

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF033018\
 Data File : BF104109.D
 Acq On : 30 Mar 2018 22:39
 Operator : JU/SJ
 Sample : J1760-09 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-032918

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-BF032818.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	2.269	27	31	37	rVB	39375	51119	3.53%	0.281%
2	5.187	522	527	537	rBV	36354	52139	3.60%	0.287%
3	5.587	591	595	619	rBV	605449	1076916	74.32%	5.928%
4	6.587	762	765	786	rBV	639967	1227990	84.74%	6.760%
5	6.728	786	789	795	rBV	984672	1140921	78.73%	6.281%
6	6.957	824	828	841	rBV	375713	576835	39.81%	3.175%
7	7.110	851	854	867	rBV	889649	954187	65.85%	5.253%
8	7.210	870	871	874	rVV2	20461	18409	1.27%	0.101%
9	7.269	878	881	887	rVB3	20200	28735	1.98%	0.158%
10	7.522	920	924	934	rBV	437512	756193	52.18%	4.163%
11	7.898	985	988	992	rVB4	11976	14825	1.02%	0.082%
12	7.969	997	1000	1002	rBV2	30057	28825	1.99%	0.159%
13	8.198	1036	1039	1042	rBV	118380	95420	6.58%	0.525%
14	8.239	1042	1046	1051	rVV	606774	739408	51.03%	4.070%
15	8.275	1051	1052	1056	rVV	94917	86946	6.00%	0.479%
16	8.310	1056	1058	1066	rVB6	26445	38471	2.65%	0.212%
17	8.428	1075	1078	1083	rVB4	20695	24704	1.70%	0.136%
18	8.516	1089	1093	1097	rVB4	34598	43478	3.00%	0.239%
19	8.633	1111	1113	1116	rBV	48487	35117	2.42%	0.193%
20	8.733	1127	1130	1132	rBV3	24903	22299	1.54%	0.123%
21	8.804	1137	1142	1147	rVV	103586	103700	7.16%	0.571%
22	8.898	1155	1158	1163	rVB4	17971	23129	1.60%	0.127%
23	9.051	1181	1184	1188	rBV3	24511	26342	1.82%	0.145%
24	9.133	1195	1198	1201	rVB2	15102	17834	1.23%	0.098%
25	9.169	1201	1204	1208	rBV3	22063	23407	1.62%	0.129%
26	9.239	1213	1216	1218	rVB	31614	27246	1.88%	0.150%
27	9.310	1224	1228	1236	rBV	1270812	1449091	100.00%	7.977%
28	9.375	1236	1239	1244	rVB	118636	122428	8.45%	0.674%
29	9.698	1289	1294	1301	rBV2	79846	123485	8.52%	0.680%
30	9.857	1316	1321	1324	rBV	19368	29062	2.01%	0.160%
31	9.904	1325	1329	1333	rVB	101689	89536	6.18%	0.493%
32	9.992	1341	1344	1357	rVB	807466	873607	60.29%	4.809%
33	10.227	1381	1384	1388	rVB4	18830	24043	1.66%	0.132%
34	10.304	1394	1397	1401	rBV6	11003	16253	1.12%	0.089%

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35	10.404	1411	1414	1416	rVB	86858	64971	4.48%	0.358%
36	10.545	1436	1438	1444	rVB3	11878	18386	1.27%	0.101%
37	10.622	1447	1451	1454	rBV	35529	39610	2.73%	0.218%
38	10.786	1476	1479	1492	rBV	583733	784605	54.14%	4.319%
39	10.874	1492	1494	1504	rVB2	72131	111471	7.69%	0.614%
40	11.327	1568	1571	1573	rBV	35464	29349	2.03%	0.162%
41	11.351	1573	1575	1579	rVB2	23288	23193	1.60%	0.128%
42	11.492	1595	1599	1601	rBV	558413	625436	43.16%	3.443%
43	11.569	1610	1612	1621	rVB	31296	56655	3.91%	0.312%
44	11.751	1638	1643	1646	rBV2	33946	34036	2.35%	0.187%
45	11.857	1657	1661	1666	rBV	432651	345648	23.85%	1.903%
46	11.992	1680	1684	1687	rBV2	38962	50110	3.46%	0.276%
47	12.021	1687	1689	1695	rVB2	23774	35134	2.42%	0.193%
48	12.104	1700	1703	1706	rBV2	27356	34374	2.37%	0.189%
49	12.316	1737	1739	1745	rVB7	12364	15446	1.07%	0.085%
50	12.545	1774	1778	1780	rBV	1455910	1312989	90.61%	7.228%
51	12.568	1780	1782	1790	rVV2	927978	926911	63.96%	5.103%
52	12.651	1792	1796	1802	rVB	167430	157430	10.86%	0.867%
53	12.710	1802	1806	1816	rBV	253675	319903	22.08%	1.761%
54	12.886	1834	1836	1839	rVB3	22817	20686	1.43%	0.114%
55	12.939	1839	1845	1851	rBV	228457	319175	22.03%	1.757%
56	13.074	1864	1868	1878	rBV	1016465	1102082	76.05%	6.067%
57	13.215	1889	1892	1894	rBV2	32273	33087	2.28%	0.182%
58	13.333	1909	1912	1915	rBV	21206	21549	1.49%	0.119%
59	13.374	1916	1919	1921	rBV2	31060	28489	1.97%	0.157%
60	13.892	2005	2007	2010	rBV3	16370	22422	1.55%	0.123%
61	13.951	2014	2017	2022	rVV2	102530	124728	8.61%	0.687%
62	14.045	2031	2033	2035	rBV2	17242	16003	1.10%	0.088%
63	14.145	2044	2050	2053	rBV2	602997	819304	56.54%	4.510%
64	15.221	2230	2233	2236	rBV	71720	112591	7.77%	0.620%
65	15.609	2297	2299	2304	rVB	51177	64328	4.44%	0.354%
66	15.680	2307	2311	2322	rVB2	302368	533507	36.82%	2.937%

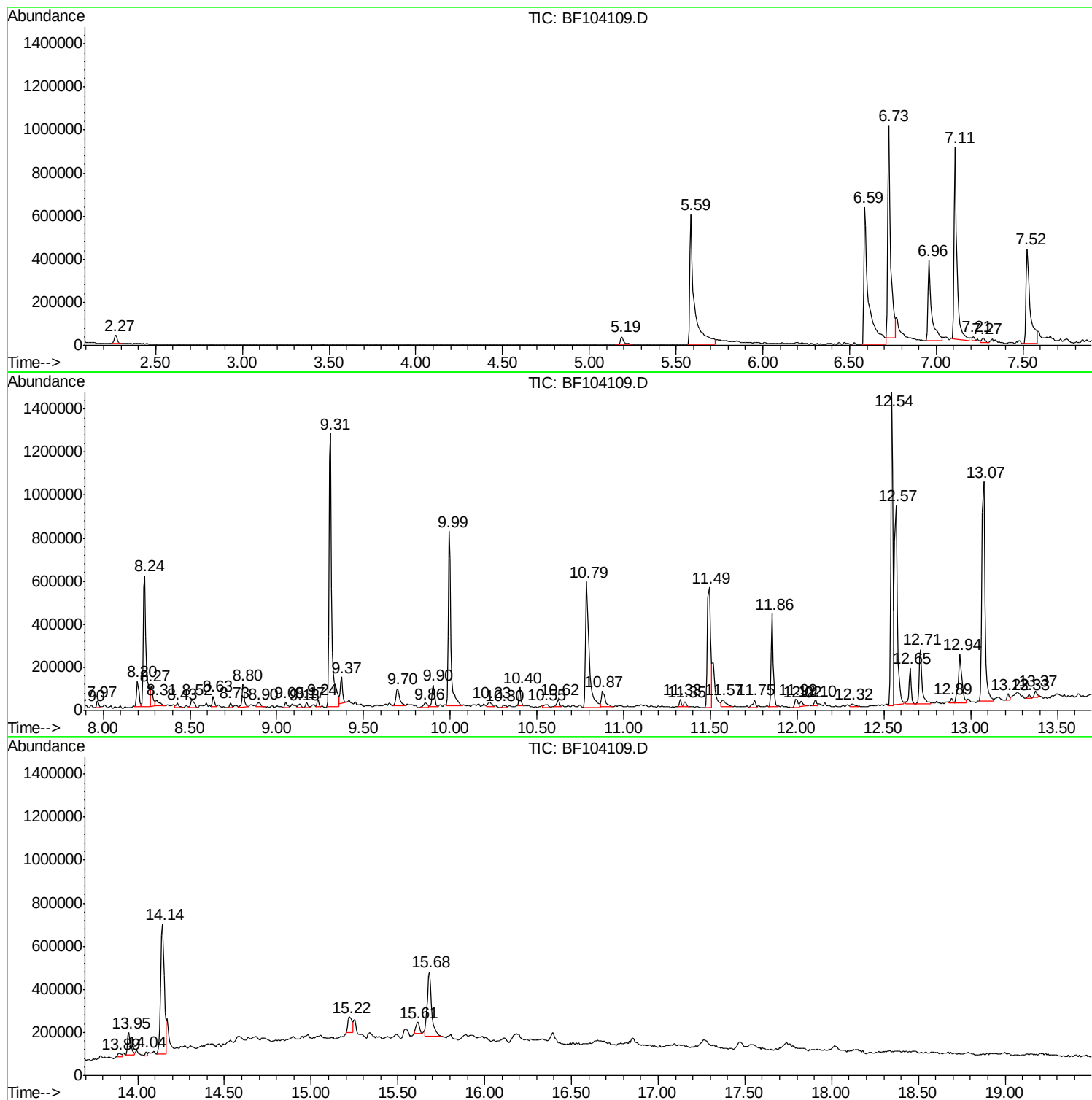
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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SU-03-032918

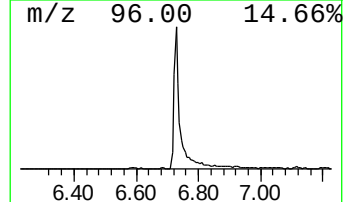
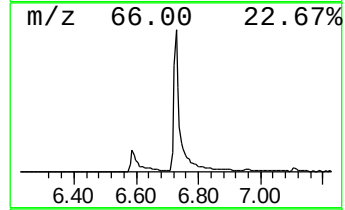
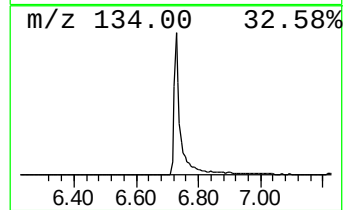
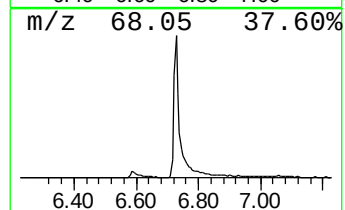
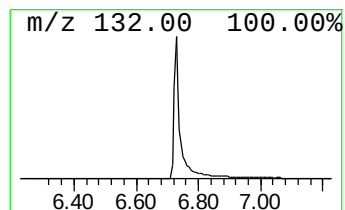
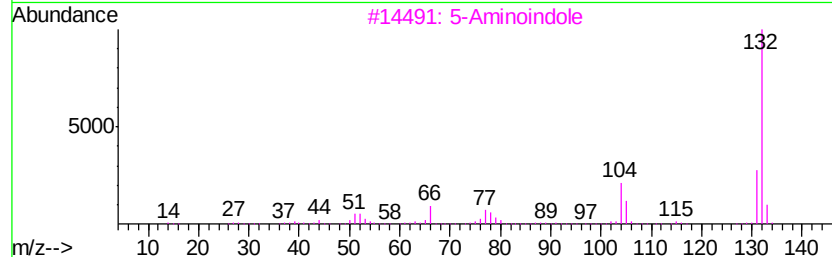
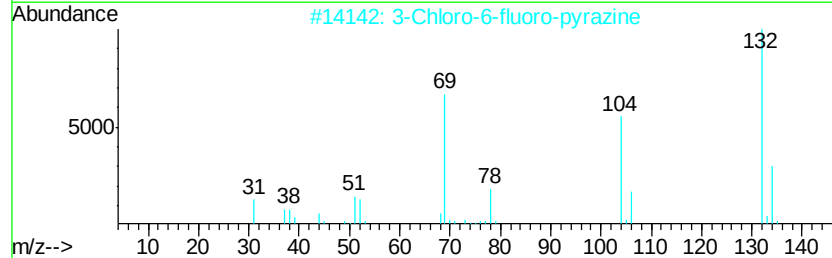
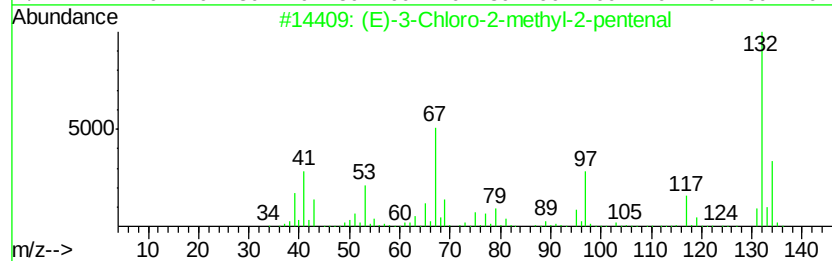
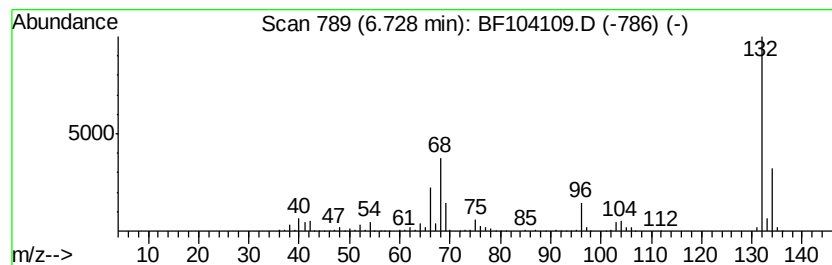
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.73 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	39.56 ng	1140920	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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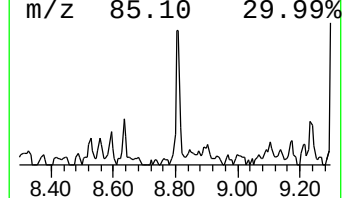
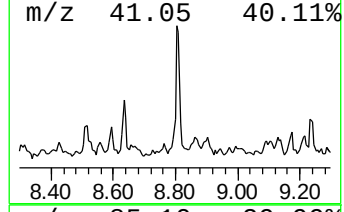
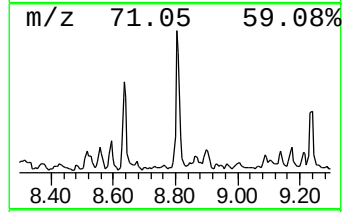
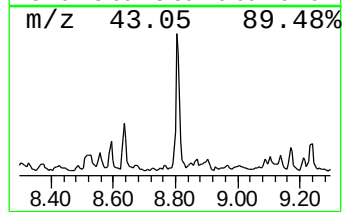
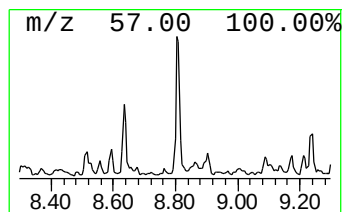
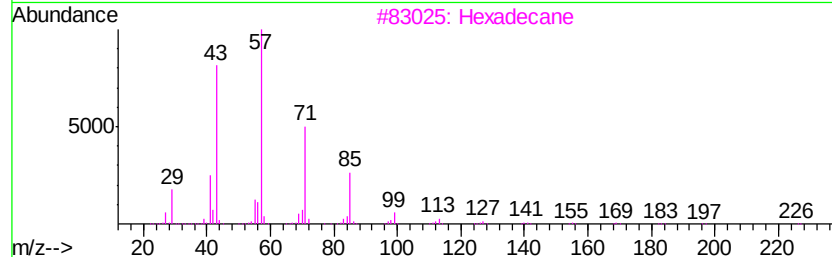
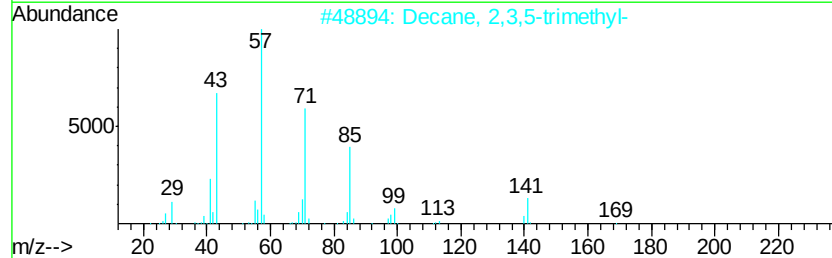
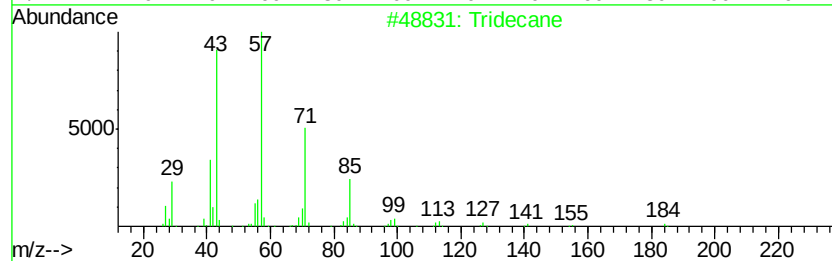
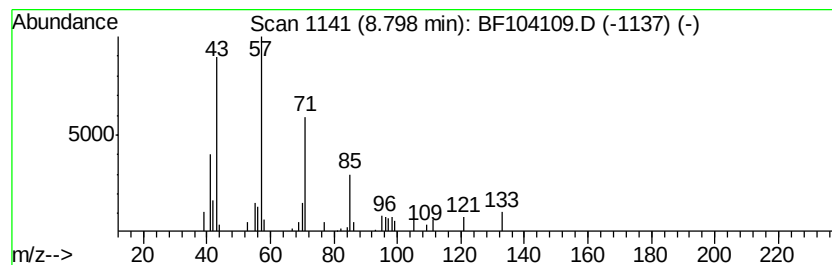
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.80	2.80 ng	103700	Naphthalene-d8	8.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
3	Hexadecane	226	C16H34	000544-76-3	86
4	Tetradecane	198	C14H30	000629-59-4	83
5	Nonane, 5-butyl-	184	C13H28	017312-63-9	81



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ALS Vial : 18 Sample Multiplier: 1

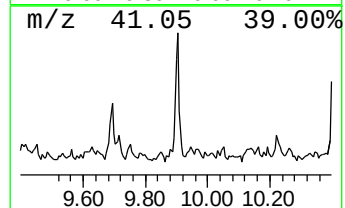
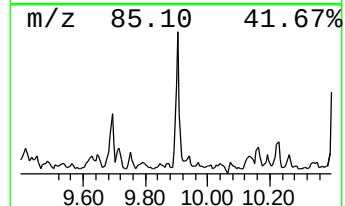
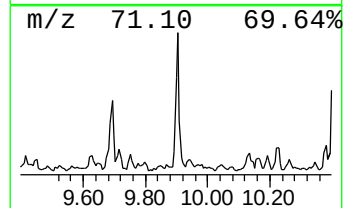
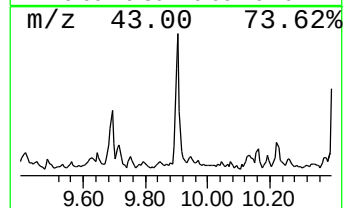
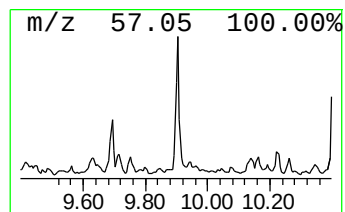
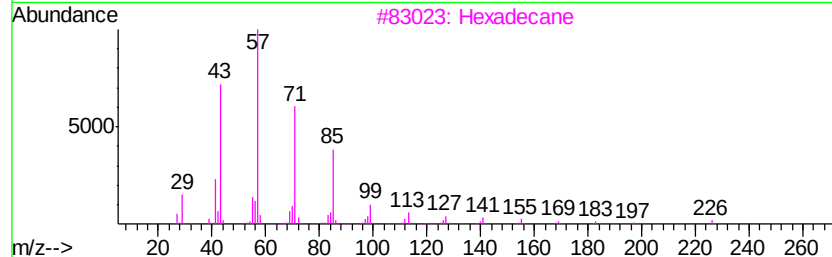
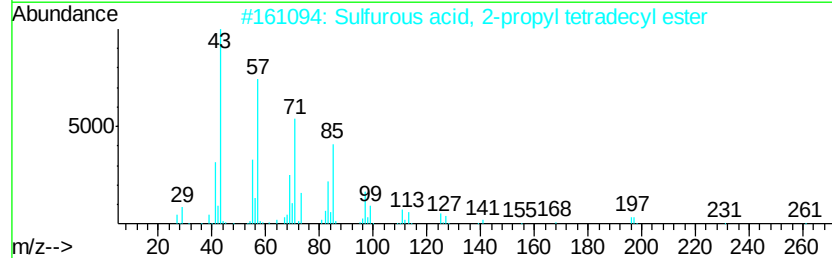
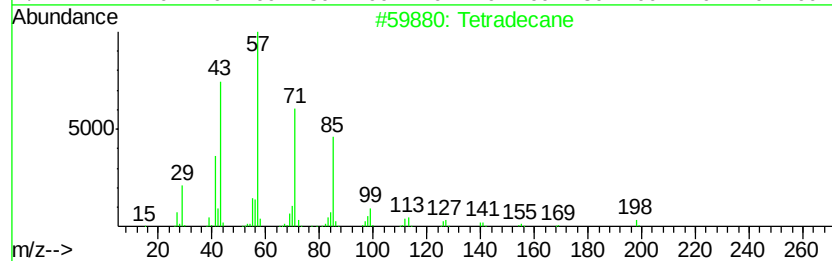
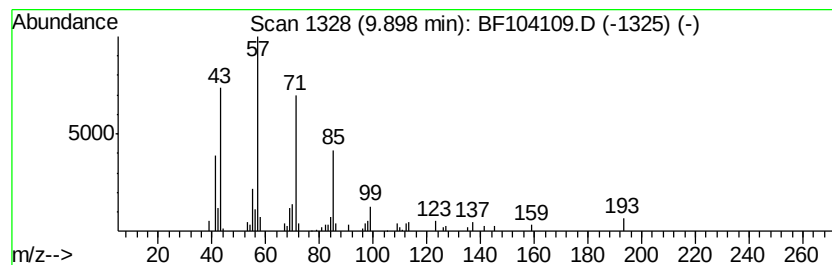
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
9.90	2.05 ng	89536	Acenaphthene-d10		9.99
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	91
2	Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Tridecane	184	C13H28	000629-50-5	90
5	Pentadecane	212	C15H32	000629-62-9	87



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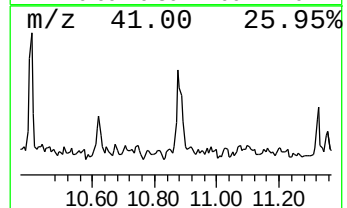
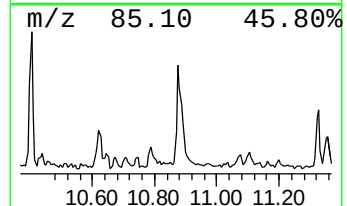
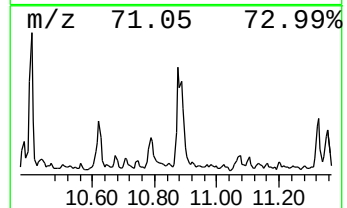
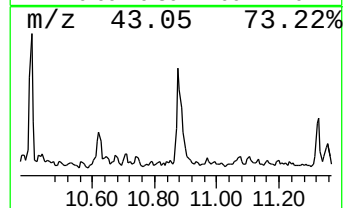
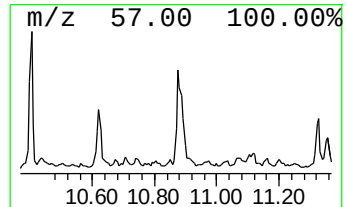
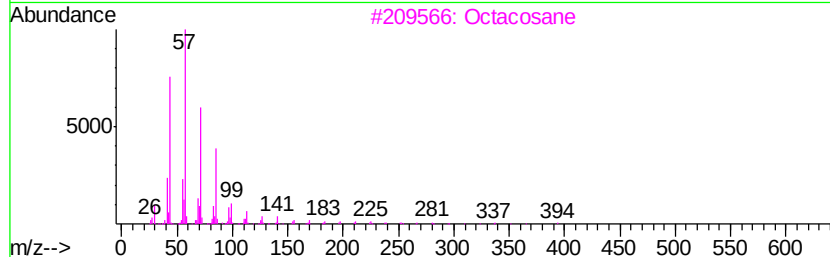
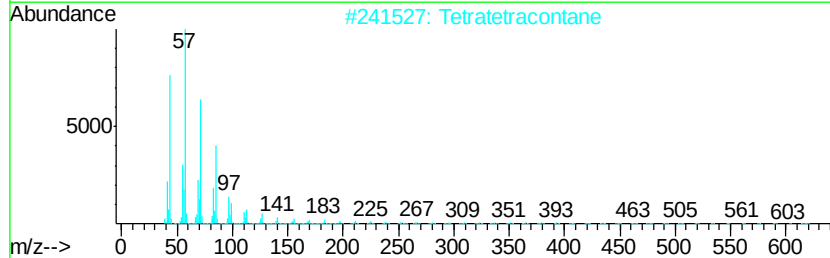
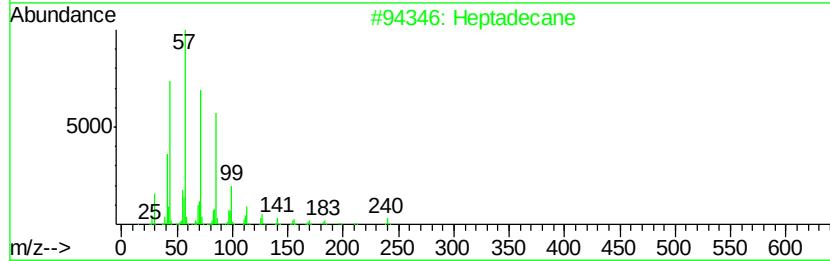
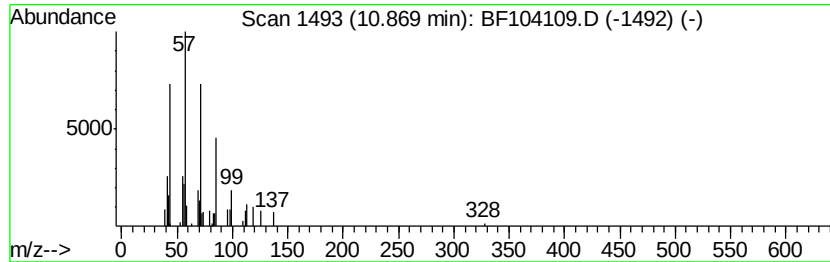
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Heptadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.87	3.56 ng	111471	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	91
2	Tetratetracontane	619	C44H90	007098-22-8	90
3	Octacosane	394	C28H58	000630-02-4	90
4	Pentadecane	212	C15H32	000629-62-9	90
5	Sulfurous acid, 2-propyl tetra...	320	C17H36O3S	1000309-12-5	90



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ALS Vial : 18 Sample Multiplier: 1

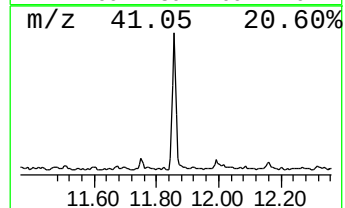
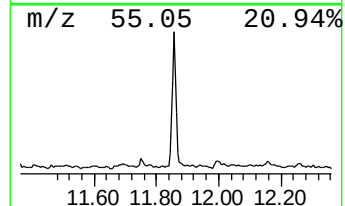
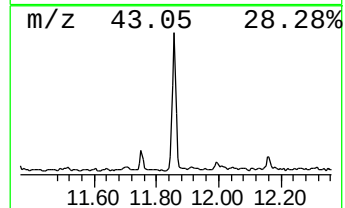
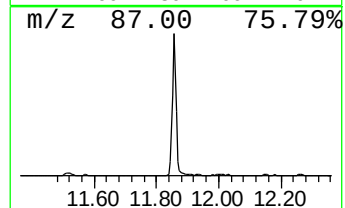
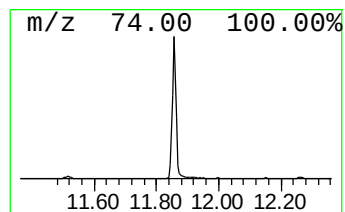
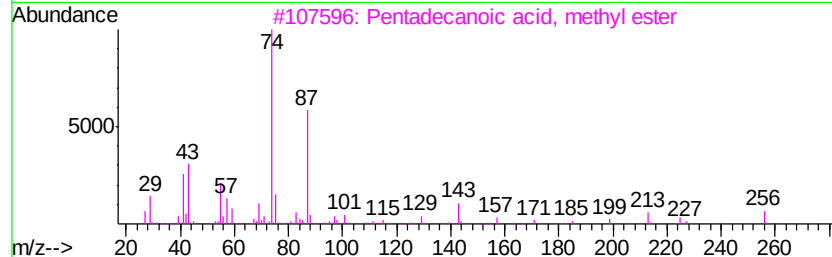
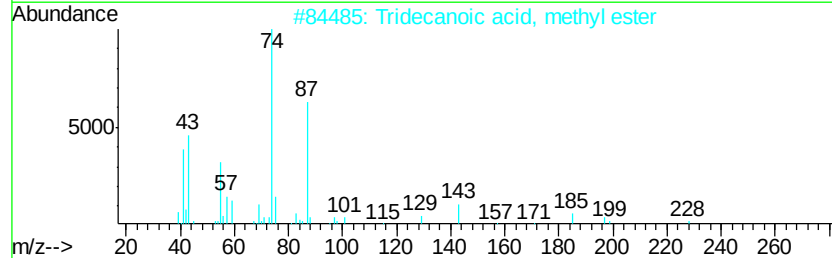
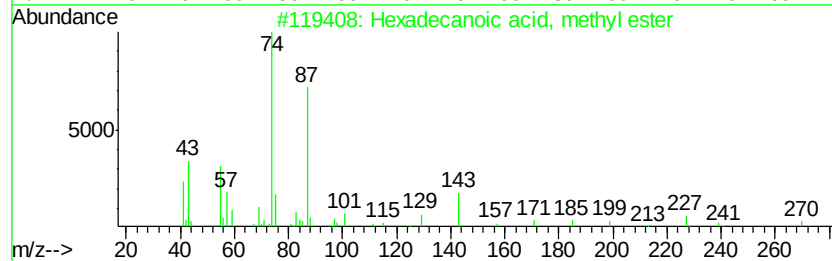
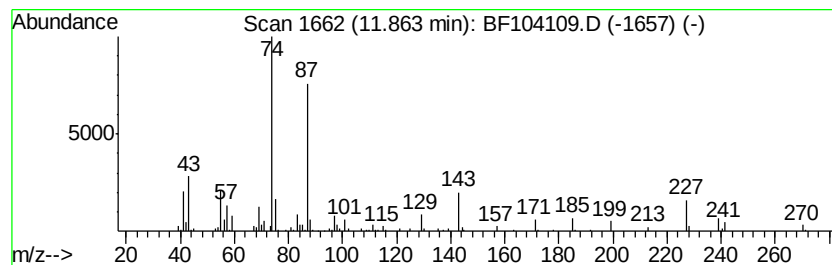
Instrument :
BNA_F
ClientSampleId :
SU-03-032918

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, methyl e... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.86	11.05 ng	345648	Phenanthrene-d10	11.49		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86
4		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	81
5		Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	80



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF033018\
Data File : BF104109.D
Acq On : 30 Mar 2018 22:39
Operator : JU/SJ
Sample : J1760-09 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

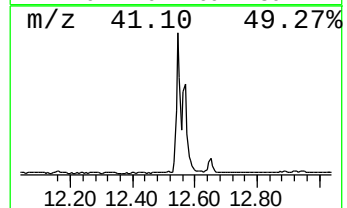
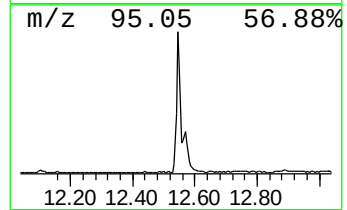
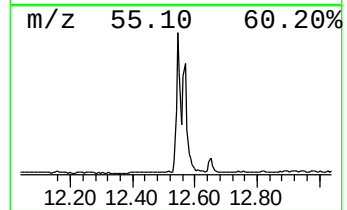
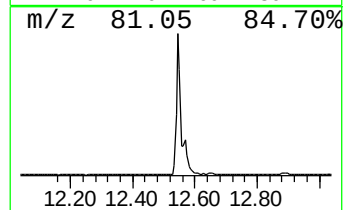
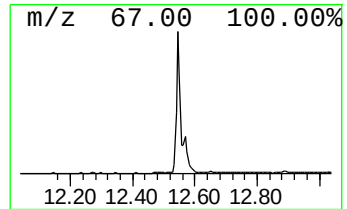
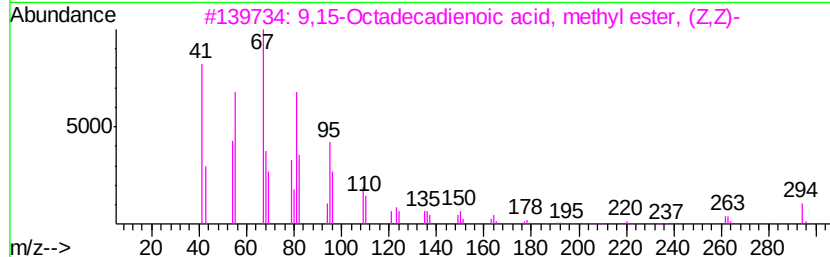
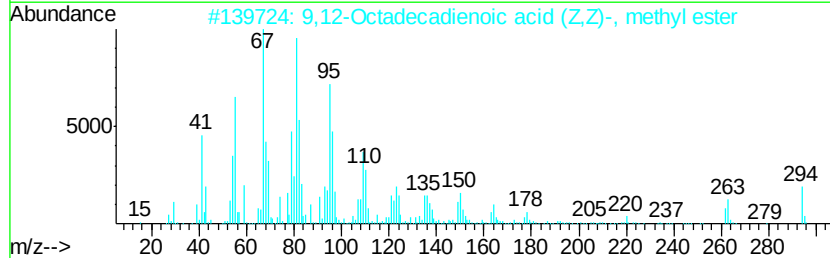
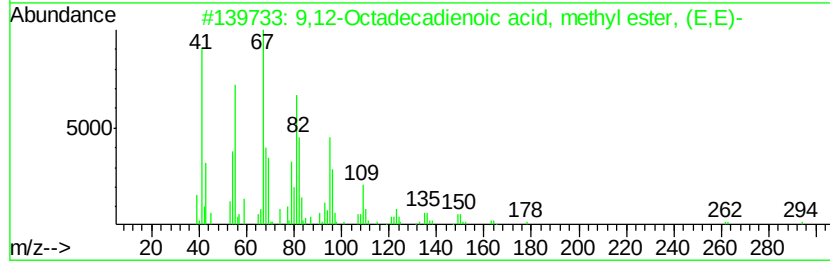
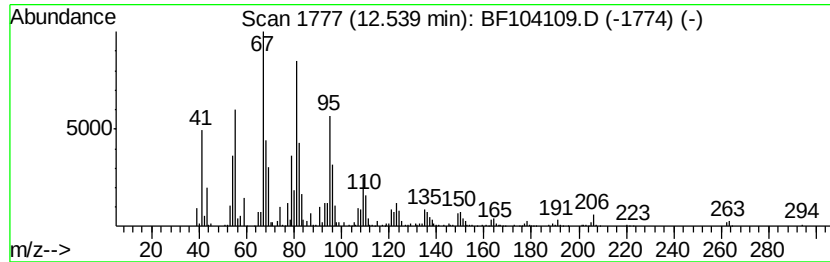
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ClientSampleId :
SU-03-032918

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.54	41.99 ng	1312990	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methv...	294	C19H34O2	002566-97-4	99
2	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3	9,15-Octadecadienoic acid, methv...	294	C19H34O2	017309-05-6	99
4	11,14-Octadecadienoic acid, meth...	294	C19H34O2	056554-61-1	98
5	Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF033018\
Data File : BF104109.D
Acq On : 30 Mar 2018 22:39
Operator : JU/SJ
Sample : J1760-09 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

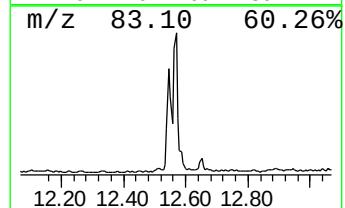
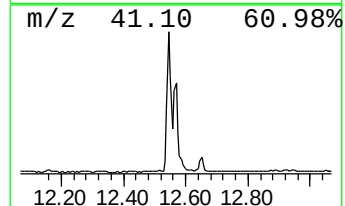
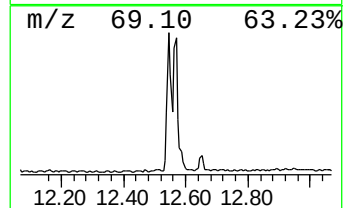
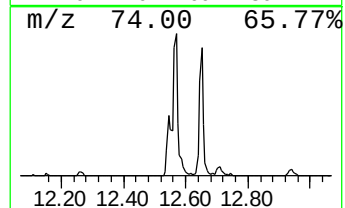
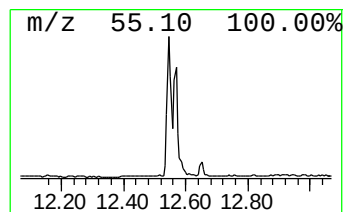
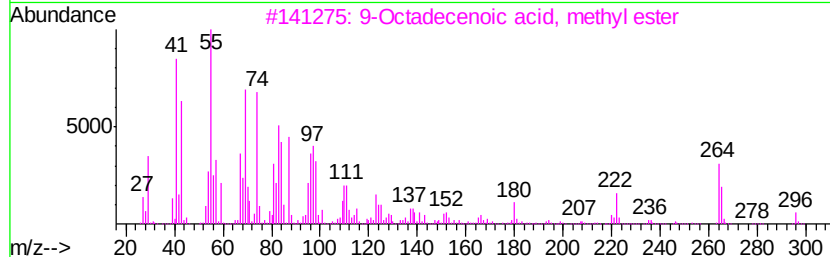
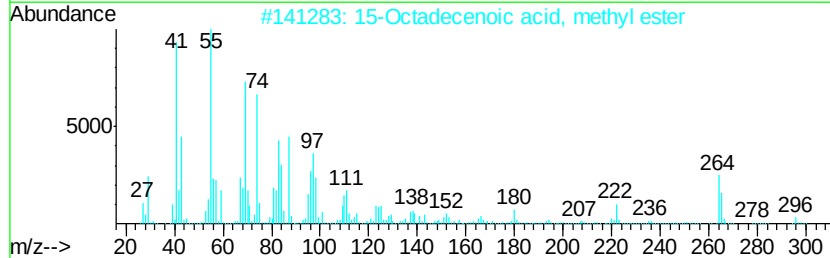
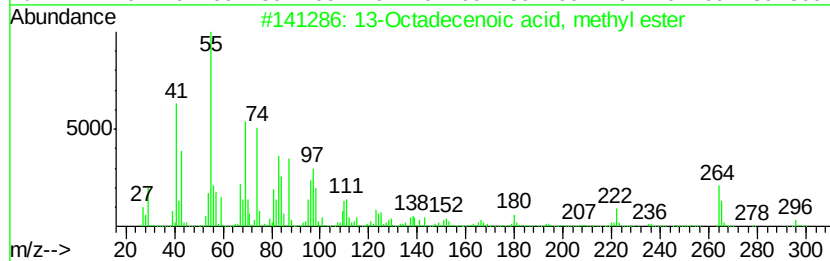
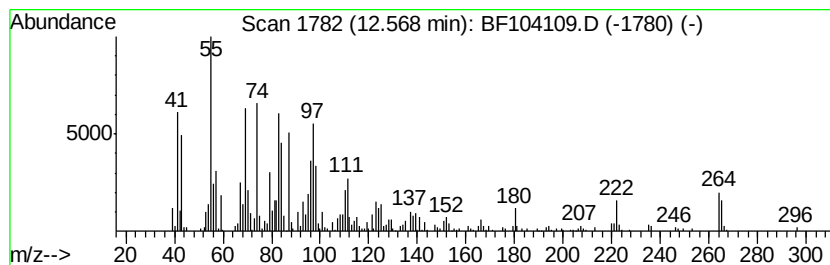
Instrument :
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ClientSampleId :
SU-03-032918

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 13-Octadecenoic acid, methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.57	29.64 ng	926911	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
2	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	99
3	9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
4	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	99
5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF033018\
 Data File : BF104109.D
 Acq On : 30 Mar 2018 22:39
 Operator : JU/SJ
 Sample : J1760-09 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

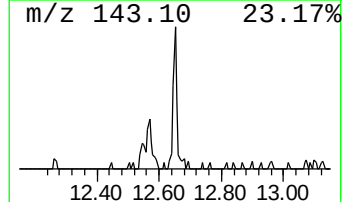
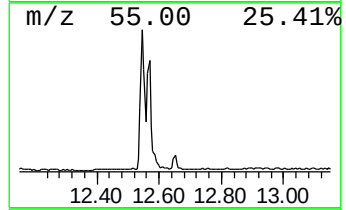
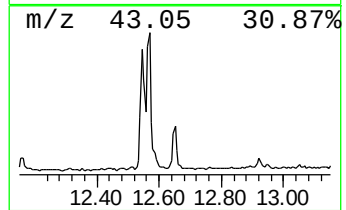
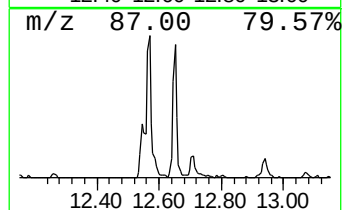
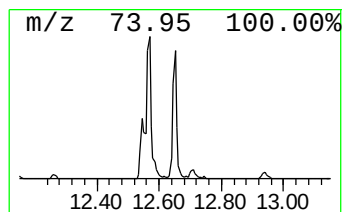
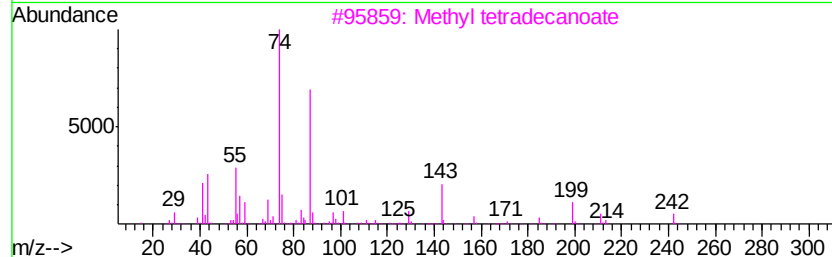
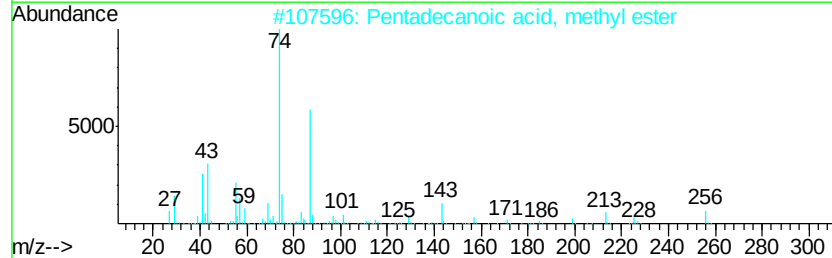
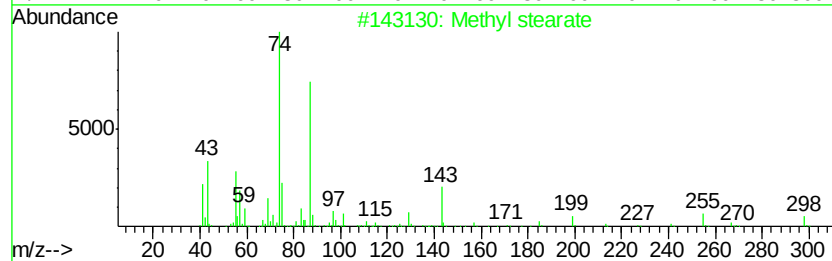
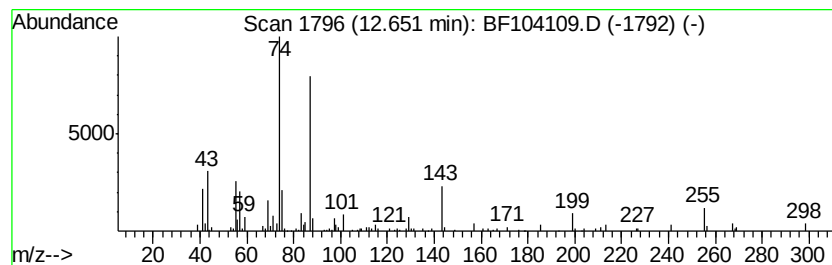
Instrument :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Methyl stearate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.65	5.03 ng	157430	Phenanthrene-d10	11.49	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	95
2	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87
3	Methyl tetradecanoate	242	C15H30O2	000124-10-7	80
4	Methyl 13-methyltetradecanoate	256	C16H32O2	1000336-31-4	72
5	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	72



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF033018\
Data File : BF104109.D
Acq On : 30 Mar 2018 22:39
Operator : JU/SJ
Sample : J1760-09 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

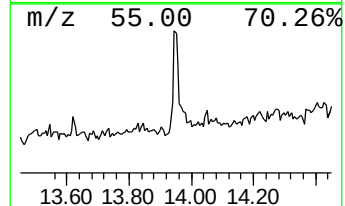
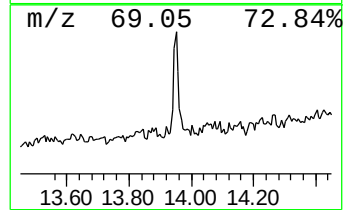
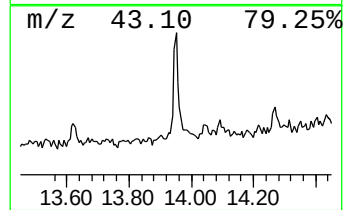
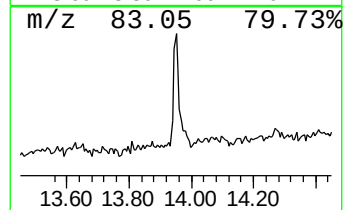
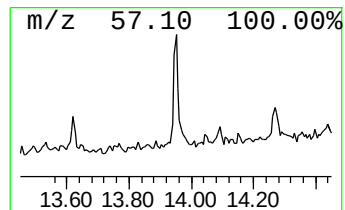
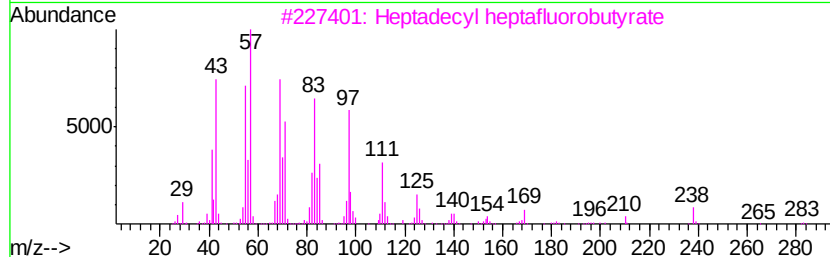
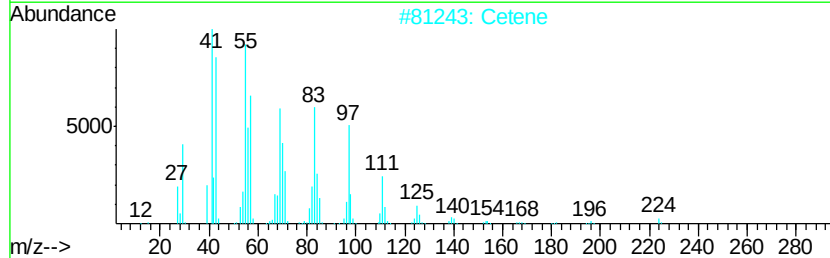
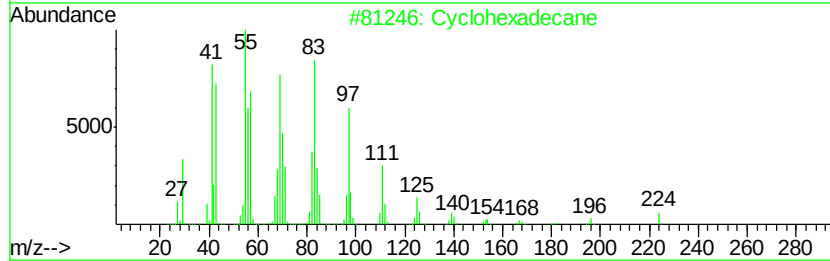
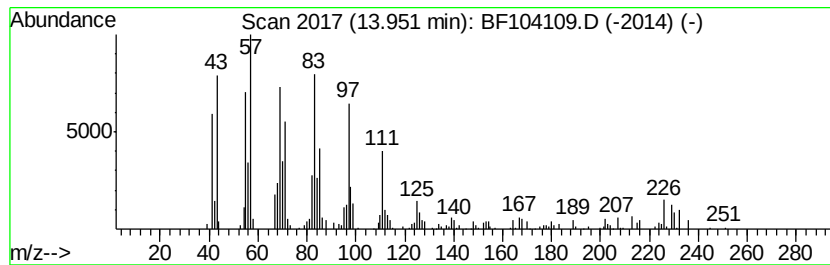
Instrument :
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ClientSampled :
SU-03-032918

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclohexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	3.04 ng	124728	Chrysene-d12	14.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 97
2		Cetene	224 C16H32	000629-73-2 96
3		Heptadecyl heptafluorobutvrate	452 C21H35F7O2	959085-66-6 93
4		Heneicosyl heptafluorobutvrate	508 C25H43F7O2	1000351-83-8 90
5		Trichloroacetic acid, pentadecyl...	372 C17H31Cl3O2	074339-53-0 87



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF033018\
Data File : BF104109.D
Acq On : 30 Mar 2018 22:39
Operator : JU/SJ
Sample : J1760-09 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-03-032918

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032818.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
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Tridecane	8.80	2.8	ng	103700	2	8.24	739408	20.0
Tetradecane	9.90	2.0	ng	89536	3	9.99	873607	20.0
Heptadecane	10.87	3.6	ng	111471	4	11.49	625436	20.0
Hexadecanoic acid...	11.86	11.1	ng	345648	4	11.49	625436	20.0
9,12-Octadecadien...	12.54	42.0	ng	1312990	4	11.49	625436	20.0
13-Octadecenoic a...	12.57	29.6	ng	926911	4	11.49	625436	20.0
Methyl stearate	12.65	5.0	ng	157430	4	11.49	625436	20.0
Cyclohexadecane	13.95	3.0	ng	124728	5	14.14	819304	20.0