

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111611.D
 Acq On : 20 Dec 2018 00:57
 Operator : JU/SJ
 Sample : J6196-11 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-121818-C

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF121918.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	3.993	319	324	331	rBV	94844	120605	1.69%	0.254%
2	5.104	509	513	518	rBV	122137	126856	1.78%	0.267%
3	5.510	575	582	585	rBV	2332406	2203344	30.96%	4.641%
4	5.751	616	623	630	rBV	146248	141018	1.98%	0.297%
5	6.510	748	752	759	rVB	2348509	2351534	33.04%	4.953%
6	6.663	774	778	781	rVB	2503536	2135634	30.01%	4.498%
7	6.722	785	788	794	rBV	165795	171275	2.41%	0.361%
8	6.892	814	817	821	rBV	1416900	1242661	17.46%	2.617%
9	7.051	840	844	847	rBV	2011726	1581314	22.22%	3.331%
10	7.451	903	912	915	rBV	1388820	1294432	18.19%	2.727%
11	8.175	1030	1035	1038	rBV	1901786	1740297	24.45%	3.666%
12	8.198	1038	1039	1042	rVB	404520	197945	2.78%	0.417%
13	8.775	1133	1137	1140	rBV	198022	158564	2.23%	0.334%
14	8.886	1150	1156	1160	rVB	181777	178395	2.51%	0.376%
15	8.986	1169	1173	1177	rBV2	105106	114809	1.61%	0.242%
16	9.251	1214	1218	1221	rBV	2708888	2124782	29.85%	4.476%
17	9.339	1229	1233	1237	rBV	264080	244063	3.43%	0.514%
18	9.510	1256	1262	1267	rBV4	74859	112720	1.58%	0.237%
19	9.592	1269	1276	1277	rBV4	72361	105093	1.48%	0.221%
20	9.639	1282	1284	1286	rBV	261683	208928	2.94%	0.440%
21	9.869	1320	1323	1325	rVB	271351	195746	2.75%	0.412%
22	9.927	1330	1333	1337	rVB	1740694	1592597	22.38%	3.355%
23	10.369	1402	1408	1410	rBV	283122	294840	4.14%	0.621%
24	10.586	1439	1445	1451	rBV	218107	305656	4.29%	0.644%
25	10.710	1463	1466	1470	rBV	1429405	1192083	16.75%	2.511%
26	10.839	1485	1488	1489	rBV	329398	264872	3.72%	0.558%
27	10.851	1489	1490	1497	rVB	257407	237268	3.33%	0.500%
28	11.286	1561	1564	1566	rBV	301488	228354	3.21%	0.481%
29	11.316	1566	1569	1574	rVB	278232	296188	4.16%	0.624%
30	11.410	1582	1585	1588	rBV	1520987	1393655	19.58%	2.936%
31	11.433	1588	1589	1592	rVB	192107	122668	1.72%	0.258%
32	11.710	1633	1636	1638	rBV	236065	209484	2.94%	0.441%
33	11.733	1638	1640	1643	rVB2	104122	89130	1.25%	0.188%
34	11.810	1649	1653	1656	rBV	3179377	2431006	34.16%	5.121%

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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-BF121918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.939	1672	1675	1678	rVB	171547	157270	2.21%	0.331%
36	12.121	1703	1706	1709	rBV	218952	201334	2.83%	0.424%
37	12.210	1719	1721	1730	rVB6	69542	92139	1.29%	0.194%
38	12.498	1766	1770	1772	rBV	7013343	7117536	100.00%	14.992%
39	12.521	1772	1774	1780	rVV	8195664	7024679	98.70%	14.796%
40	12.604	1784	1788	1790	rBV	1028689	989452	13.90%	2.084%
41	12.621	1790	1791	1793	rVV	364334	224515	3.15%	0.473%
42	12.645	1793	1795	1798	rVB	139886	125421	1.76%	0.264%
43	12.851	1825	1830	1832	rBV	347685	356272	5.01%	0.750%
44	12.880	1832	1835	1837	rVB	168805	141799	1.99%	0.299%
45	12.992	1851	1854	1858	rBV	2015905	1787058	25.11%	3.764%
46	13.245	1892	1897	1900	rBV2	199053	303975	4.27%	0.640%
47	13.327	1908	1911	1915	rVB	111089	101457	1.43%	0.214%
48	13.574	1951	1953	1956	rBV	148910	130223	1.83%	0.274%
49	13.892	2004	2007	2013	rBV3	170743	279873	3.93%	0.590%
50	13.992	2022	2024	2028	rBV	147230	152512	2.14%	0.321%
51	14.045	2029	2033	2036	rBV	1607910	1565220	21.99%	3.297%
52	14.221	2060	2063	2067	rVB2	169071	169853	2.39%	0.358%
53	14.527	2112	2115	2117	rVB	179098	158501	2.23%	0.334%
54	14.833	2165	2167	2170	rBV	150973	150849	2.12%	0.318%
55	15.521	2280	2284	2287	rVB	735623	837565	11.77%	1.764%

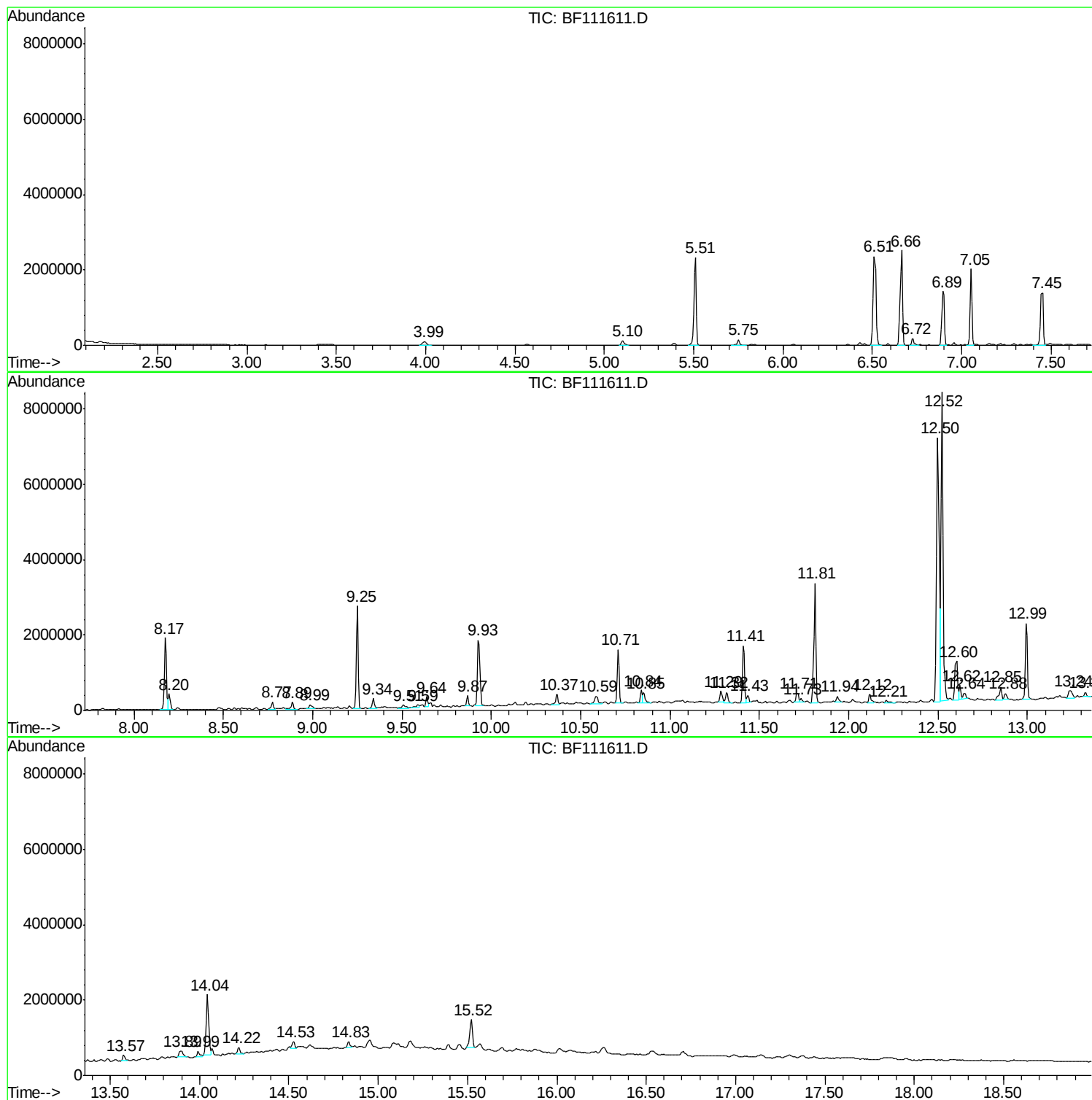
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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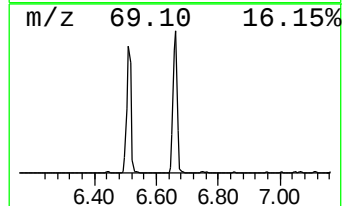
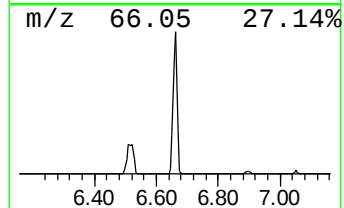
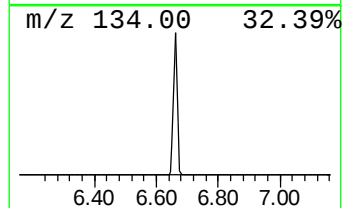
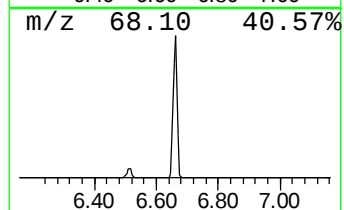
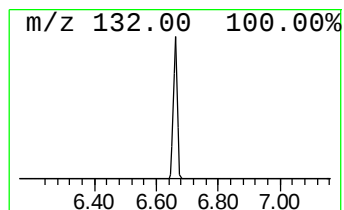
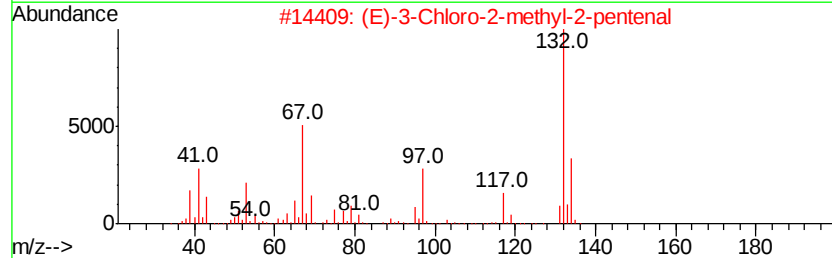
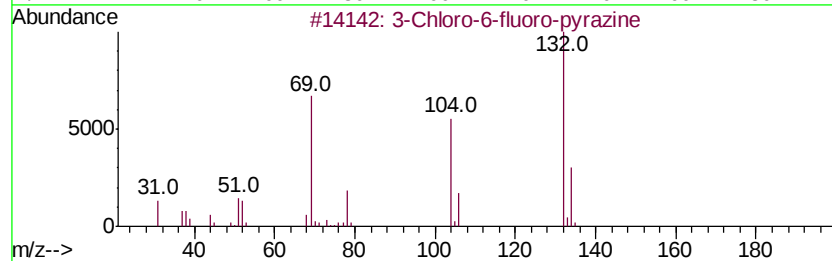
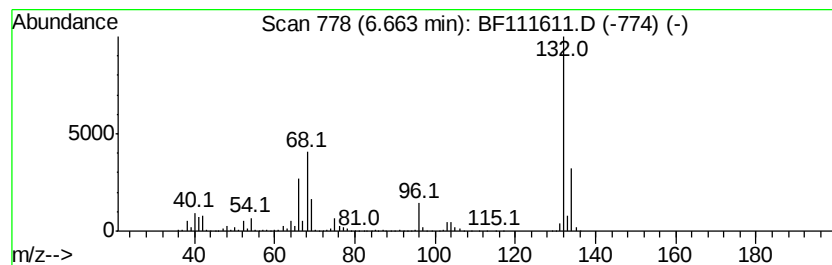
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.66 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	34.37 ng	2135630	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Xenon	132	Xe	007440-63-3	9



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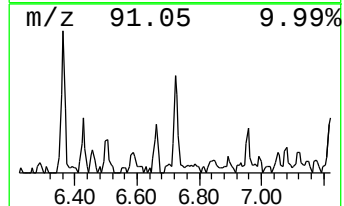
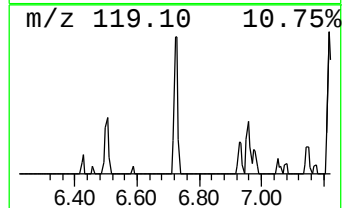
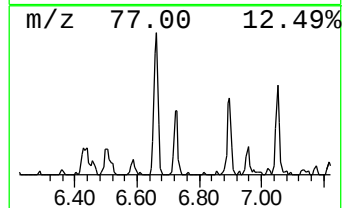
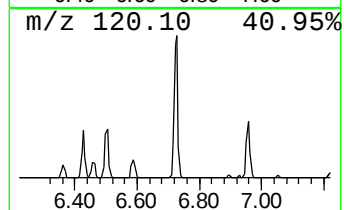
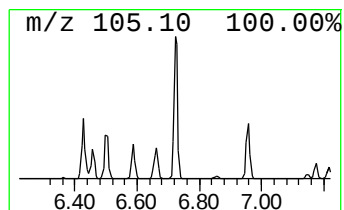
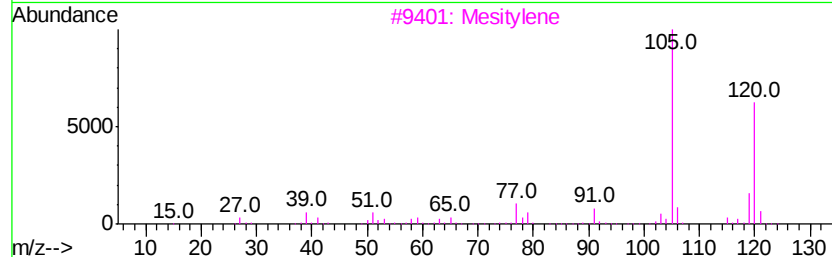
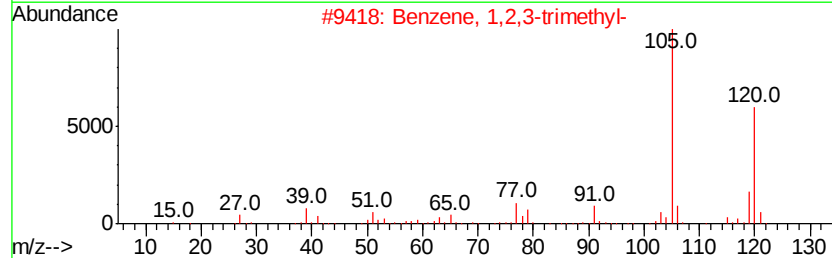
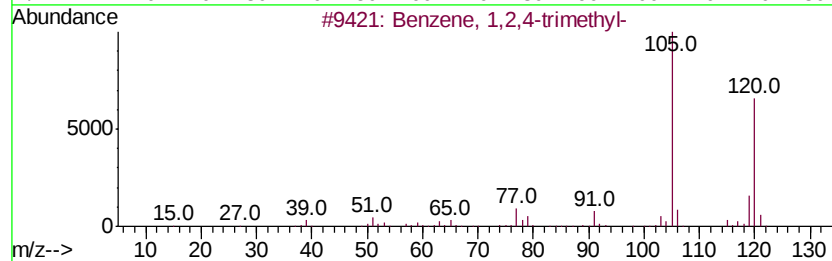
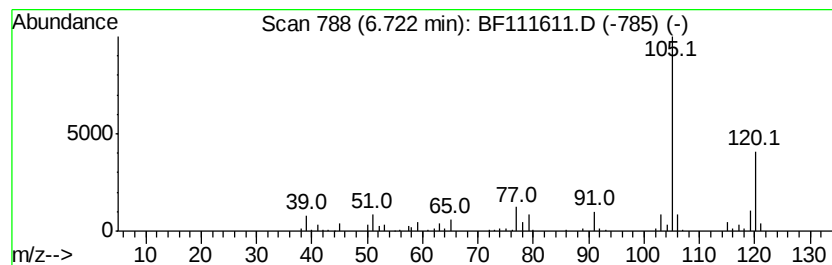
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2,4-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.72	2.76 ng	171275	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3		Mesitylene	120	C9H12	000108-67-8	94
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87



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ALS Vial : 15 Sample Multiplier: 1

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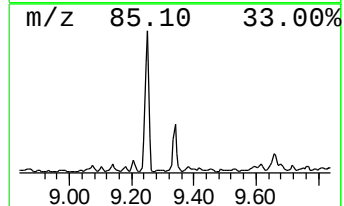
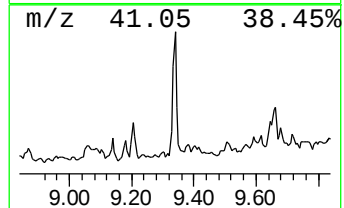
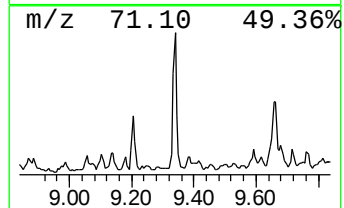
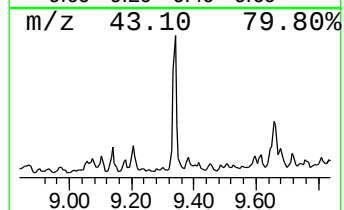
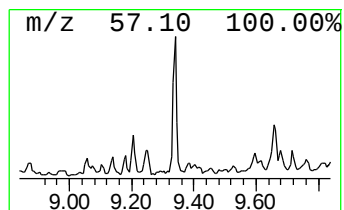
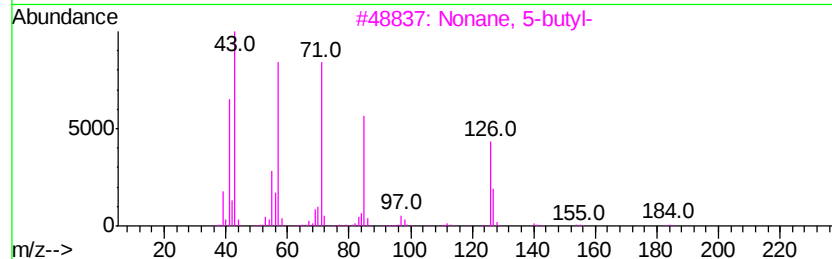
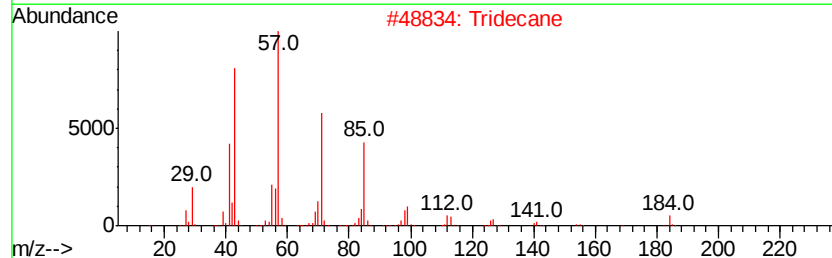
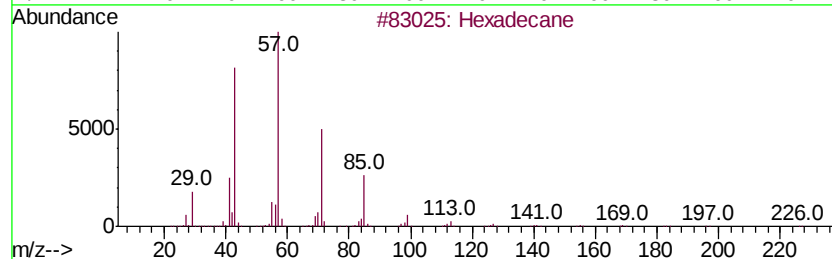
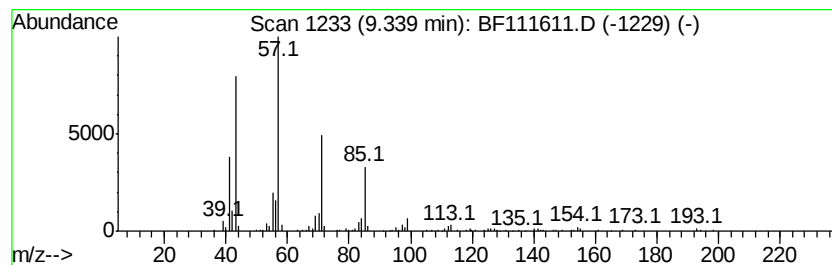
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	3.06 ng	244063	Acenaphthene-d10	9.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	90
2		Tridecane	184	C13H28	000629-50-5	83
3		Nonane, 5-butyl-	184	C13H28	017312-63-9	72
4		Decane, 3-bromo-	220	C10H21Br	030571-71-2	72
5		Undecane, 3-methyl-	170	C12H26	001002-43-3	72



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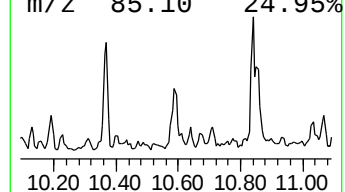
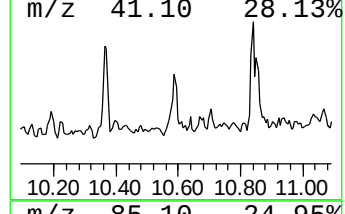
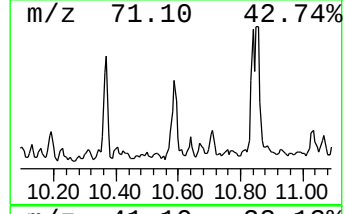
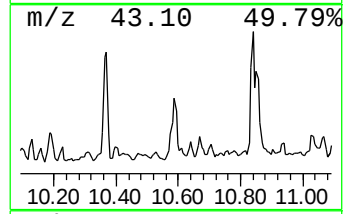
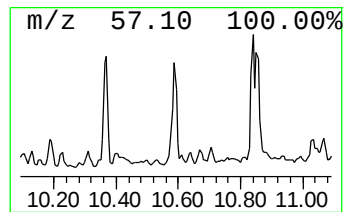
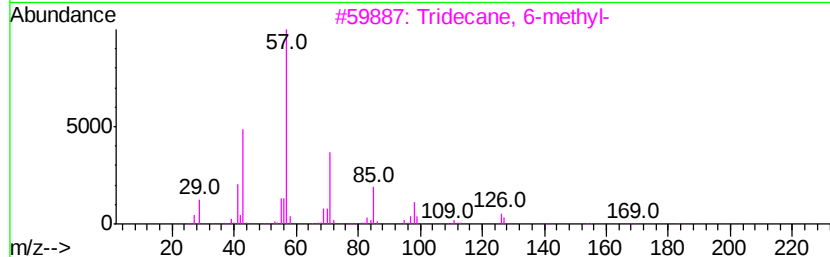
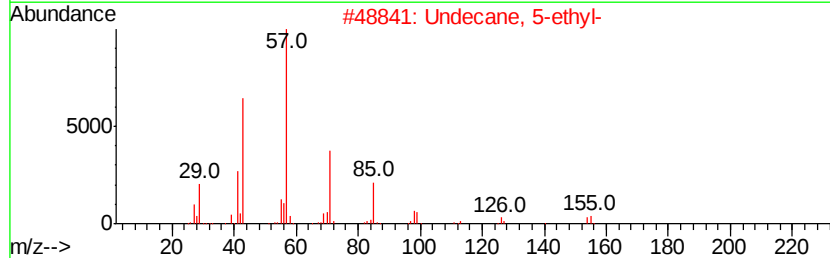
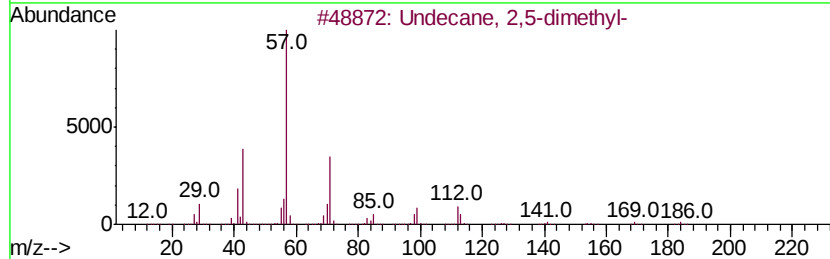
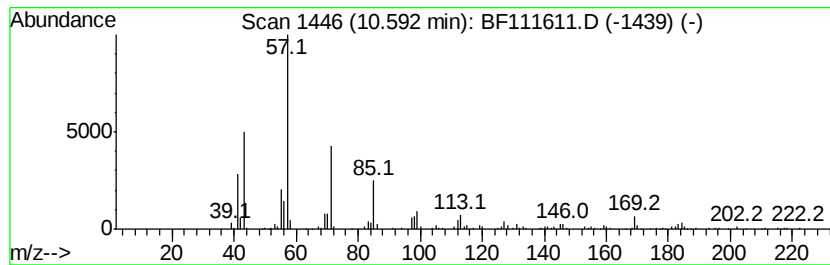
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Undecane, 2,5-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.59	3.84 ng	305656	Acenaphthene-d10	9.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	81
2	Undecane, 5-ethyl-	184	C13H28	017453-94-0	72
3	Tridecane, 6-methyl-	198	C14H30	013287-21-3	72
4	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	68
5	Heptacosane	380	C27H56	000593-49-7	64



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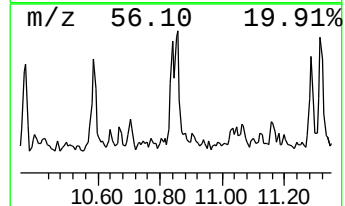
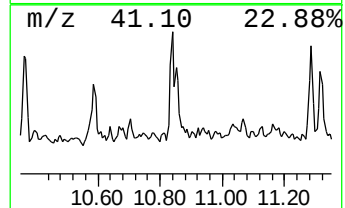
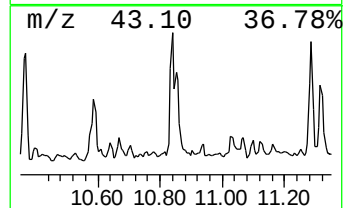
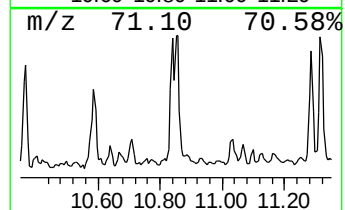
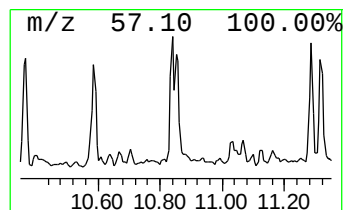
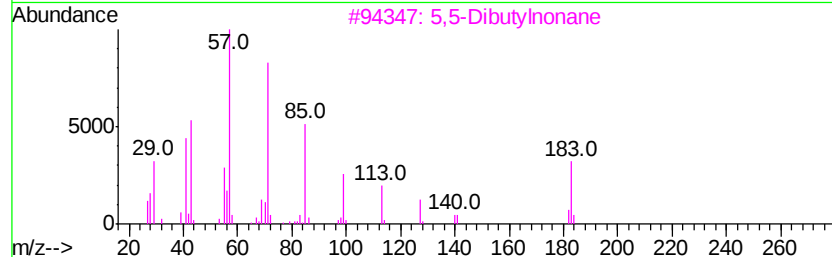
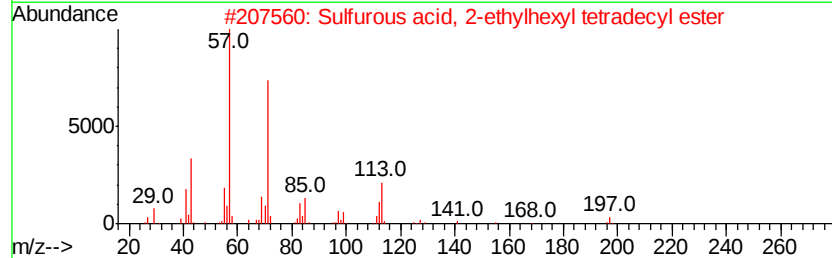
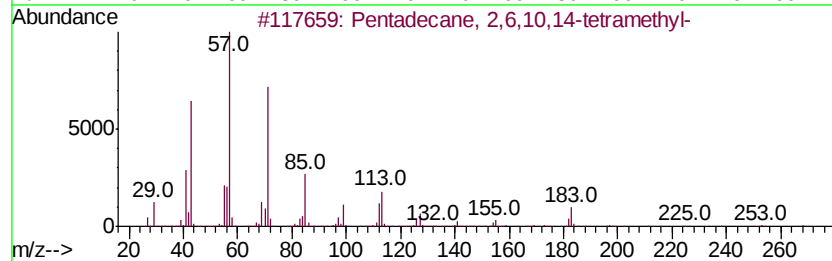
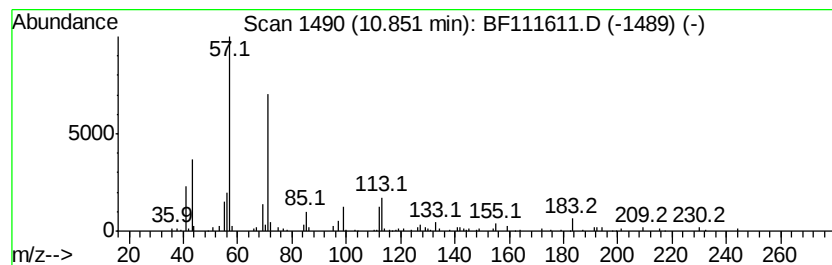
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecane, 2,6,10,14-tetr... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	3.40 ng	237268	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72
2		Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	64
3		5,5-Dibutylnonane	240	C17H36	006008-17-9	59
4		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	59
5		Tetradecane	198	C14H30	000629-59-4	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
Data File : BF111611.D
Acq On : 20 Dec 2018 00:57
Operator : JU/SJ
Sample : J6196-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-01-121818-C

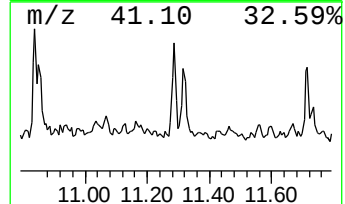
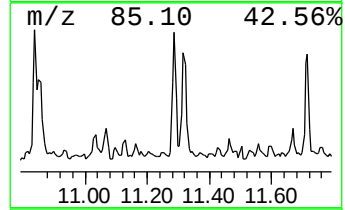
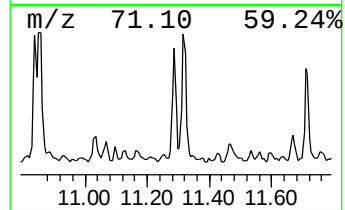
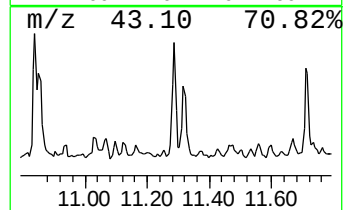
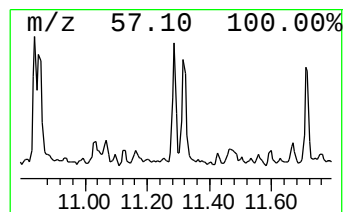
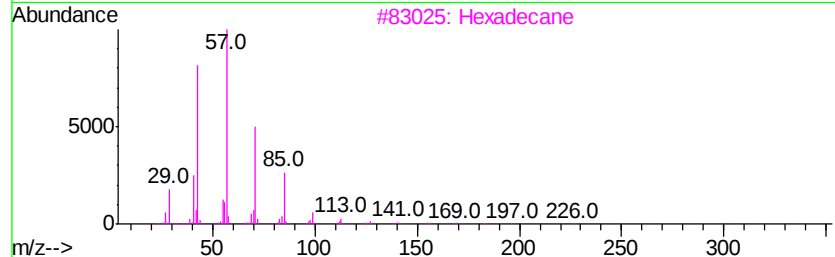
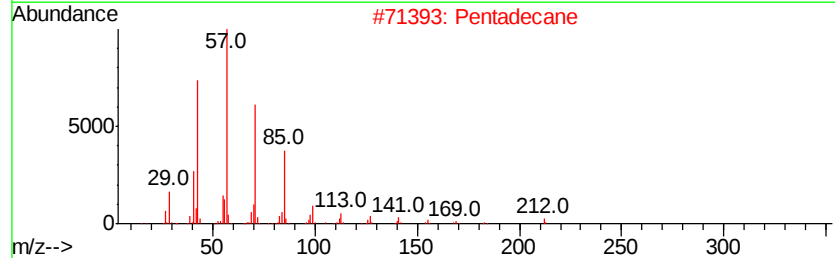
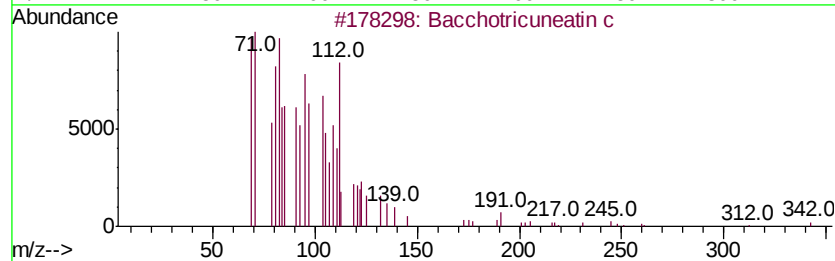
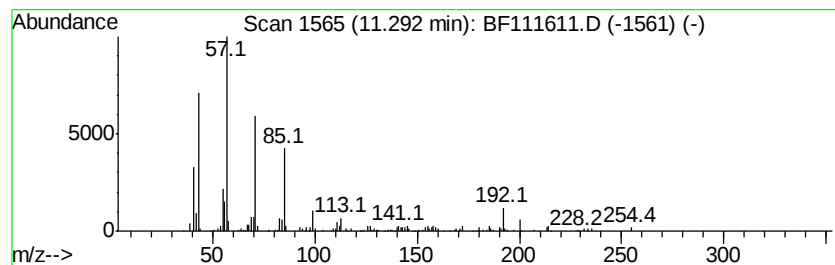
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Bacchotricuneatin c Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.29	3.28 ng	228354	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bacchotricuneatin c	342	C20H22O5	066563-30-2	99
2		Pentadecane	212	C15H32	000629-62-9	87
3		Hexadecane	226	C16H34	000544-76-3	86
4		Sulfurous acid, 2-propyl tridecy...	306	C16H34O3S	1000309-12-4	80
5		Eicosane	282	C20H42	000112-95-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
Data File : BF111611.D
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Sample : J6196-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

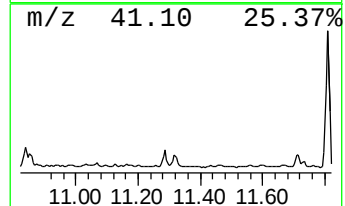
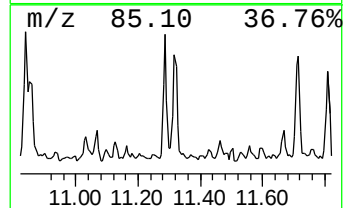
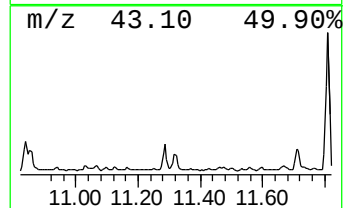
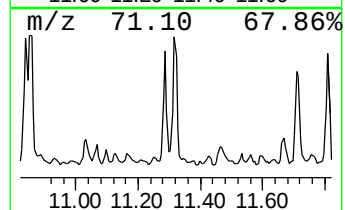
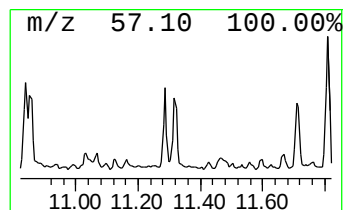
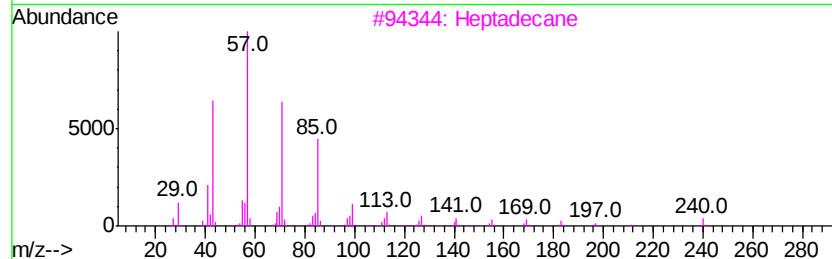
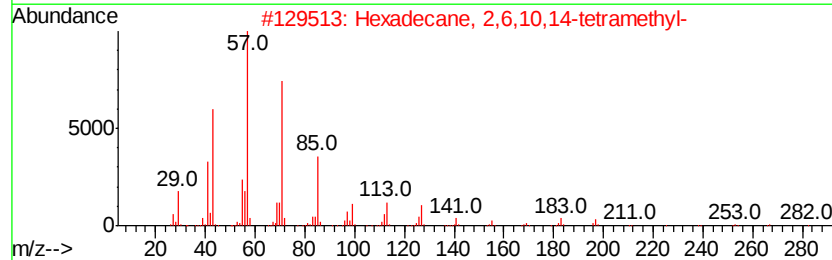
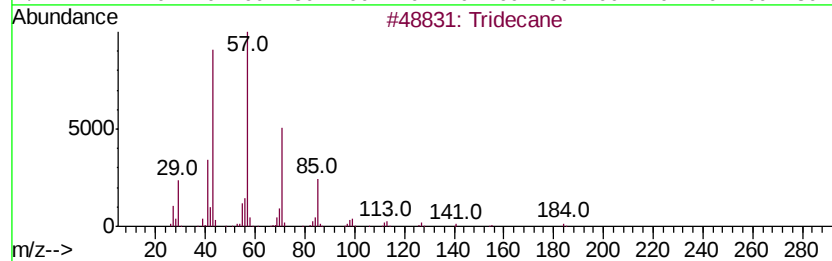
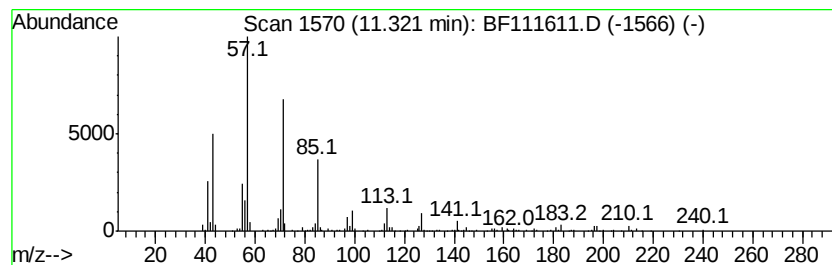
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.32	4.25 ng	296188	Phenanthrene-d10		11.41
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	91
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	87
3	Heptadecane	240	C17H36	000629-78-7	86
4	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	86
5	Heptacosane	380	C27H56	000593-49-7	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
Data File : BF111611.D
Acq On : 20 Dec 2018 00:57
Operator : JU/SJ
Sample : J6196-11 2X
Misc :
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Instrument :
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ClientSampleId :
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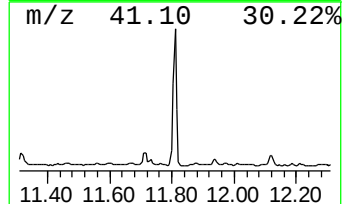
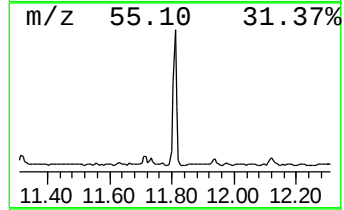
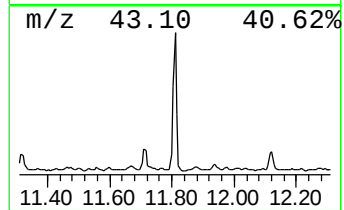
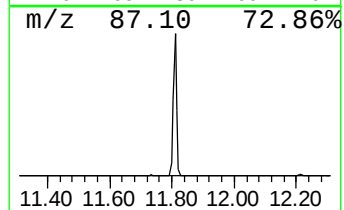
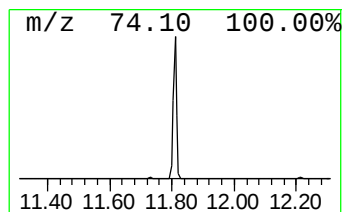
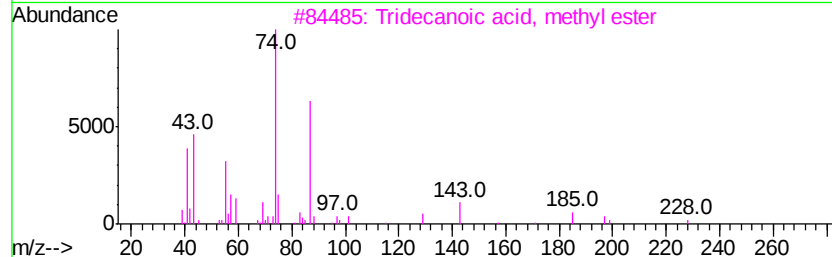
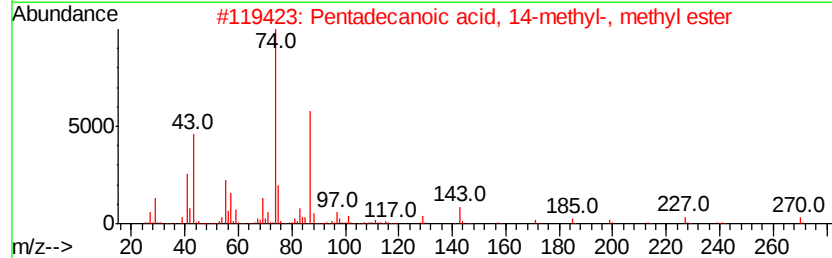
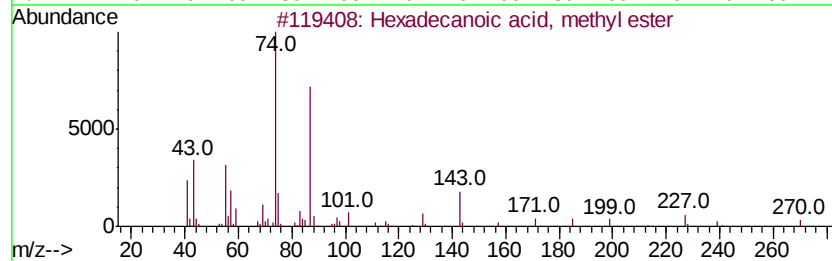
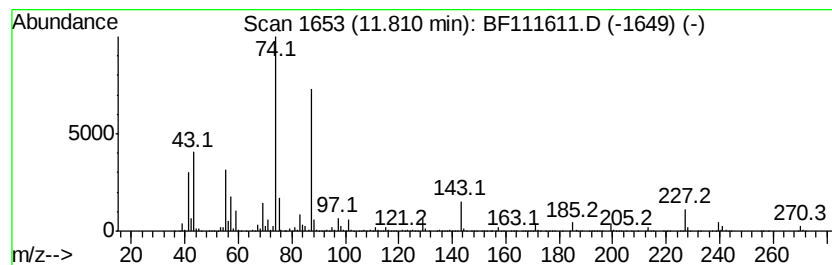
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	34.89 ng	2431010	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
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Sample : J6196-11 2X
Misc :
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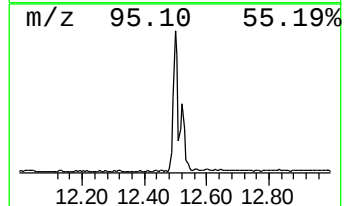
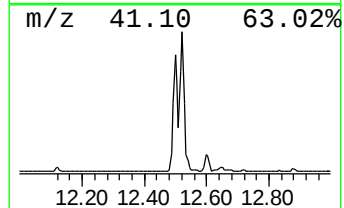
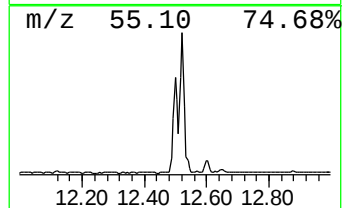
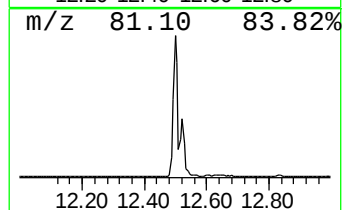
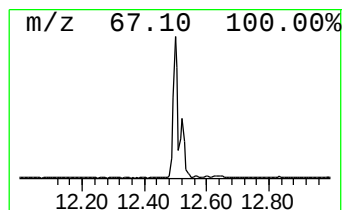
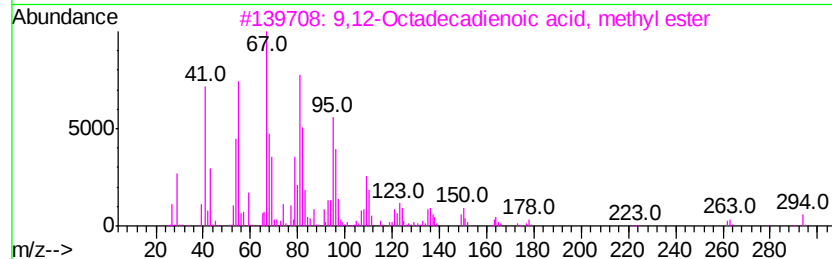
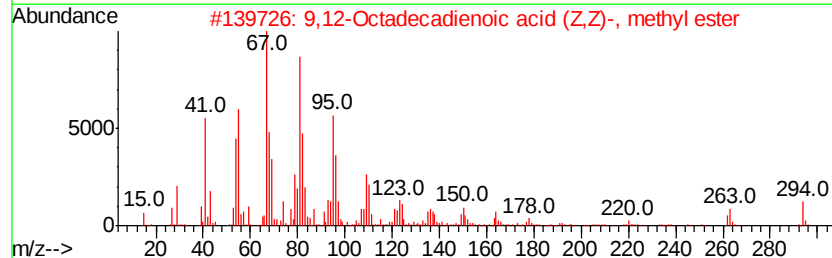
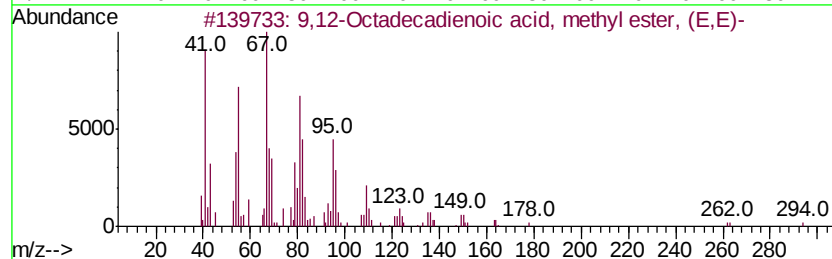
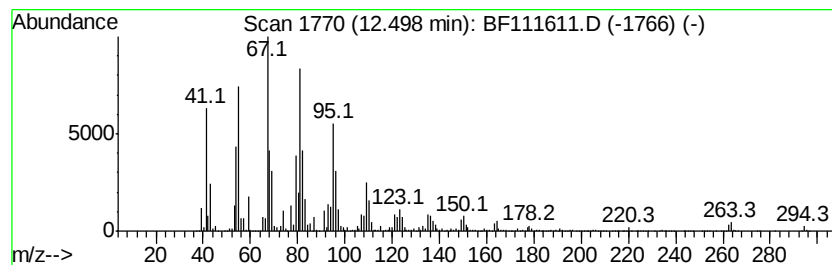
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.50	102.14 ng	7117540	Phenanthrene-d10	11.41		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99	
2	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99	
3	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99	
4	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99	
5	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
Data File : BF111611.D
Acq On : 20 Dec 2018 00:57
Operator : JU/SJ
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Misc :
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Instrument :
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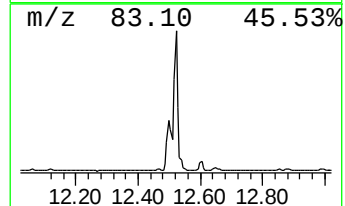
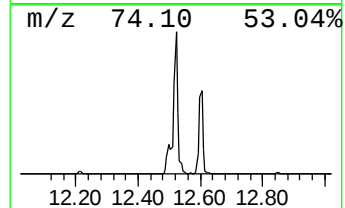
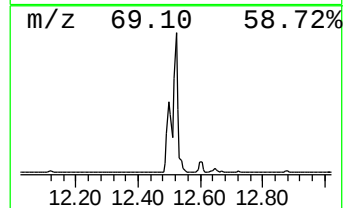
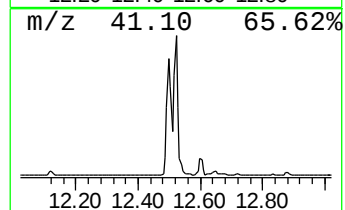
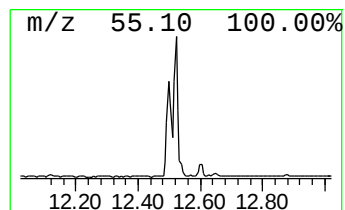
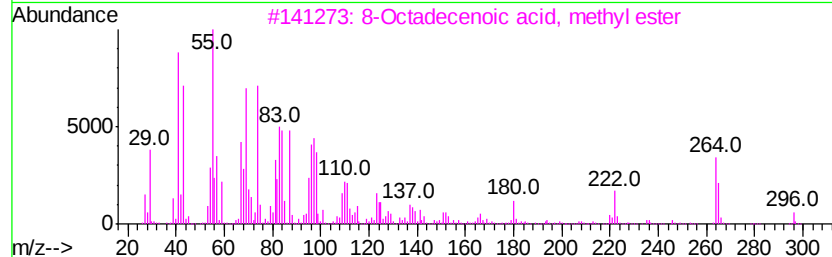
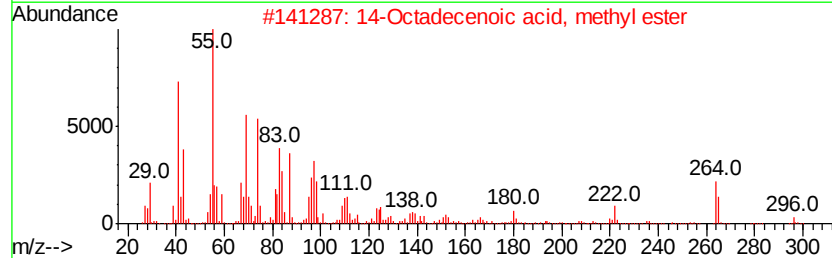
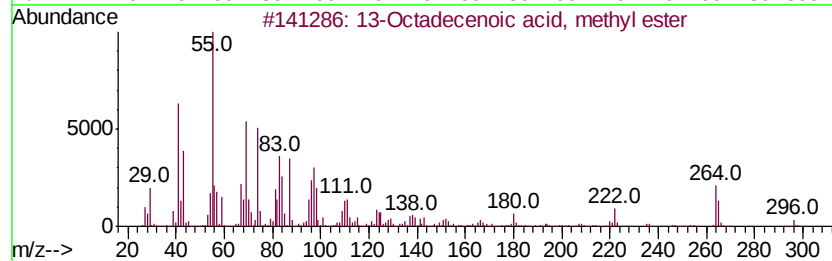
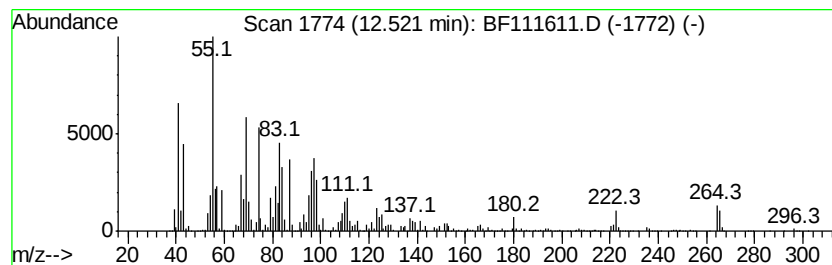
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 13-Octadecenoic acid, methyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	100.81 ng	7024680	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
2		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	99
3		8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99
4		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
5		6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
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Misc :
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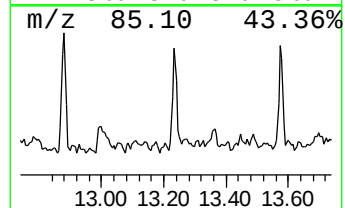
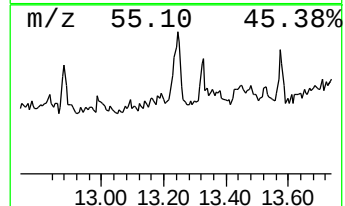
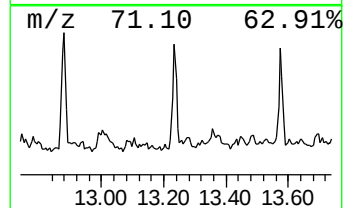
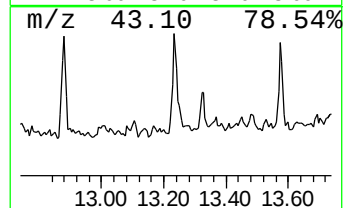
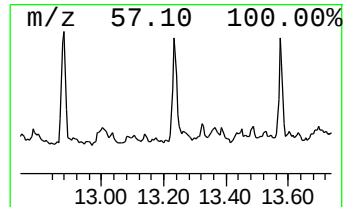
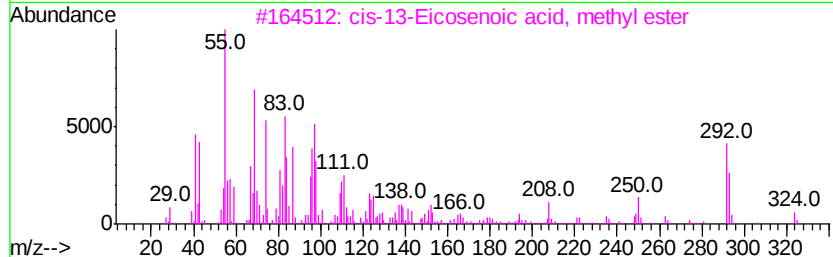
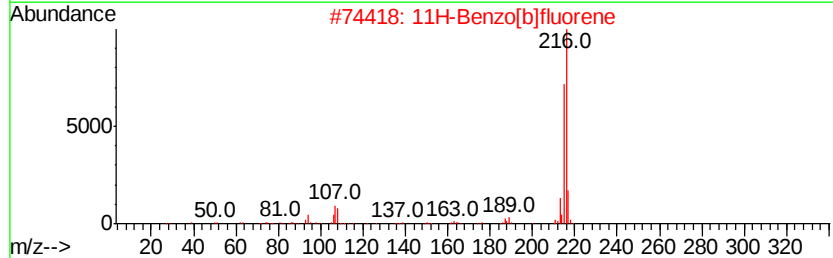
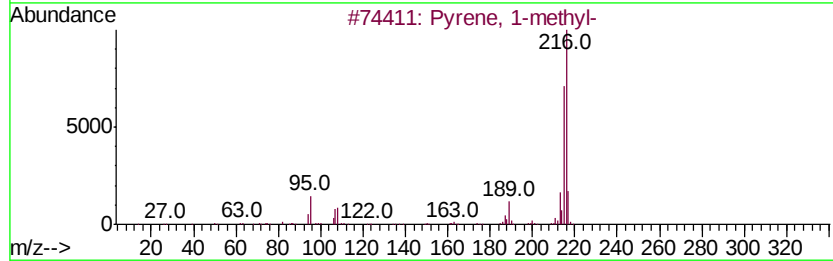
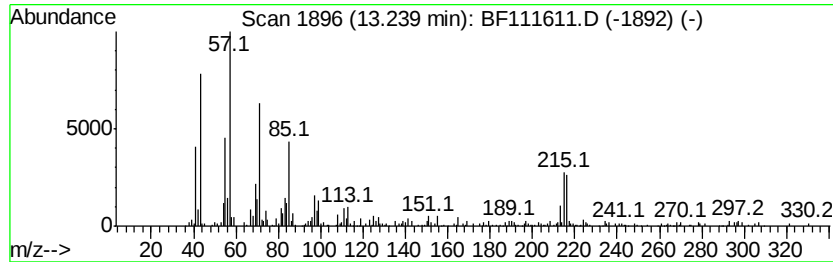
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	3.88 ng	303975	Chrysene-d12	14.04
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene, 1-methyl-	216 C17H12	002381-21-7 64
2		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 60
3		cis-13-Eicosenoic acid, methyl e...	324 C21H40O2	1000333-52-1 59
4		11-Eicosenoic acid, methyl ester	324 C21H40O2	003946-08-5 53
5		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
Data File : BF111611.D
Acq On : 20 Dec 2018 00:57
Operator : JU/SJ
Sample : J6196-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-121818-C

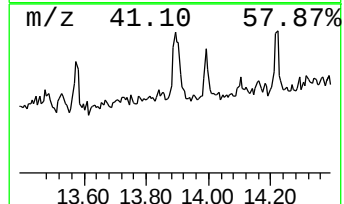
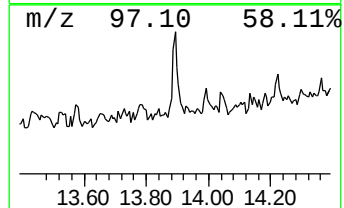
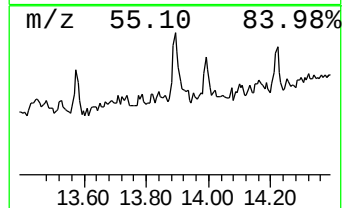
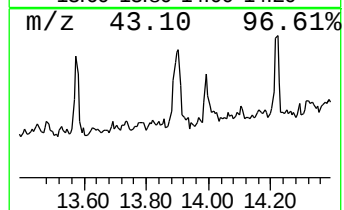
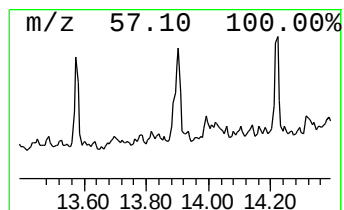
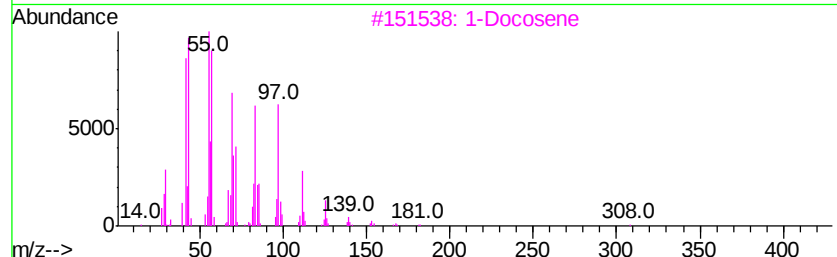
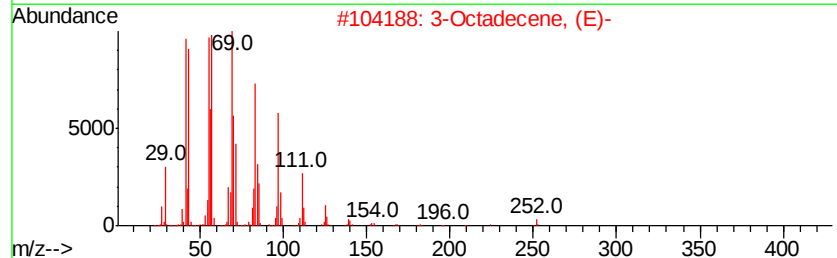
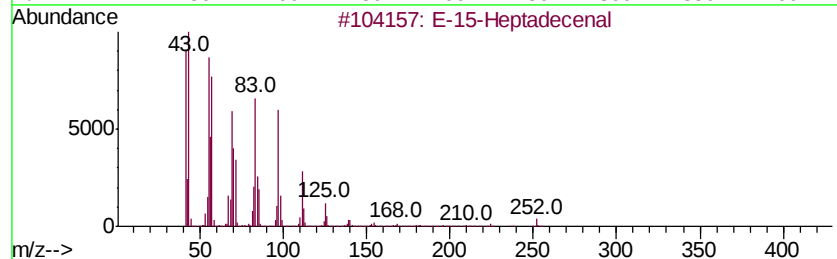
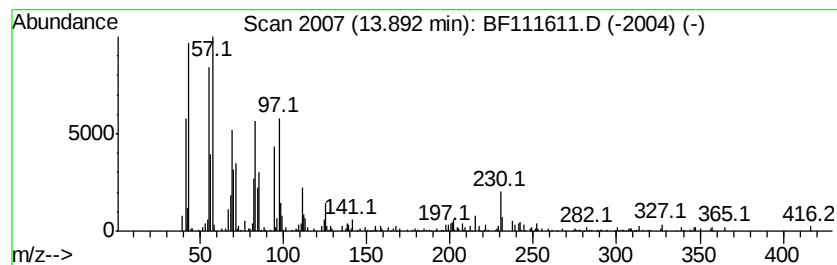
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 E-15-Heptadecenal Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	3.58 ng	279873	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	E-15-Heptadecenal	252	C17H32O	1000130-97-9	91
2		3-Octadecene, (E)-	252	C18H36	007206-19-1	83
3		1-Docosene	308	C22H44	001599-67-3	83
4		10-Heneicosene (c,t)	294	C21H42	095008-11-0	83
5		1-Octadecene	252	C18H36	000112-88-9	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF121918\
Data File : BF111611.D
Acq On : 20 Dec 2018 00:57
Operator : JU/SJ
Sample : J6196-11 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-121818-C

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.66	6.66	34.4	ng	2135630	1	6.90	1242660	20.0
Benzene, 1,2,4-tr...	6.72	2.8	ng	171275	1	6.90	1242660	20.0
Hexadecane	9.34	3.1	ng	244063	3	9.93	1592600	20.0
Undecane, 2,5-dim...	10.59	3.8	ng	305656	3	9.93	1592600	20.0
Pentadecane, 2,6,...	10.85	3.4	ng	237268	4	11.41	1393660	20.0
Bacchotricuneatin c	11.29	3.3	ng	228354	4	11.41	1393660	20.0
Tridecane	11.32	4.3	ng	296188	4	11.41	1393660	20.0
Hexadecanoic acid...	11.81	34.9	ng	2431010	4	11.41	1393660	20.0
9,12-Octadecadien...	12.50	102.1	ng	7117540	4	11.41	1393660	20.0
13-Octadecenoic a...	12.52	100.8	ng	7024680	4	11.41	1393660	20.0
Pyrene, 1-methyl-	13.24	3.9	ng	303975	5	14.04	1565220	20.0
E-15-Heptadecenal	13.89	3.6	ng	279873	5	14.04	1565220	20.0