LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070217\
Data File : BF096420.D
Acq On : 2 Jul 2017 17:08
Operator : SJ/MA
Sample : I3876-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CAN-SB-0203040506-0-11

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-BF070117.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	4.928	478	483	493	rBV	500965	640899	21.92%	2.451%
2	5.351	549	555	558	rBV	2793776	2862247	97.91%	10.945%
3	6.369	723	728	731	rBV	2613909	2923325	100.00%	11.178%
4	6.504	746	751	754	rBV	2954968	2805668	95.98%	10.728%
5	6.734	786	790	793	rVB	853212	685807	23.46%	2.622%
6	6.893	813	817	820	rVB	1839302	1679198	57.44%	6.421%
7	7.293	876	885	888	rBV	1823680	1734445	59.33%	6.632%
8	7.340	888	893	897	rVV4	21365	34492	1.18%	0.132%
9	7.387	897	901	905	rVB3	59643	73661	2.52%	0.282%
10	7.545	926	928	932	rVB3	40976	36121	1.24%	0.138%
11	7.663	945	948	952	rVB4	59187	60773	2.08%	0.232%
12	7.722	952	958	960	rBV3	57134	79509	2.72%	0.304%
13	7.851	977	980	984	rBV2	27202	32567	1.11%	0.125%
14	7.910	988	990	997	rVB4	34728	57687	1.97%	0.221%
15	8.016	1003	1008	1011	rBV	1107009	972088	33.25%	3.717%
16	8.092	1019	1021	1024	rVB	75802	52233	1.79%	0.200%
17	8.234	1042	1045	1047	rBV	61514	48332	1.65%	0.185%
18	8.310	1055	1058	1061	rVB4	47566	67552	2.31%	0.258%
19	8.345	1061	1064	1066	rBV4	44775	45136	1.54%	0.173%
20	8.451	1078	1082	1084	rBV	123510	137680	4.71%	0.526%
21	8.528	1091	1095	1098	rBV4	38618	48325	1.65%	0.185%
22	8.569	1099	1102	1106	rVV4	72900	85604	2.93%	0.327%
23	8.710	1124	1126	1130	rVB3	80791	102686	3.51%	0.393%
24	8.763	1130	1135	1139	rBV4	56140	90149	3.08%	0.345%
25	8.904	1157	1159	1162	rBV3	39035	41561	1.42%	0.159%
26	8.934	1162	1164	1168	rVB4	38203	45100	1.54%	0.172%
27	8.981	1168	1172	1173	rBV4	39039	56674	1.94%	0.217%
28	9.016	1175	1178	1180	rVV2	67128	75044	2.57%	0.287%
29	9.051	1181	1184	1186	rVV	171183	149698	5.12%	0.572%
30	9.092	1186	1191	1194	rVB	1995117	1752619	59.95%	6.702%
31	9.181	1202	1206	1211	rBV5	128435	194168	6.64%	0.742%
32	9.481	1254	1257	1260	rBV	729728	653104	22.34%	2.497%
33	9.504	1260	1261	1264	rVB	199219	131768	4.51%	0.504%
34	9.539	1264	1267	1271	rVB3	77709	97939	3.35%	0.374%

LSC Area Percent Report

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Instrument : BNA_F
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.592	1274	1276	1278	rVB2	53662	37670	1.29%	0.144%
36	9.639	1282	1284	1289	rBV4	69429	111102	3.80%	0.425%
37	9.686	1289	1292	1295	rVB	219525	185416	6.34%	0.709%
38	9.716	1295	1297	1301	rBV3	53032	73260	2.51%	0.280%
39	9.769	1301	1306	1309	rBV	895571	925259	31.65%	3.538%
40	9.816	1312	1314	1316	rBV2	40288	29635	1.01%	0.113%
41	9.886	1324	1326	1329	rVB3	46808	40279	1.38%	0.154%
42	10.039	1349	1352	1355	rBV2	71240	70299	2.40%	0.269%
43	10.145	1367	1370	1372	rBV3	35201	33407	1.14%	0.128%
44	10.181	1374	1376	1379	rVV	53446	43344	1.48%	0.166%
45	10.210	1379	1381	1383	rVB	55359	37173	1.27%	0.142%
46	10.310	1396	1398	1400	rBV2	41294	38686	1.32%	0.148%
47	10.433	1415	1419	1423	rVB	65651	81304	2.78%	0.311%
48	10.551	1435	1439	1442	rBV	1358817	1218145	41.67%	4.658%
49	10.681	1458	1461	1463	rBV	86359	89409	3.06%	0.342%
50	11.128	1535	1537	1541	rBV2	55497	46040	1.57%	0.176%
51	11.245	1553	1557	1559	rBV	800567	669140	22.89%	2.559%
52	11.269	1559	1561	1564	rVB	243911	187835	6.43%	0.718%
53	11.316	1564	1569	1573	rVB	71077	67153	2.30%	0.257%
54	11.775	1644	1647	1652	rVV2	60072	71256	2.44%	0.272%
55	11.857	1657	1661	1663	rBV	50556	55992	1.92%	0.214%
56	12.451	1758	1762	1766	rBV	323078	263059	9.00%	1.006%
57	12.675	1795	1800	1804	rBV	269482	299641	10.25%	1.146%
58	12.827	1822	1826	1829	rVB	1270934	1111925	38.04%	4.252%
59	12.898	1834	1838	1841	rBV3	36660	39731	1.36%	0.152%
60	13.010	1854	1857	1859	rVB2	32238	33370	1.14%	0.128%
61	13.110	1871	1874	1877	rBV	37549	38992	1.33%	0.149%
62	13.727	1974	1979	1986	rBV2	140599	179332	6.13%	0.686%
63	13.869	1998	2003	2006	rBV2	636406	726061	24.84%	2.776%
64	13.898	2006	2008	2014	rVB	126777	113375	3.88%	0.434%
65	14.880	2171	2175	2178	rBV	88058	126341	4.32%	0.483%
66	15.163	2220	2223	2228	rBV	54129	64311	2.20%	0.246%
67	15.221	2230	2233	2239	rBV	63247	71311	2.44%	0.273%
68	15.280	2239	2243	2248	rBV	337898	418987	14.33%	1.602%

Sum of corrected areas: 26152099

LSC Report - Integrated Chromatogram

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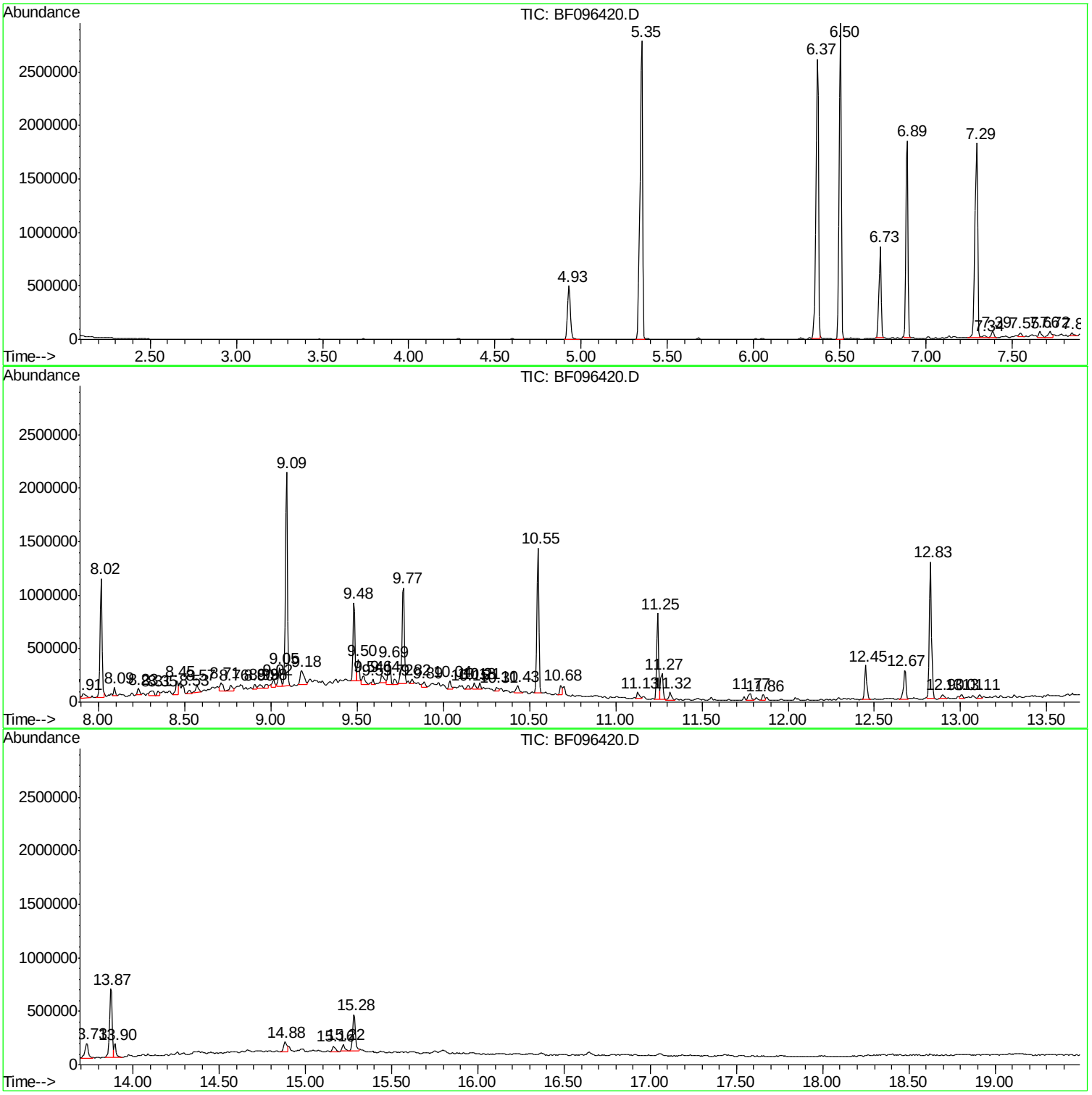
Instrument :
BNA_F
ClientSampleId :
CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070217\
Data File : BF096420.D
Acq On : 2 Jul 2017 17:08
Operator : SJ/MA
Sample : I3876-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

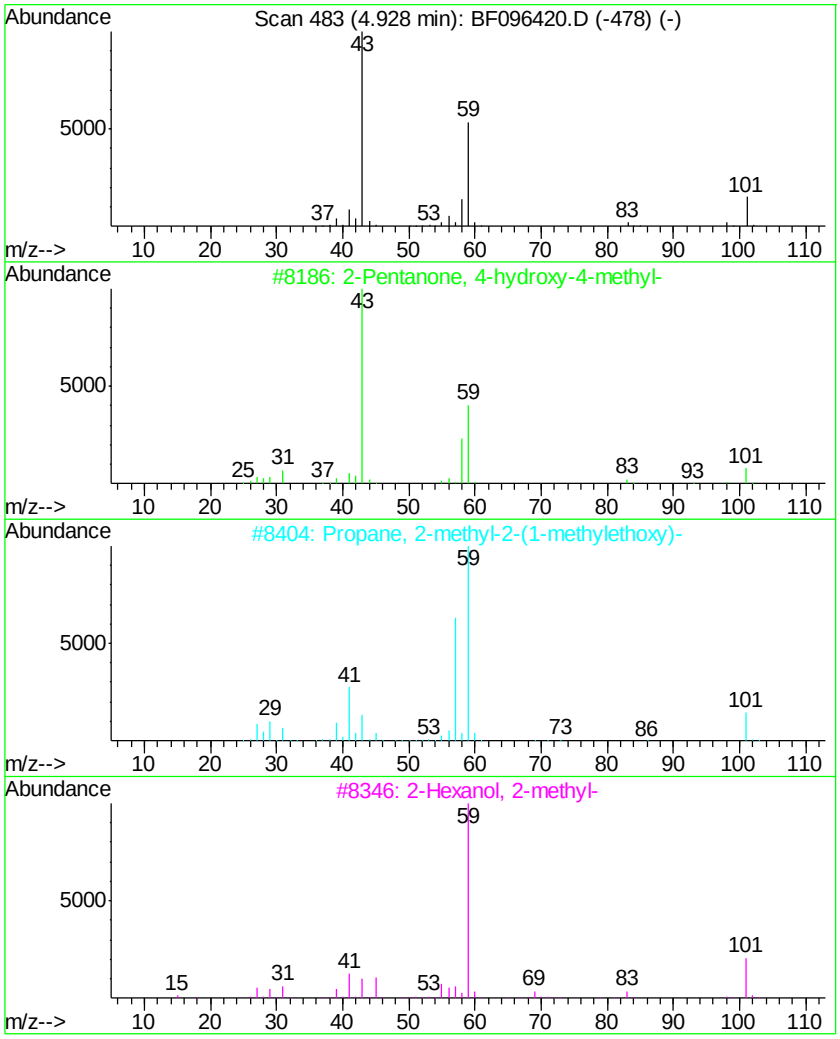
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

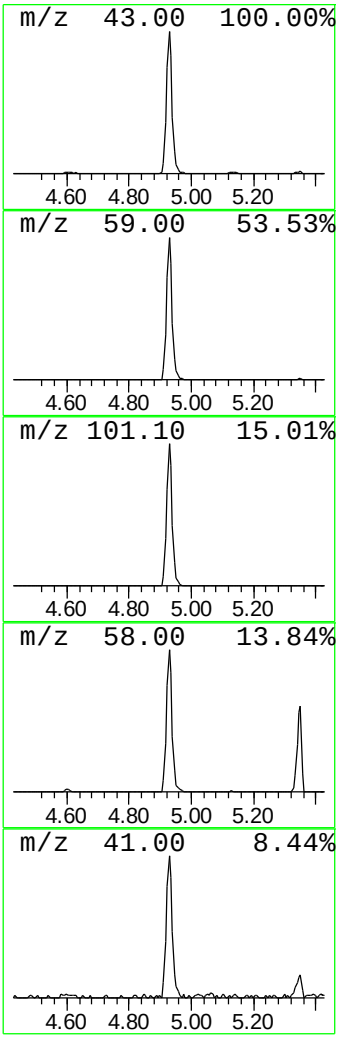
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	18.69 ng	640899	1,4-Dichlorobenzene-d4	6.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23





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Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
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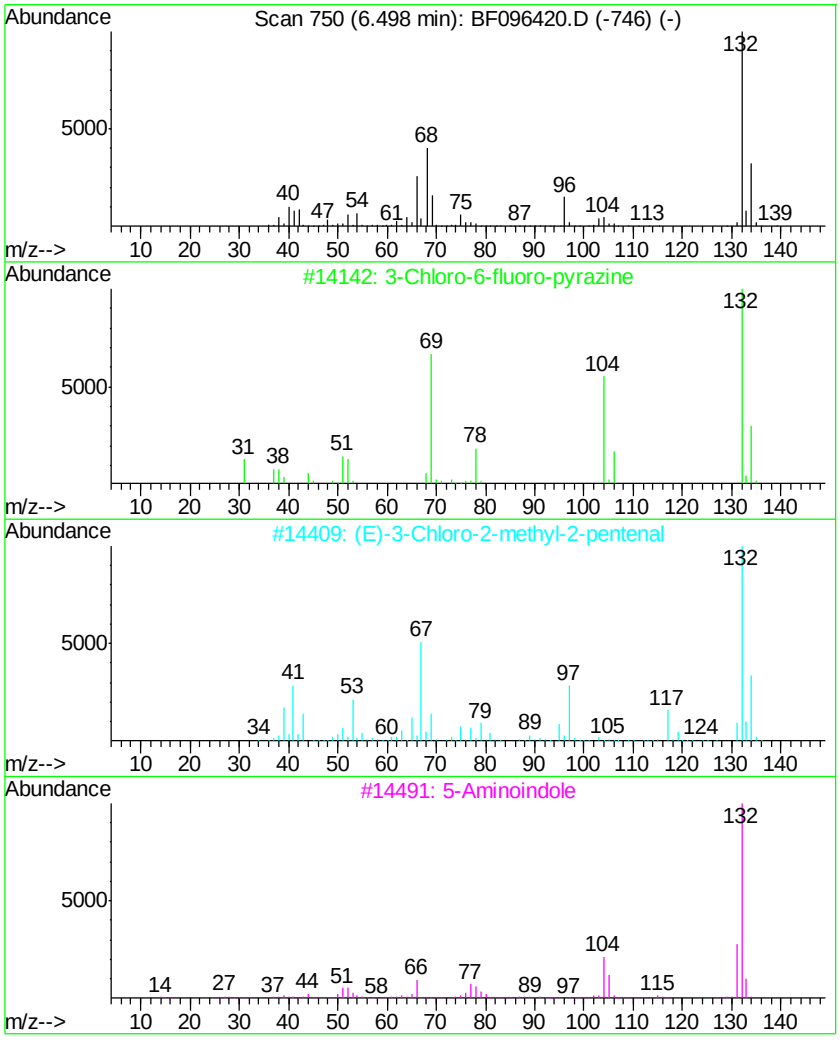
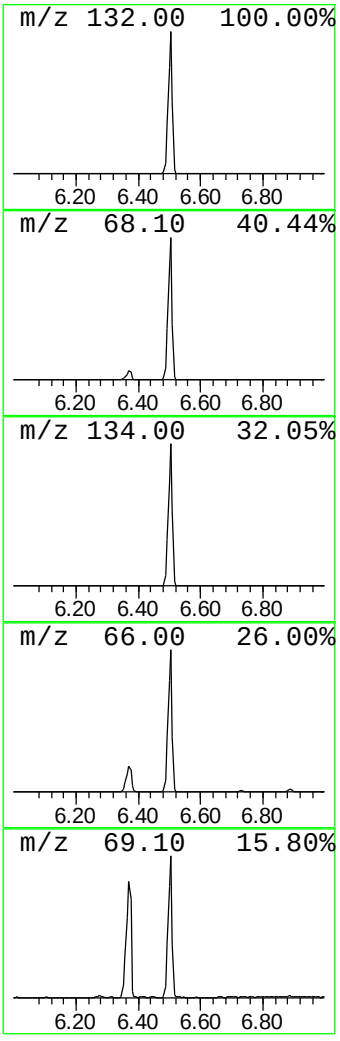
****

Peak Number 2 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	81.82 ng	2805670	1,4-Dichlorobenzene-d4	6.73

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	16
3	5-Aminoindole	132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9





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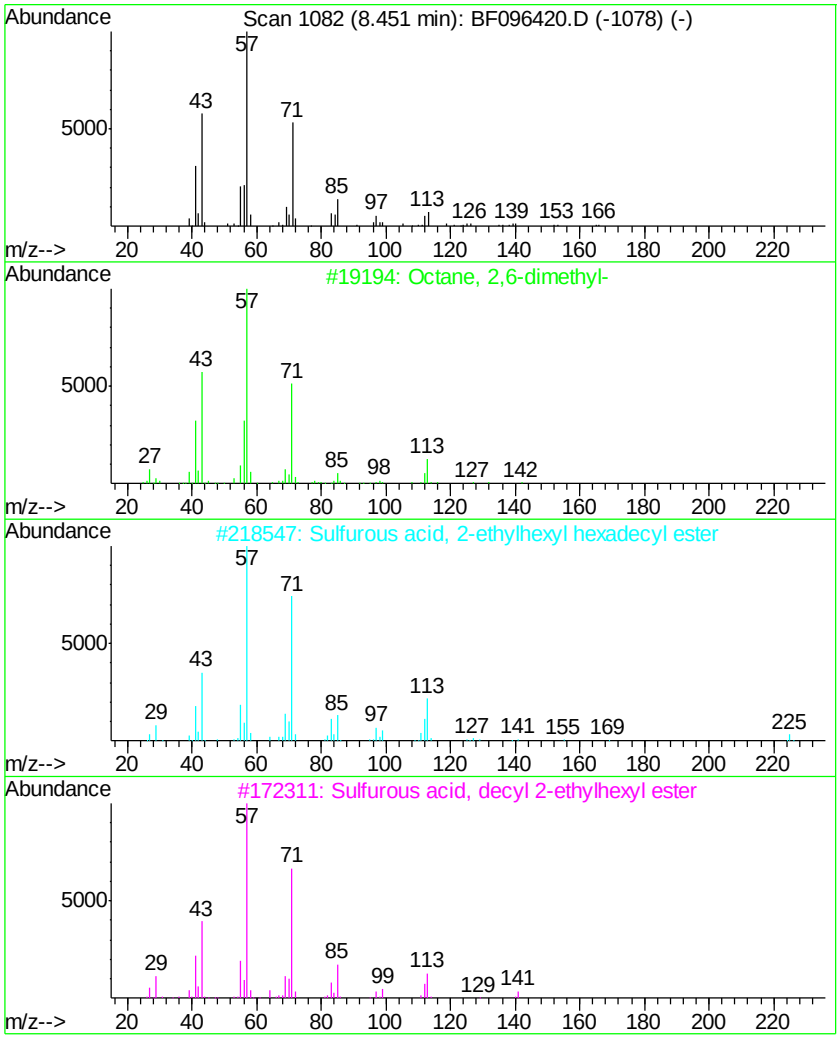
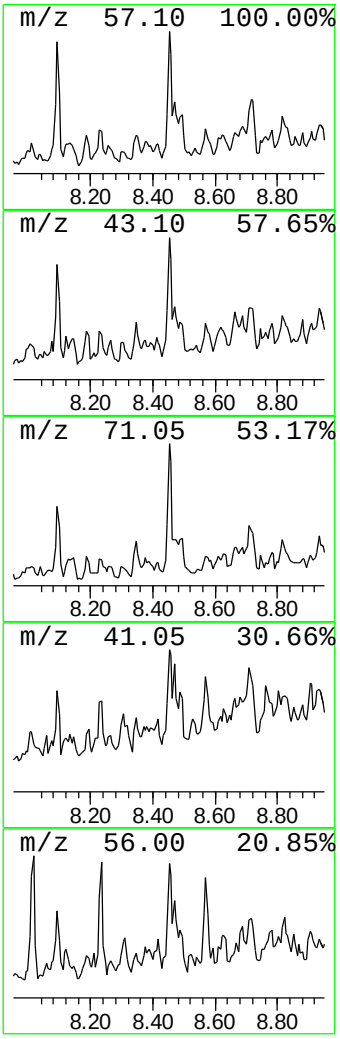
****

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.45	2.83 ng	137680	Naphthalene-d8	8.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
2	Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	64
3	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	64
4	Sulfurous acid, 2-ethylhexyl oct...	446	C26H54O3S	1000309-20-1	64
5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	64





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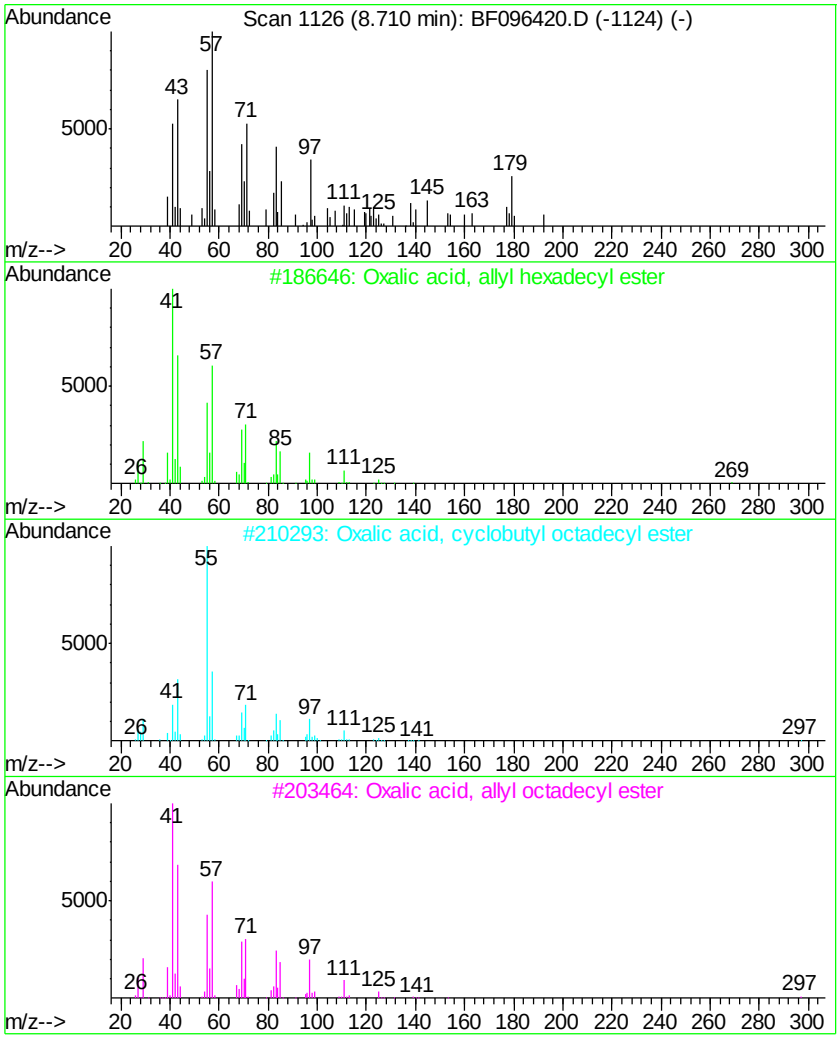
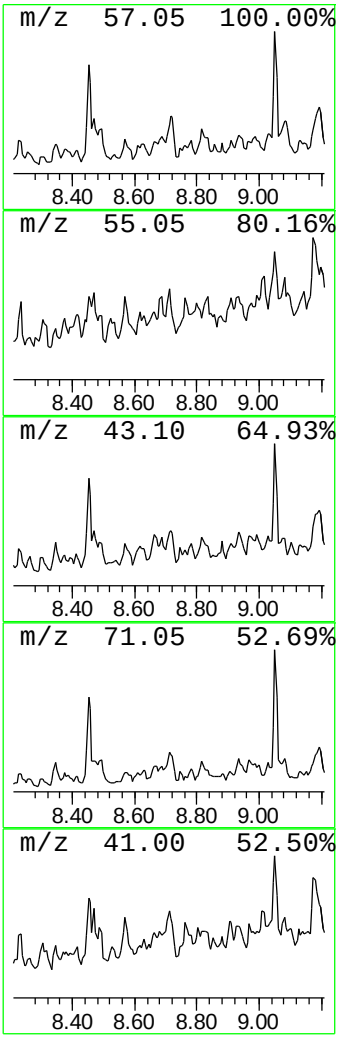
Peak Number 5 Oxalic acid, allyl hexadecy... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.71	2.11 ng	102686	Naphthalene-d8	8.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	52
2	Oxalic acid, cyclobutyl octadecy...	396	C24H44O4	1000309-70-8	52
3	Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	50
4	1-Hexacosanol	382	C26H54O	000506-52-5	46
5	Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	35





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TIC Library : C:\DATABASE\NIST11.L
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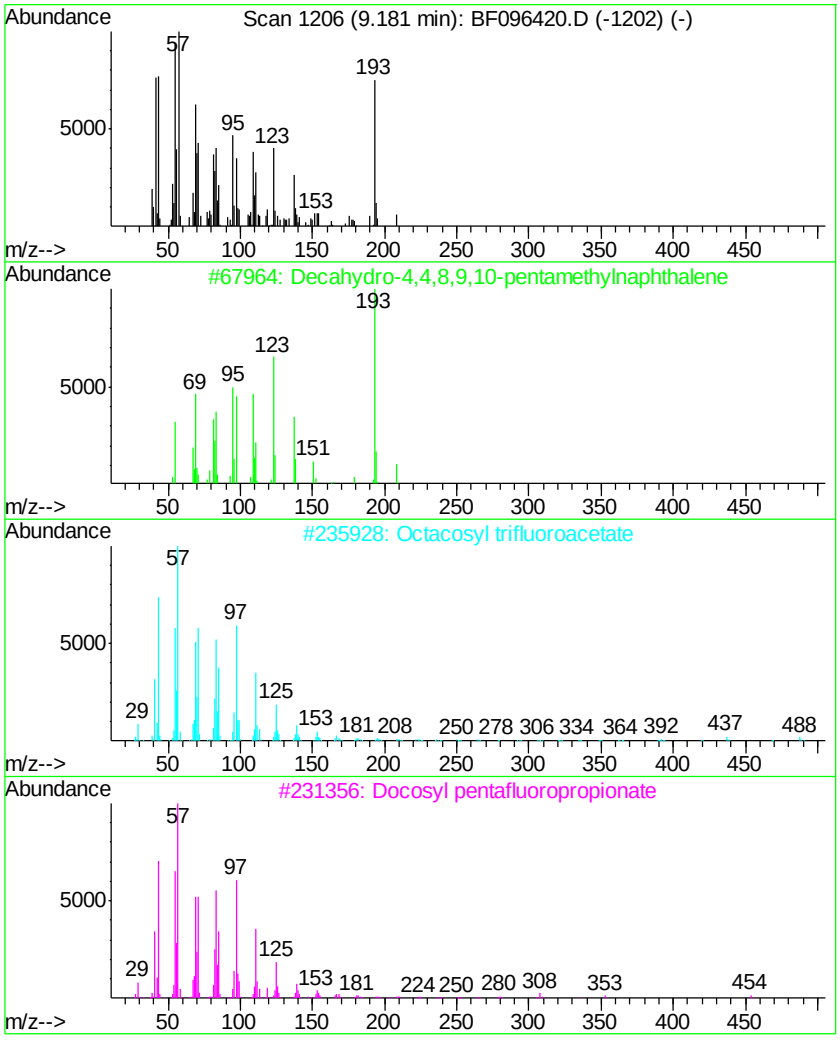
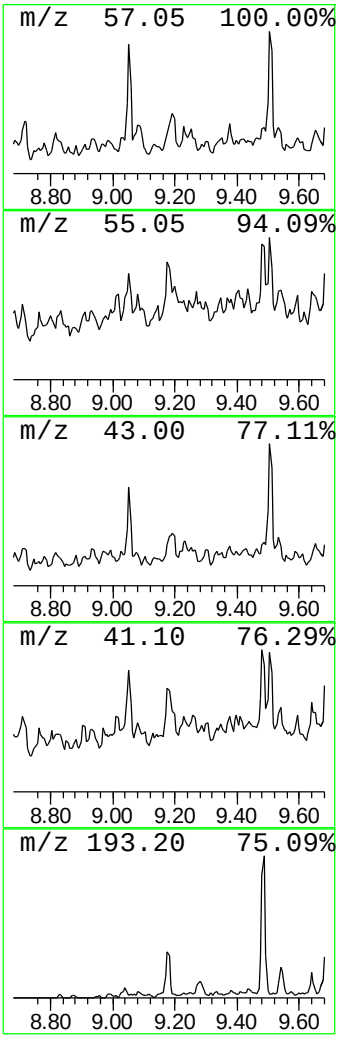
Peak Number 6 Decahydro-4,4,8,9,10-pentam... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.18	4.20 ng	194168	Acenaphthene-d10	9.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	86
2	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	20
3	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	20
4	Triacontyl acetate	480	C32H64O2	041755-58-2	18
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	18





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

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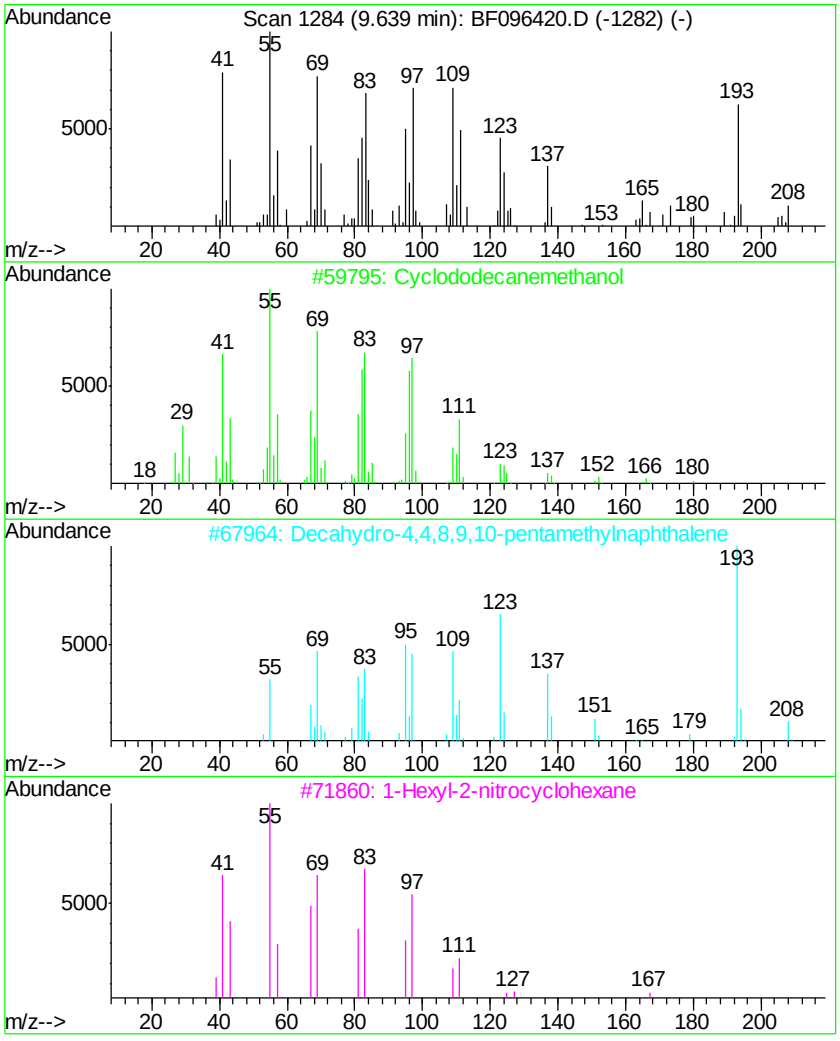
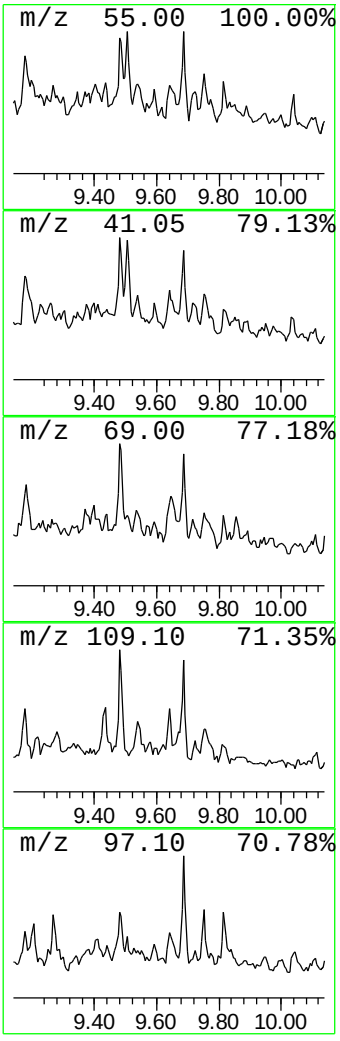
Peak Number 8 unknown9.64 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.64	2.40 ng	111102	Acenaphthene-d10	9.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclododecanemethanol	198	C13H26O	001892-12-2	47
2	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	43
3	1-Hexyl-2-nitrocyclohexane	213	C12H23NO2	118252-04-3	35
4	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	35
5	Octane, 1,8-dibromo-	270	C8H16Br2	004549-32-0	27





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Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

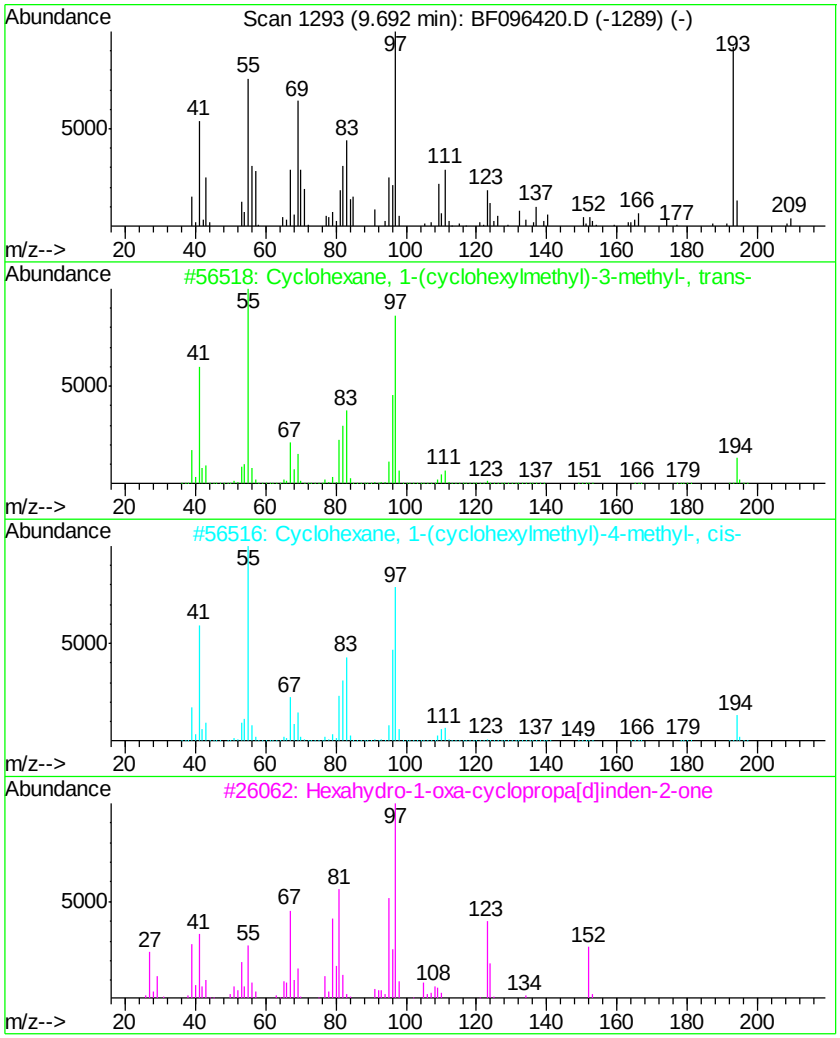
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

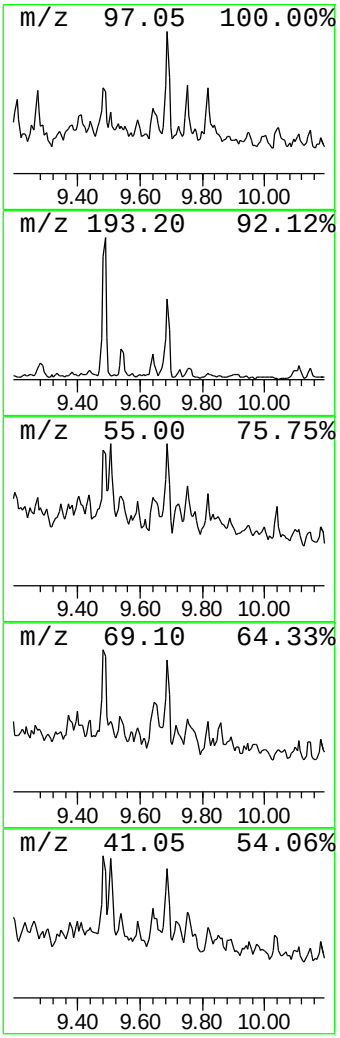
****

Peak Number 9 unknown9.69 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	4.01 ng	185416	Acenaphthene-d10	9.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-95-9	38
2	Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-97-1	30
3	Hexahydro-1-oxa-cyclopropa[d]ind...	152	C9H12O2	037643-23-5	27
4	Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-96-0	25
5	Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-98-2	18





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070217\
Data File : BF096420.D
Acq On : 2 Jul 2017 17:08
Operator : SJ/MA
Sample : I3876-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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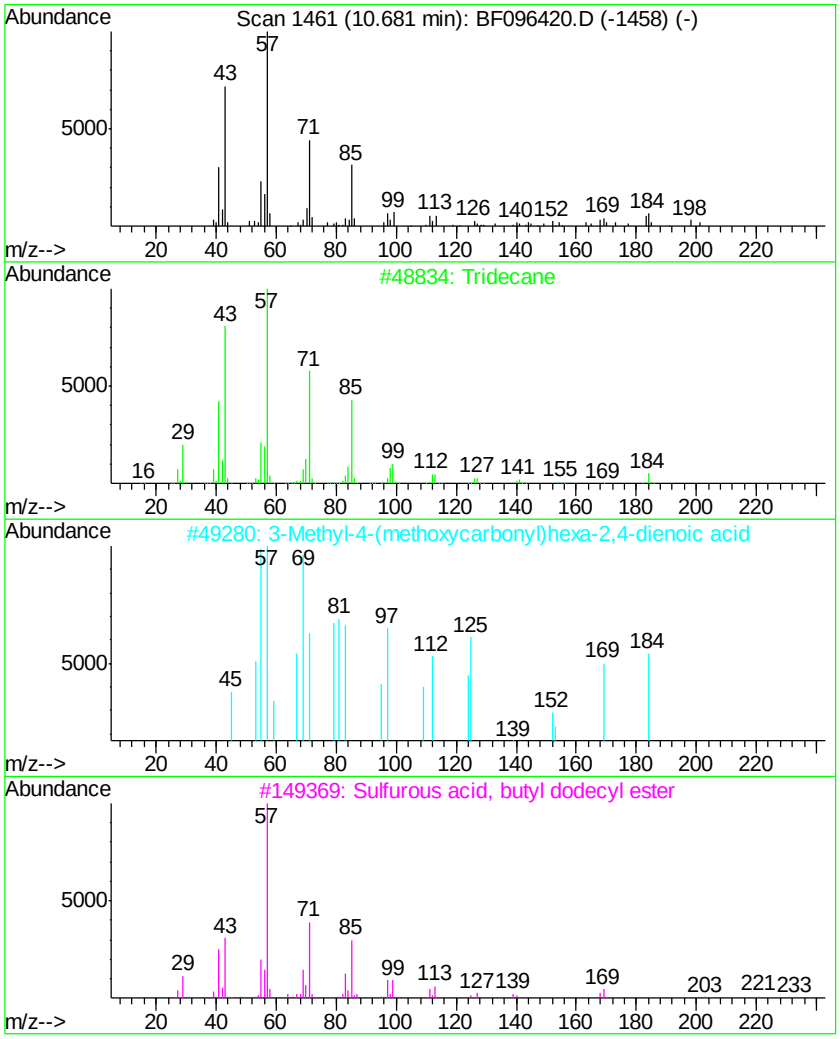
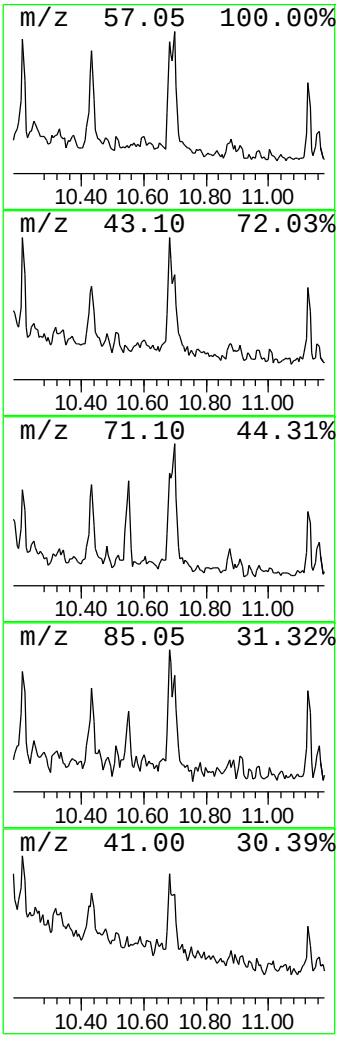
Peak Number 10 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	2.67 ng	89409	Phenanthrene-d10	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	3-Methyl-4-(methoxycarbonyl)hexa...	184	C9H12O4	1000104-10-8	90
3	Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	80
4	Dodecane	170	C12H26	000112-40-3	74
5	5-Ethyldecane	170	C12H26	017302-36-2	72





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Operator : SJ/MA
Sample : I3876-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
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TIC Library : C:\DATABASE\NIST11.L
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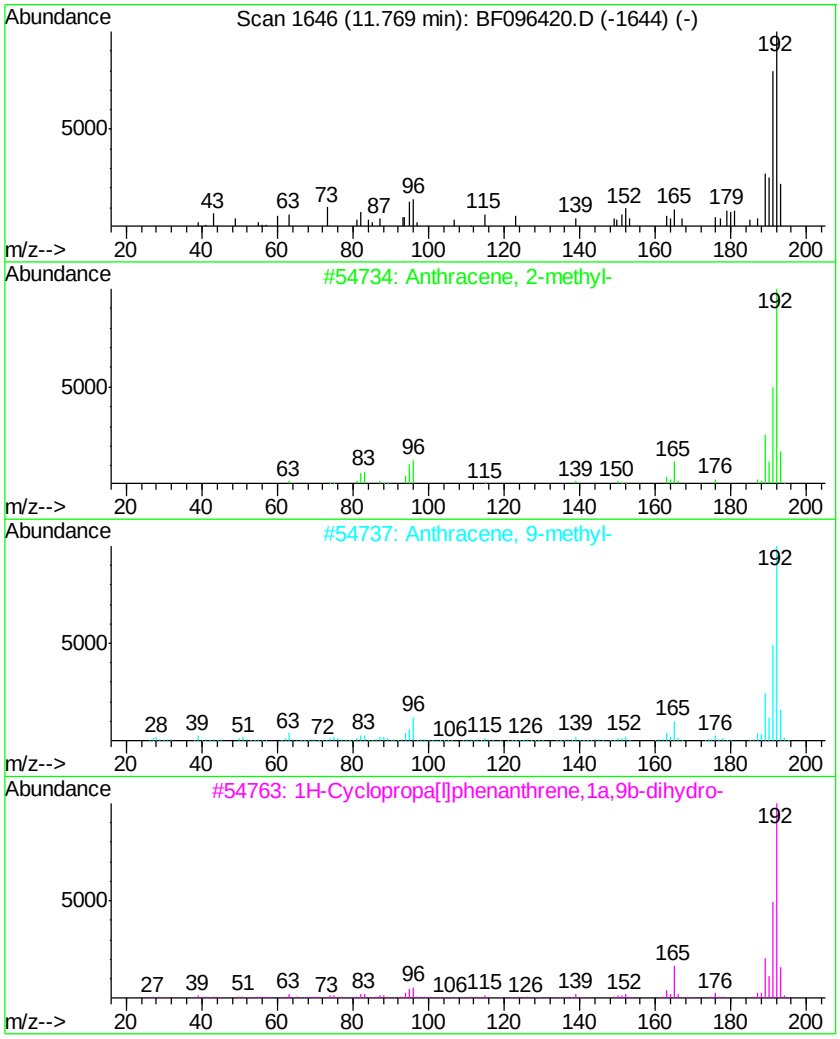
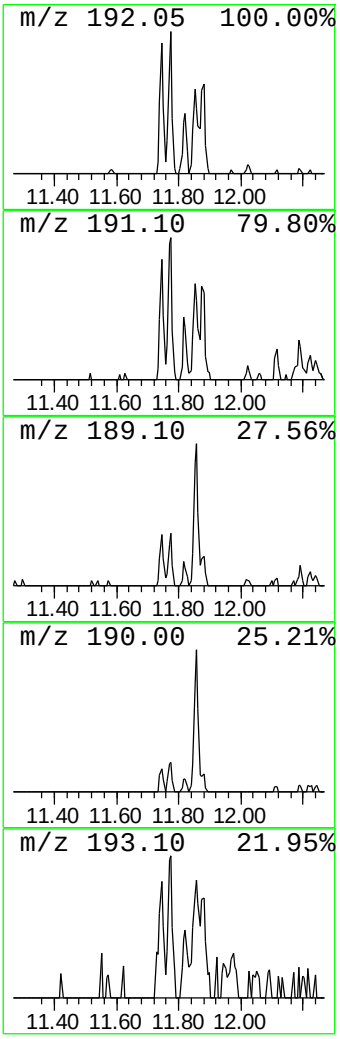
Peak Number 11 Anthracene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	2.13 ng	71256	Phenanthrene-d10	11.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	76
2	Anthracene, 9-methyl-	192	C15H12	000779-02-2	70
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	64
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	64
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	64





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Sample : I3876-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
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TIC Integration Parameters: LSCINT.P

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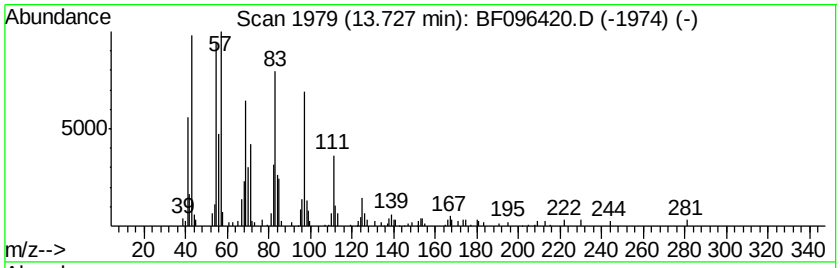
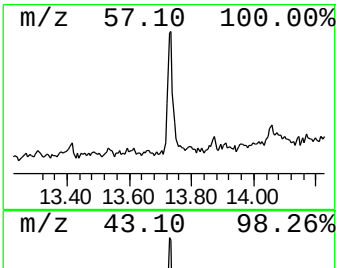
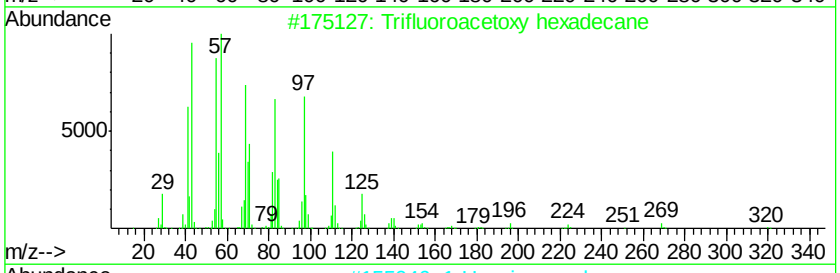
Peak Number 12 Trifluoroacetoxy hexadecane Concentration Rank 4

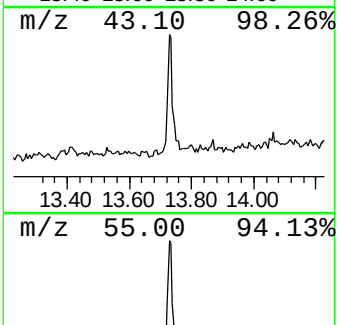
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	4.94 ng	179332	Chrysene-d12	13.87

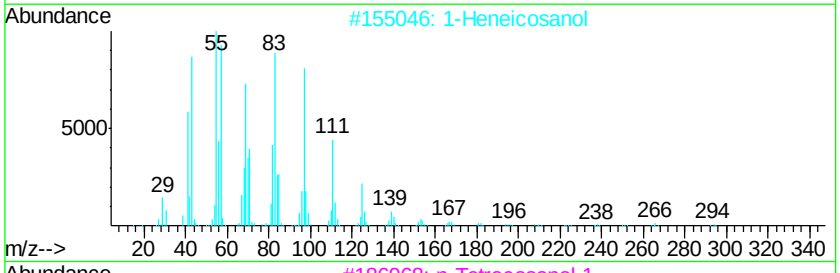
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
4	1-Docosene	308	C22H44	001599-67-3	91
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90

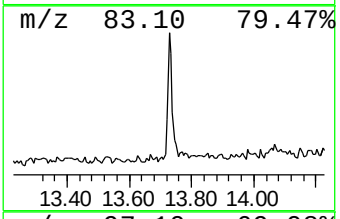


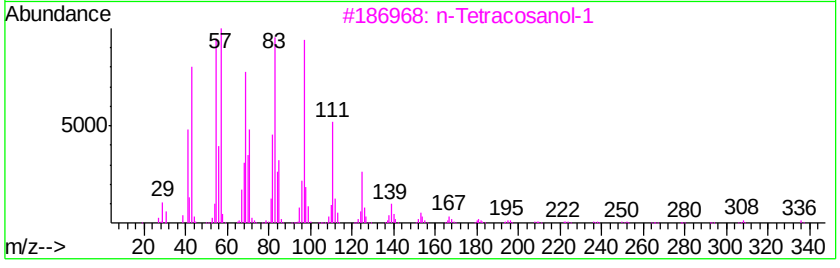












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Operator : SJ/MA
Sample : I3876-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CAN-SB-0203040506-0-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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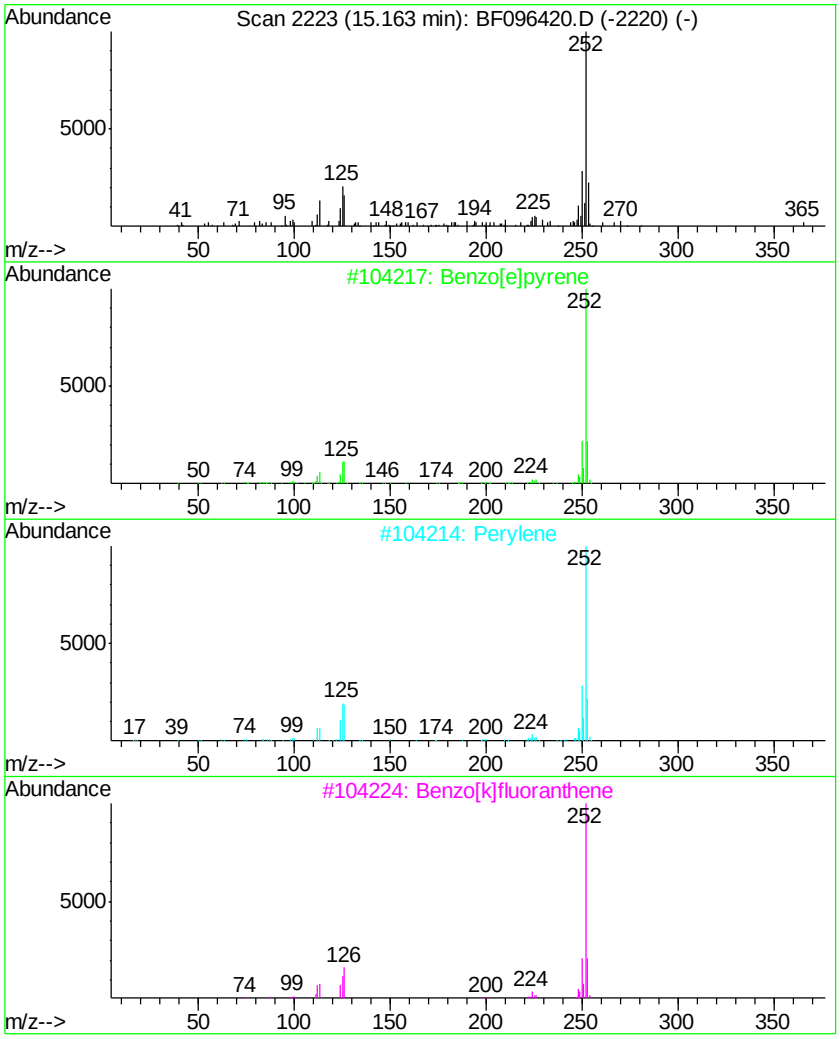
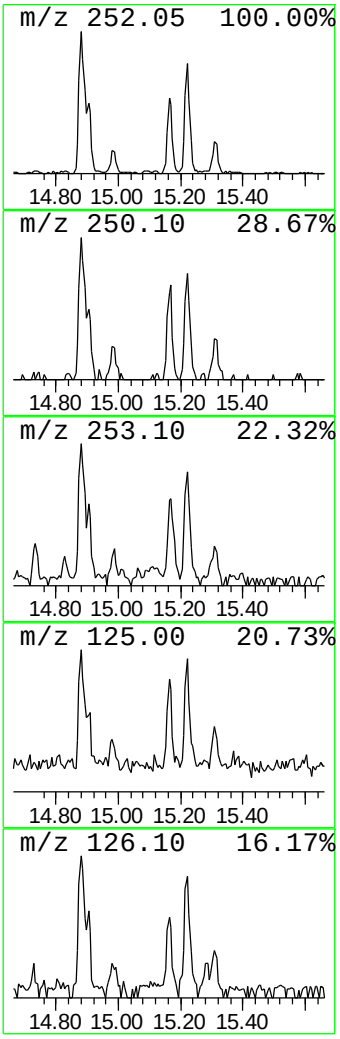
Peak Number 13 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	3.07 ng	64311	Perylene-d12	15.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	90
2	Perylene	252	C20H12	000198-55-0	90
3	Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
4	Benzo[a]pyrene	252	C20H12	000050-32-8	86
5	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	76





Tentatively Identified Compound (LSC) summary

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Operator : SJ/MA
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Misc :
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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...	4.93	18.7	ng	640899	1	6.73	685807	20.0
unknown6.50	6.50	81.8	ng	2805670	1	6.73	685807	20.0
Octane, 2,6-dimet...	8.45	2.8	ng	137680	2	8.02	972088	20.0
Oxalic acid, ally...	8.71	2.1	ng	102686	2	8.02	972088	20.0
Decahydro-4,4,8,9...	9.18	4.2	ng	194168	3	9.77	925259	20.0
unknown9.64	9.64	2.4	ng	111102	3	9.77	925259	20.0
unknown9.69	9.69	4.0	ng	185416	3	9.77	925259	20.0
Tridecane	10.68	2.7	ng	89409	4	11.25	669140	20.0
Anthracene, 2-met...	11.77	2.1	ng	71256	4	11.25	669140	20.0
Trifluoroacetoxy ...	13.73	4.9	ng	179332	5	13.87	726061	20.0
Benzo[e]pyrene	15.16	3.1	ng	64311	6	15.29	418987	20.0