LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
Data File : BF116748.D
Acq On : 1 Sep 2019 17:48
Operator : HP/JU
Sample : K4544-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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	5.451	566	572	576	rBV	1314832	1274929	33.54%	4.860%
2	6.469	739	745	752	rBV	1531645	1485340	39.08%	5.663%
3	6.610	763	769	773	rBV	1564530	1446670	38.06%	5.515%
4	6.845	805	809	812	rBV	532045	437796	11.52%	1.669%
5	6.998	831	835	839	rBV	1198690	1101594	28.98%	4.200%
6	7.398	896	903	906	rBV	991223	890260	23.42%	3.394%
7	8.128	1018	1027	1029	rBV	700937	658481	17.32%	2.510%
8	9.175	1199	1205	1206	rBV2	167450	217278	5.72%	0.828%
9	9.198	1206	1209	1213	rVB	2026303	1727540	45.45%	6.586%
10	9.286	1221	1224	1229	rBV5	117201	173902	4.58%	0.663%
11	9.592	1274	1276	1278	rBV	548532	470820	12.39%	1.795%
12	9.757	1301	1304	1307	rBV2	375268	484056	12.74%	1.845%
13	9.798	1308	1311	1314	rVB	792262	669476	17.61%	2.552%
14	9.839	1315	1318	1320	rBV	268279	277201	7.29%	1.057%
15	9.880	1320	1325	1330	rBV2	1076007	1341599	35.30%	5.115%
16	9.933	1331	1334	1336	rBV2	302002	421758	11.10%	1.608%
17	10.045	1351	1353	1355	rBV2	258569	252064	6.63%	0.961%
18	10.151	1368	1371	1373	rBV2	642351	600183	15.79%	2.288%
19	10.227	1382	1384	1387	rBV	500597	396772	10.44%	1.513%
20	10.263	1387	1390	1392	rBV3	595217	717128	18.87%	2.734%
21	10.304	1394	1397	1399	rVV	1487037	1405784	36.99%	5.359%
22	10.833	1483	1487	1490	rVB	3333175	3534307	92.99%	13.474%
23	11.304	1564	1567	1570	rVB	3226296	3800776	100.00%	14.490%
24	13.986	2021	2023	2025	rBV	602544	496904	13.07%	1.894%
25	15.033	2198	2201	2208	rVB	427766	788380	20.74%	3.006%
26	15.333	2249	2252	2258	rVB	249371	313899	8.26%	1.197%
27	15.392	2259	2262	2266	rVB	246902	266020	7.00%	1.014%
28	15.457	2269	2273	2276	rBV	453358	579721	15.25%	2.210%

Sum of corrected areas: 26230638

LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
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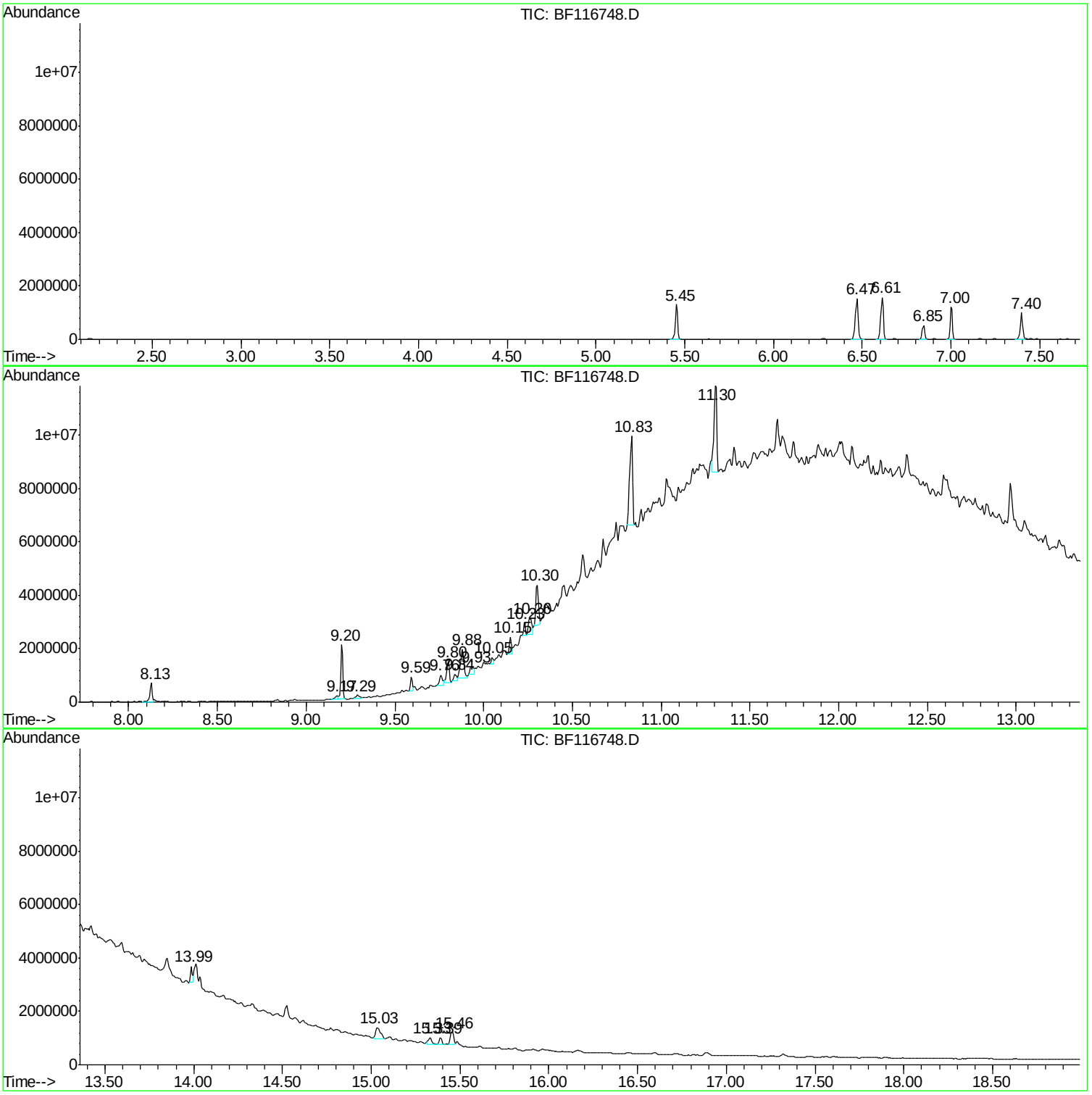
Sample : K4544-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
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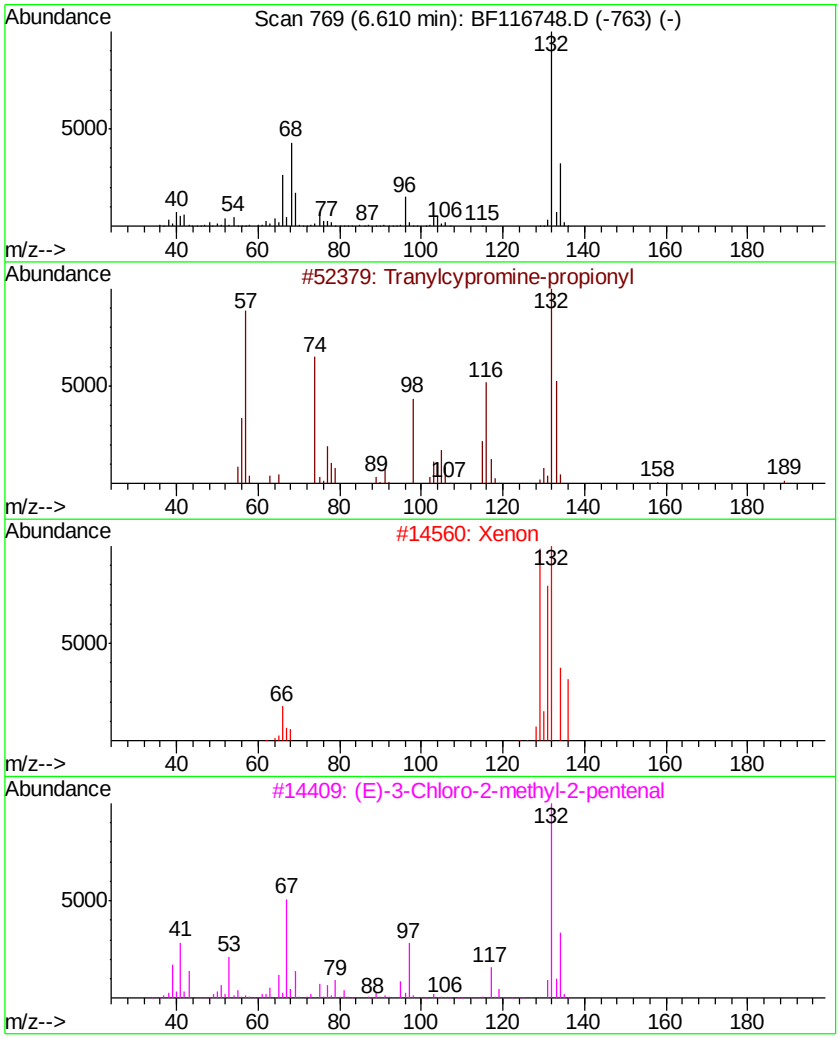
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

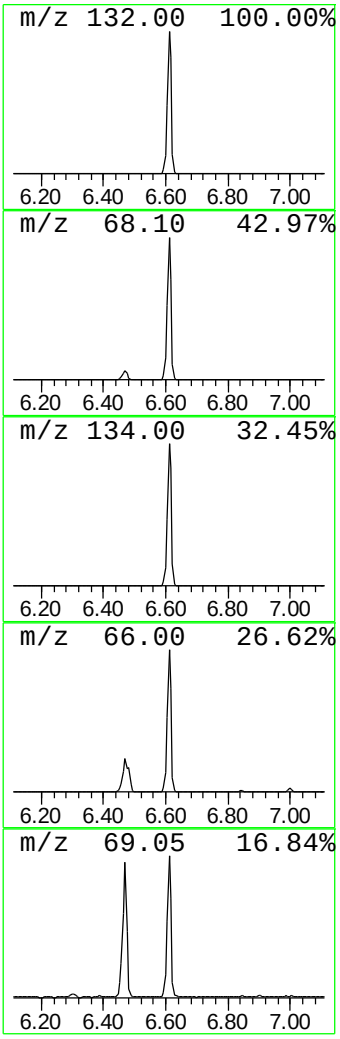
****

Peak Number 1 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	66.09 ng	1446670	1,4-Dichlorobenzene-d4	6.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
2	Xenon	132	Xe	007440-63-3	9
3	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
Data File : BF116748.D
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Operator : HP/JU
Sample : K4544-19
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Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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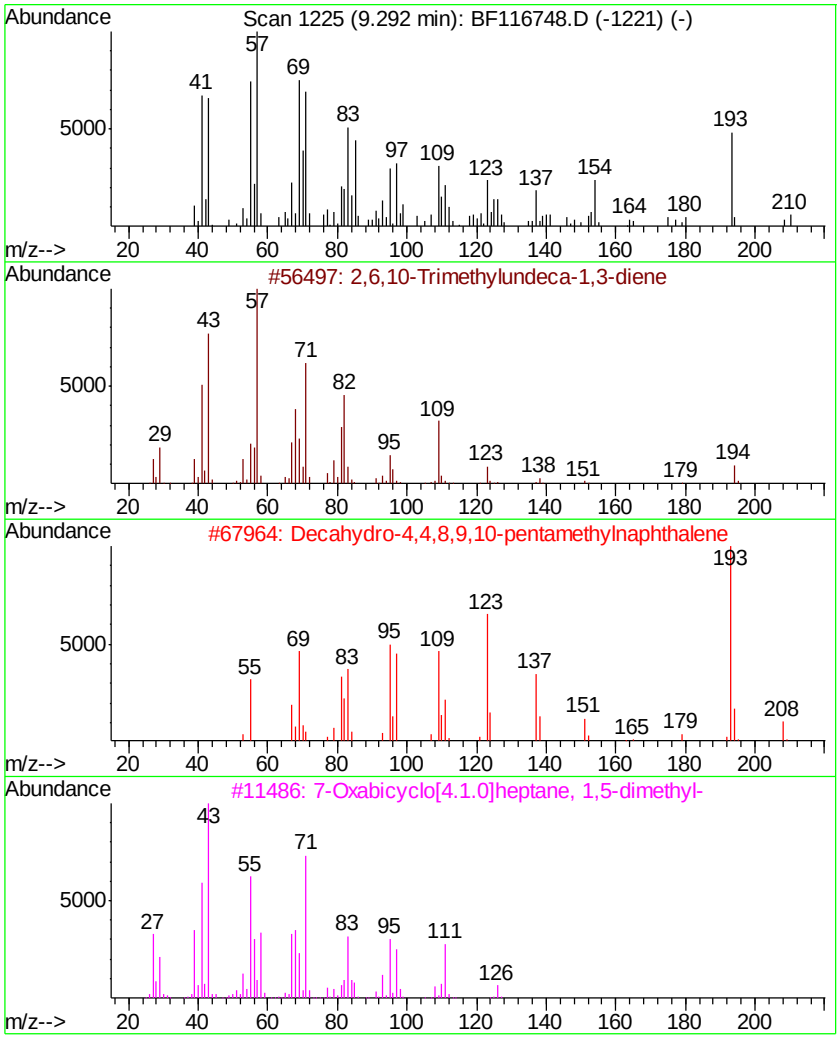
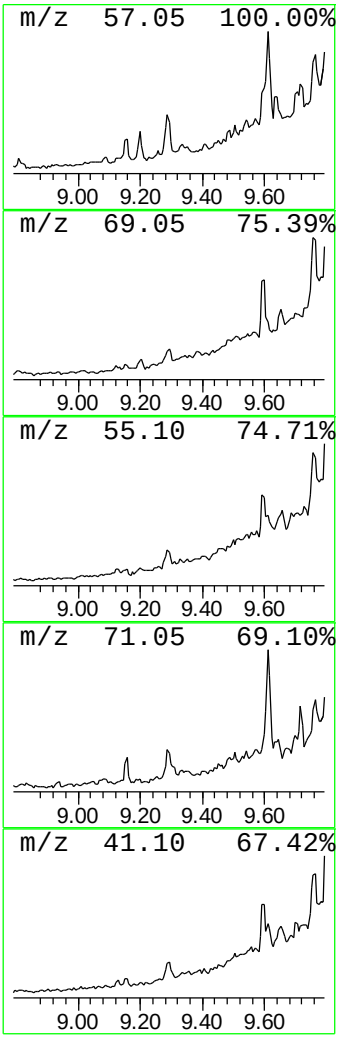
Peak Number 3 unknown9.29 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.29	2.59 ng	173902	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Trimethylundeca-1,3-diene	194	C14H26	1000222-09-1	48
2	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	46
3	7-Oxabicyclo[4.1.0]heptane, 1,5-...	126	C8H14O	162239-52-3	38
4	Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	25
5	2-Butanone, 4-(2,6,6-trimethyl-2...	194	C13H22O	039721-65-8	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
Data File : BF116748.D
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Operator : HP/JU
Sample : K4544-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
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TIC Library : C:\DATABASE\NIST11.L
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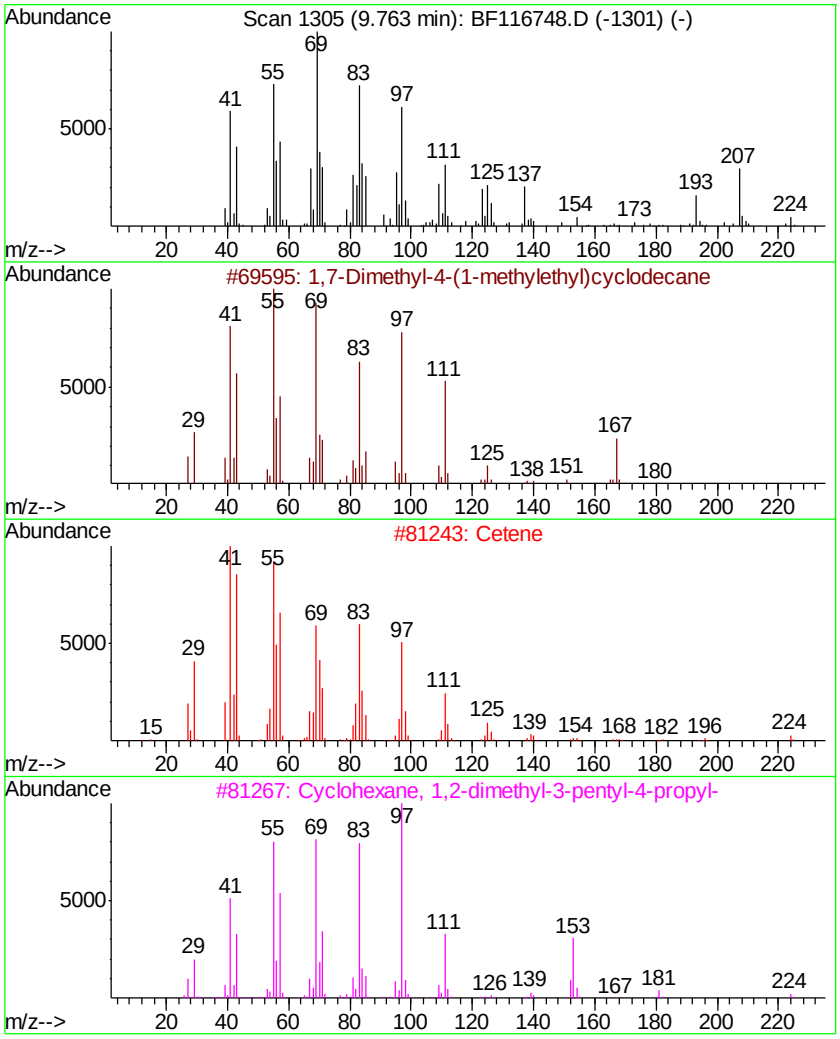
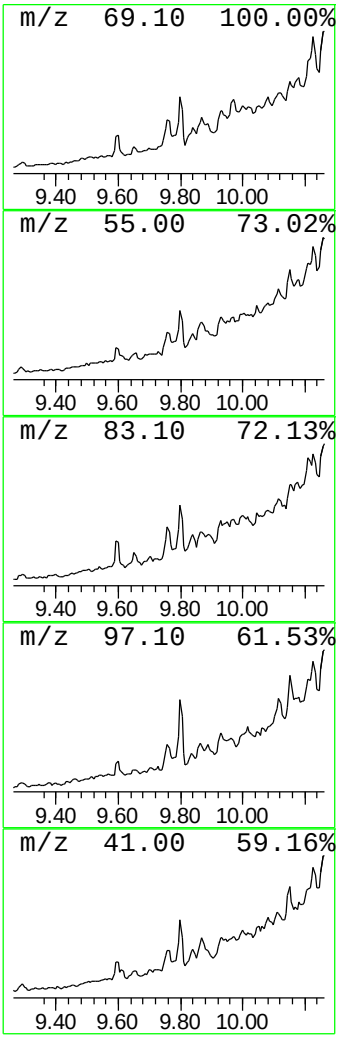
Peak Number 4 unknown9.76 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	7.22 ng	484056	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,7-Dimethyl-4-(1-methylethyl)cy...	210	C15H30	000645-10-3	43
2	Cetene	224	C16H32	000629-73-2	38
3	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	38
4	Crotyl methacrylate	140	C8H12O2	007376-45-6	30
5	Cyclohexane, 1,1'-(2-propyl-1,3-...	250	C18H34	055030-21-2	27





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
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Misc :
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Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
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TIC Library : C:\DATABASE\NIST11.L
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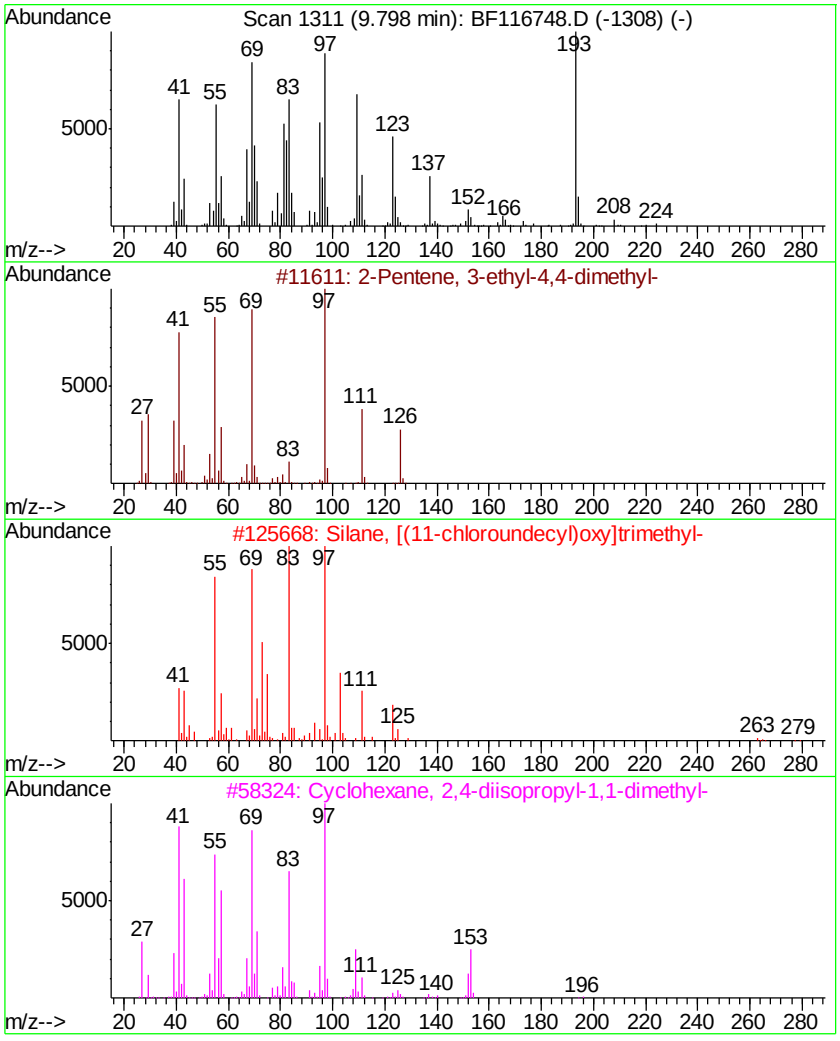
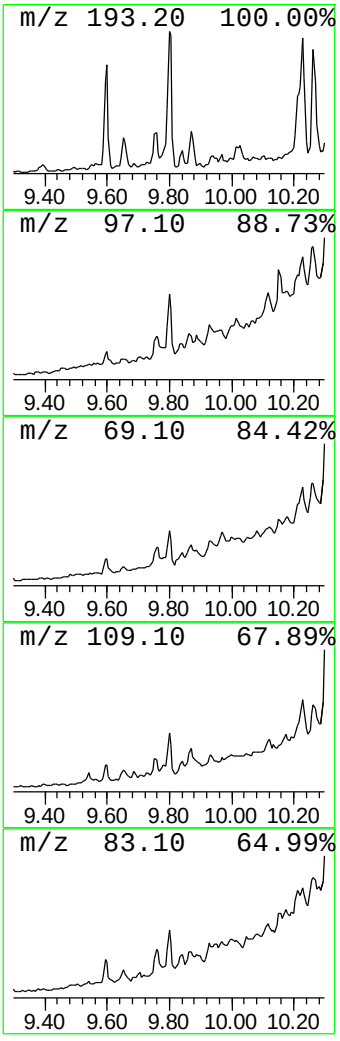
****

Peak Number 5 unknown9.80 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.80	9.98 ng	669476	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	27
2	Silane, [(11-chloroundecyl)oxy]t...	278	C14H31ClOSi	026305-82-8	27
3	Cyclohexane, 2,4-diisopropyl-1,1...	196	C14H28	1000149-60-5	27
4	1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	25
5	4-Hepten-3-one, 4-methyl-	126	C8H14O	022319-31-9	25





Library Search Compound Report

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Sample : K4544-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
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TIC Library : C:\DATABASE\NIST11.L
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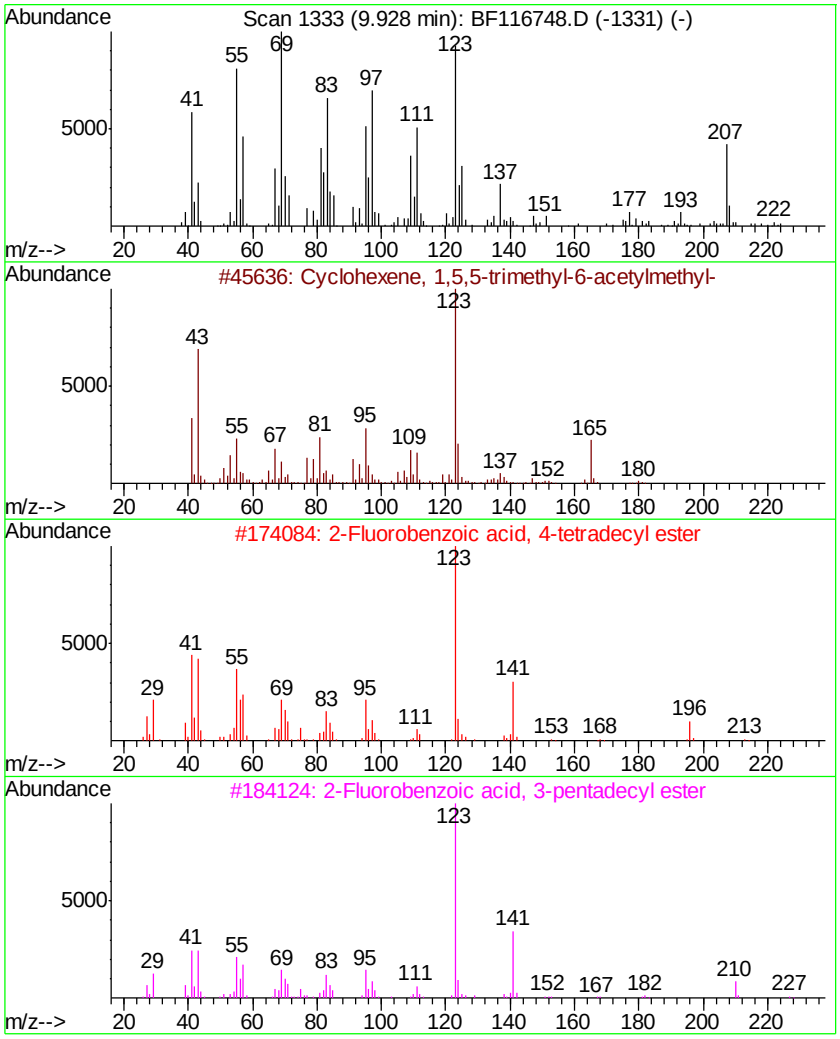
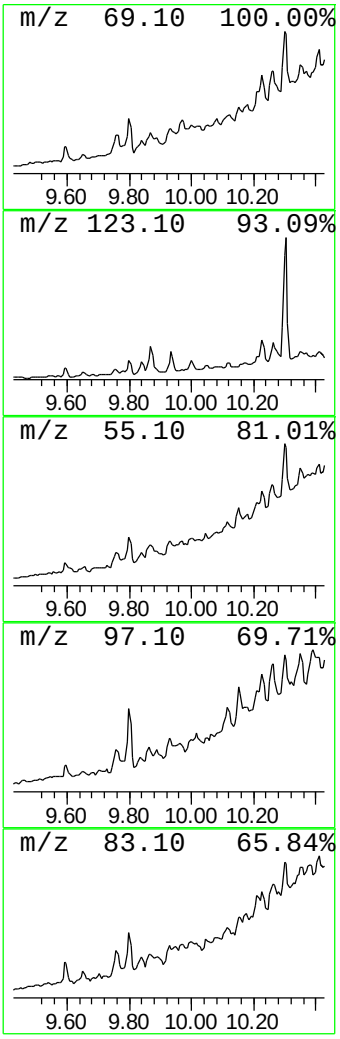
****

Peak Number 6 Cyclohexene, 1,5,5-trimethy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	6.29 ng	421758	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1,5,5-trimethyl-6-a...	180	C12H20O	211563-96-1	64
2	2-Fluorobenzoic acid, 4-tetradec...	336	C21H33FO2	1000280-61-4	50
3	2-Fluorobenzoic acid, 3-pentadec...	350	C22H35FO2	1000280-61-6	50
4	3,3,5,5-Tetramethylcyclohexanol	156	C10H20O	002650-40-0	47
5	4-Fluorobenzoic acid, 6-ethyl-3-...	280	C17H25FO2	1000279-07-4	47





Library Search Compound Report

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Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.M
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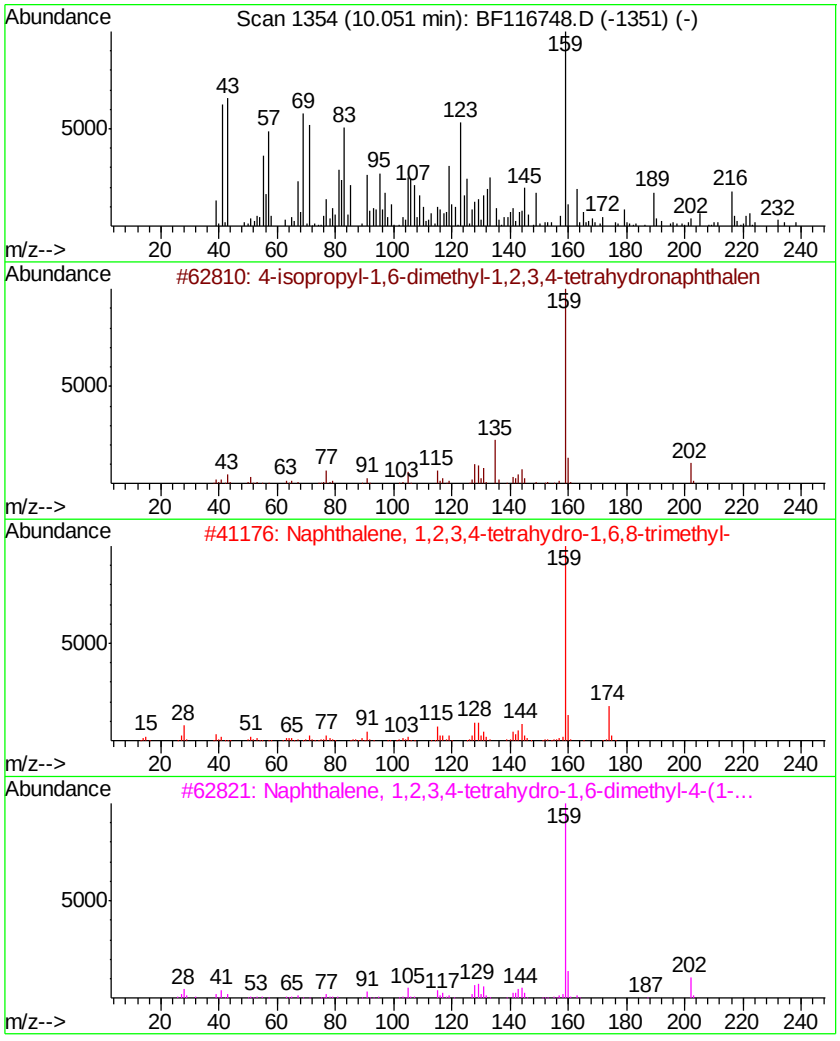
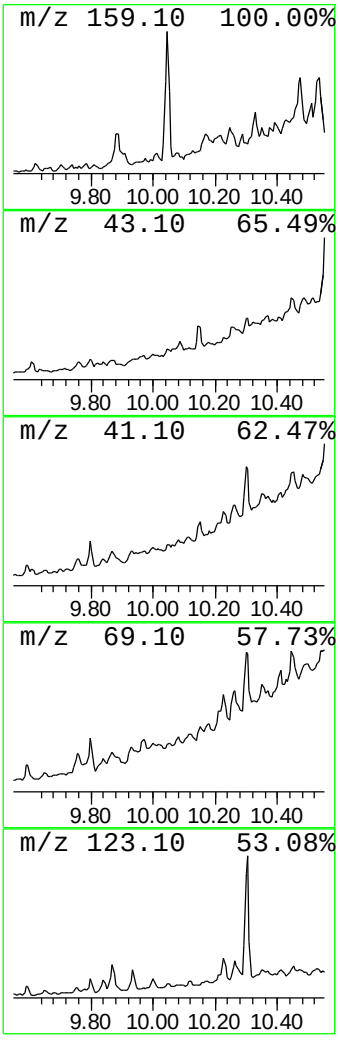
****

Peak Number 7 4-isopropyl-1,6-dimethyl-1,... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.05	3.76 ng	252064	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-isopropyl-1,6-dimethyl-1,2,3,4...	202	C15H22	1000378-99-6	55
2	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	46
3	Naphthalene, 1,2,3,4-tetrahydro-...	202	C15H22	000483-77-2	46
4	Quinoline, 6-methoxy-	159	C10H9NO	005263-87-6	46
5	Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	46





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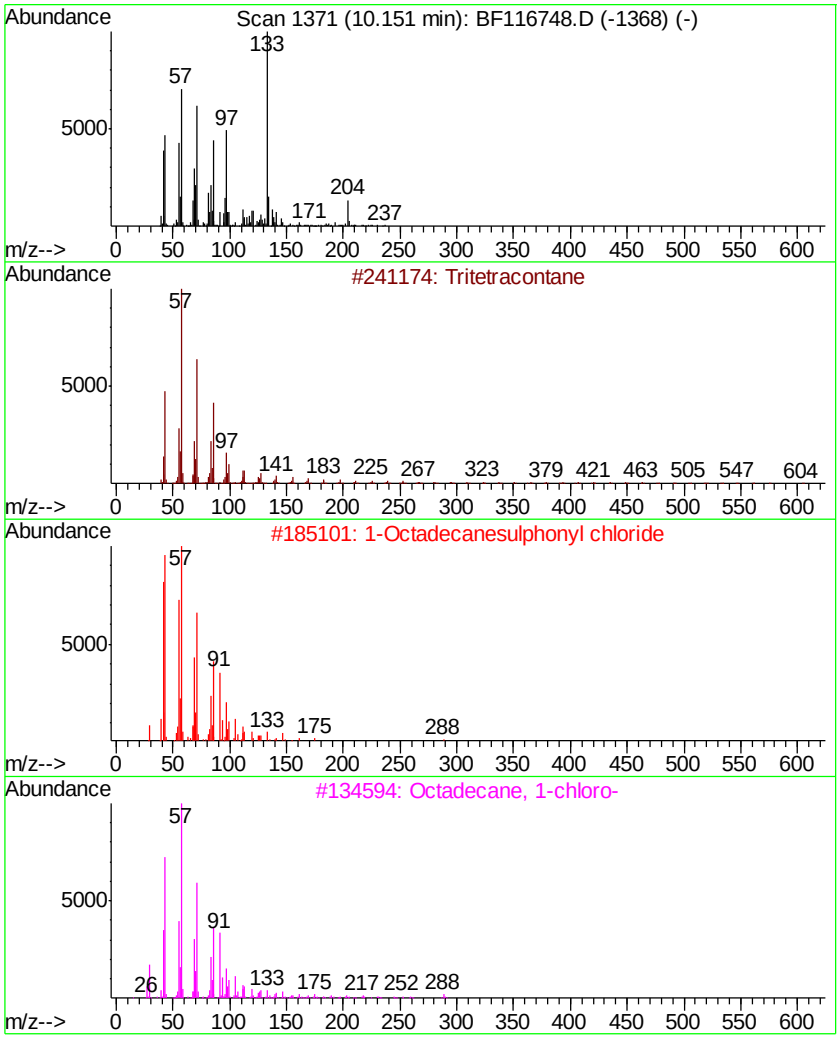
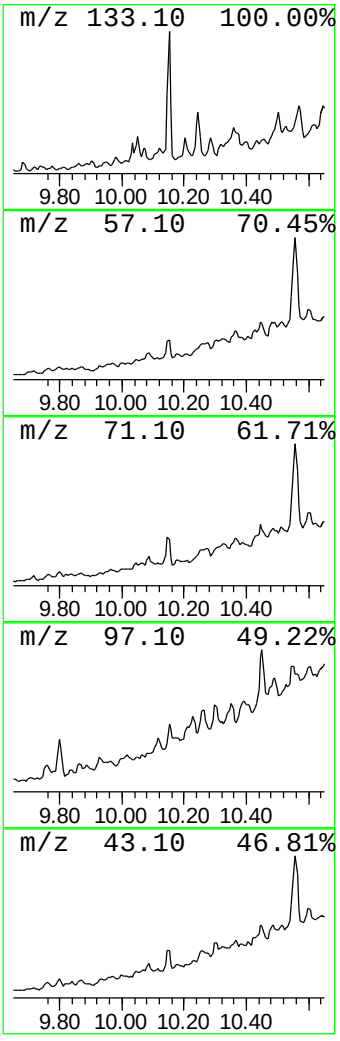
Peak Number 8 unknown10.15 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.15	8.95 ng	600183	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tritetracontane	605	C43H88	007098-21-7	27
2	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	25
3	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	22
4	Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	22
5	Nonadecane	268	C19H40	000629-92-5	22





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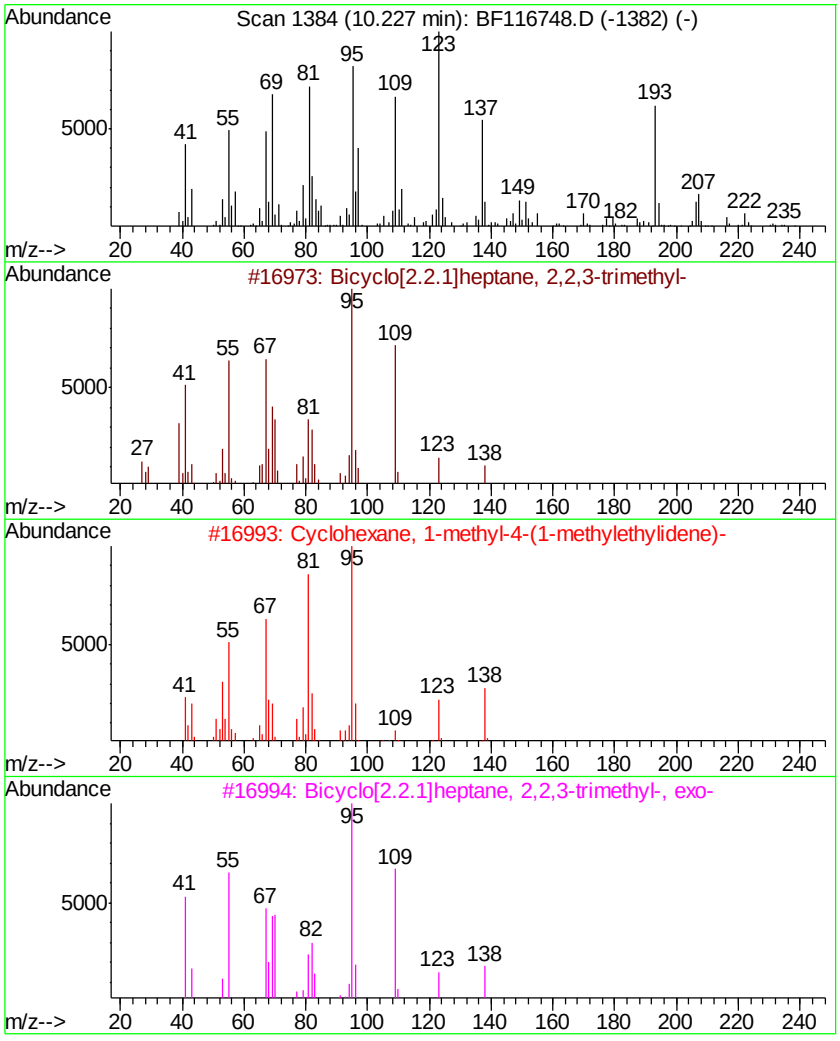
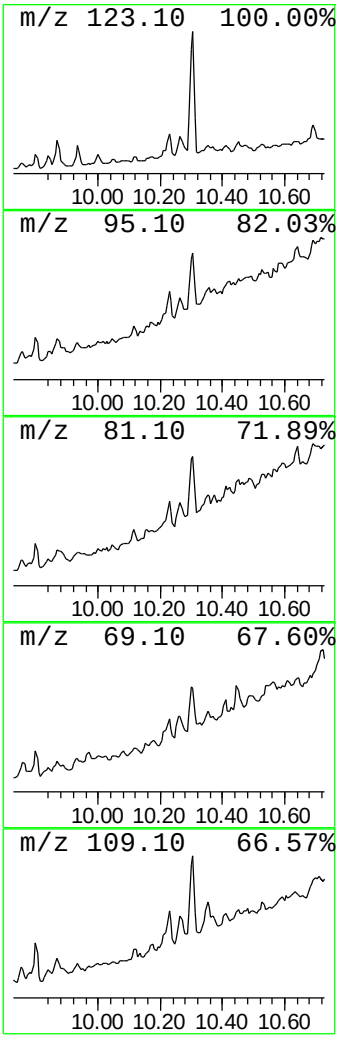
Peak Number 9 unknown10.23 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.23	5.91 ng	396772	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	46
2	Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-27-2	38
3	Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-41-8	38
4	2-(4-Fluorobenzoylamido)thiazole	222	C10H7FN2OS	313376-92-0	35
5	Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	000554-59-6	30





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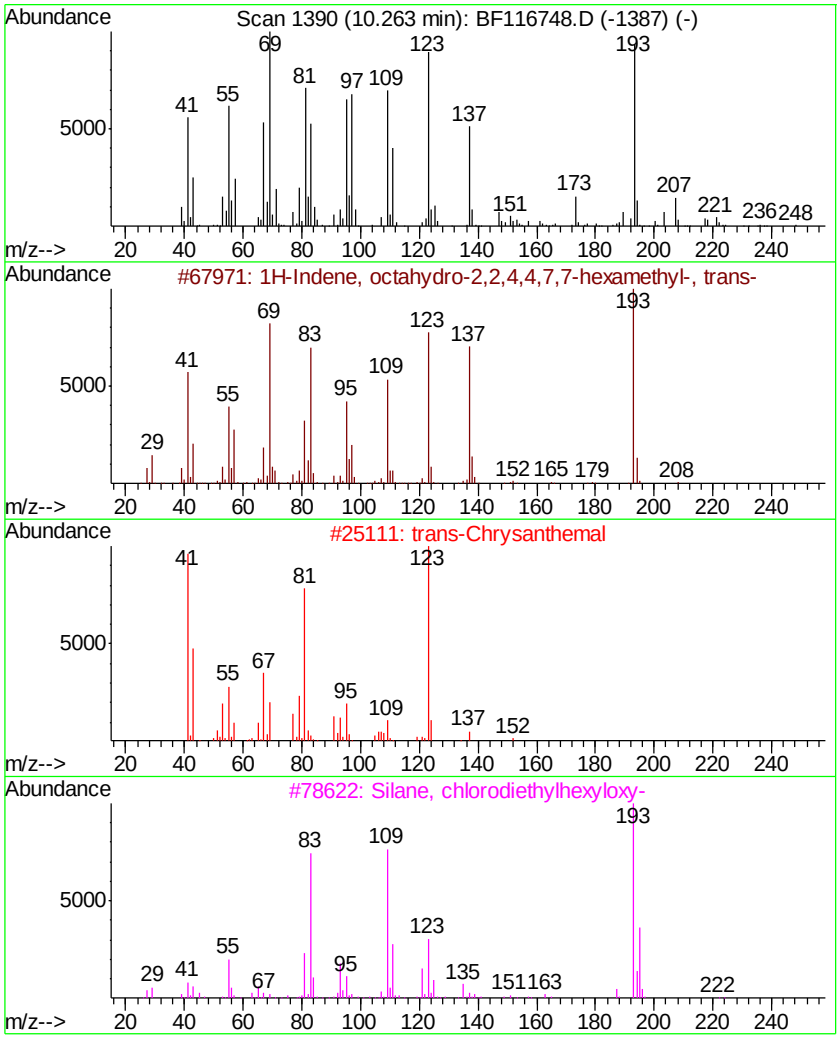
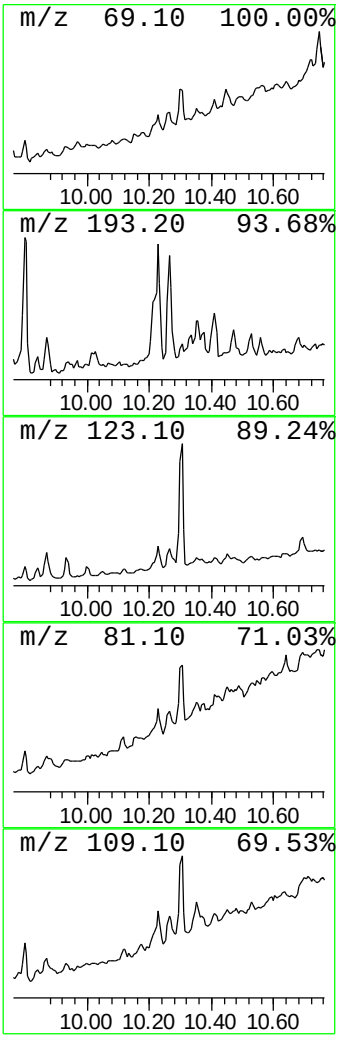
Peak Number 10 unknown10.26 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	10.69 ng	717128	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	35
2	trans-Chrysanthemal	152	C10H16O	020104-05-6	15
3	Silane, chlorodiethylhexyloxy-	222	C10H23ClOSi	1000363-74-4	14
4	Octane, 1,8-dibromo-	270	C8H16Br2	004549-32-0	14
5	3,4-Diethoxy-N-(4,5,6,7-tetrahyd...	346	C18H22N2O3S	1000311-37-1	14





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Data File : BF116748.D
Acq On : 1 Sep 2019 17:48
Operator : HP/JU
Sample : K4544-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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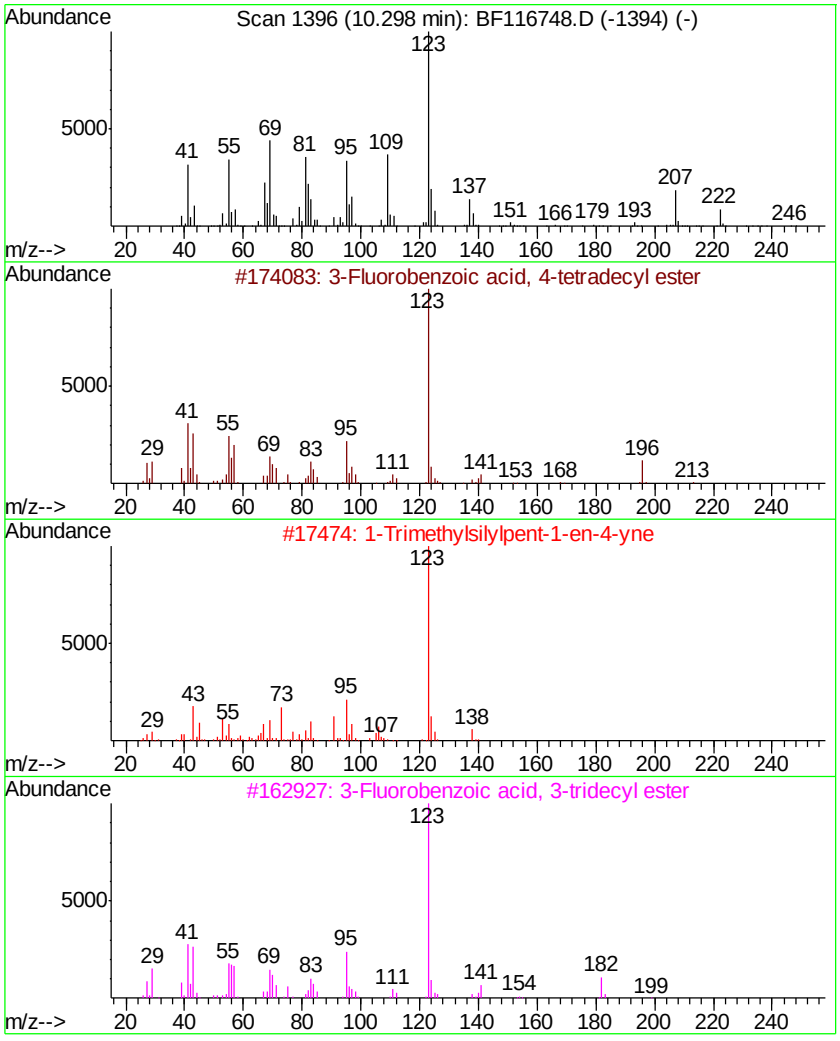
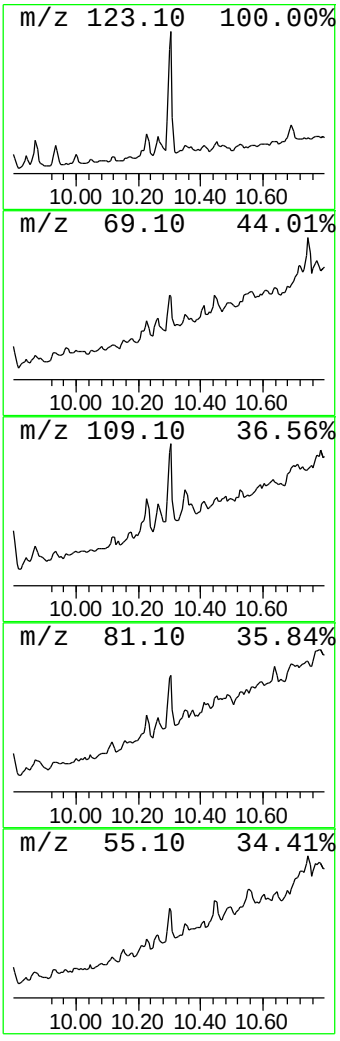
Peak Number 11 unknown10.30 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	20.96 ng	1405780	Acenaphthene-d10	9.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Fluorobenzoic acid, 4-tetradec...	336	C21H33FO2	1000280-60-6	47
2	1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	47
3	3-Fluorobenzoic acid, 3-tridecyl...	322	C20H31FO2	1000280-60-3	47
4	4-Fluorobenzoic acid, 5-pentadec...	350	C22H35FO2	1000280-68-6	47
5	2-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19FO2	1000278-98-1	43





Library Search Compound Report

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Instrument : BNA_F
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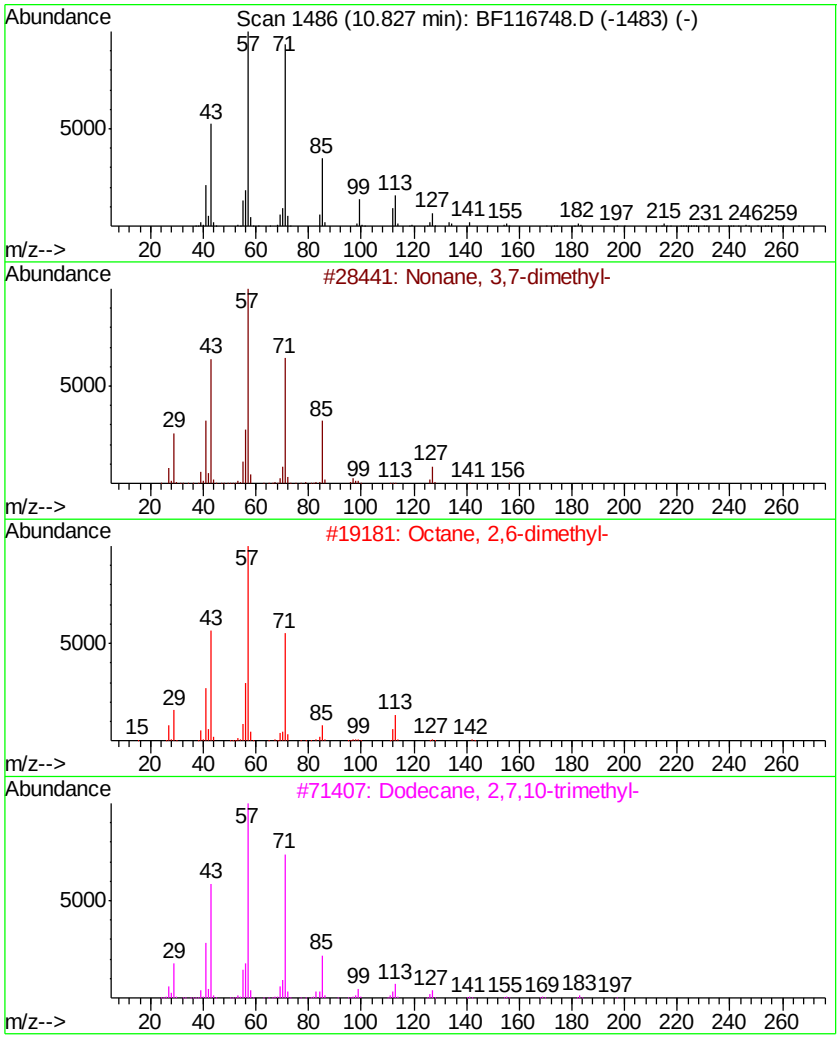
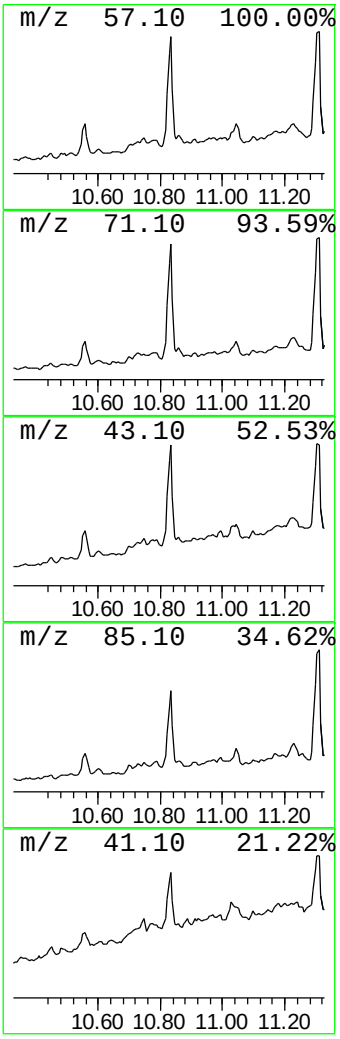
Peak Number 12 Nonane, 3,7-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.83	18.60 ng	3534310	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	81
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	80
3	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
4	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	78
5	Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	72





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Data File : BF116748.D
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Operator : HP/JU
Sample : K4544-19
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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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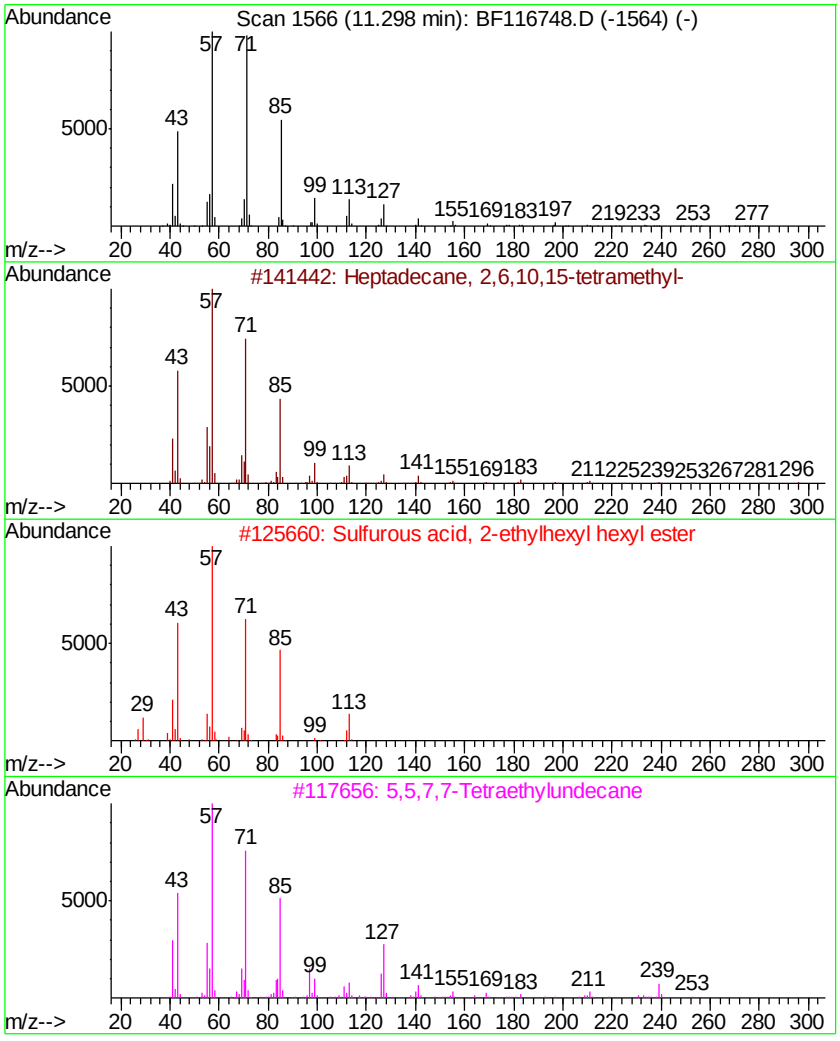
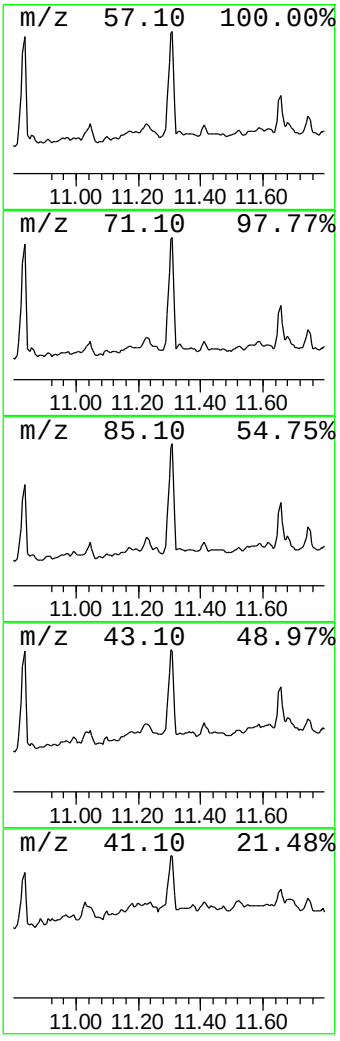
Peak Number 13 Heptadecane, 2,6,10,15-tetr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	20.00 ng	3800780	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
2	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	80
3	5,5,7,7-Tetraethylundecane	268	C19H40	1000360-42-0	72
4	Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	72
5	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	64





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Instrument : BNA_F
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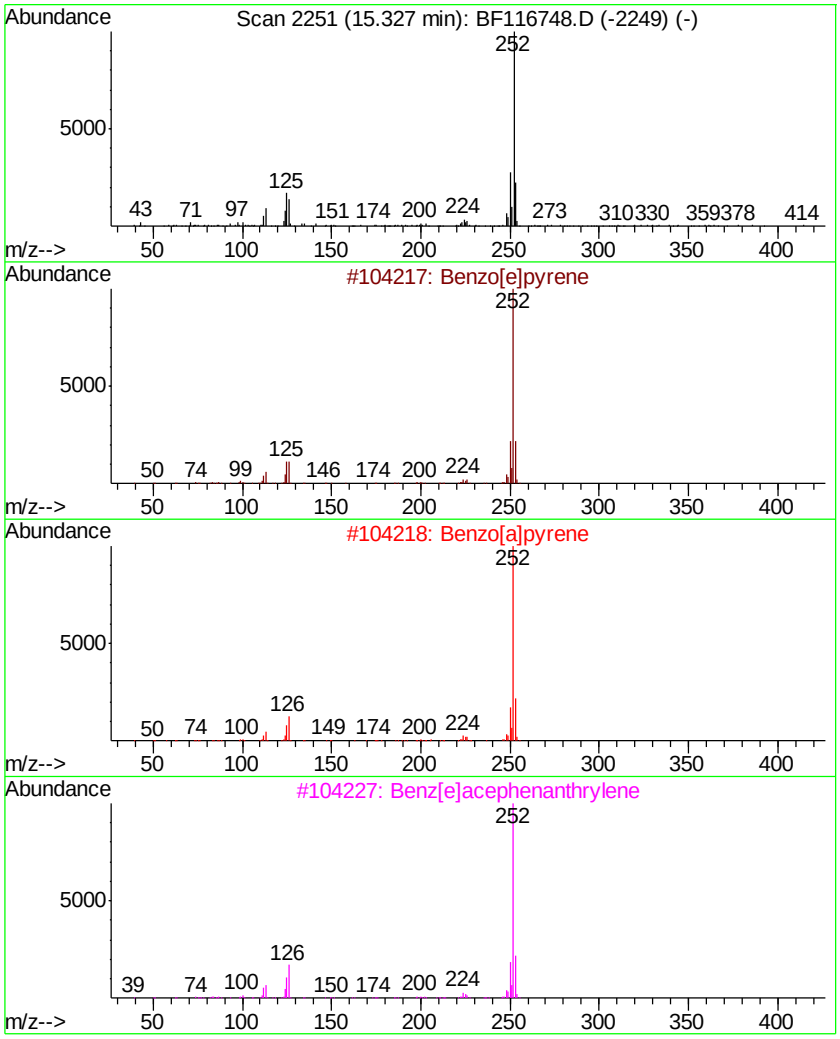
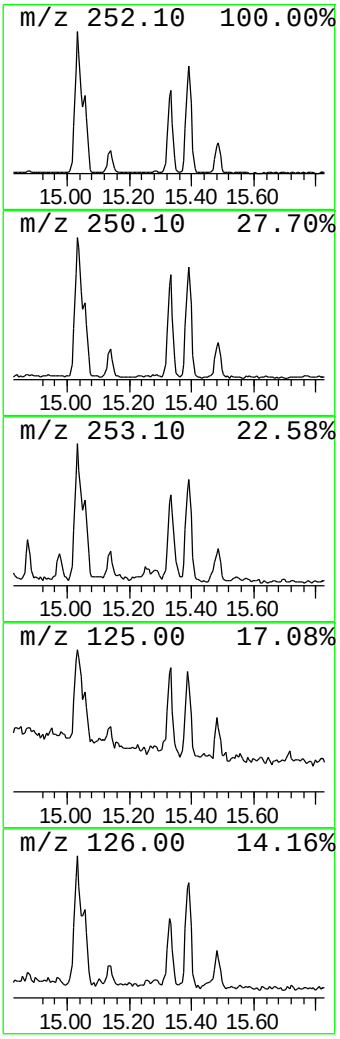
Peak Number 14 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	10.83 ng	313899	Perylene-d12	15.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	93
2	Benzo[a]pyrene	252	C20H12	000050-32-8	90
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
4	Benzo[j]fluoranthene	252	C20H12	000205-82-3	89
5	Perylene	252	C20H12	000198-55-0	89





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF090119\
Data File : BF116748.D
Acq On : 1 Sep 2019 17:48
Operator : HP/JU
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : T16-17-E1-E4-(0-2.25)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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
unknown6.61	6.61	66.1	ng	1446670	1	6.85	437796	20.0
unknown9.29	9.29	2.6	ng	173902	3	9.88	1341600	20.0
unknown9.76	9.76	7.2	ng	484056	3	9.88	1341600	20.0
unknown9.80	9.80	10.0	ng	669476	3	9.88	1341600	20.0
Cyclohexene, 1,5,...	9.93	6.3	ng	421758	3	9.88	1341600	20.0
4-isopropyl-1,6-d...	10.05	3.8	ng	252064	3	9.88	1341600	20.0
unknown10.15	10.15	8.9	ng	600183	3	9.88	1341600	20.0
unknown10.23	10.23	5.9	ng	396772	3	9.88	1341600	20.0
unknown10.26	10.26	10.7	ng	717128	3	9.88	1341600	20.0
unknown10.30	10.30	21.0	ng	1405780	3	9.88	1341600	20.0
Nonane, 3,7-dimet...	10.83	18.6	ng	3534310	4	11.39	3800780	20.0
Heptadecane, 2,6,...	11.30	20.0	ng	3800780	4	11.39	3800780	20.0
Benzo[e]pyrene	15.33	10.8	ng	313899	6	15.46	579721	20.0