LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110540.D
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

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-BF102318.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	2.234	22	25	33	rVB	53209	83052	9.15%	1.047%
2	5.575	590	593	607	rBV	163022	200964	22.14%	2.534%
3	6.581	760	764	784	rBV	148429	248311	27.35%	3.132%
4	6.728	785	789	796	rBV	222331	225712	24.87%	2.847%
5	6.957	824	828	838	rBV	428432	401159	44.19%	5.059%
6	7.110	851	854	865	rBV	182352	176381	19.43%	2.224%
7	7.522	921	924	937	rBV	68895	106443	11.73%	1.342%
8	8.239	1042	1046	1069	rVB	431911	525829	57.93%	6.631%
9	8.810	1140	1143	1149	rBV	13600	13120	1.45%	0.165%
10	8.957	1163	1168	1175	rBV	15985	26732	2.94%	0.337%
11	9.057	1182	1185	1189	rBV	24115	24983	2.75%	0.315%
12	9.310	1225	1228	1237	rBV	186178	215878	23.78%	2.723%
13	9.375	1237	1239	1243	rVB	19652	19715	2.17%	0.249%
14	9.586	1269	1275	1279	rBV2	9824	14726	1.62%	0.186%
15	9.657	1282	1287	1289	rVV	17322	19691	2.17%	0.248%
16	9.704	1289	1295	1302	rVB3	22820	42396	4.67%	0.535%
17	9.863	1318	1322	1326	rBV2	19330	20857	2.30%	0.263%
18	9.910	1326	1330	1334	rVB	22997	23460	2.58%	0.296%
19	9.998	1342	1345	1349	rBV	496791	445948	49.13%	5.624%
20	10.033	1349	1351	1359	rVB	36009	38018	4.19%	0.479%
21	10.204	1375	1380	1385	rBV3	12281	27229	3.00%	0.343%
22	10.404	1411	1414	1417	rBV	27197	23907	2.63%	0.302%
23	10.551	1436	1439	1445	rVB	30948	34195	3.77%	0.431%
24	10.627	1445	1452	1455	rBV3	13676	22100	2.43%	0.279%
25	10.792	1476	1480	1488	rBV	95642	111826	12.32%	1.410%
26	10.880	1492	1495	1500	rBV	36723	42279	4.66%	0.533%
27	11.098	1524	1532	1535	rBV6	8960	17054	1.88%	0.215%
28	11.327	1568	1571	1574	rBV	31476	25614	2.82%	0.323%
29	11.392	1579	1582	1585	rVB	11072	10615	1.17%	0.134%
30	11.492	1595	1599	1601	rBV	477233	428269	47.18%	5.401%
31	11.516	1601	1603	1609	rVB	212663	192871	21.25%	2.432%
32	11.568	1609	1612	1616	rBV	43468	45623	5.03%	0.575%
33	11.757	1641	1644	1646	rVB	29680	26593	2.93%	0.335%
34	11.857	1658	1661	1665	rVB	333268	255948	28.20%	3.228%

LSC Area Percent Report

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Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

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-BF102318.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.992	1682	1684	1687	rBV	36844	40156	4.42%	0.506%
36	12.021	1687	1689	1692	rVV	39640	37176	4.10%	0.469%
37	12.074	1695	1698	1700	rVV2	16367	15426	1.70%	0.195%
38	12.110	1700	1704	1706	rVV2	43225	54950	6.05%	0.693%
39	12.133	1706	1708	1710	rVV	22603	20101	2.21%	0.254%
40	12.163	1710	1713	1717	rVB	33434	26400	2.91%	0.333%
41	12.286	1732	1734	1737	rBV	14431	14262	1.57%	0.180%
42	12.321	1738	1740	1744	rVB4	10597	12389	1.36%	0.156%
43	12.439	1758	1760	1763	rVB2	13613	11348	1.25%	0.143%
44	12.468	1763	1765	1767	rBV3	14132	12375	1.36%	0.156%
45	12.545	1774	1778	1780	rBV	1049178	907736	100.00%	11.448%
46	12.568	1780	1782	1790	rVV2	674486	646644	71.24%	8.155%
47	12.651	1793	1796	1802	rVB	141166	130073	14.33%	1.640%
48	12.710	1802	1806	1810	rBV	239810	224885	24.77%	2.836%
49	12.863	1830	1832	1834	rBV2	13482	11875	1.31%	0.150%
50	12.886	1834	1836	1839	rVB2	42364	34616	3.81%	0.437%
51	12.921	1839	1842	1843	rBV	78557	63914	7.04%	0.806%
52	12.939	1843	1845	1851	rVV	195121	215126	23.70%	2.713%
53	12.992	1851	1854	1857	rVB	22657	23530	2.59%	0.297%
54	13.068	1864	1867	1870	rBV	220650	195521	21.54%	2.466%
55	13.145	1877	1880	1881	rBV2	14275	14142	1.56%	0.178%
56	13.210	1889	1891	1894	rBV3	23612	26263	2.89%	0.331%
57	13.251	1894	1898	1899	rVV3	36509	46663	5.14%	0.588%
58	13.274	1900	1902	1909	rVB2	51665	65599	7.23%	0.827%
59	13.374	1916	1919	1921	rBV2	32586	28661	3.16%	0.361%
60	13.468	1932	1935	1937	rBV2	21884	25319	2.79%	0.319%
61	13.621	1958	1961	1963	rBV	53139	45245	4.98%	0.571%
62	13.945	2014	2016	2020	rBV3	68807	69440	7.65%	0.876%
63	14.139	2045	2049	2052	rBV2	397524	448708	49.43%	5.659%
64	14.262	2068	2070	2073	rVB	64422	51123	5.63%	0.645%
65	14.568	2119	2122	2128	rBV2	66093	74424	8.20%	0.939%
66	15.668	2306	2309	2314	rVB2	155659	197749	21.78%	2.494%

Sum of corrected areas: 7929337

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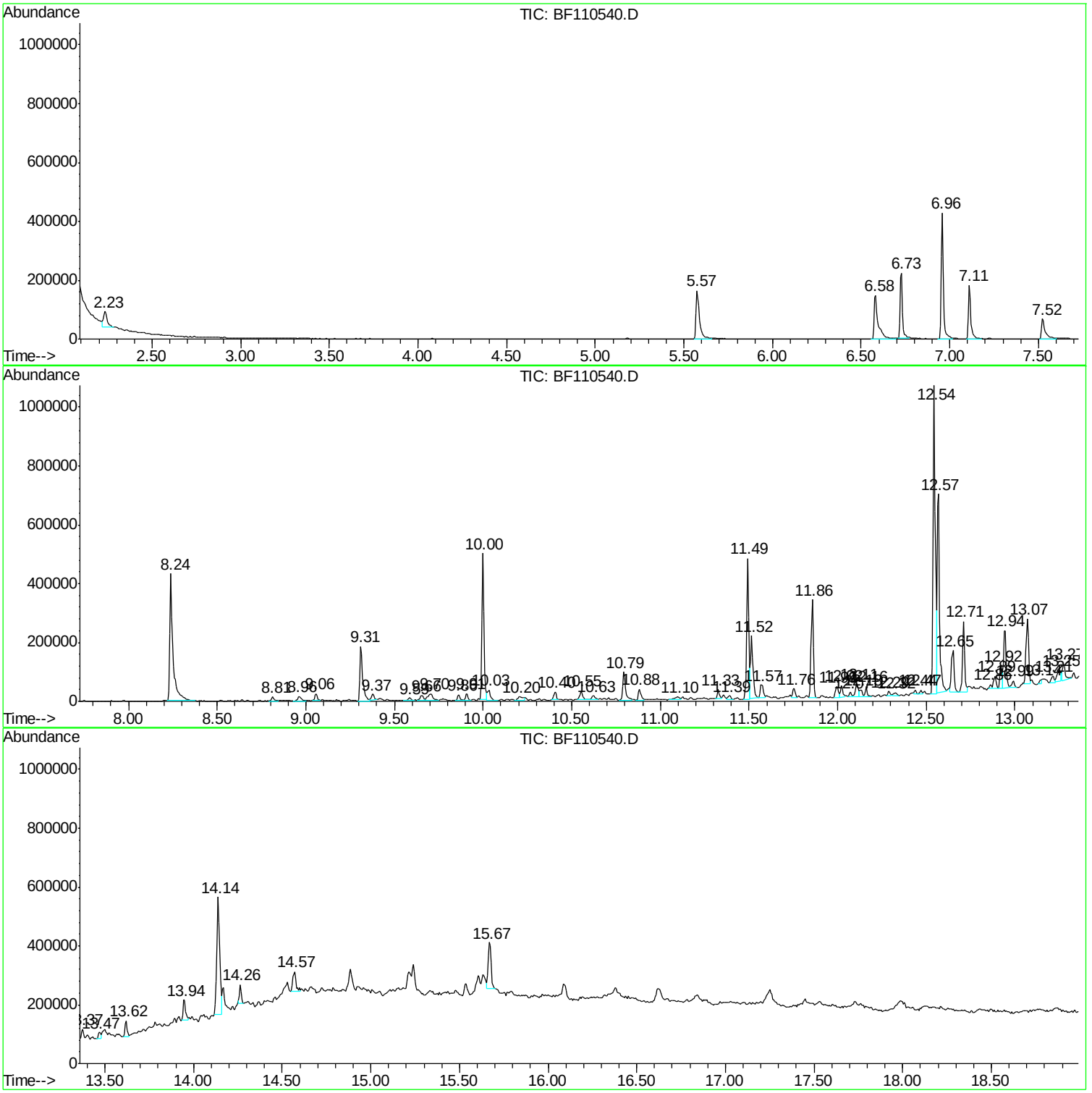
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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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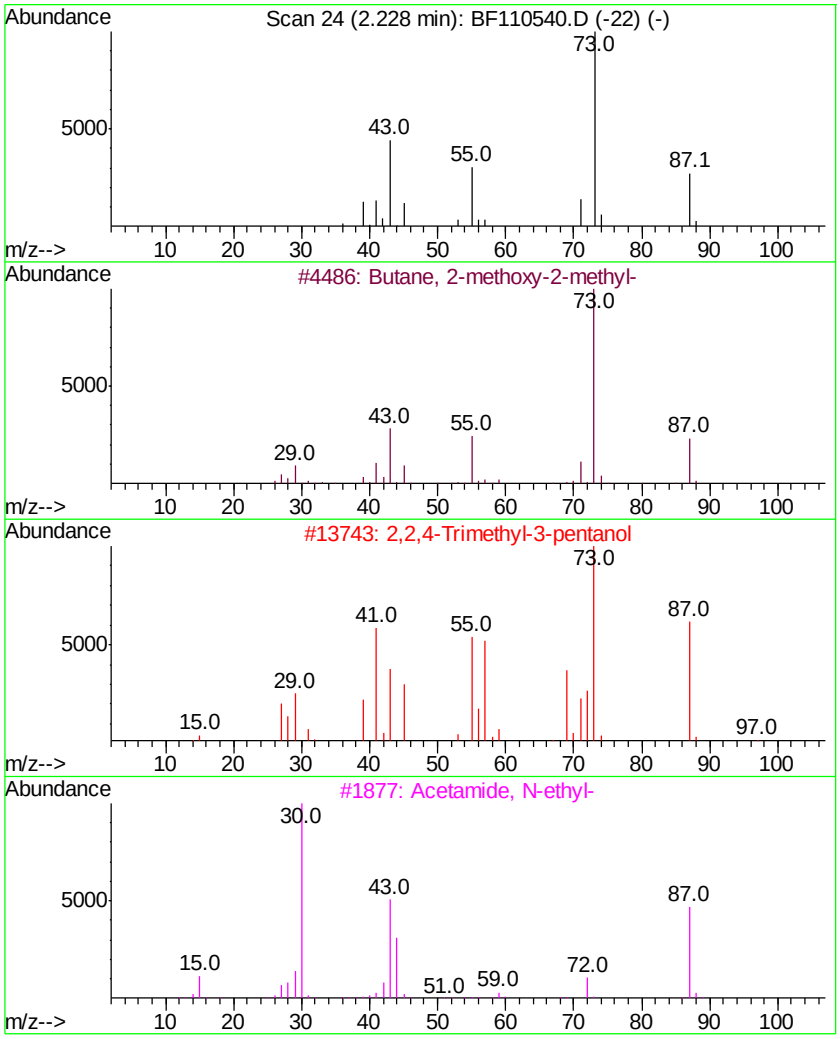
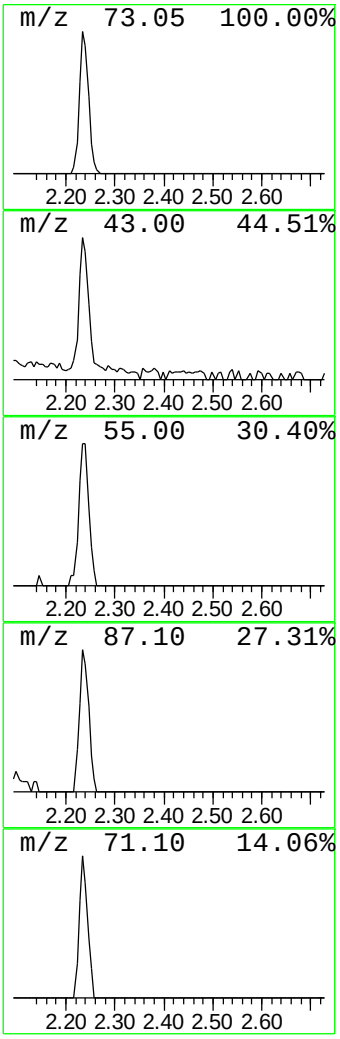
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	4.14 ng	83052	1,4-Dichlorobenzene-d4	6.96

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	39
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	25
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110540.D
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

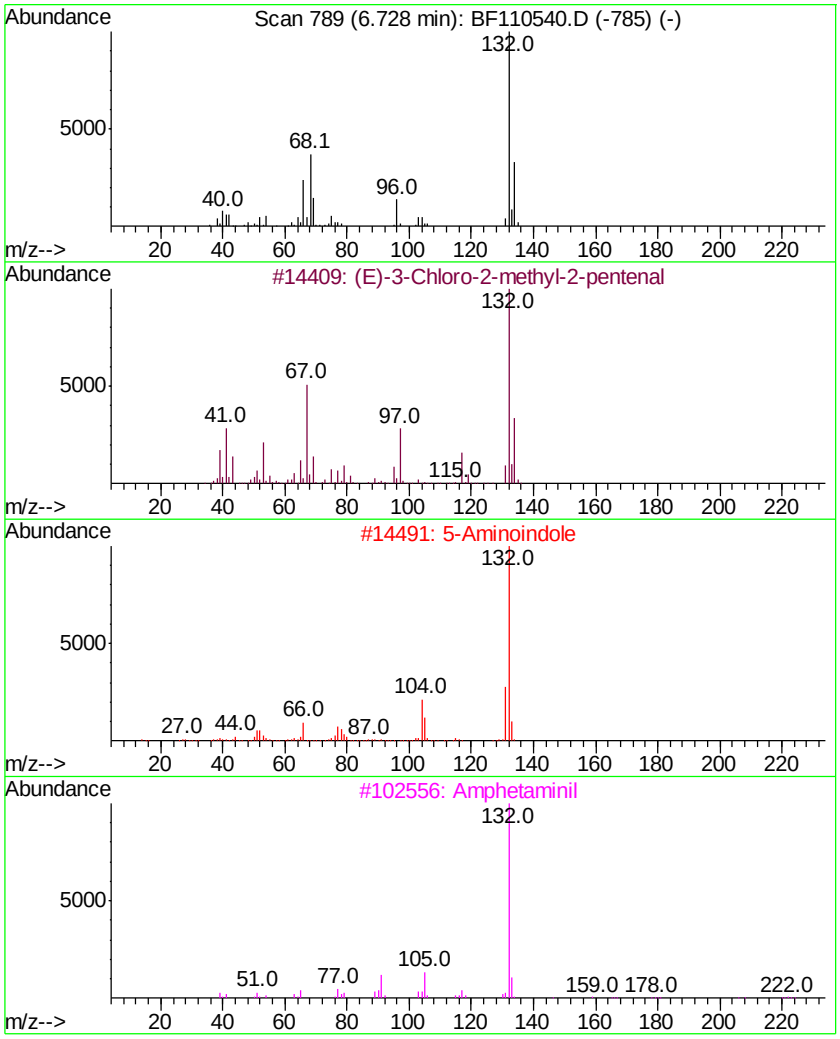
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

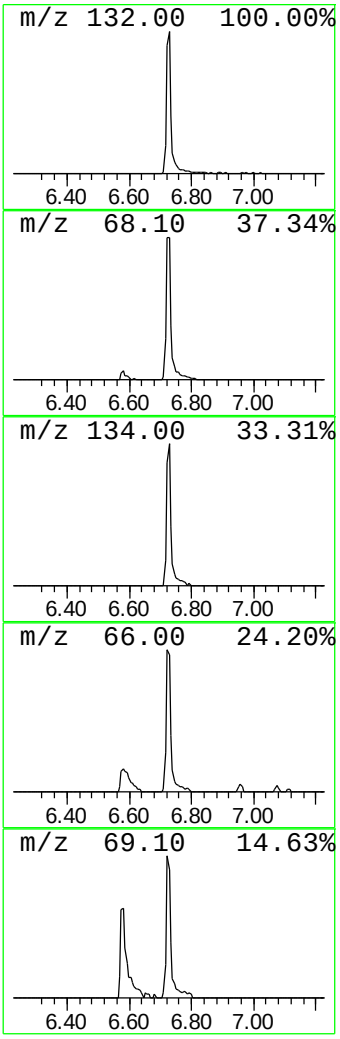
****

Peak Number 2 unknown6.73 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	11.25 ng	225712	1,4-Dichlorobenzene-d4	6.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2	5-Aminoindole	132	C8H8N2	005192-03-0	16
3	Amphetaminil	250	C17H18N2	017590-01-1	12
4	N-(-phenylisopropyl)benzaldimine	223	C16H17N	1000379-01-3	12
5	N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	12





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
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Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

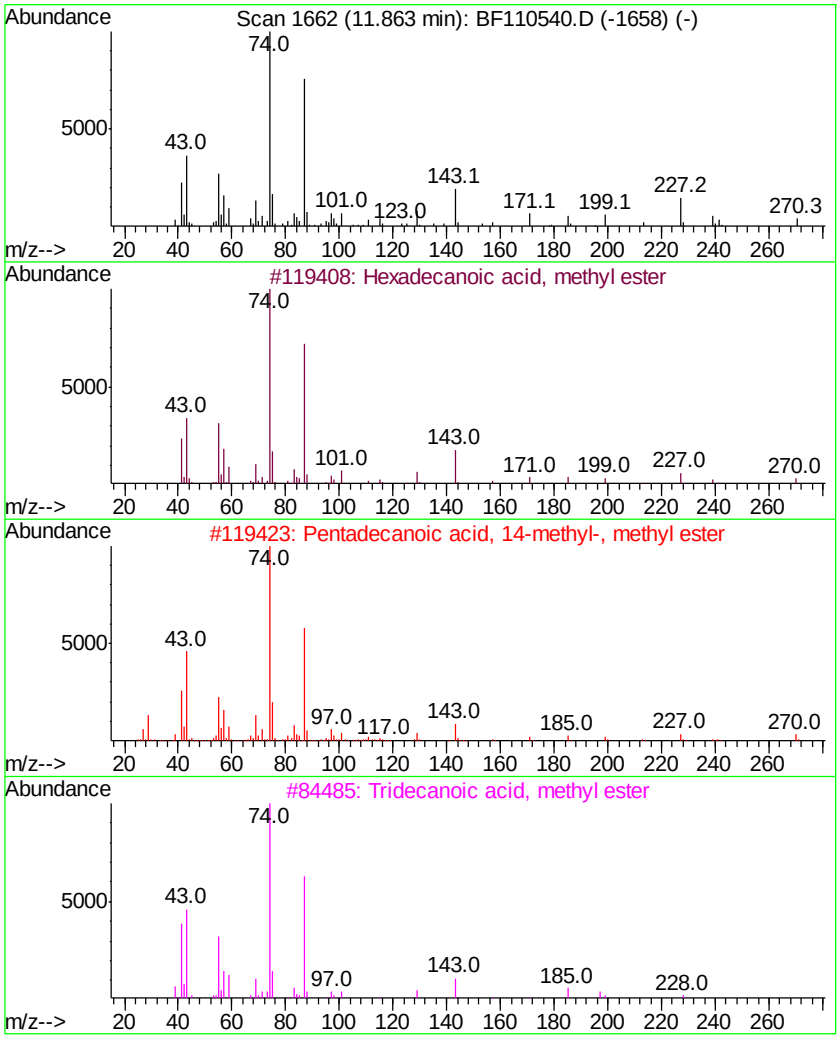
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

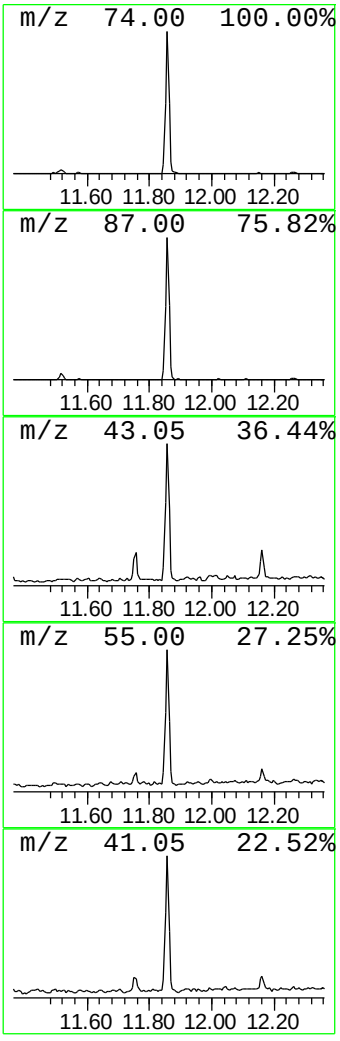
****

Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	11.95 ng	255948	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4	Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
5	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86





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Data File : BF110540.D
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

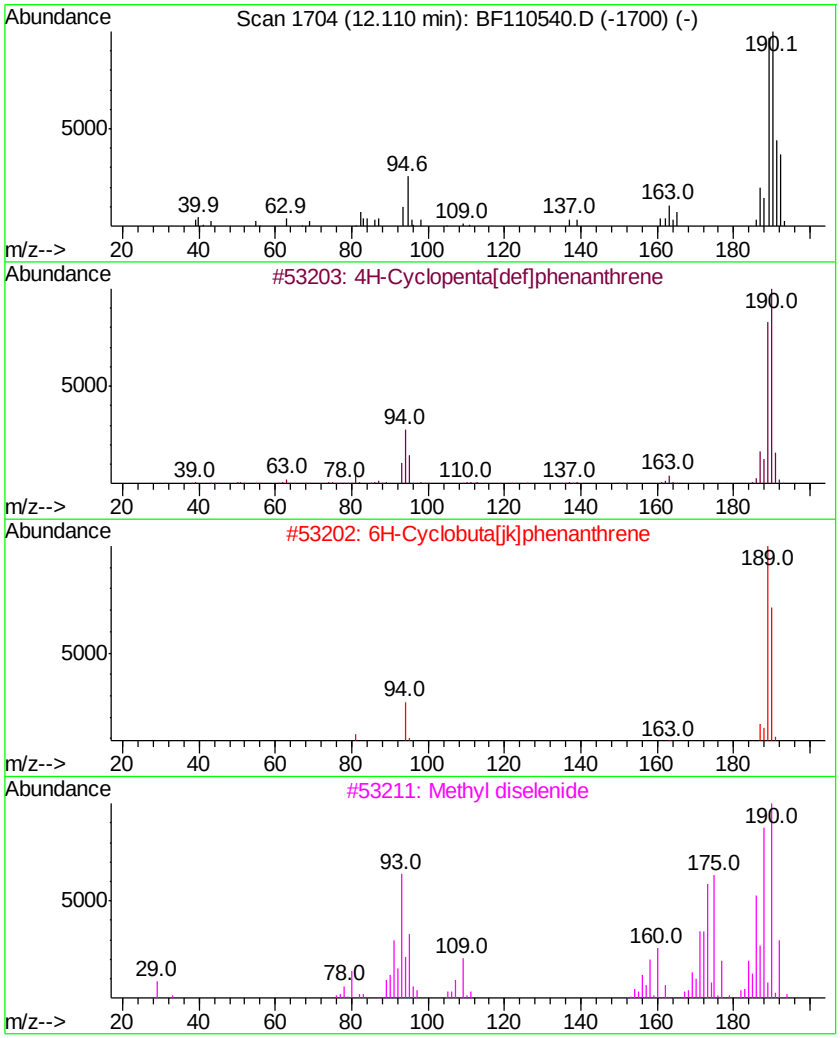
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

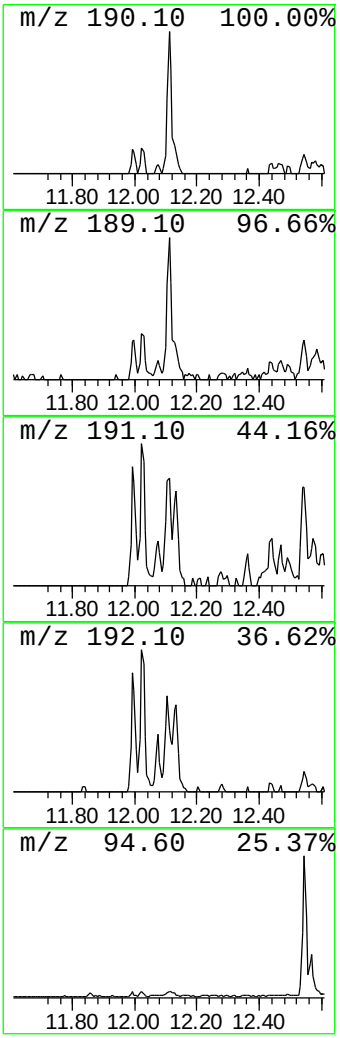
****

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	2.57 ng	54950	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	74
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3	Methyl diselenide	190	C2H6Se2	007101-31-7	38
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	33
5	5-Thiazolecarboxamide, 4-amino-2...	189	C5H7N3OS2	1000338-08-9	27





Library Search Compound Report

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Operator : JU/SJ
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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

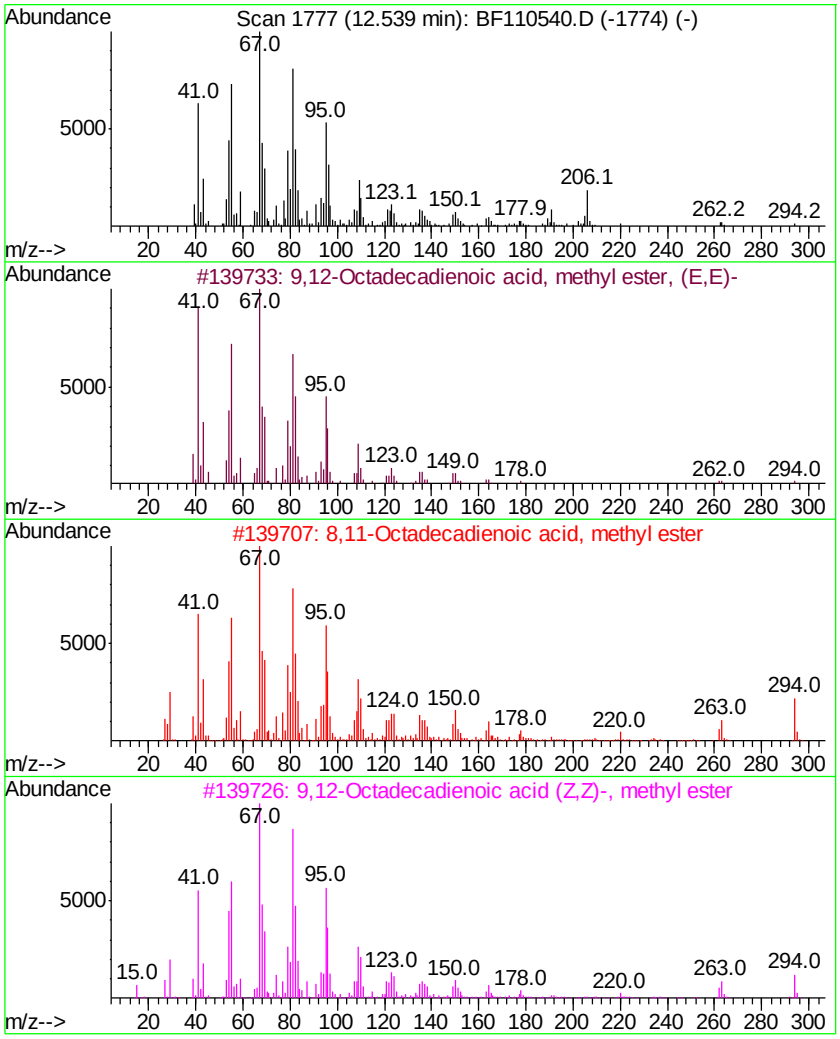
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

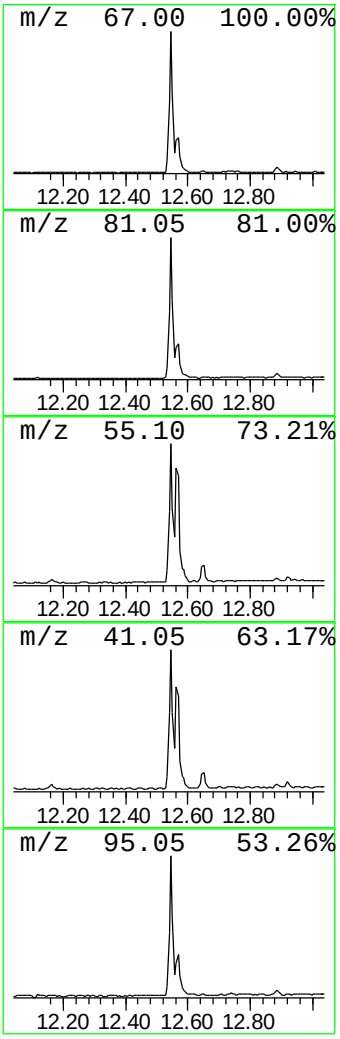
****

Peak Number 6 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	42.39 ng	907736	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

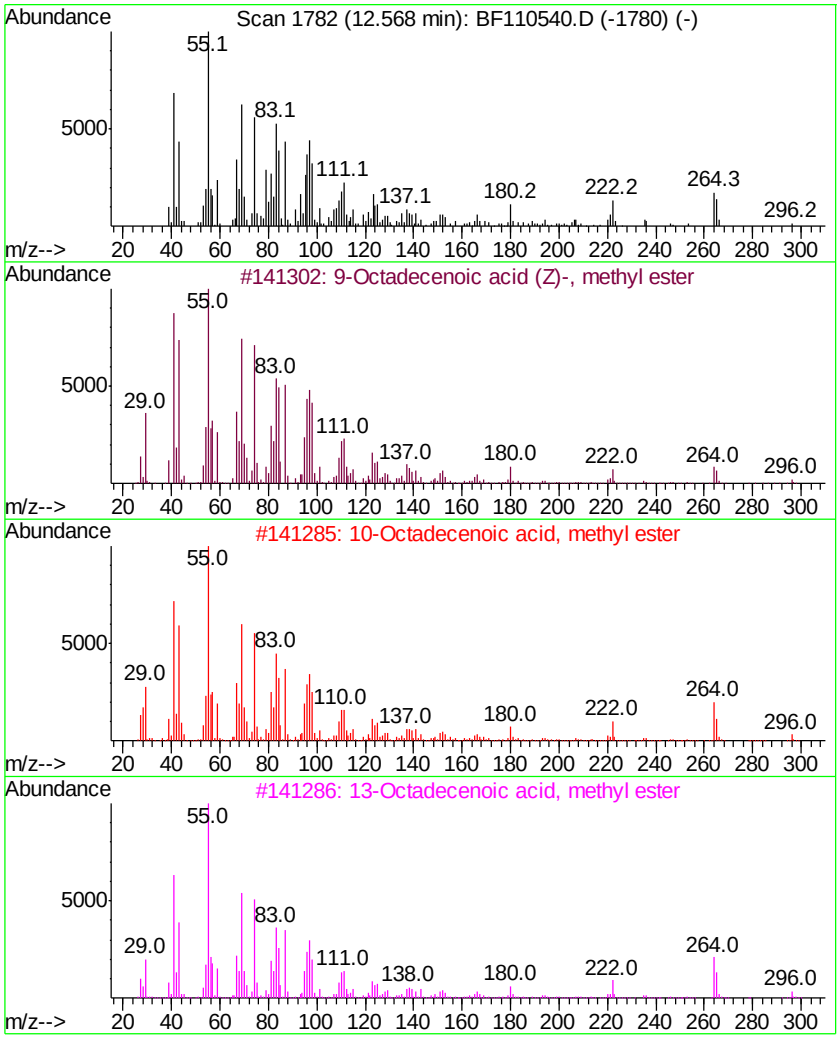
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

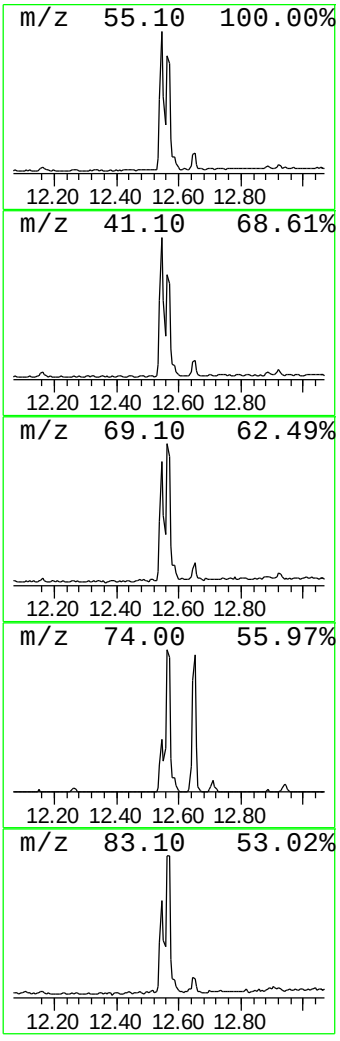
****

Peak Number 7 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	30.20 ng	646644	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
4	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99
5	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110540.D
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

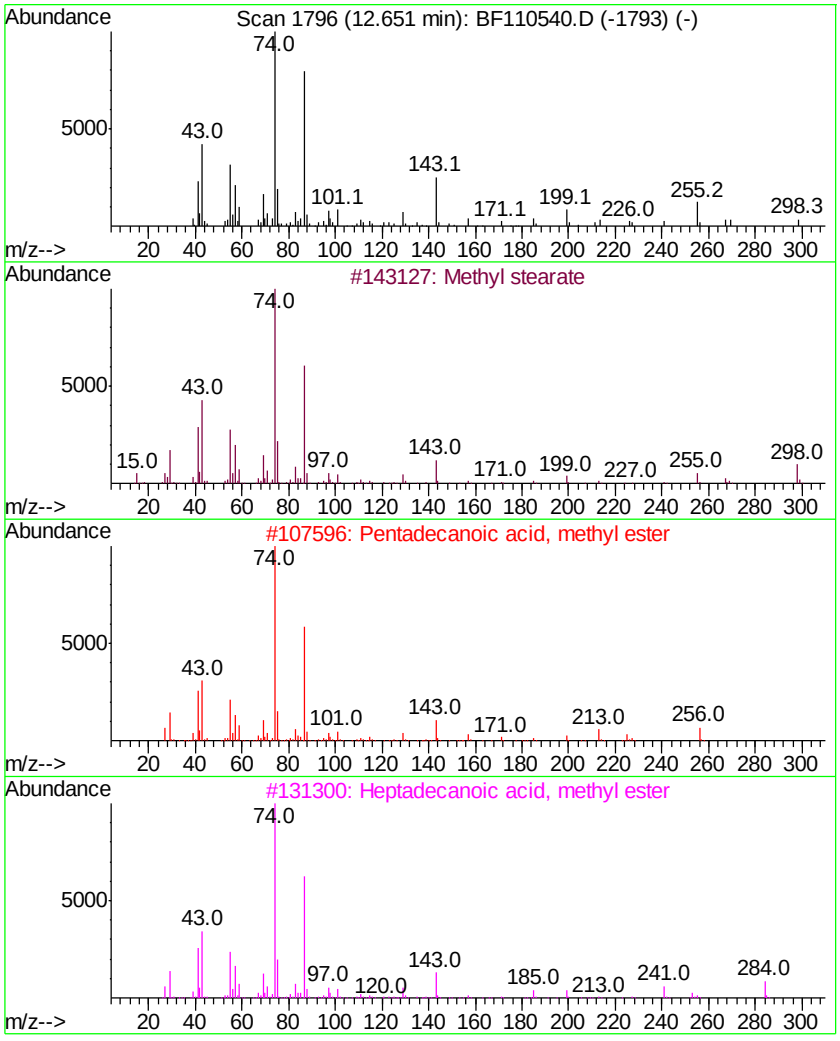
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

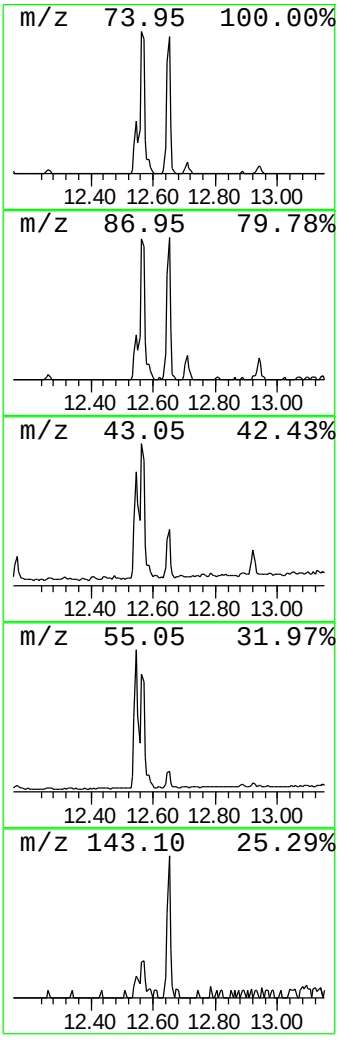
****

Peak Number 8 Methyl stearate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	6.07 ng	130073	Phenanthrene-d10	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	97
2	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	81
3	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	80
4	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	74
5	Methyl tetradecanoate	242	C15H30O2	000124-10-7	72





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Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

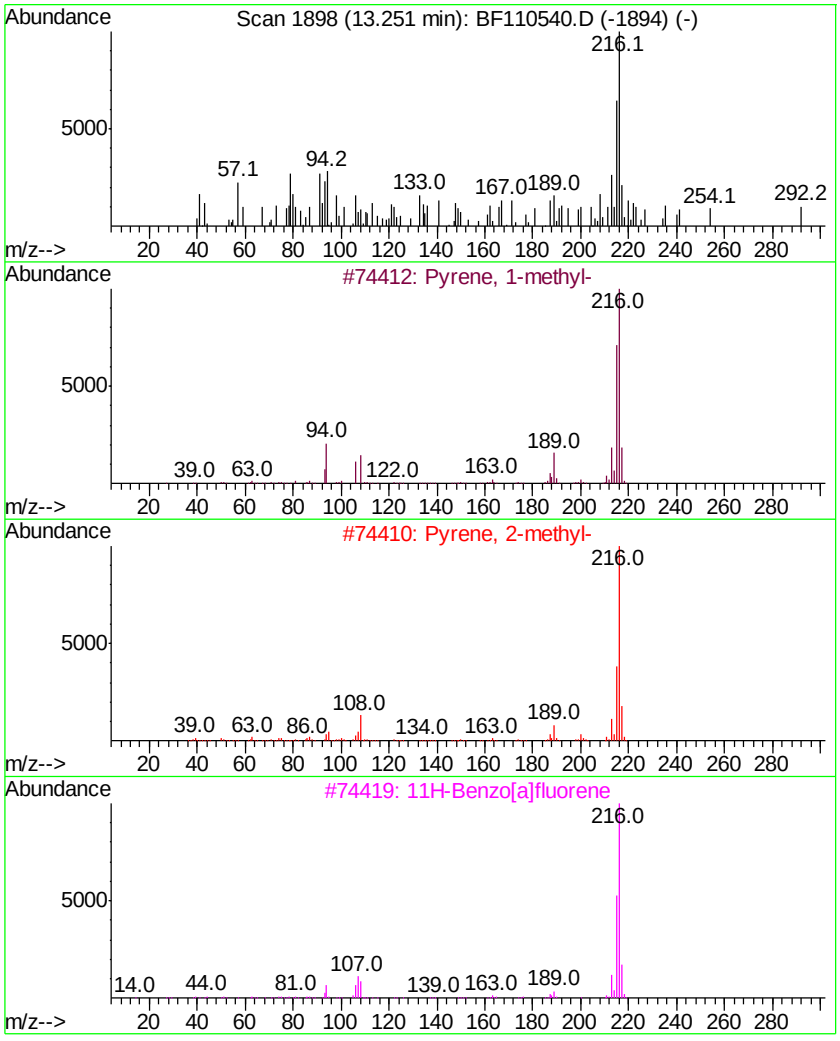
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

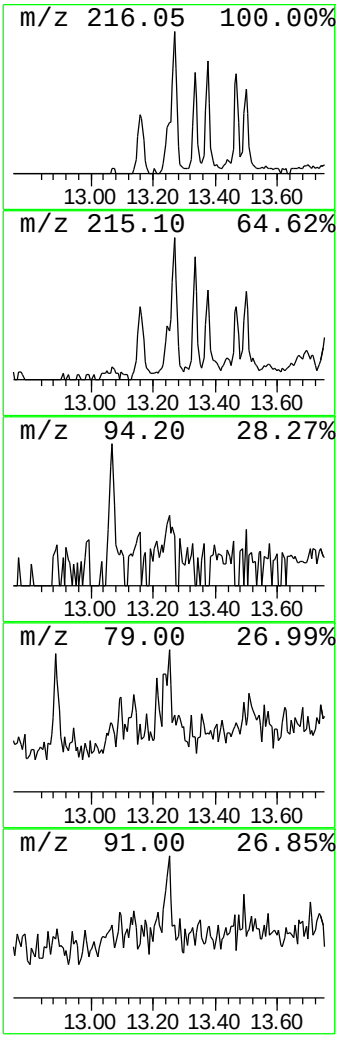
****

Peak Number 9 Pyrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.25	2.08 ng	46663	Chrysene-d12	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	64
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	53
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	53
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	52
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	49





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Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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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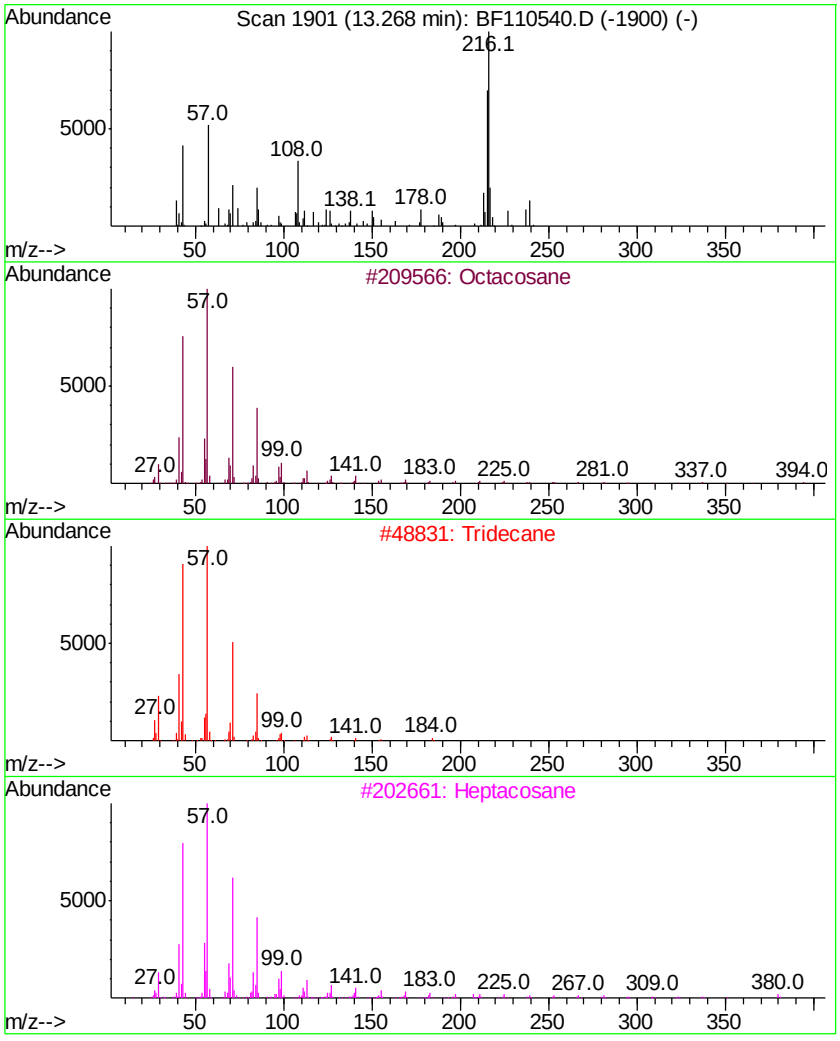
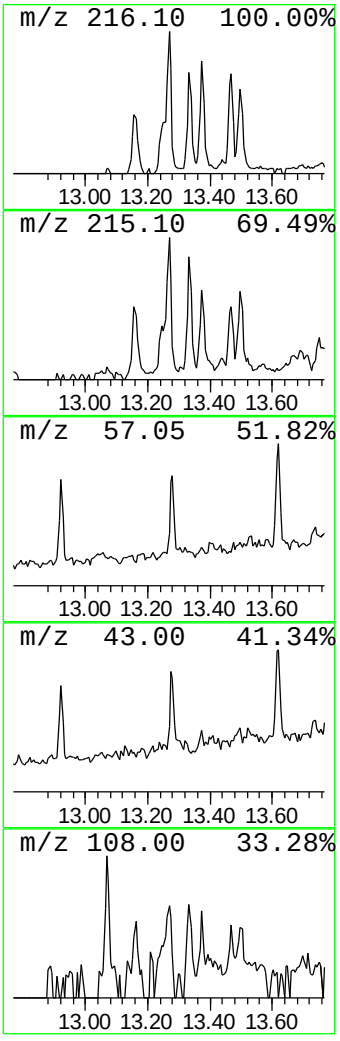
Peak Number 10 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.27	2.92 ng	65599	Chrysene-d12	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	58
2	Tridecane	184	C13H28	000629-50-5	58
3	Heptacosane	380	C27H56	000593-49-7	52
4	Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	52
5	Hentriacontane	437	C31H64	000630-04-6	52





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110618\
Data File : BF110540.D
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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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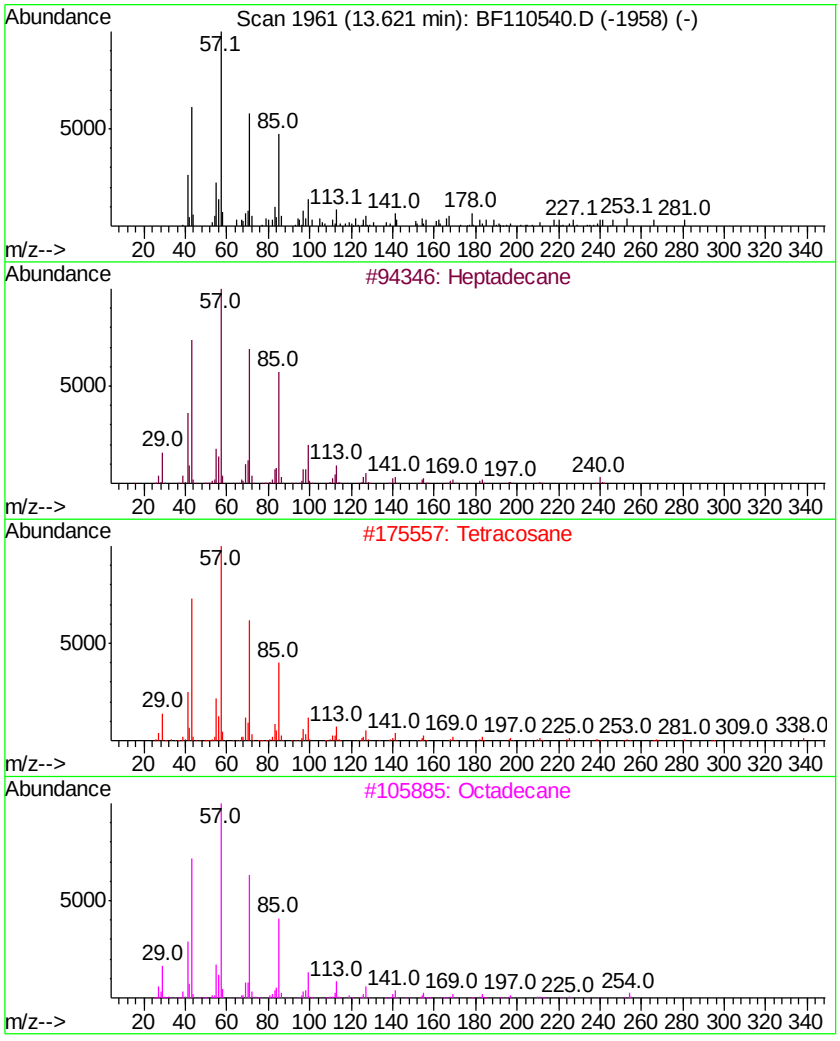
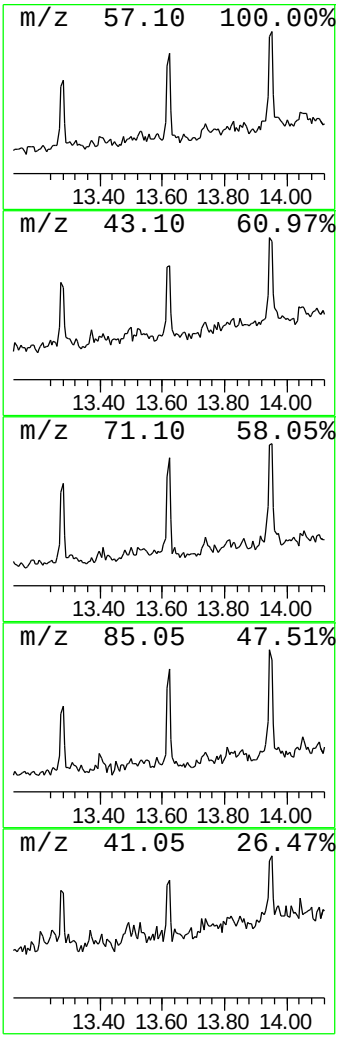
Peak Number 11 Heptadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.62	2.02 ng	45245	Chrysene-d12	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	89
2	Tetracosane	338	C24H50	000646-31-1	74
3	Octadecane	254	C18H38	000593-45-3	72
4	Docosane	310	C22H46	000629-97-0	72
5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	72





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Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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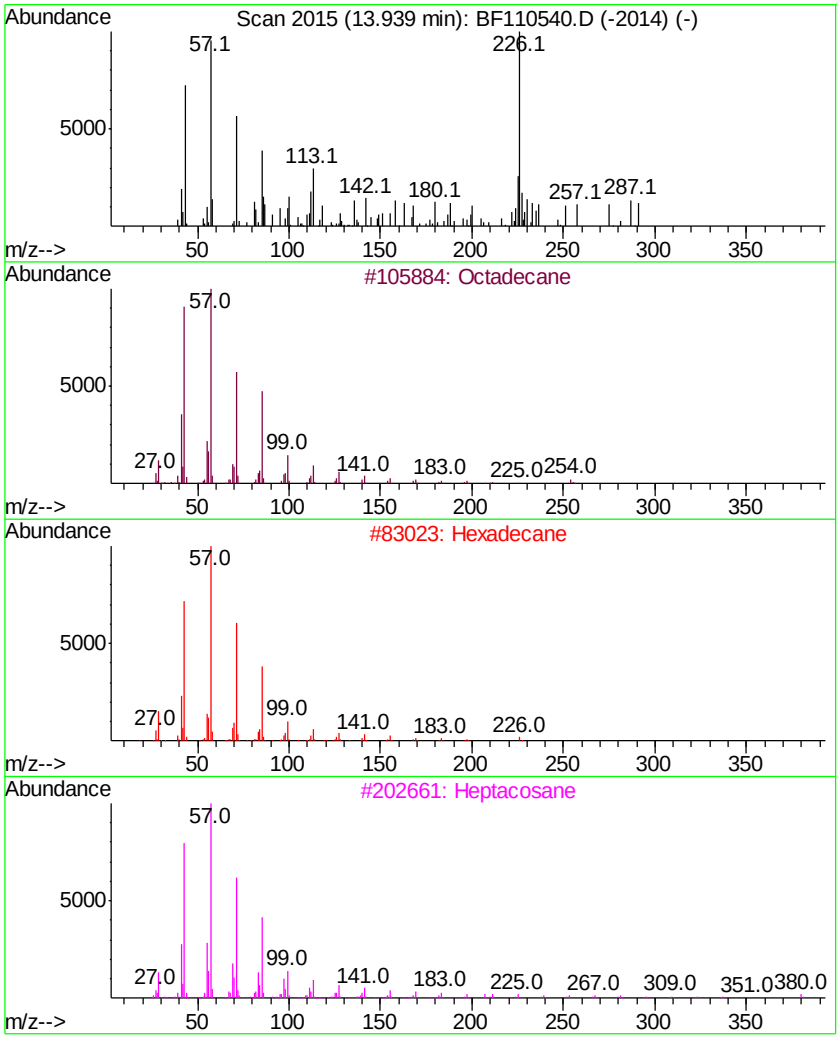
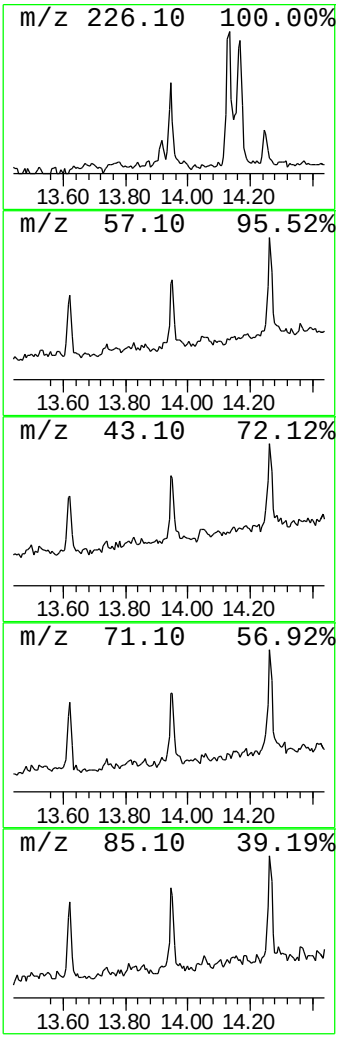
Peak Number 12 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	3.10 ng	69440	Chrysene-d12	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	64
2	Hexadecane	226	C16H34	000544-76-3	62
3	Heptacosane	380	C27H56	000593-49-7	46
4	Docosane, 11-butyl-	366	C26H54	013475-76-8	43
5	Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	43





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Operator : JU/SJ
Sample : J5765-01 5X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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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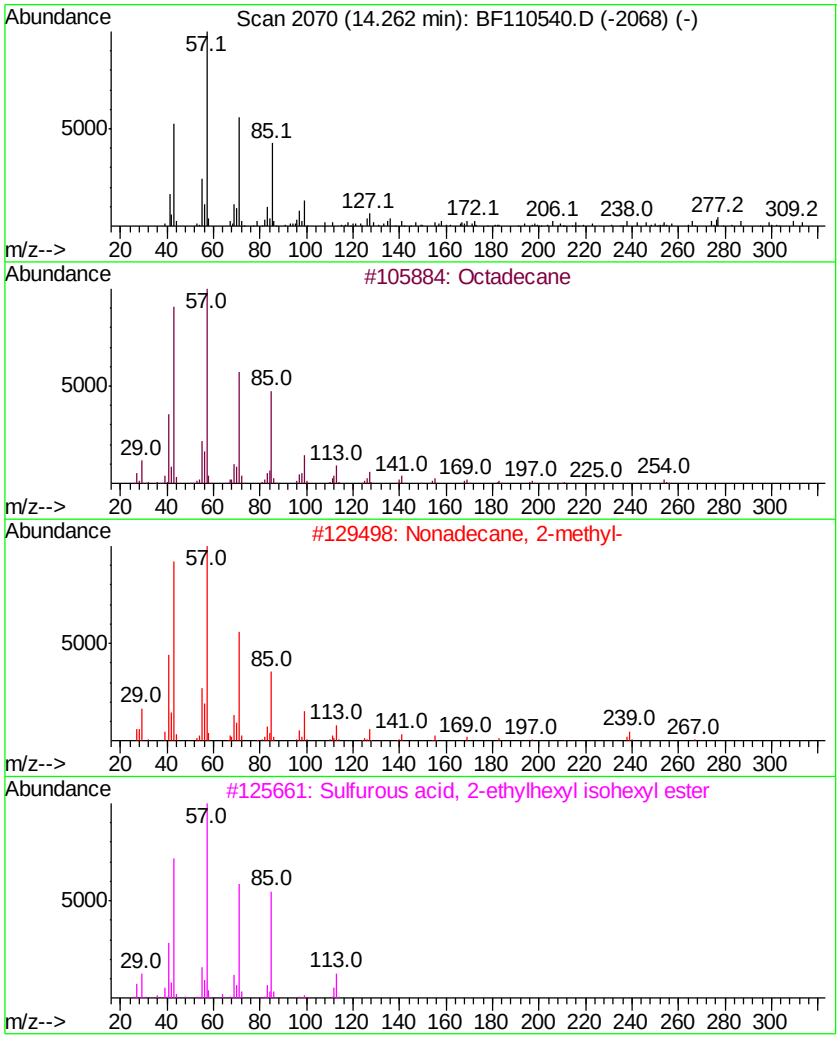
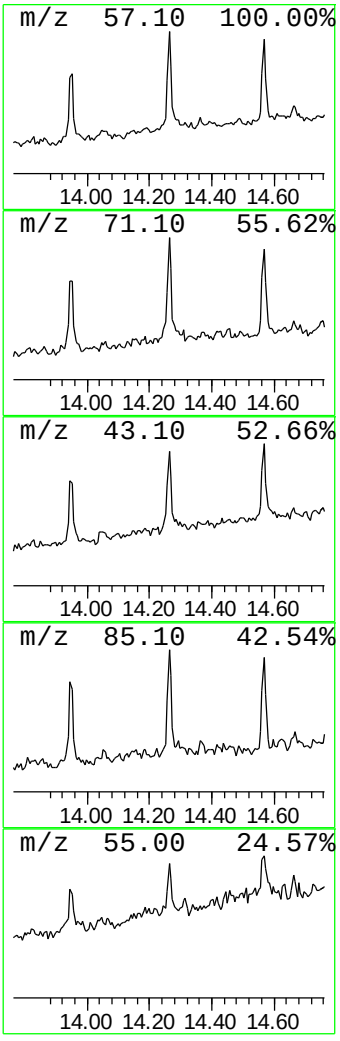
Peak Number 13 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	2.28 ng	51123	Chrysene-d12	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	87
2	Nonadecane, 2-methyl-	282	C20H42	001560-86-7	72
3	Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	72
4	Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	64
5	8-Hydroxy-2,2,8-trimethyldeca-5,...	210	C13H22O2	083040-92-0	64





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Misc :
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Instrument : BNA_F
ClientSampleId : NB-08-110518

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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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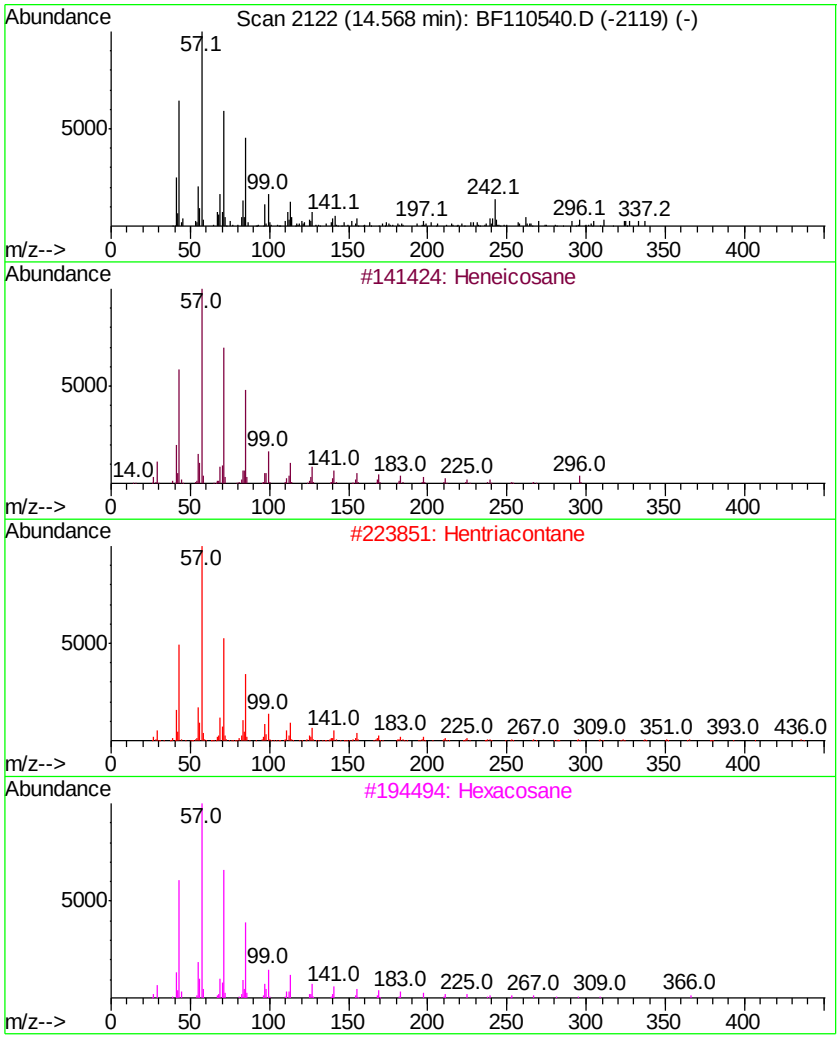
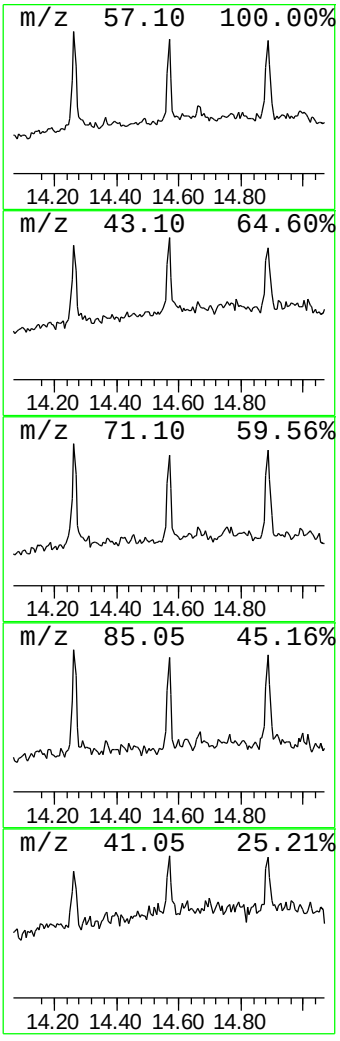
Peak Number 14 Heneicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	3.32 ng	74424	Chrysene-d12	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	93
2	Hentriacontane	437	C31H64	000630-04-6	83
3	Hexacosane	366	C26H54	000630-01-3	83
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	83
5	Triacontane	422	C30H62	000638-68-6	80





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110618\
Data File : BF110540.D
Acq On : 7 Nov 2018 2:37
Operator : JU/SJ
Sample : J5765-01 5X
Misc :
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Instrument : BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102318.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
Butane, 2-methoxy...	2.23	4.1	ng	83052	1	6.96	401159	20.0
unknown6.73	6.73	11.3	ng	225712	1	6.96	401159	20.0
Hexadecanoic acid...	11.86	11.9	ng	255948	4	11.49	428269	20.0
4H-Cyclopenta[def...	12.11	2.6	ng	54950	4	11.49	428269	20.0
9,12-Octadecadien...	12.54	42.4	ng	907736	4	11.49	428269	20.0
9-Octadecenoic ac...	12.57	30.2	ng	646644	4	11.49	428269	20.0
Methyl stearate	12.65	6.1	ng	130073	4	11.49	428269	20.0
Pyrene, 1-methyl-	13.25	2.1	ng	46663	5	14.14	448708	20.0
Octacosane	13.27	2.9	ng	65599	5	14.14	448708	20.0
Heptadecane	13.62	2.0	ng	45245	5	14.14	448708	20.0
Hexadecane	13.94	3.1	ng	69440	5	14.14	448708	20.0
Octadecane	14.26	2.3	ng	51123	5	14.14	448708	20.0
Heneicosane	14.57	3.3	ng	74424	5	14.14	448708	20.0