LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
Data File : BF089676.D
Acq On : 8 Aug 2016 23:30
Operator : UM/SJ
Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : WC2-GRAB

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-BF080416.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	4.627	263	266	286	rVB	100246	246236	20.91%	1.235%
2	5.302	323	325	348	rBV	494640	967887	82.20%	4.856%
3	6.959	468	470	477	rVV	647626	988111	83.92%	4.957%
4	7.073	477	480	486	rVV	675528	1110675	94.33%	5.572%
5	7.153	486	487	492	rVB	18224	25451	2.16%	0.128%
6	7.416	508	510	517	rBV	132820	239306	20.32%	1.201%
7	7.519	517	519	522	rVB	8366	13684	1.16%	0.069%
8	7.656	528	531	543	rBV	628681	821801	69.80%	4.123%
9	7.965	555	558	561	rVB	10467	14356	1.22%	0.072%
10	8.330	587	590	600	rBV	410133	607267	51.58%	3.047%
11	8.673	616	620	622	rBV2	6587	14653	1.24%	0.074%
12	9.451	685	688	694	rBV	243560	378569	32.15%	1.899%
13	10.616	788	790	795	rBB	10471	17516	1.49%	0.088%
14	10.765	801	803	808	rBB	12409	19271	1.64%	0.097%
15	11.222	840	843	846	rBV	695173	1052040	89.35%	5.278%
16	11.736	886	888	891	rBV2	39739	52204	4.43%	0.262%
17	11.919	901	904	911	rBV	204167	269585	22.90%	1.353%
18	12.045	912	915	919	rVB	23910	39496	3.35%	0.198%
19	12.262	932	934	937	rBV	255100	409518	34.78%	2.055%
20	12.319	937	939	943	rVB	78112	100083	8.50%	0.502%
21	12.617	963	965	970	rBV	28268	53888	4.58%	0.270%
22	12.834	978	984	985	rBV2	4368	13179	1.12%	0.066%
23	13.131	1005	1010	1011	rBV	19746	30836	2.62%	0.155%
24	13.154	1011	1012	1018	rVV	57197	96726	8.21%	0.485%
25	13.268	1018	1022	1023	rVV3	7497	16654	1.41%	0.084%
26	13.440	1035	1037	1039	rBV	13990	24568	2.09%	0.123%
27	13.554	1045	1047	1053	rBV2	429785	622748	52.89%	3.124%
28	13.897	1075	1077	1082	rBV	24486	44089	3.74%	0.221%
29	14.057	1088	1091	1093	rBV2	10918	22633	1.92%	0.114%
30	14.102	1093	1095	1098	rVB2	7739	12105	1.03%	0.061%
31	14.285	1109	1111	1113	rBV	8874	14390	1.22%	0.072%
32	14.377	1117	1119	1122	rVV	22482	33375	2.83%	0.167%
33	14.491	1127	1129	1134	rVB	34290	45957	3.90%	0.231%
34	14.628	1139	1141	1142	rBV	15105	17332	1.47%	0.087%

LSC Area Percent Report

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Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : WC2-GRAB

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-BF080416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.662	1142	1144	1145	rVV	236986	316650	26.89%	1.589%
36	14.708	1145	1148	1150	rVV	752771	1177440	100.00%	5.907%
37	14.788	1152	1155	1161	rVV	132405	232507	19.75%	1.166%
38	15.085	1178	1181	1187	rBV	30692	64487	5.48%	0.324%
39	15.200	1189	1191	1193	rVB	11641	15660	1.33%	0.079%
40	15.314	1198	1201	1205	rVB2	9382	22656	1.92%	0.114%
41	15.463	1211	1214	1216	rBV	70665	91210	7.75%	0.458%
42	15.497	1216	1217	1222	rVV	72456	113828	9.67%	0.571%
43	15.577	1222	1224	1225	rVV	25578	34004	2.89%	0.171%
44	15.611	1225	1227	1229	rVV	100588	153057	13.00%	0.768%
45	15.680	1229	1233	1237	rVB2	74869	178552	15.16%	0.896%
46	15.908	1251	1253	1255	rVV	45514	66527	5.65%	0.334%
47	15.954	1255	1257	1261	rVB	41219	61246	5.20%	0.307%
48	16.125	1270	1272	1273	rBV	10253	14583	1.24%	0.073%
49	16.171	1273	1276	1277	rVV	11303	16941	1.44%	0.085%
50	16.263	1282	1284	1286	rBV	25699	39822	3.38%	0.200%
51	16.308	1286	1288	1293	rVV4	19510	48223	4.10%	0.242%
52	16.377	1293	1294	1299	rVB2	36579	63551	5.40%	0.319%
53	16.468	1299	1302	1305	rBV	842447	1144564	97.21%	5.742%
54	16.571	1309	1311	1313	rBV2	36292	61617	5.23%	0.309%
55	16.606	1313	1314	1317	rVB	71168	62315	5.29%	0.313%
56	16.663	1317	1319	1324	rBV3	25873	65988	5.60%	0.331%
57	16.777	1326	1329	1334	rVV	778358	1119852	95.11%	5.618%
58	16.857	1334	1336	1338	rVB2	18416	25053	2.13%	0.126%
59	16.960	1343	1345	1347	rBV	36847	44305	3.76%	0.222%
60	17.028	1348	1351	1354	rVV	568304	689523	58.56%	3.459%
61	17.097	1354	1357	1360	rVB2	29875	57571	4.89%	0.289%
62	17.143	1360	1361	1363	rVB	11173	12336	1.05%	0.062%
63	17.211	1363	1367	1368	rBV2	32340	62455	5.30%	0.313%
64	17.234	1368	1369	1373	rVB	52287	67436	5.73%	0.338%
65	17.326	1373	1377	1379	rBV	43220	58163	4.94%	0.292%
66	17.371	1379	1381	1389	rBV3	60203	156759	13.31%	0.786%
67	17.486	1389	1391	1393	rVB	26639	29259	2.48%	0.147%
68	17.520	1393	1394	1398	rBV	13917	26027	2.21%	0.131%
69	17.691	1407	1409	1412	rVB	77802	113168	9.61%	0.568%
70	17.886	1424	1426	1427	rBV2	12015	21699	1.84%	0.109%
71	17.909	1427	1428	1431	rVV	37358	56467	4.80%	0.283%

LSC Area Percent Report

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Acq On : 8 Aug 2016 23:30
Operator : UM/SJ
Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : WC2-GRAB

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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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	18.046	1437	1440	1441	rBV	44294	82845	7.04%	0.416%
73	18.137	1446	1448	1451	rVV2	40552	61237	5.20%	0.307%
74	18.194	1451	1453	1455	rVV	50406	79624	6.76%	0.399%
75	18.240	1455	1457	1458	rVV	12662	17533	1.49%	0.088%
76	18.297	1460	1462	1464	rVV	38347	68404	5.81%	0.343%
77	18.354	1464	1467	1469	rVV2	425255	776338	65.93%	3.895%
78	18.400	1469	1471	1474	rVV	359841	511728	43.46%	2.567%
79	18.491	1477	1479	1482	rVB	55099	73765	6.26%	0.370%
80	18.640	1490	1492	1494	rVB2	26755	35750	3.04%	0.179%
81	18.686	1494	1496	1499	rBV2	24441	37528	3.19%	0.188%
82	18.869	1506	1512	1514	rBV2	61616	134645	11.44%	0.676%
83	18.983	1520	1522	1526	rBV4	39441	85499	7.26%	0.429%
84	19.280	1546	1548	1552	rBV4	28327	61540	5.23%	0.309%
85	19.463	1562	1564	1567	rBV	44151	73032	6.20%	0.366%
86	19.623	1576	1578	1584	rBV	364826	787261	66.86%	3.950%
87	19.737	1586	1588	1591	rVB	62443	79855	6.78%	0.401%
88	19.806	1592	1594	1596	rBV	36457	59016	5.01%	0.296%
89	19.909	1601	1603	1607	rBV	183228	265594	22.56%	1.332%
90	19.977	1607	1609	1612	rBV	247882	345957	29.38%	1.736%
91	20.034	1612	1614	1615	rBV	161879	248461	21.10%	1.247%
92	20.492	1652	1654	1657	rVB	120547	163532	13.89%	0.820%
93	21.200	1714	1716	1721	rVB2	149099	284589	24.17%	1.428%
94	21.349	1727	1729	1731	rVV	22948	34131	2.90%	0.171%
95	21.395	1731	1733	1737	rVB	31757	54750	4.65%	0.275%
96	21.543	1743	1746	1751	rBV	115968	207902	17.66%	1.043%
97	21.737	1761	1763	1769	rVB2	26068	68693	5.83%	0.345%
98	23.269	1893	1897	1906	rBV	20287	81340	6.91%	0.408%
99	23.440	1908	1912	1928	rVB3	15994	105084	8.92%	0.527%
100	24.138	1970	1973	1976	rBV3	8675	24927	2.12%	0.125%

Sum of corrected areas: 19932250

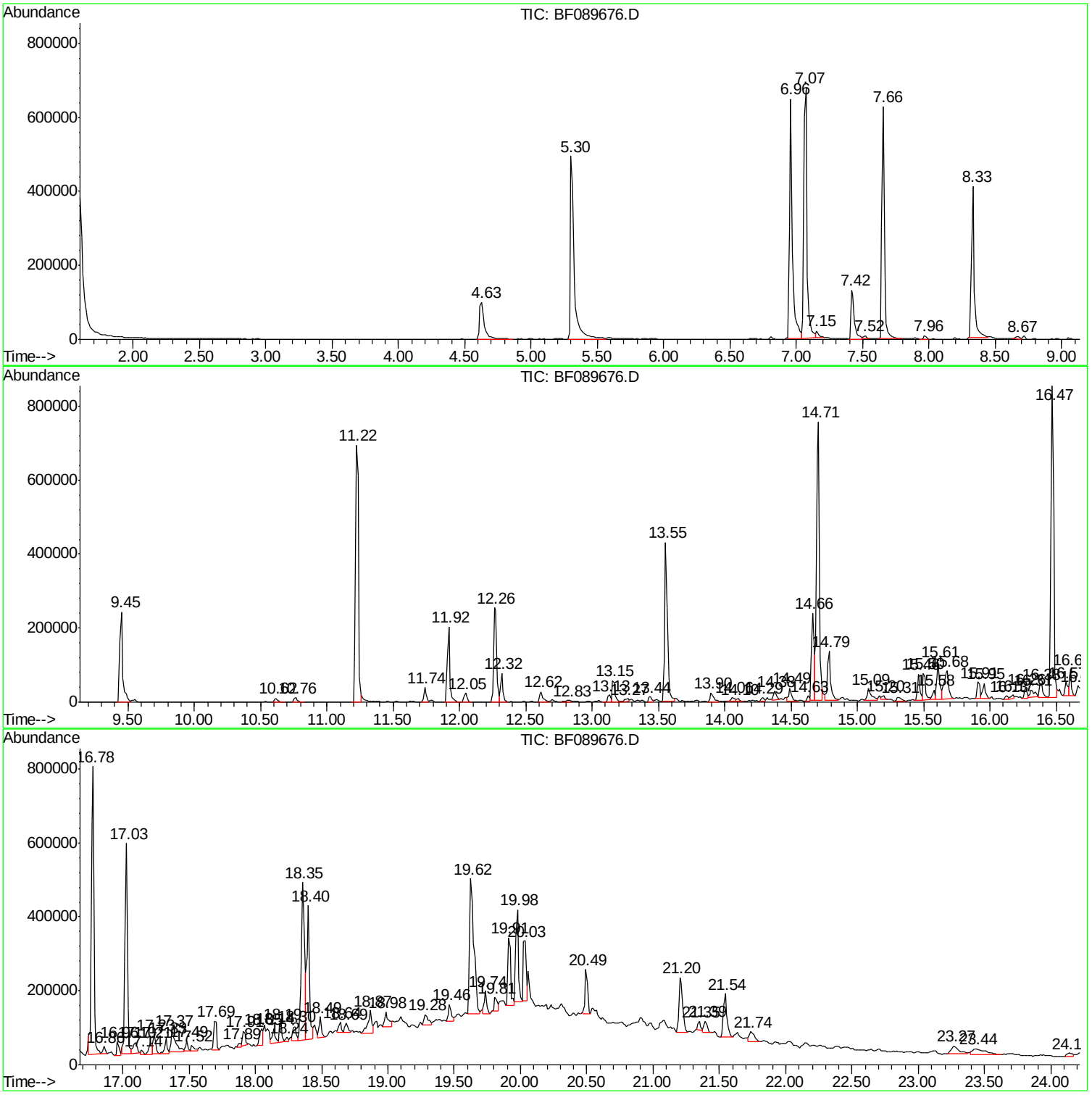
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
Data File : BF089676.D
Acq On : 8 Aug 2016 23:30
Operator : UM/SJ
Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : WC2-GRAB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.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\BF080816\
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

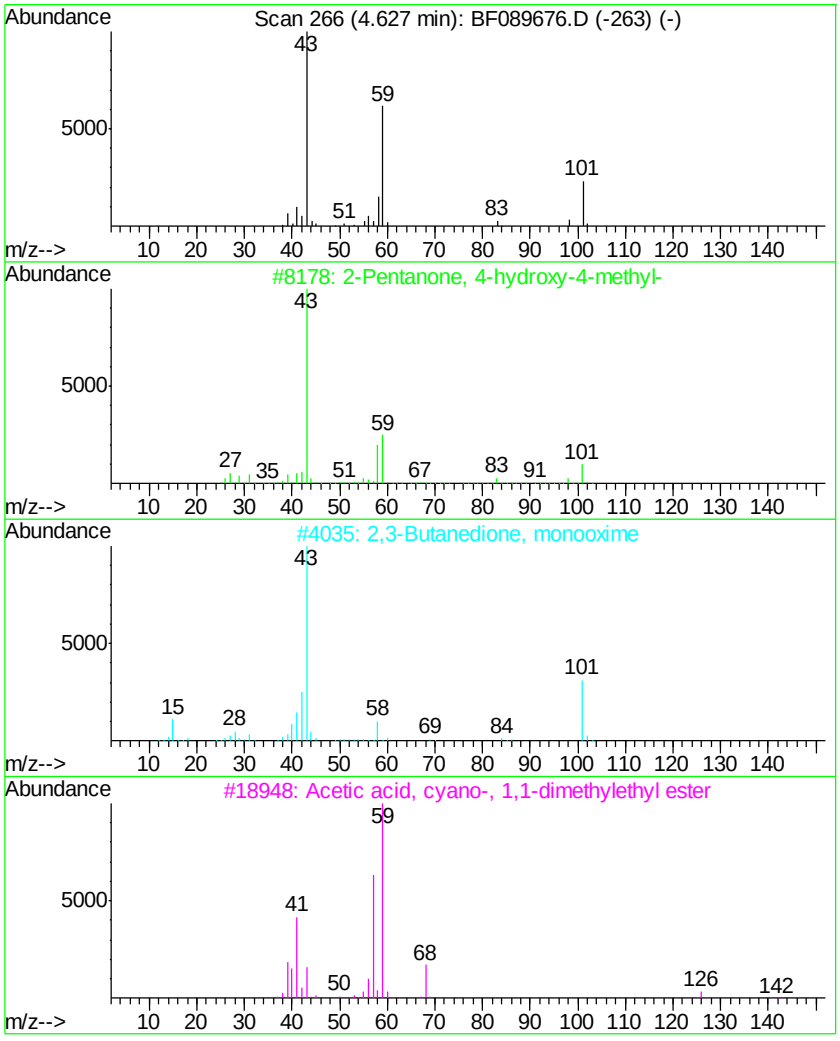
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

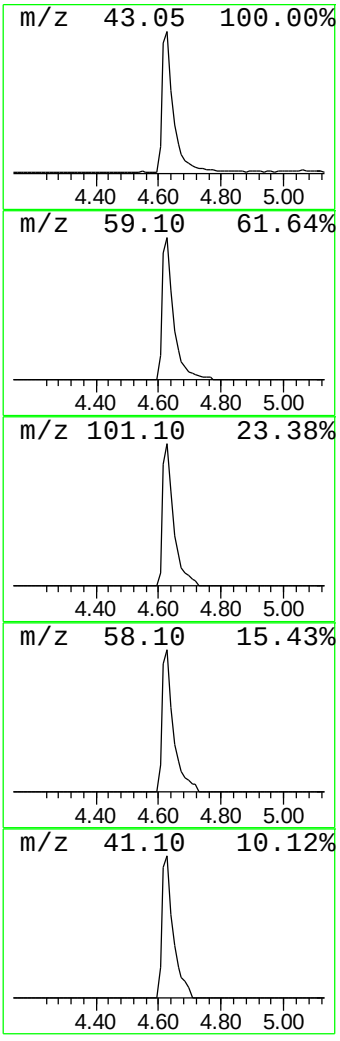
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.63	20.58 ng	246236	1,4-Dichlorobenzene-d4	7.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
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Sample : H4331-12
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Instrument : BNA_F
ClientSampleId : WC2-GRAB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

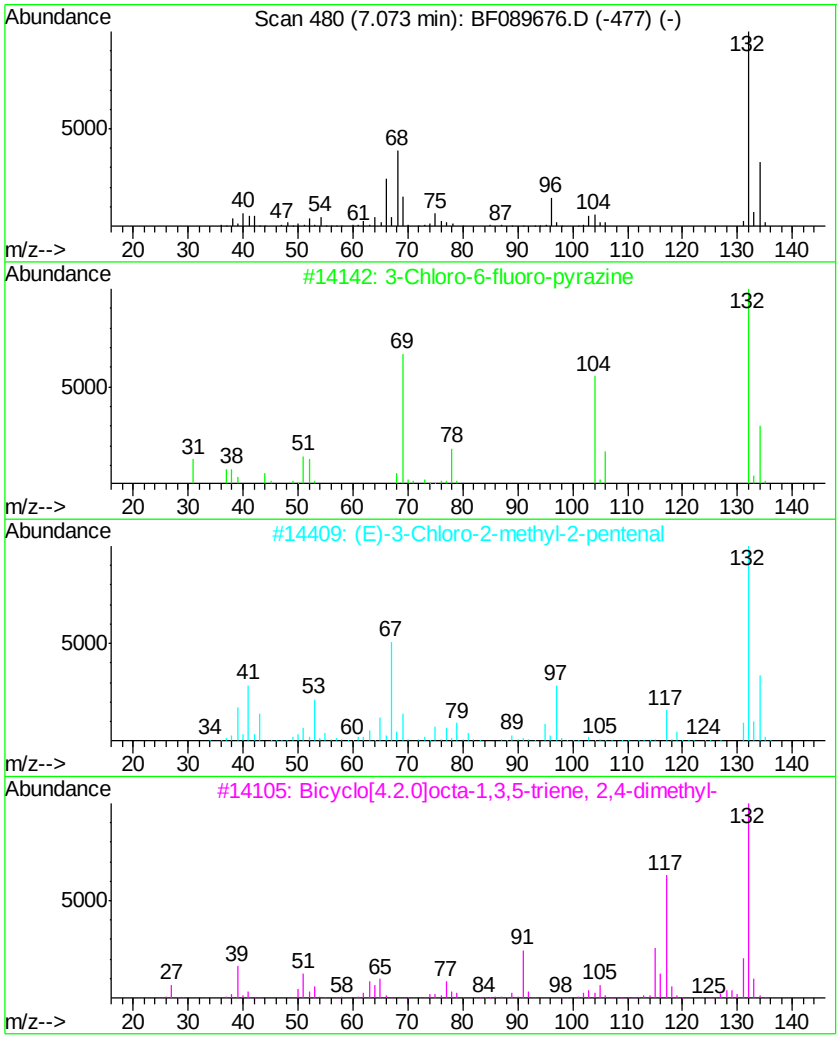
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

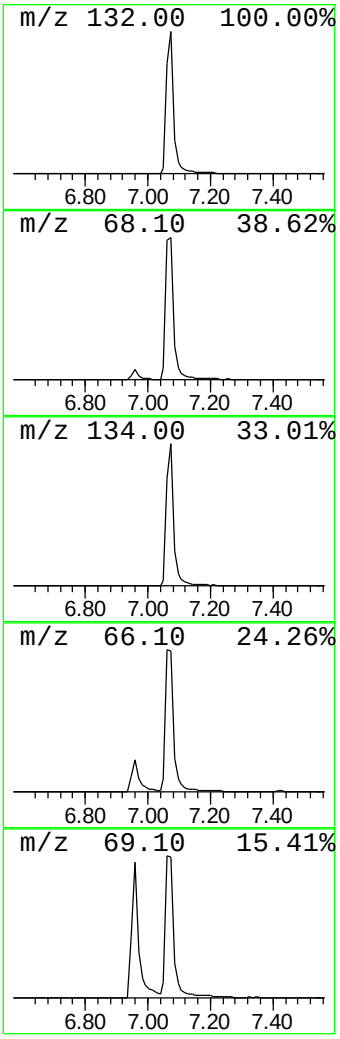
****

Peak Number 2 unknown7.07 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	92.82 ng	1110680	1,4-Dichlorobenzene-d4	7.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

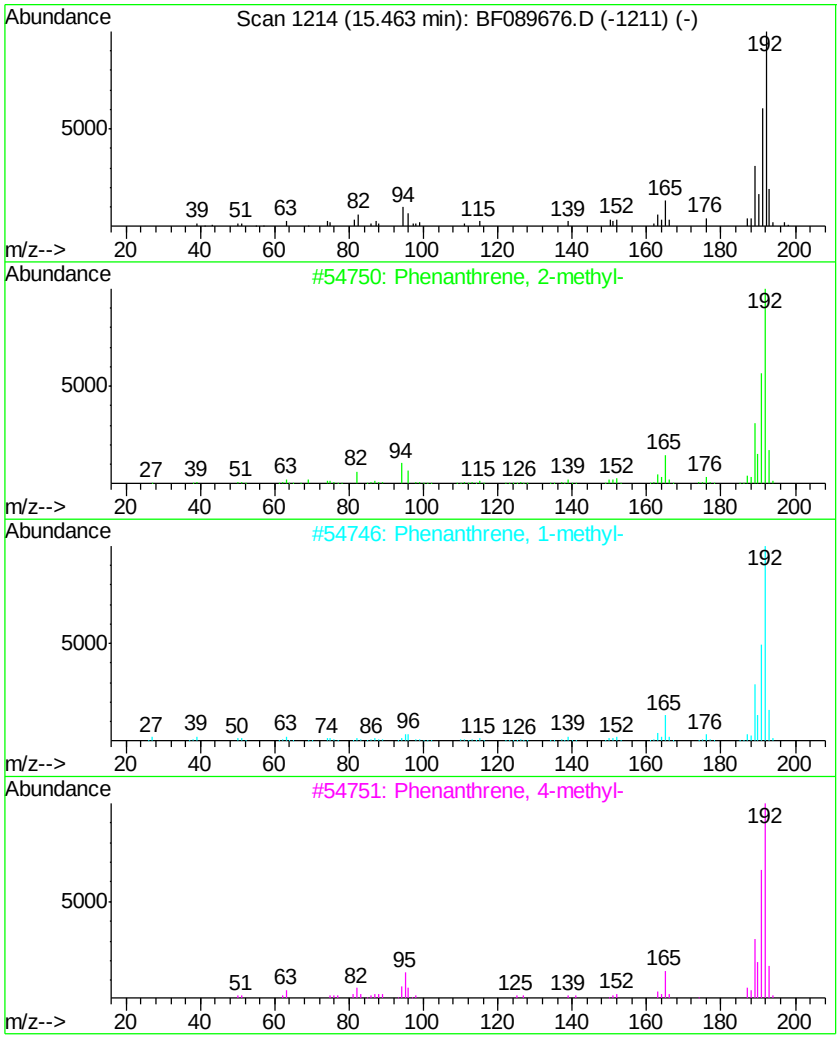
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

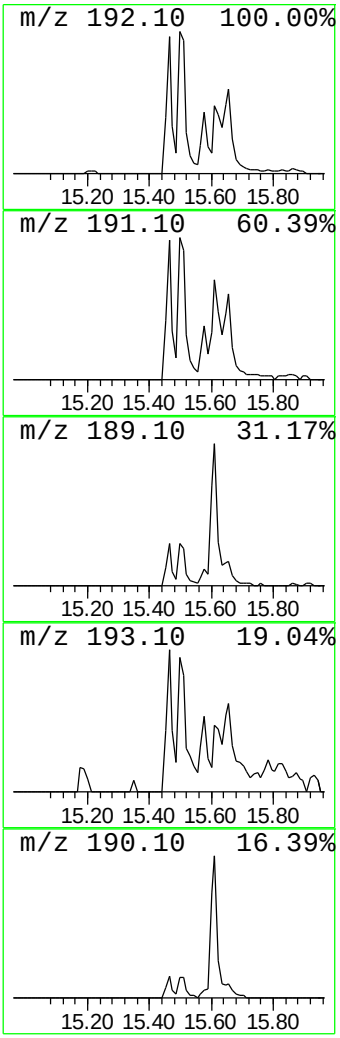
****

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	5.76 ng	91210	Phenanthrene-d10	14.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
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Operator : UM/SJ
Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.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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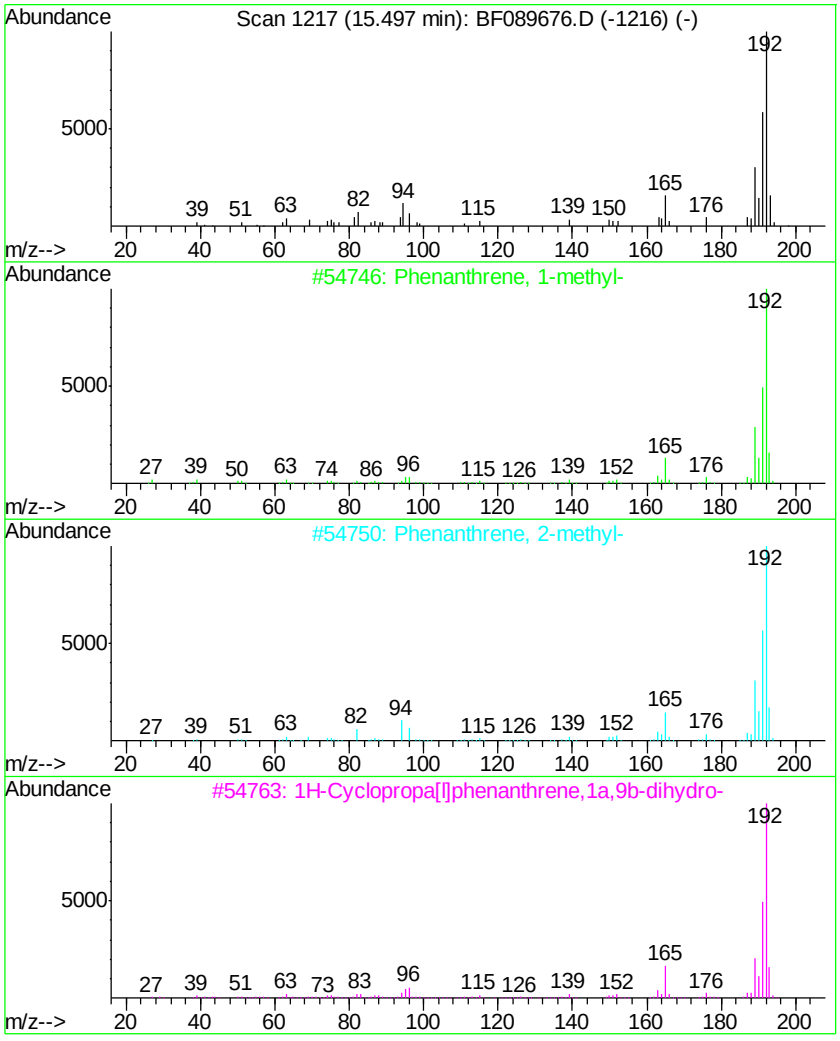
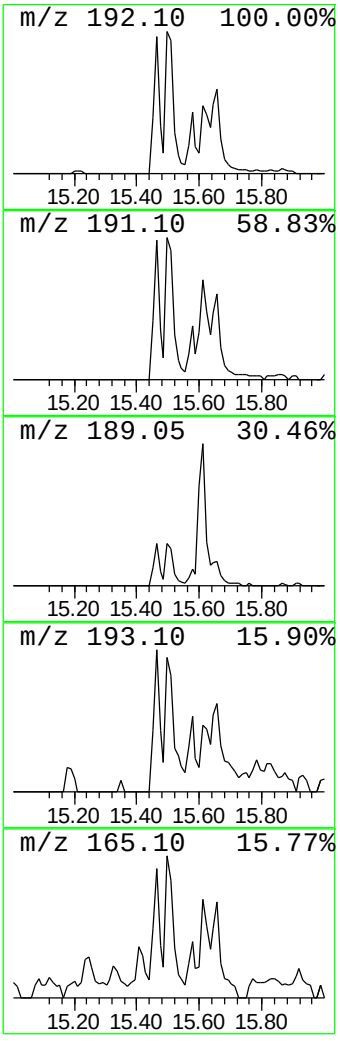
Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.50	7.19 ng	113828	Phenanthrene-d10	14.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
Data File : BF089676.D
Acq On : 8 Aug 2016 23:30
Operator : UM/SJ
Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : WC2-GRAB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.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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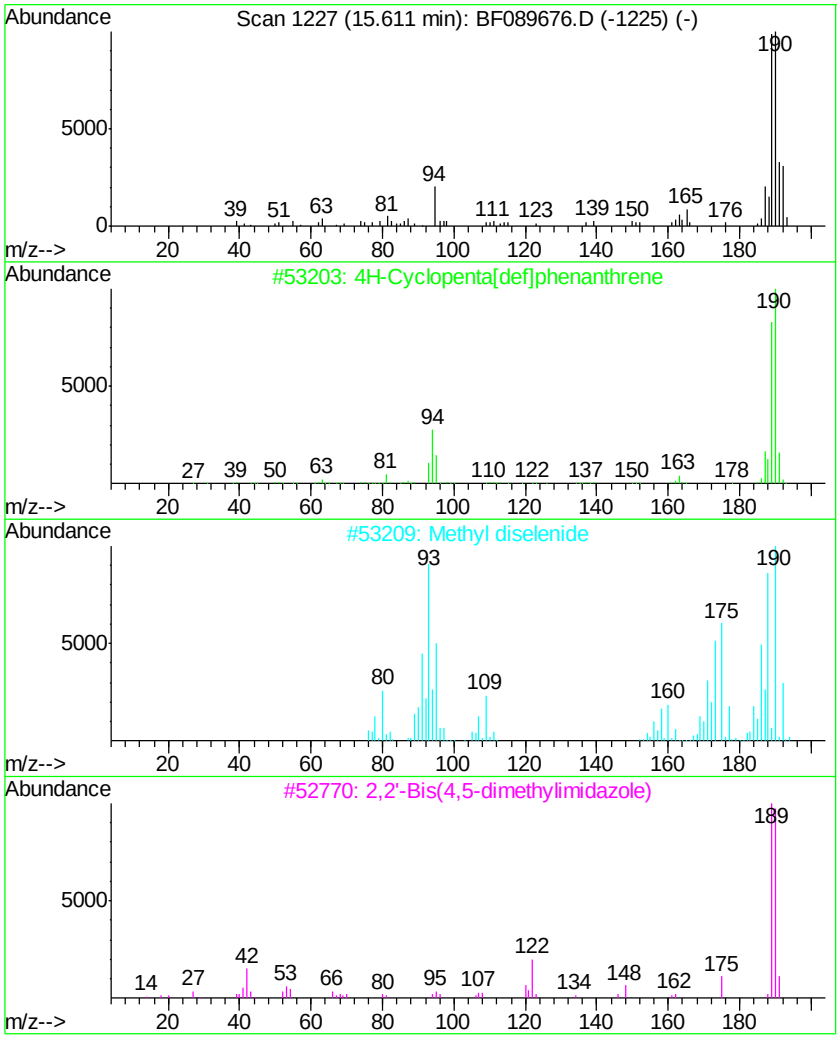
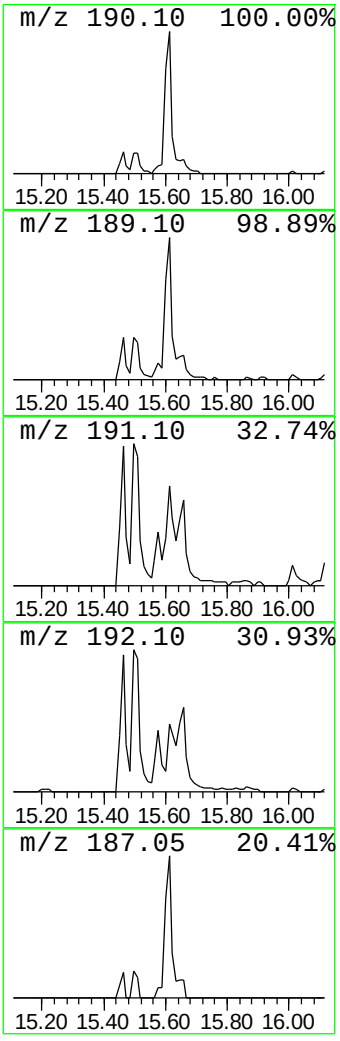
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	9.67 ng	153057	Phenanthrene-d10	14.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2	Methyl diselenide	190	C2H6Se2	007101-31-7	49
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	47
4	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	45
5	Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	28





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
Data File : BF089676.D
Acq On : 8 Aug 2016 23:30
Operator : UM/SJ
Sample : H4331-12
Misc :
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Instrument : BNA_F
ClientSampleId : WC2-GRAB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.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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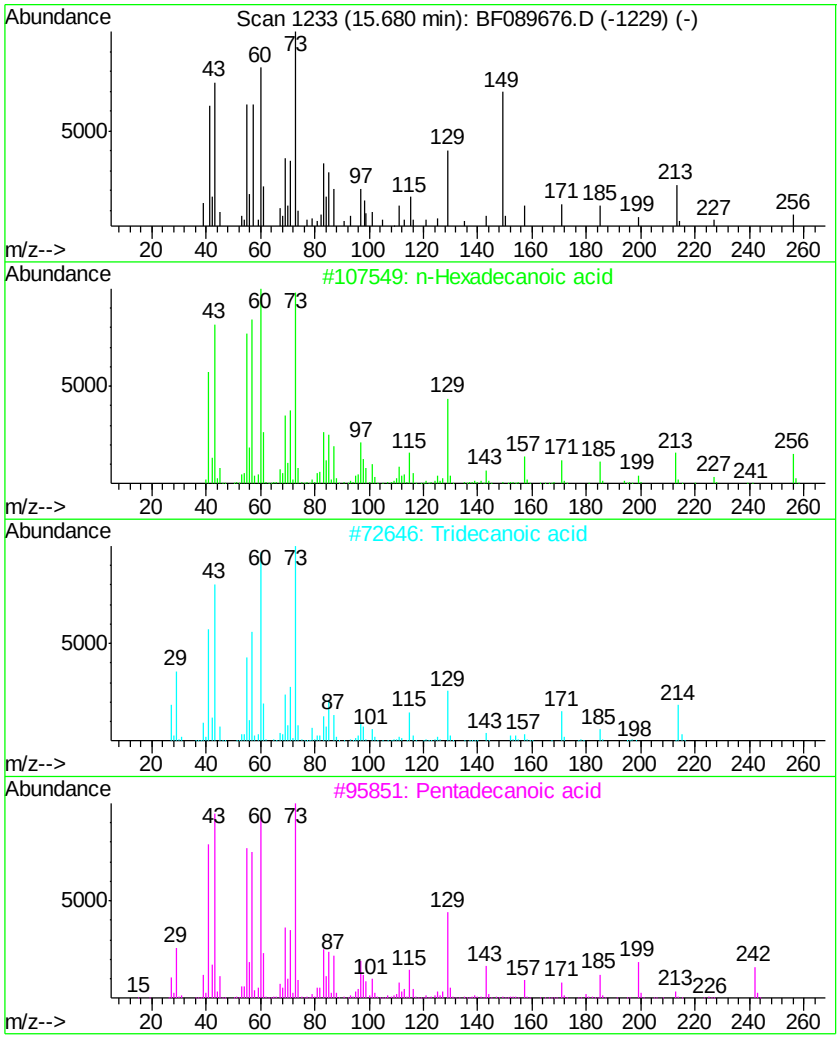
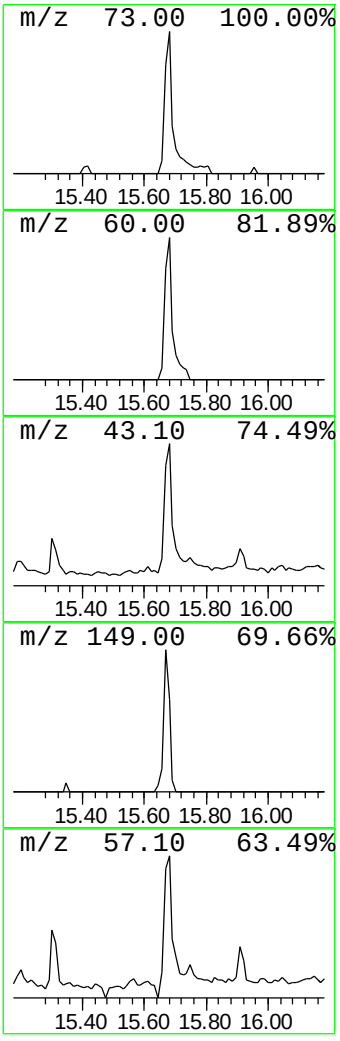
Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	11.28 ng	178552	Phenanthrene-d10	14.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	76
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
5	n-Decanoic acid	172	C10H20O2	000334-48-5	47





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ClientSampleId : WC2-GRAB

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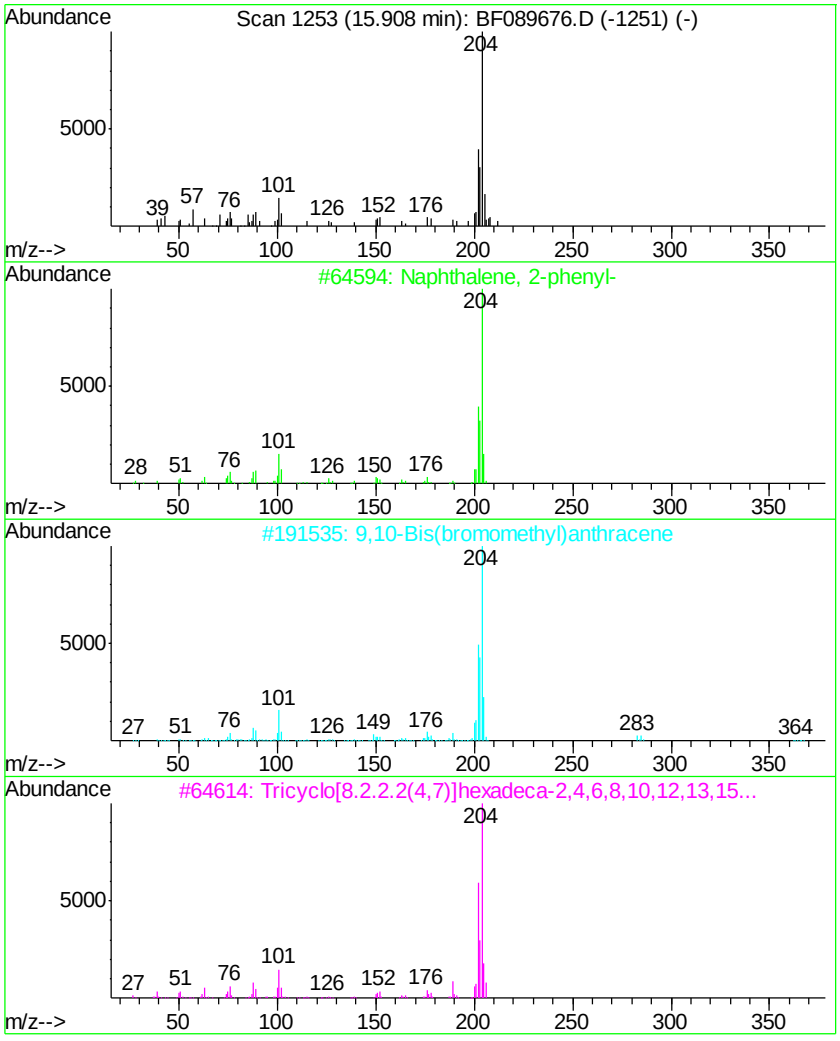
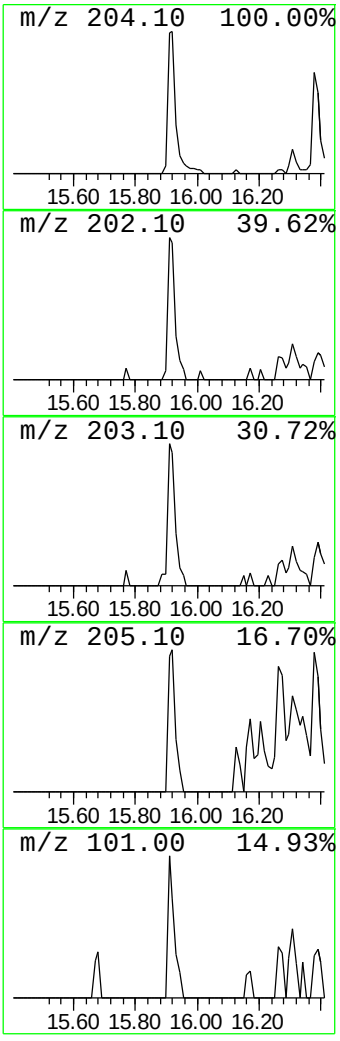
Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	4.20 ng	66527	Phenanthrene-d10	14.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	87
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	86
4	5H-Dibenzo[a,d]cycloheptene, 5-m...	204	C16H12	002975-79-3	78
5	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	60





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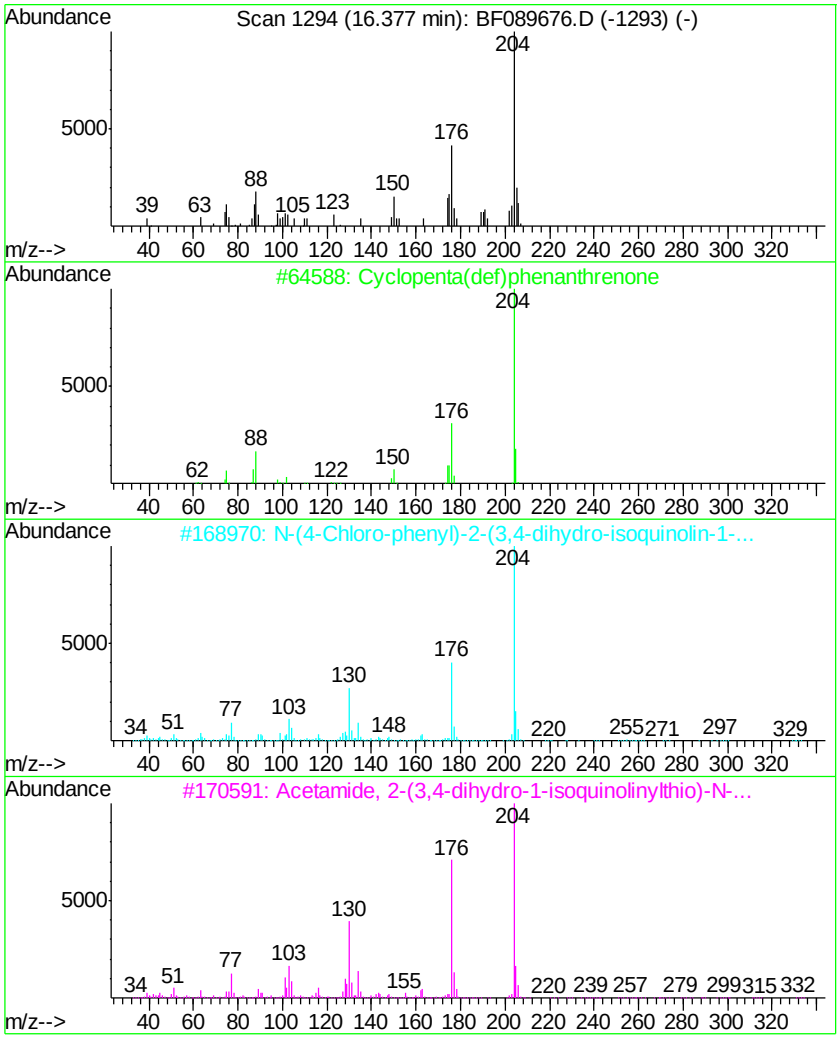
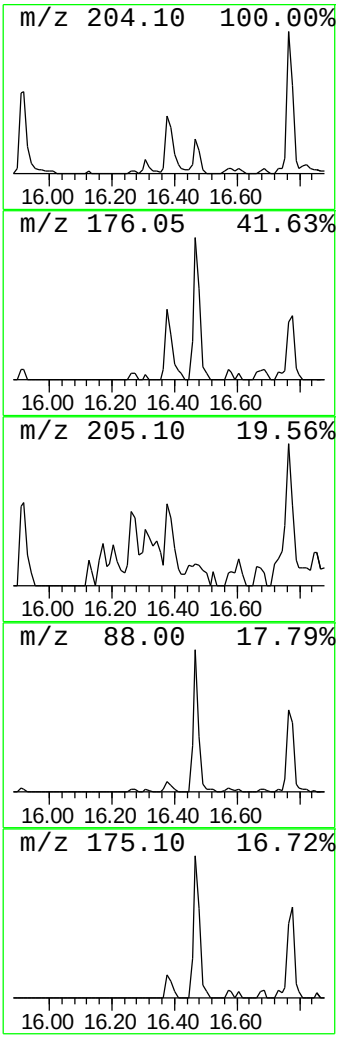
Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	4.01 ng	63551	Phenanthrene-d10	14.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	90
2	N-(4-Chloro-phenyl)-2-(3,4-dihyd...	330	C17H15ClN2OS	1000296-34-3	64
3	Acetamide, 2-(3,4-dihydro-1-isoq...	332	C17H14F2N2OS	1000270-76-1	59
4	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	52
5	6-Cyanophenanthridine	204	C14H8N2	002176-28-5	49





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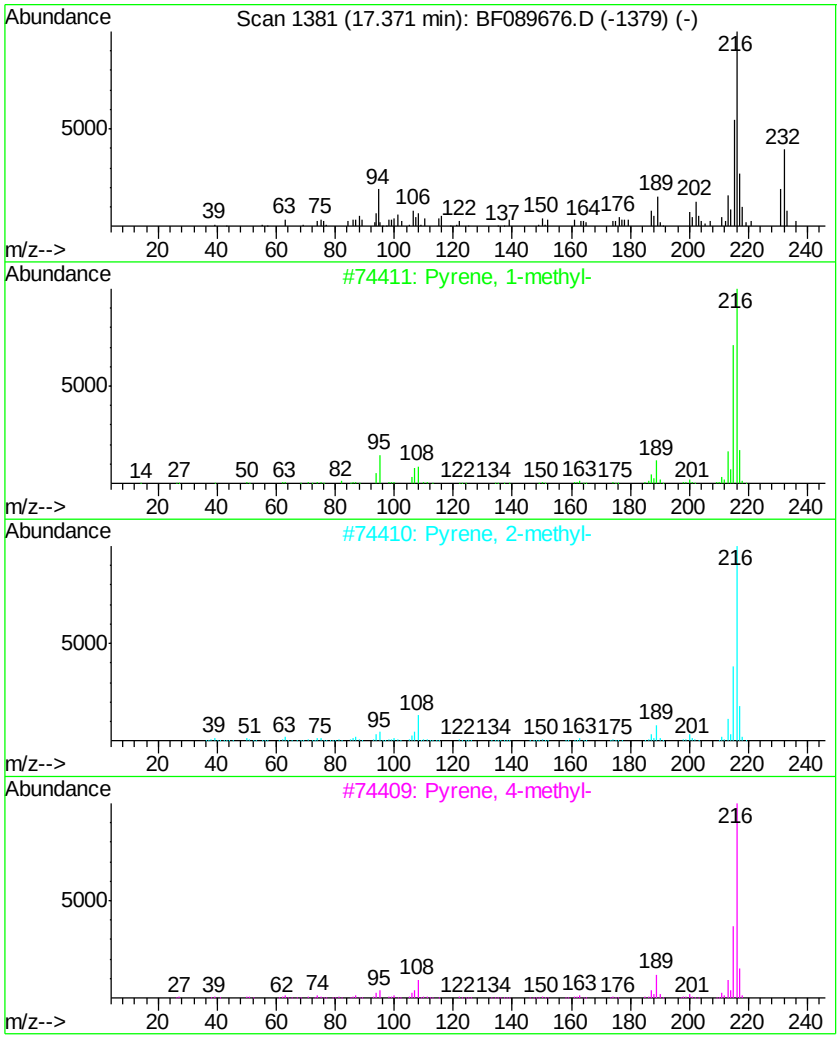
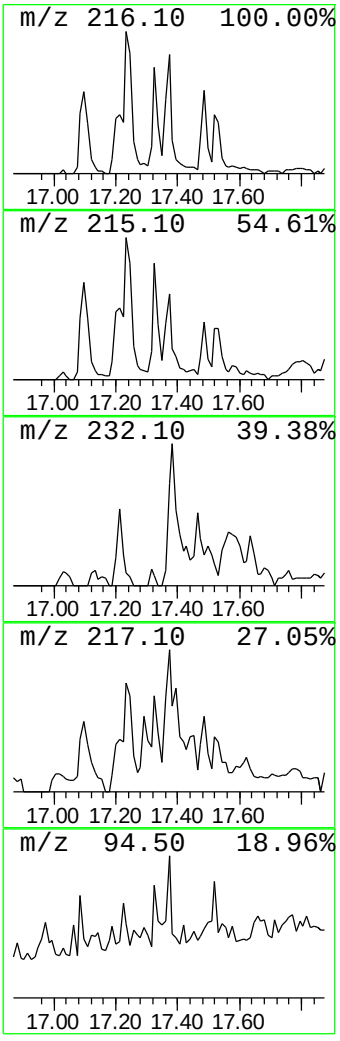
Peak Number 10 Pyrene, 1-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	4.04 ng	156759	Chrysene-d12	18.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	60
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	53
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	49
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	46
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	46





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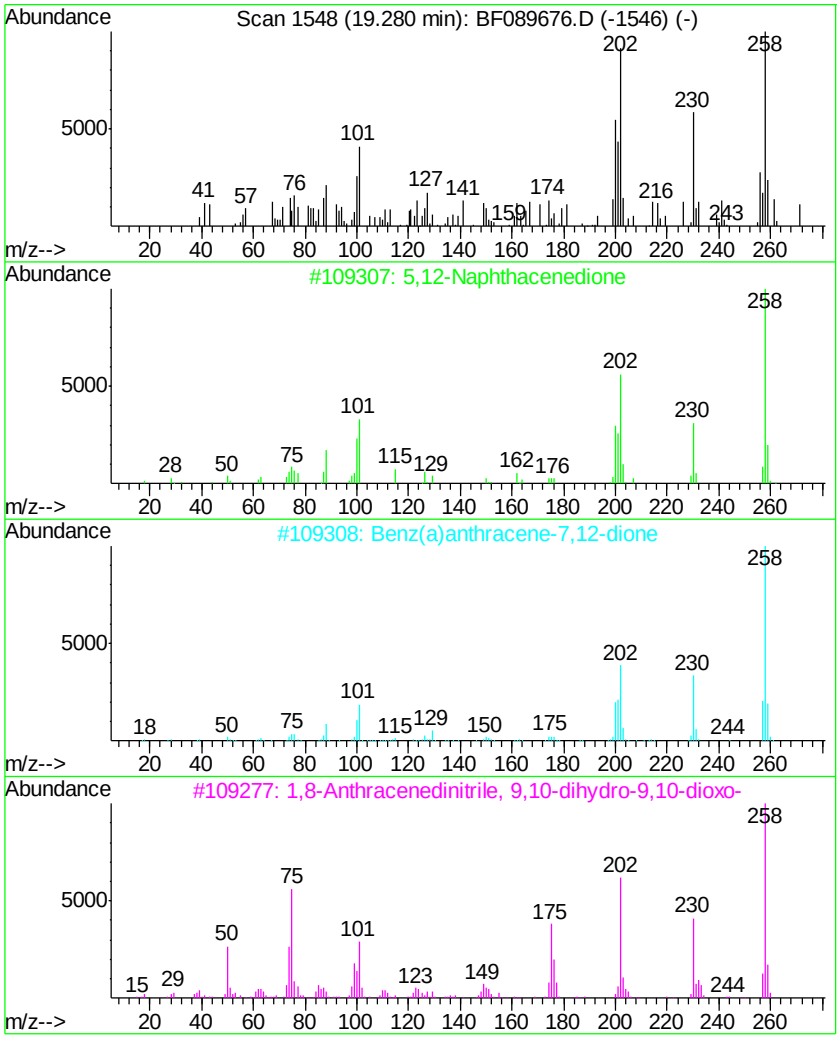
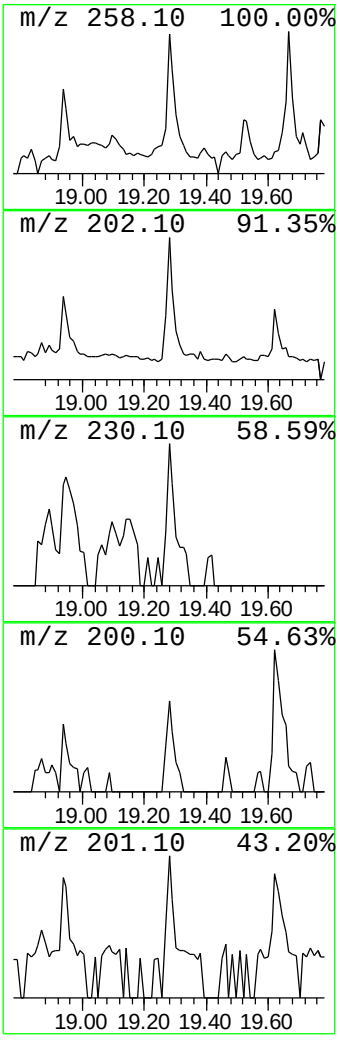
Peak Number 11 5,12-Naphthacenedione Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.28	4.95 ng	61540	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,12-Naphthacenedione	258	C18H10O2	001090-13-7	93
2	Benz(a)anthracene-7,12-dione	258	C18H10O2	002498-66-0	93
3	1,8-Anthracenedinitrile, 9,10-di...	258	C16H6N2O2	090866-50-5	86
4	1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	50
5	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	50





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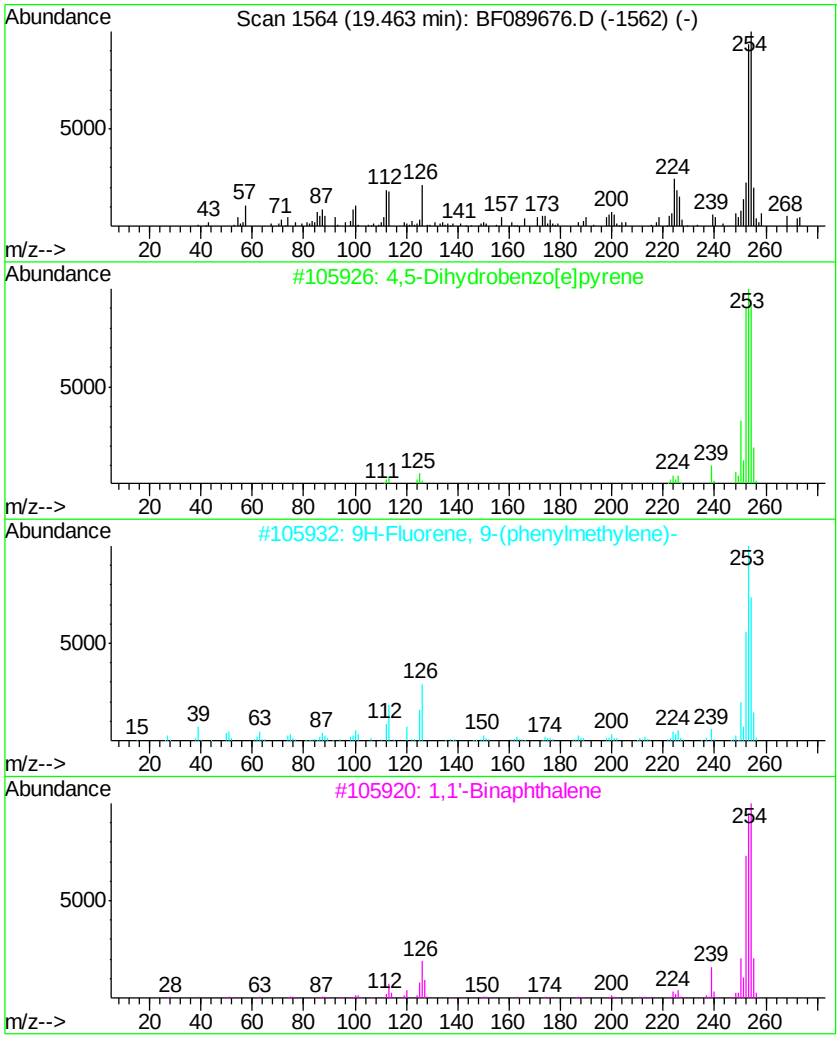
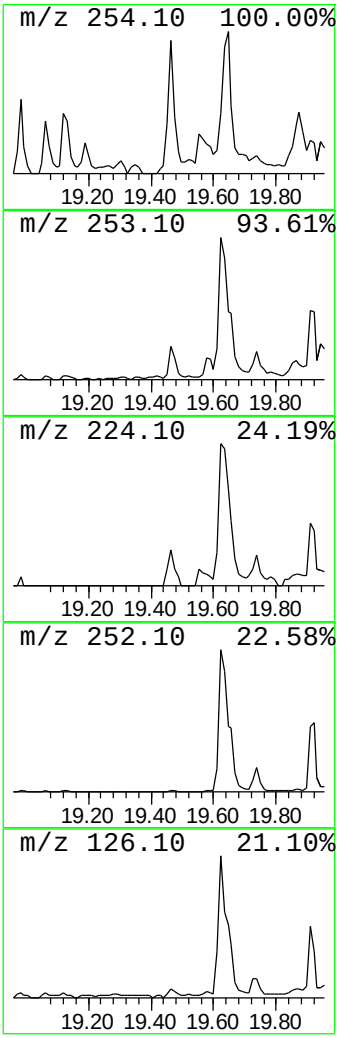
Peak Number 12 4,5-Dihydrobenzo[e]pyrene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.46	5.88 ng	73032	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,5-Dihydrobenzo[e]pyrene	254	C20H14	095676-42-9	76
2	9H-Fluorene, 9-(phenylmethylene)-	254	C20H14	001836-87-9	74
3	1,1'-Binaphthalene	254	C20H14	000604-53-5	72
4	4,6'-Biazulenyl	254	C20H14	094154-49-1	72
5	Benzo[a]pyrene, 4,5-dihydro-	254	C20H14	057652-66-1	70





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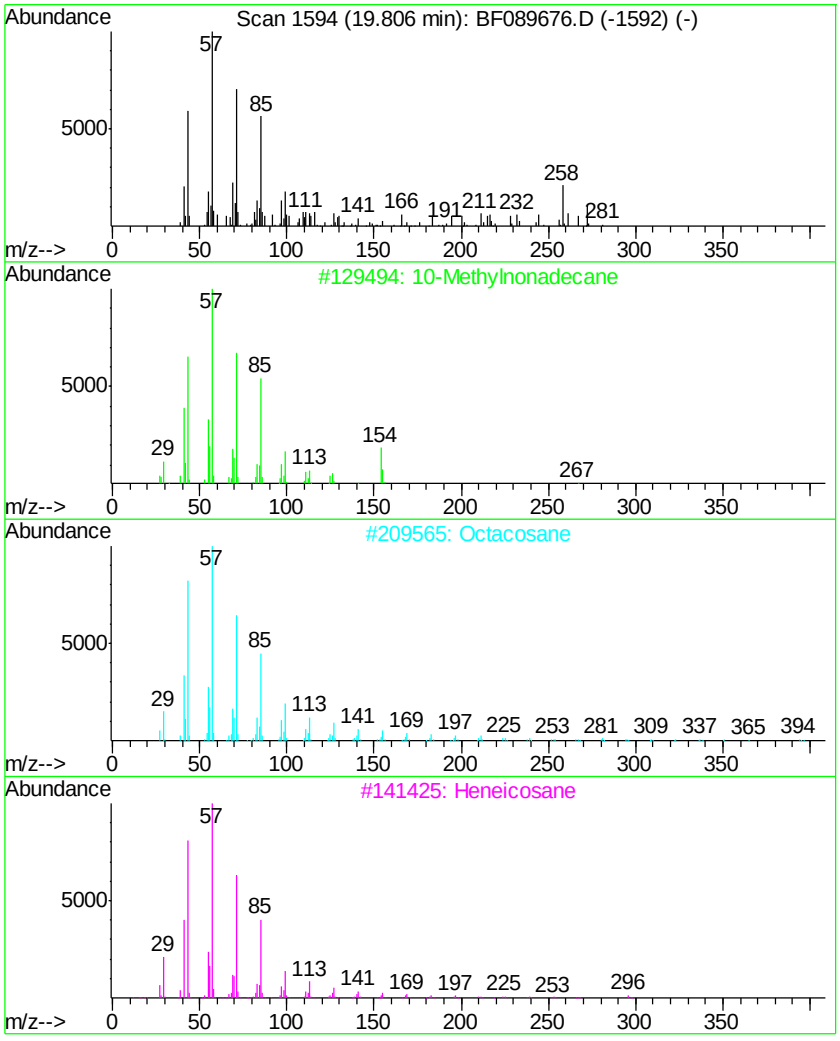
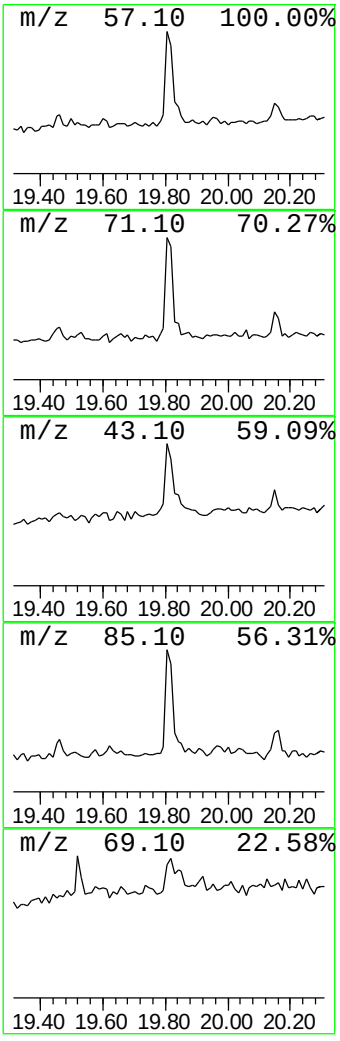
Peak Number 14 unknown19.81 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.81	4.75 ng	59016	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane	282	C20H42	056862-62-5	46
2	Octacosane	394	C28H58	000630-02-4	43
3	Heneicosane	296	C21H44	000629-94-7	43
4	Decane, 1-iodo-	268	C10H21I	002050-77-3	43
5	Pentacosane	352	C25H52	000629-99-2	43





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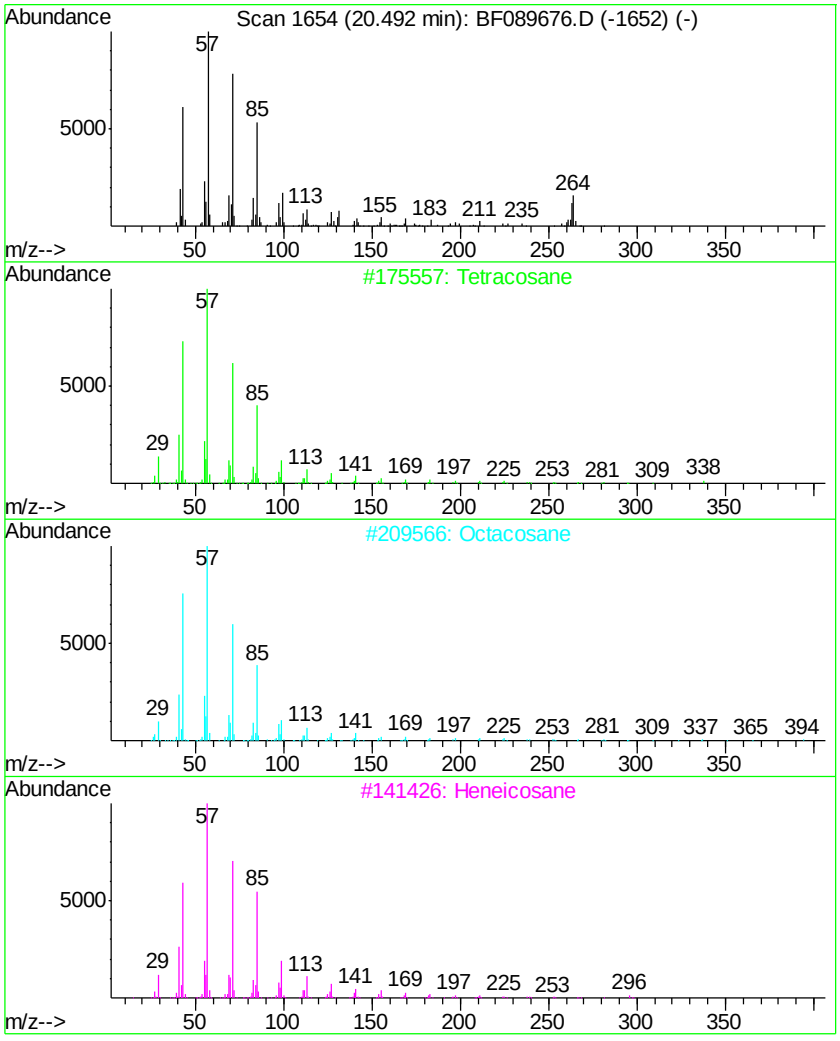
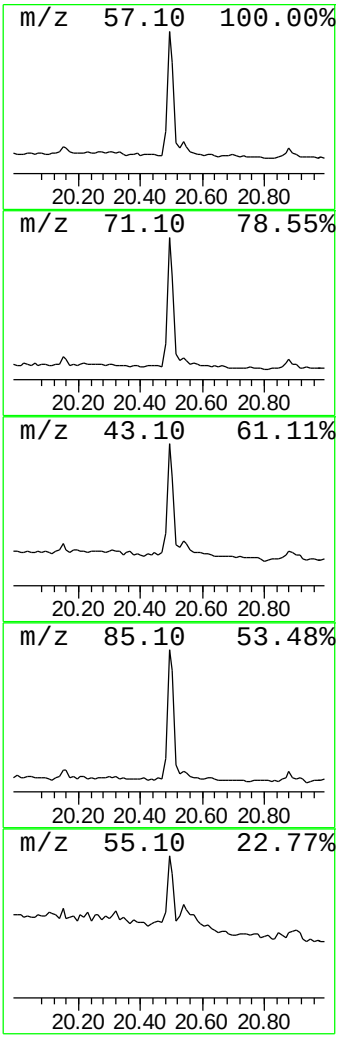
Peak Number 16 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.49	13.16 ng	163532	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	87
2	Octacosane	394	C28H58	000630-02-4	87
3	Heneicosane	296	C21H44	000629-94-7	87
4	Hexacosane	366	C26H54	000630-01-3	87
5	Heptacosane	380	C27H56	000593-49-7	83





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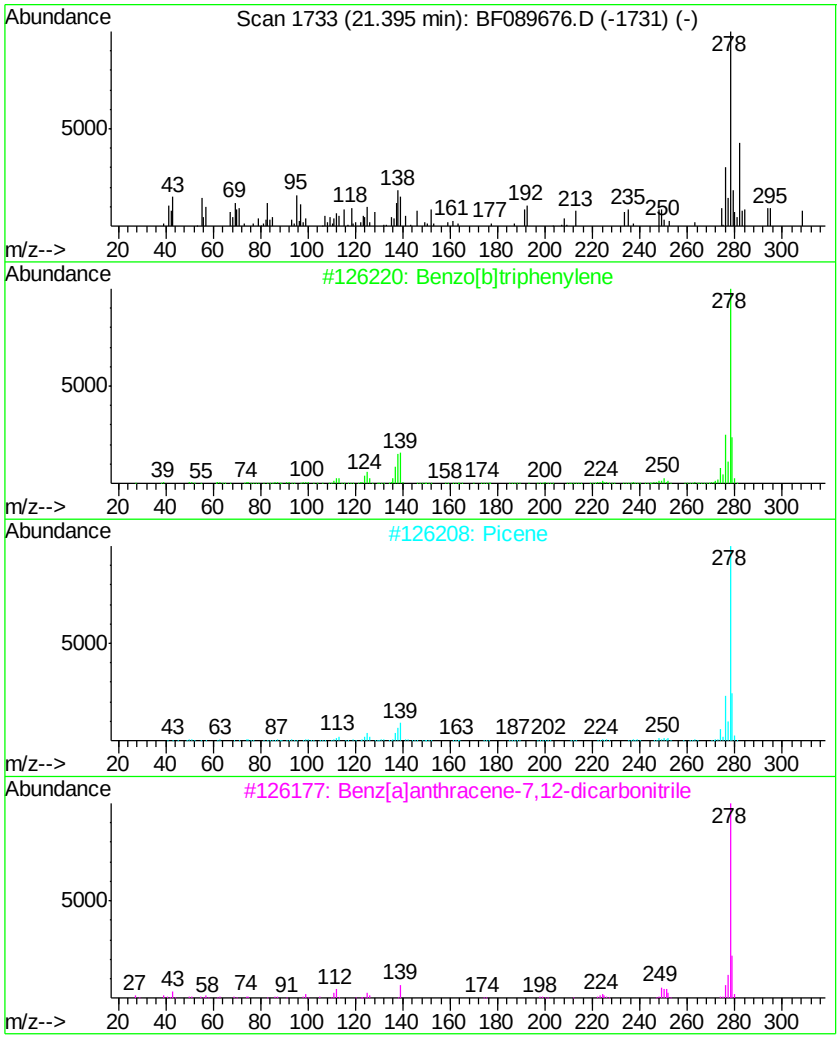
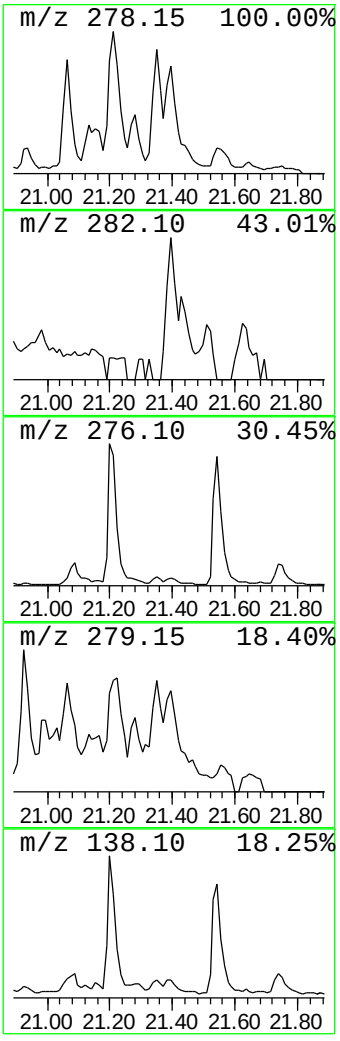
Peak Number 17 unknown21.39 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.39	4.41 ng	54750	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]triphenylene	278	C22H14	000215-58-7	46
2	Picene	278	C22H14	000213-46-7	41
3	Benz[a]anthracene-7,12-dicarboni...	278	C20H10N2	035215-32-8	38
4	Dibenz[a,h]anthracene	278	C22H14	000053-70-3	38
5	Estra-1,3,5,7,9,15-hexaen-17-one...	278	C19H18O2	056588-53-5	38





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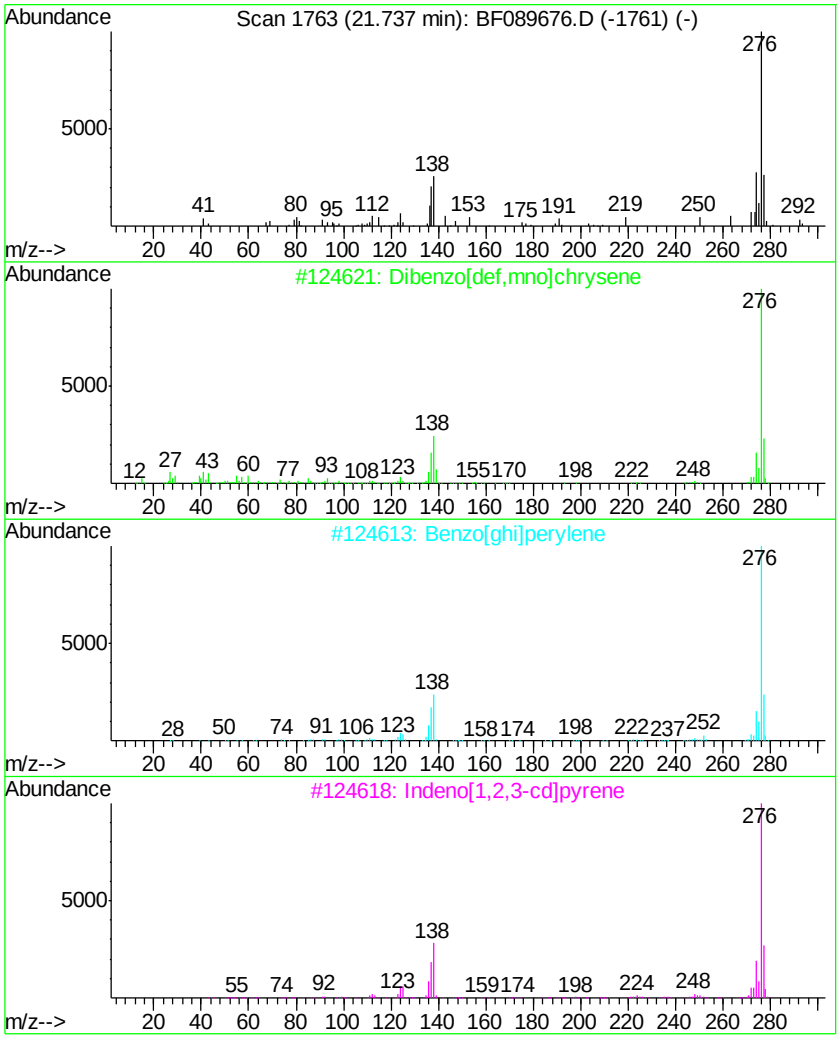
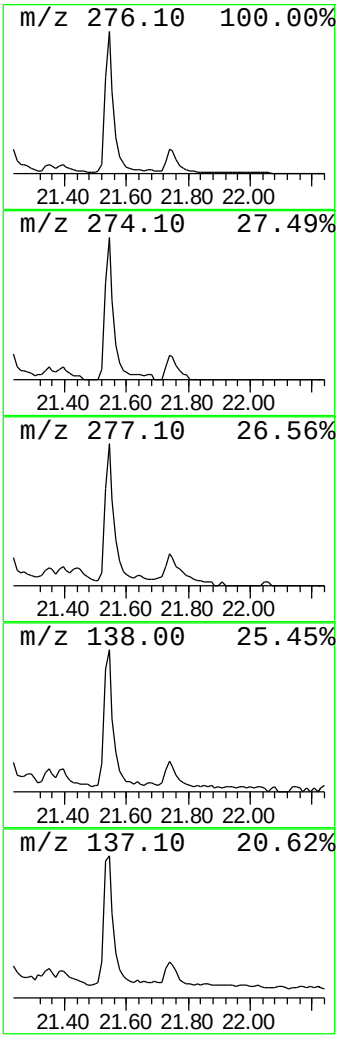
Peak Number 18 Dibenzo[def,mno]chrysene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.74	5.53 ng	68693	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	94
2	Benzo[ghi]perylene	276	C22H12	000191-24-2	93
3	Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	91
4	Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	81
5	Naphthalen-1-yl-(3-nitro-benzyli...	276	C17H12N2O2	200421-53-0	43





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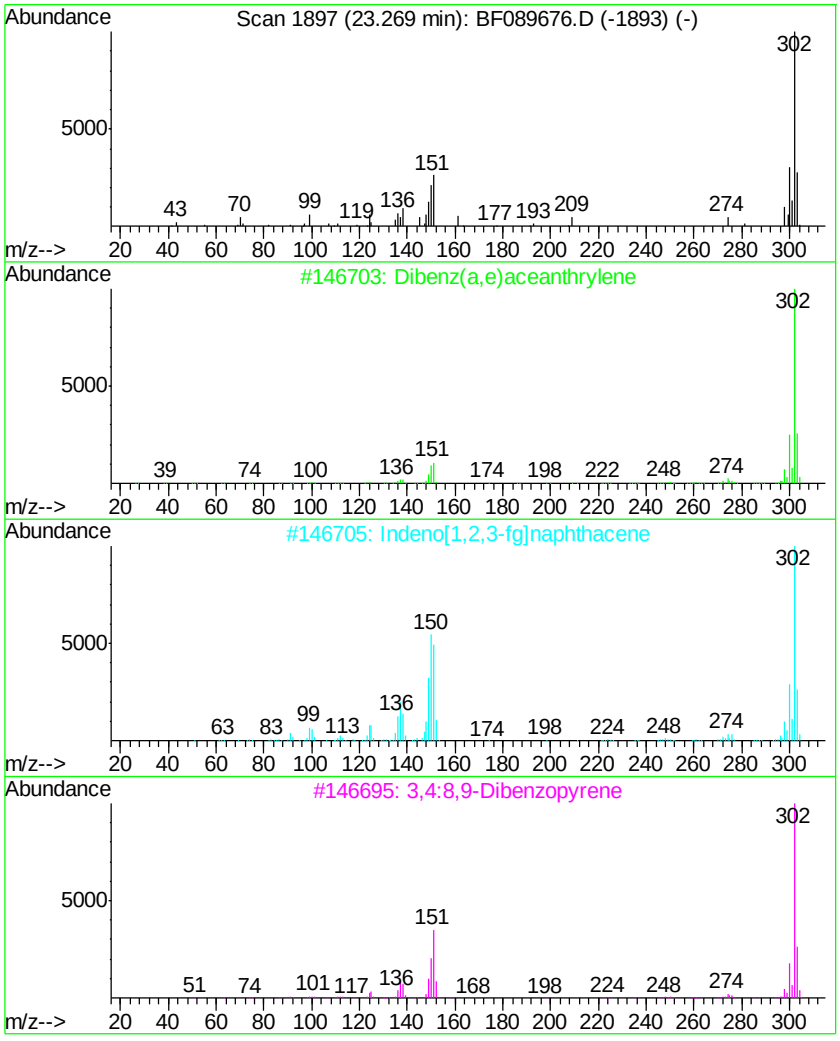
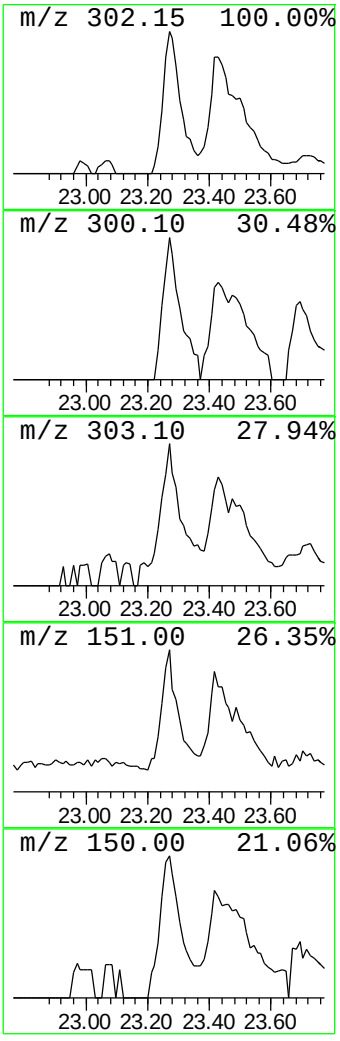
Peak Number 19 Dibenz(a,e)aceanthrylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.27	6.55 ng	81340	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	87
2	Indeno[1,2,3-fg]naphthacene	302	C24H14	000203-11-2	87
3	3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	74
4	1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	74
5	Dibenzo(a,i)pyrene	302	C24H14	000189-55-9	74





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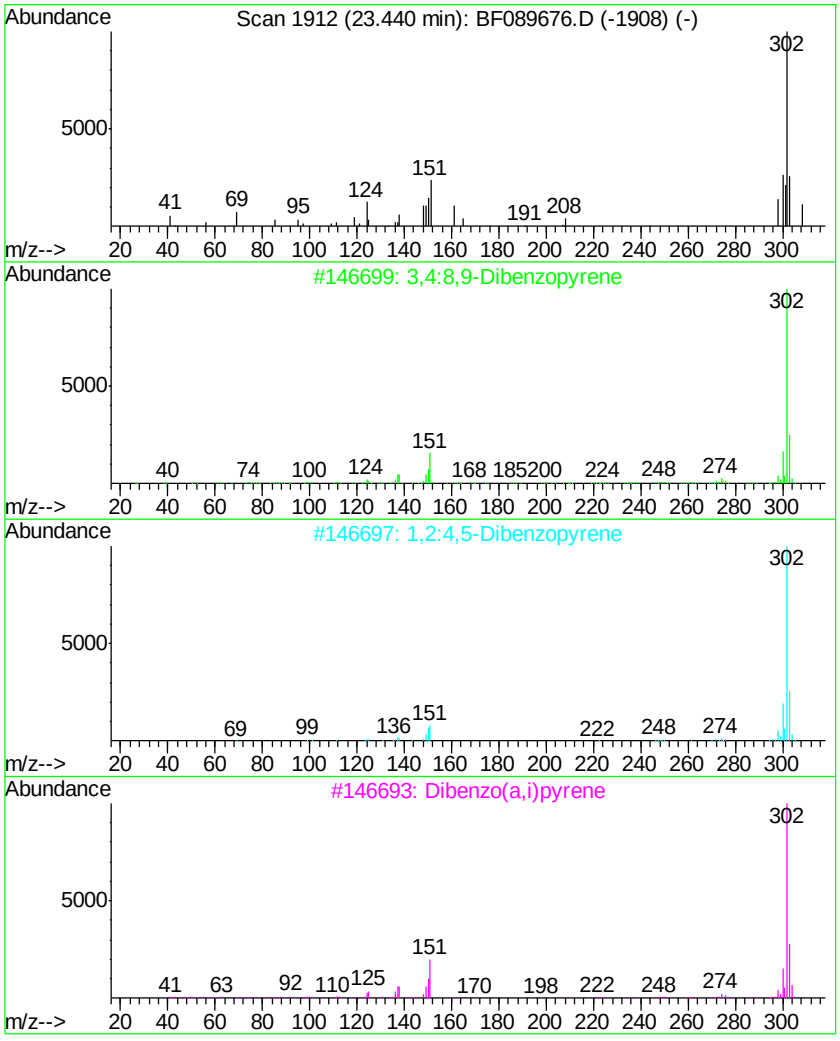
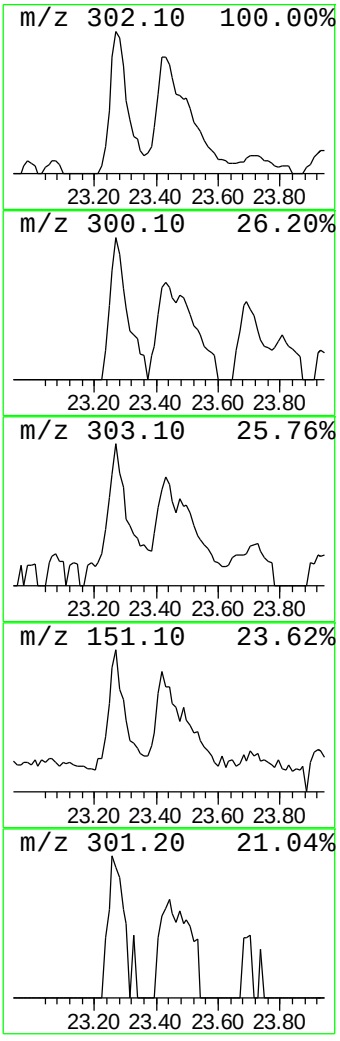
Peak Number 20 3,4:8,9-Dibenzopyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.44	8.46 ng	105084	Perylene-d12	20.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	52
2	1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	50
3	Dibenzo(a,i)pyrene	302	C24H14	000189-55-9	50
4	1H-Imidazo[4,5-b]phenazine, 2-(2...	302	C17H10N4S	1000319-15-8	47
5	1,2,9,10-Dibenzopyrene	302	C24H14	000191-30-0	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080816\
Data File : BF089676.D
Acq On : 8 Aug 2016 23:30
Operator : UM/SJ
Sample : H4331-12
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : WC2-GRAB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.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
2-Pentanone, 4-hy...	4.63	20.6	ng	246236	1	7.43	239306	20.0
unknown7.07	7.07	92.8	ng	1110680	1	7.43	239306	20.0
Phenanthrene, 2-m...	15.46	5.8	ng	91210	4	14.66	316650	20.0
Phenanthrene, 1-m...	15.50	7.2	ng	113828	4	14.66	316650	20.0
4H-Cyclopenta[def...	15.61	9.7	ng	153057	4	14.66	316650	20.0
n-Hexadecanoic acid	15.68	11.3	ng	178552	4	14.66	316650	20.0
Naphthalene, 2-ph...	15.91	4.2	ng	66527	4	14.66	316650	20.0
Cyclopenta(def)ph...	16.38	4.0	ng	63551	4	14.66	316650	20.0
Pyrene, 1-methyl-	17.37	4.0	ng	156759	5	18.37	776338	20.0
5,12-Naphthacened...	19.28	5.0	ng	61540	6	20.03	248461	20.0
4,5-Dihydrobenzo[...	19.46	5.9	ng	73032	6	20.03	248461	20.0
unknown19.81	19.81	4.8	ng	59016	6	20.03	248461	20.0
Tetracosane	20.49	13.2	ng	163532	6	20.03	248461	20.0
unknown21.39	21.39	4.4	ng	54750	6	20.03	248461	20.0
Dibenzo[def,mno]c...	21.74	5.5	ng	68693	6	20.03	248461	20.0
Dibenz(a,e)aceant...	23.27	6.5	ng	81340	6	20.03	248461	20.0
3,4:8,9-Dibenzopy...	23.44	8.5	ng	105084	6	20.03	248461	20.0