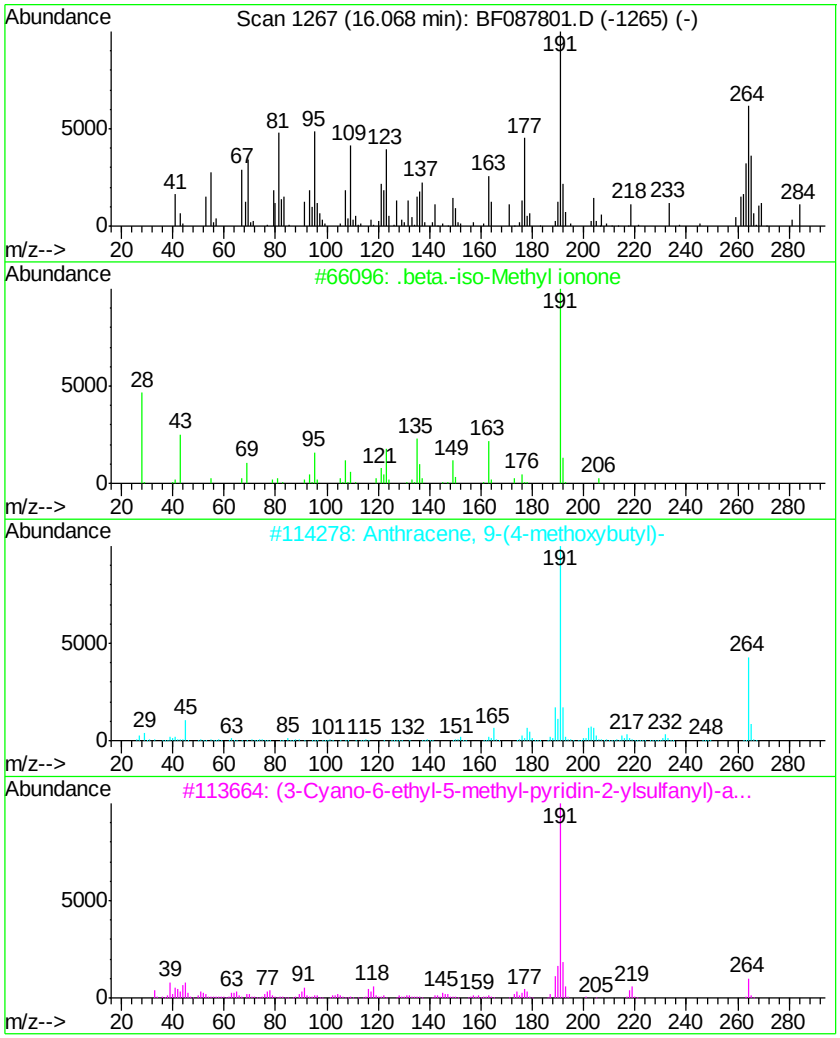
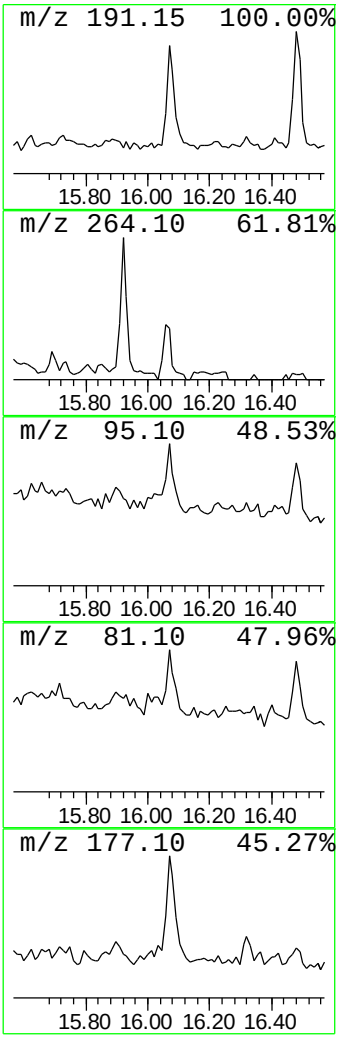
Peak Number 17 unknown16.07 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	2.61 ng	103941	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	32
2	Anthracene, 9-(4-methoxybutyl)-	264	C19H20O	144000-76-0	30
3	(3-Cyano-6-ethyl-5-methyl-pyridi...	264	C13H16N2O2S	1000300-45-4	27
4	1H-Pyrazole-4-acetic acid, 1-(4-...	264	C11H16N6O2	1000337-51-2	27
5	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	25





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ClientSampleId : SS-3

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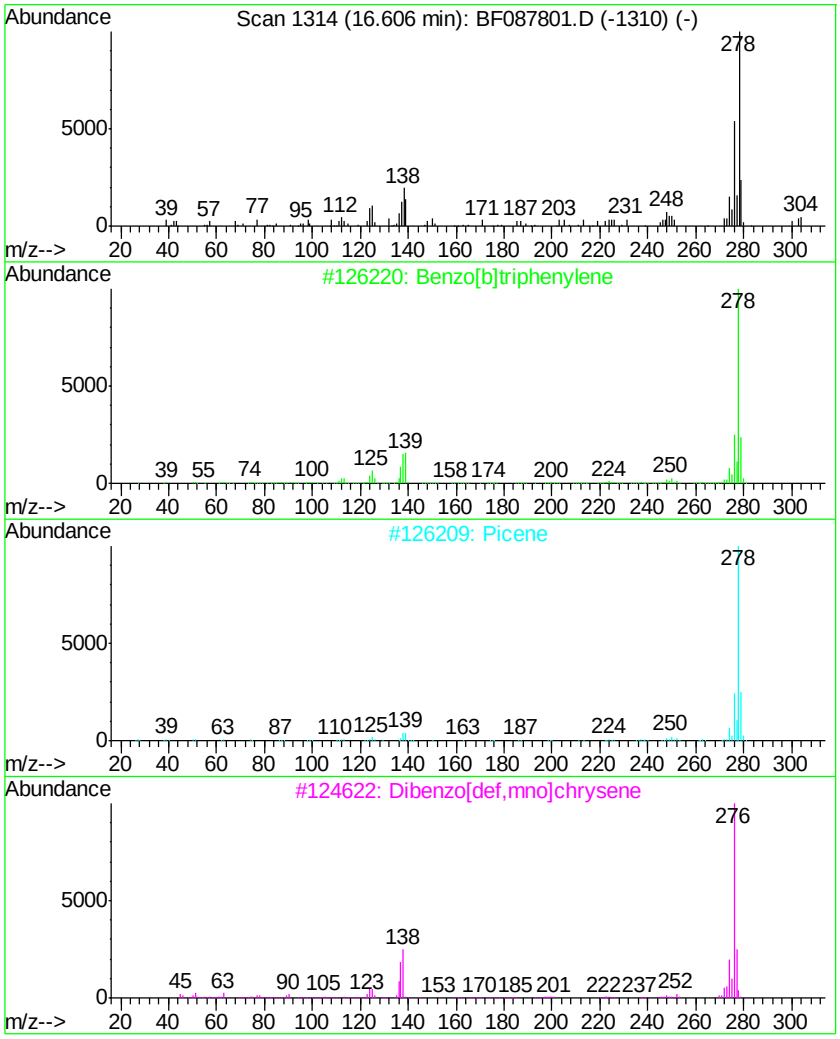
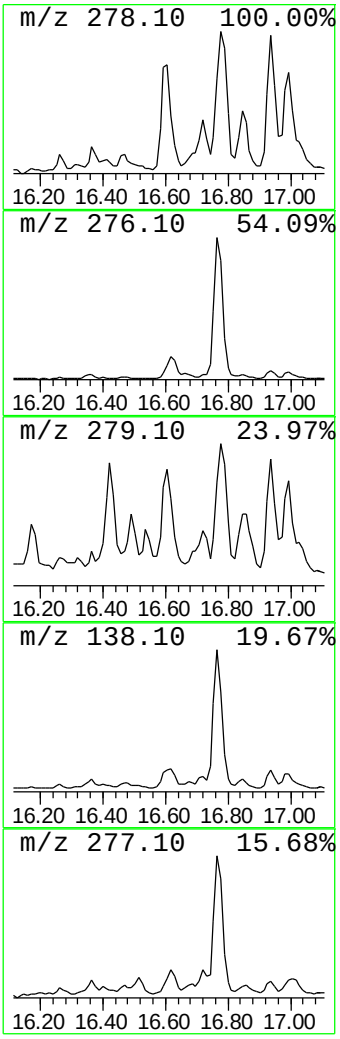
Peak Number 18 Benzo[b]triphenylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.61	3.51 ng	139452	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]triphenylene	278	C22H14	000215-58-7	64
2	Picene	278	C22H14	000213-46-7	58
3	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	49
4	Pyrene, 1-phenyl-	278	C22H14	005101-27-9	49
5	Benz[a]anthracene-7,12-dicarboni...	278	C20H10N2	035215-32-8	49





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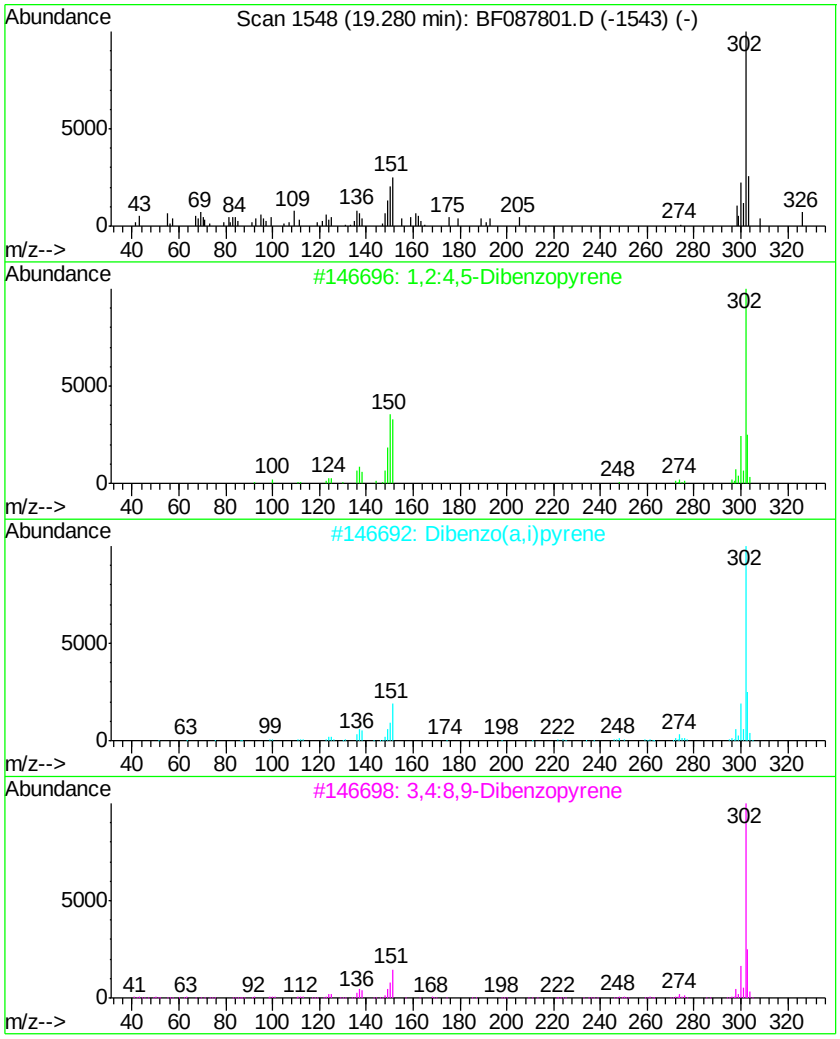
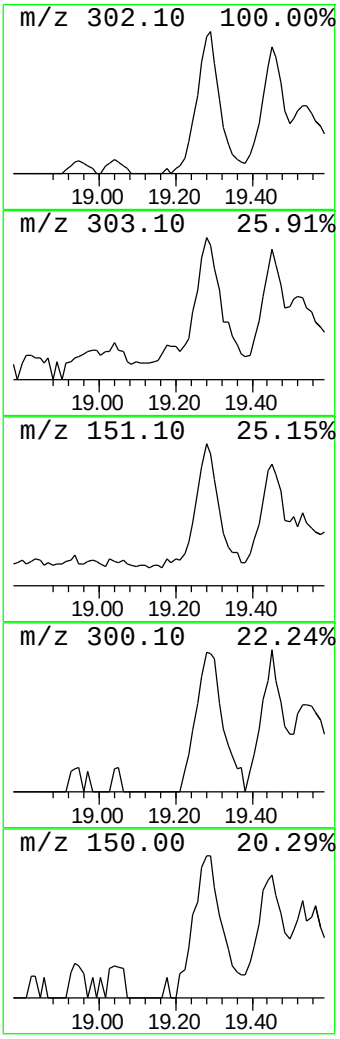
Peak Number 20 1,2:4,5-Dibenzopyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.28	4.46 ng	177139	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	81
2	Dibenzo(a,i)pyrene	302	C24H14	000189-55-9	68
3	3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	59
4	Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	50
5	Dibenzo[fg,op]naphthacene	302	C24H14	000192-51-8	50





Tentatively Identified Compound (LSC) summary

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TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Propane, 2,2-dime...	1.69	227.1	ng	6434120	1	6.84	566752	20.0
Propane, 1,1-dime...	2.09	12.0	ng	339973	1	6.84	566752	20.0
2-Pentanone, 4-hy...	5.03	8.6	ng	243168	1	6.84	566752	20.0
unknown6.60	6.60	85.8	ng	2430650	1	6.84	566752	20.0
unknown11.22	11.22	4.4	ng	166555	4	11.35	764693	20.0
Phenanthrene, 2-m...	11.87	3.2	ng	121270	4	11.35	764693	20.0
4H-Cyclopenta[def...	11.97	7.6	ng	289915	4	11.35	764693	20.0
Naphthalene, 2-ph...	12.15	3.5	ng	135182	4	11.35	764693	20.0
Phenanthrene, 2,5...	12.40	2.8	ng	107636	4	11.35	764693	20.0
11H-Benzo[a]fluorene	13.11	5.0	ng	299970	5	13.98	1211460	20.0
Benzo[e]pyrene	15.28	9.6	ng	380615	6	15.39	795177	20.0
unknown15.73	15.73	3.0	ng	118397	6	15.39	795177	20.0
unknown16.07	16.07	2.6	ng	103941	6	15.39	795177	20.0
Benzo[b]triphenylene	16.61	3.5	ng	139452	6	15.39	795177	20.0
1,2:4,5-Dibenzopy...	19.28	4.5	ng	177139	6	15.39	795177	20.0