LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

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-BF022218.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	2.152	5	11	25	rVB	996008	1360927	35.80%	3.649%
2	5.104	510	513	528	rVB	192987	254376	6.69%	0.682%
3	5.492	575	579	599	rBV	3481550	3801491	100.00%	10.193%
4	6.504	746	751	754	rBV	3419353	3484377	91.66%	9.343%
5	6.639	770	774	778	rBV	3632359	3624678	95.35%	9.719%
6	6.869	809	813	819	rBV	1241602	1013724	26.67%	2.718%
7	7.028	835	840	843	rBV	3053980	2815661	74.07%	7.550%
8	7.434	904	909	912	rBV	2464633	2408664	63.36%	6.459%
9	8.151	1027	1031	1043	rBV	1498363	1337221	35.18%	3.586%
10	9.228	1209	1214	1217	rBV	3497644	3330238	87.60%	8.930%
11	9.298	1223	1226	1229	rVB	64236	53255	1.40%	0.143%
12	9.616	1276	1280	1288	rVB	288471	291693	7.67%	0.782%
13	9.827	1312	1316	1320	rBV	155396	124614	3.28%	0.334%
14	9.910	1326	1330	1333	rBV	1299363	1270966	33.43%	3.408%
15	10.327	1398	1401	1403	rVB	164485	132321	3.48%	0.355%
16	10.545	1432	1438	1440	rBV	37638	48114	1.27%	0.129%
17	10.698	1460	1464	1475	rBV	1818954	1804341	47.46%	4.838%
18	10.798	1478	1481	1485	rBV	106425	92054	2.42%	0.247%
19	11.398	1579	1583	1585	rBV	1132581	999040	26.28%	2.679%
20	11.421	1585	1587	1592	rVV	323743	280562	7.38%	0.752%
21	11.474	1592	1596	1601	rVB	43041	45770	1.20%	0.123%
22	11.910	1664	1670	1672	rBV2	66711	106280	2.80%	0.285%
23	11.927	1672	1673	1678	rVB2	70103	48509	1.28%	0.130%
24	12.010	1684	1687	1690	rBV2	65876	79617	2.09%	0.213%
25	12.227	1721	1724	1728	rVB2	43349	43304	1.14%	0.116%
26	12.610	1785	1789	1793	rBV	488348	434847	11.44%	1.166%
27	12.792	1817	1820	1823	rVB	556032	475216	12.50%	1.274%
28	12.845	1823	1829	1831	rBV	409797	445152	11.71%	1.194%
29	12.863	1831	1832	1835	rVB	98529	73682	1.94%	0.198%
30	12.980	1848	1852	1855	rBV	2285737	2008004	52.82%	5.384%
31	13.063	1861	1866	1869	rBV2	33363	53177	1.40%	0.143%
32	13.274	1898	1902	1906	rBV2	36840	44690	1.18%	0.120%
33	13.492	1936	1939	1943	rBV	374216	341575	8.99%	0.916%
34	13.539	1943	1947	1949	rBV2	60802	53844	1.42%	0.144%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

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-BF022218.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.680	1968	1971	1978	rVB	44150	47085	1.24%	0.126%
36	13.868	1999	2003	2006	rBV2	335394	391371	10.30%	1.049%
37	14.045	2028	2033	2035	rBV2	910752	1047896	27.57%	2.810%
38	14.068	2035	2037	2044	rVB	224501	212551	5.59%	0.570%
39	14.186	2054	2057	2065	rBV2	91758	109878	2.89%	0.295%
40	14.498	2106	2110	2114	rBV4	102902	147149	3.87%	0.395%
41	14.804	2158	2162	2165	rVB2	101126	135067	3.55%	0.362%
42	14.880	2173	2175	2181	rVB3	58903	74866	1.97%	0.201%
43	14.939	2181	2185	2190	rVB4	42776	67926	1.79%	0.182%
44	15.098	2207	2212	2215	rBV	182926	284661	7.49%	0.763%
45	15.145	2218	2220	2223	rVB	85050	77640	2.04%	0.208%
46	15.204	2228	2230	2241	rVB2	42070	66820	1.76%	0.179%
47	15.404	2260	2264	2270	rVB	111896	153310	4.03%	0.411%
48	15.468	2271	2275	2280	rBV	143336	190543	5.01%	0.511%
49	15.533	2281	2286	2291	rBV	703409	938657	24.69%	2.517%
50	15.974	2356	2361	2365	rVB2	59452	94041	2.47%	0.252%
51	16.033	2367	2371	2380	rBV6	37141	85283	2.24%	0.229%
52	16.480	2443	2447	2451	rBV	37705	62973	1.66%	0.169%
53	17.045	2537	2543	2547	rBV2	45668	93592	2.46%	0.251%
54	17.080	2547	2549	2555	rVB4	29107	39083	1.03%	0.105%
55	17.498	2616	2620	2631	rVB	37761	79196	2.08%	0.212%
56	17.615	2634	2640	2648	rVB	43831	112796	2.97%	0.302%

Sum of corrected areas: 37294368

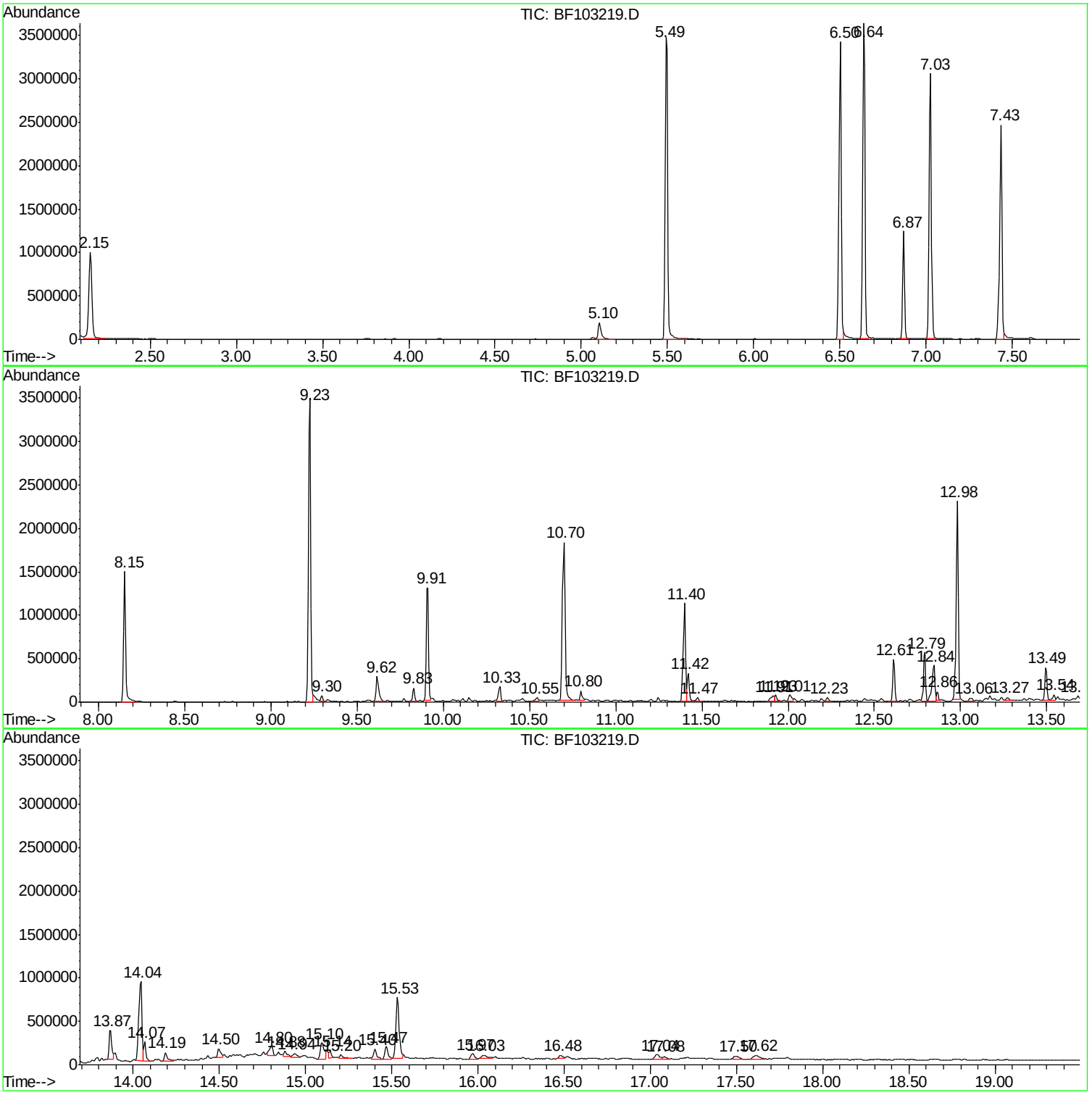
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

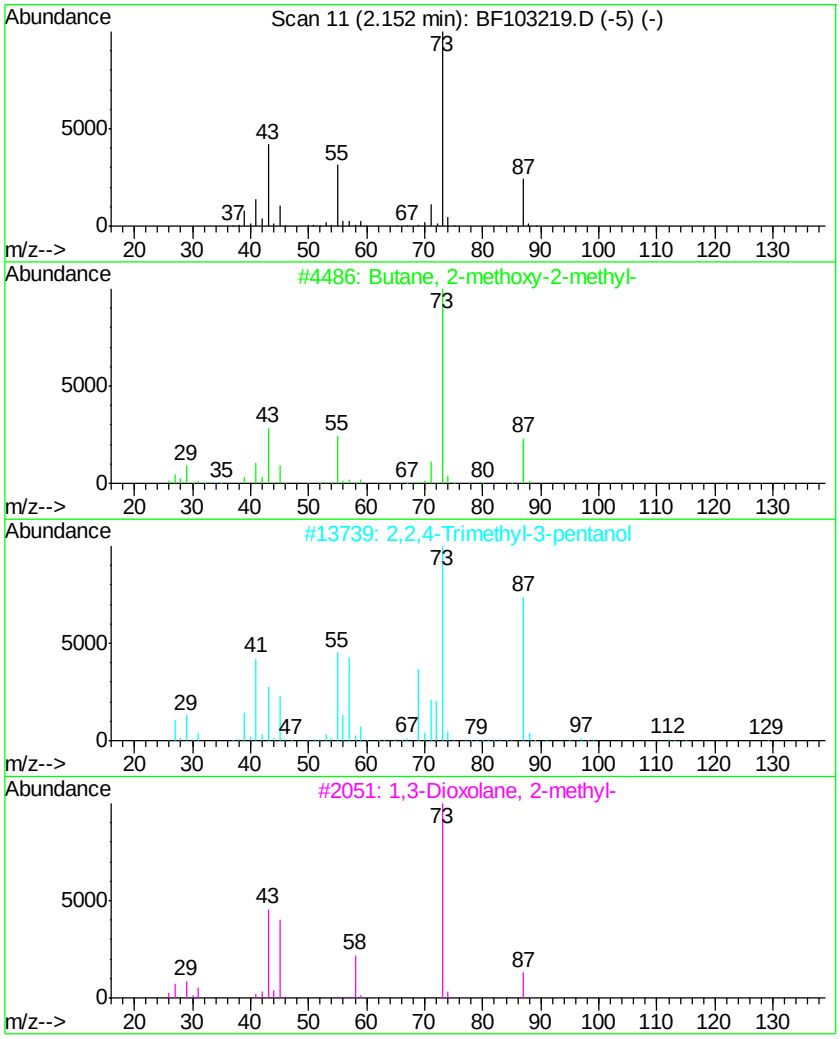
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

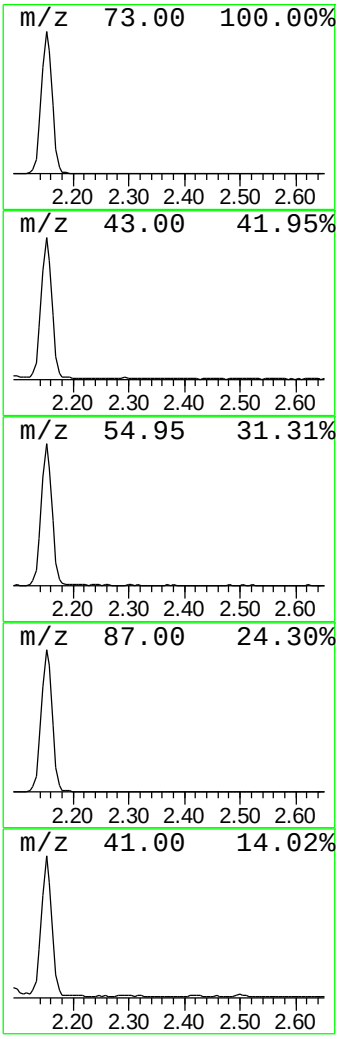
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.15	26.85 ng	1360930	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	10





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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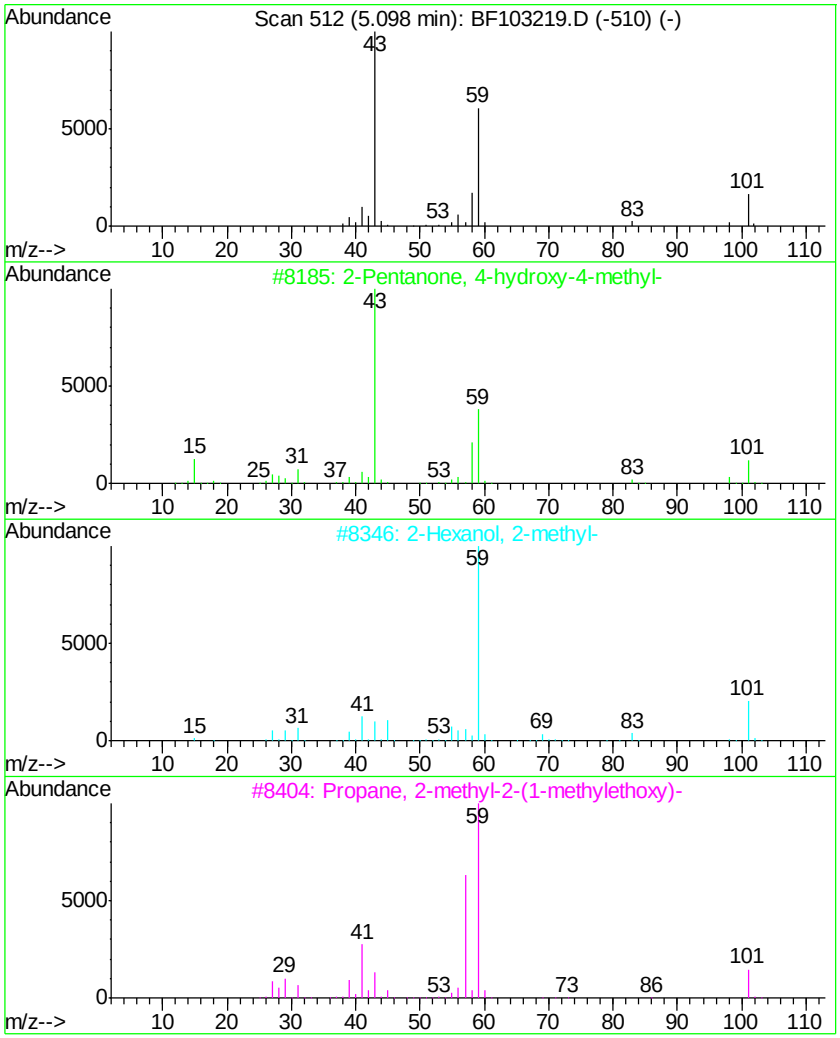
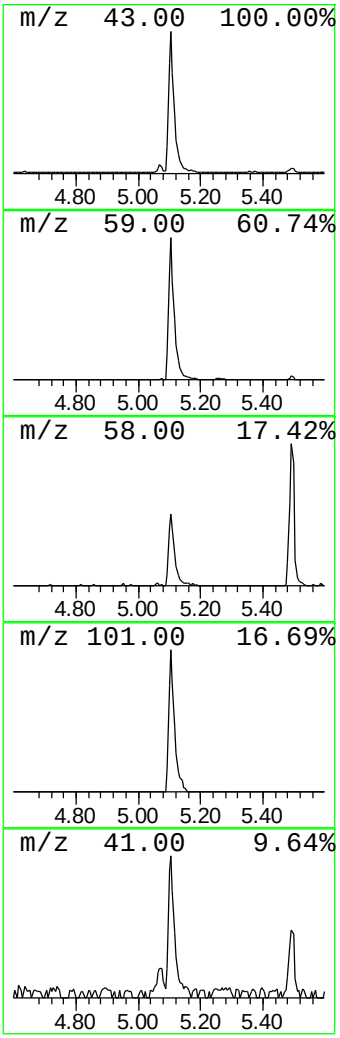
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	5.02 ng	254376	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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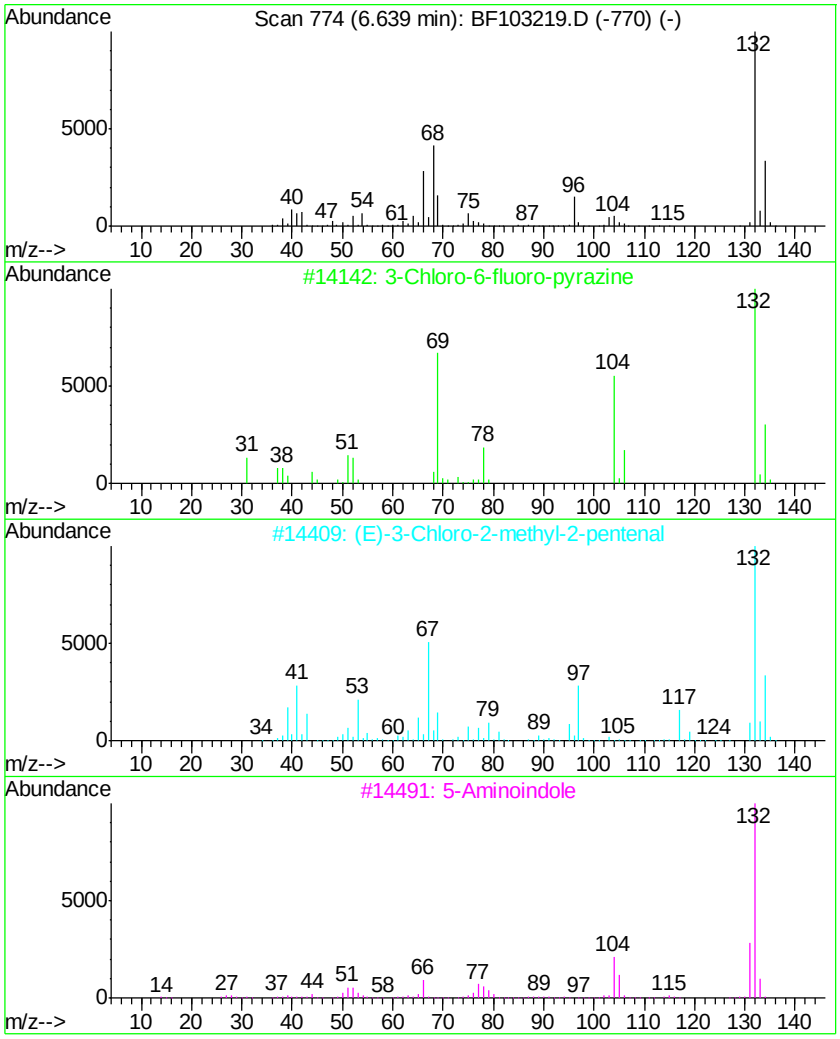
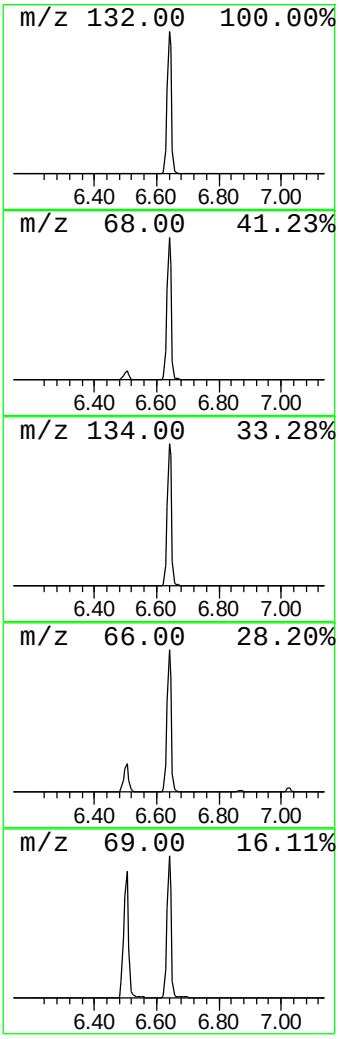
Peak Number 3 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	71.51 ng	3624680	1,4-Dichlorobenzene-d4	6.87

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	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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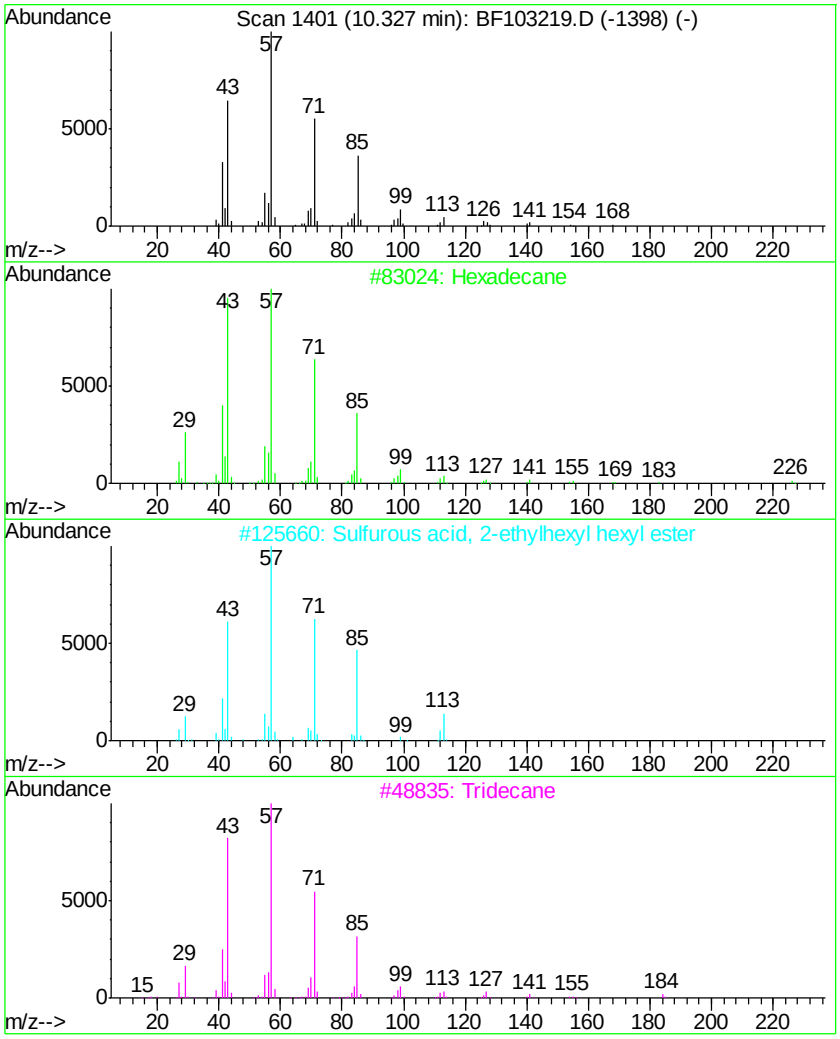
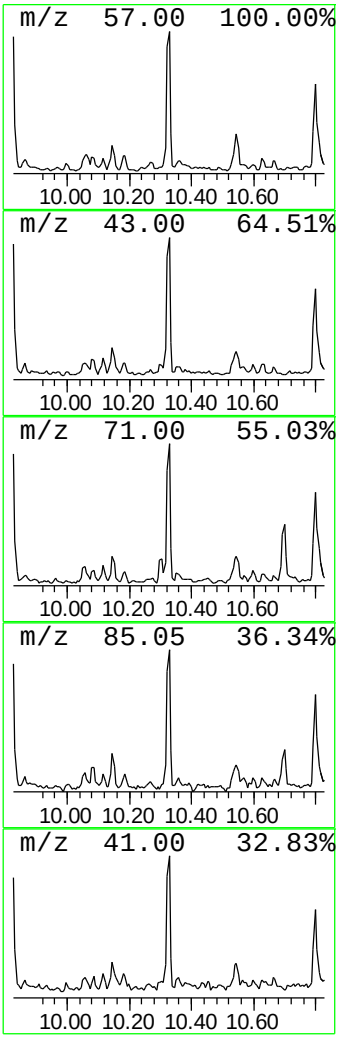
Peak Number 5 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	2.08 ng	132321	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	86
3	Tridecane	184	C13H28	000629-50-5	86
4	Heptadecane	240	C17H36	000629-78-7	86
5	Nonane, 5-butyl-	184	C13H28	017312-63-9	80





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Acq On : 26 Feb 2018 15:27
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Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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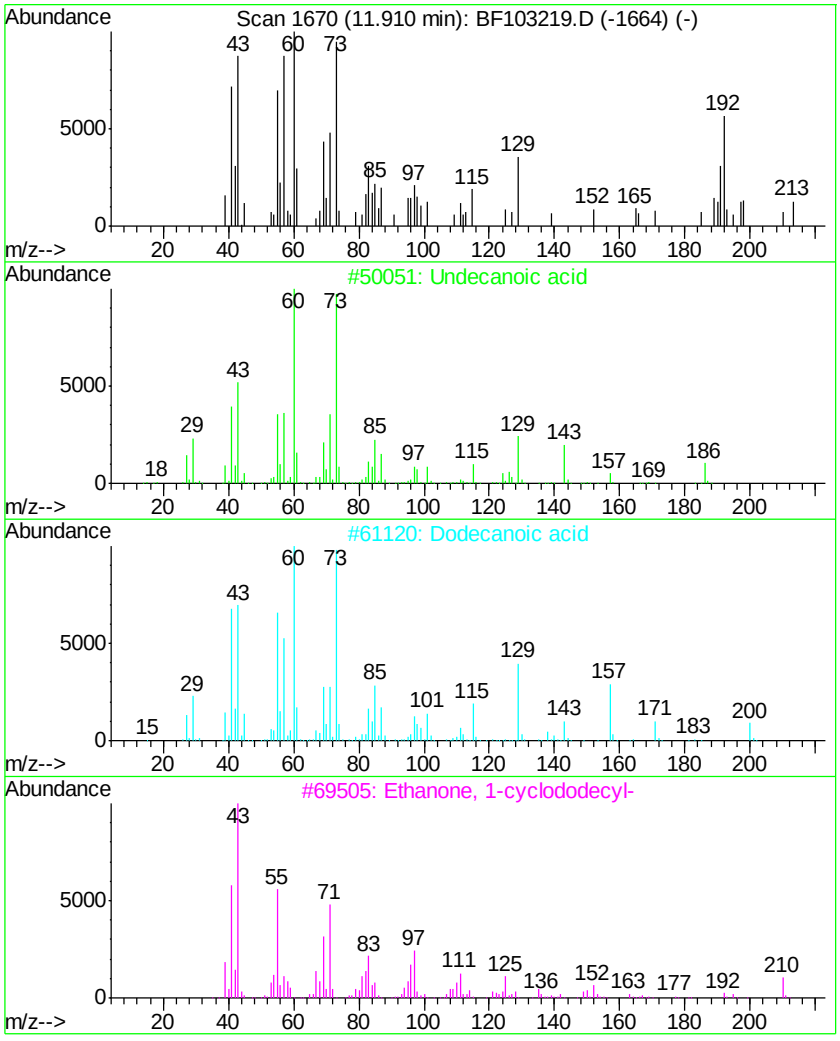
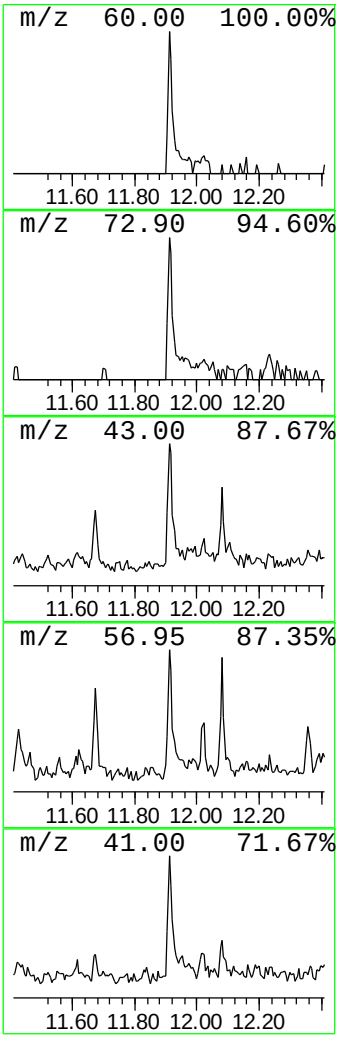
Peak Number 6 Undecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	2.13 ng	106280	Phenanthrene-d10	11.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecanoic acid	186	C11H22O2	000112-37-8	53
2	Dodecanoic acid	200	C12H24O2	000143-07-7	47
3	Ethanone, 1-cyclododecyl-	210	C14H26O	028925-00-0	25
4	1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	11
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	11





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Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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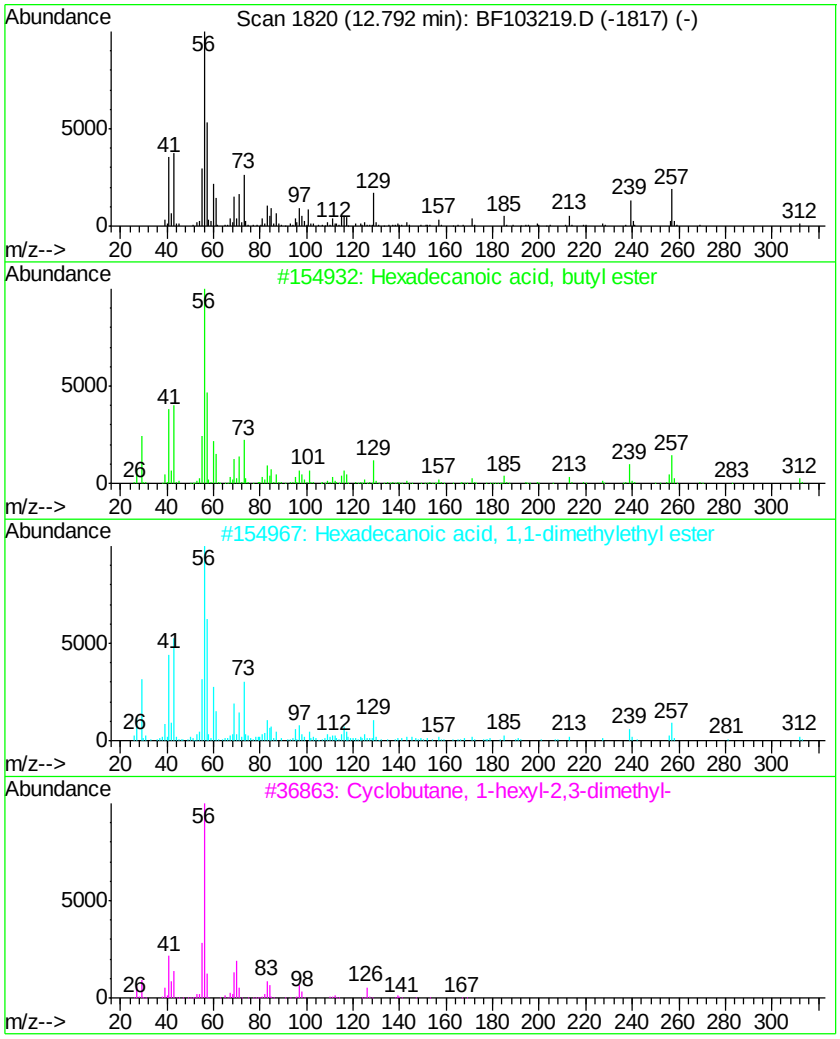
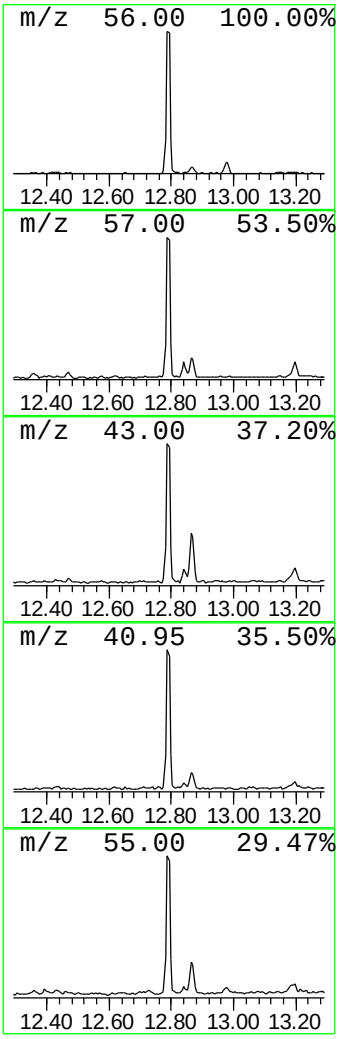
Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	9.07 ng	475216	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	80
3	Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	35
4	Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	35
5	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	32





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Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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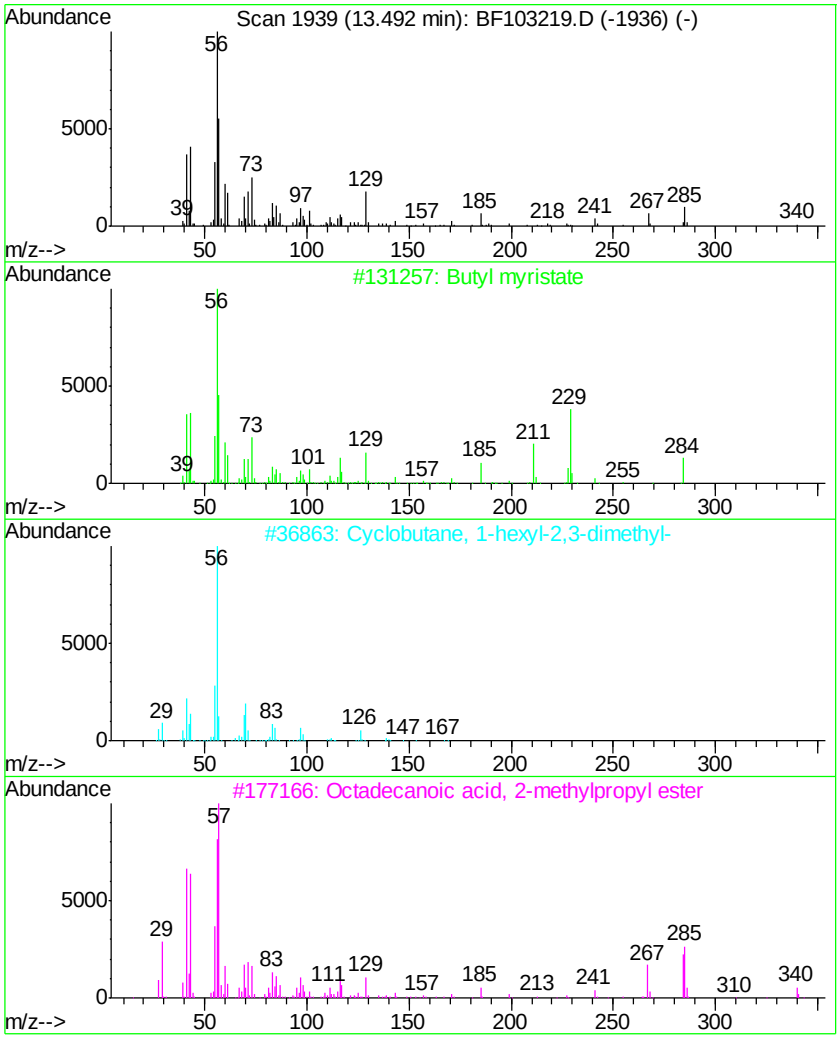
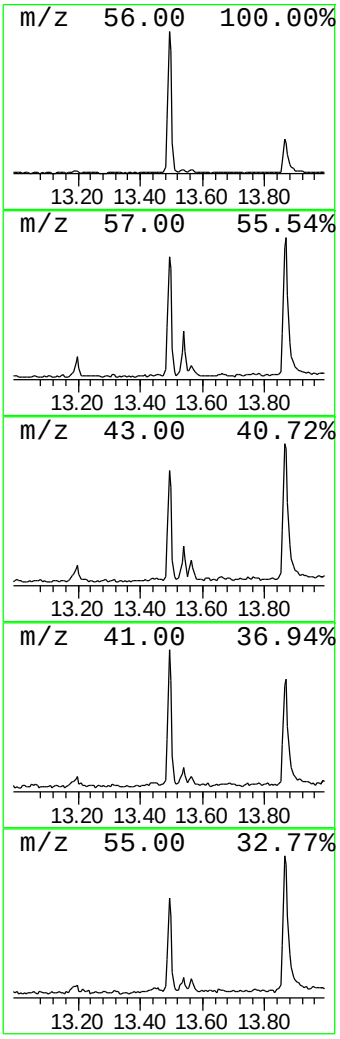
Peak Number 8 Butyl myristate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	6.52 ng	341575	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl myristate	284	C18H36O2	000110-36-1	86
2	Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	38
3	Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	37
4	Pentadecane, 8-methylene-	224	C16H32	055668-09-2	37
5	Trifluoroacetic acid, 4-methylpe...	198	C8H13F3O2	000327-71-9	37





Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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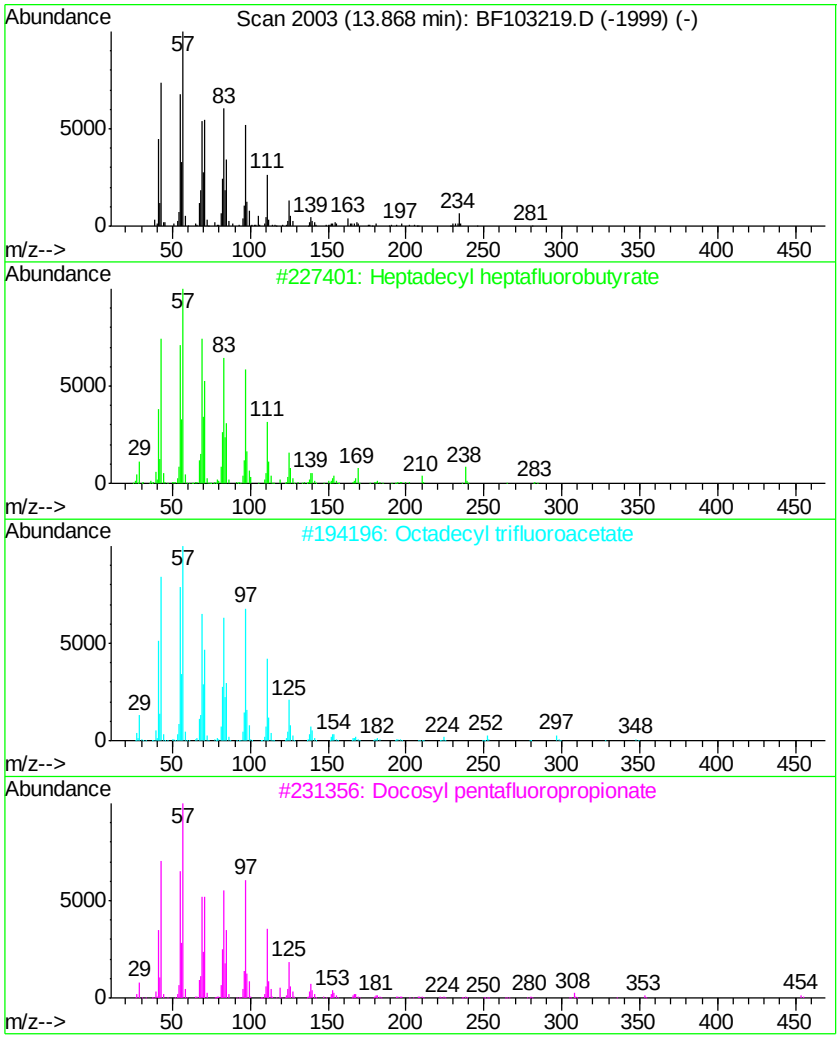
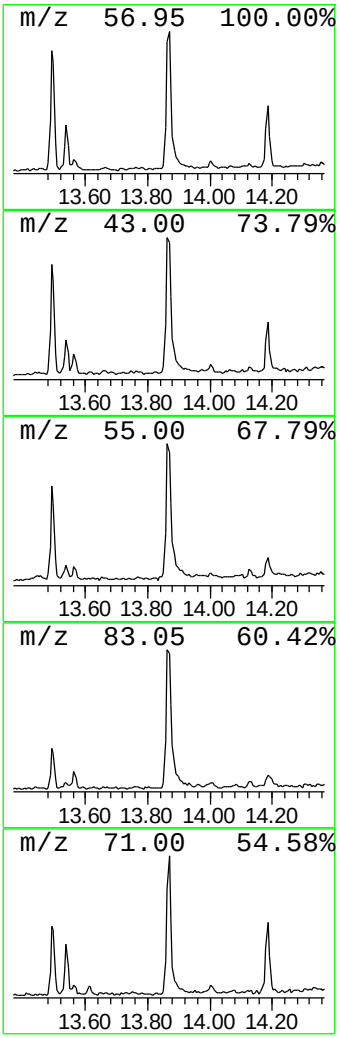
Peak Number 9 Heptadecyl heptafluorobutyrate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	7.47 ng	391371	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
2	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
3	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
4	Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91
5	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91





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Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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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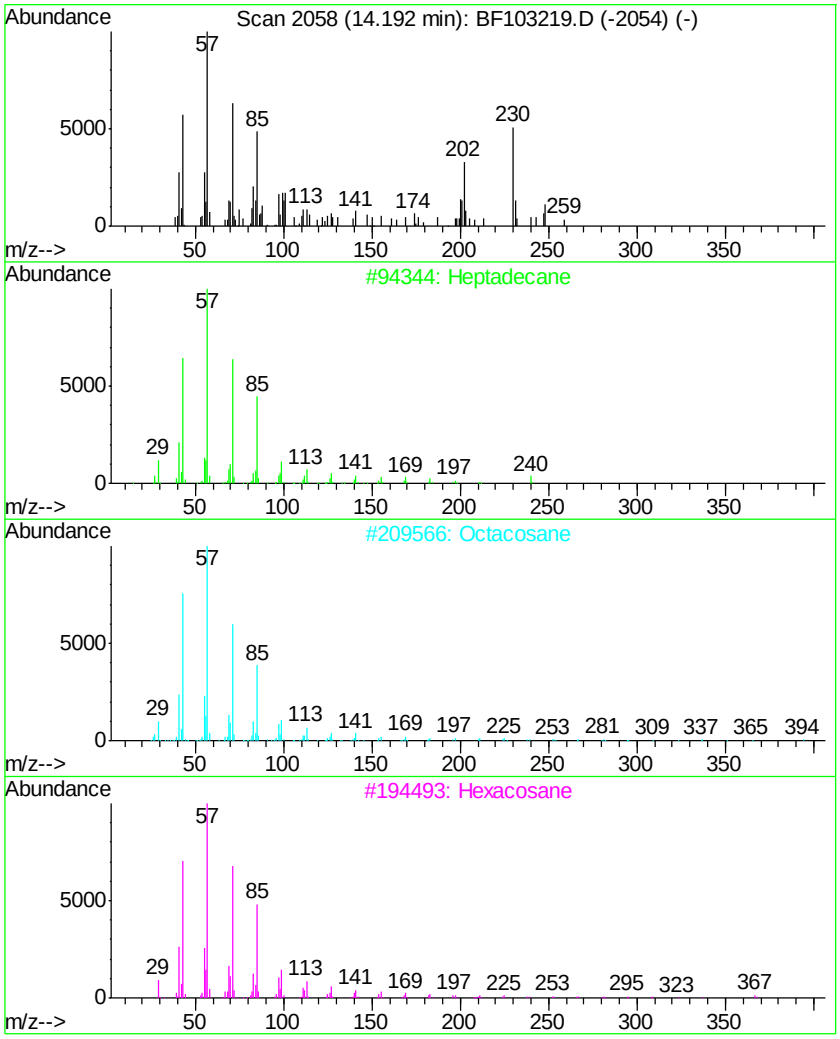
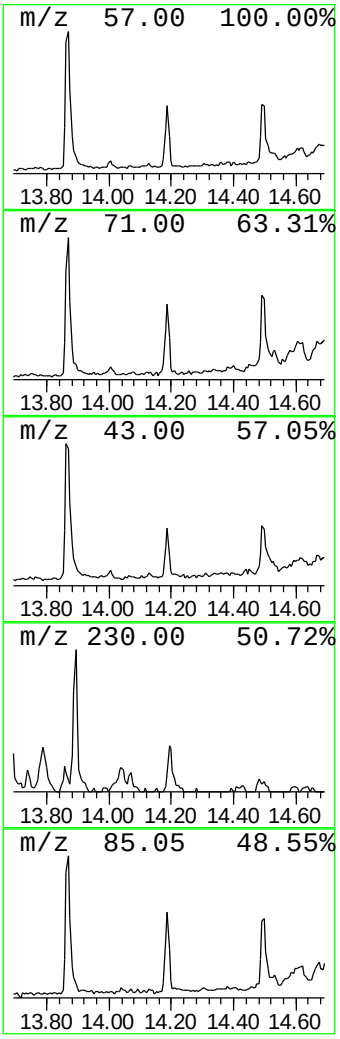
Peak Number 10 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.19	2.10 ng	109878	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	90
2	Octacosane	394	C28H58	000630-02-4	81
3	Hexacosane	366	C26H54	000630-01-3	81
4	Tetracosane	338	C24H50	000646-31-1	74
5	Heptacosane	380	C27H56	000593-49-7	72





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Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
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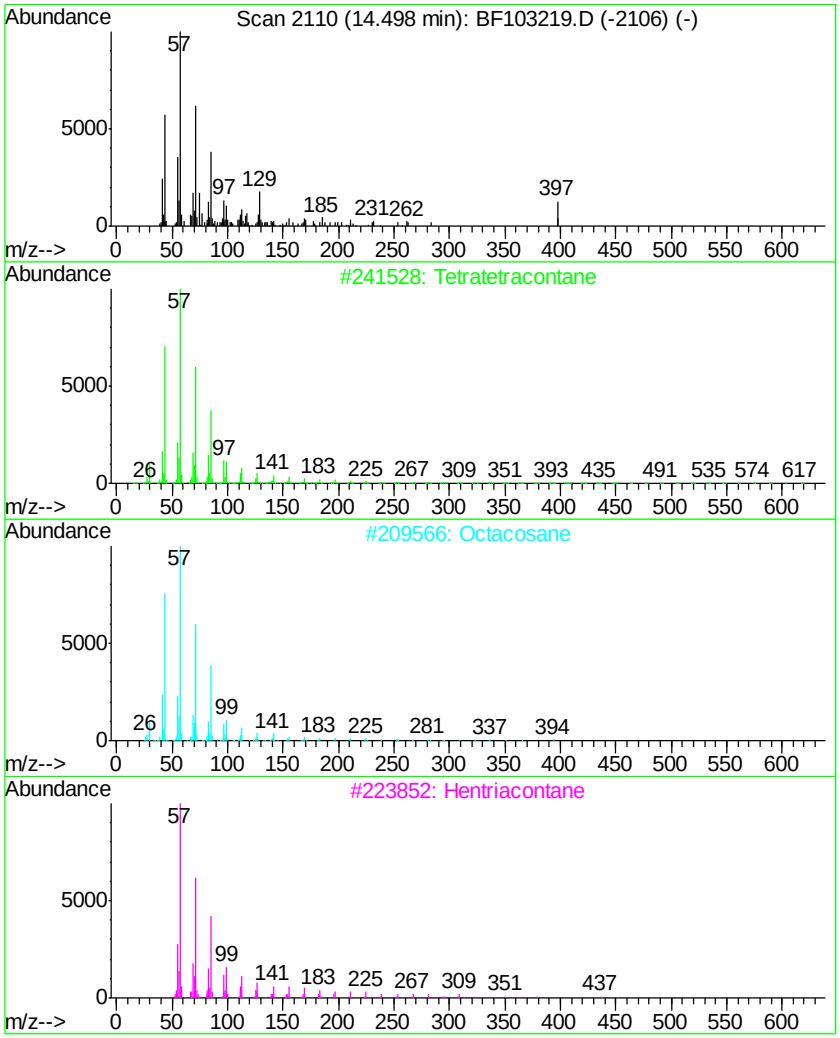
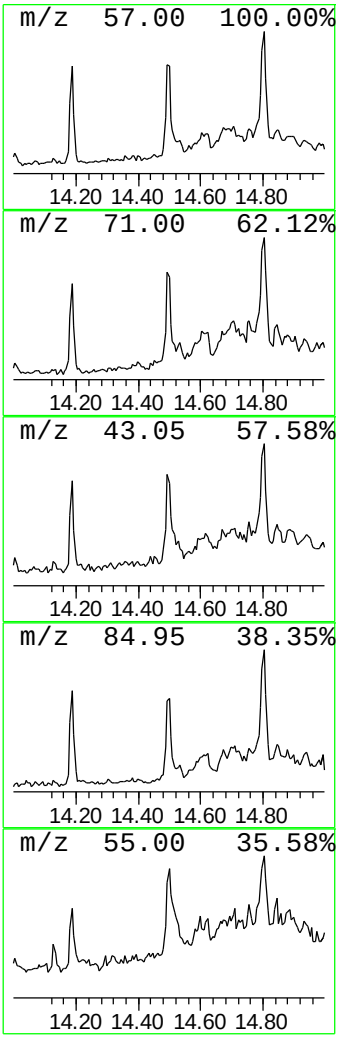
Peak Number 11 Tetratetracontane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	2.81 ng	147149	Chrysene-d12	14.04

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratetracontane	619	C44H90	007098-22-8	81
2	Octacosane	394	C28H58	000630-02-4	81
3	Hentriacontane	437	C31H64	000630-04-6	81
4	Hexacosane	366	C26H54	000630-01-3	70
5	Tetracosane	338	C24H50	000646-31-1	68





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Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
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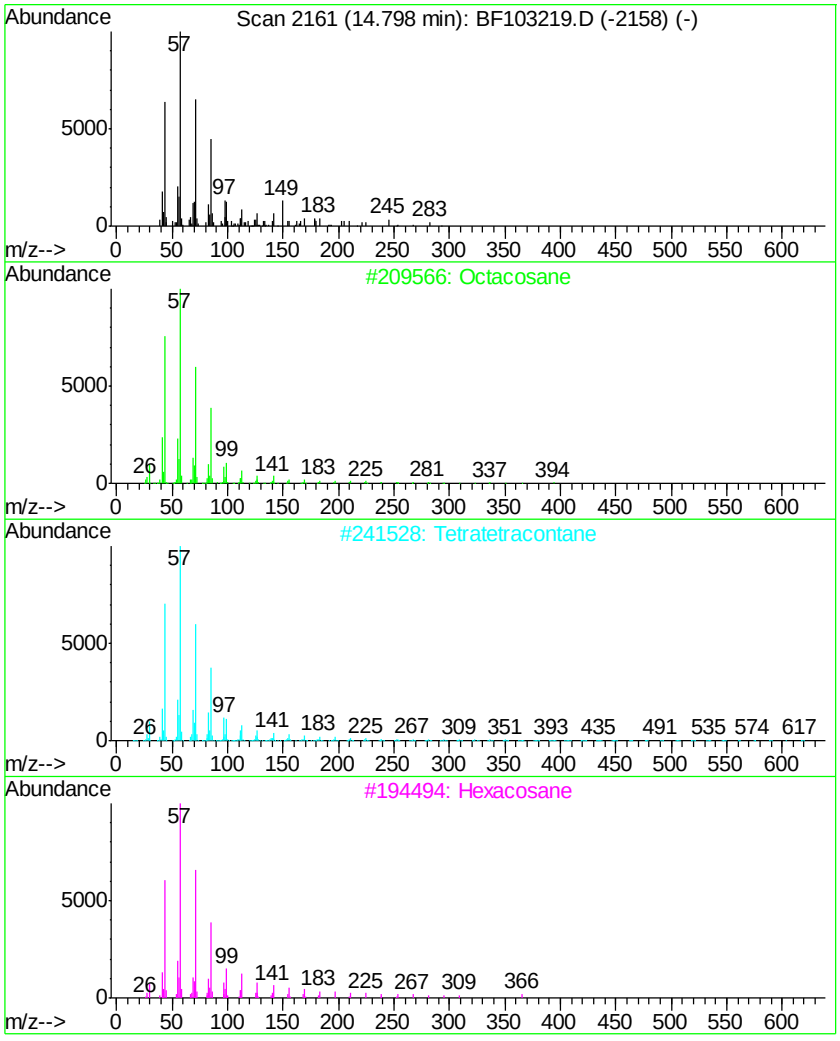
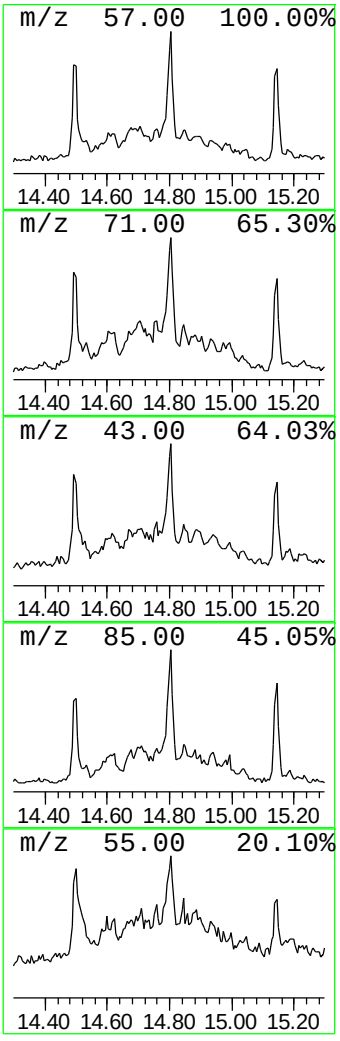
Peak Number 12 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	2.88 ng	135067	Perylene-d12	15.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Tetratetracontane	619	C44H90	007098-22-8	91
3	Hexacosane	366	C26H54	000630-01-3	90
4	Triacontane	422	C30H62	000638-68-6	90
5	Heneicosane	296	C21H44	000629-94-7	87





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Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

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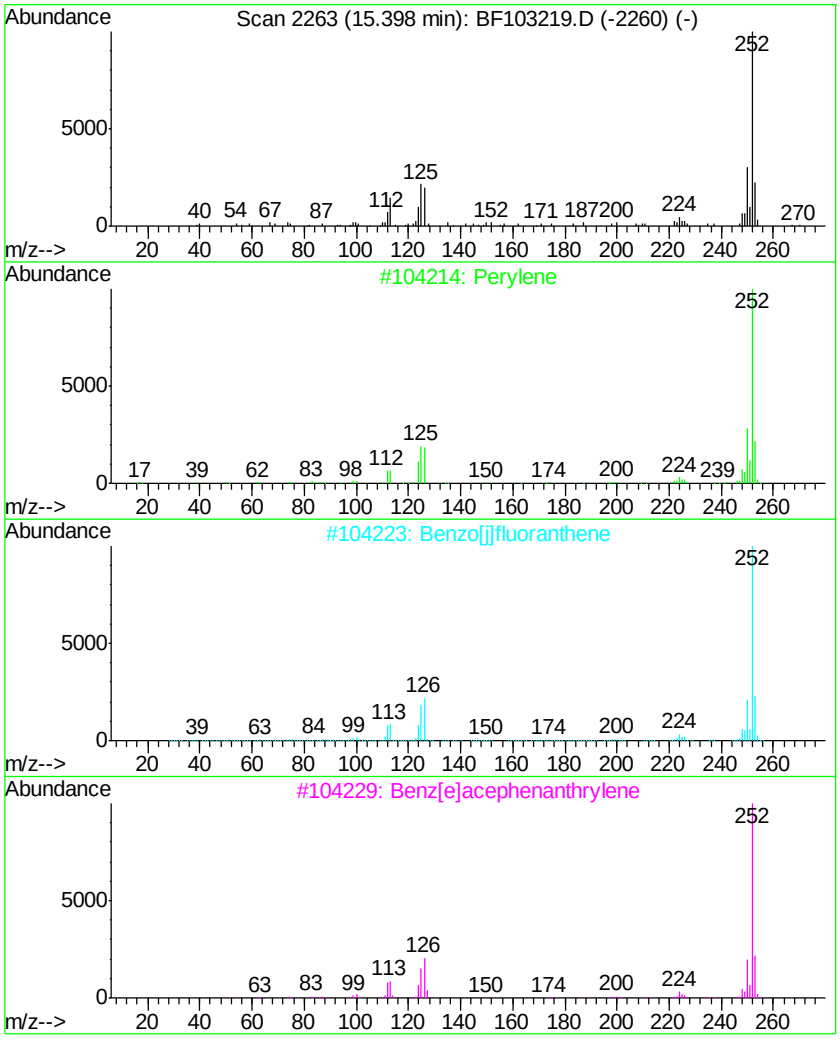
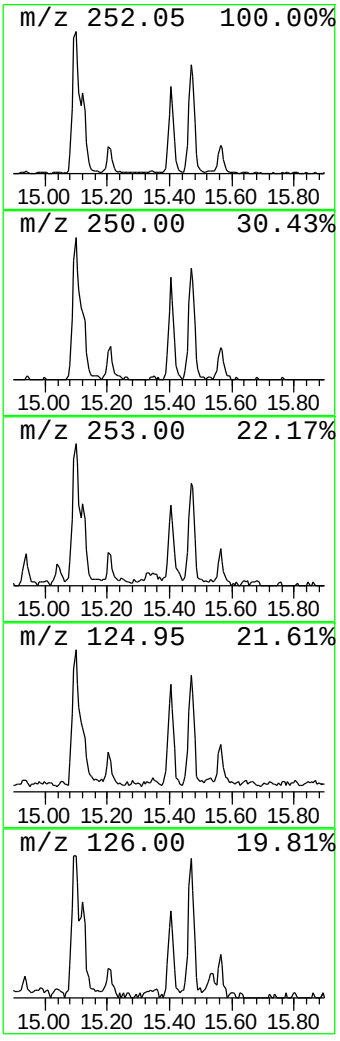
Peak Number 13 Perylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	3.27 ng	153310	Perylene-d12	15.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Perylene	252	C20H12	000198-55-0	99
2	Benzo[j]fluoranthene	252	C20H12	000205-82-3	98
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
4	Benzo[e]pyrene	252	C20H12	000192-97-2	96
5	Benzo[a]pyrene	252	C20H12	000050-32-8	94





Library Search Compound Report

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Data File : BF103219.D
Acq On : 26 Feb 2018 15:27
Operator : SJ/JU
Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

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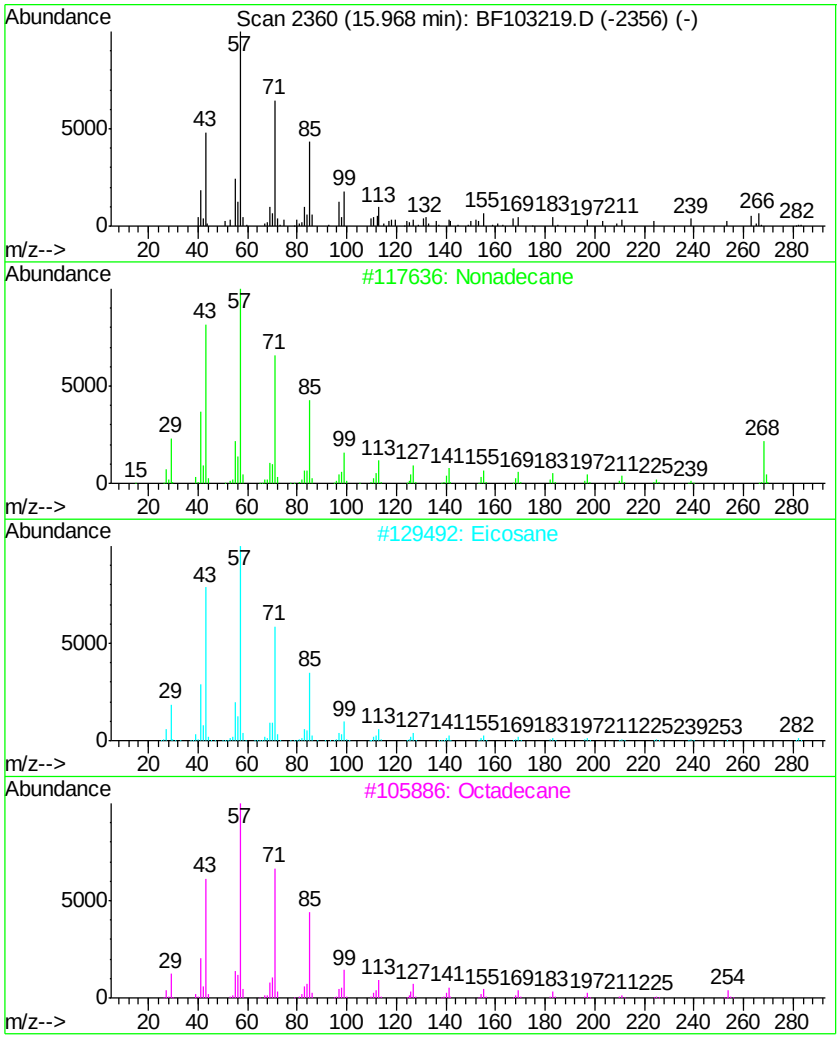
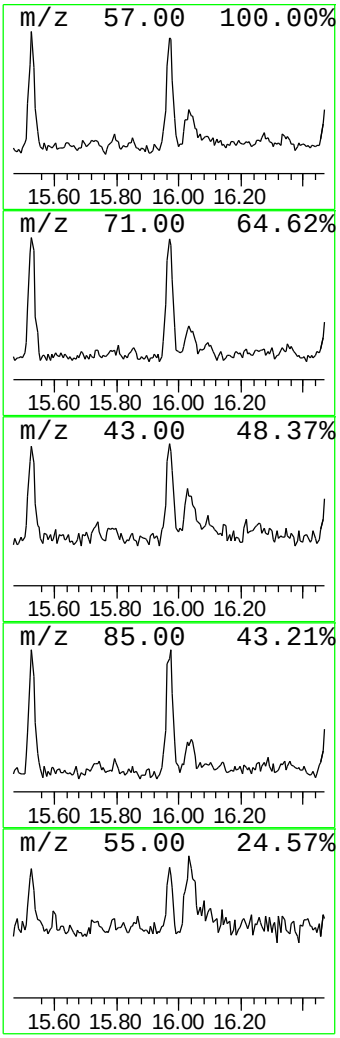
Peak Number 14 Nonadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	2.00 ng	94041	Perylene-d12	15.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	90
2	Eicosane	282	C20H42	000112-95-8	81
3	Octadecane	254	C18H38	000593-45-3	74
4	Docosane, 11-decyl-	451	C32H66	055401-55-3	74
5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	74





Library Search Compound Report

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Sample : J1660-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RT-2342

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

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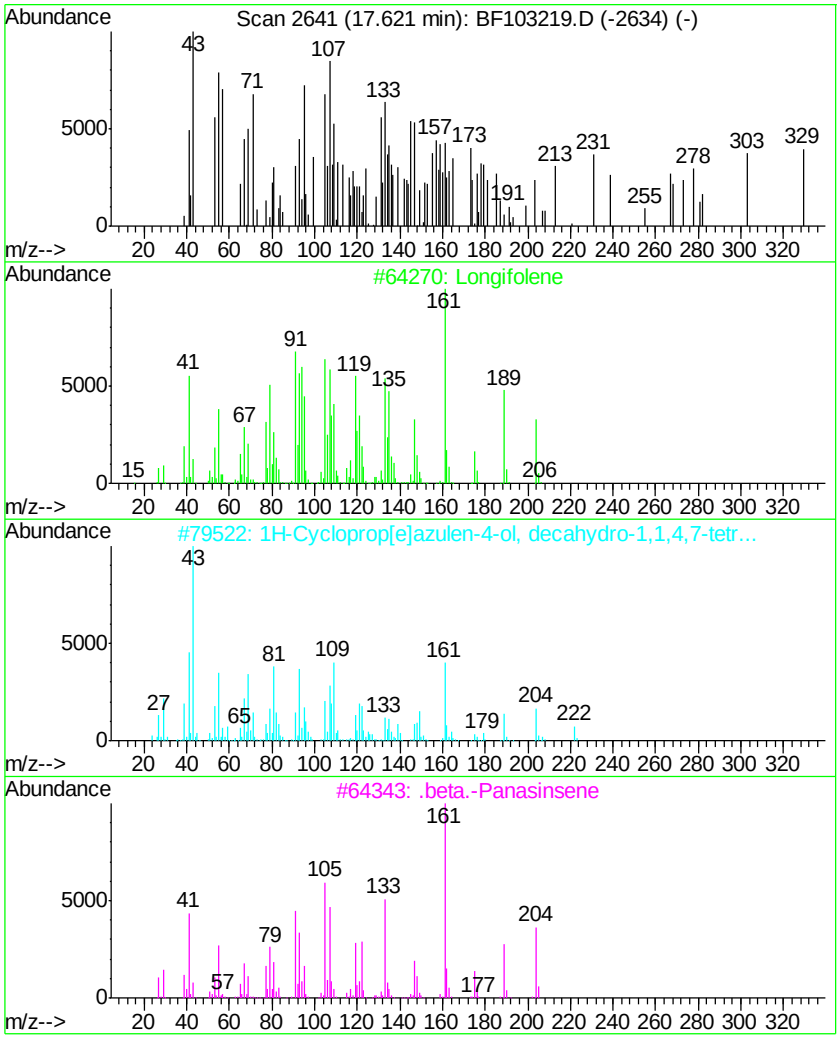
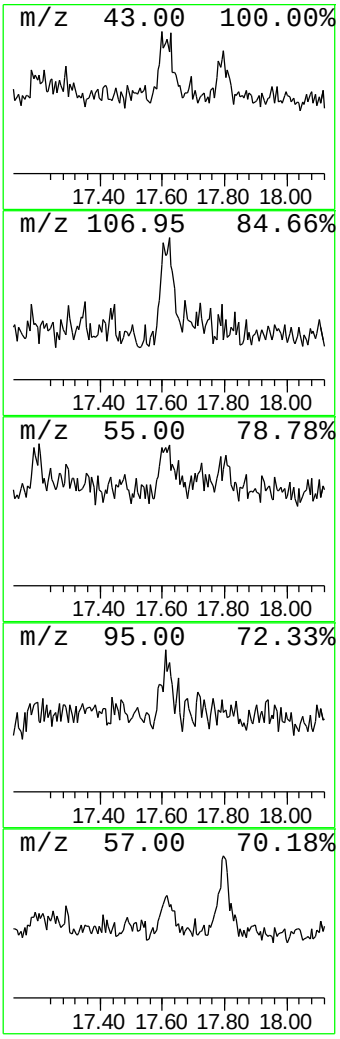
Peak Number 15 unknown17.62 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.62	2.40 ng	112796	Perylene-d12	15.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Longifolene	204	C15H24	000475-20-7	32
2	1H-Cycloprop[e]azulen-4-ol, deca...	222	C15H26O	000552-02-3	25
3	.beta.-Panasinsene	204	C15H24	1000159-39-0	16
4	Ethanone, 1-(1,2,3,4,7,7a-hexahy...	246	C17H26O	032388-58-2	16
5	Guaia-1(10),11-diene	204	C15H24	1000374-19-7	12





Tentatively Identified Compound (LSC) summary

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Operator : SJ/JU
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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
Butane, 2-methoxy...	2.15	26.9	ng	1360930	1	6.87	1013720	20.0
2-Pentanone, 4-hy...	5.10	5.0	ng	254376	1	6.87	1013720	20.0
unknown6.64	6.64	71.5	ng	3624680	1	6.87	1013720	20.0
Hexadecane	10.33	2.1	ng	132321	3	9.90	1270970	20.0
Undecanoic acid	11.91	2.1	ng	106280	4	11.40	999040	20.0
Hexadecanoic acid...	12.79	9.1	ng	475216	5	14.04	1047900	20.0
Butyl myristate	13.49	6.5	ng	341575	5	14.04	1047900	20.0
Heptadecyl heptaf...	13.87	7.5	ng	391371	5	14.04	1047900	20.0
Heptadecane	14.19	2.1	ng	109878	5	14.04	1047900	20.0
Tetratetracontane	14.50	2.8	ng	147149	5	14.04	1047900	20.0
Octacosane	14.80	2.9	ng	135067	6	15.53	938657	20.0
Perylene	15.40	3.3	ng	153310	6	15.53	938657	20.0
Nonadecane	15.97	2.0	ng	94041	6	15.53	938657	20.0
unknown17.62	17.62	2.4	ng	112796	6	15.53	938657	20.0