

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
 Data File : BF103940.D
 Acq On : 26 Mar 2018 22:07
 Operator : JU/SJ
 Sample : J1760-07 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-032318-B

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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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	2.287	30	34	49	rVB	67132	103091	1.08%	0.209%
2	5.593	592	596	612	rBV	1793724	1612382	16.90%	3.270%
3	6.593	761	766	779	rBV	1769867	1672421	17.53%	3.392%
4	6.740	787	791	795	rBV	1985080	1697911	17.80%	3.444%
5	6.975	827	831	834	rBV	1034204	831511	8.72%	1.686%
6	7.128	853	857	861	rBV	1548996	1278279	13.40%	2.593%
7	7.534	922	926	929	rBV	1186090	982311	10.30%	1.992%
8	8.257	1045	1049	1051	rBV	1347355	1063278	11.15%	2.157%
9	9.069	1182	1187	1190	rBV2	144181	155528	1.63%	0.315%
10	9.328	1227	1231	1235	rBV	2102517	1715436	17.98%	3.479%
11	9.398	1240	1243	1246	rVB	156895	136609	1.43%	0.277%
12	9.586	1272	1275	1281	rVB5	83099	124603	1.31%	0.253%
13	9.669	1281	1289	1291	rBV2	139034	206468	2.16%	0.419%
14	9.716	1295	1297	1302	rVV	201002	239801	2.51%	0.486%
15	9.786	1305	1309	1314	rVB3	87855	118579	1.24%	0.241%
16	9.875	1320	1324	1327	rBV2	127407	121921	1.28%	0.247%
17	9.928	1330	1333	1337	rVB	284315	220767	2.31%	0.448%
18	10.016	1345	1348	1351	rVV	1358835	1090689	11.43%	2.212%
19	10.045	1351	1353	1356	rVB	169551	140081	1.47%	0.284%
20	10.216	1379	1382	1385	rVB2	126993	120522	1.26%	0.244%
21	10.251	1385	1388	1392	rVB3	112126	105033	1.10%	0.213%
22	10.328	1397	1401	1404	rBV3	95181	118762	1.25%	0.241%
23	10.428	1412	1418	1421	rBV	383369	352363	3.69%	0.715%
24	10.563	1438	1441	1448	rVB	217990	223265	2.34%	0.453%
25	10.645	1451	1455	1458	rBV3	125887	162393	1.70%	0.329%
26	10.804	1479	1482	1486	rBV	1085176	993555	10.42%	2.015%
27	10.904	1496	1499	1508	rVB	363541	457199	4.79%	0.927%
28	11.110	1529	1534	1536	rBV4	64320	112340	1.18%	0.228%
29	11.351	1572	1575	1578	rBV	316629	243249	2.55%	0.493%
30	11.380	1578	1580	1582	rVV2	114033	106462	1.12%	0.216%
31	11.510	1598	1602	1604	rBV	1123543	1020793	10.70%	2.070%
32	11.533	1604	1606	1610	rVV	1372009	1085592	11.38%	2.202%
33	11.586	1612	1615	1618	rVV	276971	257106	2.70%	0.521%
34	11.780	1645	1648	1650	rBV	219067	173703	1.82%	0.352%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

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:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.886	1662	1666	1669	rBV	4052175	3604754	37.79%	7.311%
36	12.010	1684	1687	1690	rBV	290014	278238	2.92%	0.564%
37	12.039	1690	1692	1695	rVB	332761	263181	2.76%	0.534%
38	12.122	1703	1706	1709	rBV2	271436	341779	3.58%	0.693%
39	12.145	1709	1710	1713	rVB	173629	106065	1.11%	0.215%
40	12.186	1714	1717	1719	rBV	175423	157803	1.65%	0.320%
41	12.586	1778	1785	1787	rBV	5502927	9538719	100.00%	19.346%
42	12.610	1787	1789	1794	rVV2	5419371	5459900	57.24%	11.074%
43	12.651	1794	1796	1798	rVV2	118964	118443	1.24%	0.240%
44	12.680	1798	1801	1804	rVB	1977860	1605528	16.83%	3.256%
45	12.727	1806	1809	1812	rBV	1080384	1016618	10.66%	2.062%
46	12.916	1838	1841	1843	rBV2	273987	234249	2.46%	0.475%
47	12.963	1843	1849	1855	rVB	1125483	1275879	13.38%	2.588%
48	13.098	1868	1872	1874	rBV	1313355	1165565	12.22%	2.364%
49	13.174	1882	1885	1891	rBV5	73012	159438	1.67%	0.323%
50	13.239	1893	1896	1898	rBV	309523	294252	3.08%	0.597%
51	13.310	1906	1908	1909	rVV	150262	117922	1.24%	0.239%
52	13.327	1909	1911	1913	rVV2	196424	154327	1.62%	0.313%
53	13.351	1913	1915	1918	rVB	121220	111803	1.17%	0.227%
54	13.404	1920	1924	1926	rBV2	230452	256921	2.69%	0.521%
55	13.516	1941	1943	1946	rBV	101218	95848	1.00%	0.194%
56	14.074	2035	2038	2042	rBV2	175498	173455	1.82%	0.352%
57	14.163	2048	2053	2056	rBV2	967961	1290059	13.52%	2.616%
58	14.186	2056	2057	2063	rVB	380461	341548	3.58%	0.693%
59	15.068	2203	2207	2216	rVB	252872	417218	4.37%	0.846%
60	15.251	2234	2238	2241	rBV	255713	409352	4.29%	0.830%
61	15.574	2290	2293	2297	rVB	183128	230587	2.42%	0.468%
62	15.639	2301	2304	2311	rVB	264789	379752	3.98%	0.770%
63	15.710	2312	2316	2320	rBV	497169	661857	6.94%	1.342%

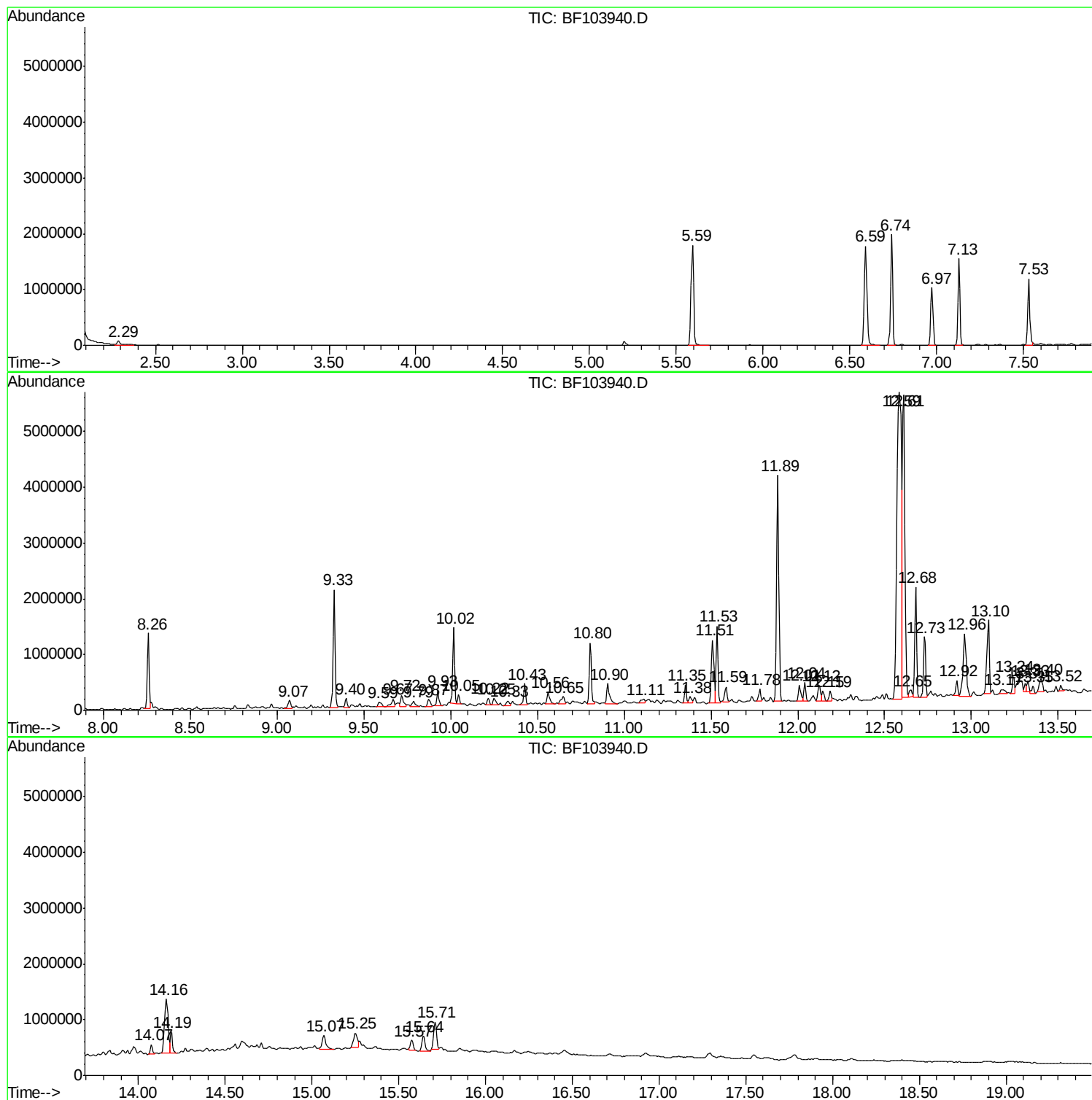
Sum of corrected areas: 49305063

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

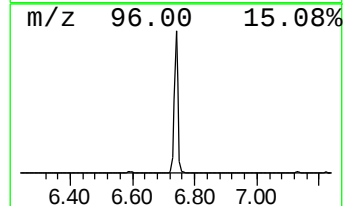
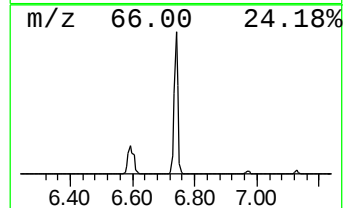
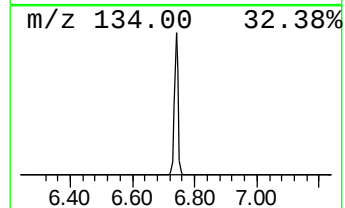
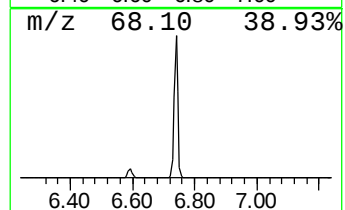
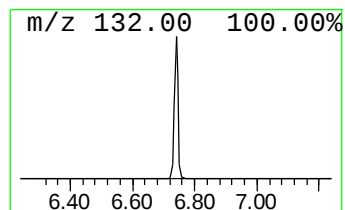
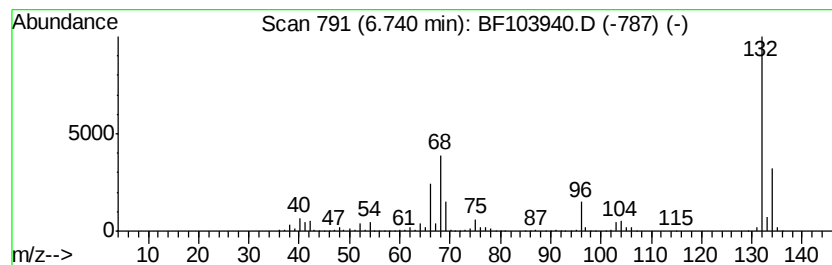
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.74 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	40.84 ng	1697910	1,4-Dichlorobenzene-d4	6.97

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	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

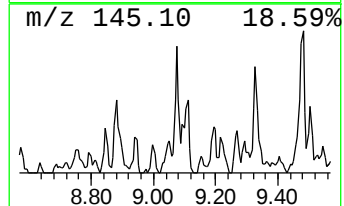
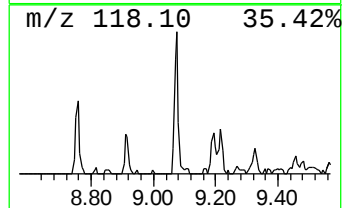
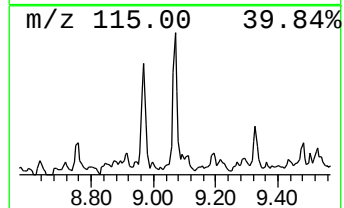
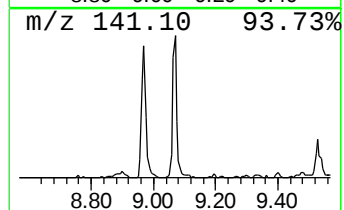
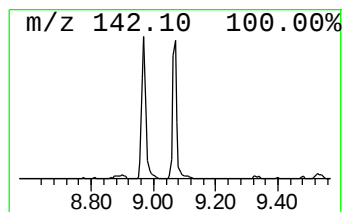
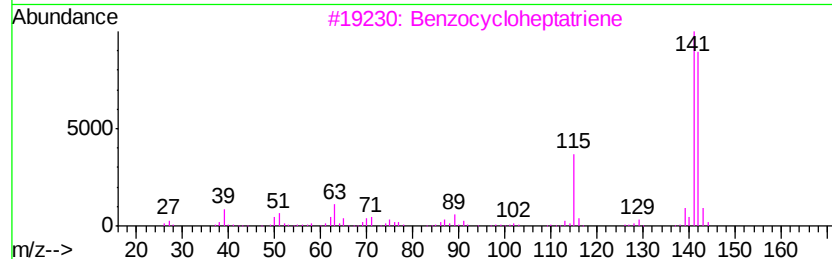
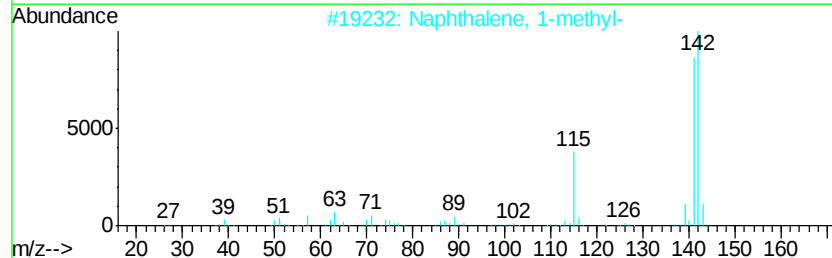
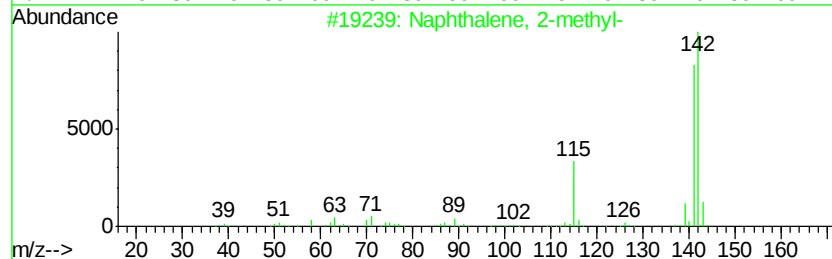
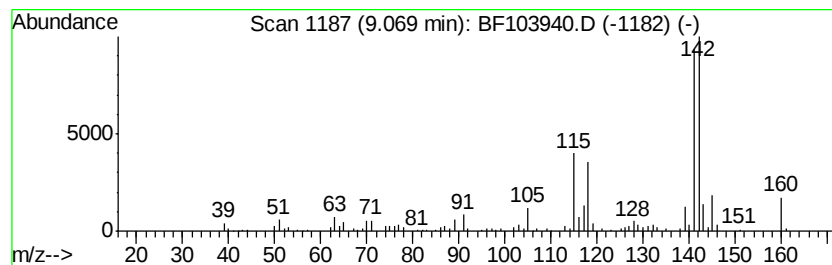
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Naphthalene, 2-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.07	2.93 ng	155528	Naphthalene-d8	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	81
4		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	81
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

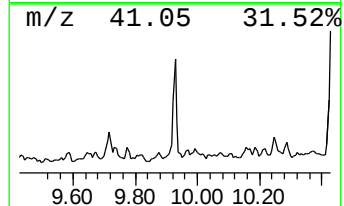
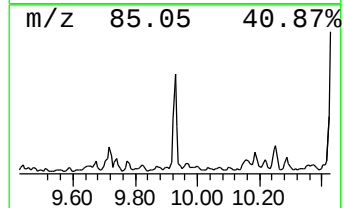
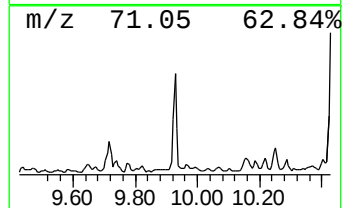
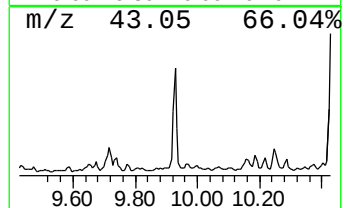
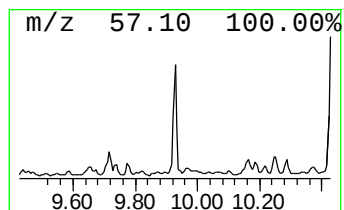
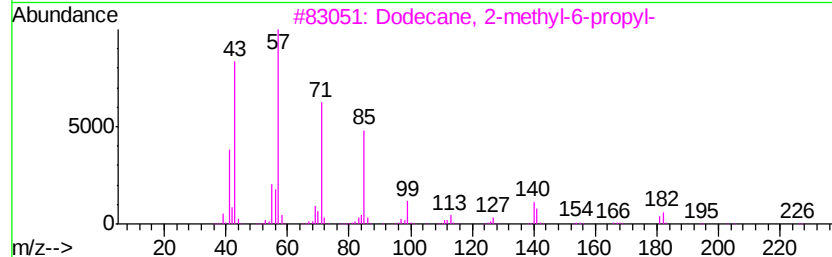
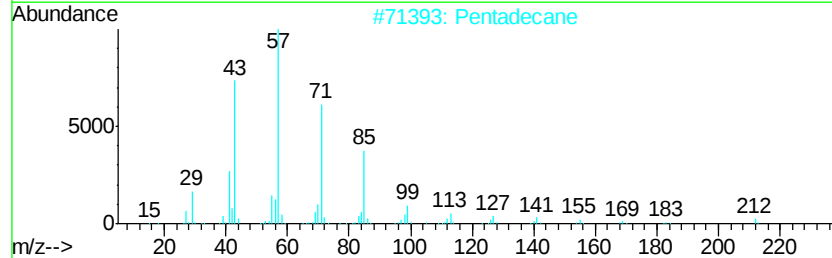
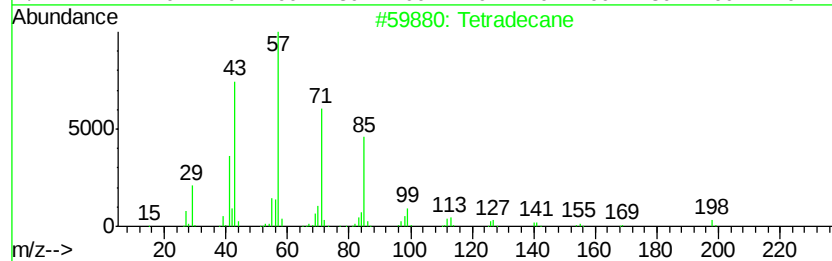
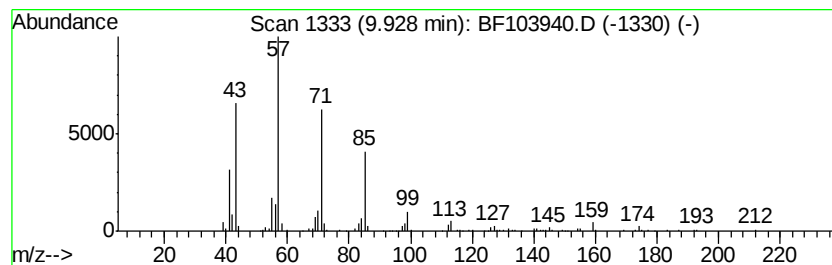
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Tetradecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	4.05 ng	220767	Acenaphthene-d10	10.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	87
2			Pentadecane	212	C15H32	000629-62-9	78
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	78
4			Hexadecane	226	C16H34	000544-76-3	72
5			Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

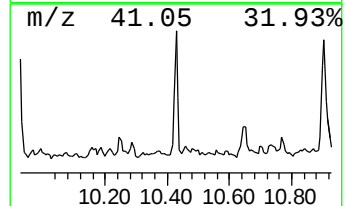
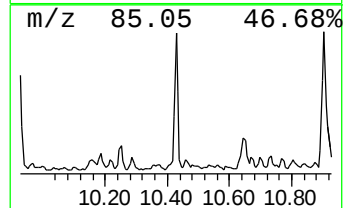
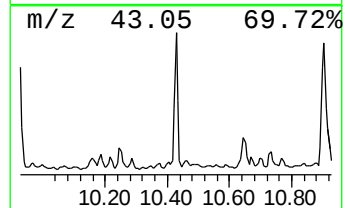
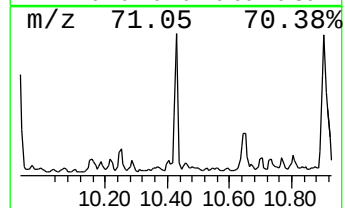
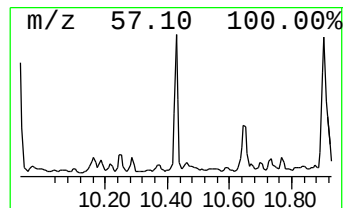
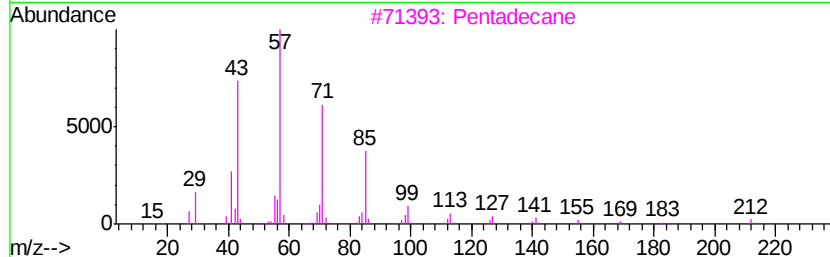
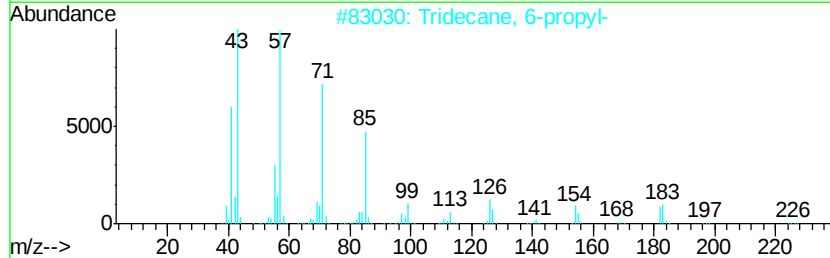
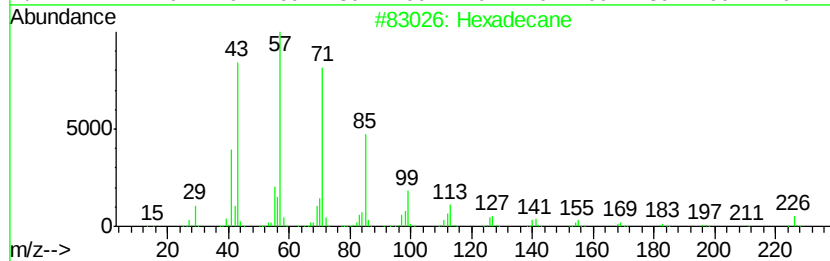
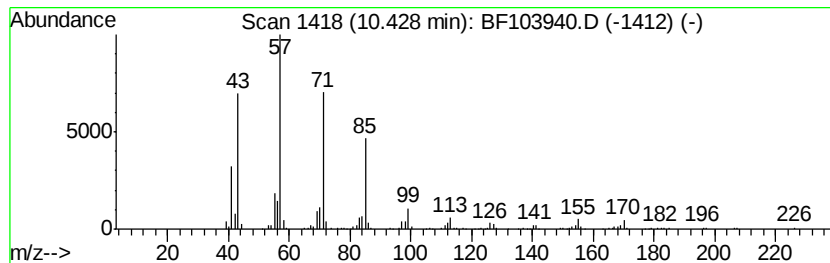
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	6.46 ng	352363	Acenaphthene-d10	10.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	93
2	Tridecane, 6-propyl-		226	C16H34	055045-10-8	86
3	Pentadecane		212	C15H32	000629-62-9	86
4	Dodecane		170	C12H26	000112-40-3	83
5	Tetradecane		198	C14H30	000629-59-4	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

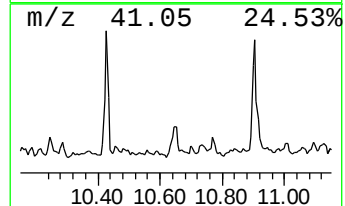
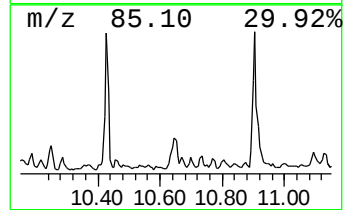
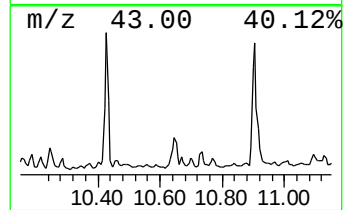
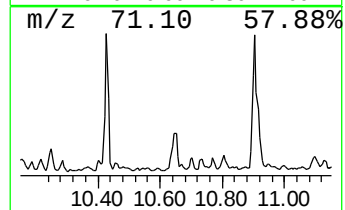
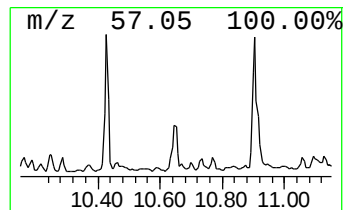
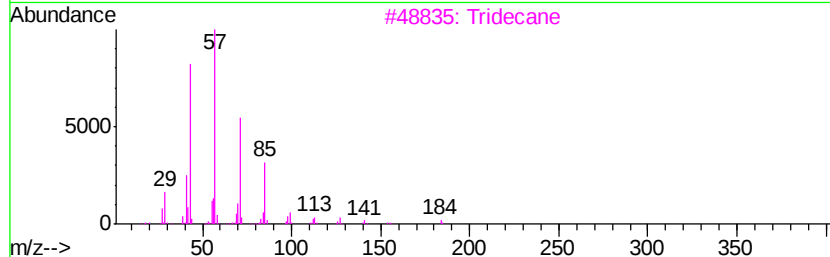
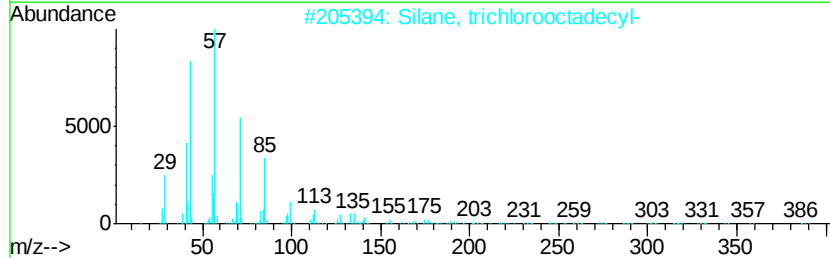
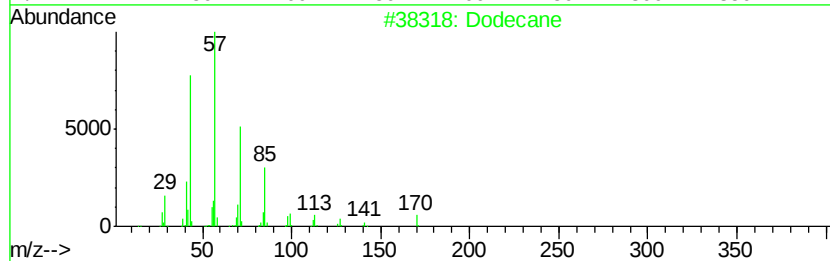
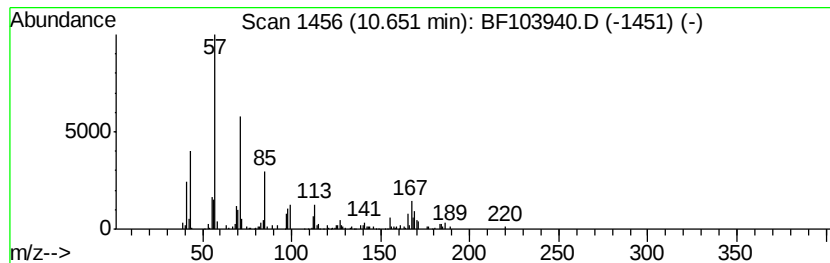
Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dodecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	2.98 ng	162393	Acenaphthene-d10	10.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	90
2	Silane, trichlorooctadecyl-	386 C18H37Cl3Si	000112-04-9	72
3	Tridecane	184 C13H28	000629-50-5	70
4	Dodecane, 1-iodo-	296 C12H25I	004292-19-7	64
5	Eicosane	282 C20H42	000112-95-8	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

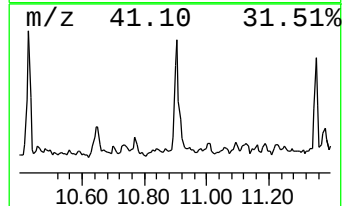
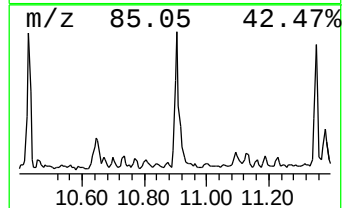
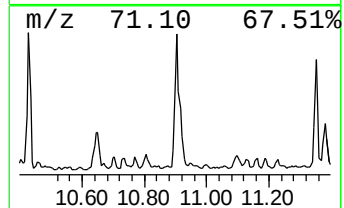
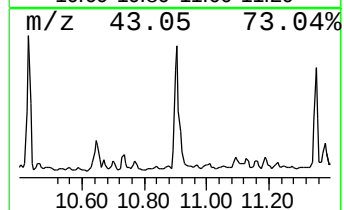
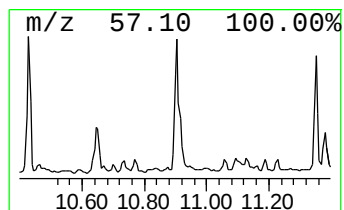
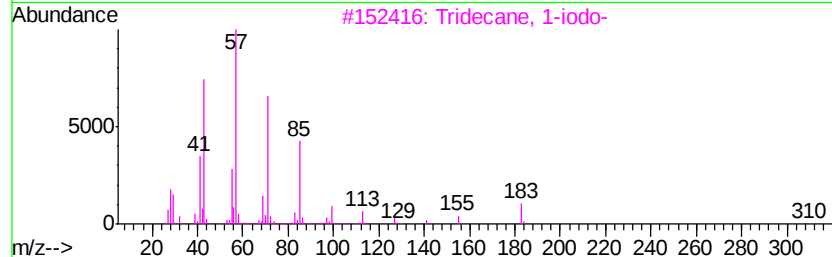
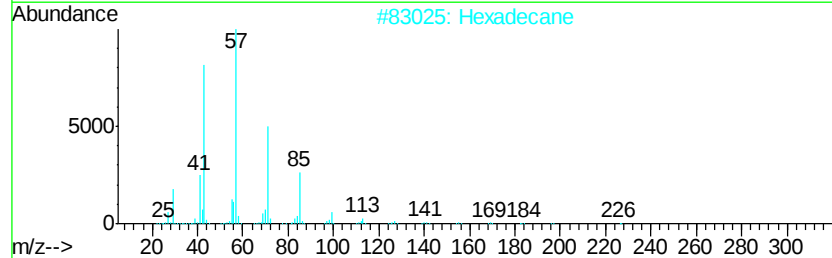
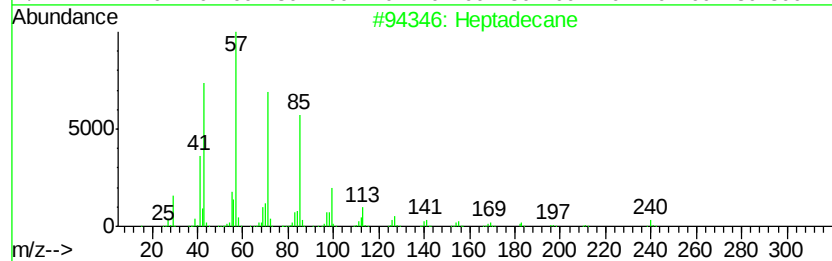
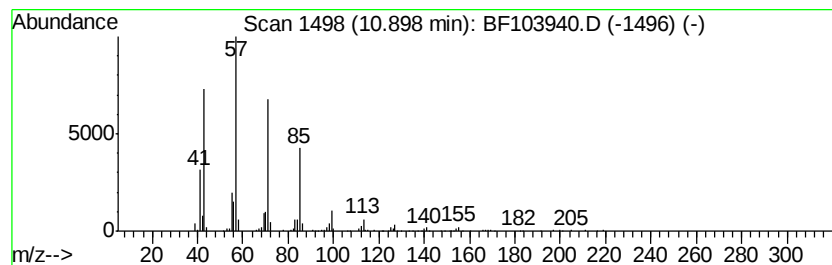
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Heptadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.90	8.96 ng	457199	Phenanthrene-d10	11.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	91
2	Hexadecane		226	C16H34	000544-76-3	90
3	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	90
4	Heptacosane, 1-chloro-		414	C27H55Cl	062016-79-9	86
5	Sulfurous acid, 2-propyl tetra...		320	C17H36O3S	1000309-12-5	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

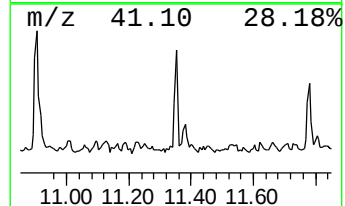
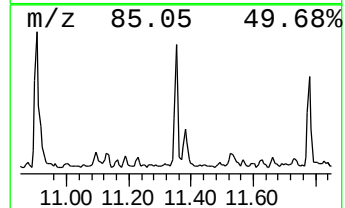
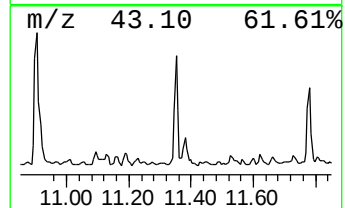
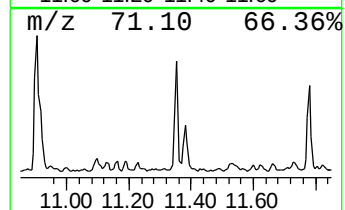
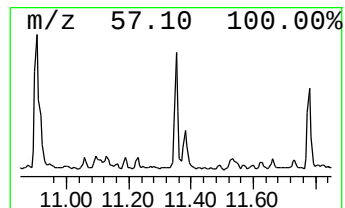
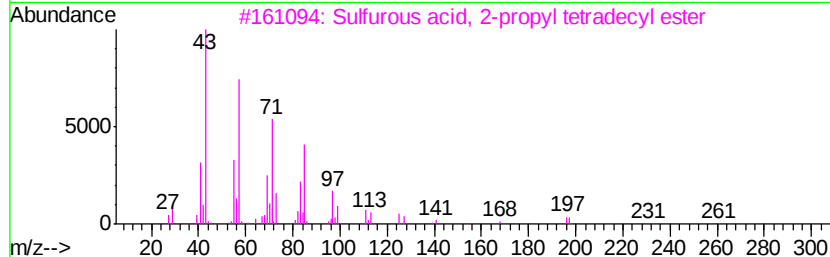
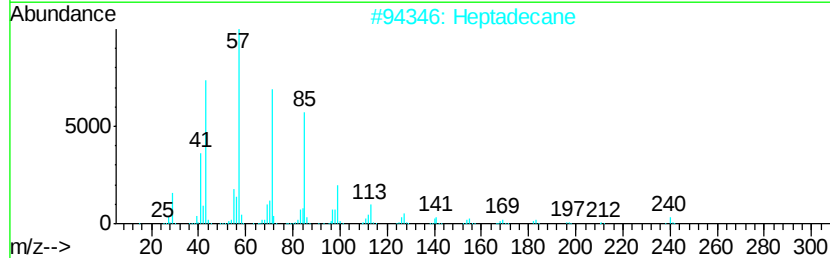
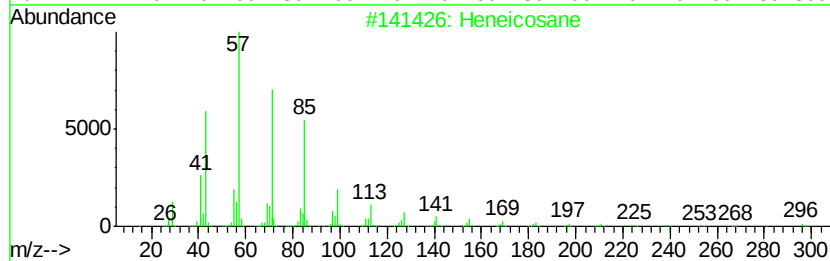
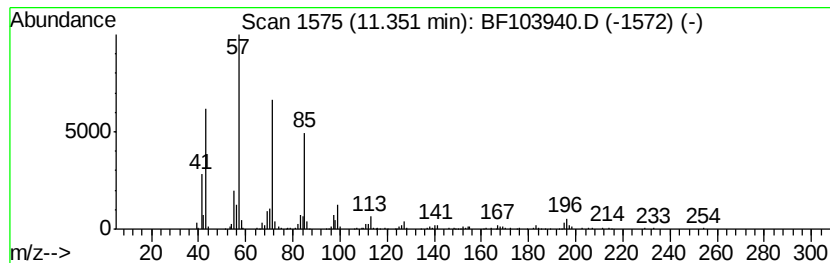
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heneicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	4.77 ng	243249	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Heptadecane	240	C17H36	000629-78-7	91
3		Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	90
4		Hexadecane	226	C16H34	000544-76-3	86
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

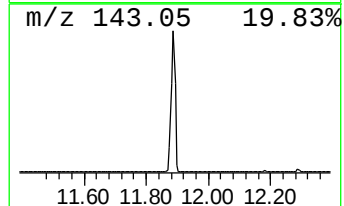
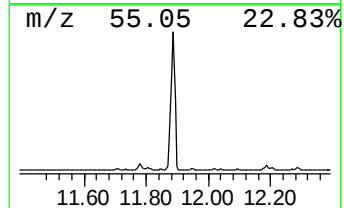
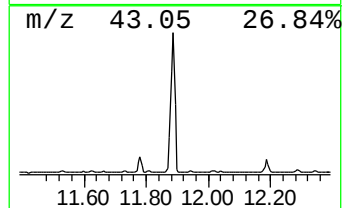
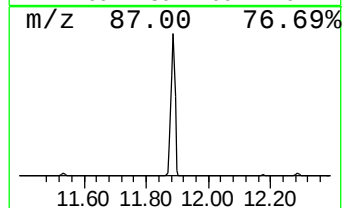
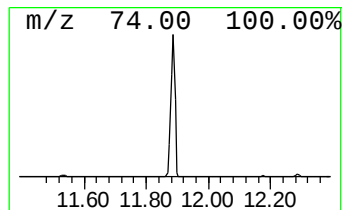
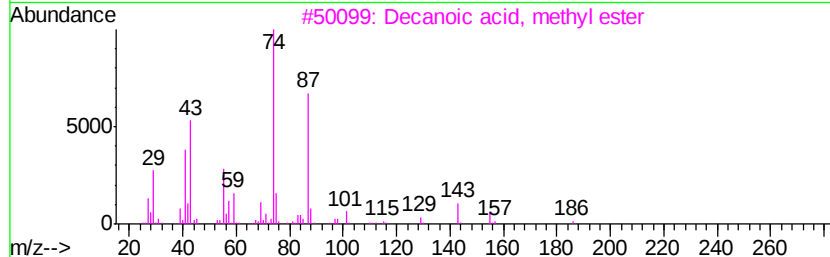
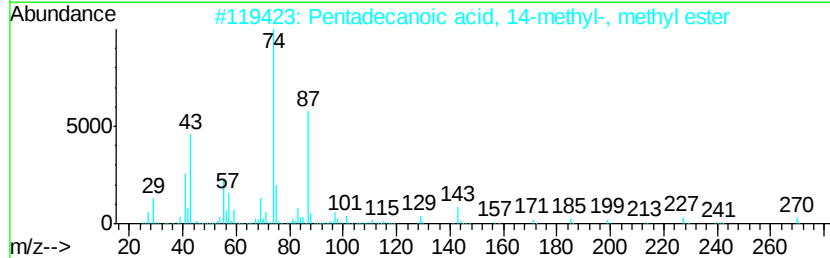
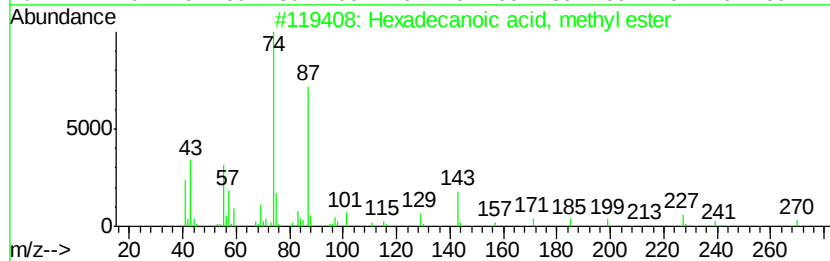
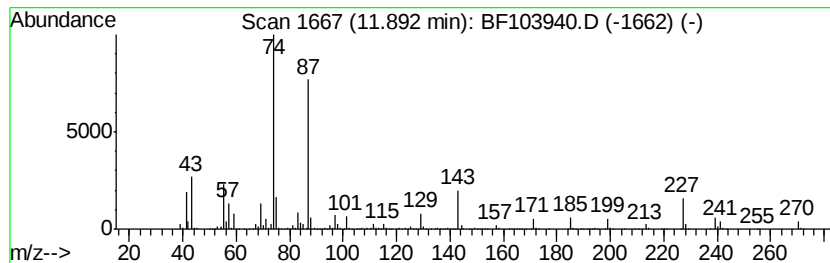
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	70.63 ng	3604750	Phenanthrene-d10	11.51

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	93
3		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

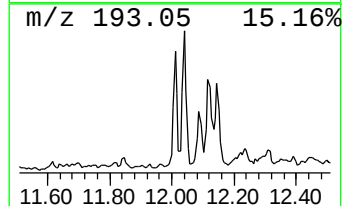
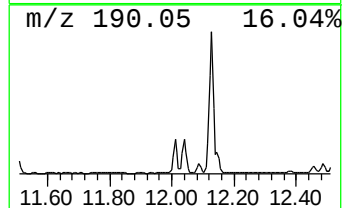
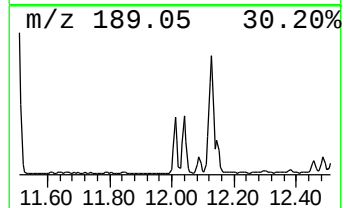
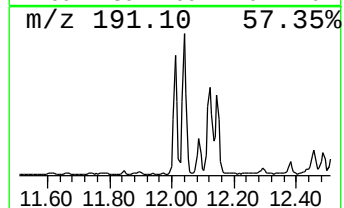
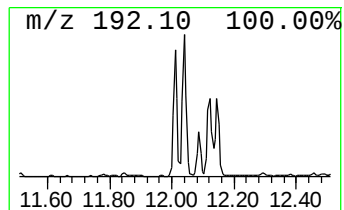
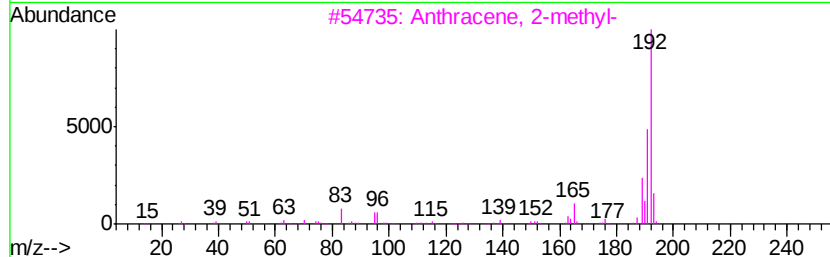
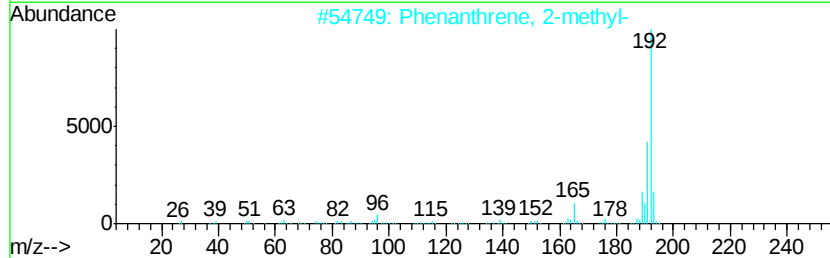
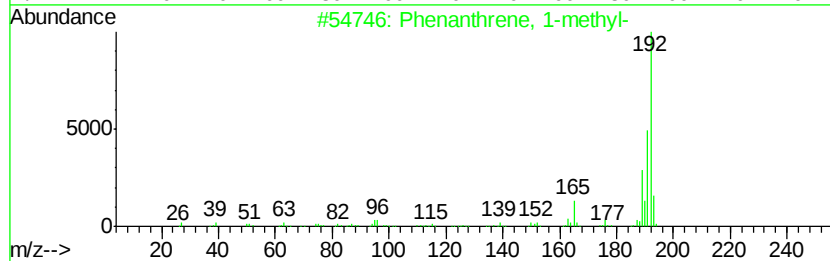
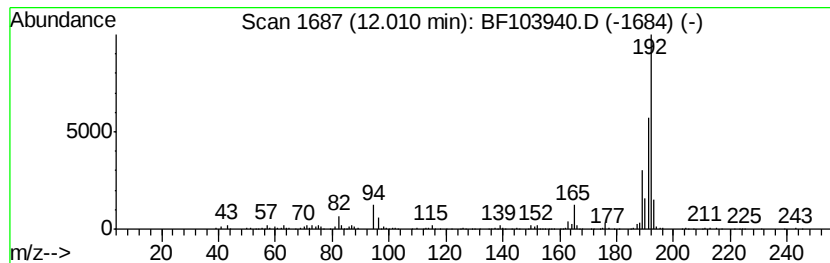
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenanthrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	5.45 ng	278238	Phenanthrene-d10	11.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	95
5			Anthracene, 9-methyl-	192	C15H12	000779-02-2	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

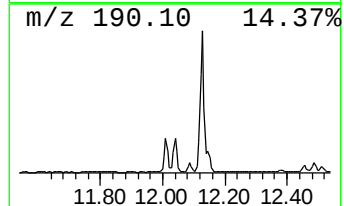
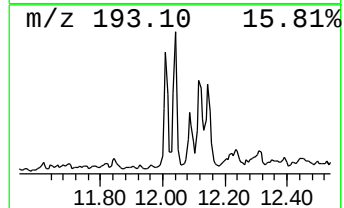
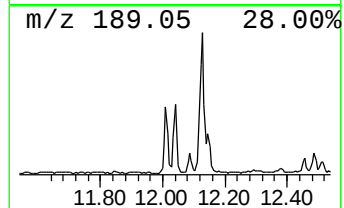
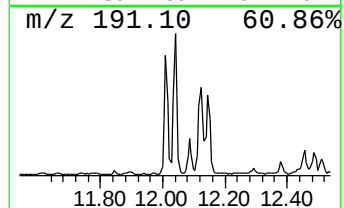
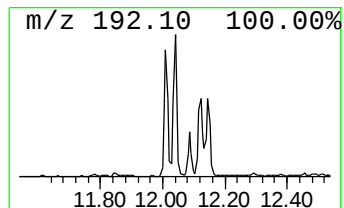
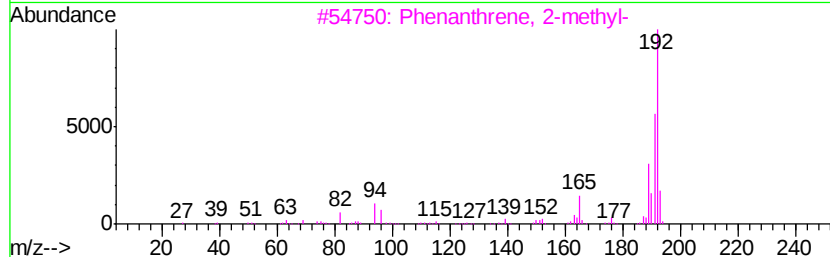
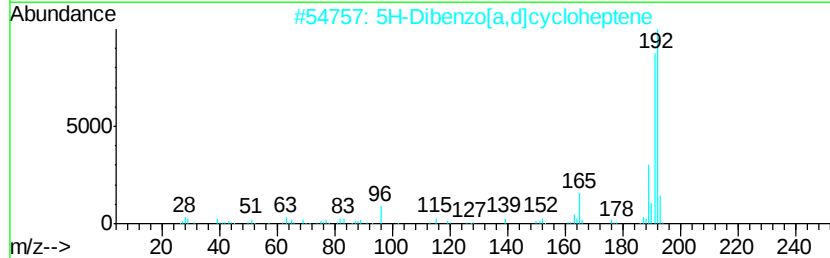
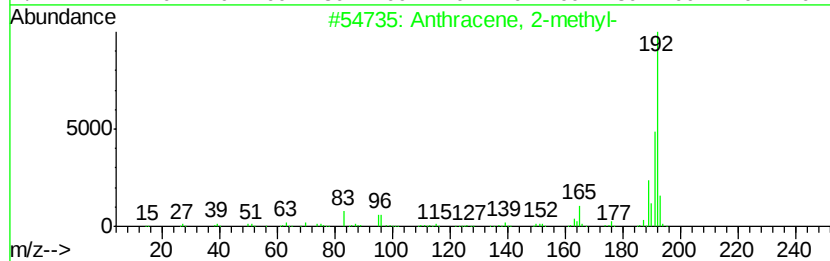
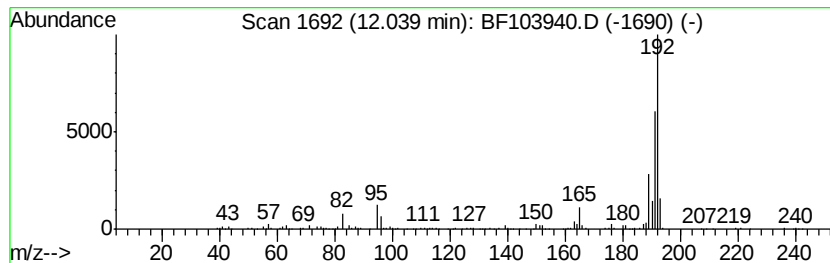
Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Anthracene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.04	5.16 ng	263181	Phenanthrene-d10	11.51	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	95
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

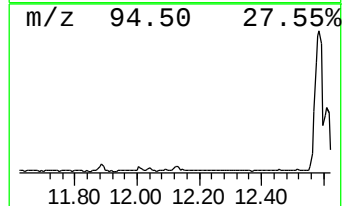
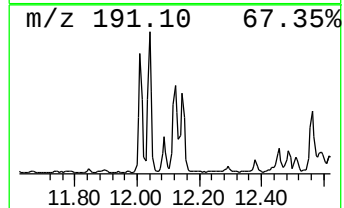
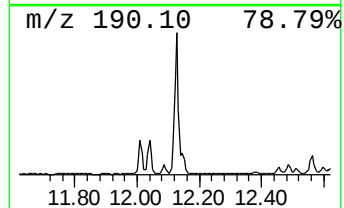
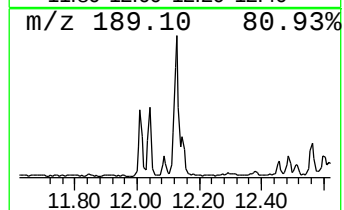
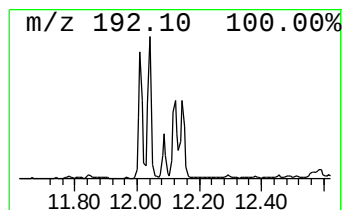
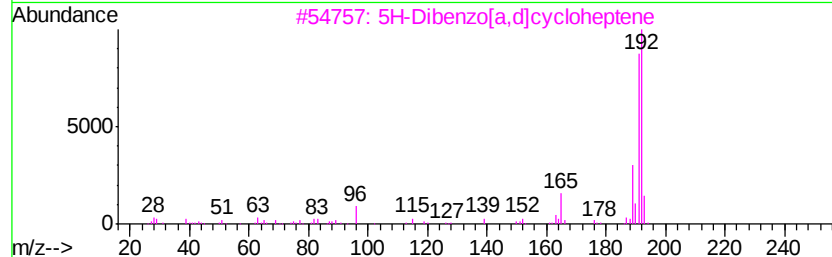
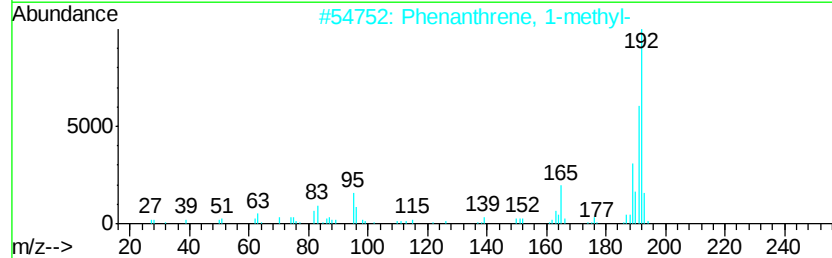
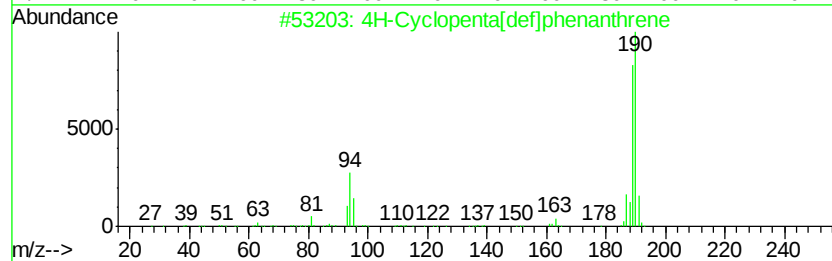
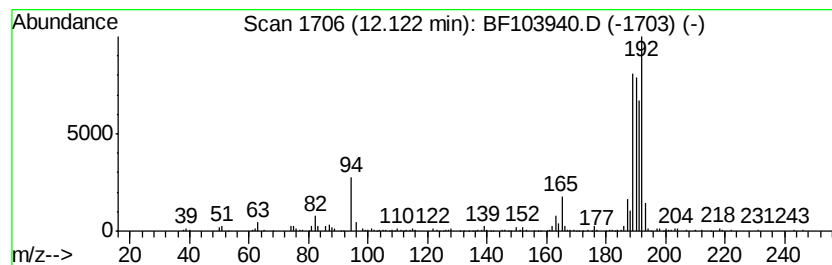
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown12.12 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	6.70 ng	341779	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
3	5H-Dibenzof[a,d]cycloheptene	192	C15H12	000256-81-5	42
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

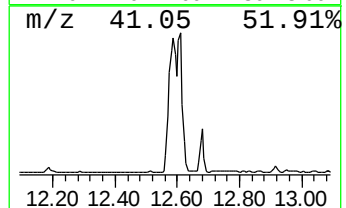
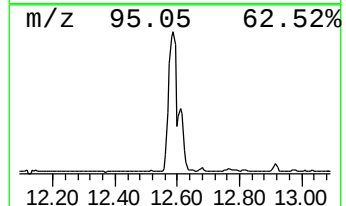
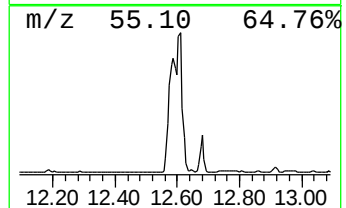
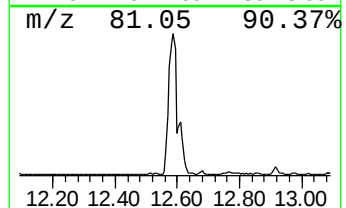
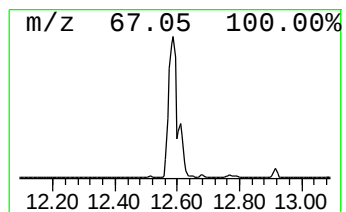
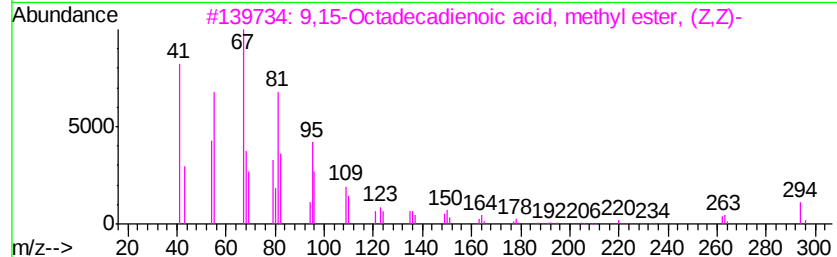
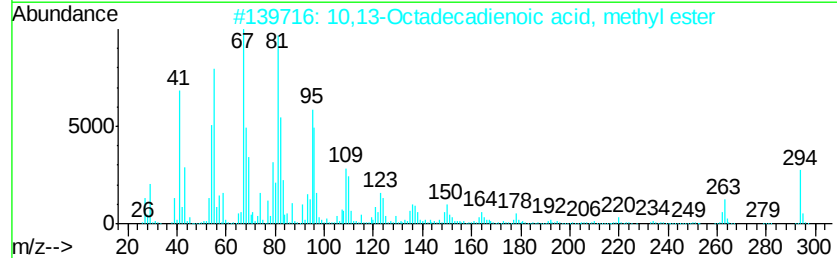
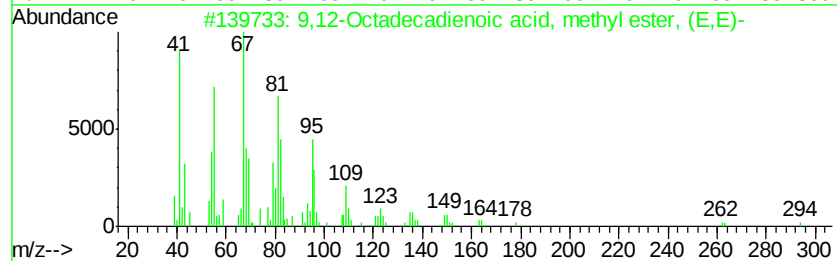
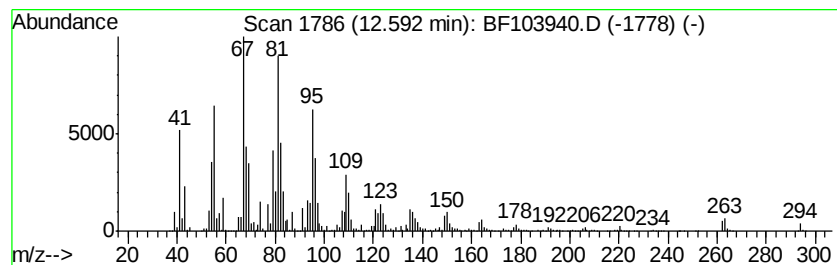
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	186.89 ng	9538720	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
3		9,15-Octadecadienoic acid, methv...	294	C19H34O2	017309-05-6	99
4		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
5		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

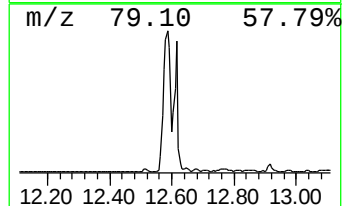
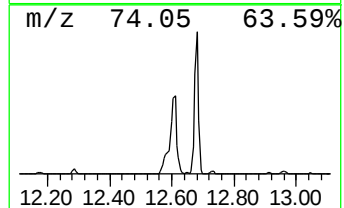
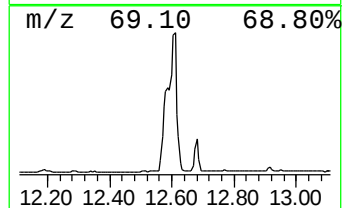
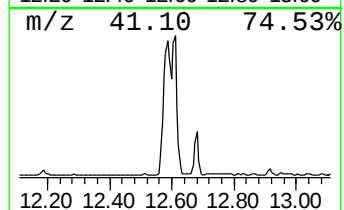
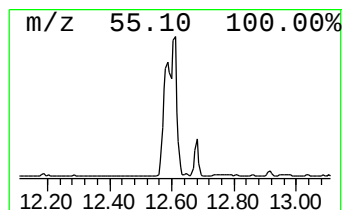
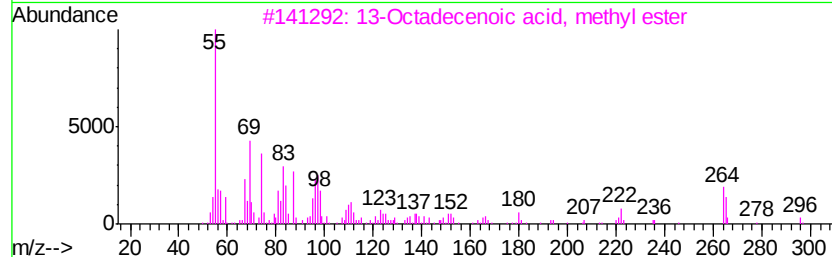
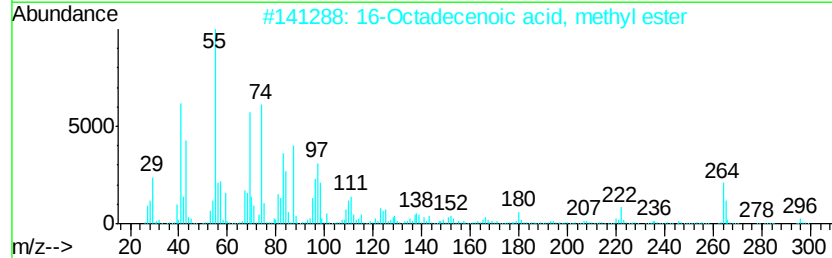
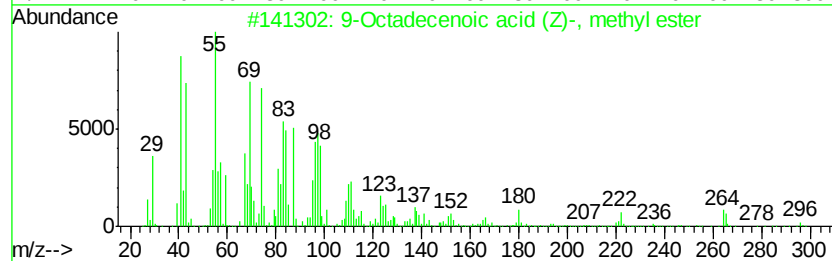
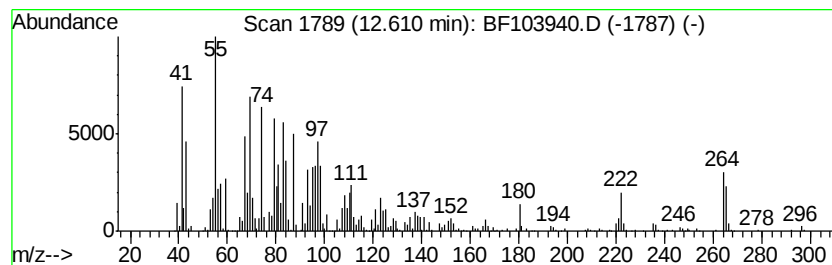
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	106.97 ng	5459900	Phenanthrene-d10	11.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
3		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	98
4		6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	96
5		15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

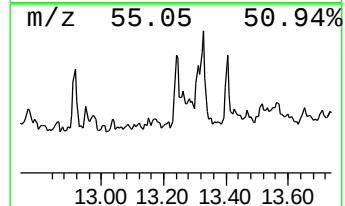
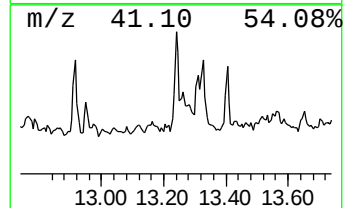
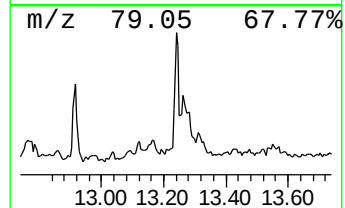
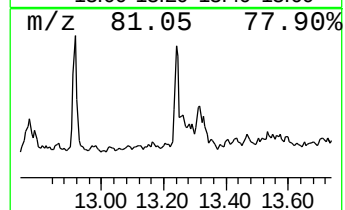
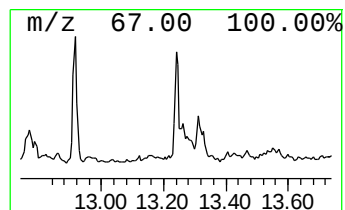
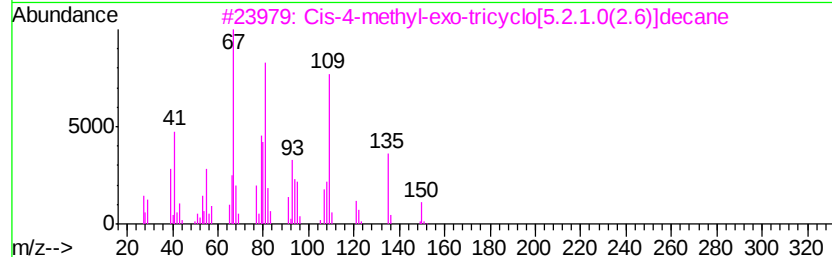
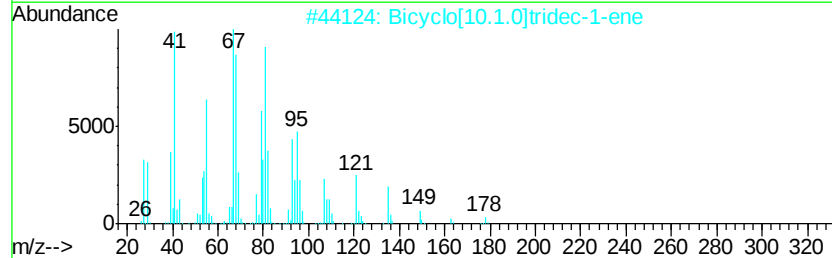
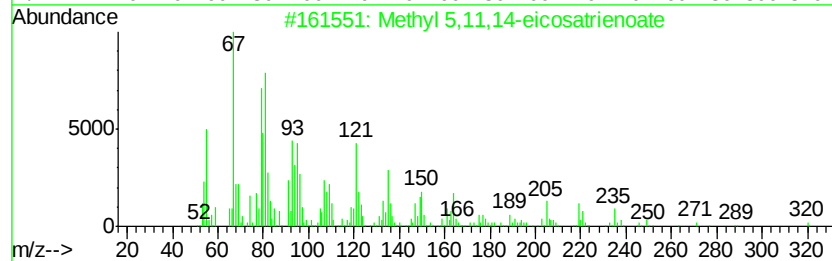
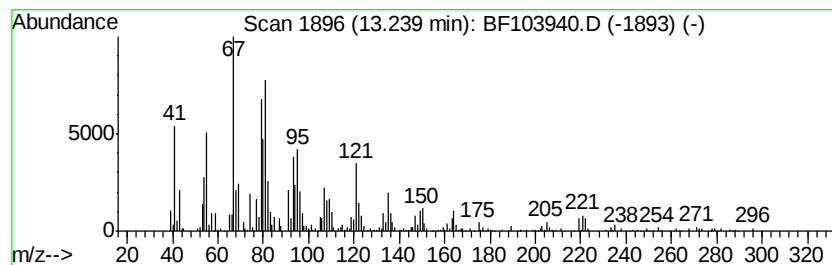
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Methyl 5,11,14-eicosatrienoate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	4.56 ng	294252	Chrysene-d12	14.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 5,11,14-eicosatrienoate	320	C21H36O2	1000336-40-2	91
2			Bicyclo[10.1.0]tridec-1-ene	178	C13H22	054766-91-5	90
3			Cis-4-methyl-exo-tricyclo[5.2.1....	150	C11H18	1000215-29-2	68
4			Cyclooctene, 4-ethenyl-	136	C10H16	001124-45-4	64
5			Bicyclo[6.1.0]non-1-ene	122	C9H14	002570-06-1	60



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

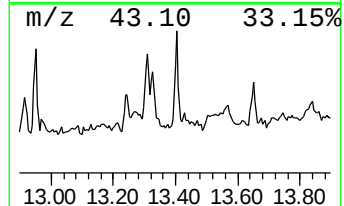
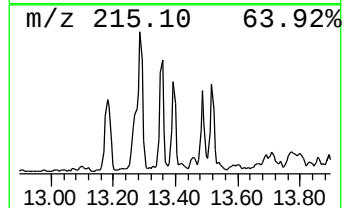
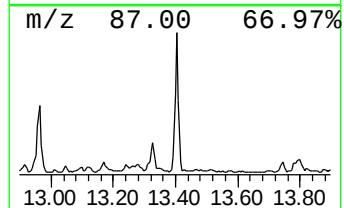
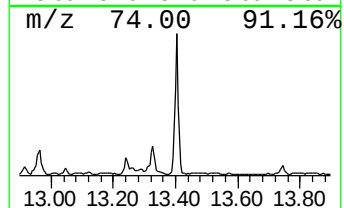
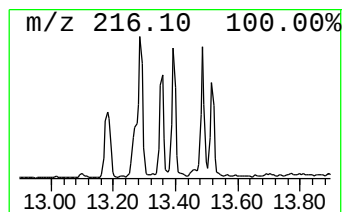
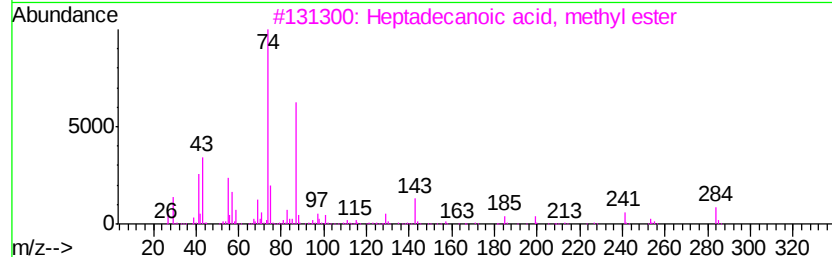
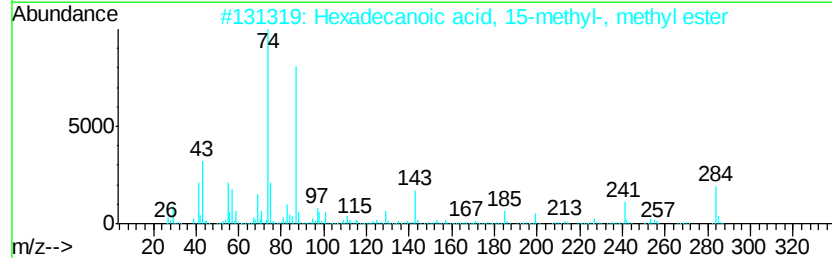
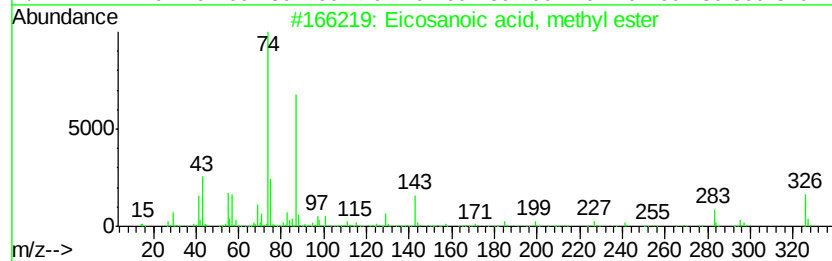
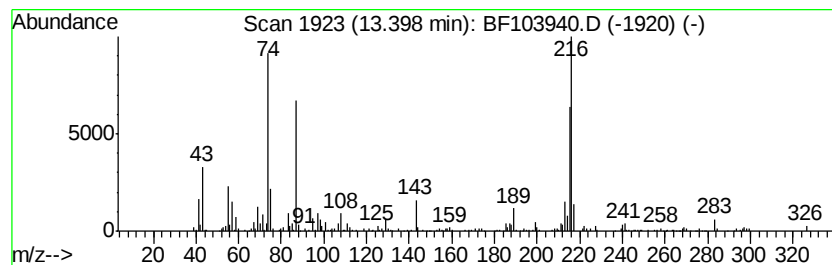
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Eicosanoic acid, methyl ester Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.40	3.98 ng	256921	Chrysene-d12	14.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	97
2		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	95
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
4		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
5		Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

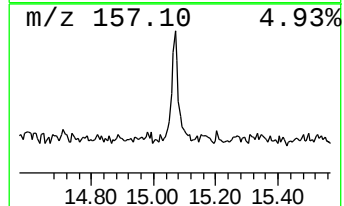
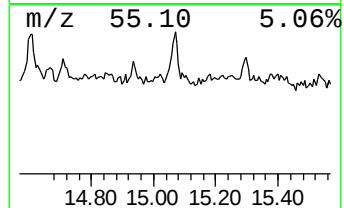
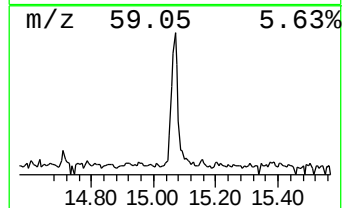
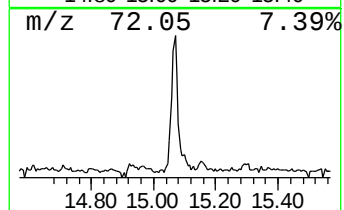
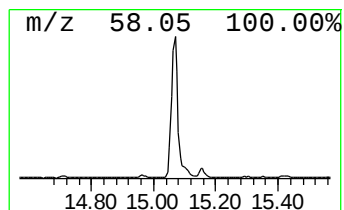
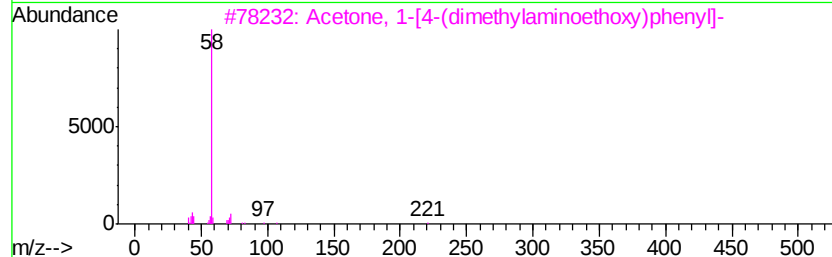
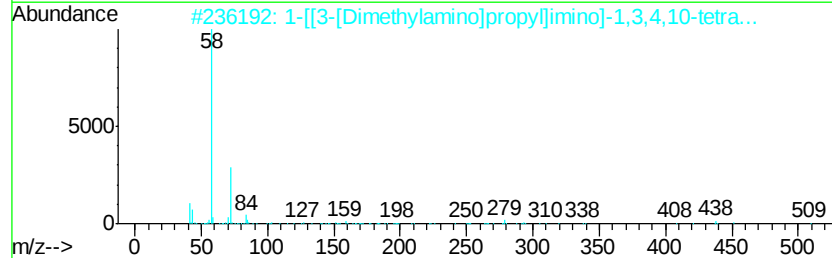
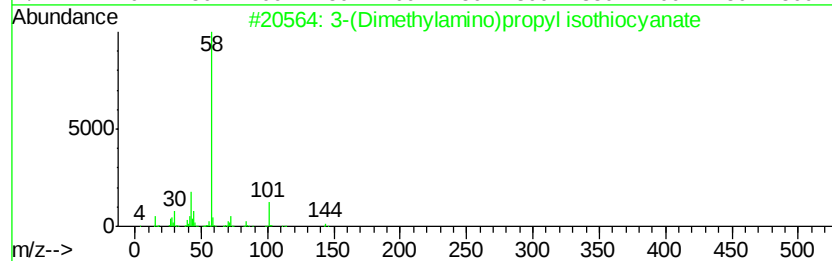
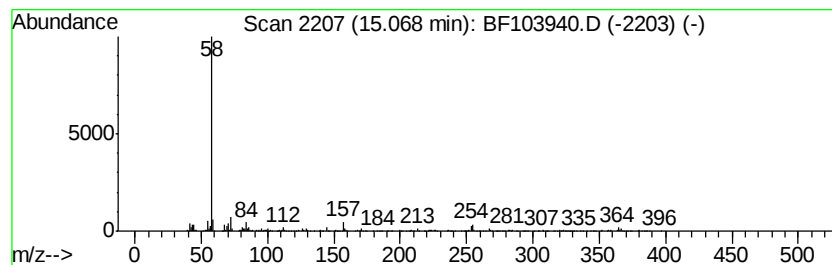
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 3-(Dimethylamino)propyl iso... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	12.61 ng	417218	Perylene-d12	15.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-(Dimethylamino)propyl isothioc...	144	C6H12N2S	027421-70-1	64
2		1-[[3-[Dimethylamino]propyl]imin...	509	C26H25F6N3O	144155-44-2	59
3		Acetone, 1-[4-(dimethylaminoetho...	221	C13H19NO2	086608-11-9	59
4		N,N-Dimethyl-2-cyclohexyloxvethy...	171	C10H21NO	071126-60-8	59
5		1-Nonanamine, N,N-dimethyl-	171	C11H25N	017373-27-2	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032618\
Data File : BF103940.D
Acq On : 26 Mar 2018 22:07
Operator : JU/SJ
Sample : J1760-07 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-032318-B

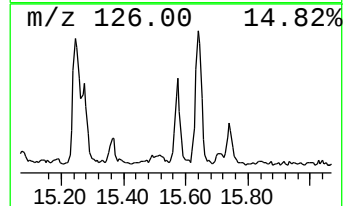
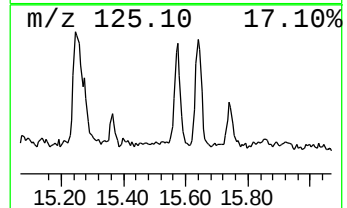
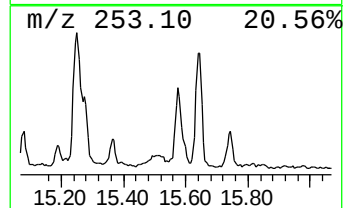
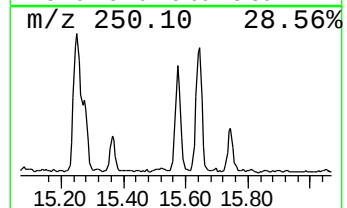
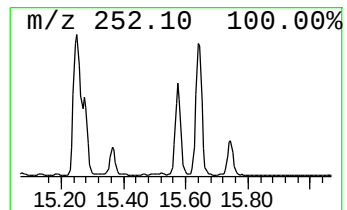
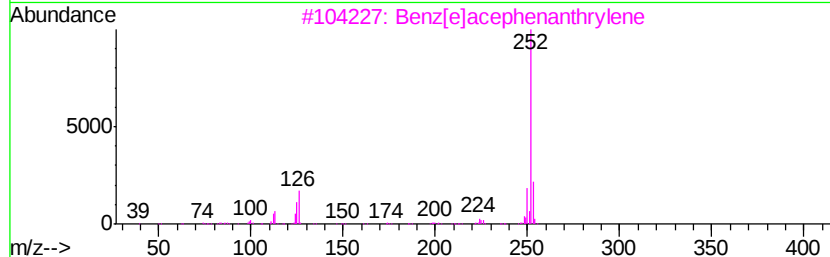
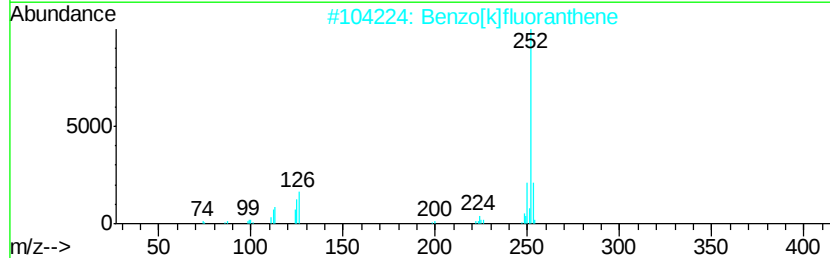
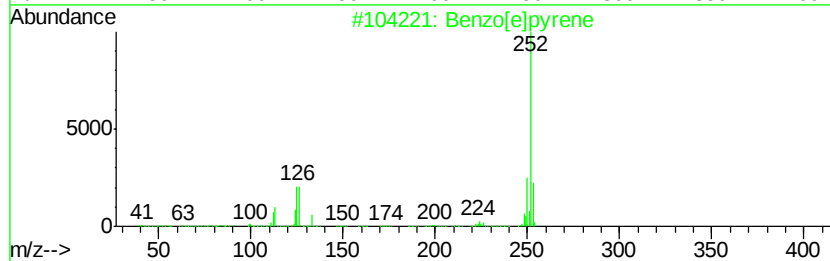
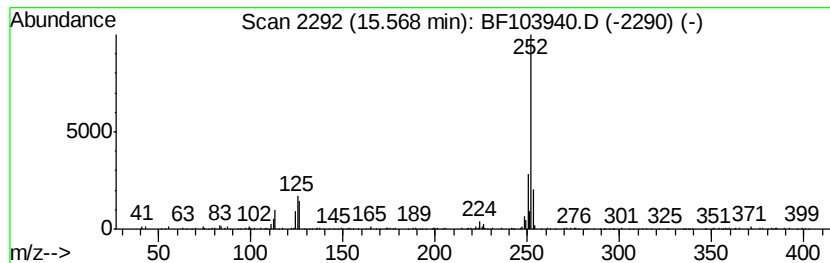
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	6.97 ng	230587	Perylene-d12	15.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
3			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	97
4			Perylene	252	C20H12	000198-55-0	94
5			Benzo[a]pyrene	252	C20H12	000050-32-8	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032618\
 Data File : BF103940.D
 Acq On : 26 Mar 2018 22:07
 Operator : JU/SJ
 Sample : J1760-07 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-032318-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031518.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
unknown6.74	6.74	40.8	ng	1697910	1	6.97	831511	20.0
Naphthalene, 2-me...	9.07	2.9	ng	155528	2	8.26	1063280	20.0
Tetradecane	9.93	4.0	ng	220767	3	10.02	1090690	20.0
Hexadecane	10.43	6.5	ng	352363	3	10.02	1090690	20.0
Dodecane	10.65	3.0	ng	162393	3	10.02	1090690	20.0
Heptadecane	10.90	9.0	ng	457199	4	11.51	1020790	20.0
Heneicosane	11.35	4.8	ng	243249	4	11.51	1020790	20.0
Hexadecanoic acid...	11.89	70.6	ng	3604750	4	11.51	1020790	20.0
Phenanthrene, 1-m...	12.01	5.5	ng	278238	4	11.51	1020790	20.0
Anthracene, 2-met...	12.04	5.2	ng	263181	4	11.51	1020790	20.0
unknown12.12	12.12	6.7	ng	341779	4	11.51	1020790	20.0
9,12-Octadecadien...	12.59	186.9	ng	9538720	4	11.51	1020790	20.0
9-Octadecenoic ac...	12.61	107.0	ng	5459900	4	11.51	1020790	20.0
Methyl 5,11,14-ei...	13.24	4.6	ng	294252	5	14.16	1290060	20.0
Eicosanoic acid, ...	13.40	4.0	ng	256921	5	14.16	1290060	20.0
3-(Dimethylamino)...	15.07	12.6	ng	417218	6	15.71	661857	20.0
Benzo[e]pyrene	15.57	7.0	ng	230587	6	15.71	661857	20.0