

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
 Data File : BF096264.D
 Acq On : 28 Jun 2017 22:04
 Operator : SJ/MA
 Sample : I3911-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-062317

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-BF062717.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	4.981	487	492	501	rVB	712694	883622	22.08%	2.348%
2	5.399	558	563	566	rBV	4049844	4001344	100.00%	10.632%
3	6.416	731	736	739	rBV	3867139	3915717	97.86%	10.404%
4	6.551	755	759	763	rBV	3779115	3745737	93.61%	9.953%
5	6.787	795	799	802	rBV	1037558	944669	23.61%	2.510%
6	6.940	821	825	829	rVB	3042834	2812172	70.28%	7.472%
7	7.187	862	867	870	rBV4	41541	41923	1.05%	0.111%
8	7.340	887	893	896	rBV	2516810	2350431	58.74%	6.245%
9	7.392	896	902	905	rVV2	77460	85815	2.14%	0.228%
10	7.434	905	909	913	rVB4	93068	98238	2.46%	0.261%
11	7.763	961	965	967	rBV2	35463	42846	1.07%	0.114%
12	7.792	967	970	974	rVV4	46618	52673	1.32%	0.140%
13	7.828	974	976	980	rVV3	47268	62945	1.57%	0.167%
14	7.875	980	984	986	rVV2	44109	49238	1.23%	0.131%
15	8.063	1012	1016	1020	rBV	1528970	1300444	32.50%	3.455%
16	8.140	1027	1029	1031	rVB	101935	73066	1.83%	0.194%
17	8.239	1042	1046	1051	rBV5	27967	53154	1.33%	0.141%
18	8.281	1051	1053	1056	rVB	76033	64350	1.61%	0.171%
19	8.363	1063	1067	1070	rBV4	63997	98860	2.47%	0.263%
20	8.457	1079	1083	1086	rVB3	56825	66505	1.66%	0.177%
21	8.498	1086	1090	1093	rBV	186868	200696	5.02%	0.533%
22	8.628	1107	1112	1115	rBV3	61342	103237	2.58%	0.274%
23	8.669	1116	1119	1122	rBV4	40573	56991	1.42%	0.151%
24	8.763	1132	1135	1139	rVB2	98245	134502	3.36%	0.357%
25	8.822	1139	1145	1147	rBV7	34596	64823	1.62%	0.172%
26	8.875	1150	1154	1157	rVB4	58122	87430	2.19%	0.232%
27	8.951	1164	1167	1169	rBV3	58184	57009	1.42%	0.151%
28	9.098	1189	1192	1195	rVV	290760	233403	5.83%	0.620%
29	9.139	1195	1199	1203	rVB	3780604	3463168	86.55%	9.202%
30	9.234	1211	1215	1219	rBV3	104858	158404	3.96%	0.421%
31	9.275	1220	1222	1227	rVV3	58593	93254	2.33%	0.248%
32	9.345	1231	1234	1238	rVB4	53024	65336	1.63%	0.174%
33	9.398	1240	1243	1246	rBV2	52502	75999	1.90%	0.202%
34	9.481	1253	1257	1259	rBV2	72482	95434	2.39%	0.254%

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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

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35	9.504	1259	1261	1263	rVV2	79077	88359	2.21%	0.235%
36	9.534	1263	1266	1268	rVV	699354	649600	16.23%	1.726%
37	9.551	1268	1269	1272	rVB	216094	159294	3.98%	0.423%
38	9.657	1286	1287	1292	rVB3	41691	43118	1.08%	0.115%
39	9.704	1292	1295	1298	rVB4	44839	50467	1.26%	0.134%
40	9.739	1298	1301	1303	rBV2	56870	61918	1.55%	0.165%
41	9.763	1303	1305	1309	rVB2	56301	49177	1.23%	0.131%
42	9.816	1309	1314	1318	rBV2	1357917	1234536	30.85%	3.280%
43	9.928	1331	1333	1337	rVB4	74013	99137	2.48%	0.263%
44	10.022	1341	1349	1352	rVV3	124120	231487	5.79%	0.615%
45	10.057	1352	1355	1357	rVV2	49842	53190	1.33%	0.141%
46	10.086	1357	1360	1363	rVV3	120884	111085	2.78%	0.295%
47	10.257	1387	1389	1392	rVB	87321	72730	1.82%	0.193%
48	10.316	1396	1399	1404	rVB6	58244	78328	1.96%	0.208%
49	10.386	1408	1411	1414	rVB5	46085	51591	1.29%	0.137%
50	10.481	1420	1427	1433	rBV	175215	229358	5.73%	0.609%
51	10.528	1433	1435	1439	rVB3	41775	48292	1.21%	0.128%
52	10.598	1443	1447	1451	rVB	1999611	1990950	49.76%	5.290%
53	10.733	1467	1470	1471	rBV	144152	141417	3.53%	0.376%
54	10.745	1471	1472	1479	rVB	160129	138219	3.45%	0.367%
55	11.180	1543	1546	1548	rBV2	74126	68753	1.72%	0.183%
56	11.210	1548	1551	1556	rVB	106155	106476	2.66%	0.283%
57	11.298	1562	1566	1568	rBV	1191091	1009925	25.24%	2.683%
58	11.604	1616	1618	1621	rBV	46437	41965	1.05%	0.112%
59	11.827	1653	1656	1661	rBV	279768	254243	6.35%	0.676%
60	12.345	1742	1744	1747	rBV2	64239	58901	1.47%	0.157%
61	12.504	1768	1771	1775	rBV	73000	86961	2.17%	0.231%
62	12.616	1787	1790	1793	rBV	85150	95850	2.40%	0.255%
63	12.880	1831	1835	1839	rBV	2822043	2823656	70.57%	7.503%
64	13.780	1985	1988	1996	rVB2	331617	340735	8.52%	0.905%
65	13.927	2009	2013	2016	rBV	987834	930736	23.26%	2.473%
66	15.363	2253	2257	2261	rVB	477490	601553	15.03%	1.598%

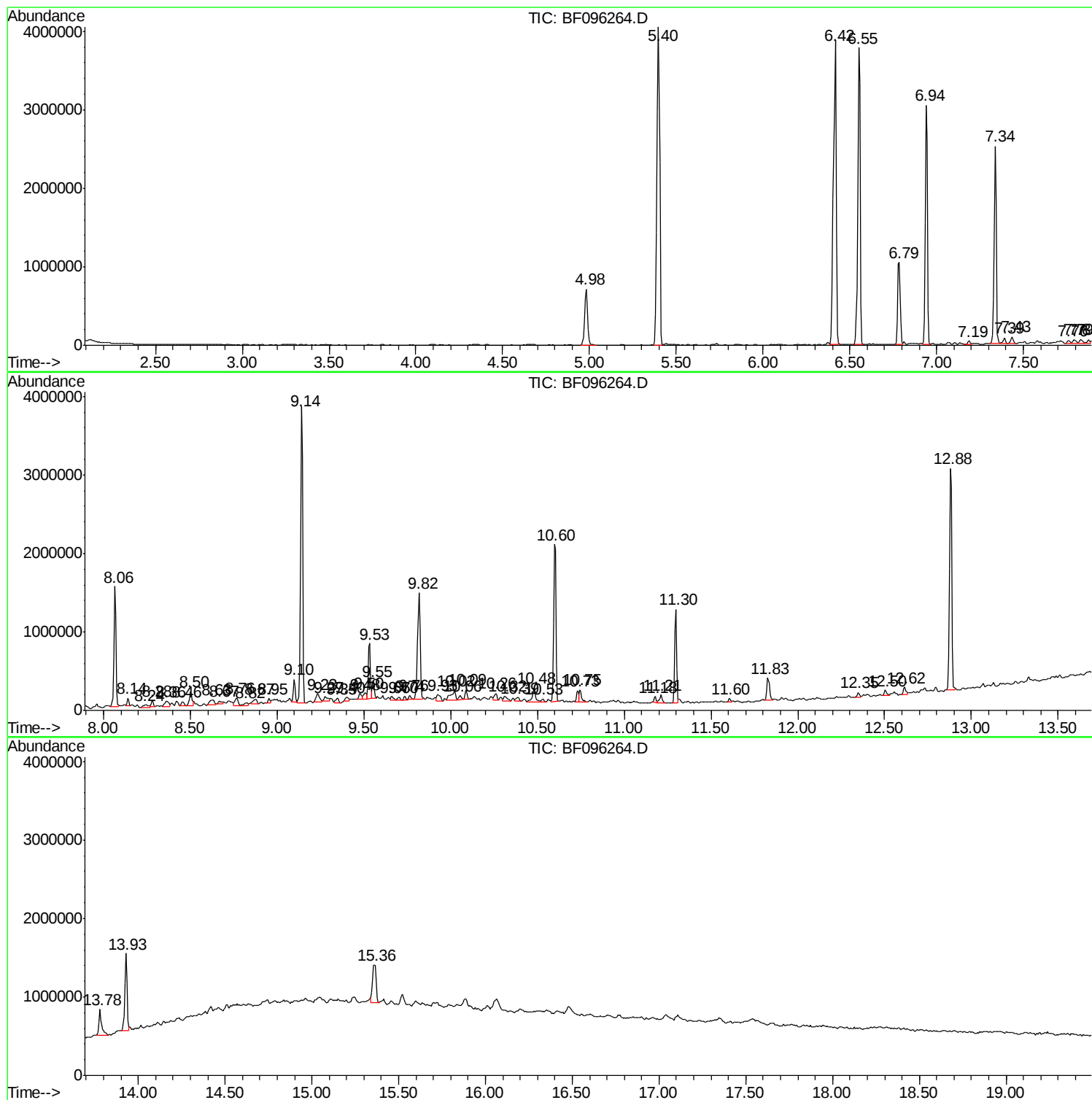
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
BNA_F
ClientSampled :
SU-04-062317

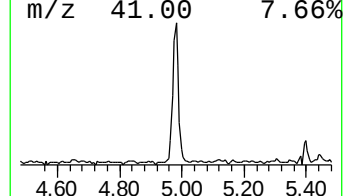
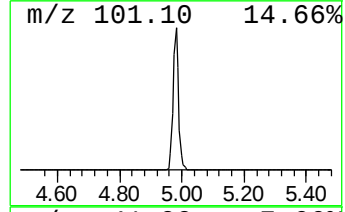
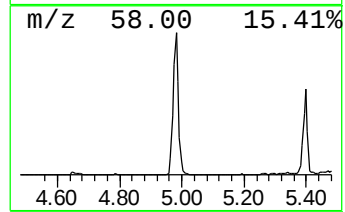
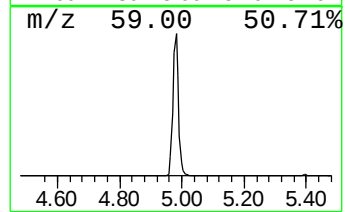
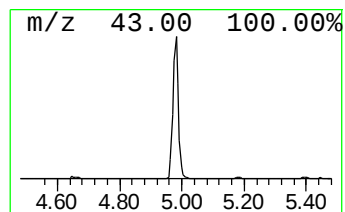
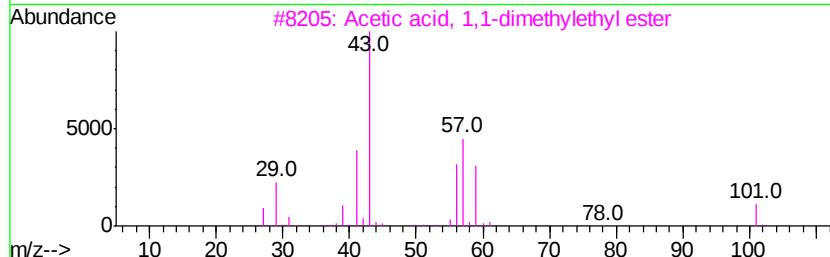
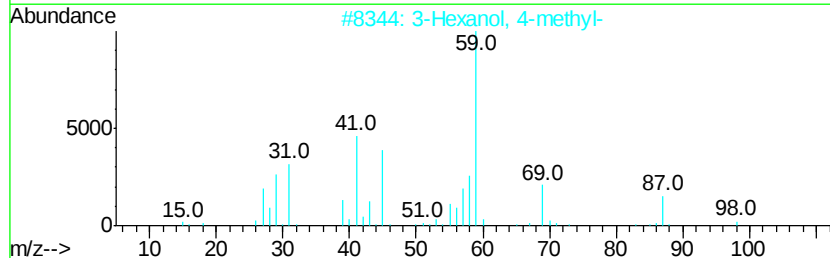
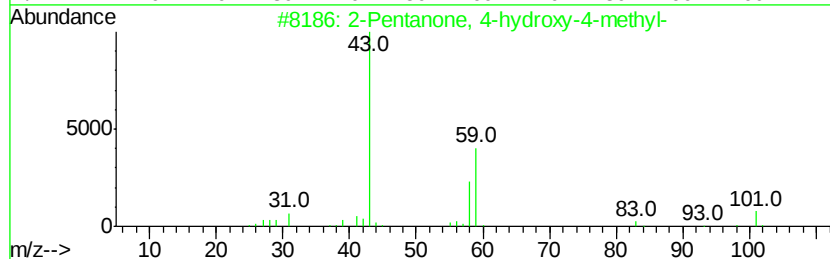
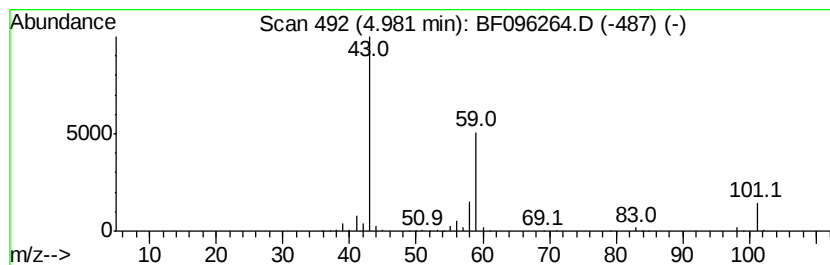
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	18.71 ng	883622	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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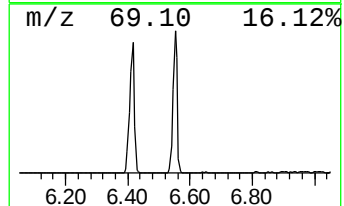
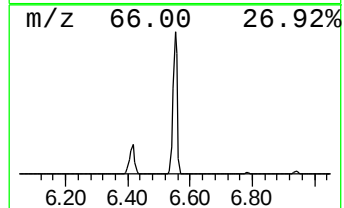
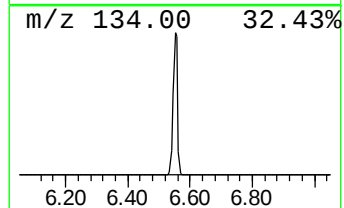
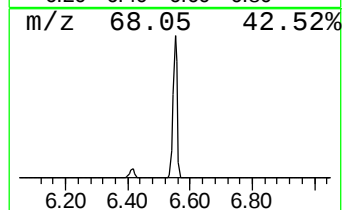
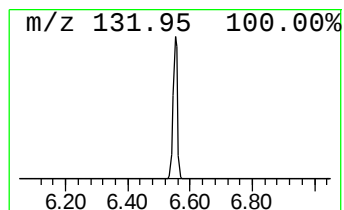
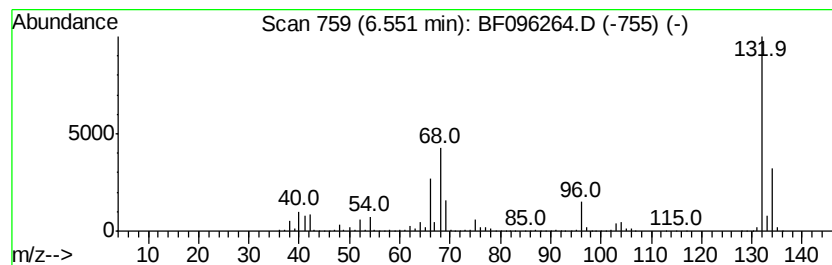
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	79.30 ng	3745740	1,4-Dichlorobenzene-d4	6.79

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	16
3	5-Aminoindole			132	C8H8N2	005192-03-0	12
4	Amphetaminil			250	C17H18N2	017590-01-1	10
5	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10



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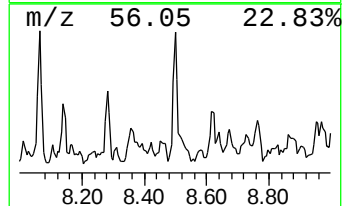
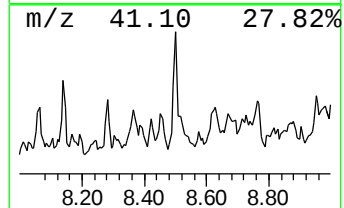
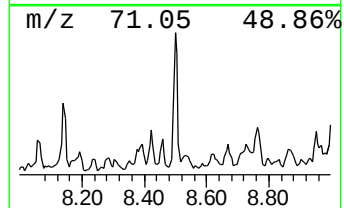
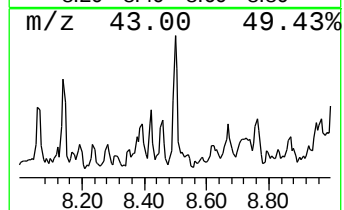
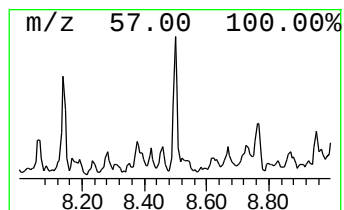
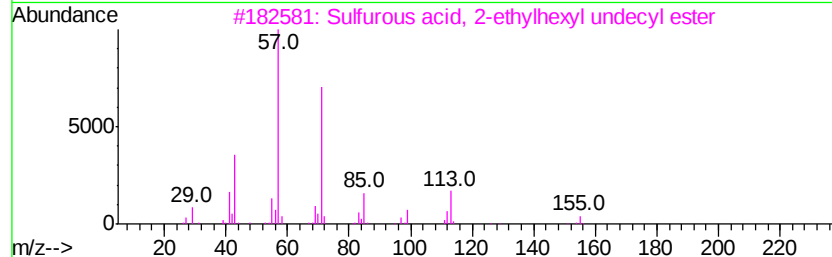
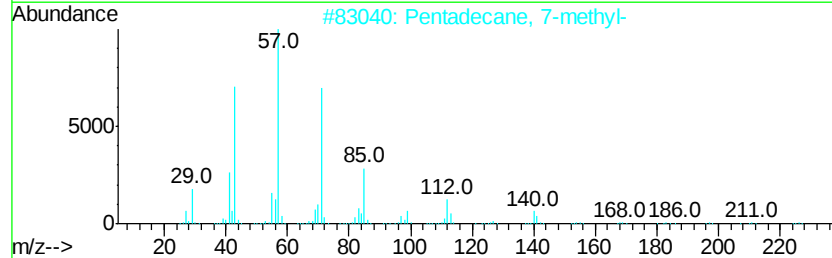
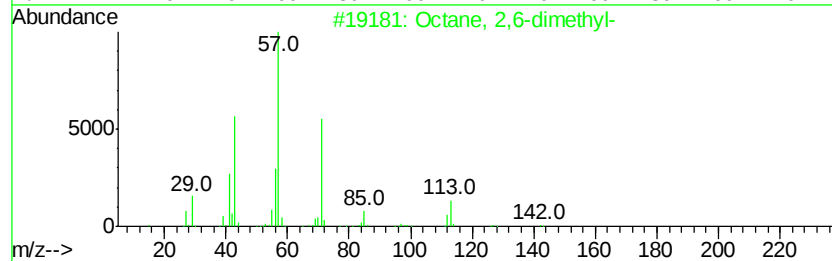
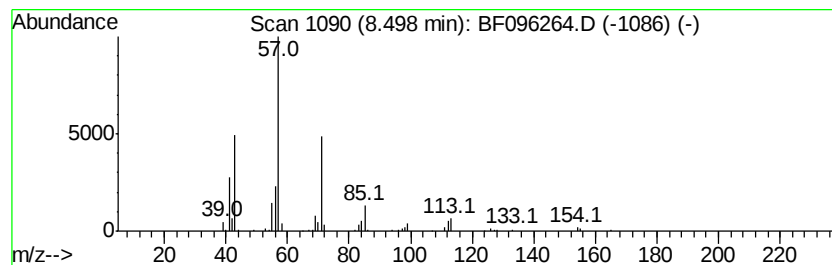
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Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.50	3.09 ng	200696	Naphthalene-d8	8.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
2		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	72
3		Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	64
4		Nonane, 3-methyl-	142	C10H22	005911-04-6	64
5		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	59



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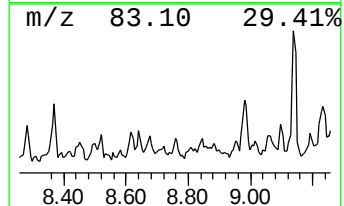
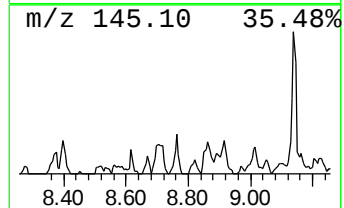
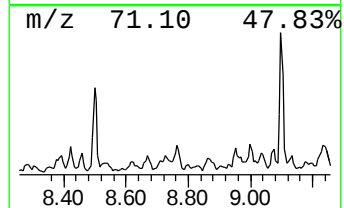
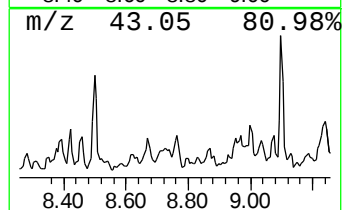
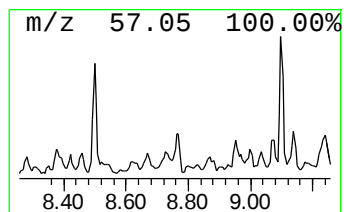
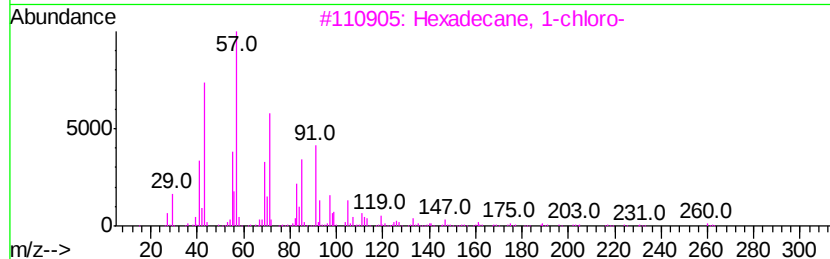
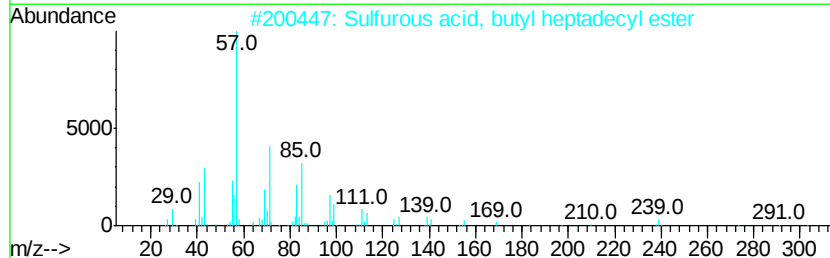
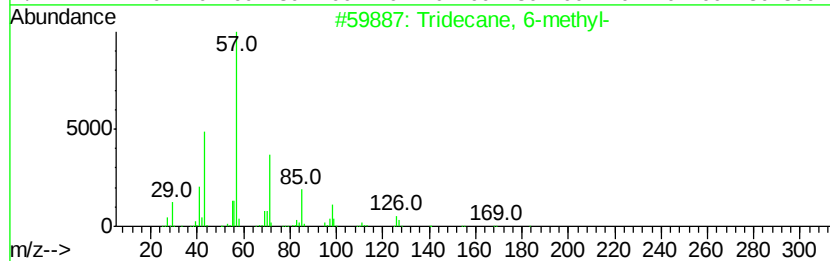
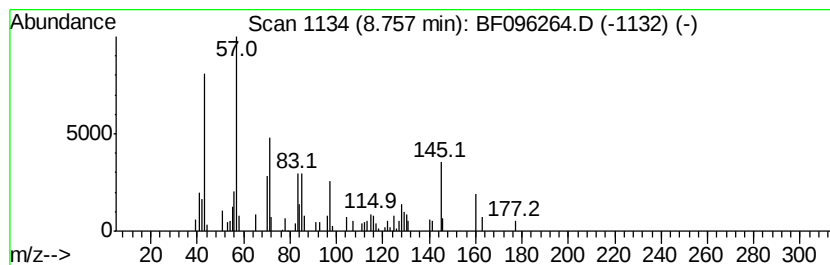
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown8.76 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.76	2.07 ng	134502	Napthalene-d8	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 6-methyl-	198	C14H30	013287-21-3	47
2		Sulfurous acid, butyl heptadecyl...	376	C21H44O3S	1000309-18-4	47
3		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	43
4		Octane, 4-ethyl-	142	C10H22	015869-86-0	43
5		Tetratetracontane	619	C44H90	007098-22-8	43



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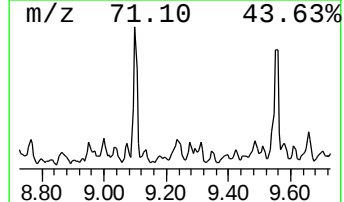
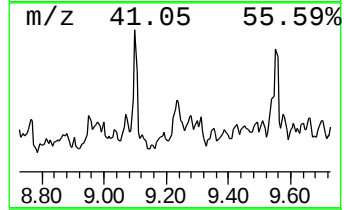
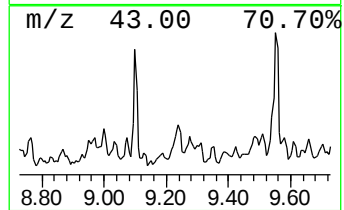
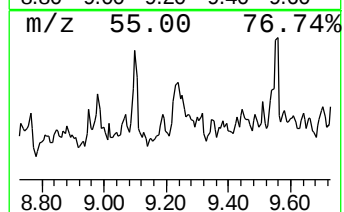
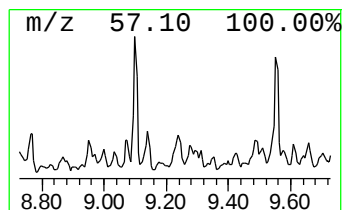
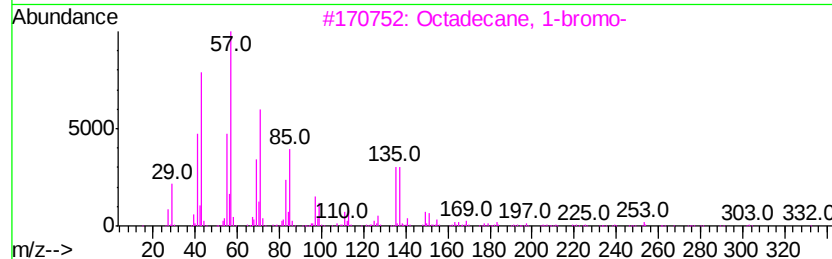
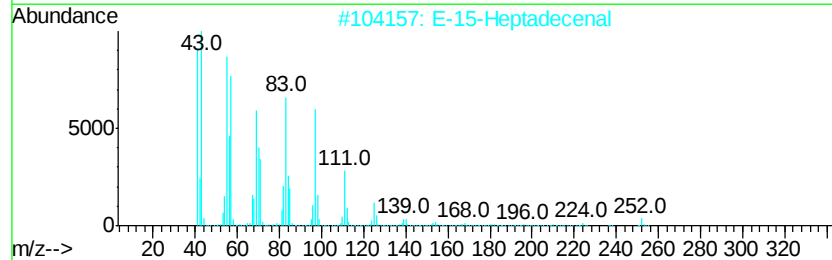
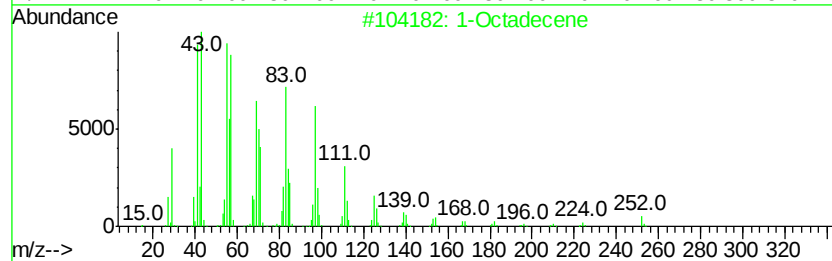
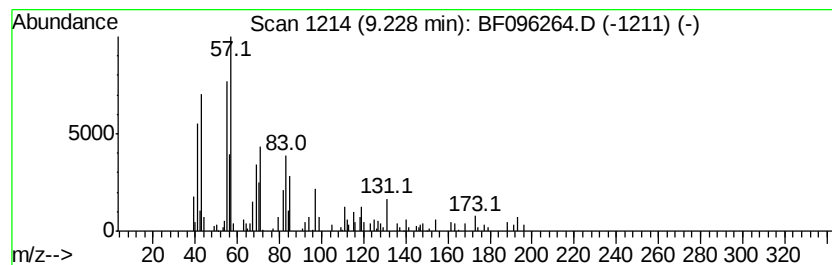
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ClientSampleId :
SU-04-062317

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown9.23 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.23	2.57 ng	158404	Acenaphthene-d10	9.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	49
2	E-15-Heptadecenal	252	C17H32O	1000130-97-9	46
3	Octadecane, 1-bromo-	332	C18H37Br	000112-89-0	43
4	Pentadecane	212	C15H32	000629-62-9	27
5	Tetradecane	198	C14H30	000629-59-4	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
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Acq On : 28 Jun 2017 22:04
Operator : SJ/MA
Sample : I3911-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

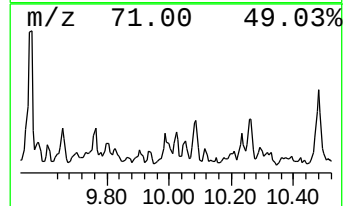
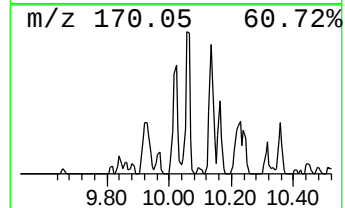
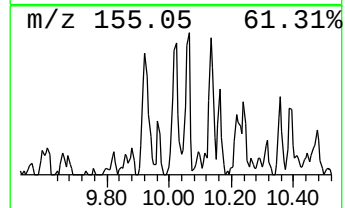
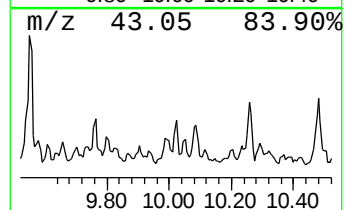
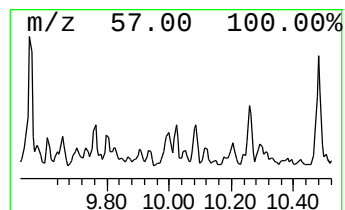
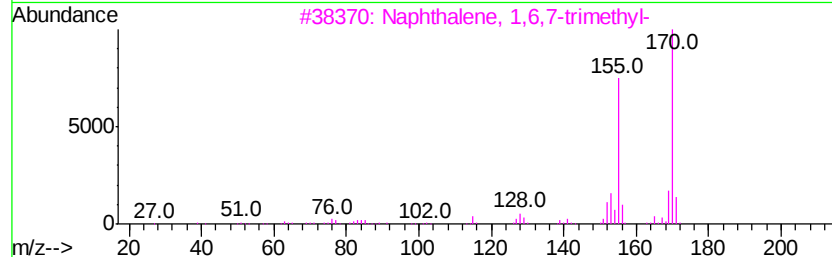
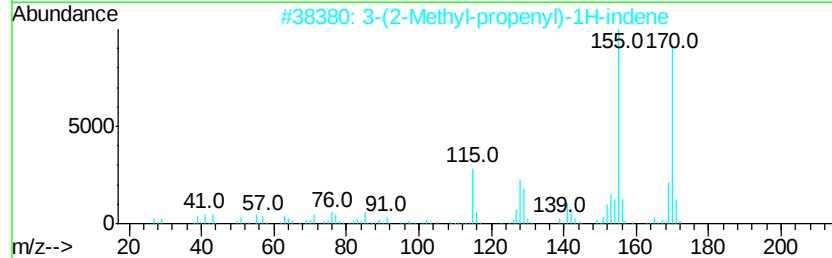
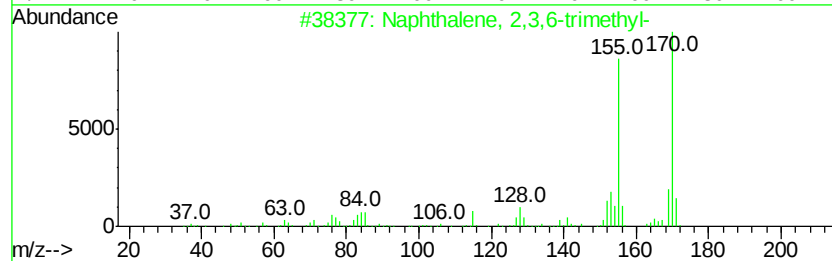
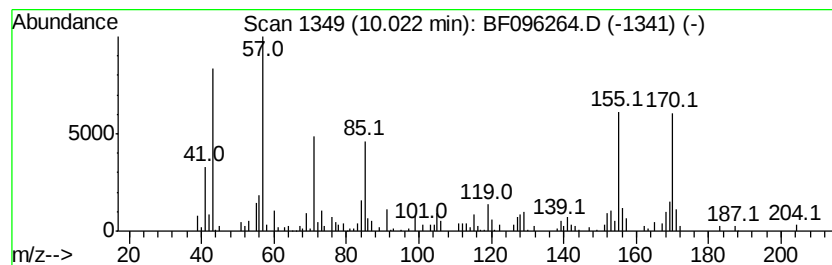
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 2,3,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.02	3.75 ng	231487	Acenaphthene-d10	9.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	64
2	3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	60
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	60
4	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	60
5	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
Data File : BF096264.D
Acq On : 28 Jun 2017 22:04
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Sample : I3911-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

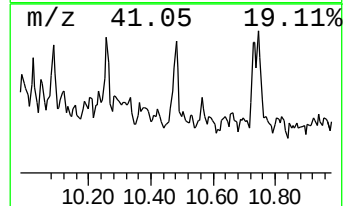
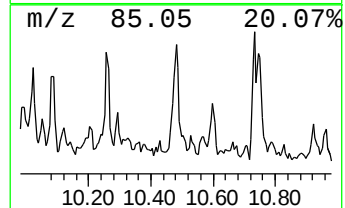
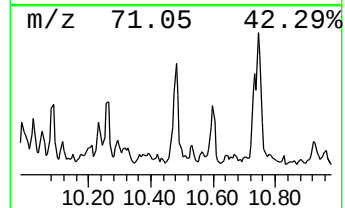
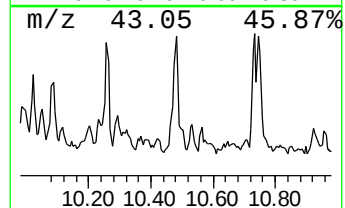
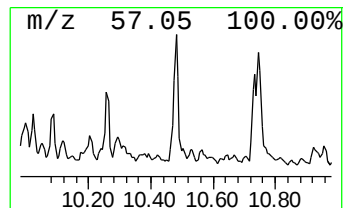
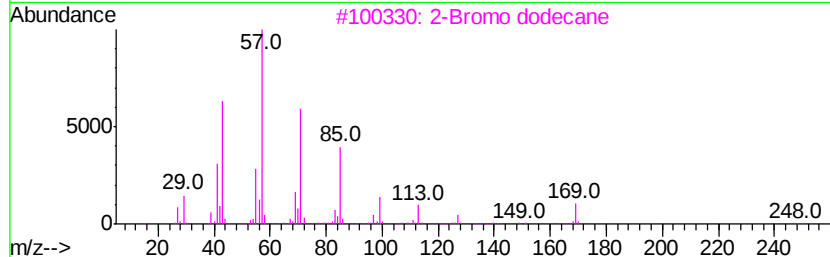
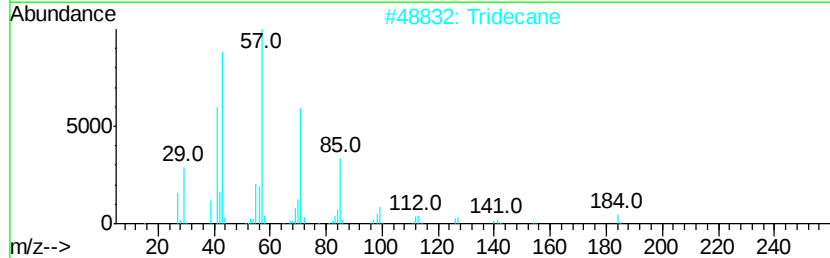
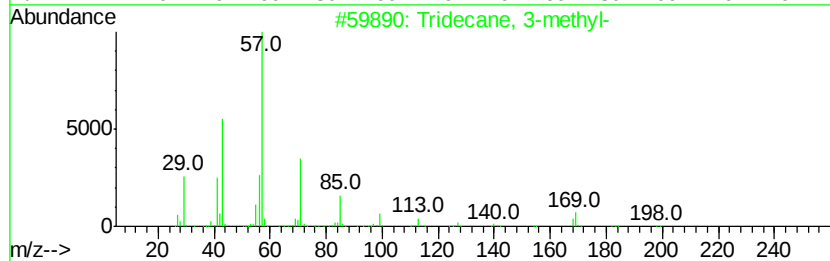
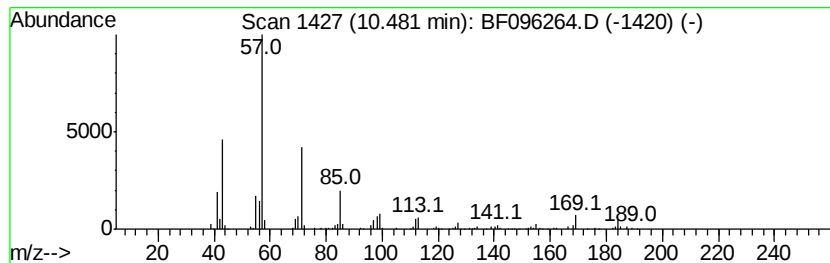
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tridecane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.48	3.72 ng	229358	Acenaphthene-d10	9.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 3-methyl-	198	C14H30	006418-41-3	72
2	Tridecane	184	C13H28	000629-50-5	68
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	64
4	Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	64
5	Heptacosane	380	C27H56	000593-49-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
Data File : BF096264.D
Acq On : 28 Jun 2017 22:04
Operator : SJ/MA
Sample : I3911-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SU-04-062317

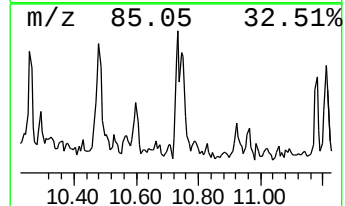
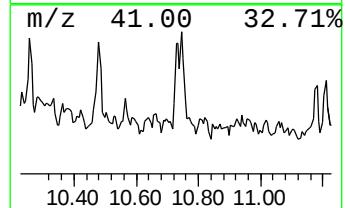
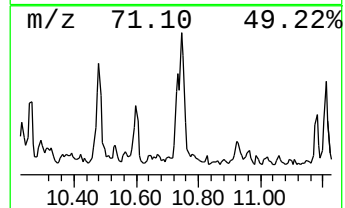
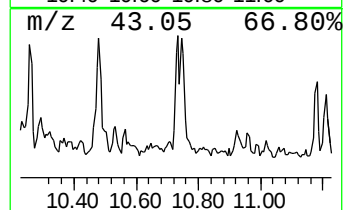
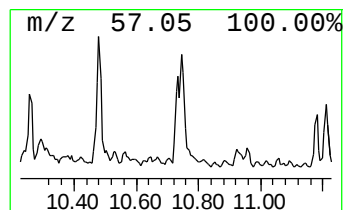
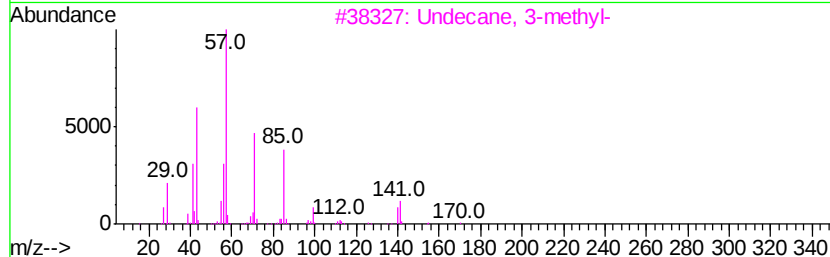
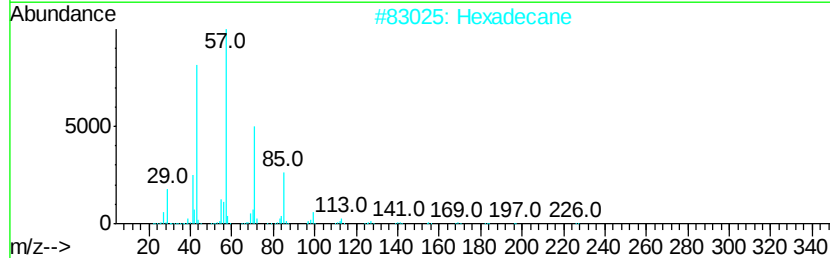
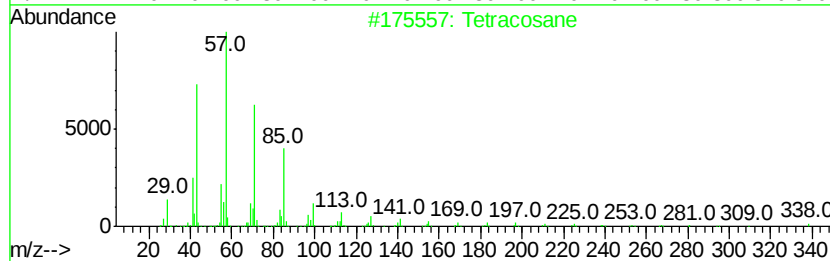
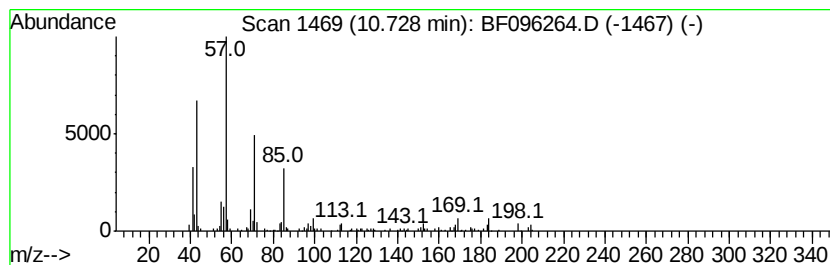
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	2.80 ng	141417	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	80
2		Hexadecane	226	C16H34	000544-76-3	72
3		Undecane, 3-methyl-	170	C12H26	001002-43-3	72
4		Tridecane	184	C13H28	000629-50-5	68
5		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
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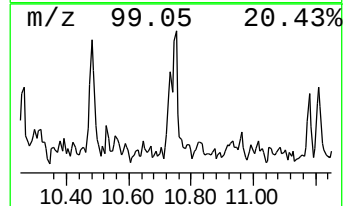
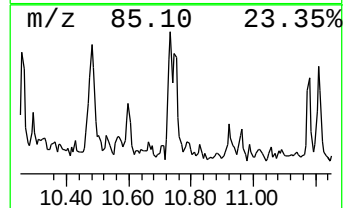
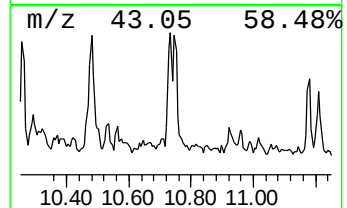
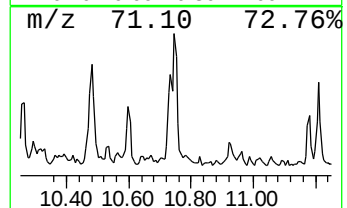
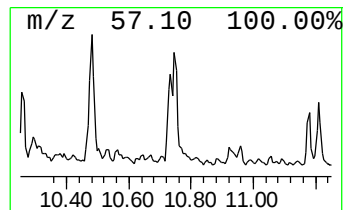
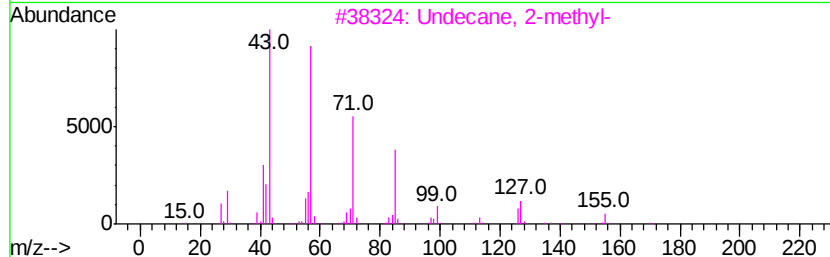
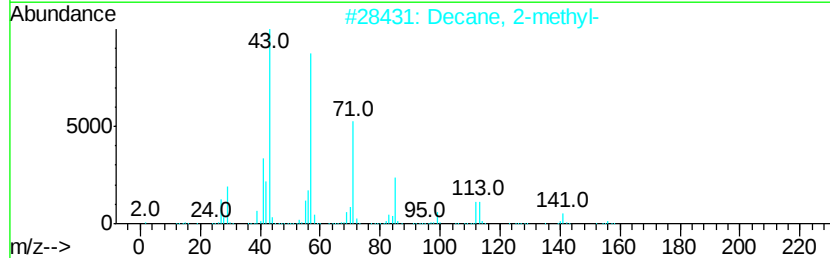
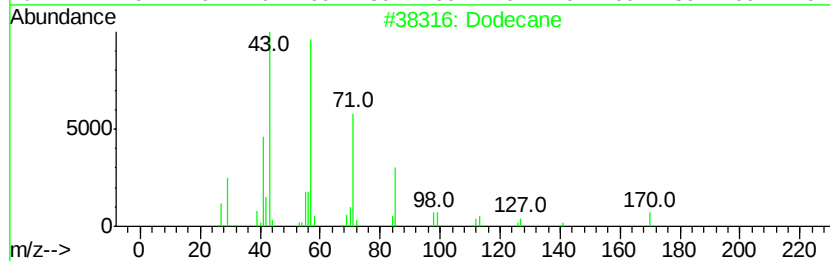
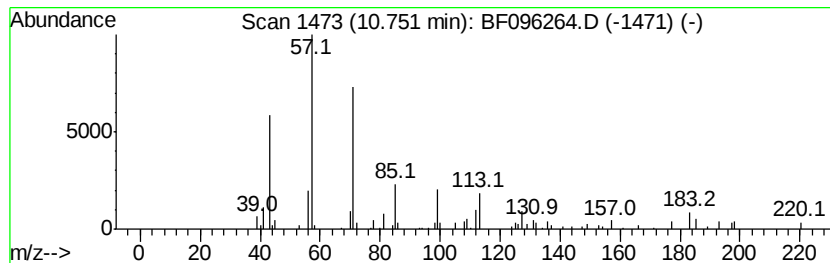
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Dodecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
10.75	2.74 ng	138219	Phenanthrene-d10		11.30		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	64
2			Decane, 2-methyl-	156	C11H24	006975-98-0	64
3			Undecane, 2-methyl-	170	C12H26	007045-71-8	59
4			Bacchotricuneatin c	342	C20H22O5	066563-30-2	58
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
Data File : BF096264.D
Acq On : 28 Jun 2017 22:04
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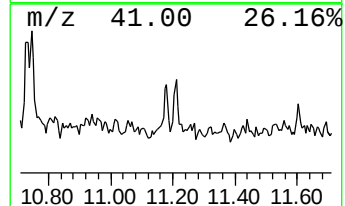
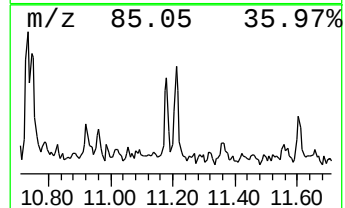
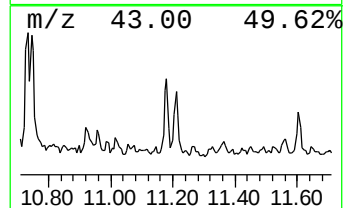
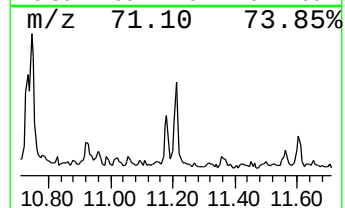
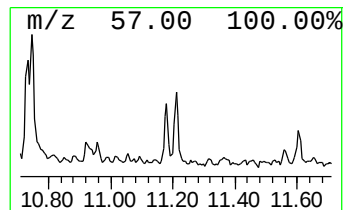
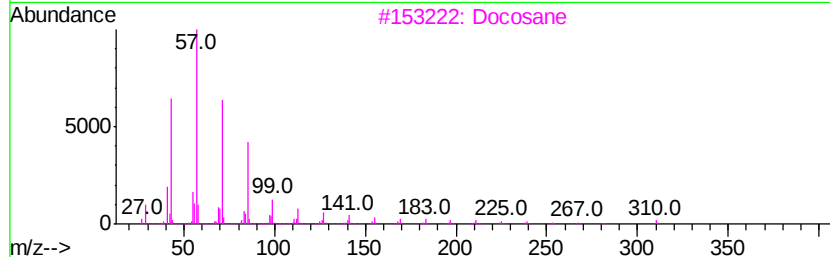
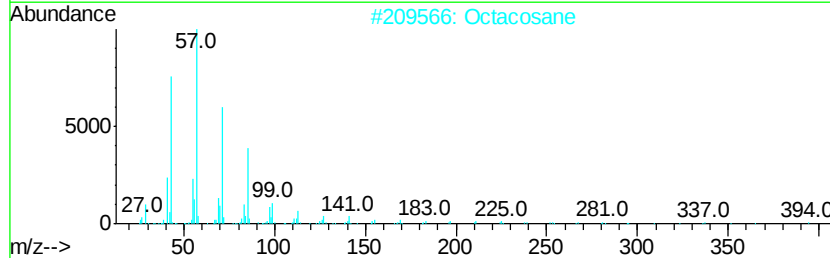
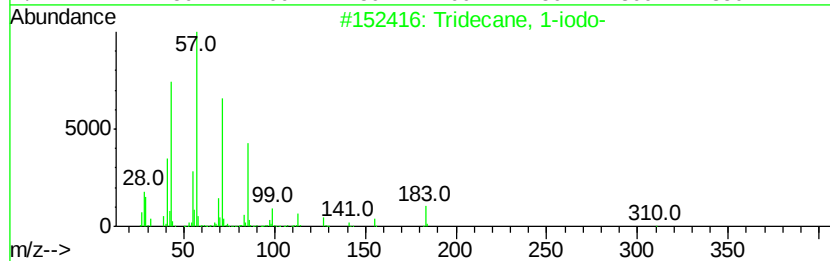
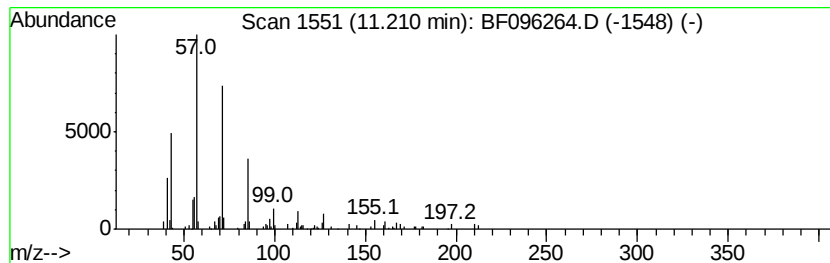
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecane, 1-iodo- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.21	2.11 ng	106476	Phenanthrene-d10	11.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86
2			Octacosane	394	C28H58	000630-02-4	78
3			Docosane	310	C22H46	000629-97-0	78
4			Heptacosane	380	C27H56	000593-49-7	78
5			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
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Acq On : 28 Jun 2017 22:04
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ALS Vial : 23 Sample Multiplier: 1

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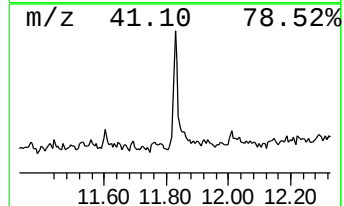
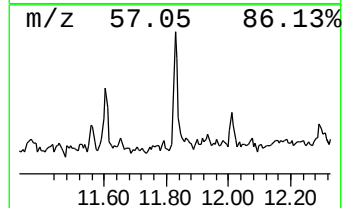
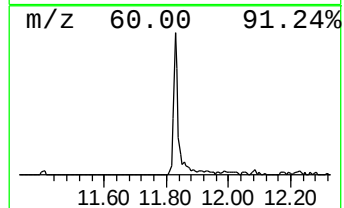
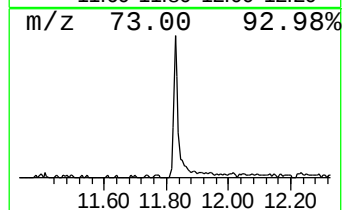
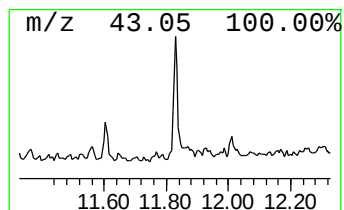
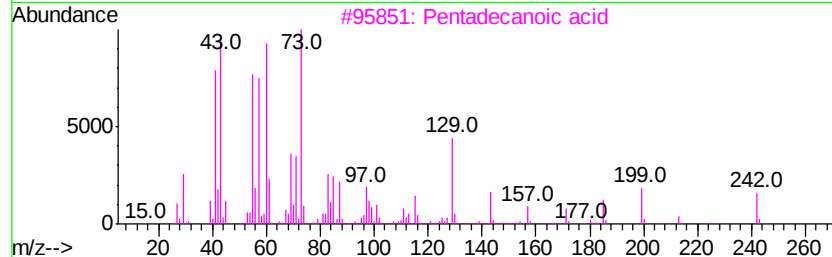
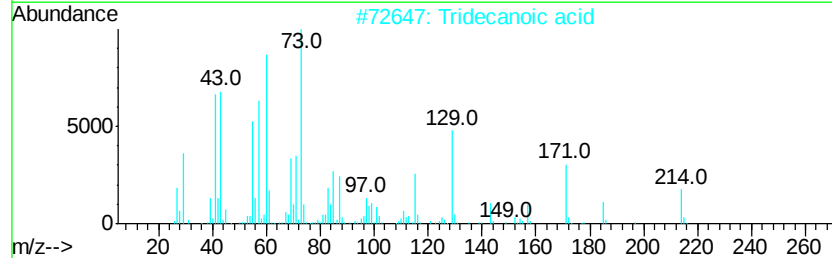
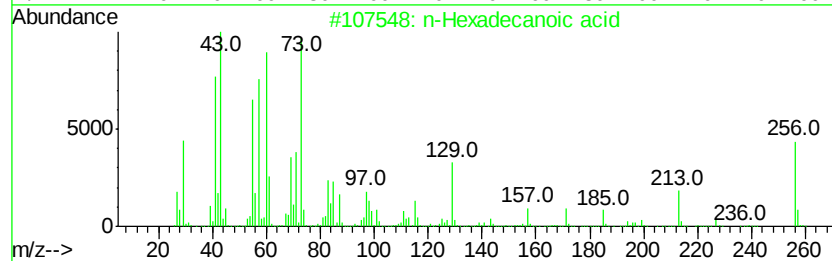
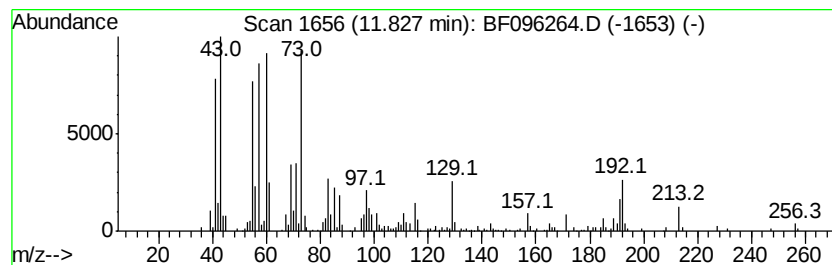
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	5.03 ng	254243	Phenanthrene-d10	11.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			Tridecanoic acid	214	C13H26O2	000638-53-9	91
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	59
5			Dodecanoic acid	200	C12H24O2	000143-07-7	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062817\
Data File : BF096264.D
Acq On : 28 Jun 2017 22:04
Operator : SJ/MA
Sample : I3911-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SU-04-062317

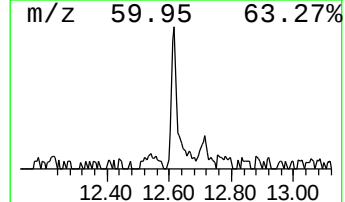
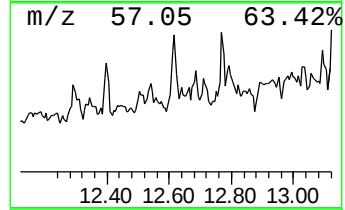
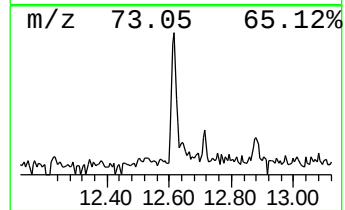
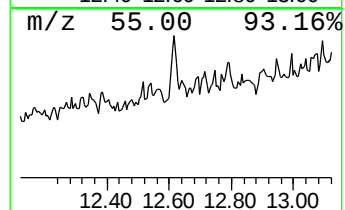
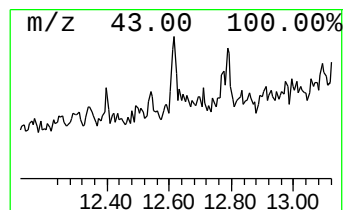
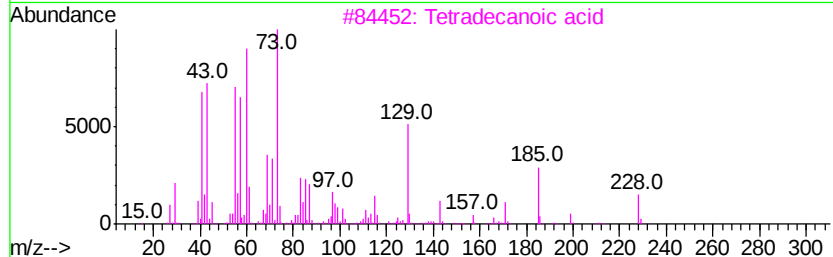
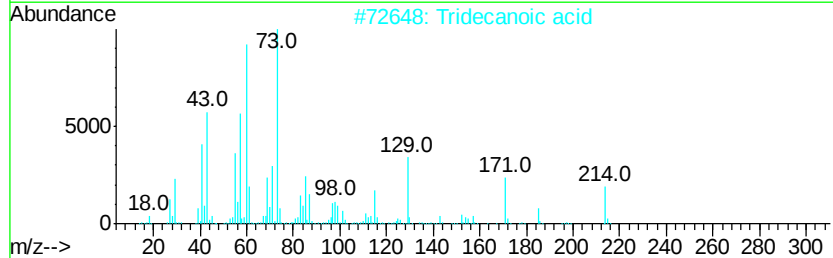
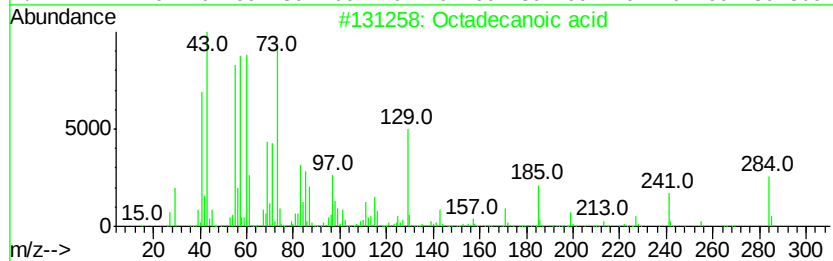
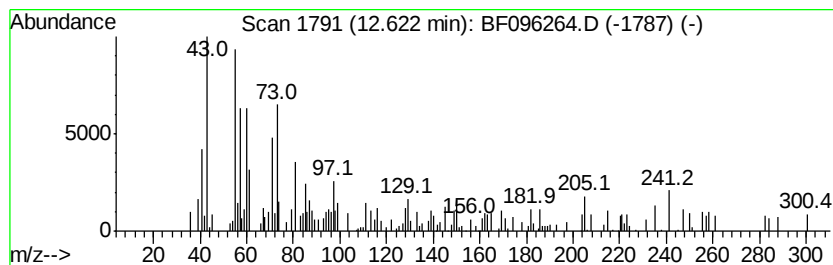
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.62	2.06 ng	95850	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2		Tridecanoic acid	214	C13H26O2	000638-53-9	62
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	60
4		Dodecanoic acid	200	C12H24O2	000143-07-7	49
5		Undecanoic acid	186	C11H22O2	000112-37-8	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
Data File : BF096264.D
Acq On : 28 Jun 2017 22:04
Operator : SJ/MA
Sample : I3911-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

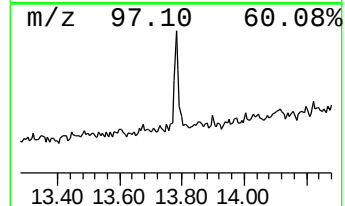
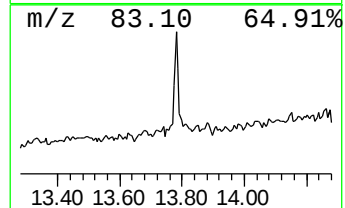
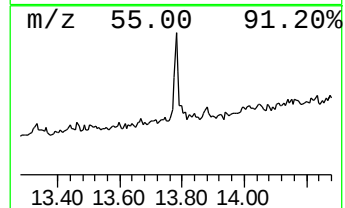
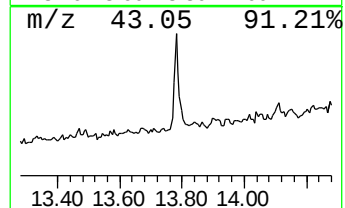
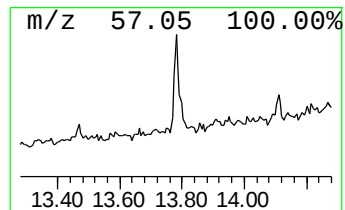
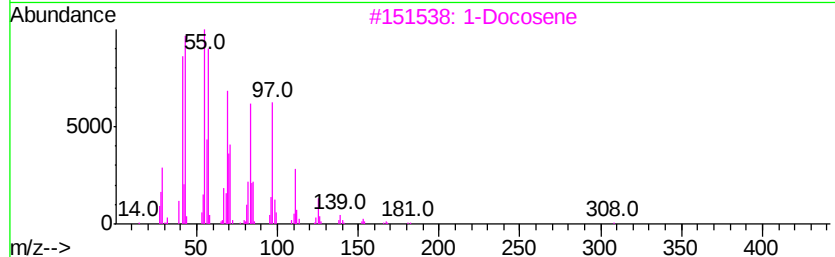
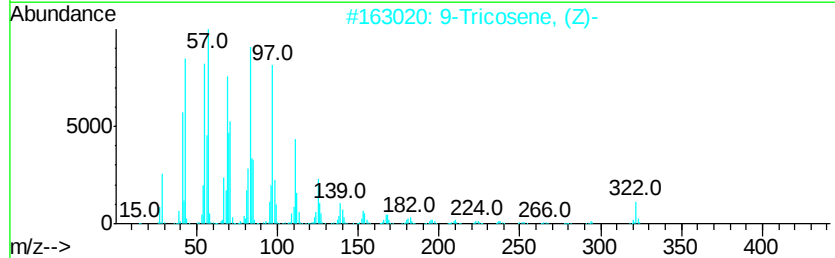
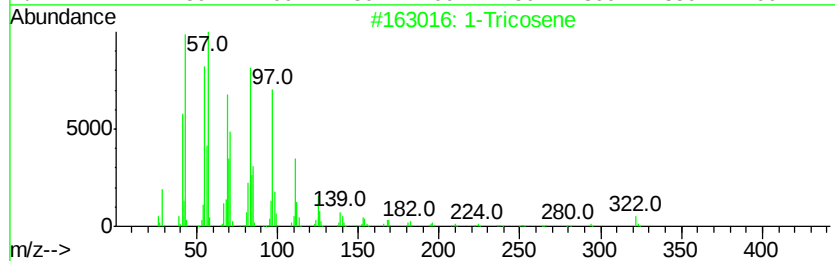
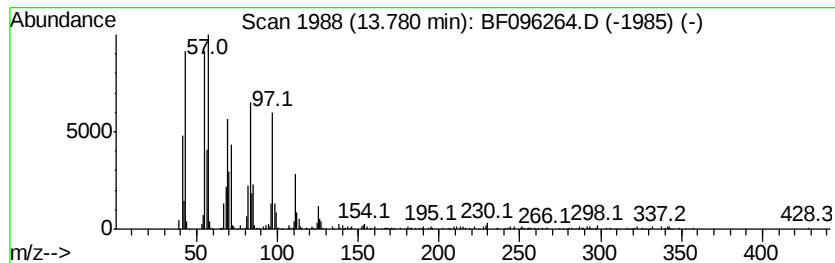
Instrument :
BNA_F
ClientSampleId :
SU-04-062317

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Tricosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.78	7.32 ng	340735	Chrysene-d12	13.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	96
2	9-Tricosene, (Z)-	322	C23H46	027519-02-4	94
3	1-Docosene	308	C22H44	001599-67-3	94
4	Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062817\
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Instrument :
 BNA_F
 ClientSampleId :
 SU-04-062317

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF062717.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...	4.98	18.7	ng	883622	1	6.79	944669	20.0
unknown6.55	6.55	79.3	ng	3745740	1	6.79	944669	20.0
Octane, 2,6-dimet...	8.50	3.1	ng	200696	2	8.06	1300440	20.0
unknown8.76	8.76	2.1	ng	134502	2	8.06	1300440	20.0
unknown9.23	9.23	2.6	ng	158404	3	9.82	1234540	20.0
Naphthalene, 2,3,...	10.02	3.8	ng	231487	3	9.82	1234540	20.0
Tridecane, 3-methyl-	10.48	3.7	ng	229358	3	9.82	1234540	20.0
Tetracosane	10.73	2.8	ng	141417	4	11.30	1009930	20.0
Dodecane	10.75	2.7	ng	138219	4	11.30	1009930	20.0
Tridecane, 1-iodo-	11.21	2.1	ng	106476	4	11.30	1009930	20.0
n-Hexadecanoic acid	11.83	5.0	ng	254243	4	11.30	1009930	20.0
Octadecanoic acid	12.62	2.1	ng	95850	5	13.93	930736	20.0
1-Tricosene	13.78	7.3	ng	340735	5	13.93	930736	20.0