LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
Data File : BF113053.D
Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

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-BF022719.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	3.716	268	277	283	rVB2	36718	59423	1.17%	0.101%
2	5.345	549	554	557	rBV	4083972	4068116	79.97%	6.892%
3	6.375	723	729	745	rVB	3735748	4349519	85.50%	7.368%
4	6.498	745	750	753	rBV	4351610	4300058	84.53%	7.285%
5	6.728	785	789	792	rBV	1313069	1220903	24.00%	2.068%
6	6.881	811	815	819	rBV	3521336	3376452	66.37%	5.720%
7	7.286	878	884	887	rBV	2681347	2768128	54.42%	4.689%
8	8.010	1002	1007	1013	rBV	1499678	1533577	30.15%	2.598%
9	9.086	1184	1190	1193	rBV	4798161	4758991	93.55%	8.062%
10	9.204	1208	1210	1215	rVB	71615	64739	1.27%	0.110%
11	9.475	1251	1256	1267	rVB	456483	398633	7.84%	0.675%
12	9.757	1299	1304	1308	rBV2	1877777	1722030	33.85%	2.917%
13	10.545	1432	1438	1441	rBV	3357360	3307907	65.03%	5.604%
14	11.233	1551	1555	1558	rBV	1929896	1929442	37.93%	3.269%
15	11.257	1558	1559	1563	rVV	418373	266425	5.24%	0.451%
16	11.310	1565	1568	1572	rVB	71182	70285	1.38%	0.119%
17	11.469	1590	1595	1602	rBV3	37422	80976	1.59%	0.137%
18	11.733	1637	1640	1643	rBV	85345	85722	1.69%	0.145%
19	11.780	1643	1648	1656	rVV2	1490190	1716814	33.75%	2.908%
20	11.845	1656	1659	1661	rVV	189413	203043	3.99%	0.344%
21	11.869	1661	1663	1668	rVB	87306	85407	1.68%	0.145%
22	12.027	1687	1690	1692	rBV	55721	54994	1.08%	0.093%
23	12.051	1692	1694	1700	rVB2	73265	76085	1.50%	0.129%
24	12.180	1711	1716	1719	rBV3	53745	68643	1.35%	0.116%
25	12.286	1731	1734	1737	rBV2	121728	137845	2.71%	0.234%
26	12.333	1740	1742	1745	rVB2	93532	91518	1.80%	0.155%
27	12.363	1745	1747	1753	rVB3	73345	90209	1.77%	0.153%
28	12.439	1756	1760	1764	rBV	1021452	979486	19.25%	1.659%
29	12.557	1776	1780	1785	rBV	655708	708436	13.93%	1.200%
30	12.668	1793	1799	1802	rBV	1025289	1073341	21.10%	1.818%
31	12.704	1802	1805	1807	rVV2	105474	128291	2.52%	0.217%
32	12.727	1807	1809	1813	rVB2	79432	91275	1.79%	0.155%
33	12.821	1820	1825	1828	rVB	4924126	5086997	100.00%	8.618%
34	12.886	1833	1836	1840	rBV2	85209	108720	2.14%	0.184%

LSC Area Percent Report

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Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

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-BF022719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.974	1848	1851	1852	rBV3	68194	68087	1.34%	0.115%
36	12.998	1852	1855	1858	rVB	197609	191854	3.77%	0.325%
37	13.063	1858	1866	1869	rBV2	202156	290420	5.71%	0.492%
38	13.098	1869	1872	1875	rBV	158667	154277	3.03%	0.261%
39	13.192	1885	1888	1891	rBV	111239	89421	1.76%	0.151%
40	13.221	1891	1893	1897	rVB2	109288	107548	2.11%	0.182%
41	13.398	1920	1923	1926	rBV	143788	116362	2.29%	0.197%
42	13.498	1938	1940	1945	rVB	103526	124323	2.44%	0.211%
43	13.610	1955	1959	1962	rBV2	122875	173545	3.41%	0.294%
44	13.639	1962	1964	1966	rVV	93063	78829	1.55%	0.134%
45	13.663	1966	1968	1972	rVV2	72754	88040	1.73%	0.149%
46	13.721	1973	1978	1985	rVV2	881389	1216769	23.92%	2.061%
47	13.863	1994	2002	2004	rBV2	2397526	2664689	52.38%	4.514%
48	13.886	2004	2006	2009	rVB	558412	429354	8.44%	0.727%
49	13.963	2014	2019	2024	rBV3	105542	152139	2.99%	0.258%
50	14.039	2029	2032	2036	rVB2	271042	272437	5.36%	0.462%
51	14.245	2063	2067	2070	rBV	116690	147154	2.89%	0.249%
52	14.280	2070	2073	2076	rVB2	89281	79295	1.56%	0.134%
53	14.339	2080	2083	2086	rBV2	319205	325981	6.41%	0.552%
54	14.610	2126	2129	2132	rBV	564639	601923	11.83%	1.020%
55	14.639	2132	2134	2141	rVB2	318232	324808	6.39%	0.550%
56	14.704	2142	2145	2150	rBV2	83648	112082	2.20%	0.190%
57	14.868	2169	2173	2183	rVB	524873	930665	18.29%	1.577%
58	14.951	2183	2187	2197	rBV2	269229	432198	8.50%	0.732%
59	15.151	2217	2221	2227	rVB	321940	373752	7.35%	0.633%
60	15.210	2227	2231	2235	rBV	418097	511334	10.05%	0.866%
61	15.268	2236	2241	2244	rVV	1434337	1793098	35.25%	3.038%
62	15.304	2244	2247	2253	rVB3	277931	412892	8.12%	0.699%
63	15.709	2312	2316	2320	rBV	181225	247062	4.86%	0.419%
64	15.762	2321	2325	2337	rVV7	142627	396095	7.79%	0.671%
65	16.174	2392	2395	2398	rBV2	64000	104981	2.06%	0.178%
66	16.209	2399	2401	2410	rVB	126022	168343	3.31%	0.285%
67	16.621	2466	2471	2481	rVB2	214868	421280	8.28%	0.714%
68	17.027	2535	2540	2550	rVB	176540	365757	7.19%	0.620%

Sum of corrected areas: 59029224

LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
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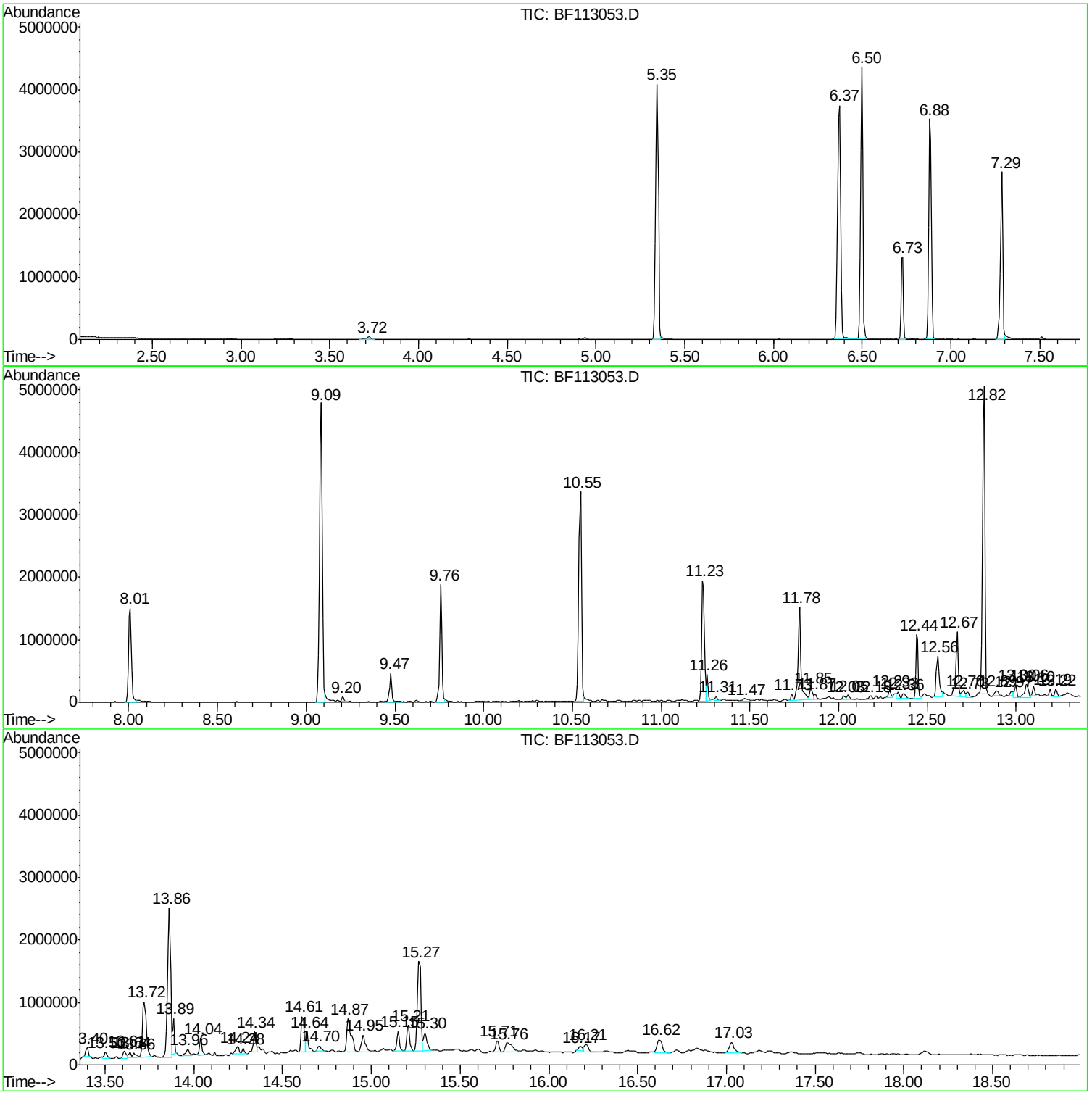
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
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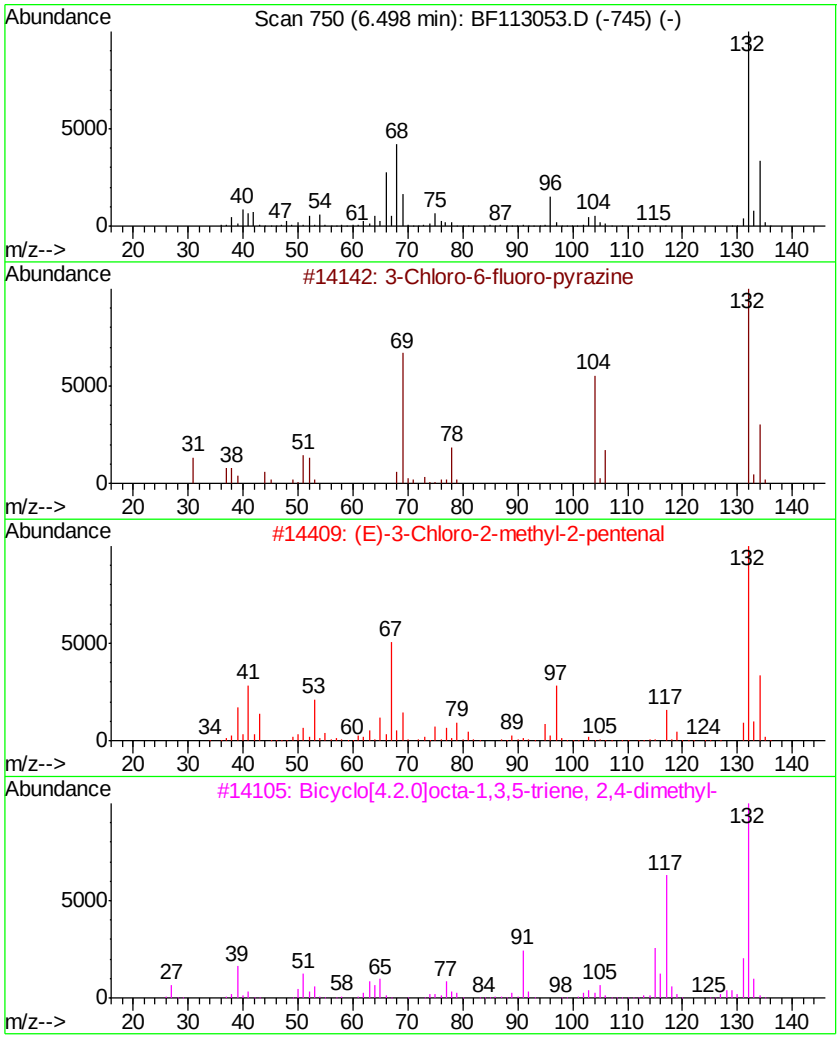
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

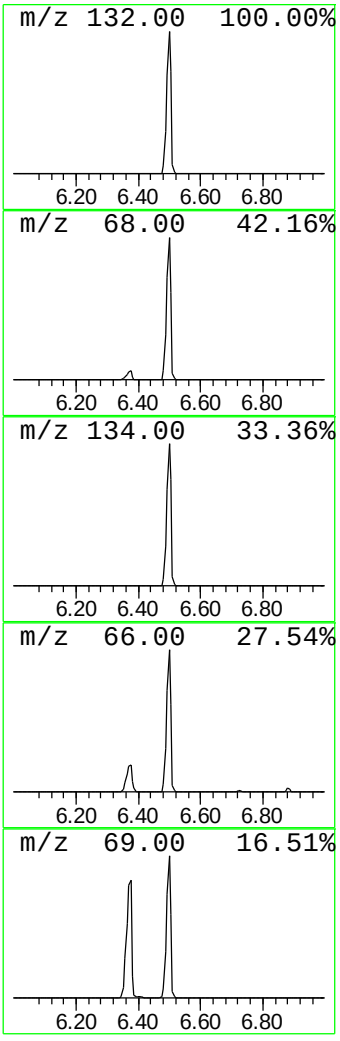
****

Peak Number 1 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	70.44 ng	4300060	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	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5	Xenon	132	Xe	007440-63-3	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
Data File : BF113053.D
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Operator : JU/SJ
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

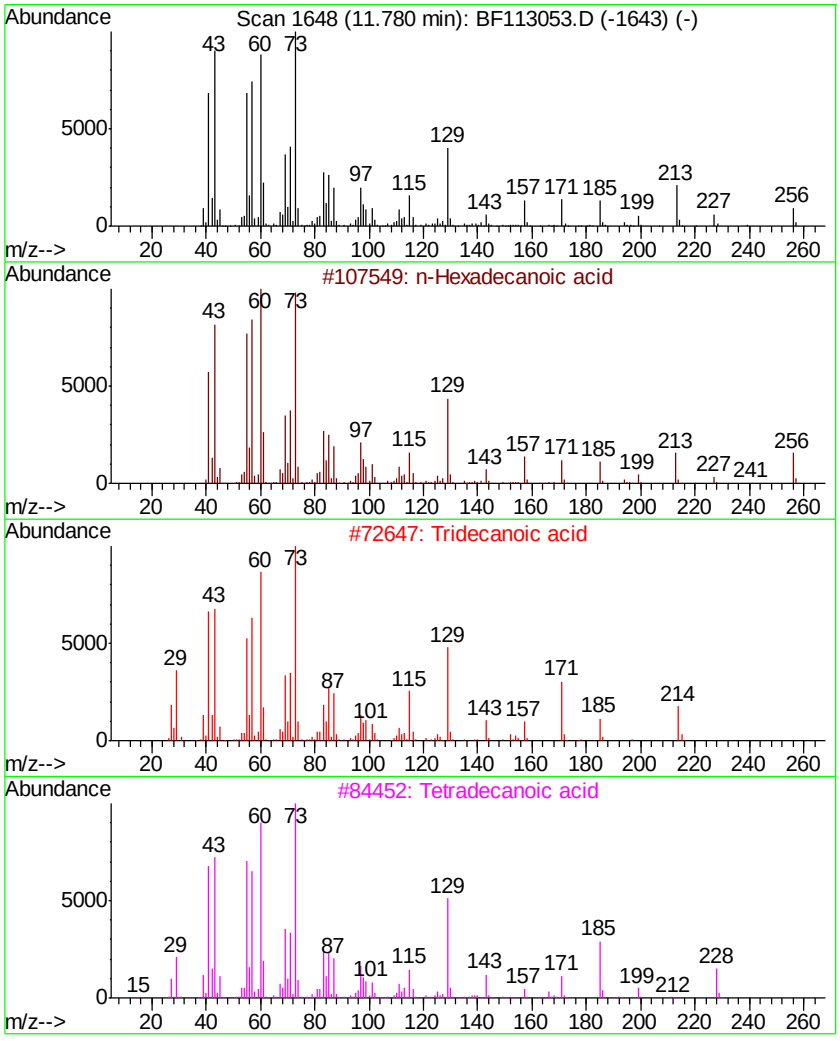
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

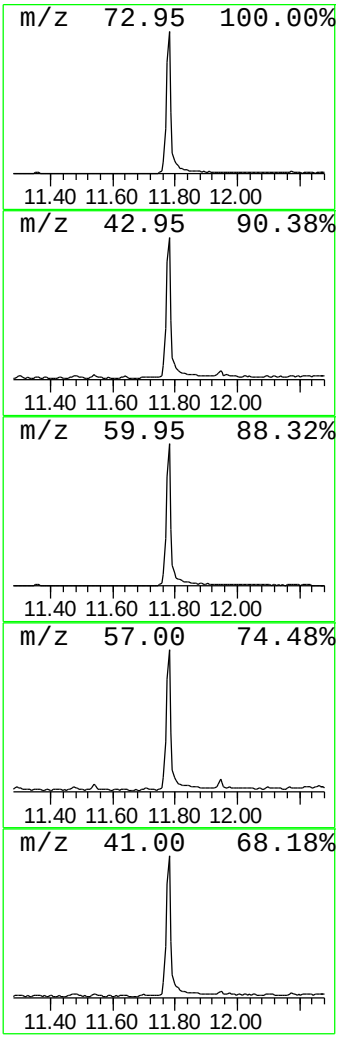
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.78	17.80 ng	1716810	Phenanthrene-d10	11.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	95
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
5	n-Decanoic acid	172	C10H20O2	000334-48-5	76





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
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Sample : K1866-03
Misc :
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Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

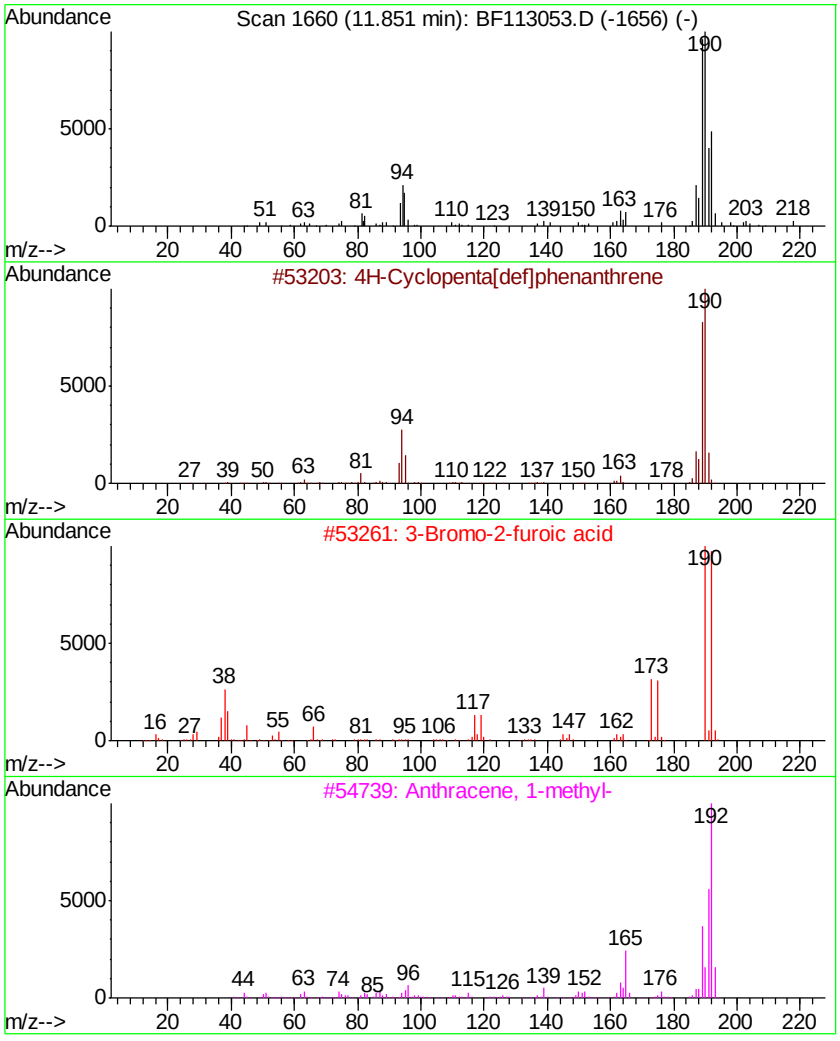
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

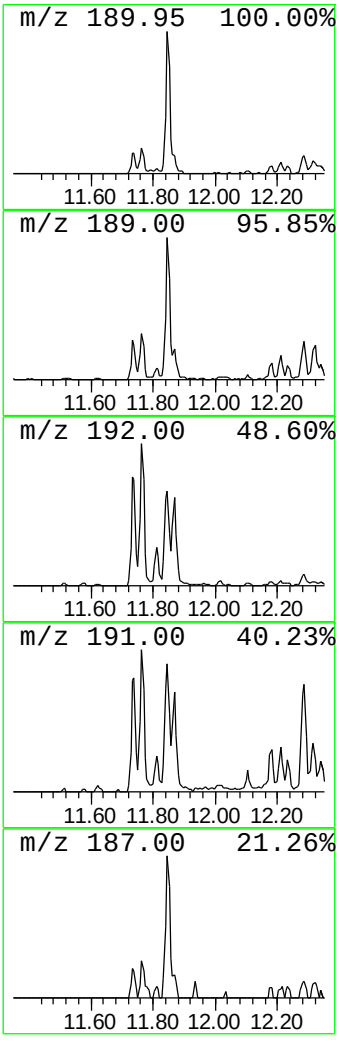
****

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	2.10 ng	203043	Phenanthrene-d10	11.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2	3-Bromo-2-furoic acid	190	C5H3BrO3	014903-90-3	27
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	25
4	1-(3-Nitro-phenyl)-pyrrolidine	192	C10H12N2O2	1000300-34-7	25
5	Benzene, 1,1'-(2-cyclopropen-1-y...	192	C15H12	022825-21-4	22





Library Search Compound Report

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TIC Library : C:\DATABASE\NIST11.L
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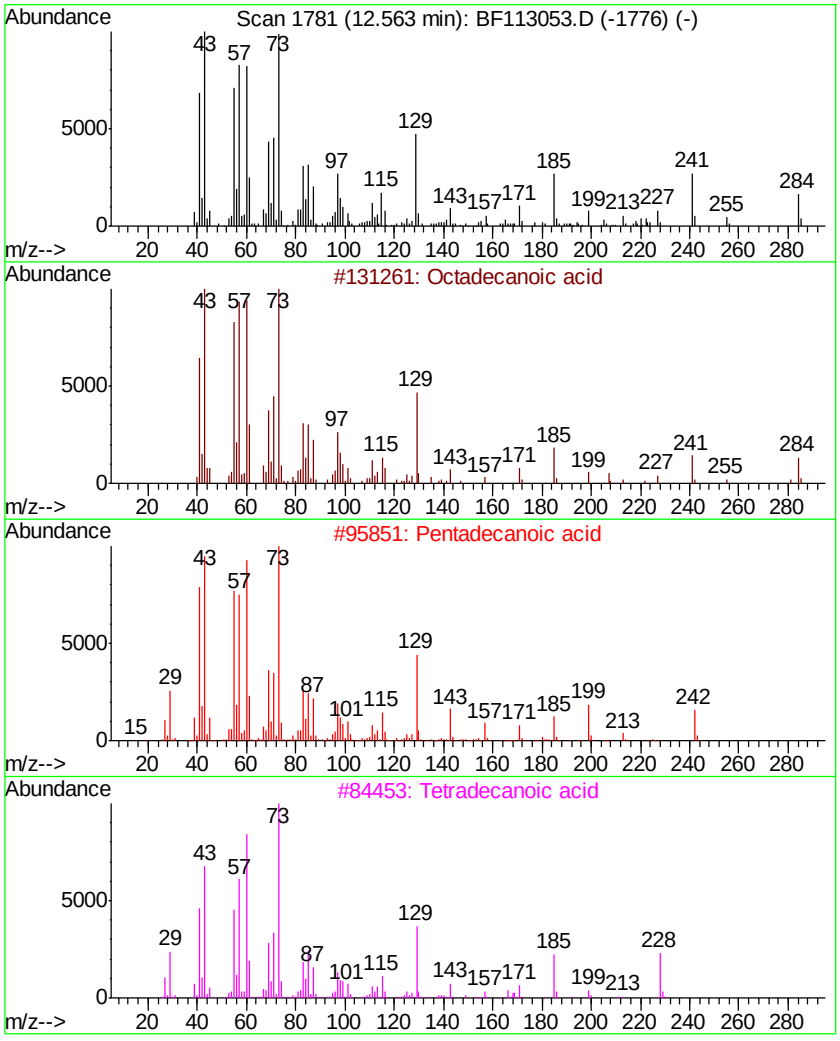
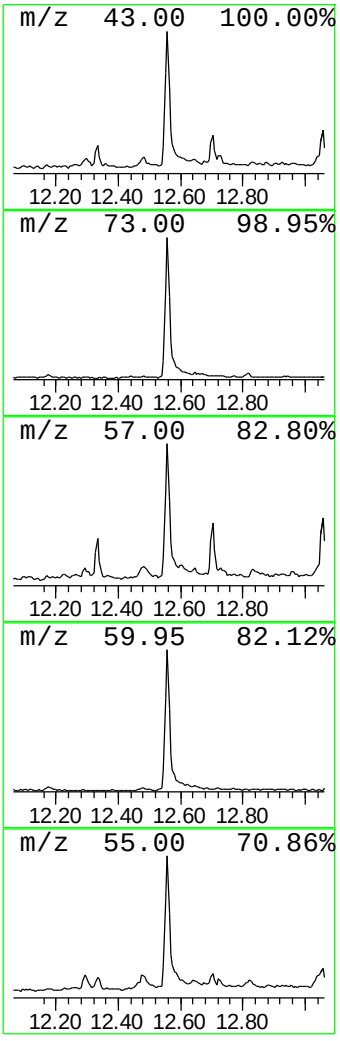
Peak Number 5 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	5.32 ng	708436	Chrysene-d12	13.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74
5	Undecanoic acid	186	C11H22O2	000112-37-8	74





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Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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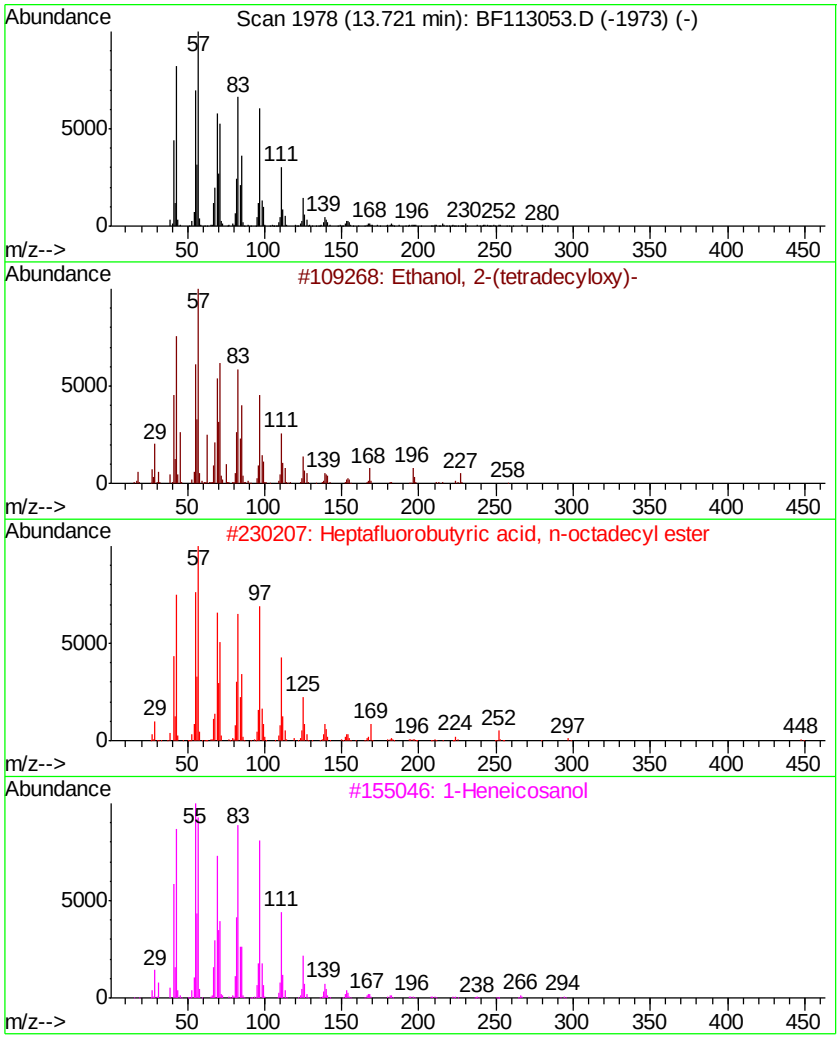
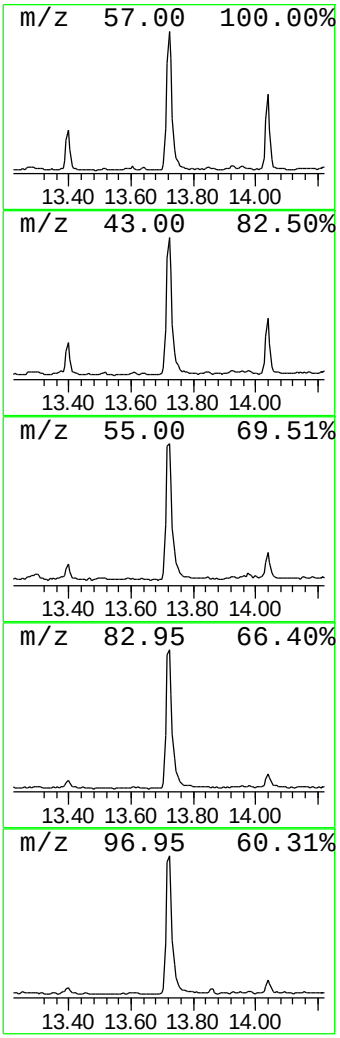
Peak Number 7 Ethanol, 2-(tetradecyloxy)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	9.13 ng	1216770	Chrysene-d12	13.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	96
2	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	93
4	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91





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Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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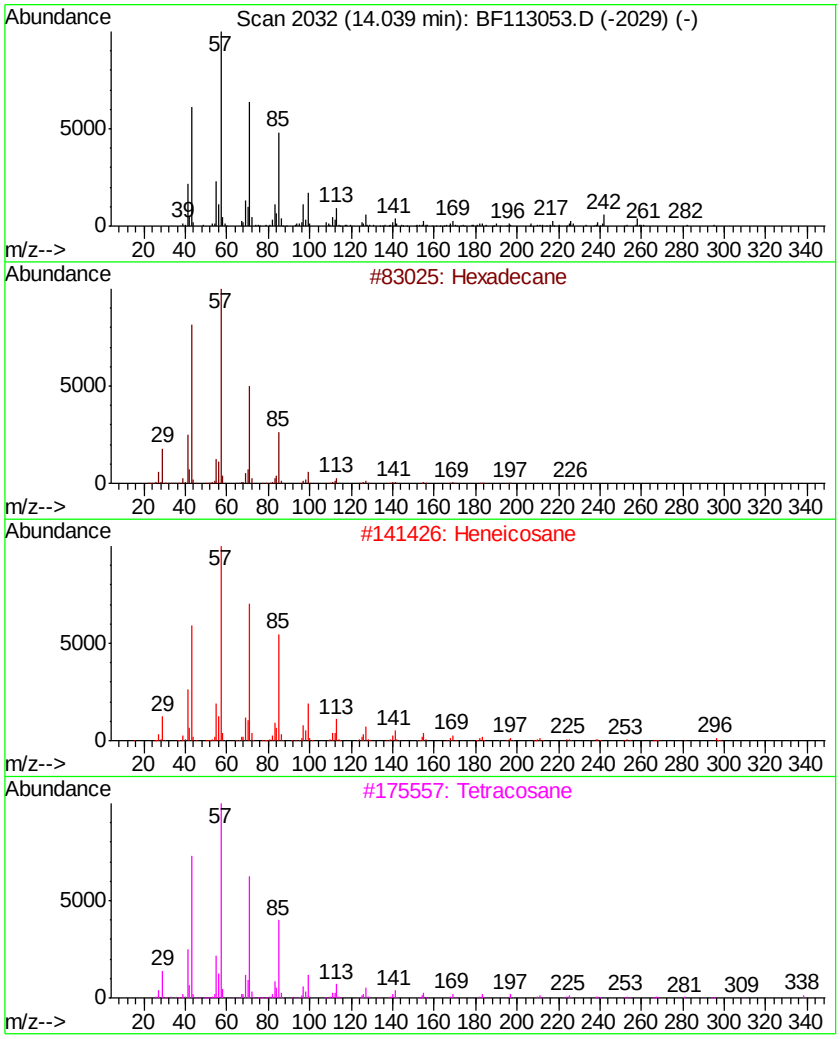
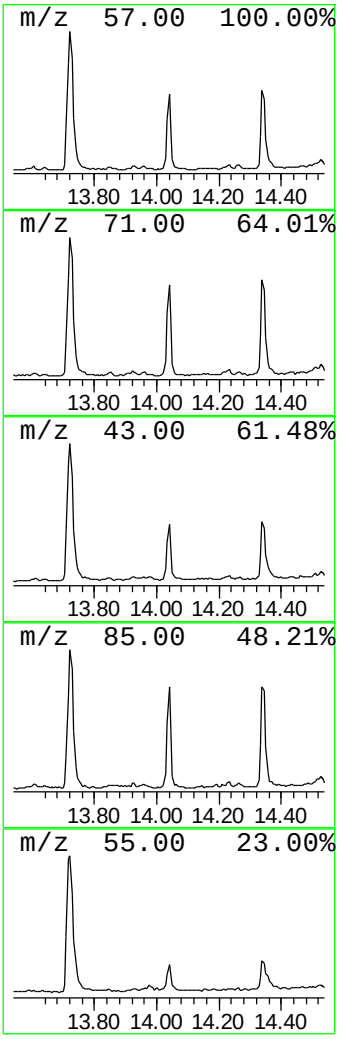
Peak Number 8 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	2.04 ng	272437	Chrysene-d12	13.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	94
2	Heneicosane	296	C21H44	000629-94-7	87
3	Tetracosane	338	C24H50	000646-31-1	87
4	Heptadecane	240	C17H36	000629-78-7	86
5	Eicosane	282	C20H42	000112-95-8	83





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
Data File : BF113053.D
Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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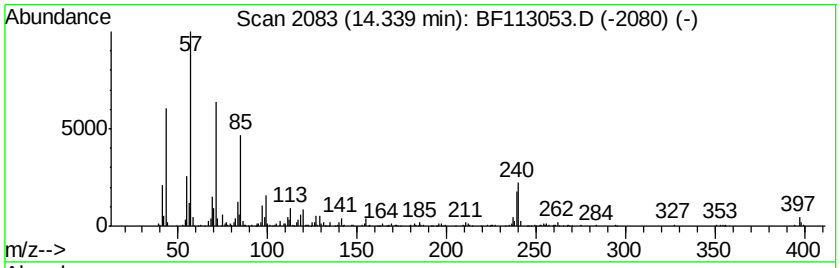
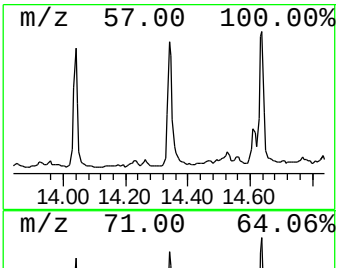
Peak Number 9 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.34	2.45 ng	325981	Chrysene-d12	13.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	94
2	Octadecane	254	C18H38	000593-45-3	86
3	Hexadecane	226	C16H34	000544-76-3	86
4	Heneicosane	296	C21H44	000629-94-7	76
5	Hexacosane	366	C26H54	000630-01-3	76





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031419\
Data File : BF113053.D
Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

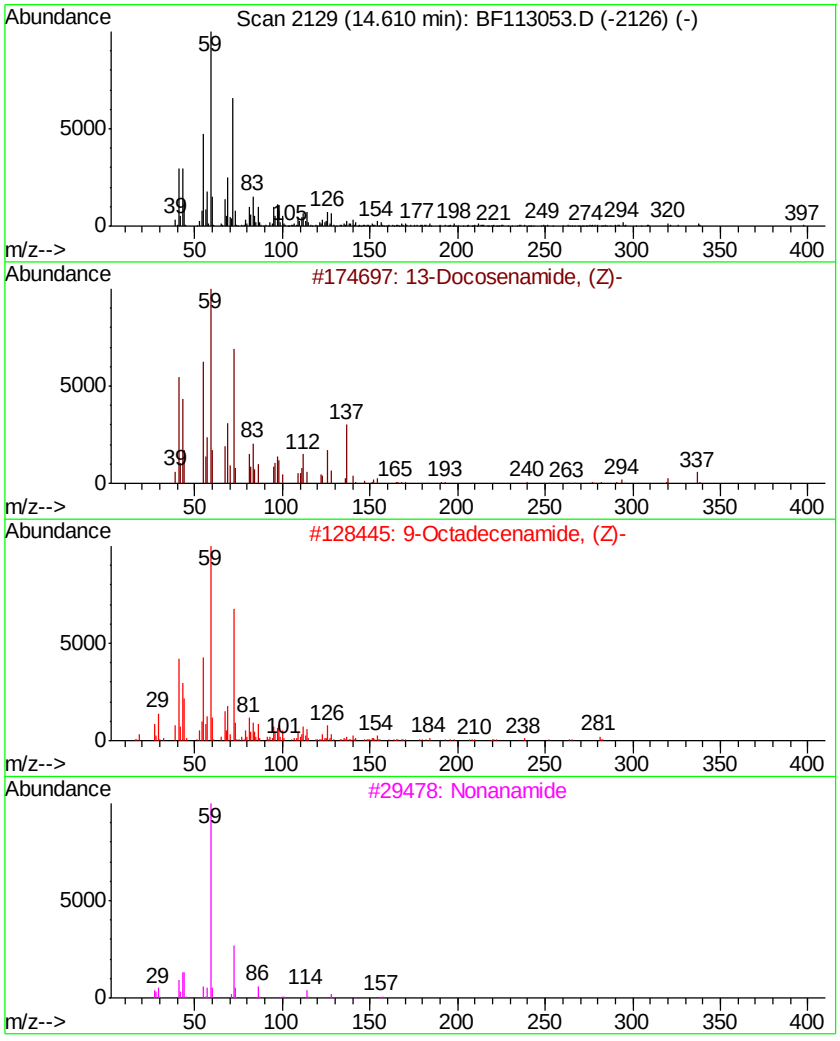
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

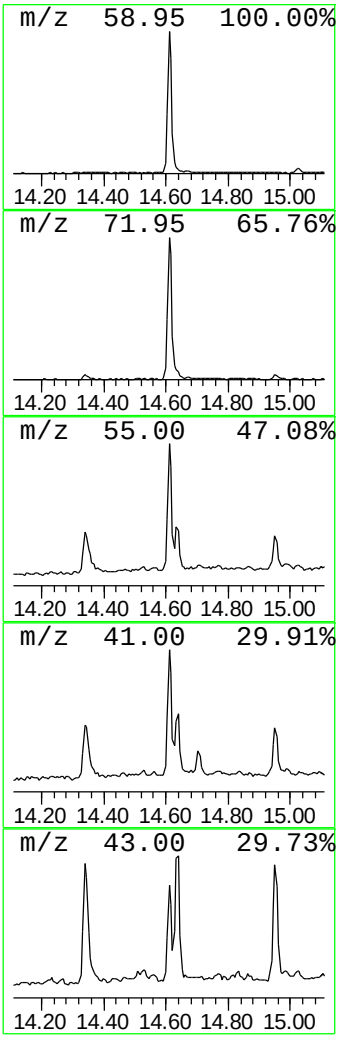
****

Peak Number 10 13-Docosenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	6.71 ng	601923	Perylene-d12	15.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	93
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
3	Nonanamide	157	C9H19NO	001120-07-6	53
4	Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	50
5	Butanamide, 3-methyl-	101	C5H11NO	000541-46-8	38





Library Search Compound Report

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Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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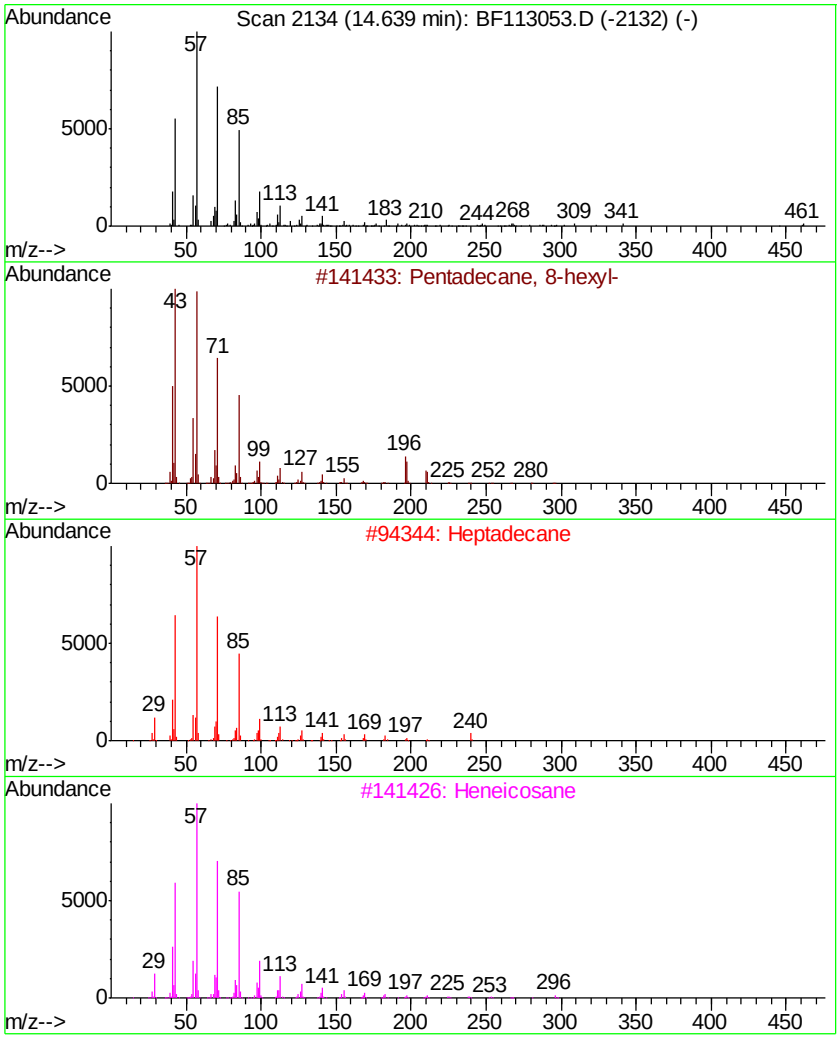
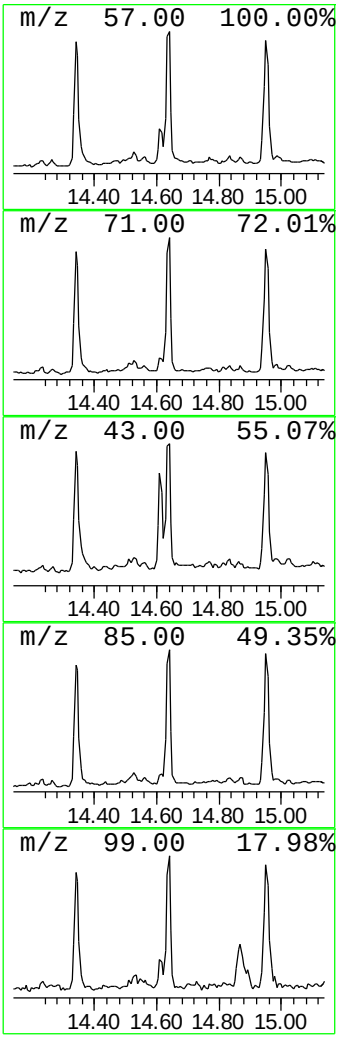
Peak Number 11 Pentadecane, 8-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	3.62 ng	324808	Perylene-d12	15.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
2	Heptadecane	240	C17H36	000629-78-7	87
3	Heneicosane	296	C21H44	000629-94-7	87
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
5	5-Butyl-5-ethylheptadecane	324	C23H48	1000360-43-0	78





Library Search Compound Report

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Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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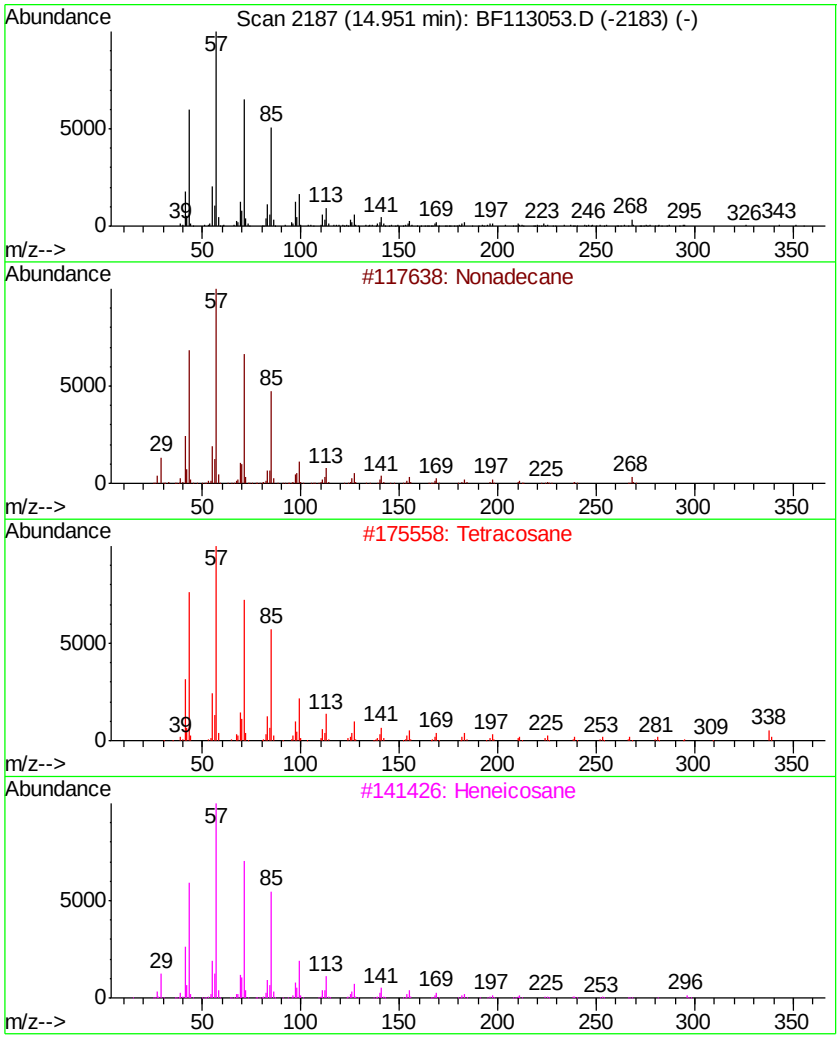
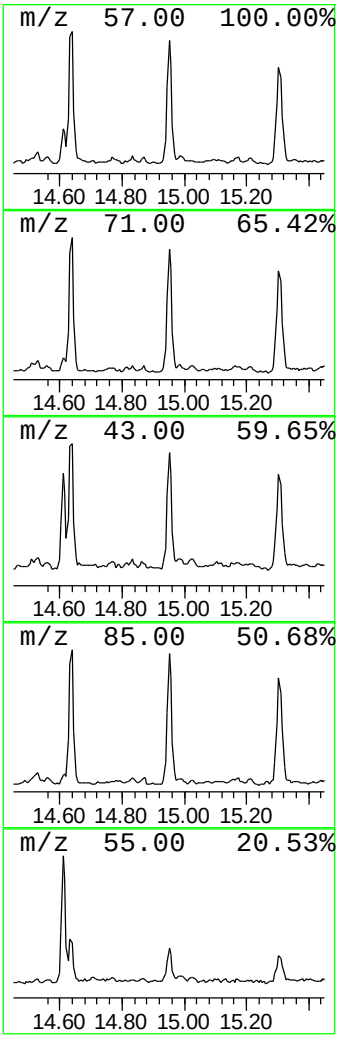
Peak Number 12 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	4.82 ng	432198	Perylene-d12	15.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	97
2	Tetracosane	338	C24H50	000646-31-1	91
3	Heneicosane	296	C21H44	000629-94-7	87
4	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	86
5	Pentacosane	352	C25H52	000629-99-2	83





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Data File : BF113053.D
Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.M
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TIC Integration Parameters: LSCINT.P

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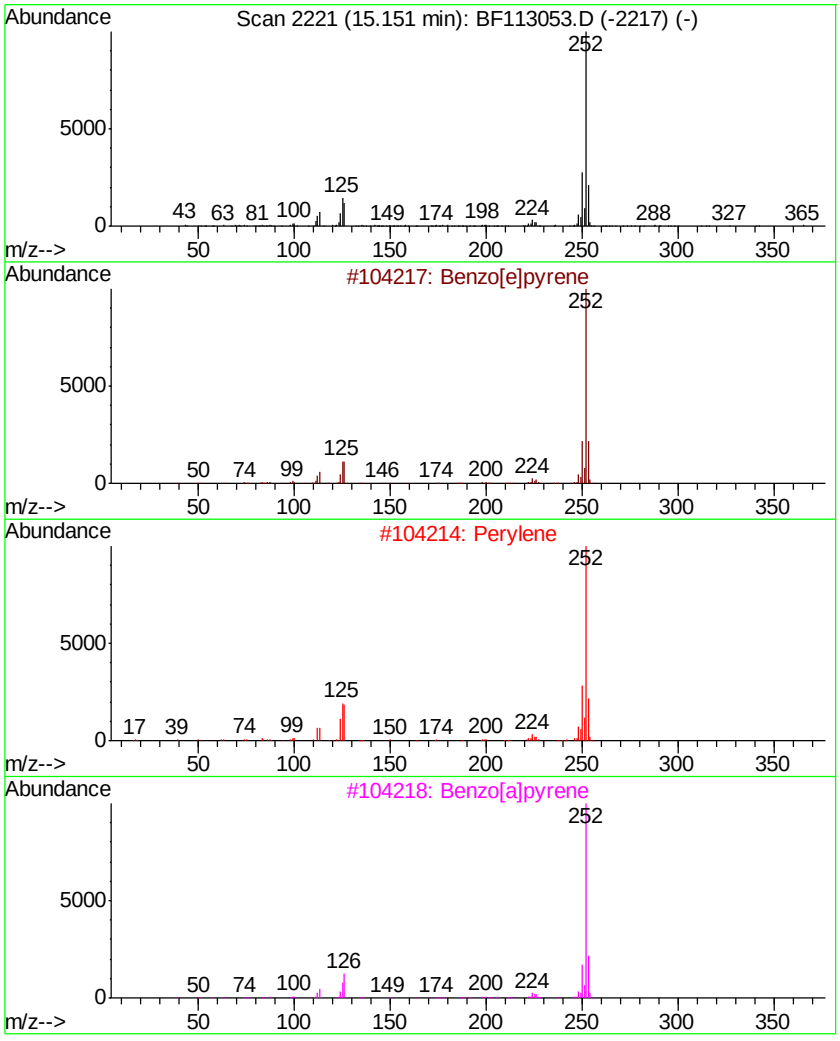
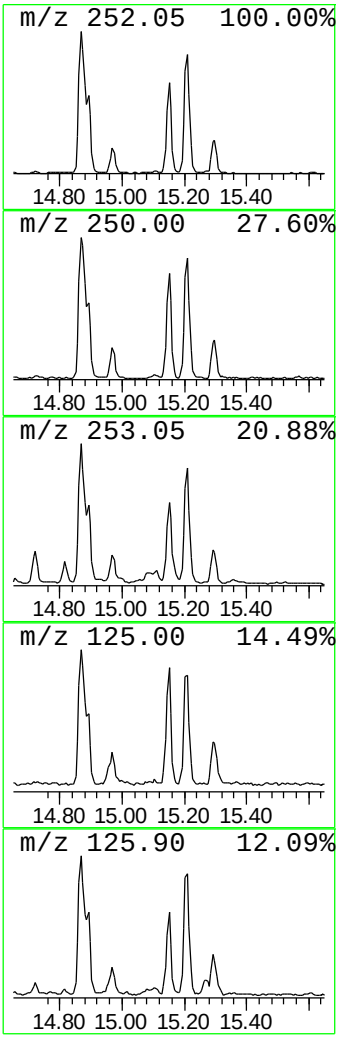
Peak Number 13 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.15	4.17 ng	373752	Perylene-d12	15.27

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





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Data File : BF113053.D
Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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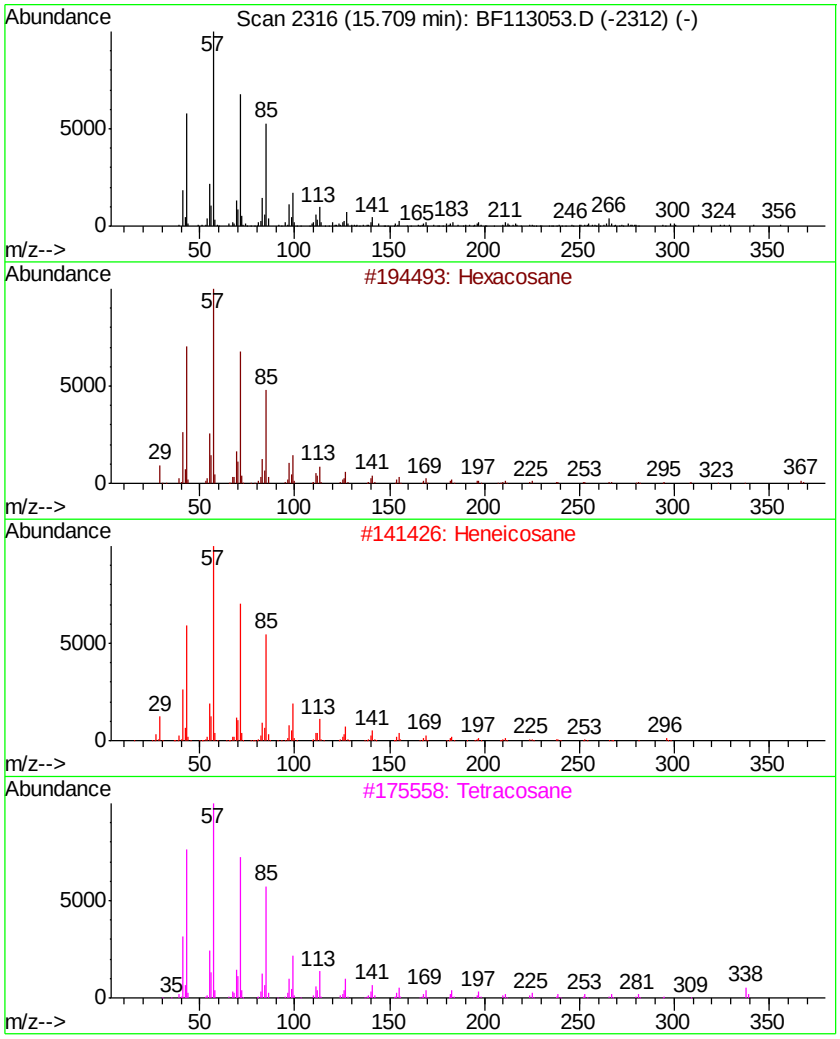
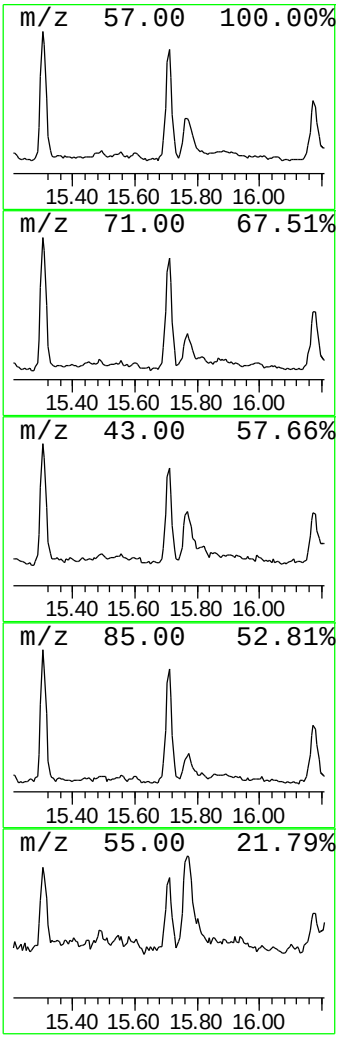
Peak Number 14 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.71	2.76 ng	247062	Perylene-d12	15.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	90
2	Heneicosane	296	C21H44	000629-94-7	90
3	Tetracosane	338	C24H50	000646-31-1	87
4	Octacosane	394	C28H58	000630-02-4	87
5	Nonadecane	268	C19H40	000629-92-5	87





Library Search Compound Report

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Misc :
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Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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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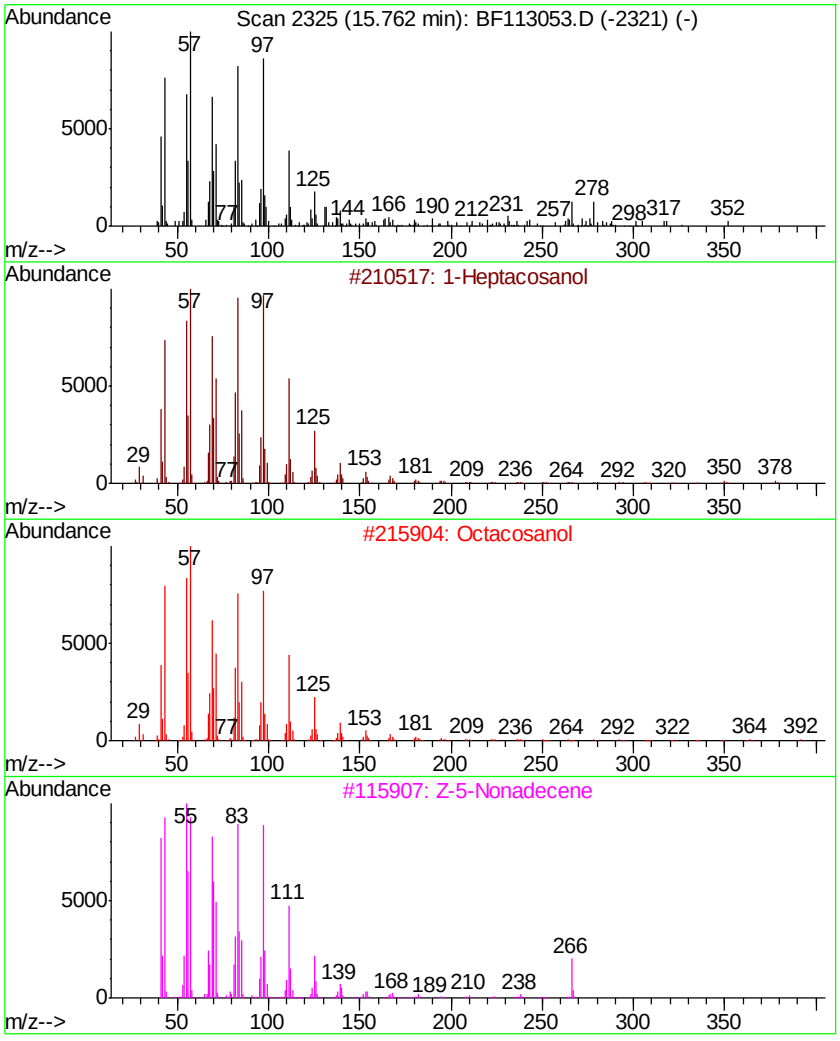
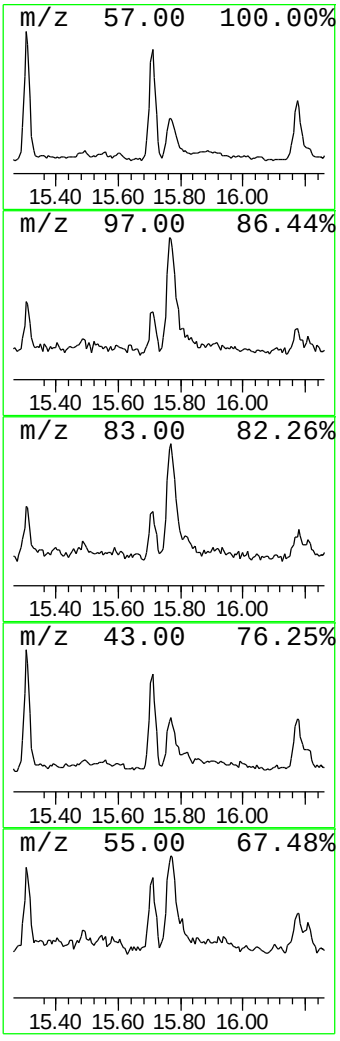
Peak Number 15 1-Heptacosanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	4.42 ng	396095	Perylene-d12	15.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	93
2	Octacosanol	410	C28H58O	000557-61-9	93
3	Z-5-Nonadecene	266	C19H38	1000131-11-8	91
4	1-Nonadecene	266	C19H38	018435-45-5	91
5	1-Docosene	308	C22H44	001599-67-3	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF031419\
Data File : BF113053.D
Acq On : 14 Mar 2019 16:50
Operator : JU/SJ
Sample : K1866-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MR-MH-03-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022719.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
unknown6.50	6.50	70.4	ng	4300060	1	6.73	1220900	20.0
n-Hexadecanoic acid	11.78	17.8	ng	1716810	4	11.23	1929440	20.0
4H-Cyclopenta[def...	11.85	2.1	ng	203043	4	11.23	1929440	20.0
Octadecanoic acid	12.56	5.3	ng	708436	5	13.86	2664690	20.0
Ethanol, 2-(tetra...	13.72	9.1	ng	1216770	5	13.86	2664690	20.0
Hexadecane	14.04	2.0	ng	272437	5	13.86	2664690	20.0
Octacosane	14.34	2.5	ng	325981	5	13.86	2664690	20.0
13-Docosenamide, ...	14.61	6.7	ng	601923	6	15.27	1793100	20.0
Pentadecane, 8-he...	14.64	3.6	ng	324808	6	15.27	1793100	20.0
Nonadecane	14.95	4.8	ng	432198	6	15.27	1793100	20.0
Benzo[e]pyrene	15.15	4.2	ng	373752	6	15.27	1793100	20.0
Hexacosane	15.71	2.8	ng	247062	6	15.27	1793100	20.0
1-Heptacosanol	15.76	4.4	ng	396095	6	15.27	1793100	20.0