

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
 Data File : BF098967.D
 Acq On : 30 Sep 2017 17:19
 Operator : SJ/JU
 Sample : I5511-01 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-092817

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-BF092317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.522	580	584	592	rBV	338238	299794	2.55%	0.743%
2	6.528	751	755	759	rVB	362077	299302	2.55%	0.742%
3	6.675	777	780	784	rVB	367017	328789	2.80%	0.815%
4	6.916	817	821	824	rVB	991399	860010	7.31%	2.131%
5	7.069	842	847	850	rBV	321951	286188	2.43%	0.709%
6	7.469	912	915	918	rVV	236078	230056	1.96%	0.570%
7	7.498	918	920	924	rVB	206334	190824	1.62%	0.473%
8	8.169	1030	1034	1036	rVV	352825	360376	3.06%	0.893%
9	8.198	1036	1039	1041	rVV	1360162	1164570	9.90%	2.886%
10	8.245	1044	1047	1050	rVB2	168495	165610	1.41%	0.410%
11	8.481	1083	1087	1091	rBV5	147189	167620	1.43%	0.415%
12	8.604	1105	1108	1111	rVB2	144699	124363	1.06%	0.308%
13	8.698	1120	1124	1126	rBV	201877	164373	1.40%	0.407%
14	8.781	1132	1138	1142	rBV	388873	397423	3.38%	0.985%
15	9.016	1175	1178	1180	rBV2	157958	124964	1.06%	0.310%
16	9.139	1196	1199	1204	rBV3	127883	154239	1.31%	0.382%
17	9.210	1208	1211	1213	rVV	164660	147918	1.26%	0.367%
18	9.269	1217	1221	1225	rVB	448632	399775	3.40%	0.991%
19	9.345	1230	1234	1237	rBV	507604	461353	3.92%	1.143%
20	9.422	1244	1247	1249	rVB	133527	120366	1.02%	0.298%
21	9.528	1262	1265	1271	rVB5	118007	132743	1.13%	0.329%
22	9.604	1271	1278	1279	rBV5	102341	187394	1.59%	0.464%
23	9.663	1284	1288	1290	rVV3	255154	301339	2.56%	0.747%
24	9.875	1321	1324	1328	rVB	585114	471844	4.01%	1.169%
25	9.951	1332	1337	1341	rVB	1126925	1087976	9.25%	2.696%
26	10.192	1375	1378	1383	rBV3	158421	208867	1.78%	0.518%
27	10.375	1404	1409	1411	rBV	616768	549523	4.67%	1.362%
28	10.498	1424	1430	1433	rBV6	60205	137960	1.17%	0.342%
29	10.592	1441	1446	1448	rBV	239791	289791	2.46%	0.718%
30	10.845	1486	1489	1494	rBV	696031	733262	6.24%	1.817%
31	11.292	1562	1565	1568	rBV	607028	482339	4.10%	1.195%
32	11.322	1568	1570	1577	rVB	228625	250123	2.13%	0.620%
33	11.439	1586	1590	1593	rBV	987117	879336	7.48%	2.179%
34	11.722	1635	1638	1640	rBV	426787	357805	3.04%	0.887%

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 Sampling : 1 Min Area: 3 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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35	11.828	1651	1656	1659	rBV	5060237	5187178	44.12%	12.854%
36	11.957	1675	1678	1685	rVB5	89563	136347	1.16%	0.338%
37	12.127	1704	1707	1716	rVB	361872	368660	3.14%	0.914%
38	12.527	1767	1775	1777	rBV2	5398891	11758279	100.00%	29.138%
39	12.616	1787	1790	1794	rVB	2929864	2635730	22.42%	6.532%
40	12.663	1794	1798	1799	rBV2	541131	697527	5.93%	1.729%
41	12.680	1799	1801	1807	rVV	1041287	1142124	9.71%	2.830%
42	12.851	1827	1830	1833	rBV	265371	233243	1.98%	0.578%
43	12.886	1833	1836	1839	rBV	307519	267184	2.27%	0.662%
44	13.022	1856	1859	1862	rVV	307697	294526	2.50%	0.730%
45	13.210	1887	1891	1894	rVB2	206727	304933	2.59%	0.756%
46	13.245	1894	1897	1898	rBV	175844	149471	1.27%	0.370%
47	13.263	1898	1900	1906	rVB2	270701	266752	2.27%	0.661%
48	13.339	1910	1913	1916	rBV	340234	269681	2.29%	0.668%
49	13.769	1982	1986	1992	rVB2	194479	292163	2.48%	0.724%
50	14.004	2023	2026	2029	rBV	355819	306345	2.61%	0.759%
51	14.080	2035	2039	2042	rBV	862838	845335	7.19%	2.095%
52	14.521	2111	2114	2123	rVB5	253334	468867	3.99%	1.162%
53	14.968	2186	2190	2202	rVB	912016	1538208	13.08%	3.812%
54	15.574	2289	2293	2298	rBV2	511514	672801	5.72%	1.667%

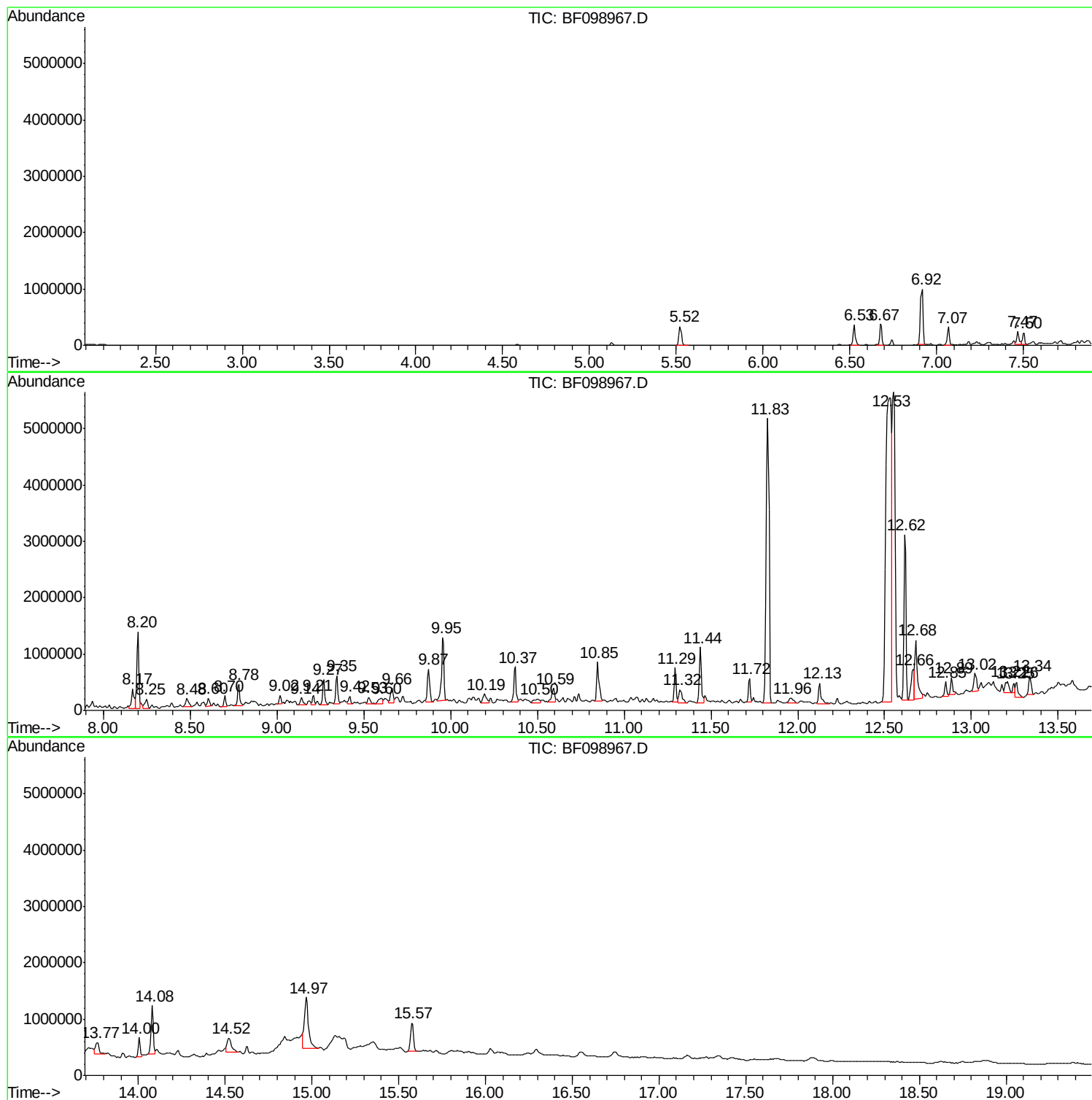
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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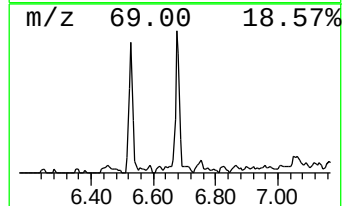
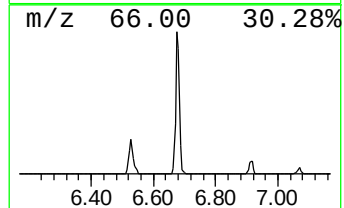
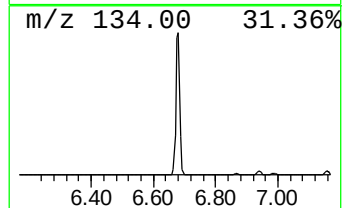
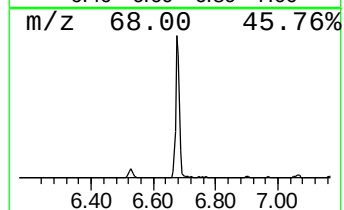
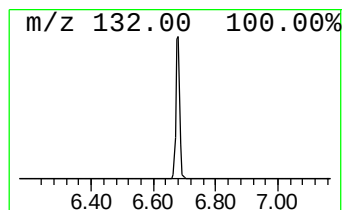
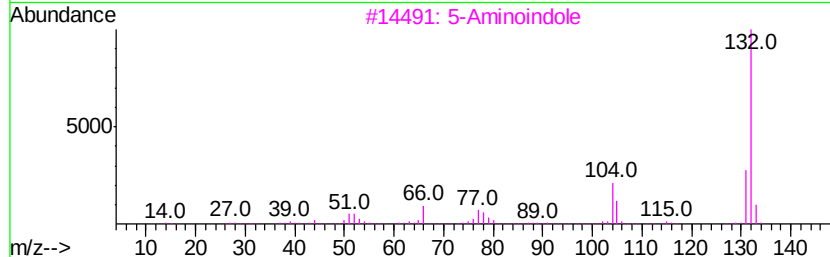
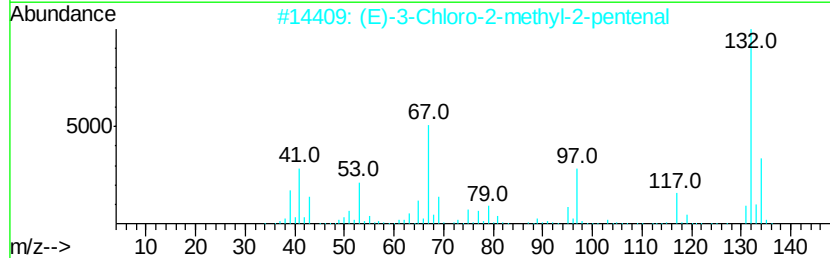
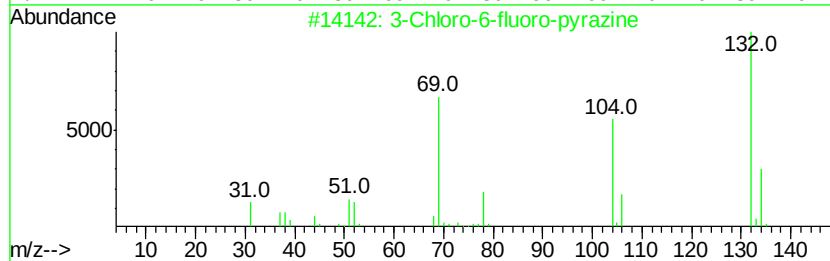
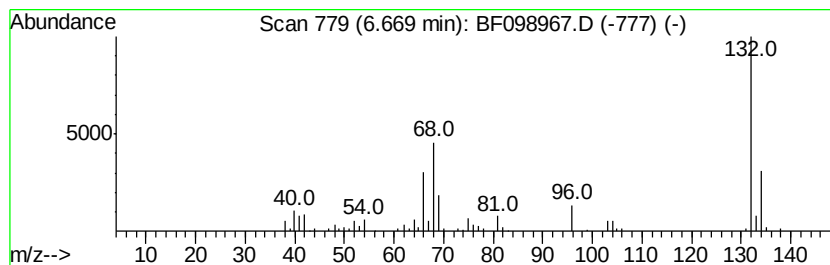
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.67 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	7.65 ng	328789	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3			5-Aminoindole	132	C8H8N2	005192-03-0	10
4			Amphetaminil	250	C17H18N2	017590-01-1	10
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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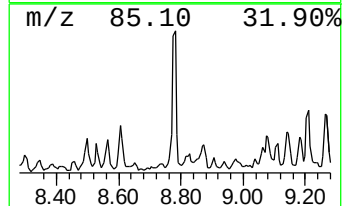
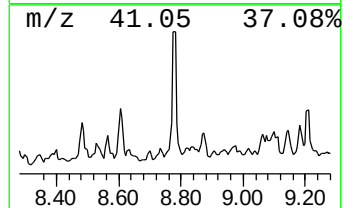
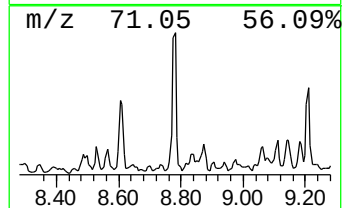
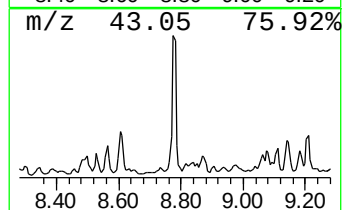
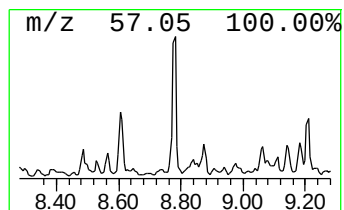
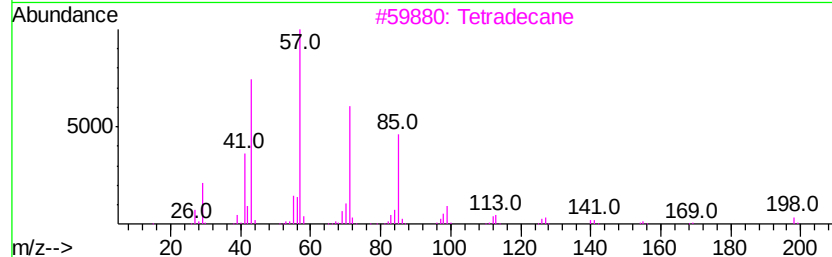
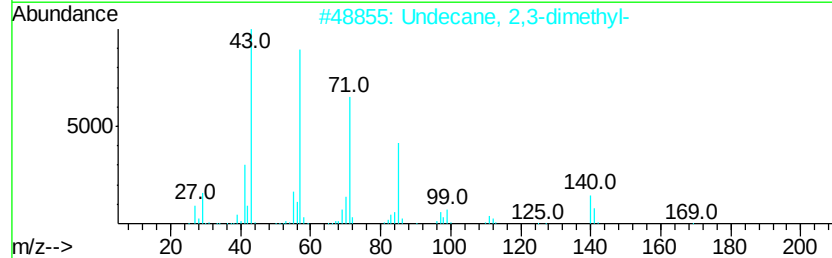
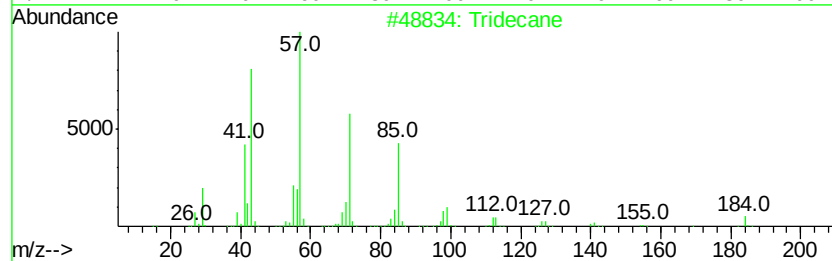
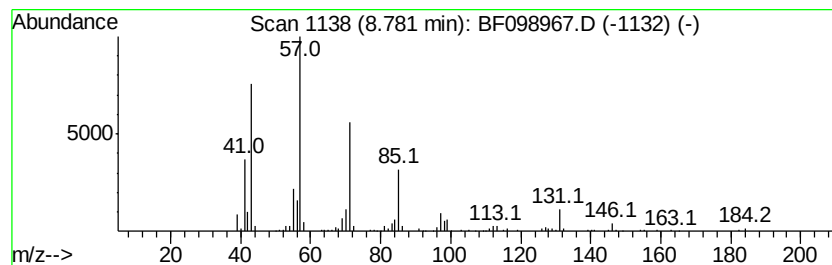
Instrument :
BNA_F
ClientSampleId :
EO-01-092817

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.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 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.78	6.83 ng	397423	Naphthalene-d8	8.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	95
2	Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	80
3	Tetradecane	198	C14H30	000629-59-4	64
4	Hexadecane	226	C16H34	000544-76-3	64
5	Pentadecane	212	C15H32	000629-62-9	64



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ALS Vial : 11 Sample Multiplier: 1

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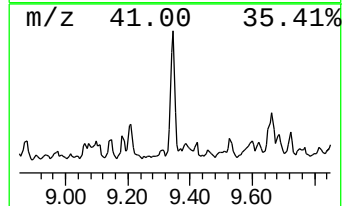
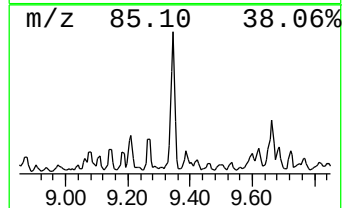
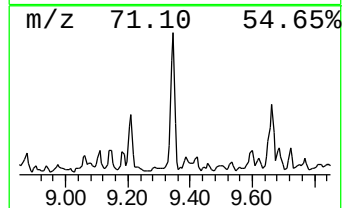
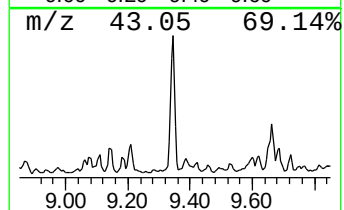
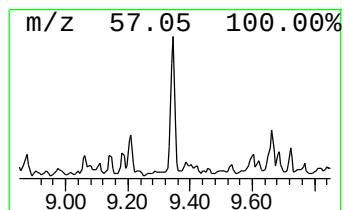
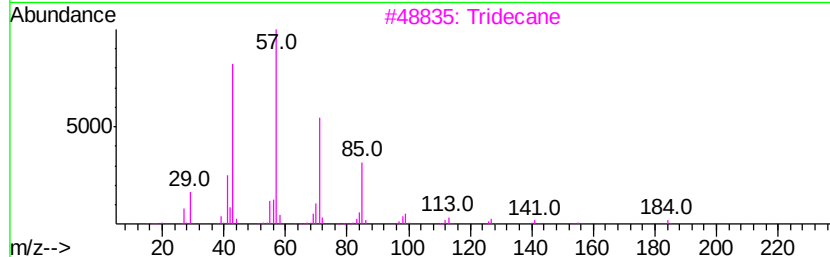
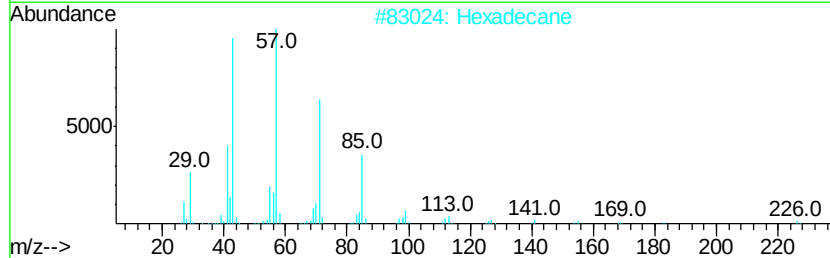
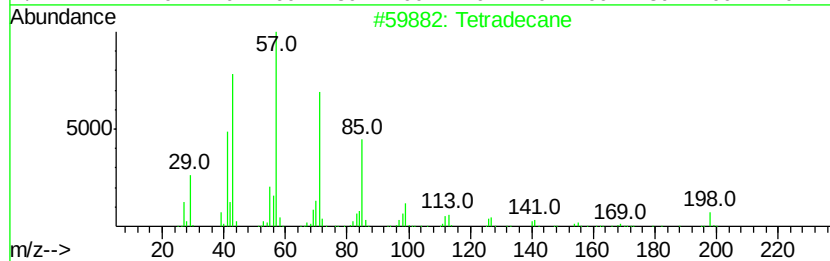
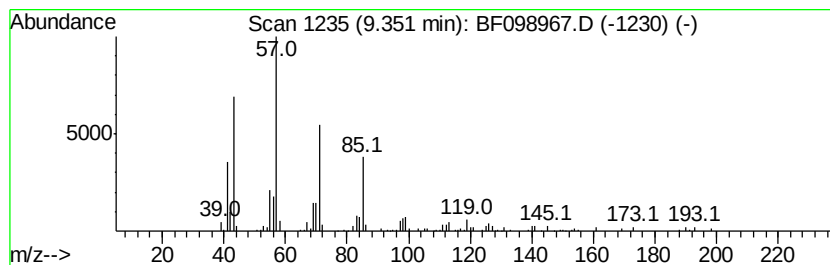
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.35	8.48 ng	461353	Acenaphthene-d10	9.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	95
2		Hexadecane	226	C16H34	000544-76-3	86
3		Tridecane	184	C13H28	000629-50-5	86
4		Tridecane, 6-methyl-	198	C14H30	013287-21-3	76
5		Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	64



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ALS Vial : 11 Sample Multiplier: 1

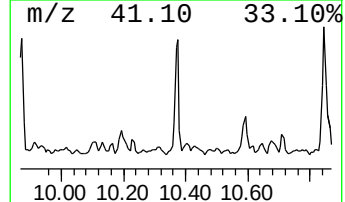
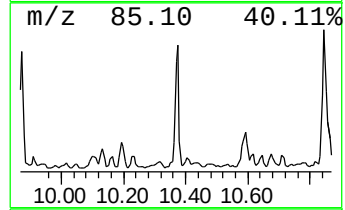
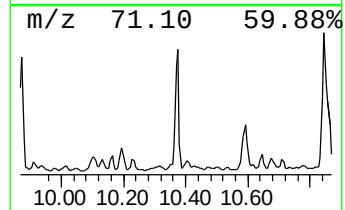
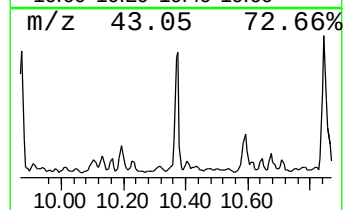
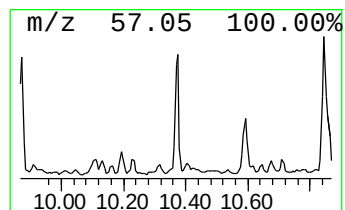
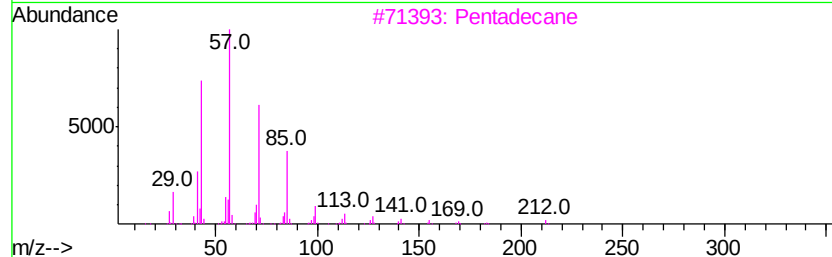
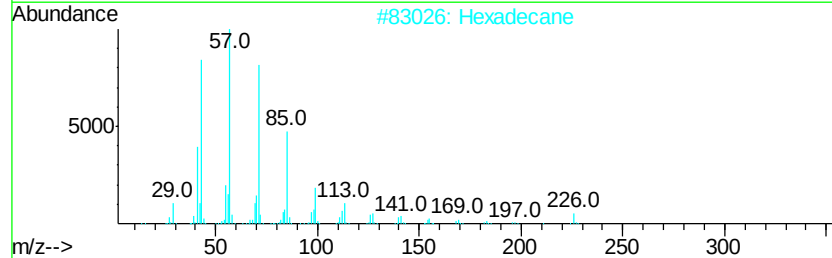
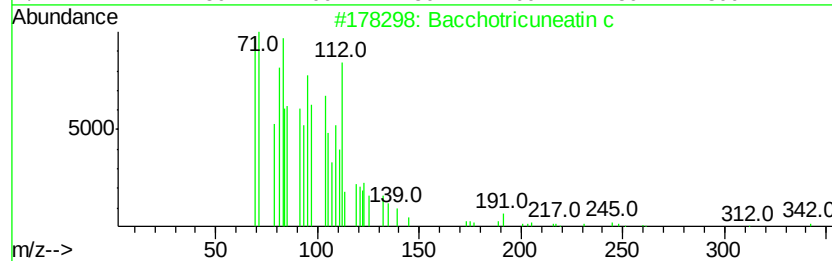
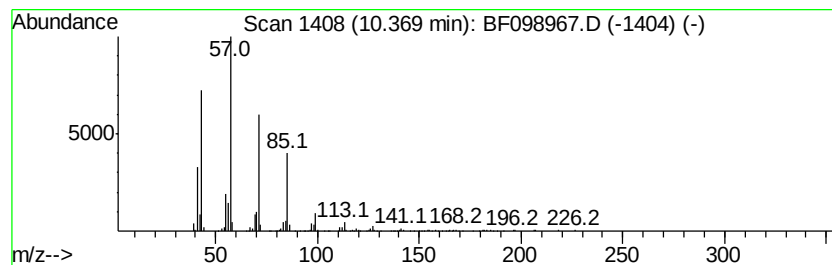
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Bacchotricuneatin c Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	10.10 ng	549523	Acenaphthene-d10	9.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Bacchotricuneatin c	342 C20H22O5	066563-30-2 96
2		Hexadecane	226 C16H34	000544-76-3 94
3		Pentadecane	212 C15H32	000629-62-9 90
4		Pentadecane, 7-methyl-	226 C16H34	006165-40-8 87
5		Tetradecane	198 C14H30	000629-59-4 86



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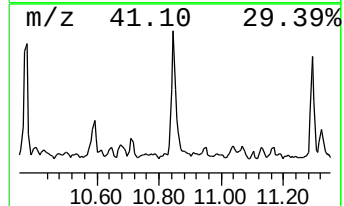
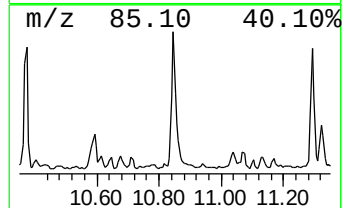
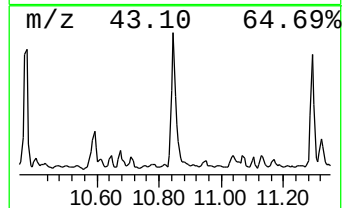
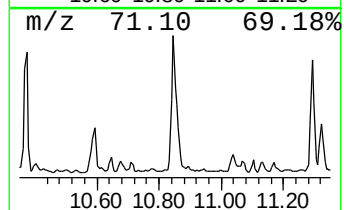
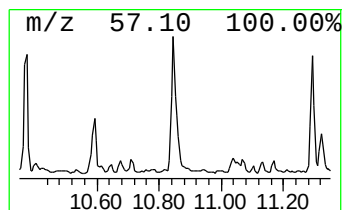
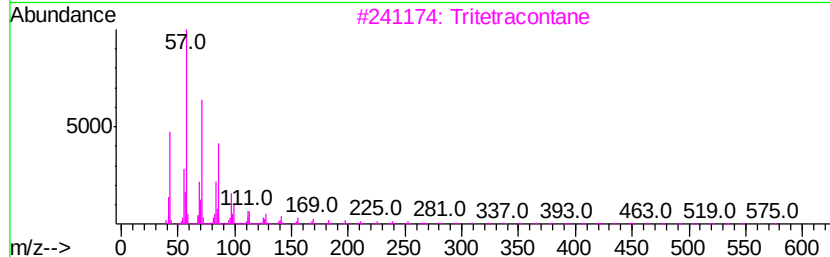
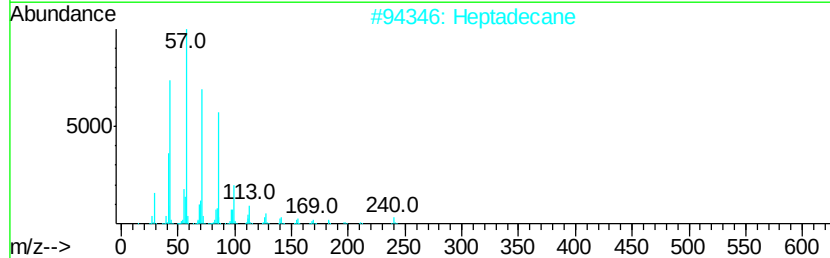
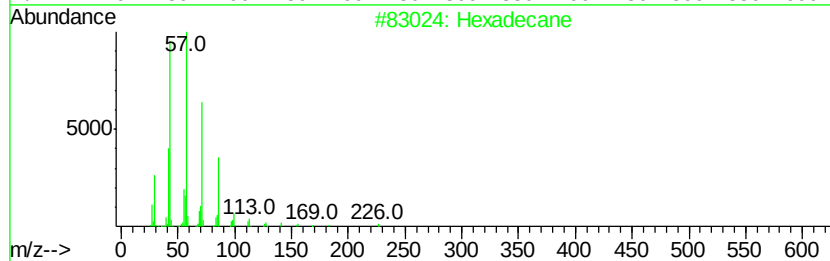
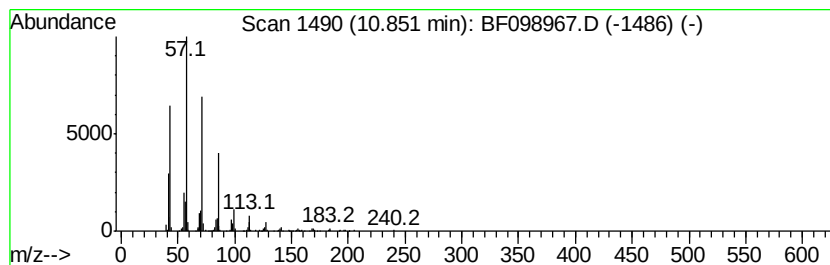
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.85	16.68 ng	733262	Phenanthrene-d10	11.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	90
2	Heptadecane		240	C17H36	000629-78-7	90
3	Tritetracontane		605	C43H88	007098-21-7	90
4	Heneicosane		296	C21H44	000629-94-7	90
5	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
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Operator : SJ/JU
Sample : I5511-01 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

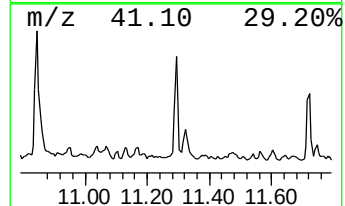
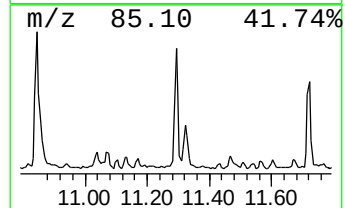
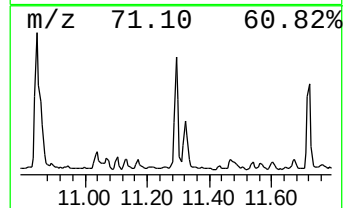
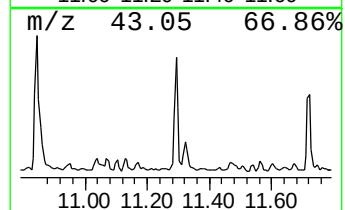
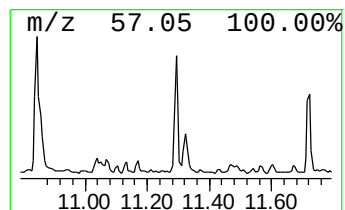
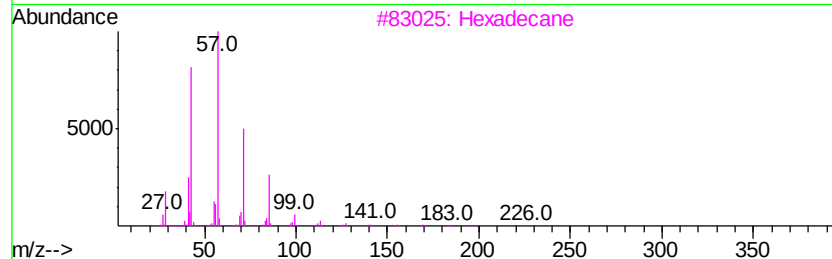
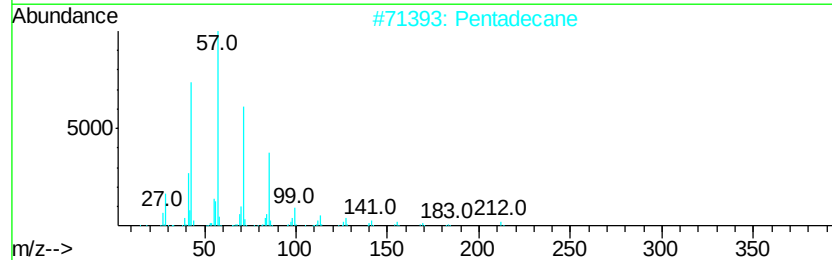
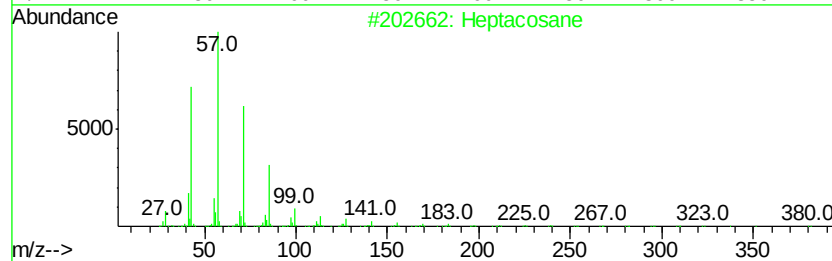
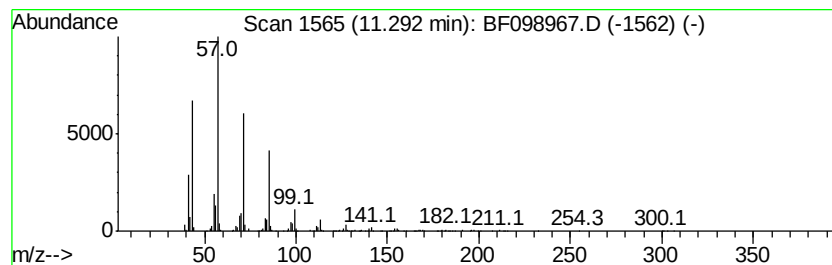
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.29	10.97 ng	482339	Phenanthrene-d10		11.44	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	91
2	Pentadecane		212	C15H32	000629-62-9	90
3	Hexadecane		226	C16H34	000544-76-3	90
4	Heneicosane		296	C21H44	000629-94-7	90
5	Heptadecane		240	C17H36	000629-78-7	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
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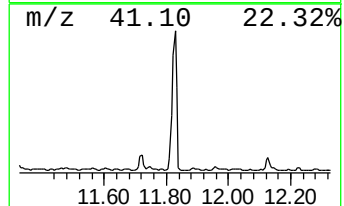
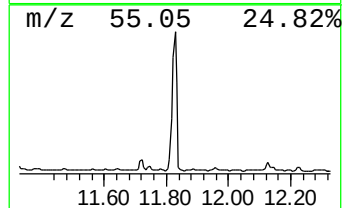
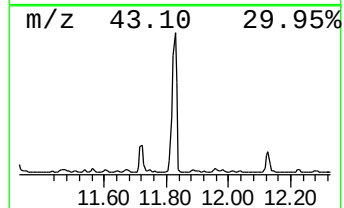
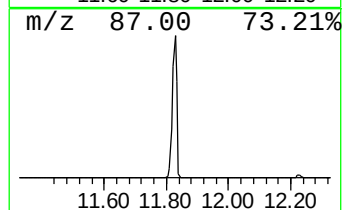
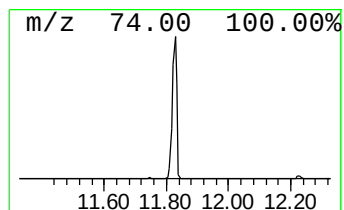
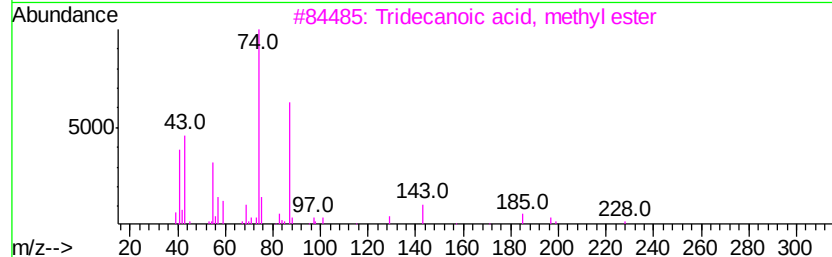
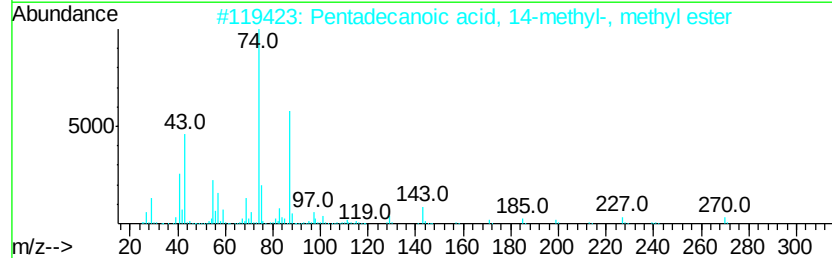
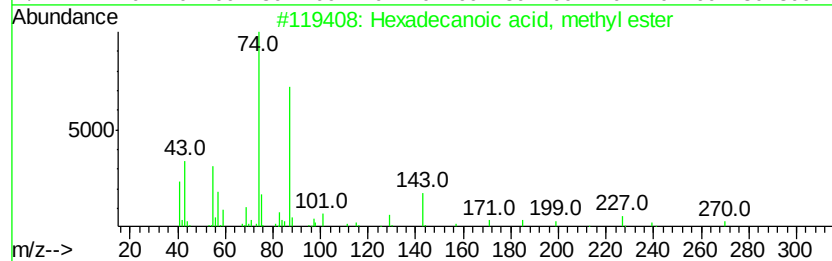
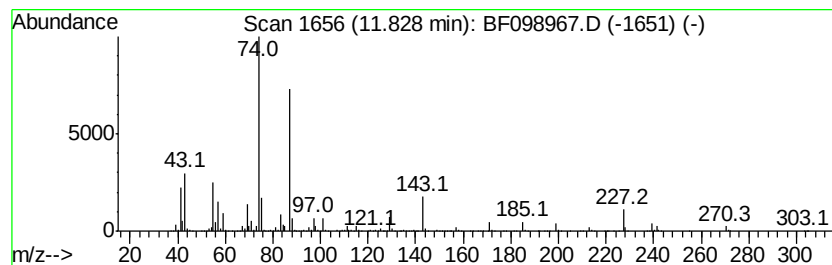
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	117.98 ng	5187180	Phenanthrene-d10	11.44

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	92
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4		Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
5		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
Data File : BF098967.D
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ALS Vial : 11 Sample Multiplier: 1

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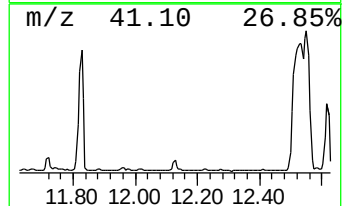
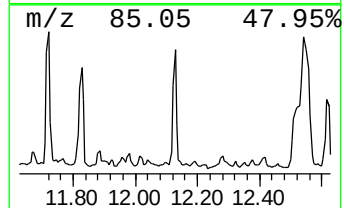
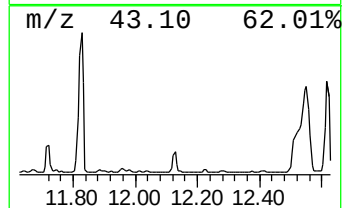
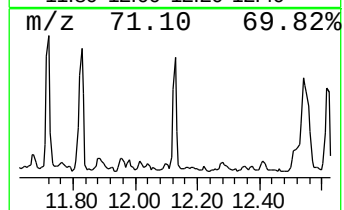
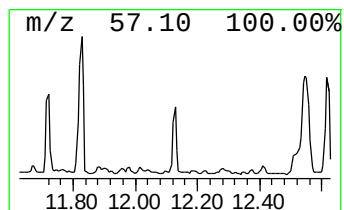
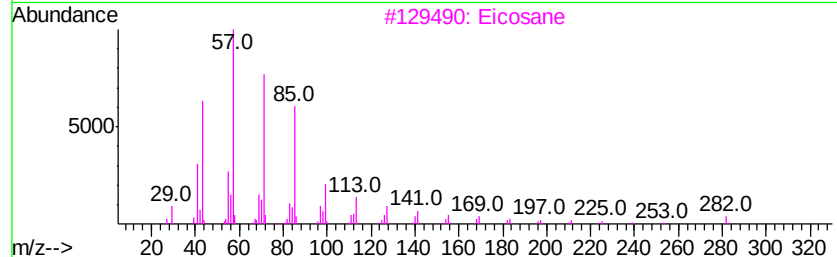
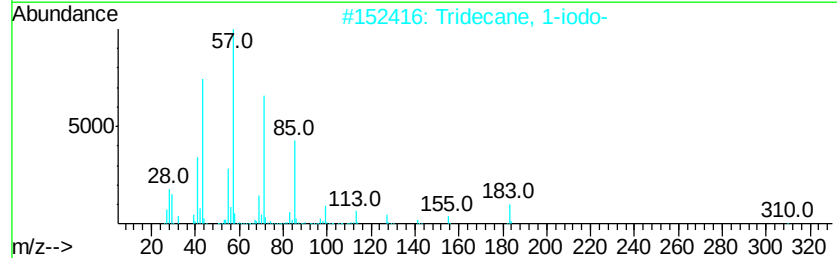
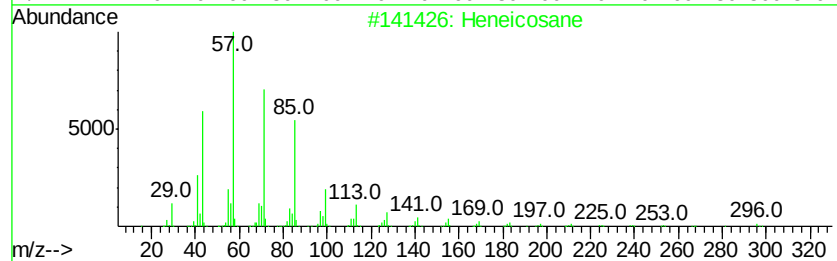
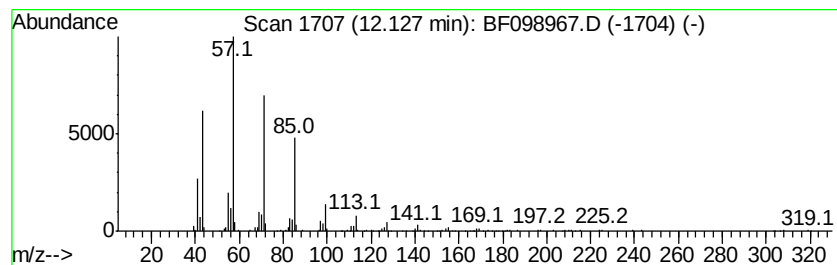
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	8.38 ng	368660	Phenanthrene-d10	11.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
3		Eicosane	282	C20H42	000112-95-8	90
4		Nonadecane	268	C19H40	000629-92-5	90
5		Hexadecane	226	C16H34	000544-76-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
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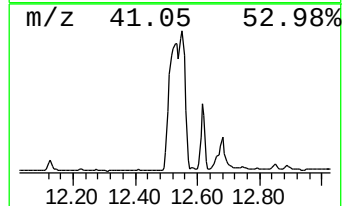
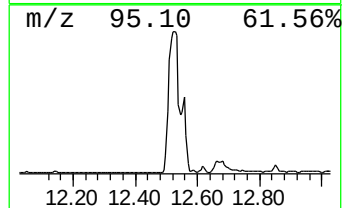
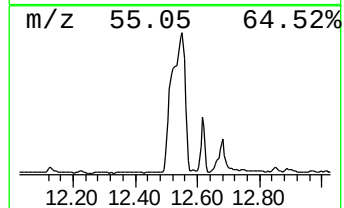
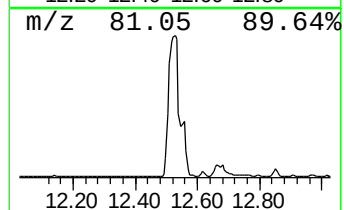
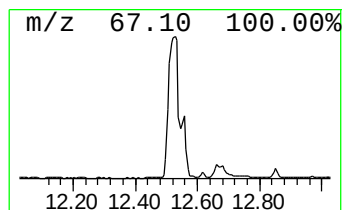
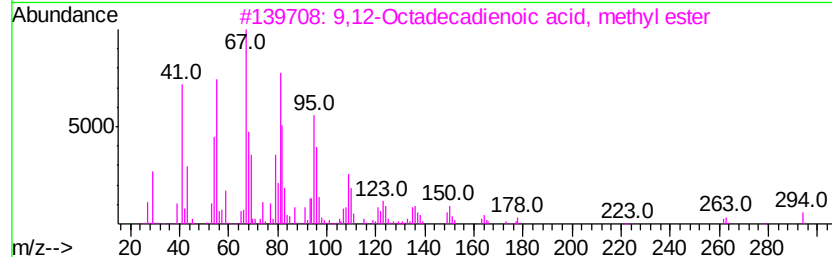
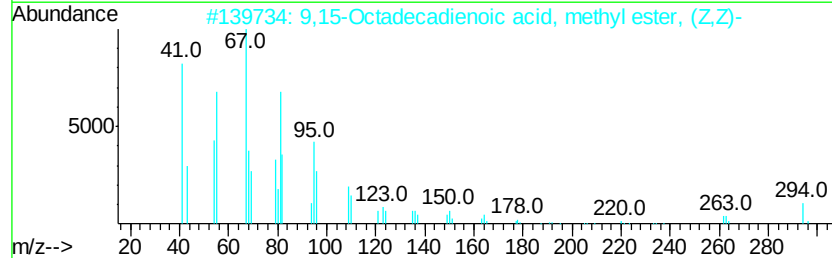
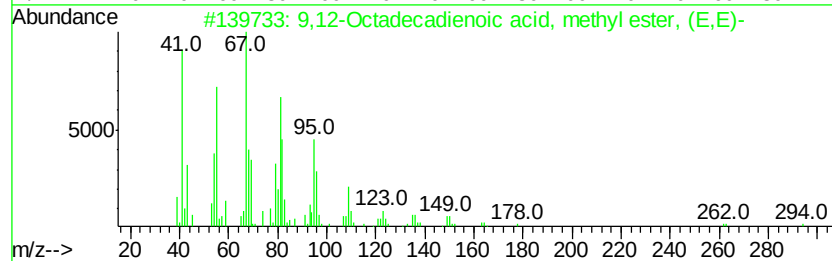
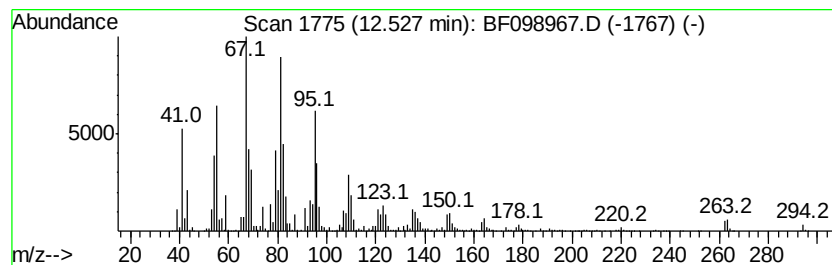
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.53	267.44 ng	11758300	Phenanthrene-d10	11.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
3		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
4		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
Data File : BF098967.D
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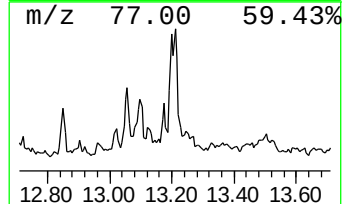
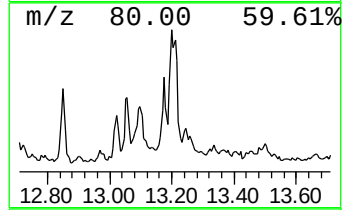
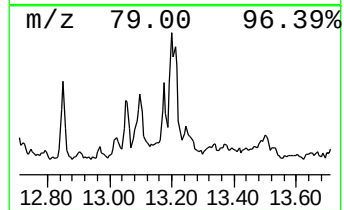
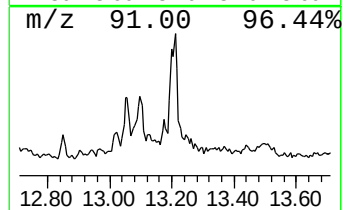
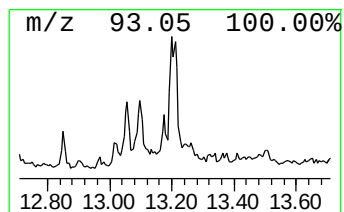
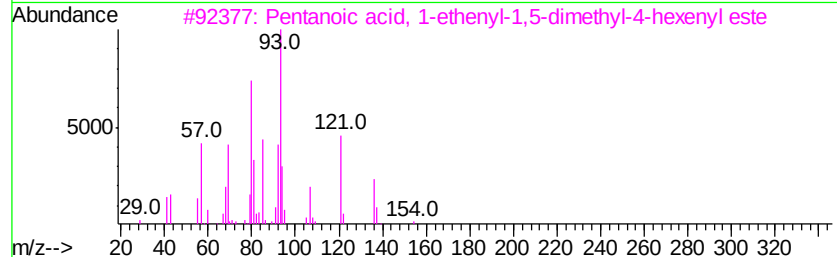
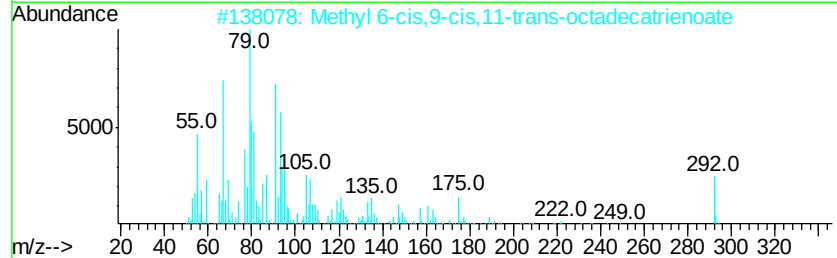
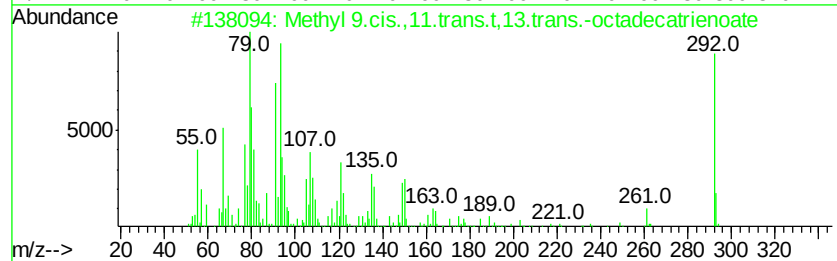
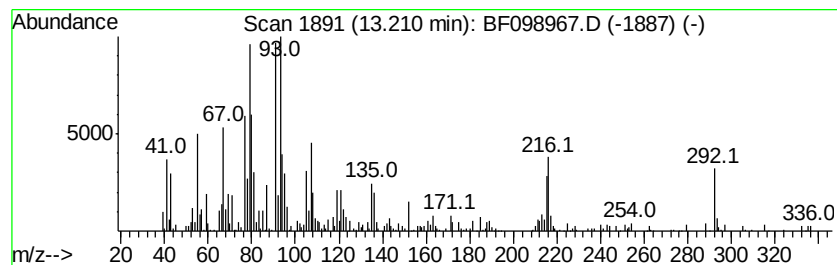
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Methyl 9.cis.,11.trans.t,13... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	7.21 ng	304933	Chrysene-d12	14.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methyl 9.cis.,11.trans.t,13.tran...	292	C19H32O2	1000336-42-6	89
2		Methyl 6-cis,9-cis,11-trans-octa...	292	C19H32O2	1000336-37-7	83
3		Pentanoic acid, 1-ethenyl-1,5-di...	238	C15H26O2	010471-96-2	43
4		Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	016022-04-1	43
5		Camphene	136	C10H16	000079-92-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
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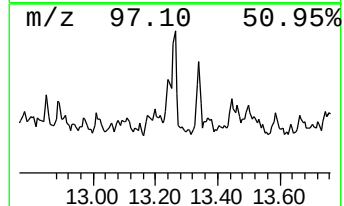
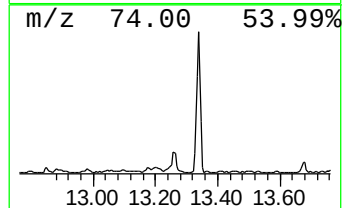
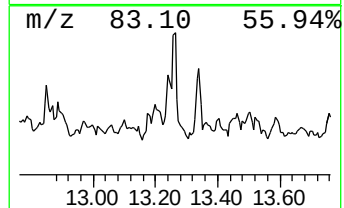
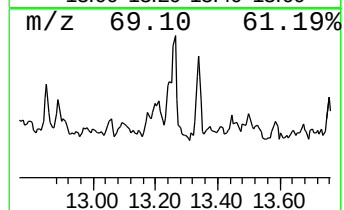
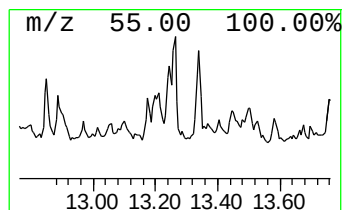
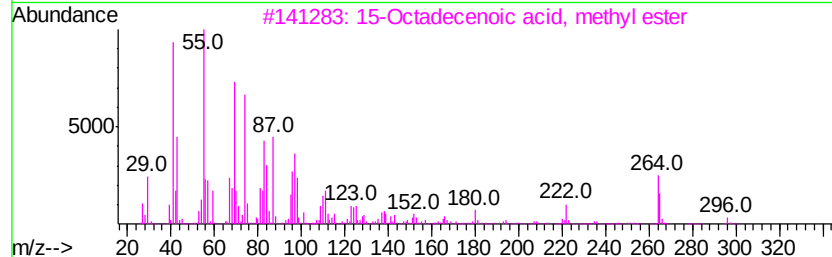
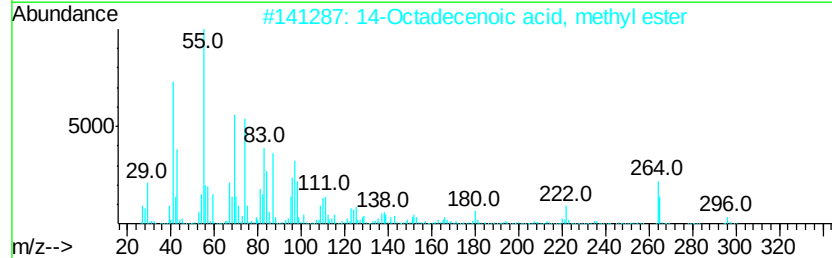
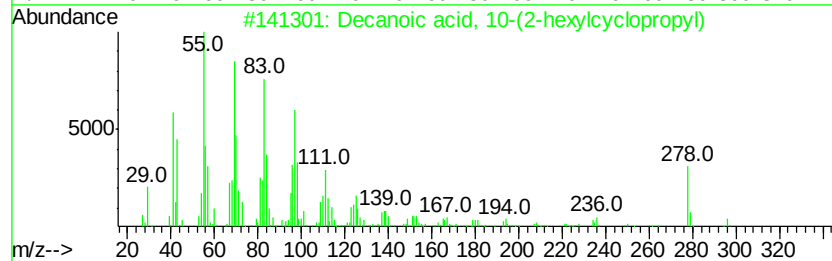
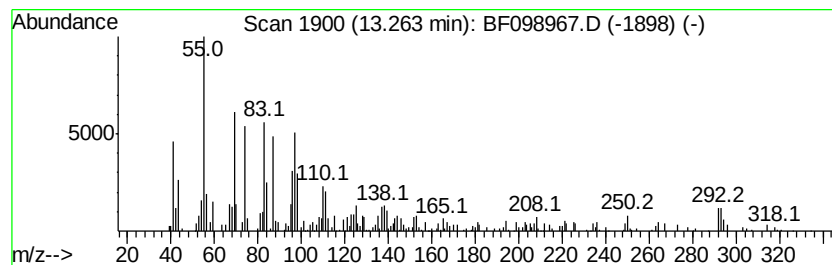
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Decanoic acid, 10-(2-hexylc... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.26	6.31 ng	266752	Chrysene-d12	14.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decanoic acid, 10-(2-hexylcyclop...	296	C19H36O2	1000197-99-5	83
2	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	83
3	15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	83
4	14-Pentadecenoic acid	240	C15H28O2	017351-34-7	78
5	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	60



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
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ALS Vial : 11 Sample Multiplier: 1

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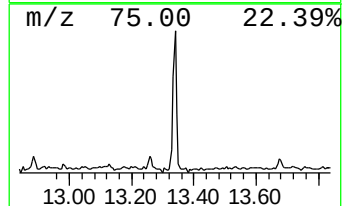
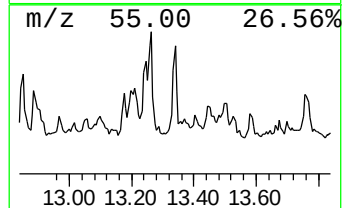
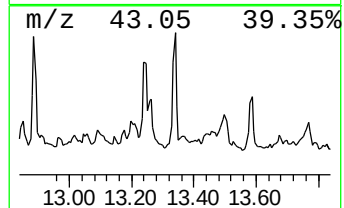
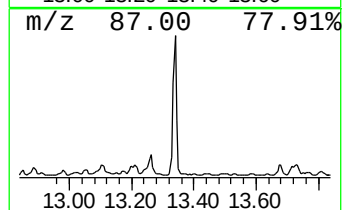
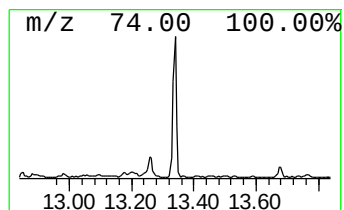
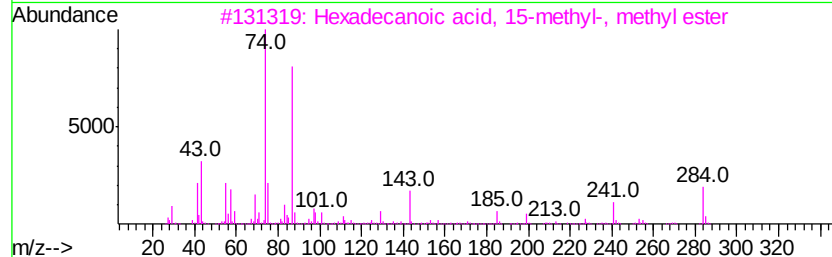
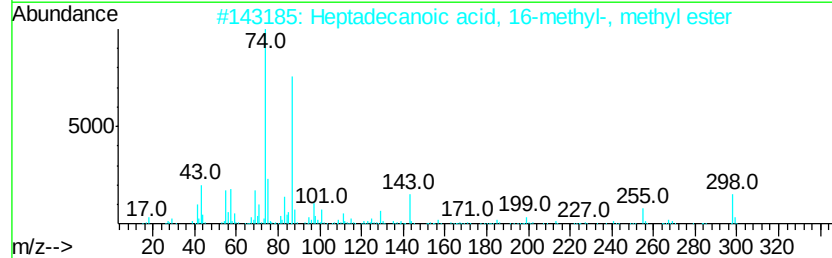
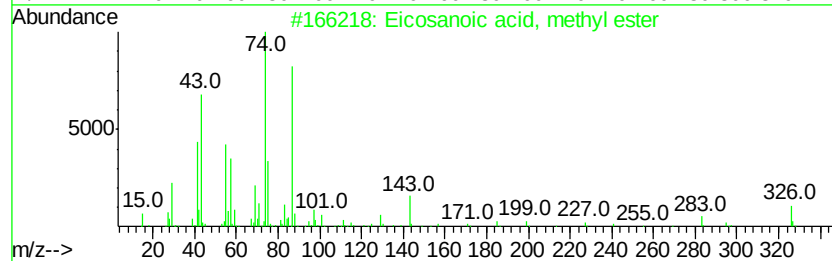
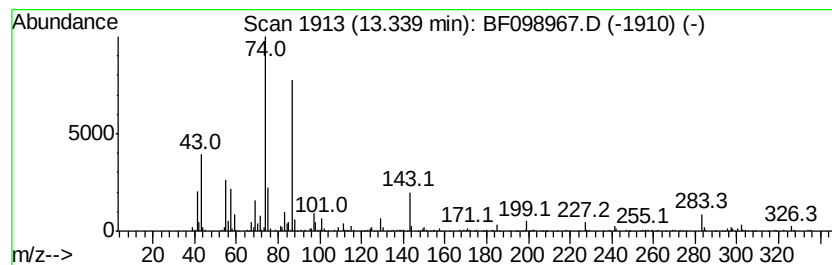
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Eicosanoic acid, methyl ester Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.34	6.38 ng	269681	Chrysene-d12	14.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	96
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	96
3			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	94
4			Methyl stearate	298	C19H38O2	000112-61-8	91
5			Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
Data File : BF098967.D
Acq On : 30 Sep 2017 17:19
Operator : SJ/JU
Sample : I5511-01 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-092817

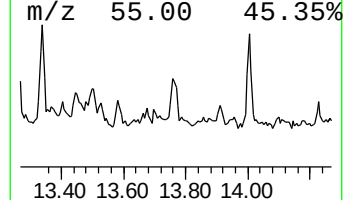
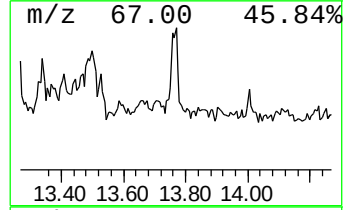
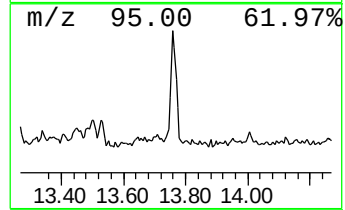
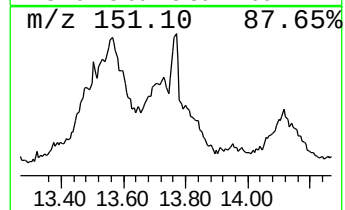
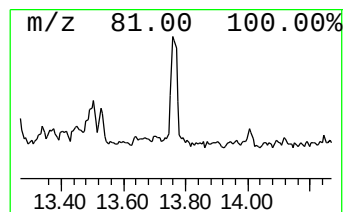
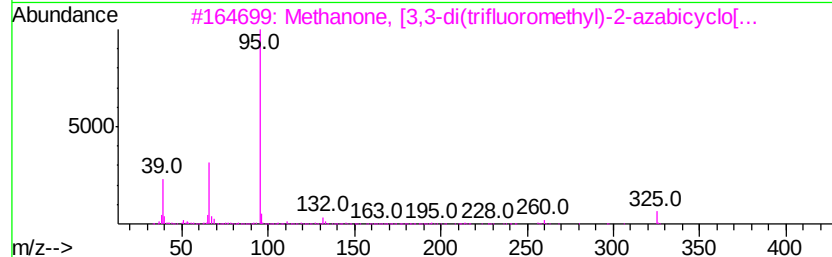
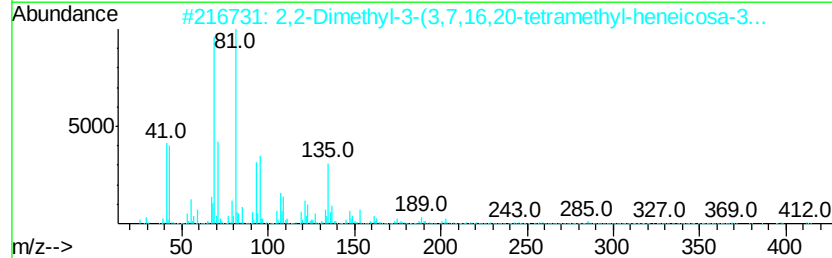
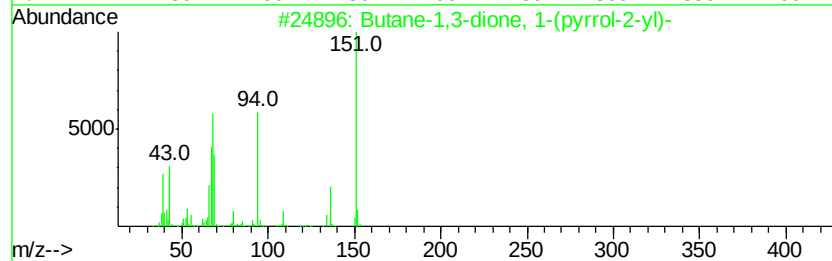
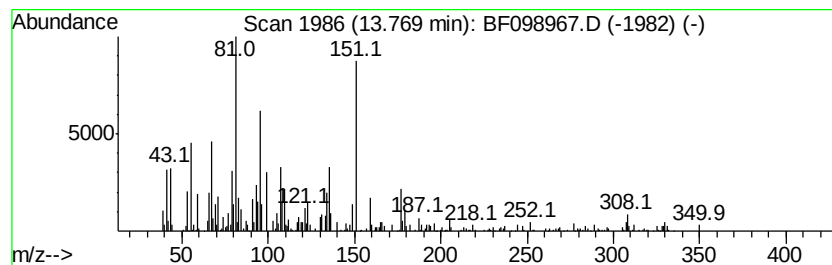
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown13.77 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	6.91 ng	292163	Chrysene-d12	14.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane-1,3-dione, 1-(pyrrol-2-yl)-	151	C8H9NO2	022441-25-4	27
2			2,2-Dimethyl-3-(3,7,16,20-tetram...	412	C29H48O	1000194-78-6	22
3			Methanone, [3,3-di(trifluorometh...	325	C13H9F6NO2	1000276-52-5	18
4			Bis-bicyclo [2.2.1] hept-2-ylketone	218	C15H22O	1000222-21-8	16
5			4,7-Methano-1H-indene, octahydro-	136	C10H16	006004-38-2	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
Data File : BF098967.D
Acq On : 30 Sep 2017 17:19
Operator : SJ/JU
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Instrument :
BNA_F
ClientSampleId :
EO-01-092817

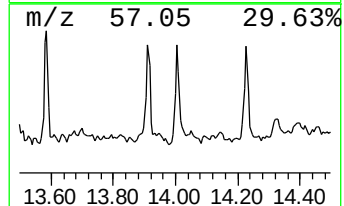
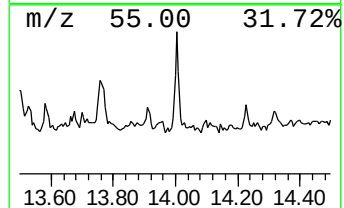
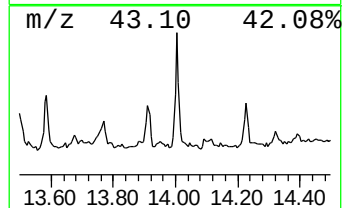
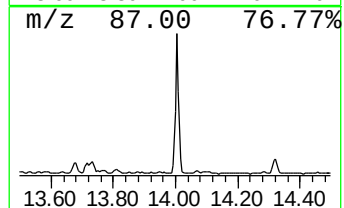
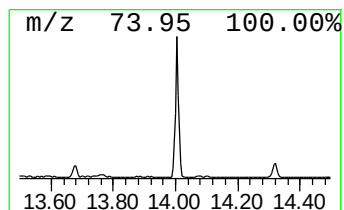
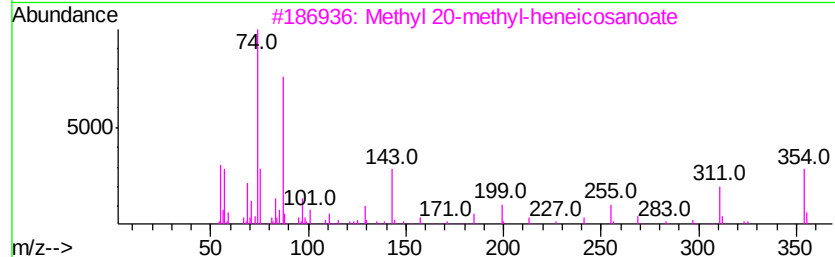
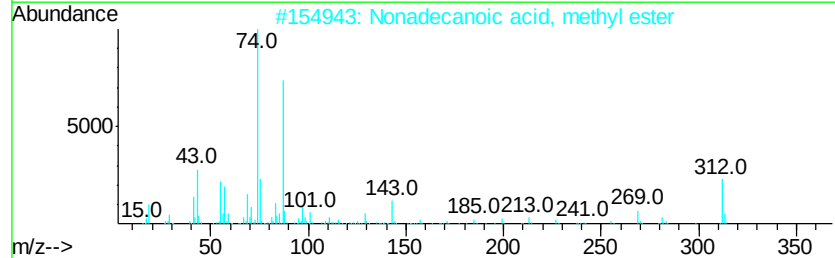
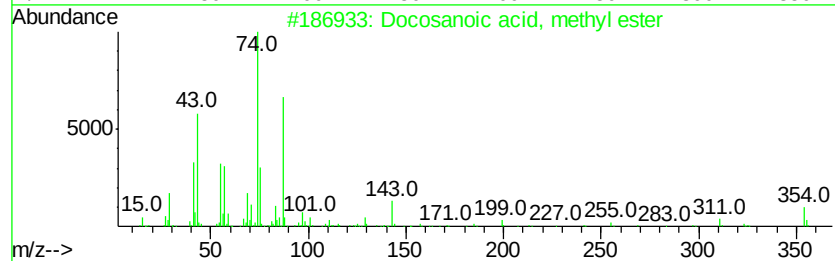
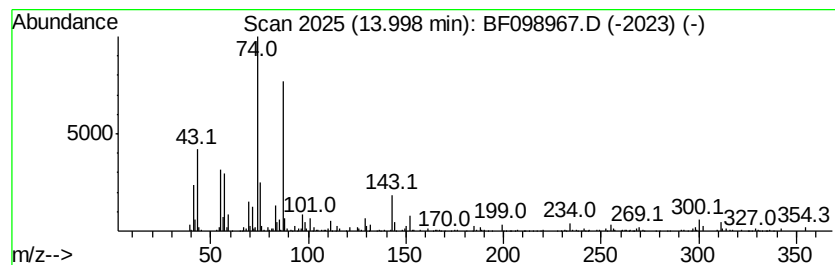
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Docosanoic acid, methyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	7.25 ng	306345	Chrysene-d12	14.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	96
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	95
3		Methyl 20-methyl-heneicosanoate	354	C23H46O2	1000336-47-4	89
4		Octadecanoic acid, 10-methyl-, m...	312	C20H40O2	002490-19-9	87
5		Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
Data File : BF098967.D
Acq On : 30 Sep 2017 17:19
Operator : SJ/JU
Sample : I5511-01 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EO-01-092817

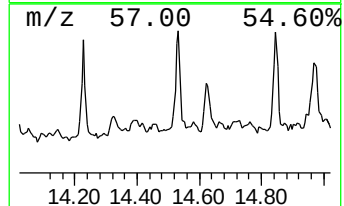
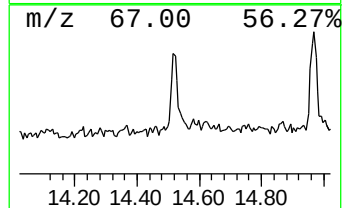
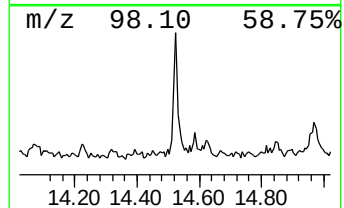
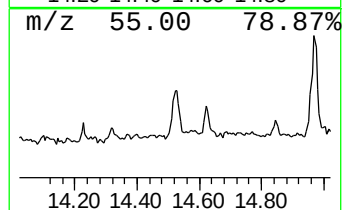
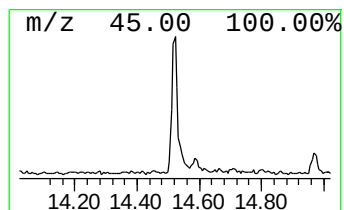
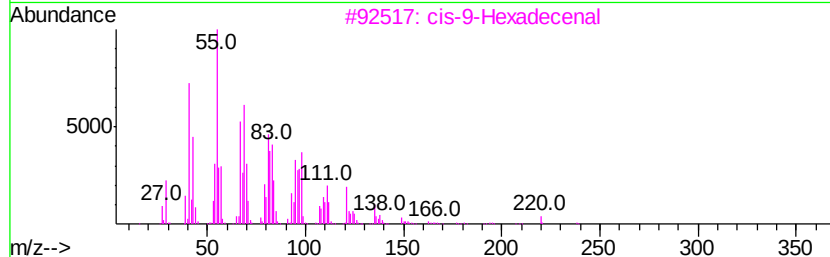
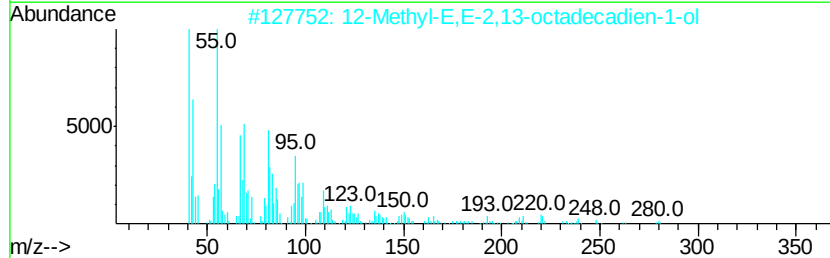
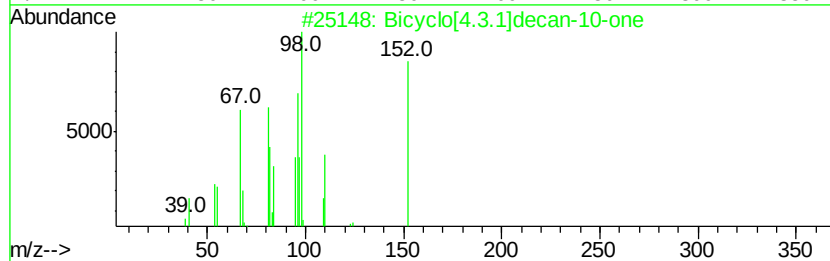
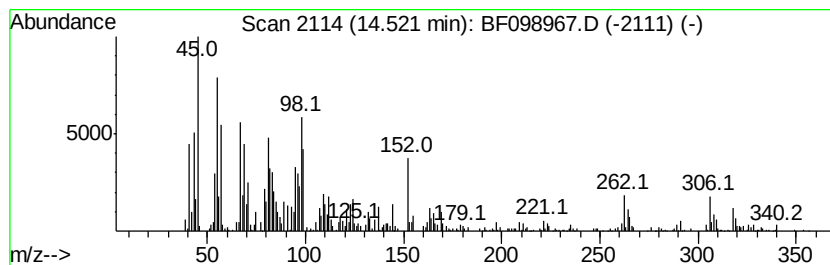
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown14.52 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	11.09 ng	468867	Chrysene-d12	14.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.3.1]decan-10-one	152	C10H16O	020440-21-5	49
2		12-Methyl-E,E-2,13-octadecadien-...	280	C19H36O	1000130-90-4	46
3		cis-9-Hexadecenal	238	C16H30O	056219-04-6	44
4		1-Decalone (cis-trans)	152	C10H16O	004832-16-0	44
5		2(1H)-Naphthalenone, octahydro-,...	152	C10H16O	016021-08-2	44



Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
Data File : BF098967.D
Acq On : 30 Sep 2017 17:19
Operator : SJ/JU
Sample : I5511-01 10X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-01-092817

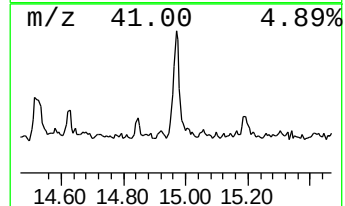
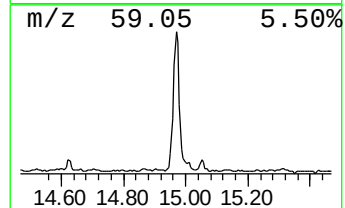
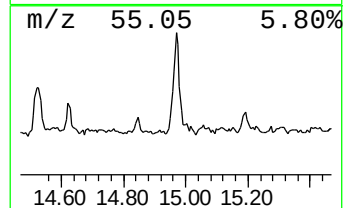
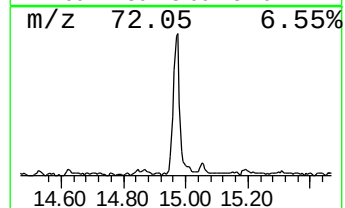
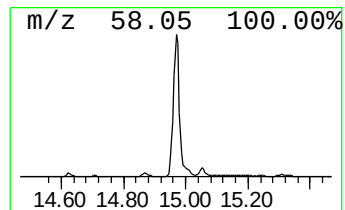
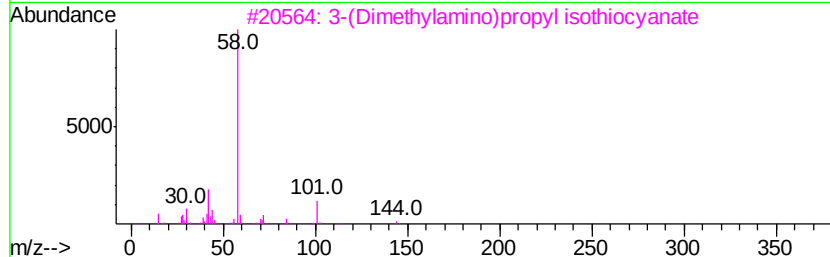
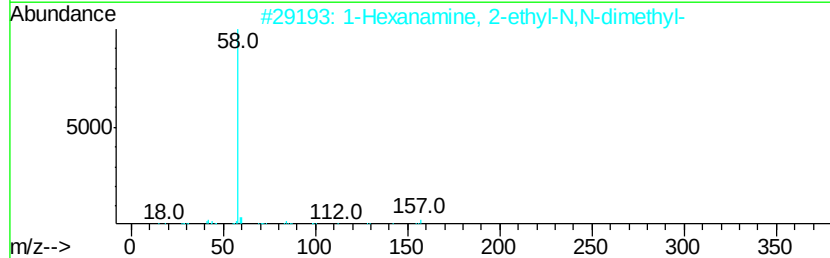
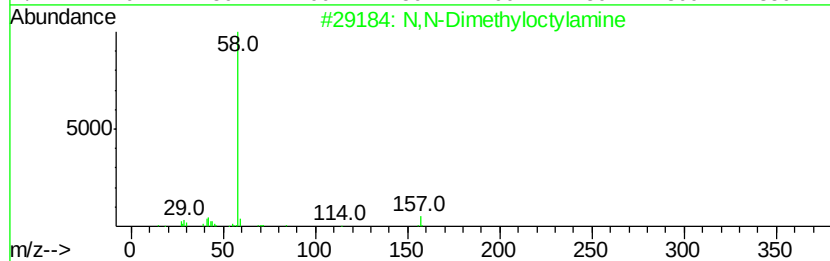
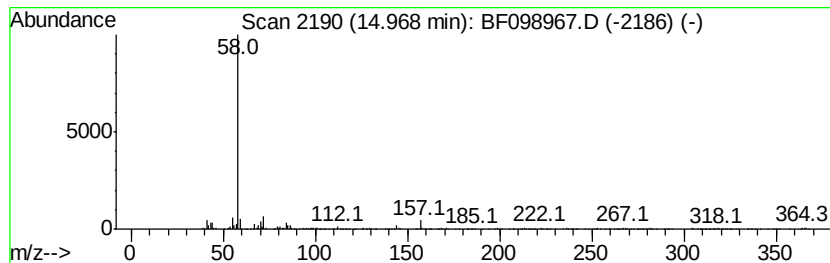
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 N,N-Dimethyloctylamine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	45.73 ng	1538210	Perylene-d12	15.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N,N-Dimethyloctylamine	157	C10H23N	007378-99-6	72
2		1-Hexanamine, 2-ethyl-N,N-dimethyl-	157	C10H23N	028056-87-3	72
3		3-(Dimethylamino)propyl isothiocy...	144	C6H12N2S	027421-70-1	72
4		N,N-Dimethyl-2-cyclohexyloxvethy...	171	C10H21NO	071126-60-8	64
5		1-Dodecanamine, N,N-dimethyl-	213	C14H31N	000112-18-5	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF093017\
 Data File : BF098967.D
 Acq On : 30 Sep 2017 17:19
 Operator : SJ/JU
 Sample : I5511-01 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-092817

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.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
unknown6.67	6.67	7.7	ng	328789	1	6.92	860010	20.0
Tridecane	8.78	6.8	ng	397423	2	8.20	1164570	20.0
Tetradecane	9.35	8.5	ng	461353	3	9.95	1087980	20.0
Bacchotricuneatin c	10.37	10.1	ng	549523	3	9.95	1087980	20.0
Hexadecane	10.85	16.7	ng	733262	4	11.44	879336	20.0
Heptacosane	11.29	11.0	ng	482339	4	11.44	879336	20.0
Hexadecanoic acid...	11.83	118.0	ng	5187180	4	11.44	879336	20.0
Heneicosane	12.13	8.4	ng	368660	4	11.44	879336	20.0
9,12-Octadecadien...	12.53	267.4	ng	11758300	4	11.44	879336	20.0
Methyl 9.cis.,11....	13.21	7.2	ng	304933	5	14.08	845335	20.0
Decanoic acid, 10...	13.26	6.3	ng	266752	5	14.08	845335	20.0
Eicosanoic acid, ...	13.34	6.4	ng	269681	5	14.08	845335	20.0
unknown13.77	13.77	6.9	ng	292163	5	14.08	845335	20.0
Docosanoic acid, ...	14.00	7.3	ng	306345	5	14.08	845335	20.0
unknown14.52	14.52	11.1	ng	468867	5	14.08	845335	20.0
N,N-Dimethyloctyl...	14.97	45.7	ng	1538210	6	15.57	672801	20.0