LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

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-BF012519.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	5.469	572	575	597	rBV	894314	932807	8.92%	1.959%
2	6.481	744	747	762	rBV	1000326	1012304	9.68%	2.126%
3	6.610	766	769	777	rBV	1109891	1073882	10.27%	2.255%
4	6.839	804	808	816	rBV	1591420	1327557	12.70%	2.788%
5	6.998	831	835	842	rBV	880403	857550	8.20%	1.801%
6	7.398	900	903	908	rBV	626307	567636	5.43%	1.192%
7	8.122	1020	1026	1032	rBV	2002834	1827736	17.48%	3.838%
8	8.628	1108	1112	1115	rBV2	111146	112613	1.08%	0.236%
9	8.710	1123	1126	1130	rBV	225534	215181	2.06%	0.452%
10	8.945	1160	1166	1168	rBV2	118962	142638	1.36%	0.300%
11	9.192	1205	1208	1213	rBV	1332168	1149763	11.00%	2.414%
12	9.275	1219	1222	1227	rBV	341786	317598	3.04%	0.667%
13	9.598	1272	1277	1279	rBV3	152441	203029	1.94%	0.426%
14	9.804	1309	1312	1317	rVB	427958	341065	3.26%	0.716%
15	9.875	1321	1324	1328	rVB	1964264	1769295	16.92%	3.715%
16	10.122	1363	1366	1371	rBV5	100285	145153	1.39%	0.305%
17	10.304	1391	1397	1400	rBV	505145	448270	4.29%	0.941%
18	10.427	1411	1418	1423	rBV3	78570	163882	1.57%	0.344%
19	10.522	1430	1434	1440	rVB	145319	213994	2.05%	0.449%
20	10.663	1456	1458	1465	rBV	563460	541865	5.18%	1.138%
21	10.774	1474	1477	1487	rVB	468751	556471	5.32%	1.169%
22	11.222	1550	1553	1556	rBV	420317	350134	3.35%	0.735%
23	11.257	1556	1559	1564	rVB2	182945	214430	2.05%	0.450%
24	11.363	1573	1577	1579	rBV	1897683	1618886	15.48%	3.400%
25	11.386	1579	1581	1588	rVB	401624	375839	3.59%	0.789%
26	11.651	1623	1626	1628	rBV	336615	289013	2.76%	0.607%
27	11.751	1639	1643	1646	rBV	4063753	3268572	31.26%	6.864%
28	11.892	1664	1667	1672	rVB3	110581	140392	1.34%	0.295%
29	12.057	1691	1695	1700	rBV	260709	263462	2.52%	0.553%
30	12.439	1752	1760	1762	rBV	9195460	10457128	100.00%	21.959%
31	12.463	1762	1764	1769	rVB2	8749635	7370519	70.48%	15.478%
32	12.539	1774	1777	1780	rVB	1669526	1304115	12.47%	2.739%
33	12.574	1780	1783	1788	rBV	458554	438423	4.19%	0.921%
34	12.810	1818	1823	1829	rVB2	477281	581898	5.56%	1.222%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

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-BF012519.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.939	1841	1845	1849	rBV	1032956	889165	8.50%	1.867%
36	13.168	1881	1884	1885	rBV	165374	139055	1.33%	0.292%
37	13.263	1897	1900	1903	rVB	143502	127152	1.22%	0.267%
38	13.510	1939	1942	1945	rBV	126191	107243	1.03%	0.225%
39	13.839	1995	1998	2003	rVB3	131368	167290	1.60%	0.351%
40	13.927	2010	2013	2017	rBV2	159405	162240	1.55%	0.341%
41	13.998	2021	2025	2028	rVV2	1400631	1534398	14.67%	3.222%
42	14.027	2028	2030	2037	rVB	184581	191500	1.83%	0.402%
43	14.151	2048	2051	2057	rBV2	141011	143251	1.37%	0.301%
44	14.457	2099	2103	2106	rBV	217509	238321	2.28%	0.500%
45	14.757	2151	2154	2159	rBV	202630	231265	2.21%	0.486%
46	14.874	2171	2174	2183	rVB	166568	253391	2.42%	0.532%
47	15.039	2199	2202	2208	rBV	147297	276891	2.65%	0.581%
48	15.092	2208	2211	2218	rVB	287380	366863	3.51%	0.770%
49	15.404	2261	2264	2270	rBV	144523	190518	1.82%	0.400%
50	15.468	2270	2275	2283	rVV	1200503	1558063	14.90%	3.272%
51	15.898	2345	2348	2353	rVB2	177278	250563	2.40%	0.526%
52	16.398	2429	2433	2438	rVB2	121875	200575	1.92%	0.421%

Sum of corrected areas: 47620844

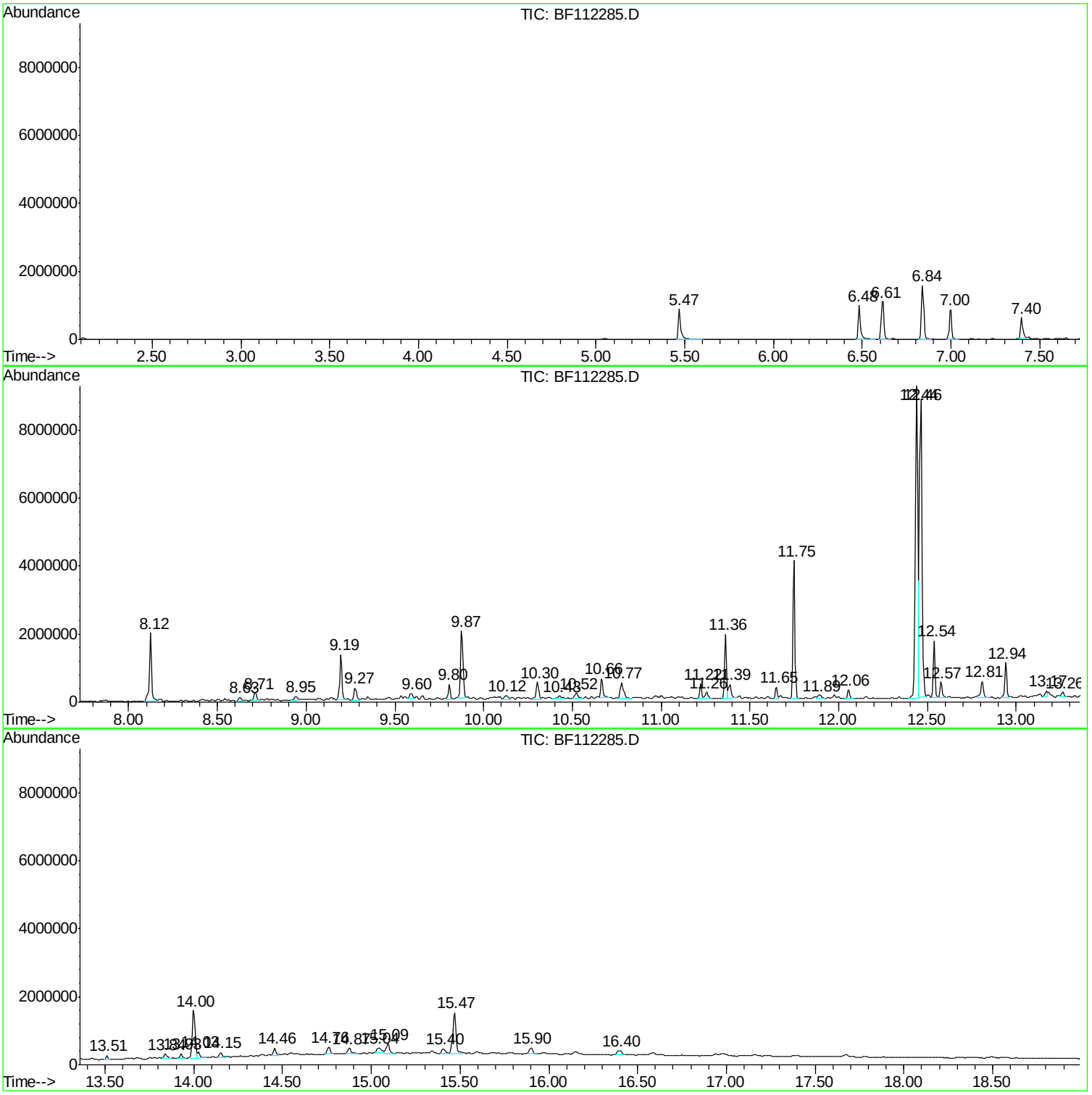
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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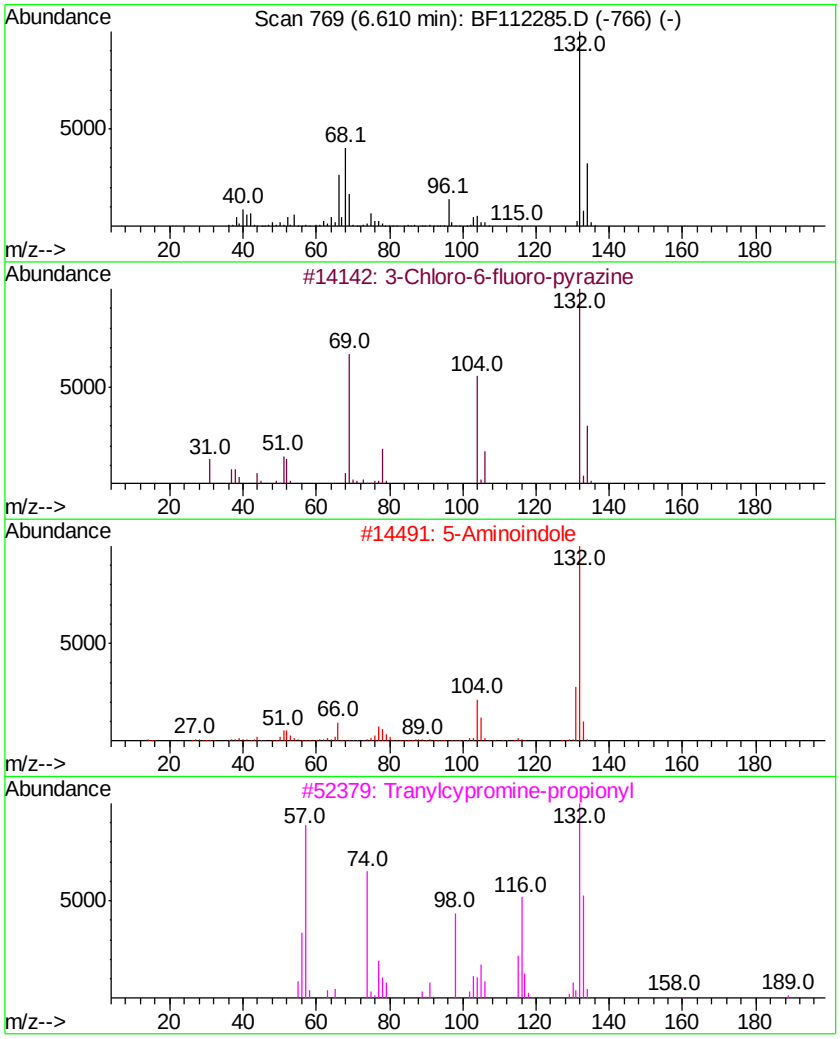
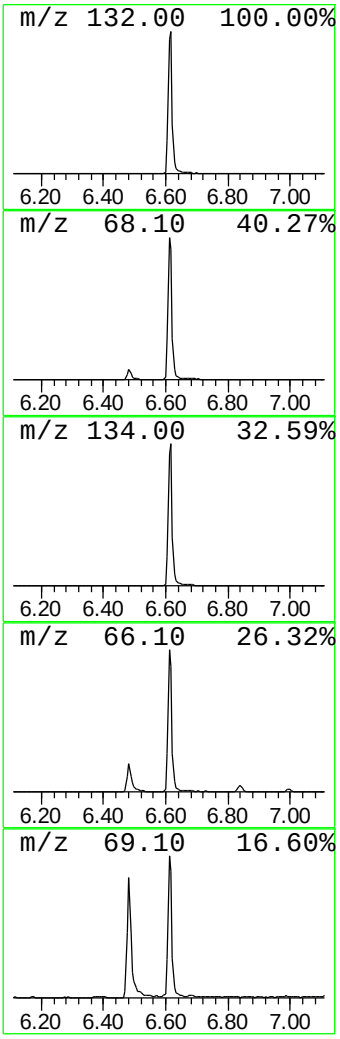
Peak Number 1 unknown6.61 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	16.18 ng	1073880	1,4-Dichlorobenzene-d4	6.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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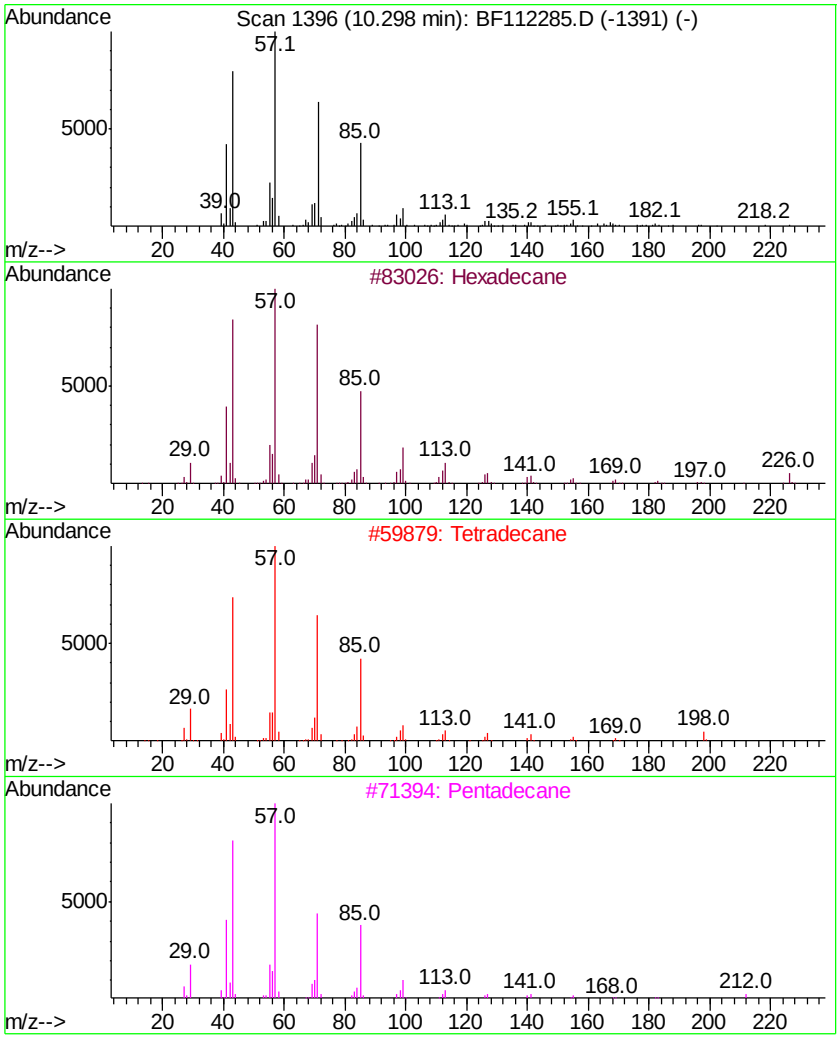
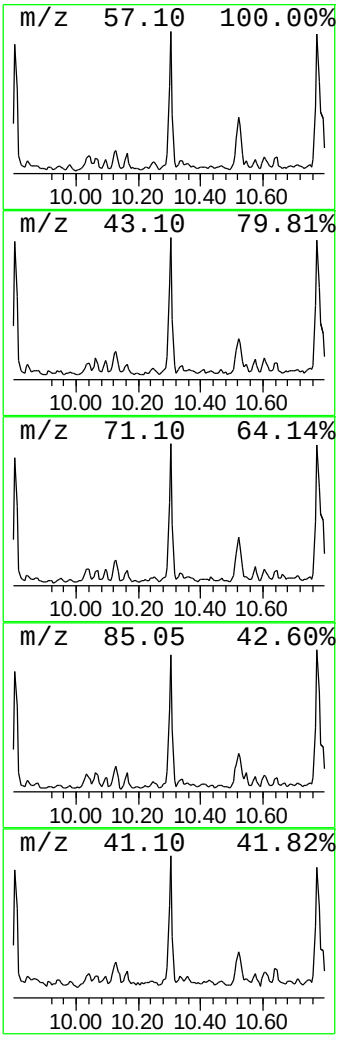
Peak Number 5 Hexadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	5.07 ng	448270	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	97
2	Tetradecane	198	C14H30	000629-59-4	91
3	Pentadecane	212	C15H32	000629-62-9	91
4	Dodecane	170	C12H26	000112-40-3	90
5	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1000309-12-5	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

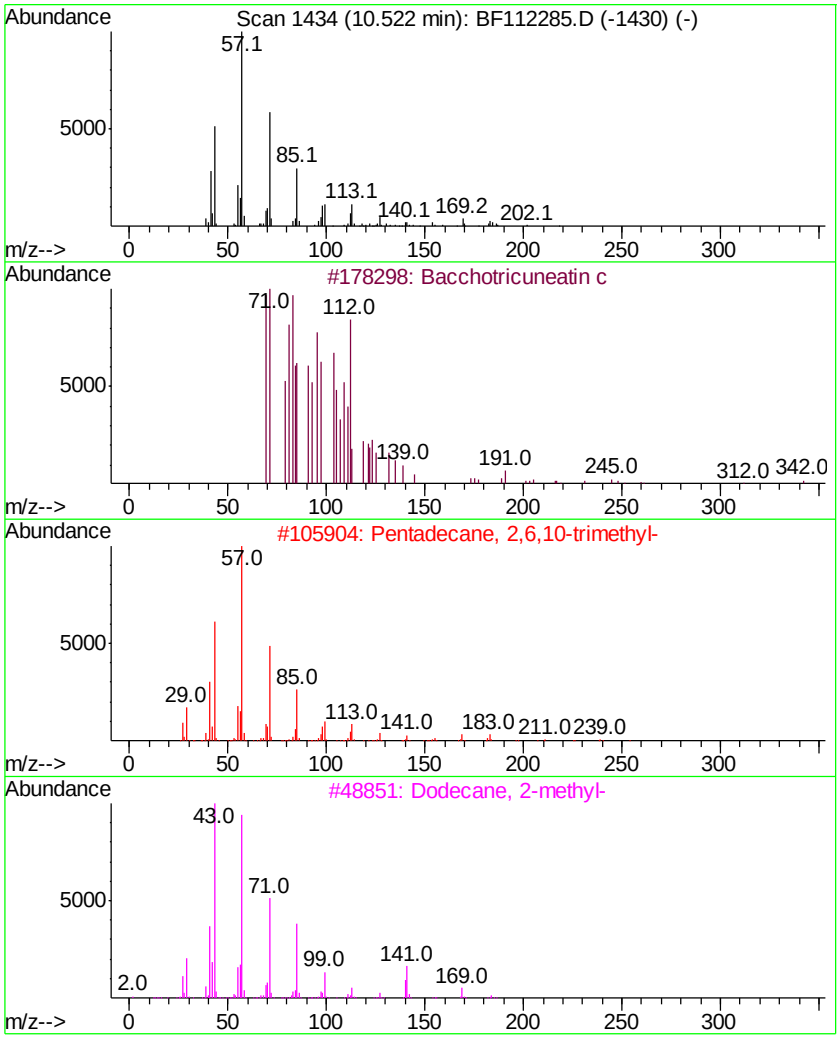
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

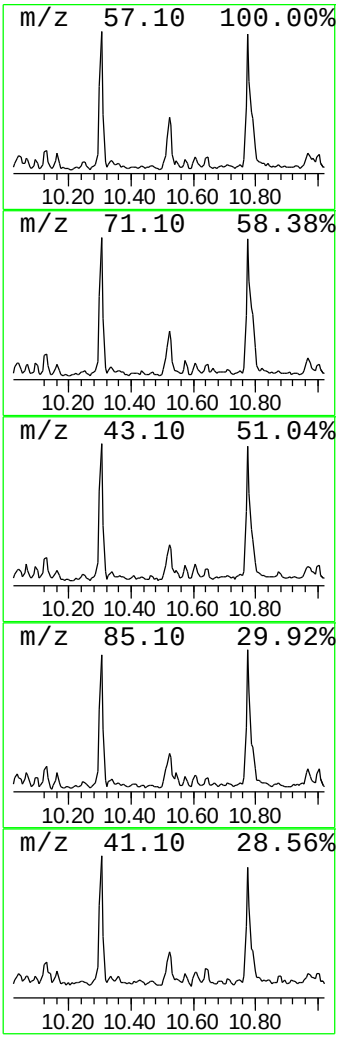
****

Peak Number 6 Bacchotricuneatin c Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.52	2.42 ng	213994	Acenaphthene-d10	9.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bacchotricuneatin c	342	C20H22O5	066563-30-2	97
2	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
3	Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
4	Tetracosane	338	C24H50	000646-31-1	72
5	Undecane, 5-methyl-	170	C12H26	001632-70-8	68





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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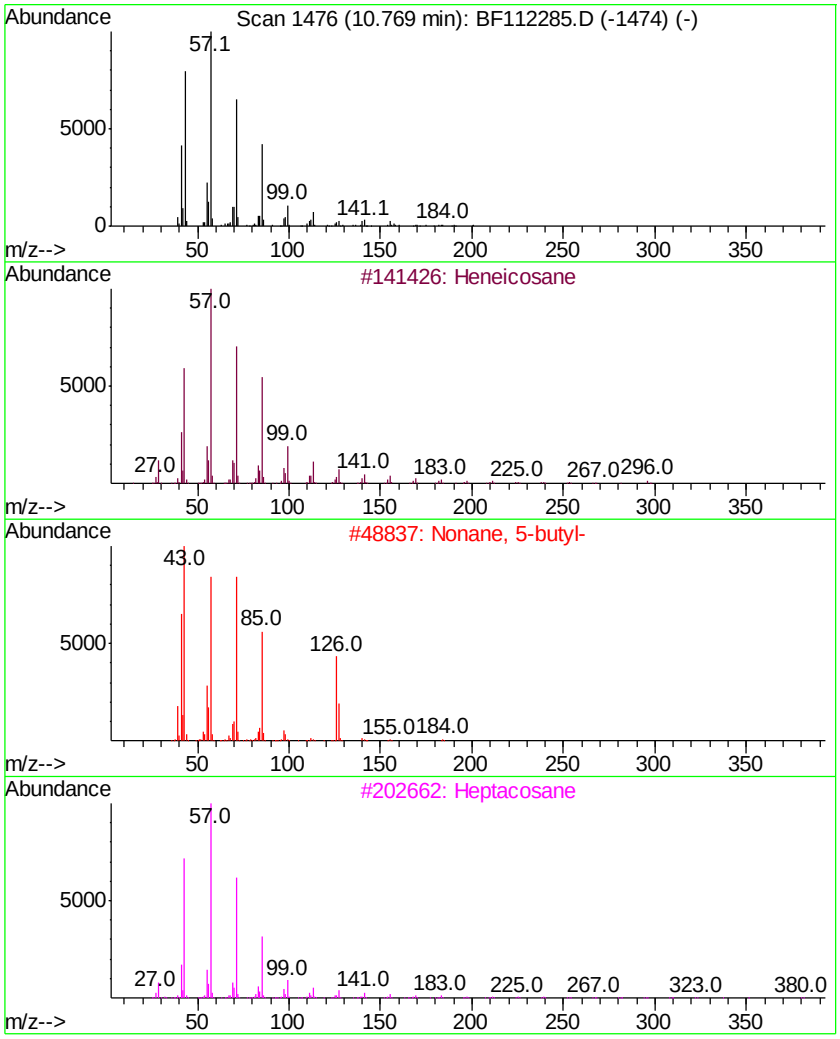
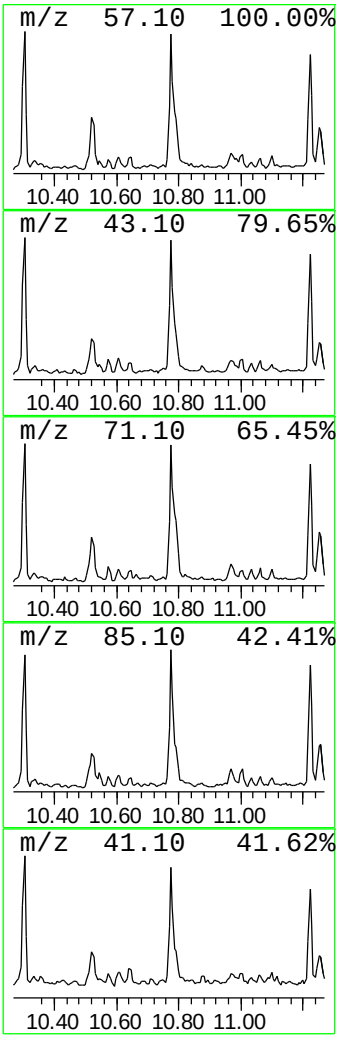
Peak Number 7 Heneicosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	6.87 ng	556471	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Nonane, 5-butyl-	184	C13H28	017312-63-9	90
3	Heptacosane	380	C27H56	000593-49-7	90
4	Tetradecane	198	C14H30	000629-59-4	87
5	Heptadecane	240	C17H36	000629-78-7	87





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Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

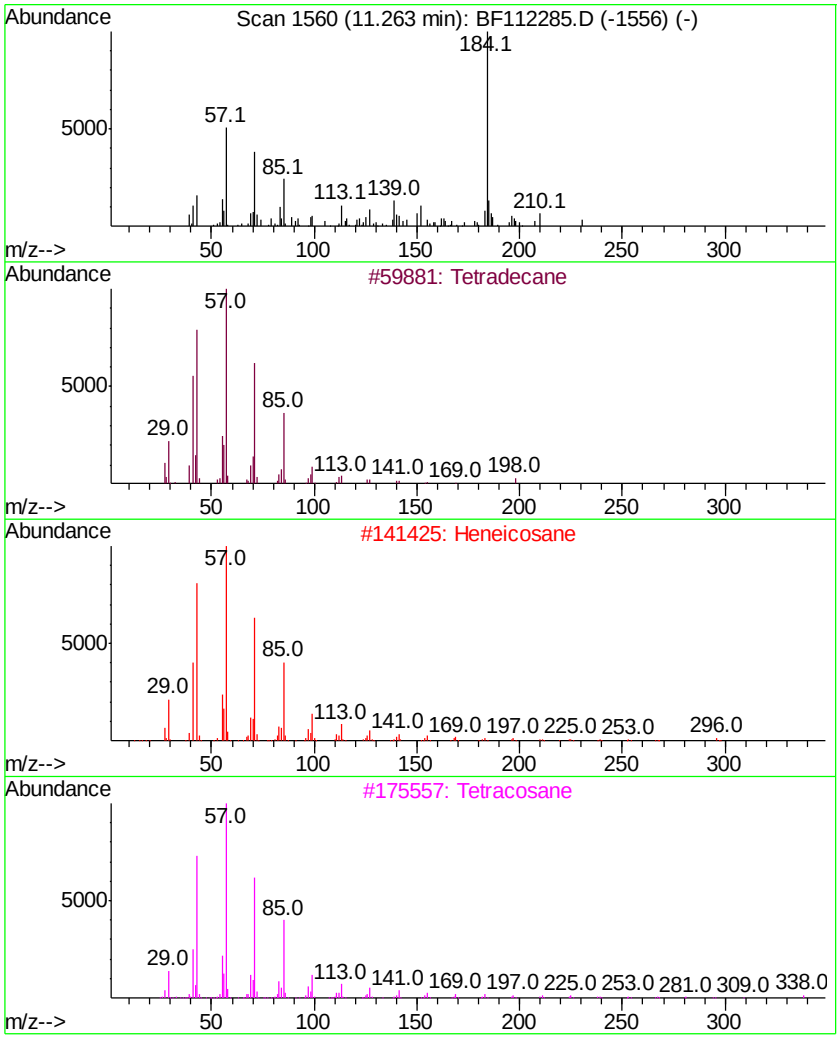
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

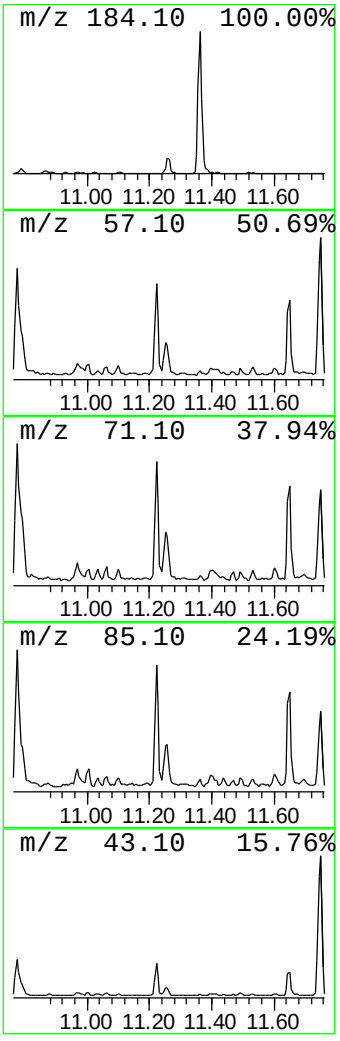
****

Peak Number 9 Tetradecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.26	2.65 ng	214430	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	64
2	Heneicosane	296	C21H44	000629-94-7	58
3	Tetracosane	338	C24H50	000646-31-1	58
4	2-Bromo dodecane	248	C12H25Br	013187-99-0	58
5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	58





Library Search Compound Report

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Misc :
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Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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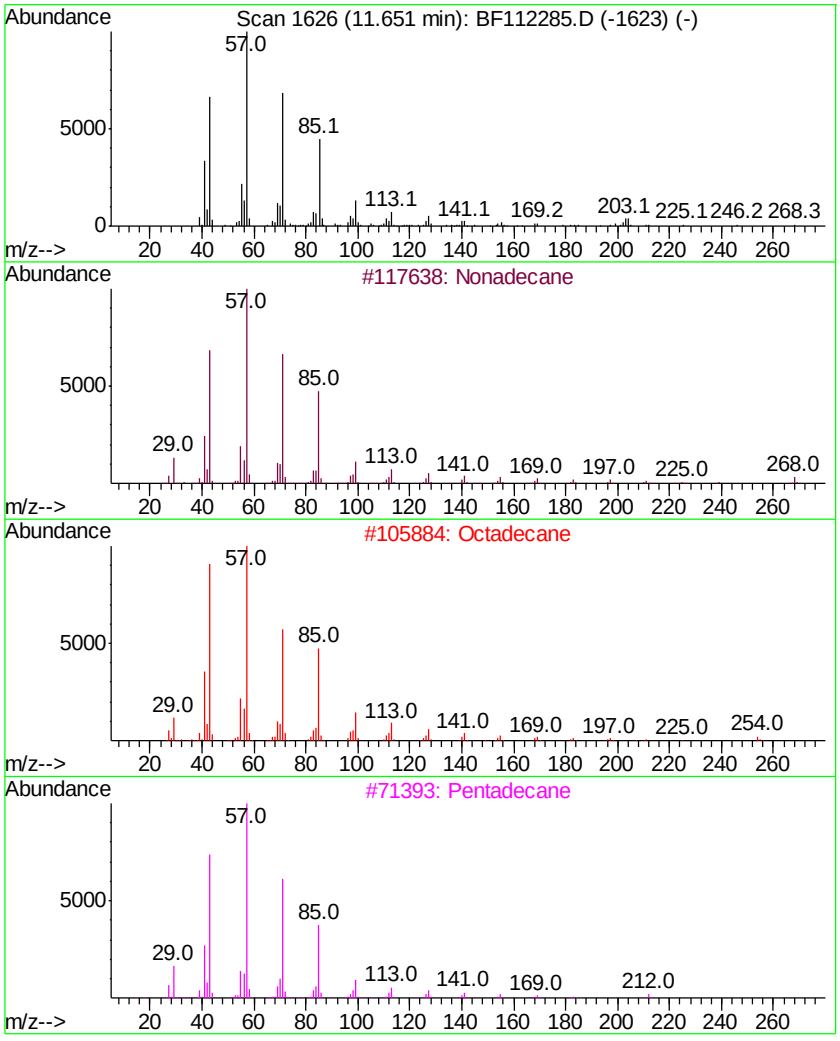
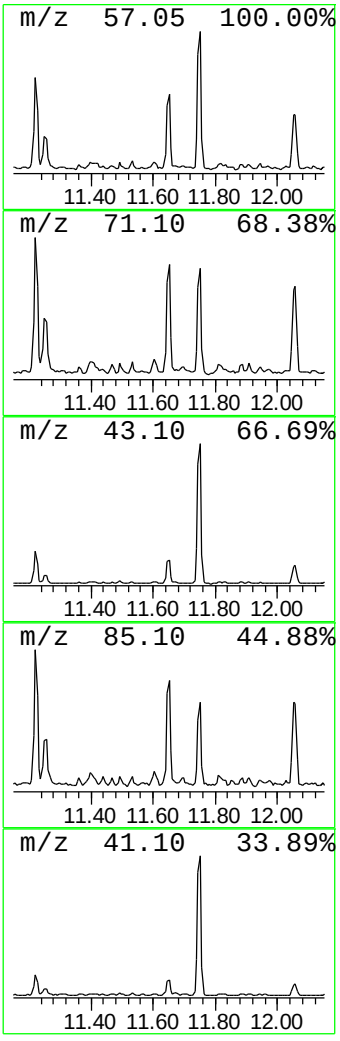
Peak Number 10 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	3.57 ng	289013	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	93
2	Octadecane	254	C18H38	000593-45-3	87
3	Pentadecane	212	C15H32	000629-62-9	87
4	Heneicosane	296	C21H44	000629-94-7	86
5	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1000309-12-5	86





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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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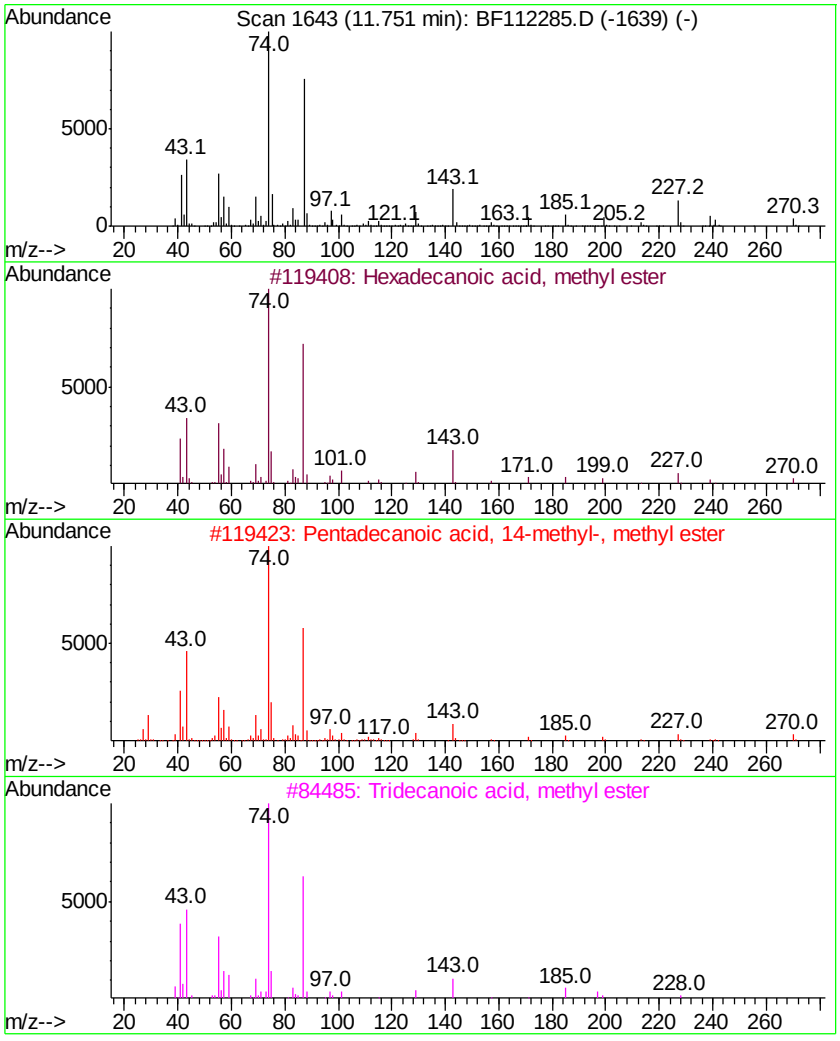
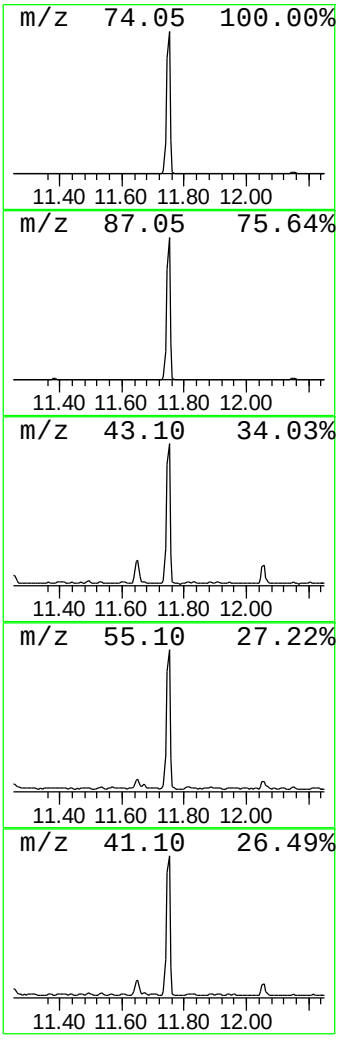
Peak Number 11 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	40.38 ng	3268570	Phenanthrene-d10	11.36

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	92
3	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5	Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	83





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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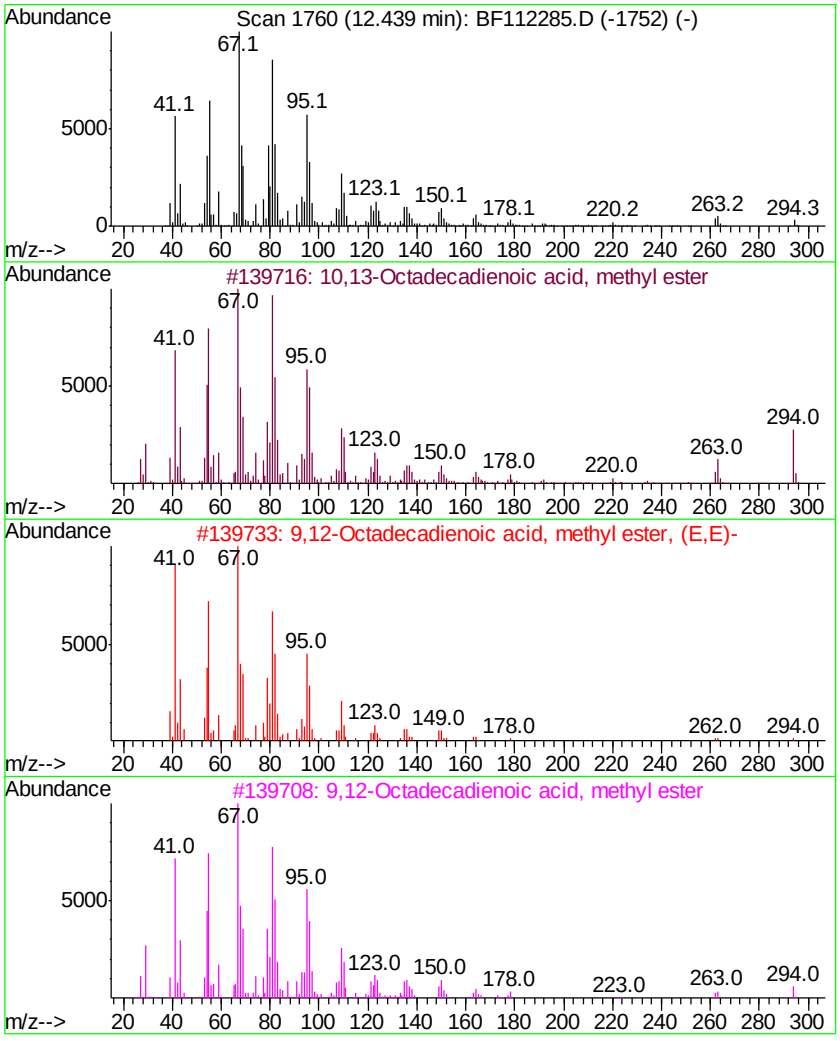
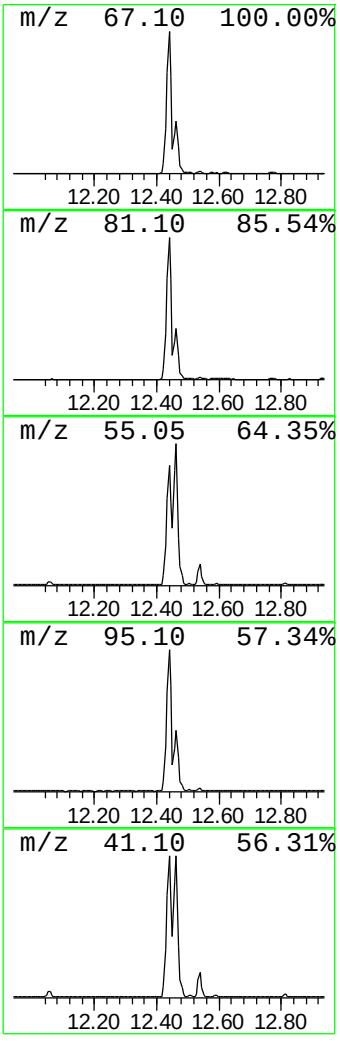
Peak Number 13 10,13-Octadecadienoic acid,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.44	129.19 ng	10457100	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
4	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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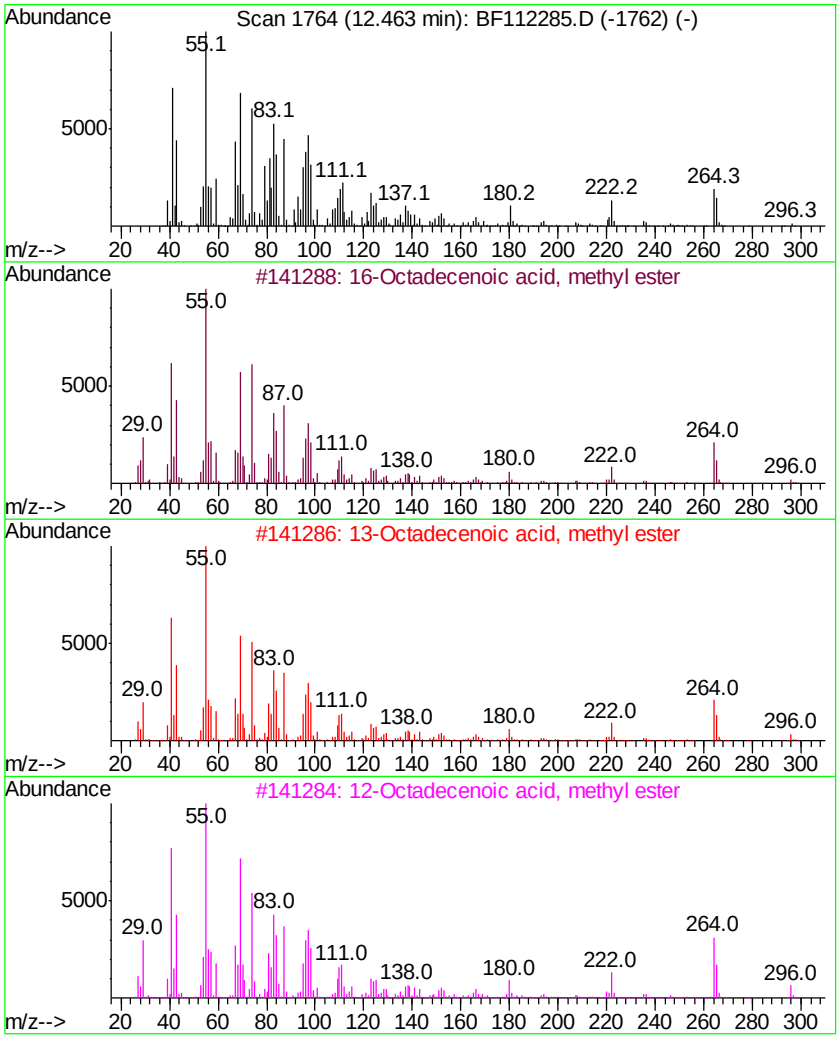
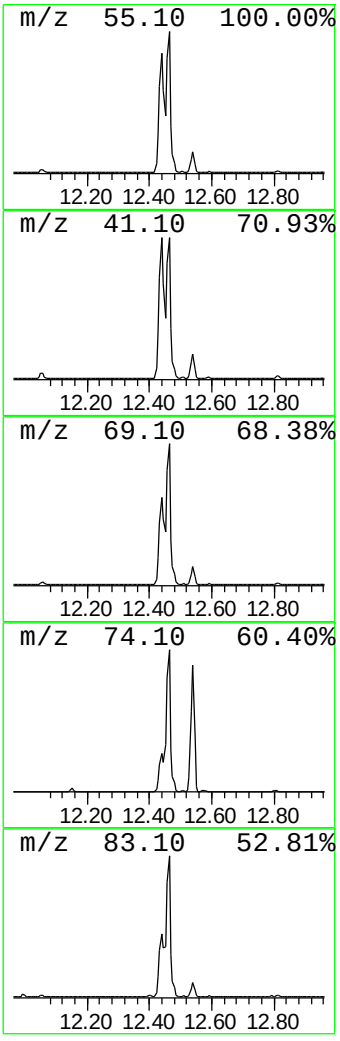
Peak Number 14 16-Octadecenoic acid, methy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	91.06 ng	7370520	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
2	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
3	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
4	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99





Library Search Compound Report

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Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.M
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TIC Library : C:\DATABASE\NIST11.L
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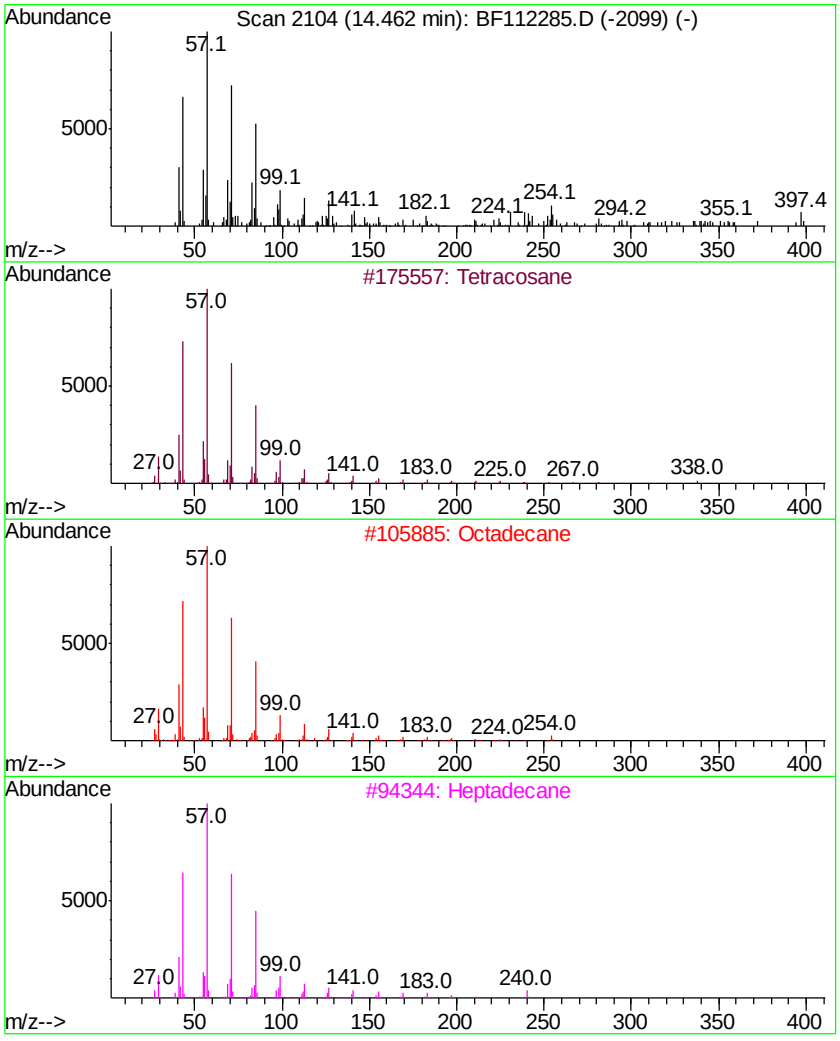
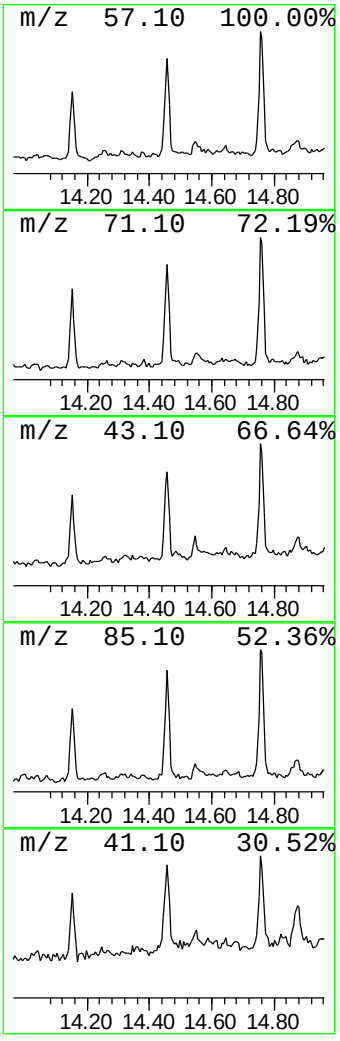
Peak Number 15 Tetracosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	3.11 ng	238321	Chrysene-d12	14.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Octadecane	254	C18H38	000593-45-3	96
3	Heptadecane	240	C17H36	000629-78-7	94
4	2-methyloctacosane	408	C29H60	1000376-72-8	90
5	Hexacosane	366	C26H54	000630-01-3	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.M
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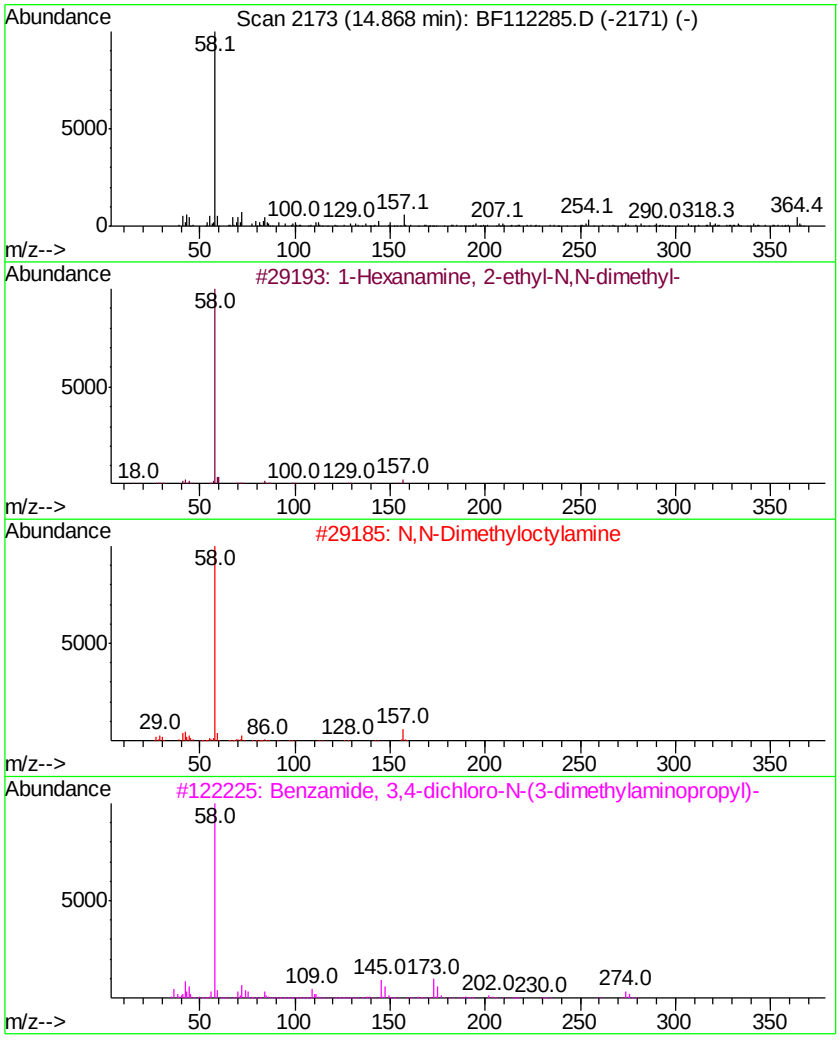
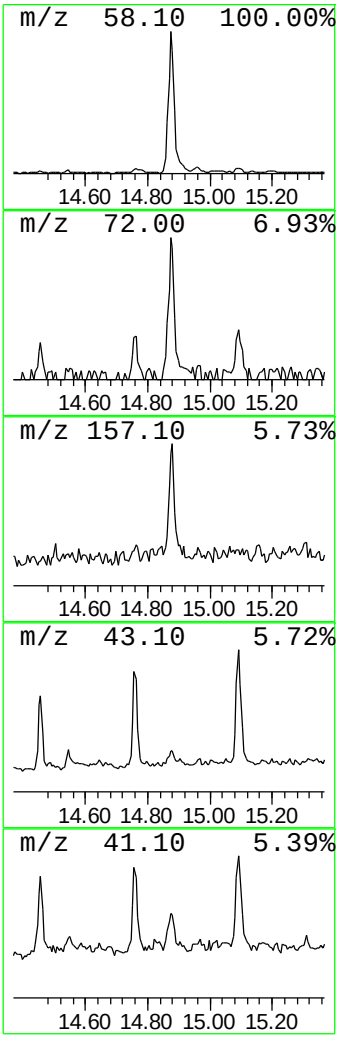
Peak Number 17 1-Hexanamine, 2-ethyl-N,N-d... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	3.25 ng	253391	Perylene-d12	15.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanamine, 2-ethyl-N,N-dimethyl-	157	C10H23N	028056-87-3	64
2	N,N-Dimethyloctylamine	157	C10H23N	007378-99-6	64
3	Benzamide, 3,4-dichloro-N-(3-dim...	274	C12H16Cl2N2O	333347-10-7	64
4	3-(Dimethylamino)propyl isothioc...	144	C6H12N2S	027421-70-1	64
5	3-Hexanamine, 3-ethyl-	129	C8H19N	056667-17-5	59





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Misc :
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Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.M
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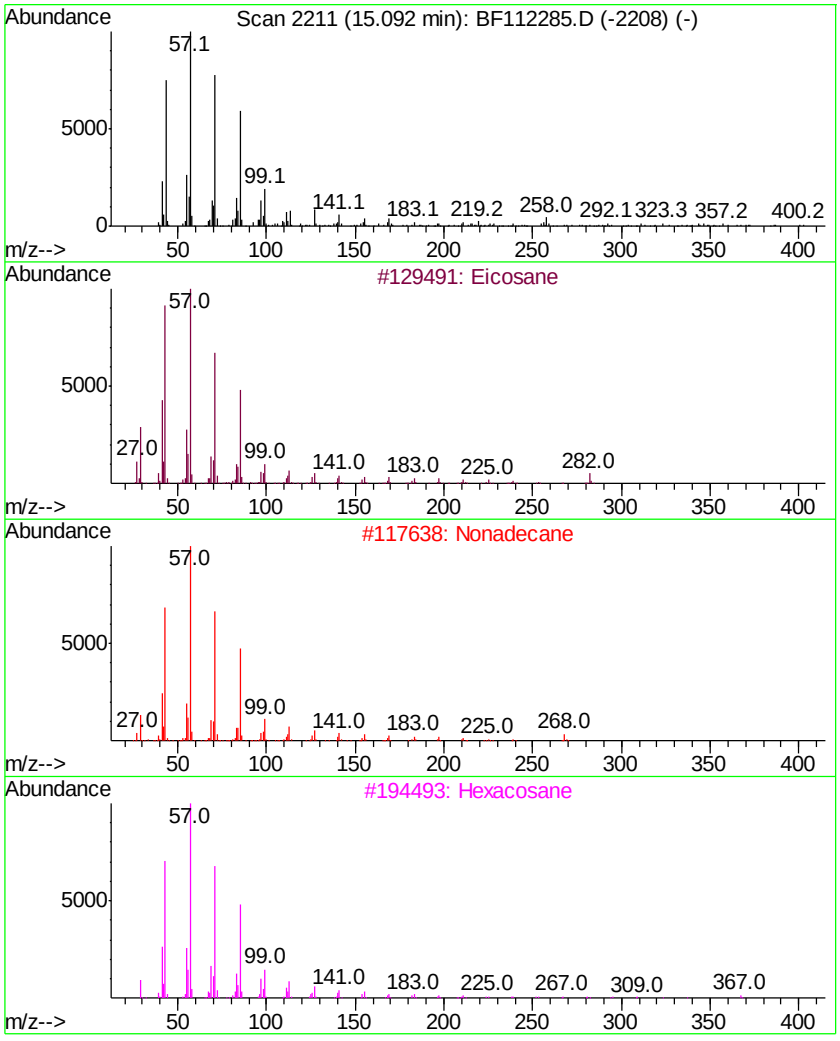
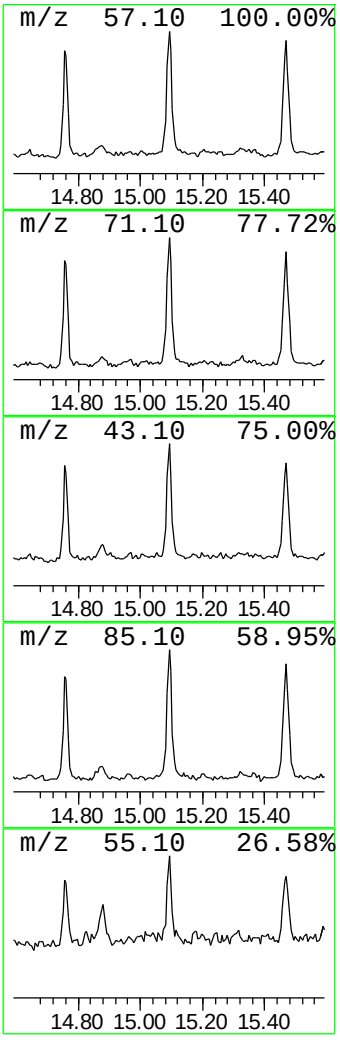
Peak Number 18 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	4.71 ng	366863	Perylene-d12	15.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	90
2	Nonadecane	268	C19H40	000629-92-5	90
3	Hexacosane	366	C26H54	000630-01-3	90
4	Heneicosane	296	C21H44	000629-94-7	87
5	Heptacosane	380	C27H56	000593-49-7	87





Library Search Compound Report

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Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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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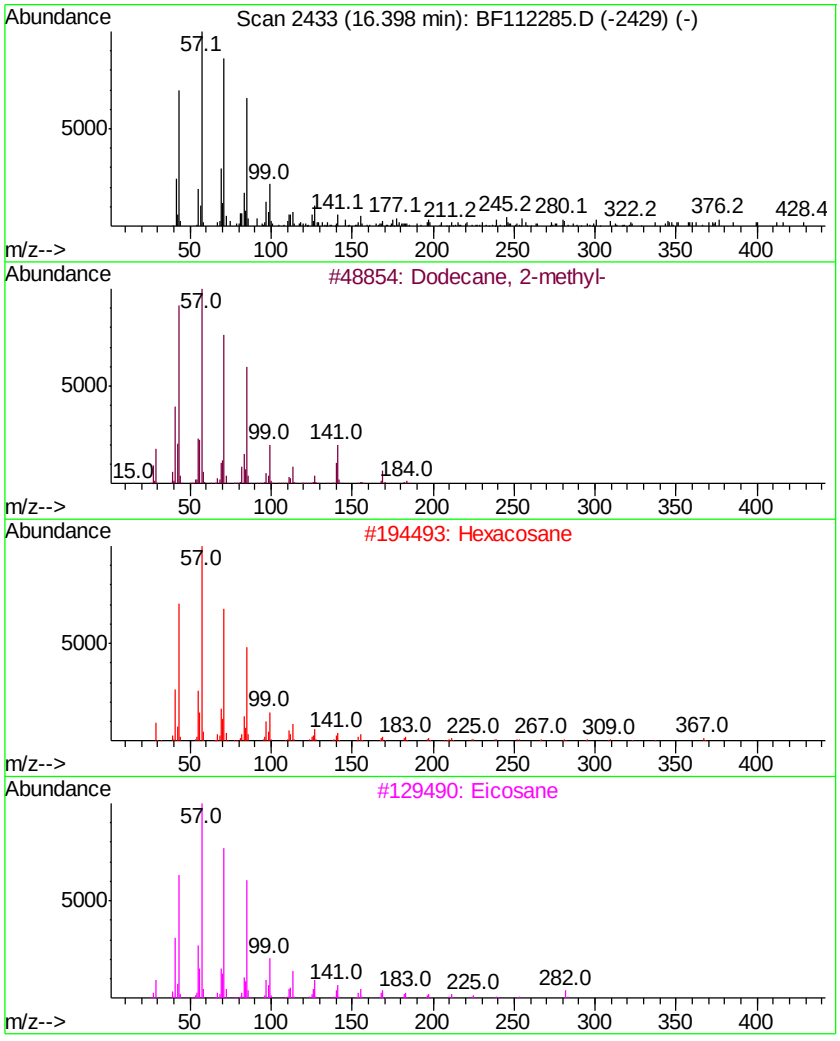
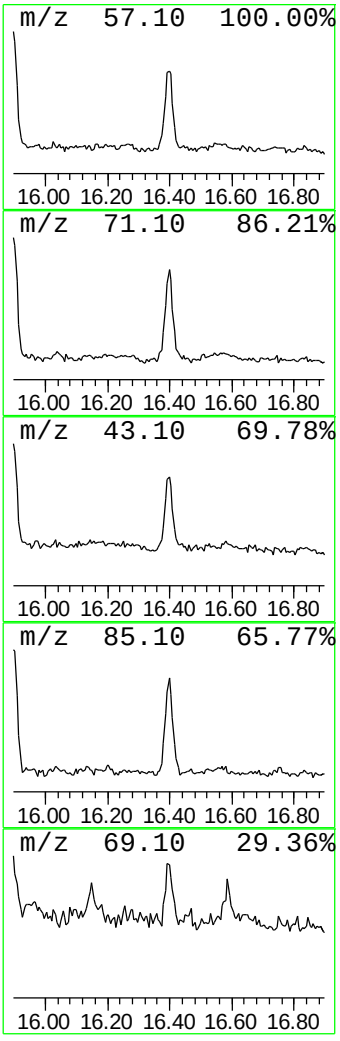
Peak Number 20 Dodecane, 2-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	2.57 ng	200575	Perylene-d12	15.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
2	Hexacosane	366	C26H54	000630-01-3	83
3	Eicosane	282	C20H42	000112-95-8	83
4	Heneicosane	296	C21H44	000629-94-7	83
5	10-Methylnonadecane	282	C20H42	056862-62-5	80





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF012819\
Data File : BF112285.D
Acq On : 28 Jan 2019 20:06
Operator : JU/SJ
Sample : K1011-11 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : OR-01-012519-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012519.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	16.2	ng	1073880	1	6.84	1327560	20.0
Hexadecane	10.30	5.1	ng	448270	3	9.87	1769300	20.0
Bacchotricuneatin c	10.52	2.4	ng	213994	3	9.87	1769300	20.0
Heneicosane	10.77	6.9	ng	556471	4	11.36	1618890	20.0
Tetradecane	11.26	2.6	ng	214430	4	11.36	1618890	20.0
Nonadecane	11.65	3.6	ng	289013	4	11.36	1618890	20.0
Hexadecanoic acid...	11.75	40.4	ng	3268570	4	11.36	1618890	20.0
10,13-Octadecadie...	12.44	129.2	ng	10457100	4	11.36	1618890	20.0
16-Octadecenoic a...	12.46	91.1	ng	7370520	4	11.36	1618890	20.0
Tetracosane	14.46	3.1	ng	238321	5	14.00	1534400	20.0
1-Hexanamine, 2-e...	14.87	3.3	ng	253391	6	15.47	1558060	20.0
Eicosane	15.09	4.7	ng	366863	6	15.47	1558060	20.0
Dodecane, 2-methyl-	16.40	2.6	ng	200575	6	15.47	1558060	20.0