

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108096.D
 Acq On : 10 Aug 2018 20:51
 Operator : JU/SJ
 Sample : J4272-15 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-080918-H

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-BF080818.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	4.713	398	404	409	rBV4	30840	45212	1.37%	0.096%
2	5.248	491	495	500	rVB	590827	600223	18.13%	1.272%
3	5.637	556	561	564	rBV	3445294	3033613	91.64%	6.429%
4	6.631	725	730	734	rBV	3654968	3283173	99.18%	6.958%
5	6.784	752	756	760	rBV	3669275	3179691	96.05%	6.739%
6	6.842	762	766	770	rVB	70645	63641	1.92%	0.135%
7	7.019	792	796	800	rBV	2252843	1784534	53.91%	3.782%
8	7.172	819	822	826	rBV	2625776	2434530	73.54%	5.160%
9	7.578	887	891	894	rBV	2418154	1968644	59.47%	4.172%
10	8.301	1010	1014	1020	rVB	2539024	2216459	66.95%	4.697%
11	8.878	1109	1112	1117	rVB3	38989	42171	1.27%	0.089%
12	9.307	1183	1185	1188	rVB	53949	41281	1.25%	0.087%
13	9.378	1193	1197	1200	rBV	4187025	3310379	100.00%	7.016%
14	9.448	1206	1209	1212	rBV3	76382	78276	2.36%	0.166%
15	9.630	1238	1240	1246	rBV3	107510	110094	3.33%	0.233%
16	9.766	1260	1263	1266	rBV	403079	338811	10.23%	0.718%
17	9.925	1288	1290	1293	rBV	69411	60657	1.83%	0.129%
18	9.978	1296	1299	1303	rVB	146221	121628	3.67%	0.258%
19	10.066	1311	1314	1318	rVV	2264637	1977442	59.73%	4.191%
20	10.101	1318	1320	1323	rVB	55596	47148	1.42%	0.100%
21	10.266	1346	1348	1351	rVB2	49750	44370	1.34%	0.094%
22	10.301	1351	1354	1359	rBV4	58224	83100	2.51%	0.176%
23	10.383	1365	1368	1370	rBV3	64410	83993	2.54%	0.178%
24	10.477	1381	1384	1387	rBV	197407	158147	4.78%	0.335%
25	10.613	1405	1407	1410	rVB2	60412	52368	1.58%	0.111%
26	10.701	1417	1422	1424	rBV	175131	189640	5.73%	0.402%
27	10.754	1429	1431	1434	rBV3	53422	55311	1.67%	0.117%
28	10.860	1445	1449	1453	rBV	1682979	1558869	47.09%	3.304%
29	10.966	1462	1467	1471	rBV2	239192	393259	11.88%	0.833%
30	11.148	1496	1498	1500	rBV2	43147	42482	1.28%	0.090%
31	11.283	1518	1521	1522	rBV2	44399	44962	1.36%	0.095%
32	11.407	1539	1542	1544	rBV	226338	184545	5.57%	0.391%
33	11.436	1544	1547	1549	rVV	228368	214056	6.47%	0.454%
34	11.460	1549	1551	1554	rVB	75682	64385	1.94%	0.136%

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ClientSampleId :
JC-03-080918-H

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

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35	11.566	1565	1569	1571	rBV	2089632	1758169	53.11%	3.726%
36	11.589	1571	1573	1576	rVV	585323	447686	13.52%	0.949%
37	11.642	1579	1582	1586	rVV	173401	162019	4.89%	0.343%
38	11.677	1586	1588	1592	rVB5	61243	58920	1.78%	0.125%
39	11.713	1592	1594	1598	rBV3	47495	67955	2.05%	0.144%
40	11.789	1604	1607	1610	rBV3	95176	99063	2.99%	0.210%
41	11.836	1612	1615	1617	rVB	185559	162381	4.91%	0.344%
42	11.936	1629	1632	1636	rVB	1171520	903029	27.28%	1.914%
43	12.072	1651	1655	1657	rBV	294728	308895	9.33%	0.655%
44	12.095	1657	1659	1663	rVB	176803	147962	4.47%	0.314%
45	12.183	1671	1674	1676	rBV2	150324	145767	4.40%	0.309%
46	12.242	1681	1684	1687	rBV	191091	148990	4.50%	0.316%
47	12.360	1702	1704	1707	rVB	61208	47614	1.44%	0.101%
48	12.630	1746	1750	1751	rBV	2745321	2401721	72.55%	5.090%
49	12.648	1751	1753	1760	rVB	2824878	2536633	76.63%	5.376%
50	12.730	1765	1767	1771	rVB	690226	606992	18.34%	1.286%
51	12.789	1773	1777	1780	rVV	1126702	1108177	33.48%	2.349%
52	12.819	1780	1782	1785	rVB2	138174	128252	3.87%	0.272%
53	12.971	1806	1808	1810	rVB2	106691	76805	2.32%	0.163%
54	13.019	1810	1816	1821	rVB	1066703	1219587	36.84%	2.585%
55	13.154	1835	1839	1842	rBV	3017775	2566835	77.54%	5.440%
56	13.177	1842	1843	1846	rVB3	83444	56606	1.71%	0.120%
57	13.366	1873	1875	1877	rVB	113270	79561	2.40%	0.169%
58	13.571	1908	1910	1914	rBV3	208507	205328	6.20%	0.435%
59	14.042	1988	1990	1996	rVB5	187548	305218	9.22%	0.647%
60	14.224	2016	2021	2023	rBV2	1813220	2073247	62.63%	4.394%
61	14.248	2023	2025	2031	rVB	439308	407373	12.31%	0.863%
62	15.795	2284	2288	2292	rVB	830889	1046017	31.60%	2.217%

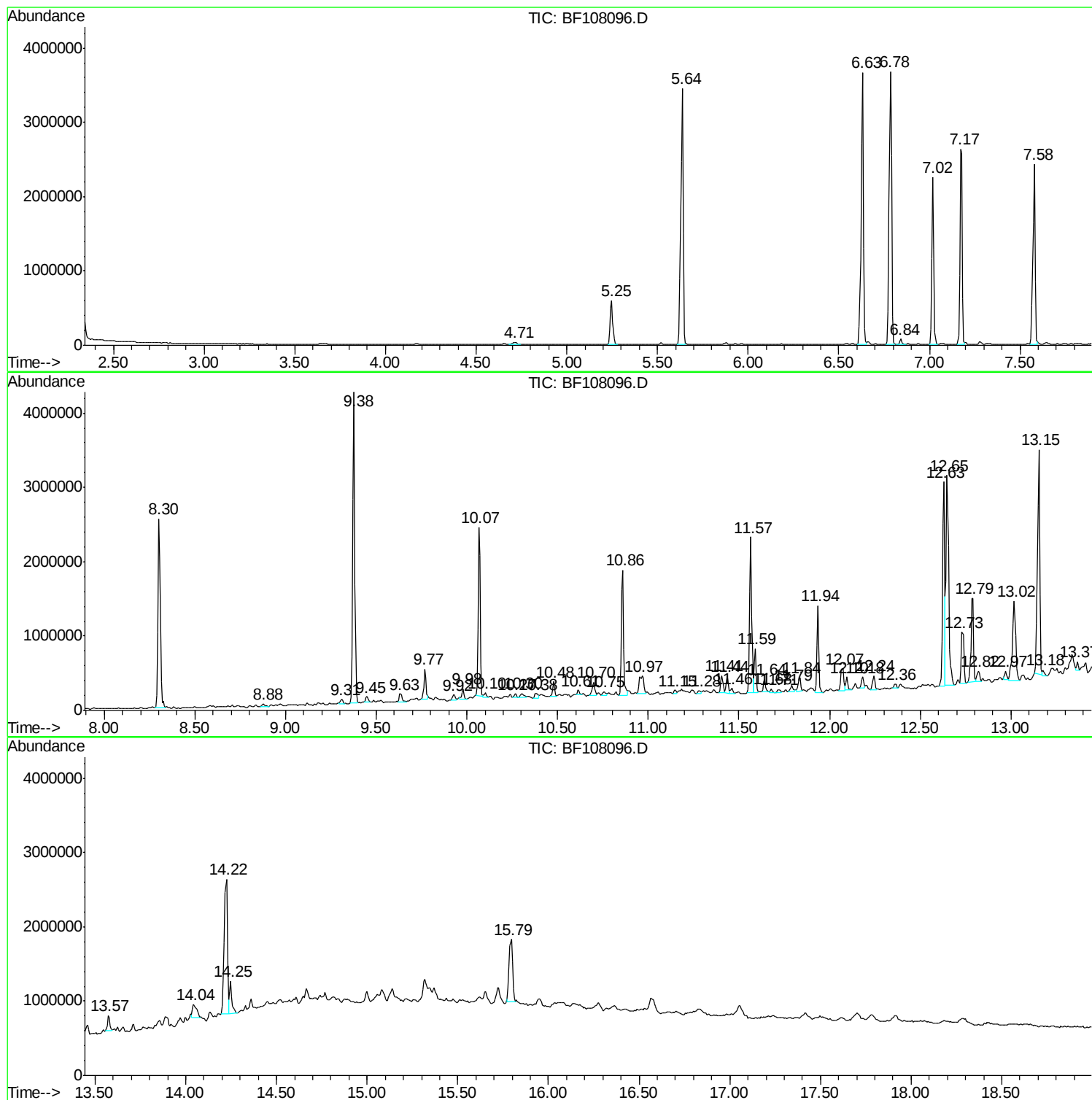
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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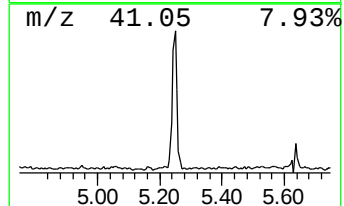
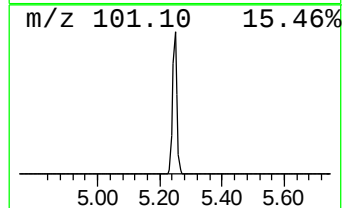
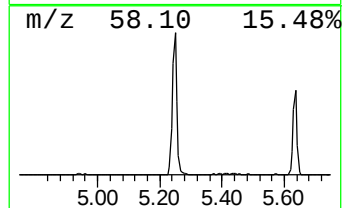
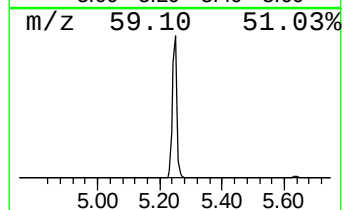
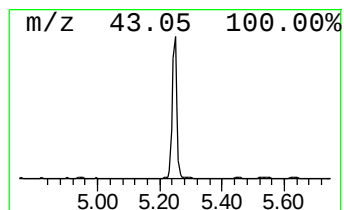
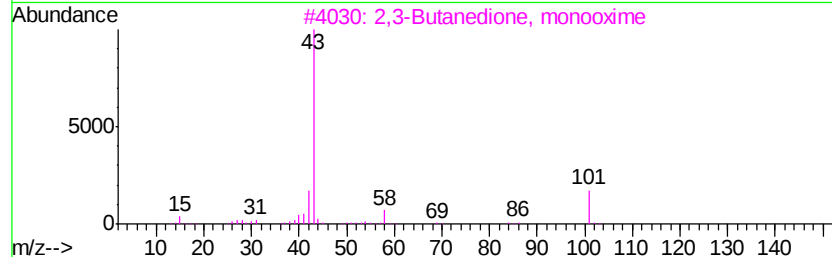
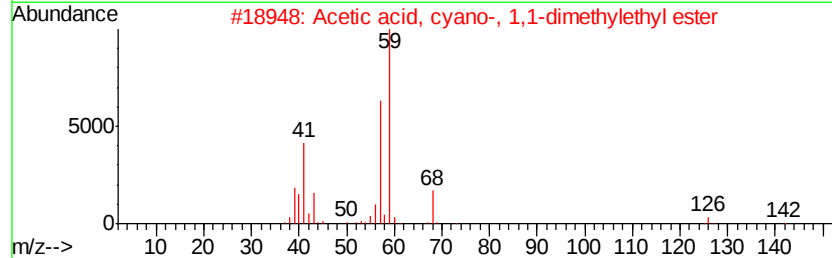
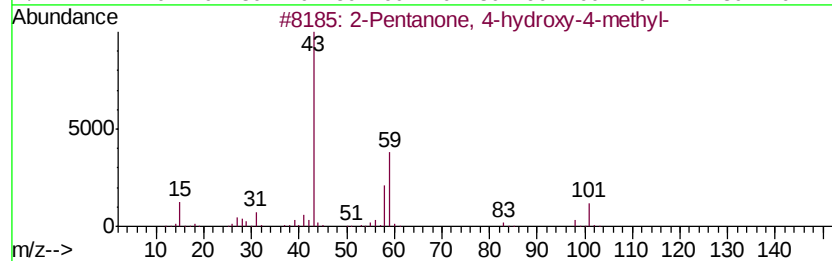
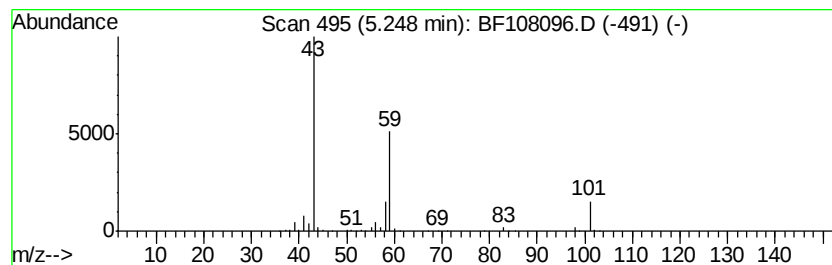
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	6.73 ng	600223	1,4-Dichlorobenzene-d4	7.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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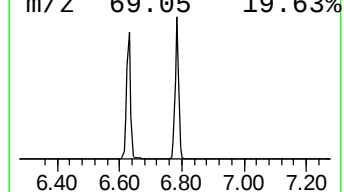
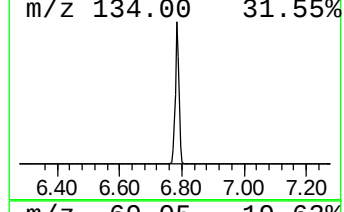
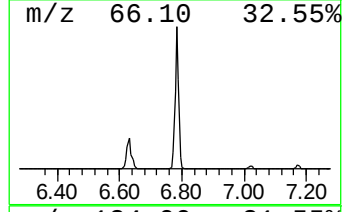
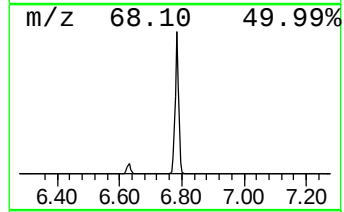
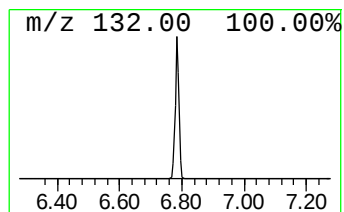
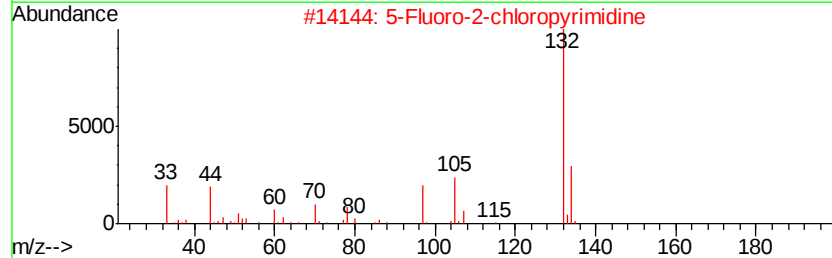
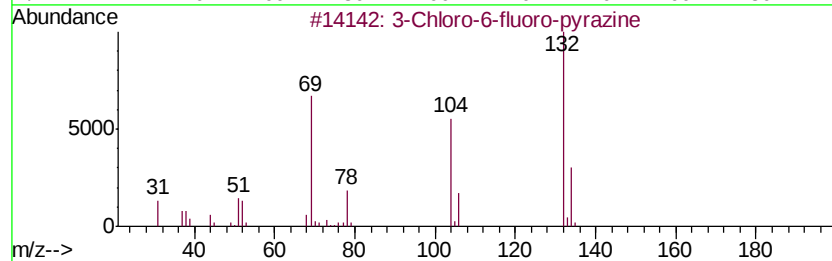
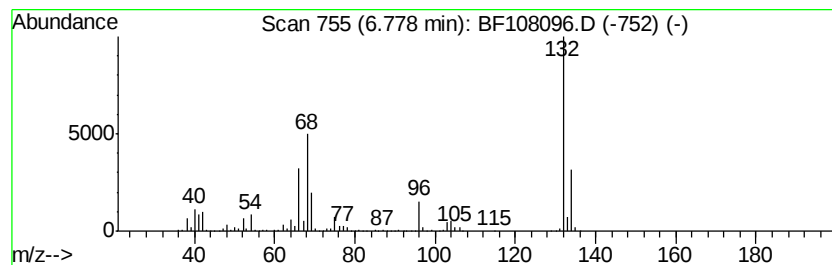
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.78 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	35.64 ng	3179690	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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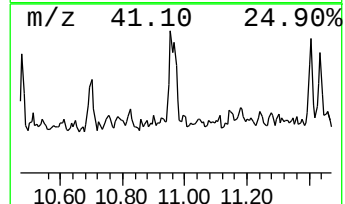
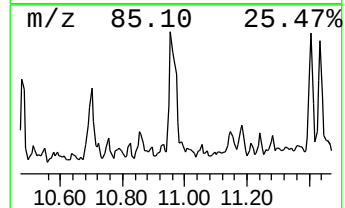
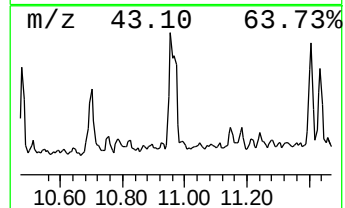
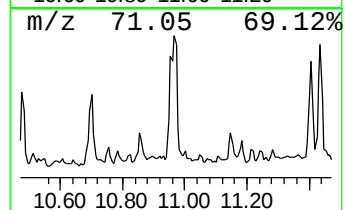
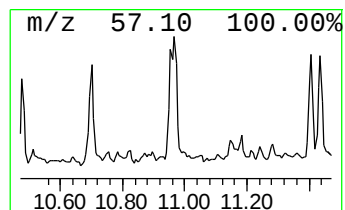
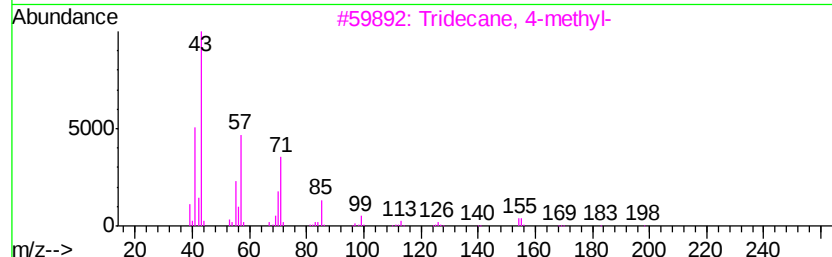
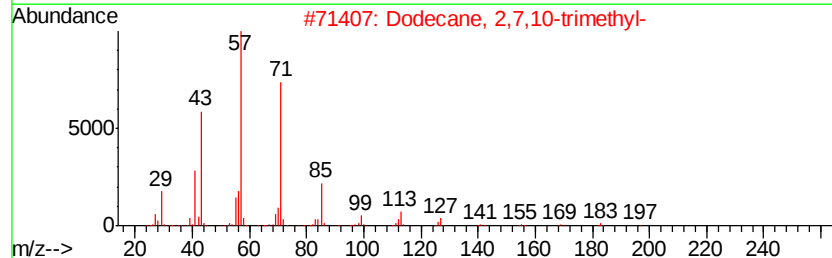
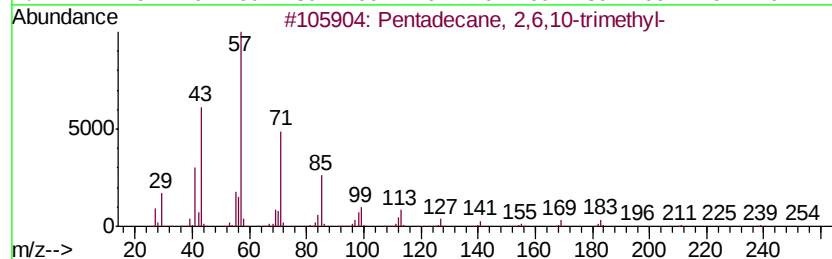
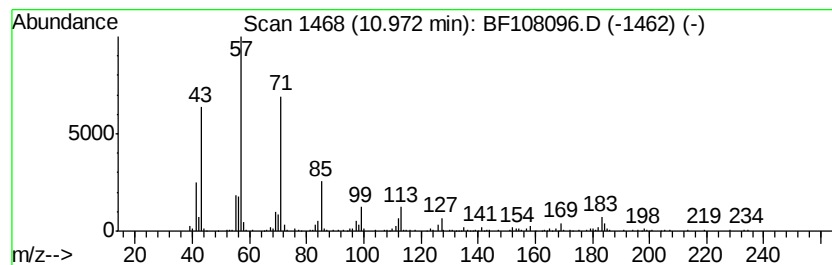
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentadecane, 2,6,10-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	4.47 ng	393259	Phenanthrene-d10	11.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
2		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	83
3		Tridecane, 4-methyl-	198	C14H30	026730-12-1	72
4		Heptacosane	380	C27H56	000593-49-7	64
5		Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64



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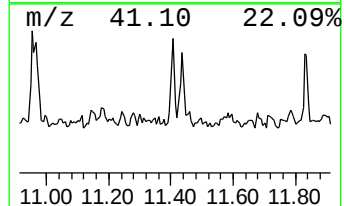
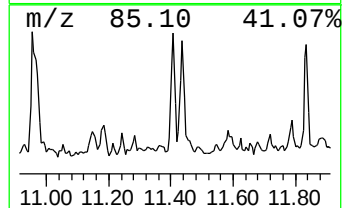
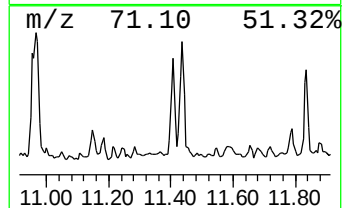
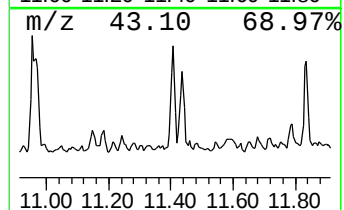
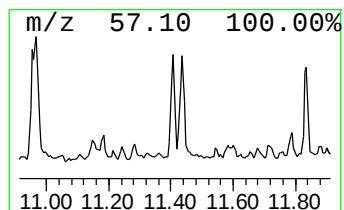
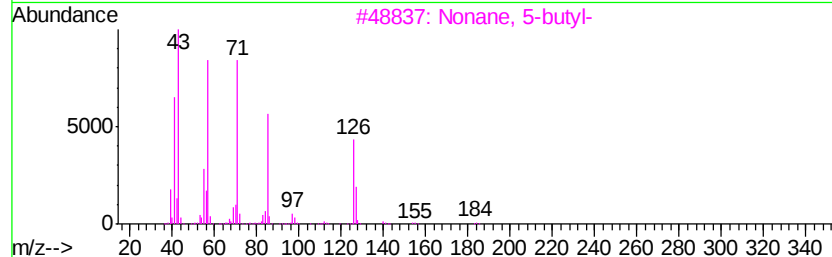
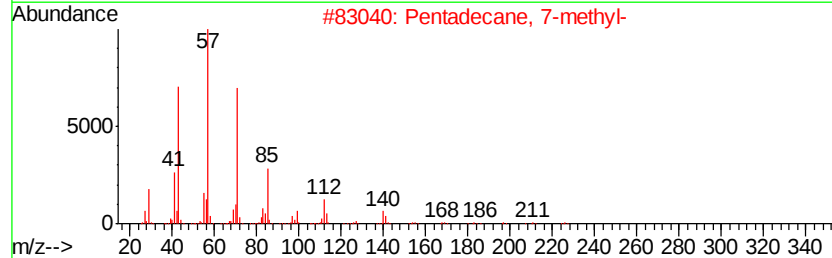
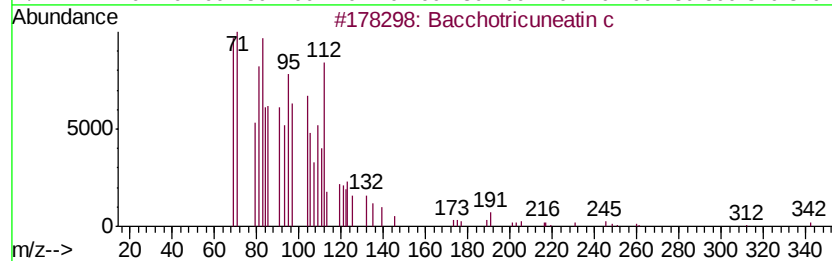
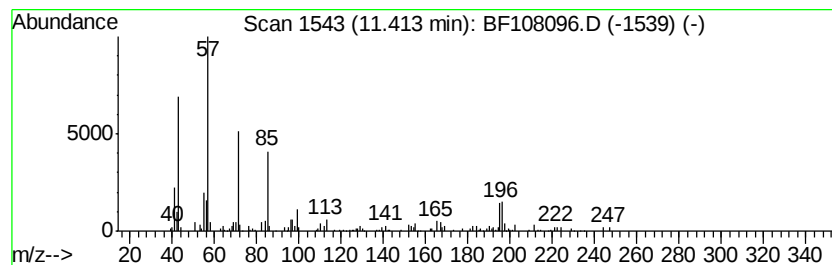
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Bacchotricuneatin c Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.41	2.10 ng	184545	Phenanthrene-d10		11.57
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bacchotricuneatin c	342	C20H22O5	066563-30-2	96
2	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	87
3	Nonane, 5-butyl-	184	C13H28	017312-63-9	86
4	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	83
5	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	81



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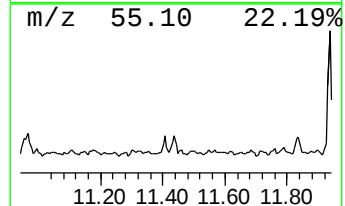
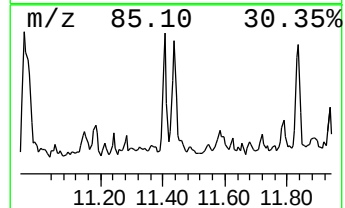
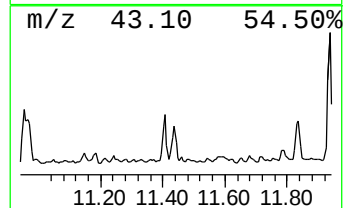
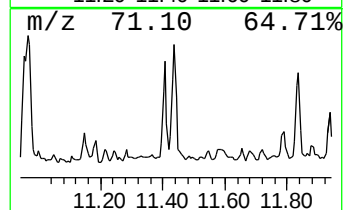
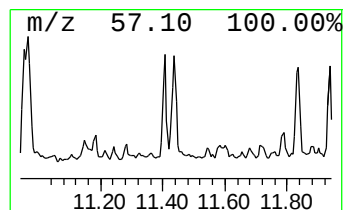
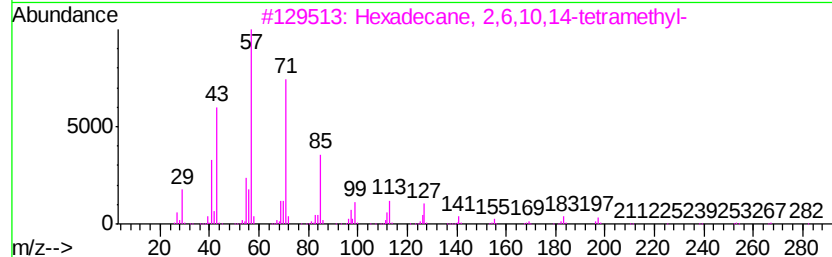
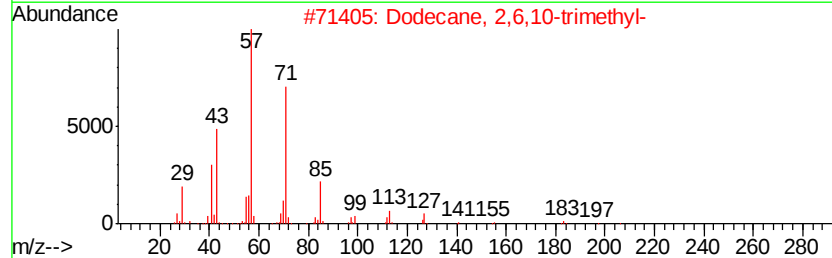
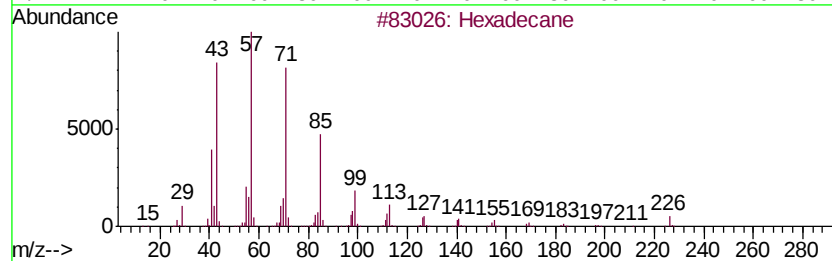
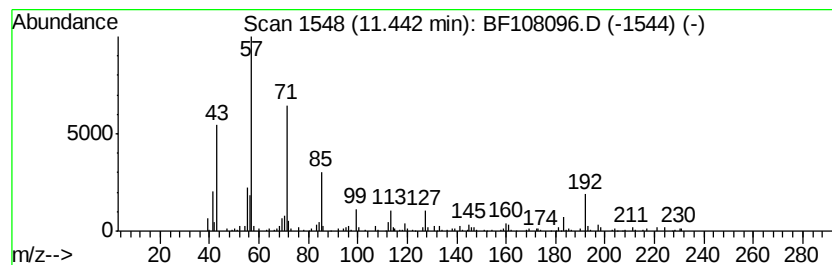
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ClientSampleId :
JC-03-080918-H

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.44	2.43 ng	214056	Phenanthrene-d10	11.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	72
2	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	72
3	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72
4	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	72
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
Data File : BF108096.D
Acq On : 10 Aug 2018 20:51
Operator : JU/SJ
Sample : J4272-15 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

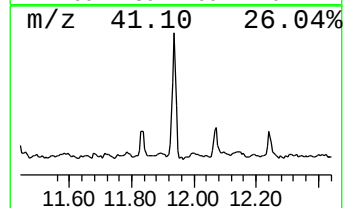
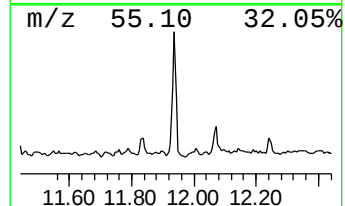
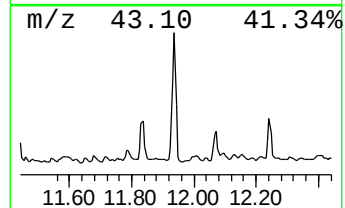
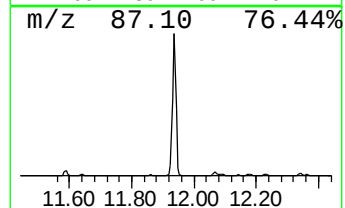
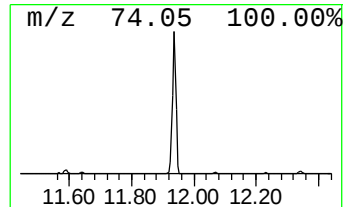
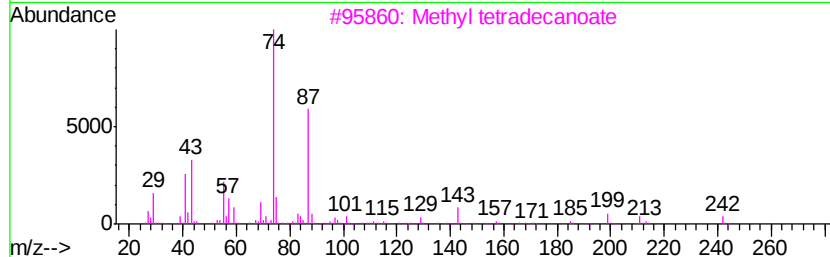
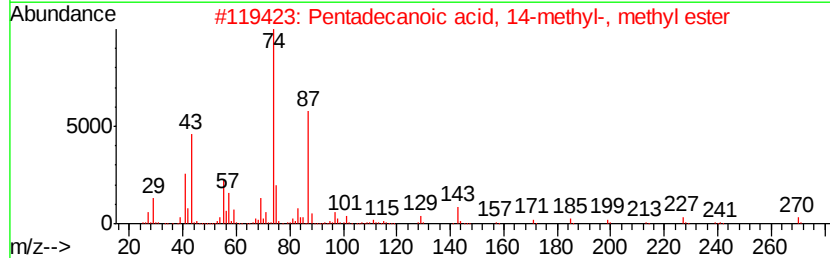
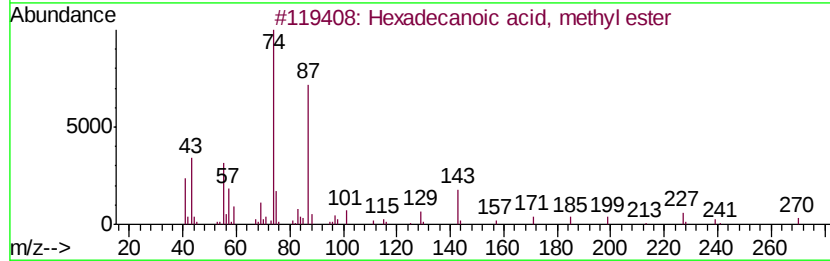
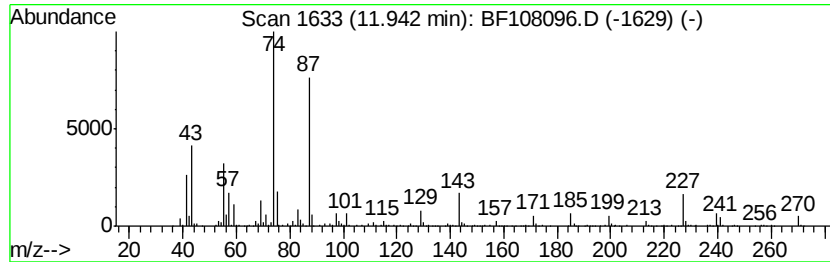
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.	
11.94	10.27 ng	903029	Phenanthrene-d10	11.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3	Methyl tetradecanoate	242	C15H30O2	000124-10-7	93
4	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	87
5	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
Data File : BF108096.D
Acq On : 10 Aug 2018 20:51
Operator : JU/SJ
Sample : J4272-15 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JC-03-080918-H

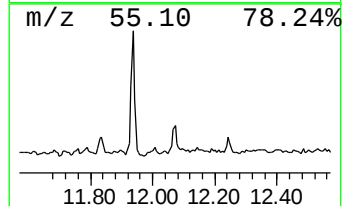
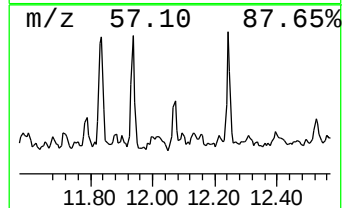
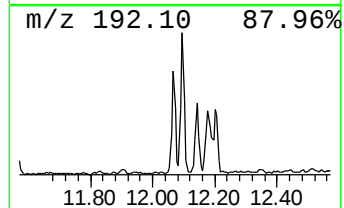
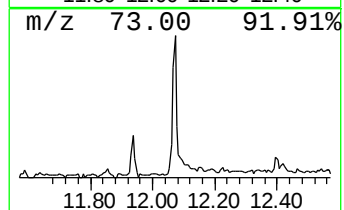
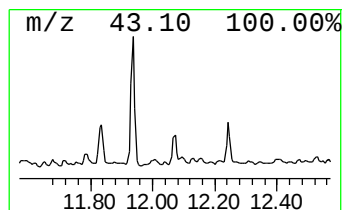
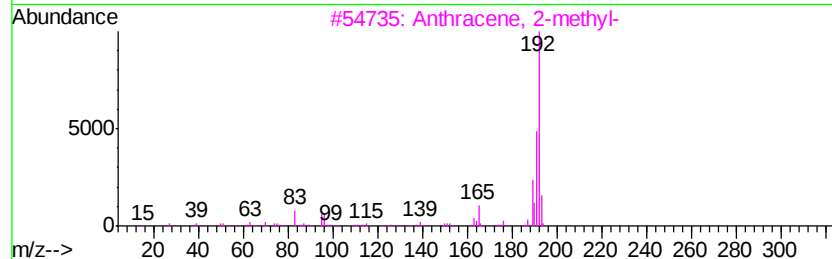
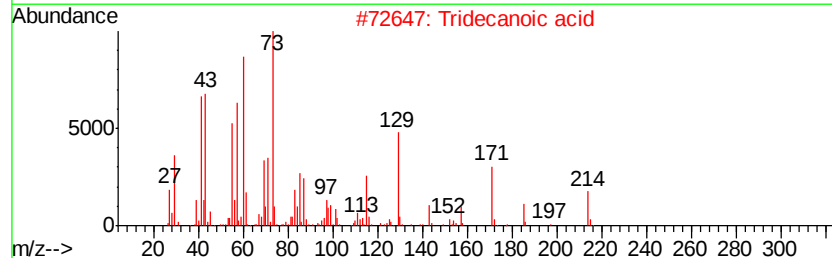
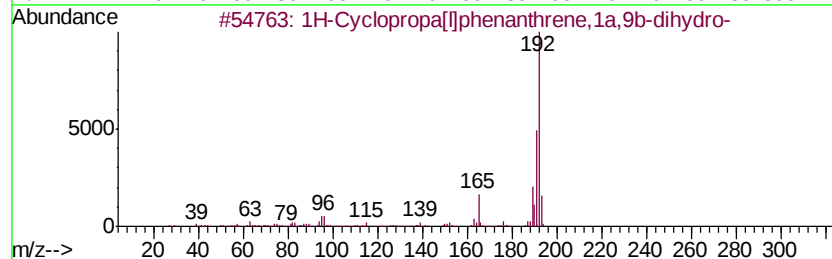
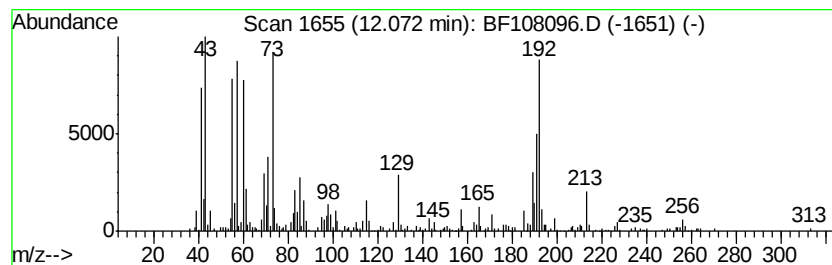
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1H-Cyclopropa[l]phenanthren... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.07	3.51 ng	308895	Phenanthrene-d10	11.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	92
2		Tridecanoic acid	214	C13H26O2	000638-53-9	83
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	78
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	70
5		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
Data File : BF108096.D
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Operator : JU/SJ
Sample : J4272-15 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JC-03-080918-H

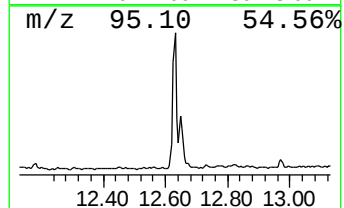
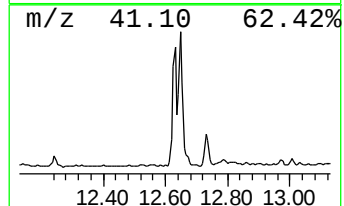
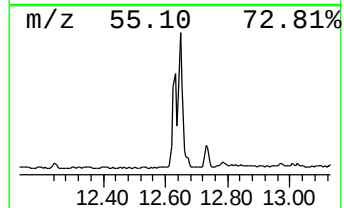
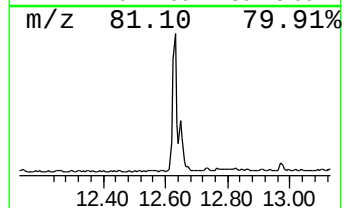
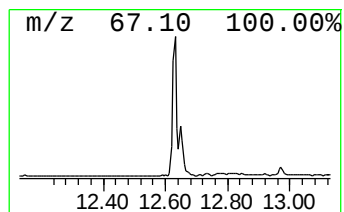
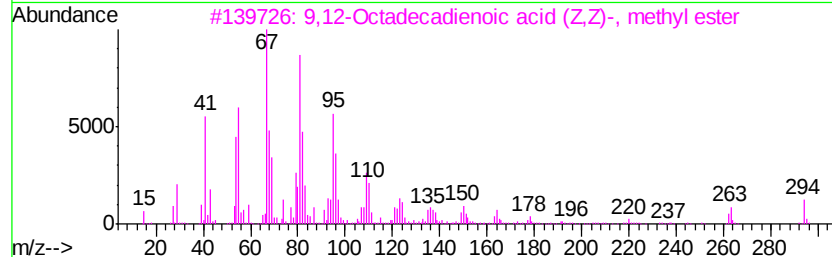
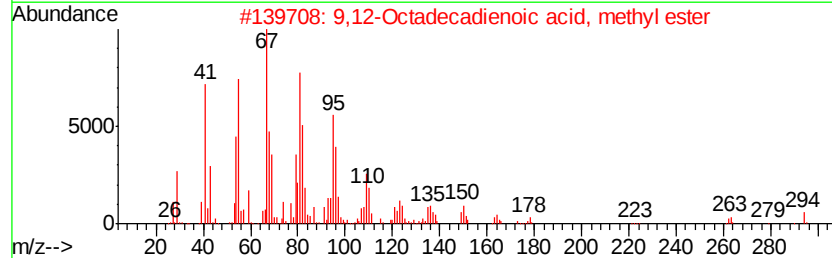
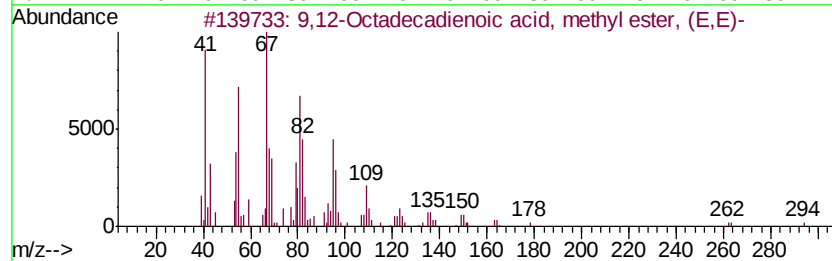
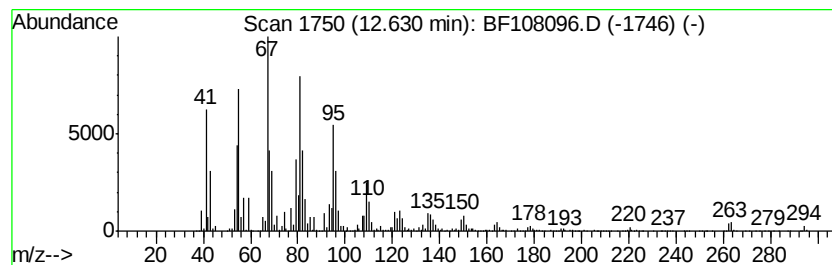
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,12-Octadecadienoic acid, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	27.32 ng	2401720	Phenanthrene-d10	11.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
3		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
4		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
Data File : BF108096.D
Acq On : 10 Aug 2018 20:51
Operator : JU/SJ
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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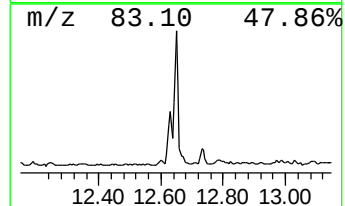
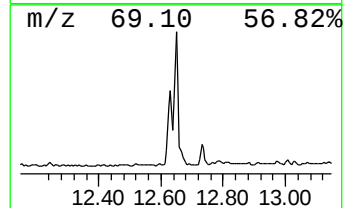
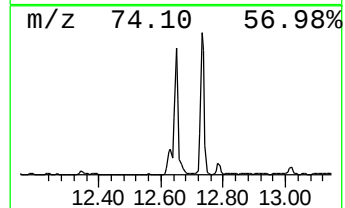
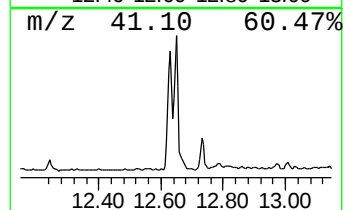
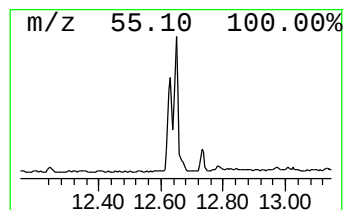
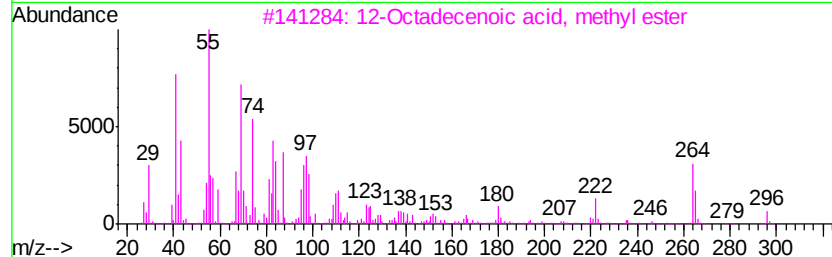
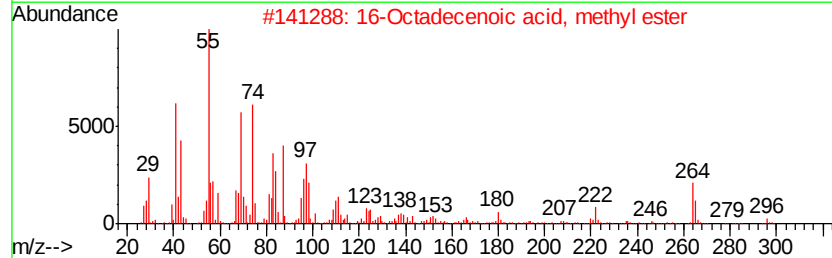
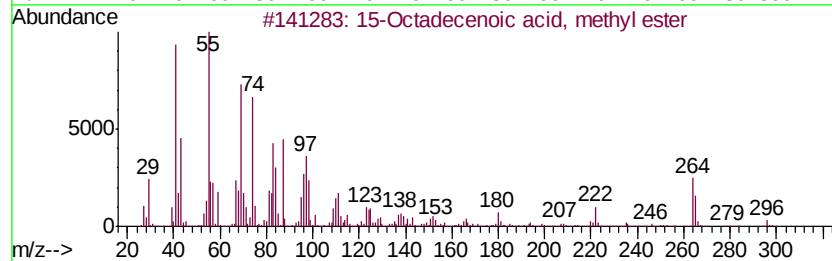
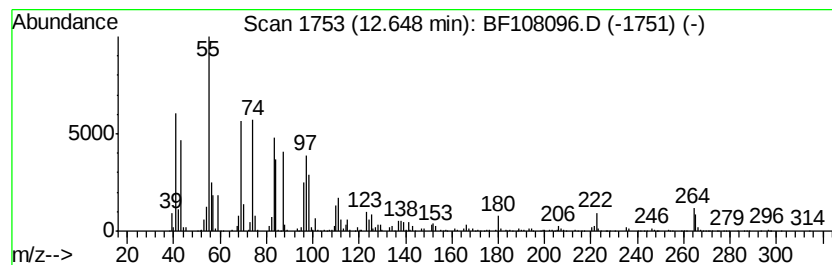
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 15-Octadecenoic acid, methyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	28.86 ng	2536630	Phenanthrene-d10	11.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	96
2		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	96
3		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	96
4		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	95
5		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
Data File : BF108096.D
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Sample : J4272-15 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

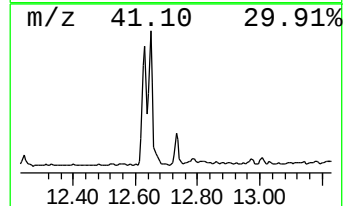
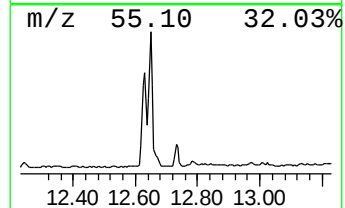
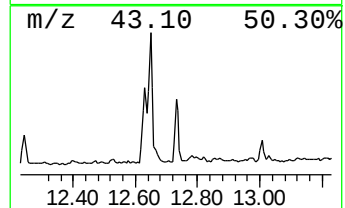
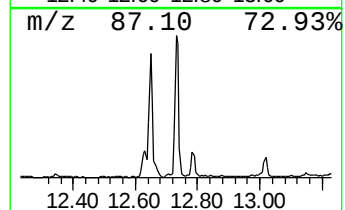
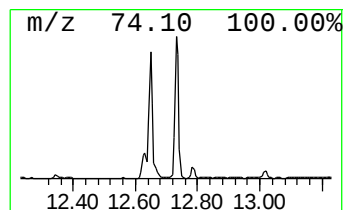
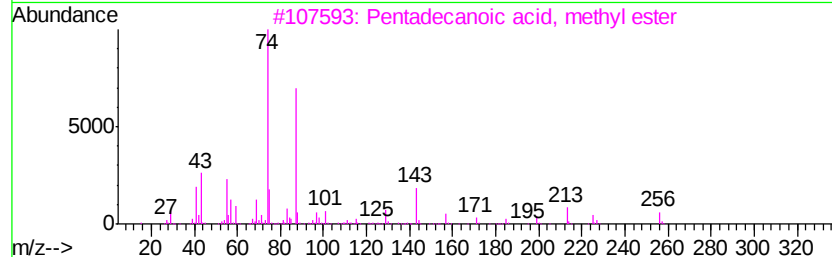
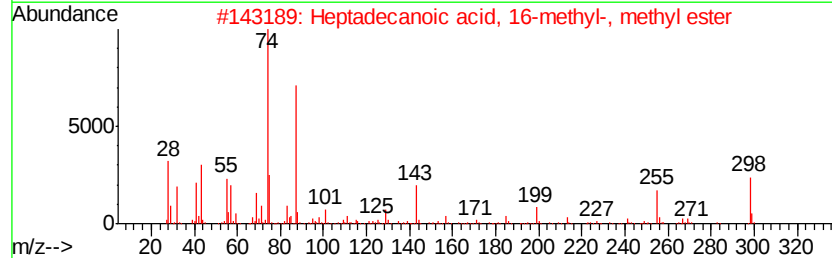
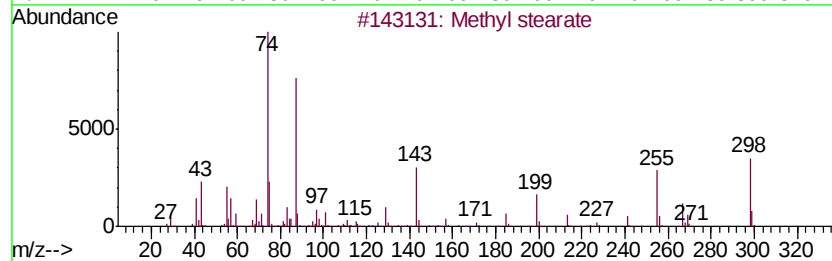
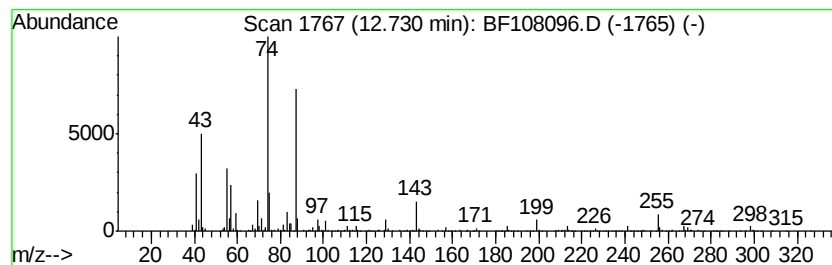
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methyl stearate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.73	6.90 ng	606992	Phenanthrene-d10		11.57	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl stearate	298	C19H38O2	000112-61-8	98
2		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	97
3		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	91
4		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	91
5		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
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JC-03-080918-H

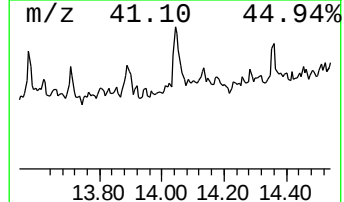
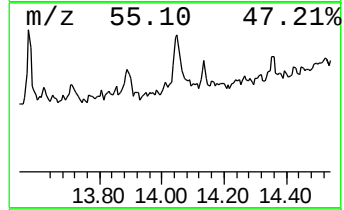
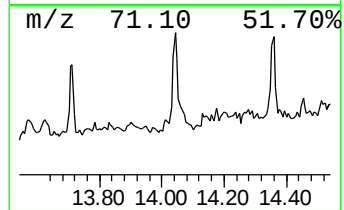
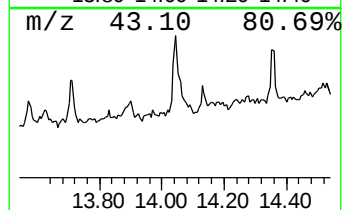
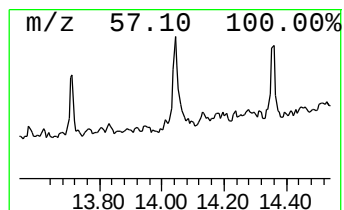
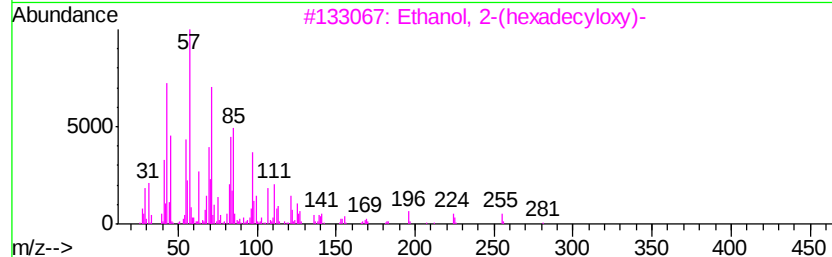
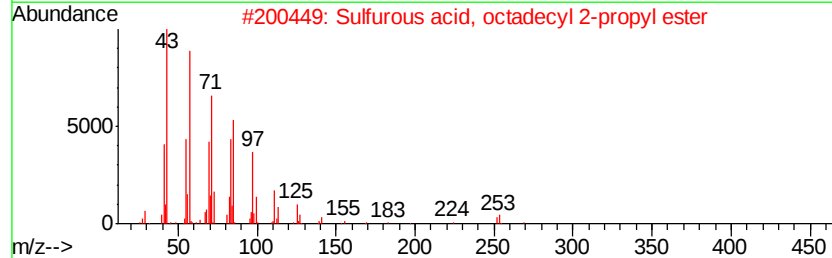
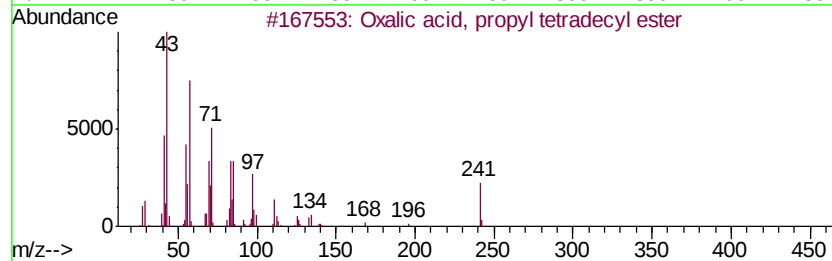
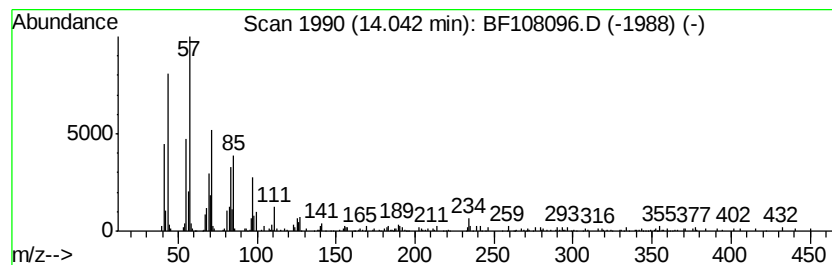
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Oxalic acid, propyl tetradecyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	2.94 ng	305218	Chrysene-d12	14.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxalic acid, propyl tetradecyl e...	328	C19H36O4	1000309-26-7	87
2		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	83
3		Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	80
4		1-Chloroeicosane	316	C20H41Cl	042217-02-7	76
5		Octacosane	394	C28H58	000630-02-4	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081018\
 Data File : BF108096.D
 Acq On : 10 Aug 2018 20:51
 Operator : JU/SJ
 Sample : J4272-15 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-080918-H

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.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
2-Pentanone, 4-hy...	5.25	6.7	ng	600223	1	7.02	1784530	20.0
unknown6.78	6.78	35.6	ng	3179690	1	7.02	1784530	20.0
Pentadecane, 2,6,...	10.97	4.5	ng	393259	4	11.57	1758170	20.0
Bacchotricuneatin c	11.41	2.1	ng	184545	4	11.57	1758170	20.0
Hexadecane	11.44	2.4	ng	214056	4	11.57	1758170	20.0
Hexadecanoic acid...	11.94	10.3	ng	903029	4	11.57	1758170	20.0
1H-Cyclopropa[l]p...	12.07	3.5	ng	308895	4	11.57	1758170	20.0
9,12-Octadecadien...	12.63	27.3	ng	2401720	4	11.57	1758170	20.0
15-Octadecenoic a...	12.65	28.9	ng	2536630	4	11.57	1758170	20.0
Methyl stearate	12.73	6.9	ng	606992	4	11.57	1758170	20.0
Oxalic acid, prop...	14.04	2.9	ng	305218	5	14.22	2073250	20.0