

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
 Data File : BF097398.D
 Acq On : 5 Aug 2017 15:41
 Operator : SJ/JU
 Sample : I4592-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 15-50-RT3460-RT4215-RT3451

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-BF072517.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	4.551	466	470	488	rBV	412311	714006	24.41%	2.698%
2	5.022	546	550	576	rBV	2264653	2755295	94.18%	10.413%
3	6.098	729	733	746	rBV	2598084	2925460	100.00%	11.056%
4	6.198	746	750	758	rVB	2733867	2654891	90.75%	10.033%
5	6.422	785	788	798	rVB	833357	755117	25.81%	2.854%
6	6.575	811	814	822	rBV	1242288	1275931	43.61%	4.822%
7	6.828	854	857	865	rBV2	17393	42676	1.46%	0.161%
8	6.998	881	886	901	rBV	1457467	1596504	54.57%	6.033%
9	7.286	930	935	944	rBV2	23127	39295	1.34%	0.149%
10	7.704	1003	1006	1013	rBV	1191221	1098957	37.57%	4.153%
11	7.763	1013	1016	1019	rVB	52753	49273	1.68%	0.186%
12	7.898	1034	1039	1042	rBV	1296796	1397079	47.76%	5.280%
13	8.104	1071	1074	1079	rBV	54824	58172	1.99%	0.220%
14	8.322	1107	1111	1114	rVB	35555	31871	1.09%	0.120%
15	8.357	1114	1117	1123	rVB2	27329	30328	1.04%	0.115%
16	8.657	1165	1168	1176	rBV2	102819	137341	4.69%	0.519%
17	8.727	1176	1180	1185	rBV2	48802	71300	2.44%	0.269%
18	8.786	1185	1190	1193	rVV	1621840	1580755	54.03%	5.974%
19	8.880	1202	1206	1211	rBV2	52195	52787	1.80%	0.199%
20	9.186	1255	1258	1263	rBV	443737	381800	13.05%	1.443%
21	9.310	1276	1279	1289	rBV	34333	51642	1.77%	0.195%
22	9.410	1291	1296	1299	rBV	71088	68175	2.33%	0.258%
23	9.451	1299	1303	1307	rVV	1140439	1044896	35.72%	3.949%
24	9.904	1377	1380	1383	rBV	82367	65246	2.23%	0.247%
25	10.121	1412	1417	1420	rBV	45050	61425	2.10%	0.232%
26	10.239	1433	1437	1439	rBV	1353817	1328523	45.41%	5.021%
27	10.263	1439	1441	1449	rVB	169311	174689	5.97%	0.660%
28	10.374	1457	1460	1467	rBV	122797	148147	5.06%	0.560%
29	10.568	1489	1493	1496	rBV5	25394	38226	1.31%	0.144%
30	10.816	1532	1535	1538	rBV	80308	75529	2.58%	0.285%
31	10.921	1549	1553	1557	rBV	1215378	1059369	36.21%	4.004%
32	10.992	1561	1565	1570	rVB6	17573	31504	1.08%	0.119%
33	11.239	1605	1607	1613	rBV	58759	53777	1.84%	0.203%
34	11.480	1645	1648	1655	rBV	351776	333116	11.39%	1.259%

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 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	11.674	1678	1681	1686	rVB	340139	306999	10.49%	1.160%
36	11.880	1710	1716	1719	rBV	31550	38787	1.33%	0.147%
37	12.127	1756	1758	1761	rBV	36481	36431	1.25%	0.138%
38	12.162	1761	1764	1770	rVB6	17333	35884	1.23%	0.136%
39	12.257	1777	1780	1791	rVV	219982	236315	8.08%	0.893%
40	12.351	1793	1796	1802	rVB2	46043	56546	1.93%	0.214%
41	12.504	1817	1822	1825	rBV	1586467	1480907	50.62%	5.597%
42	12.751	1857	1864	1868	rBV6	21427	37422	1.28%	0.141%
43	13.074	1913	1919	1921	rBV3	24113	43546	1.49%	0.165%
44	13.133	1925	1929	1932	rVB	27163	37898	1.30%	0.143%
45	13.304	1955	1958	1963	rVB4	29810	44974	1.54%	0.170%
46	13.421	1974	1978	1985	rVB2	78341	115023	3.93%	0.435%
47	13.539	1994	1998	2007	rVB	675699	664133	22.70%	2.510%
48	13.915	2059	2062	2067	rBV	78382	71887	2.46%	0.272%
49	14.539	2165	2168	2174	rVB	25639	36949	1.26%	0.140%
50	14.880	2222	2226	2230	rBV	524020	596711	20.40%	2.255%
51	15.262	2287	2291	2295	rBV2	33443	46247	1.58%	0.175%
52	15.656	2353	2358	2366	rBV2	29255	53145	1.82%	0.201%
53	15.756	2369	2375	2383	rBV	78196	132637	4.53%	0.501%
54	16.115	2430	2436	2442	rVB4	24255	50431	1.72%	0.191%
55	16.274	2457	2463	2469	rVB2	36845	67833	2.32%	0.256%
56	16.650	2523	2527	2536	rVB5	17767	39972	1.37%	0.151%
57	17.292	2628	2636	2642	rBV8	17956	47067	1.61%	0.178%

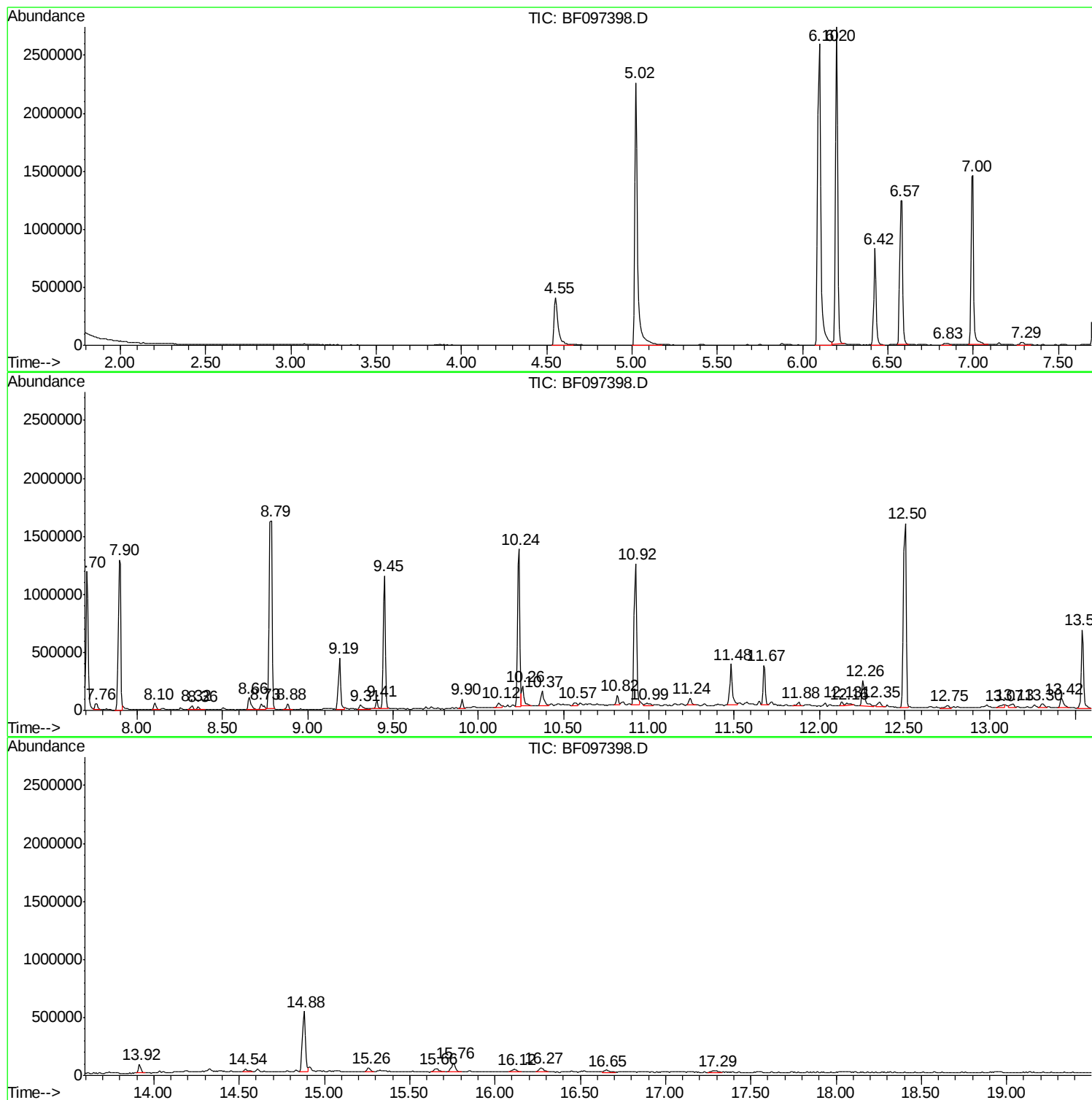
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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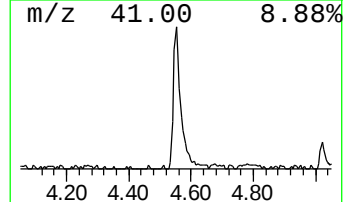
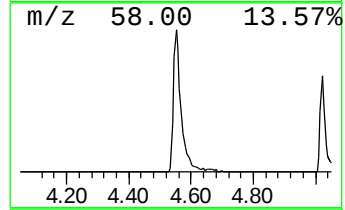
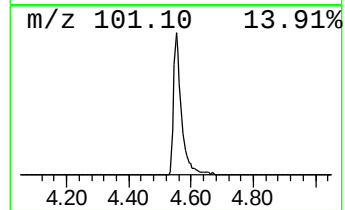
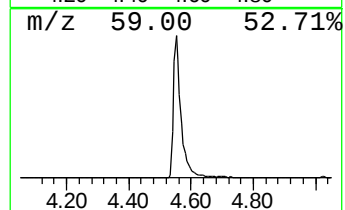
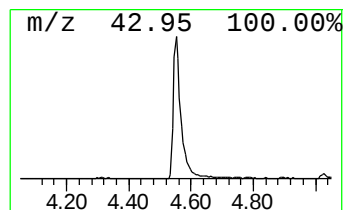
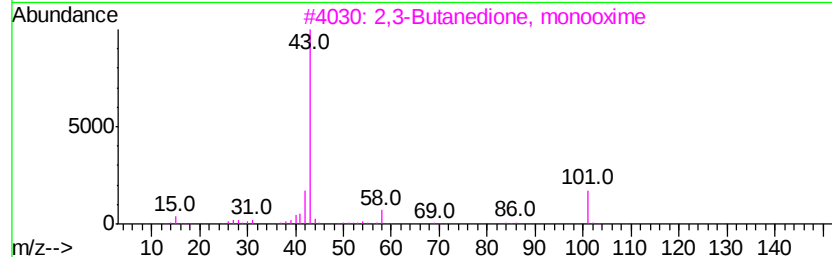
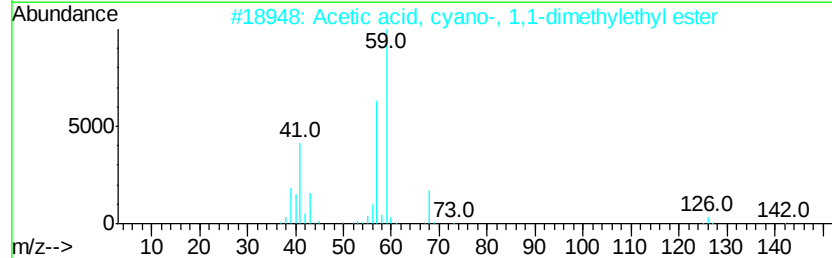
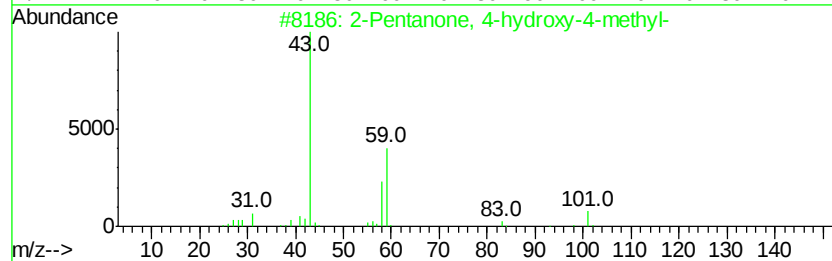
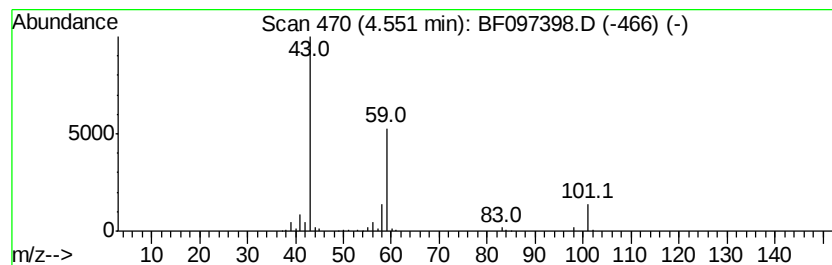
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.55	18.91 ng	714006	1,4-Dichlorobenzene-d4	6.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		1,2-Butanediol	90	C4H10O2	000584-03-2	9



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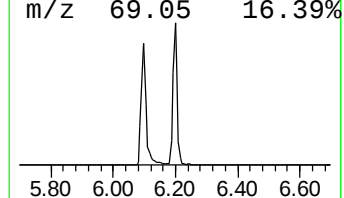
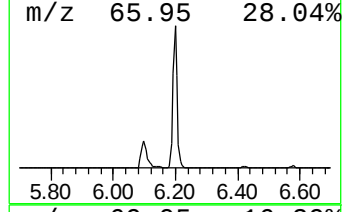
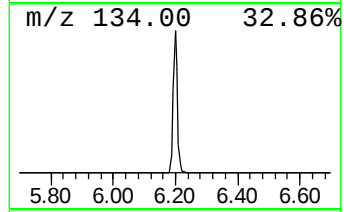
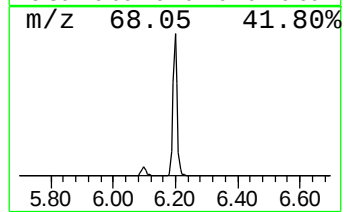
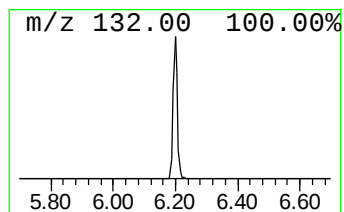
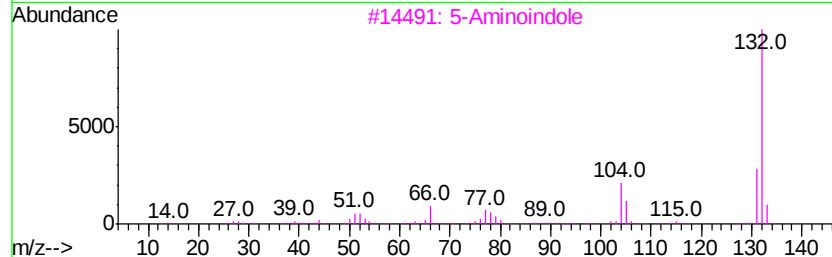
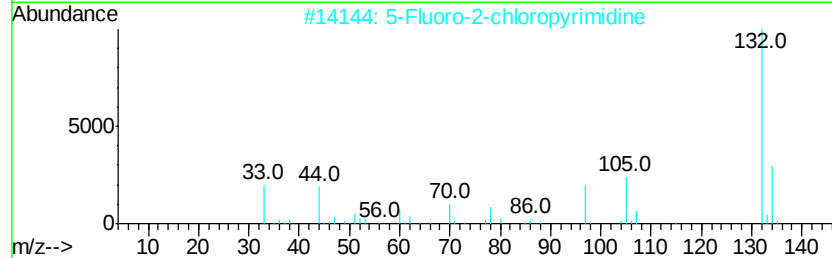
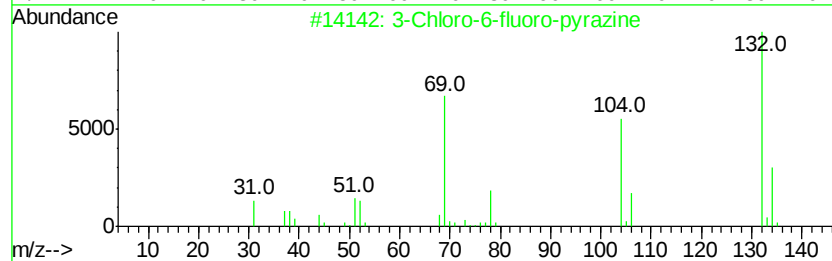
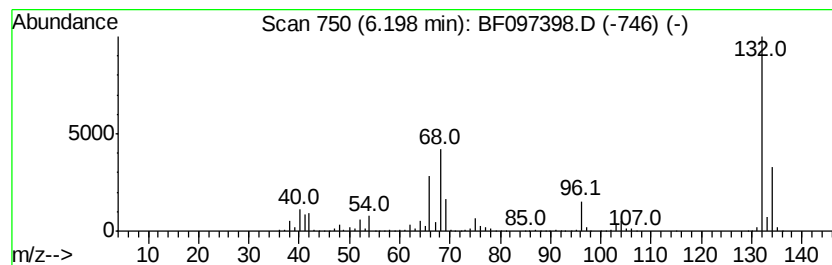
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.20	70.32 ng	2654890	1,4-Dichlorobenzene-d4	6.42

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		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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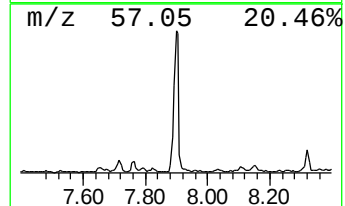
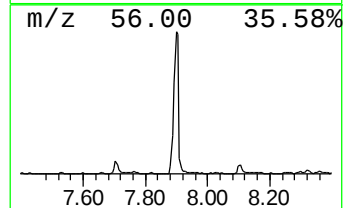
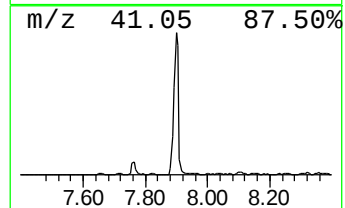
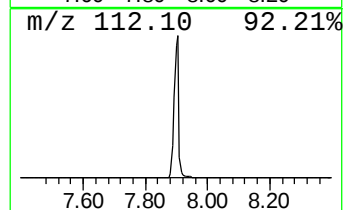
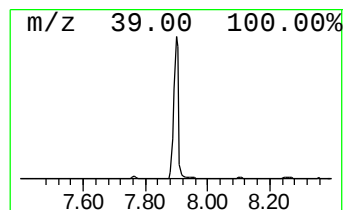
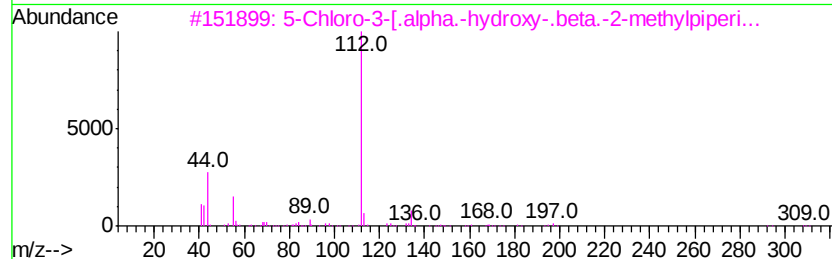
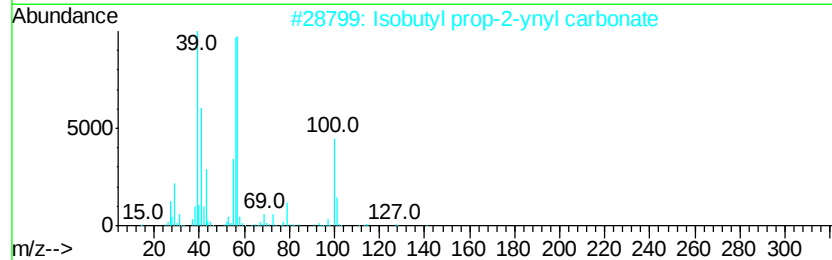
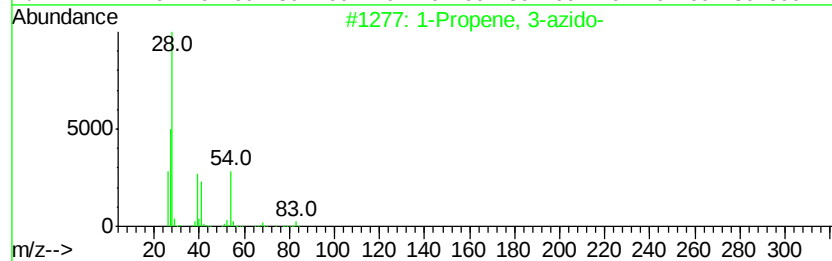
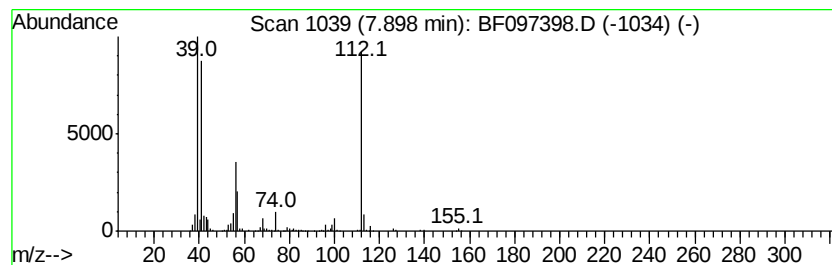
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.90 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	25.43 ng	1397080	Naphthalene-d8	7.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propene, 3-azido-	83	C3H5N3	000821-13-6	36
2		Isobutyl prop-2-ynyl carbonate	156	C8H12O3	1000372-90-0	36
3		5-Chloro-3-[.alpha.-hydroxy-.bet...	309	C16H20ClNOS	1000251-01-5	28
4		N-Methyl-tetrahydro-solasodine	431	C28H49NO2	1000256-11-3	25
5		2-Furancarboxylic acid	112	C5H4O3	000088-14-2	9



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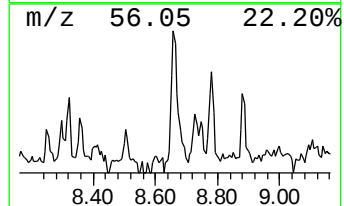
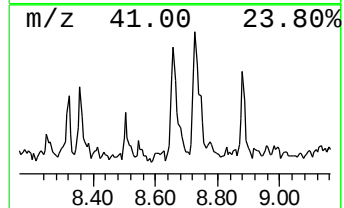
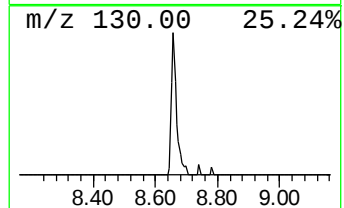
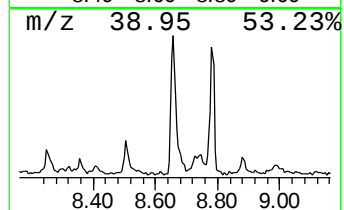
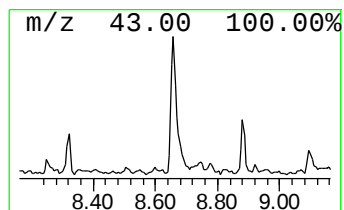
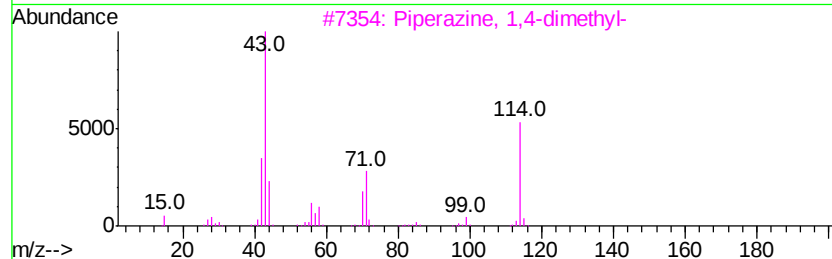
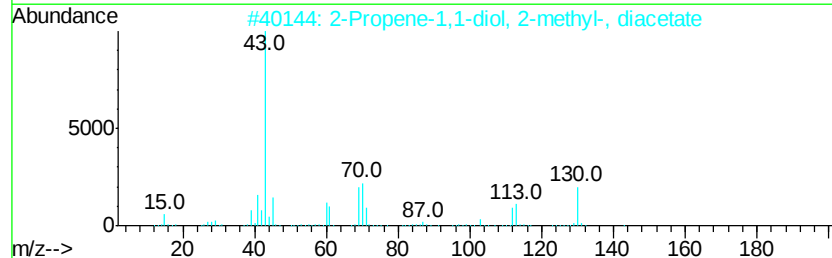
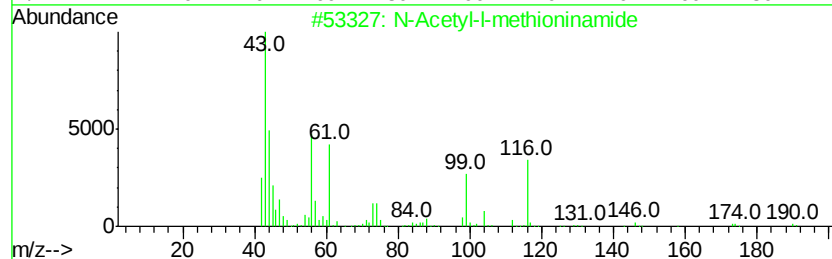
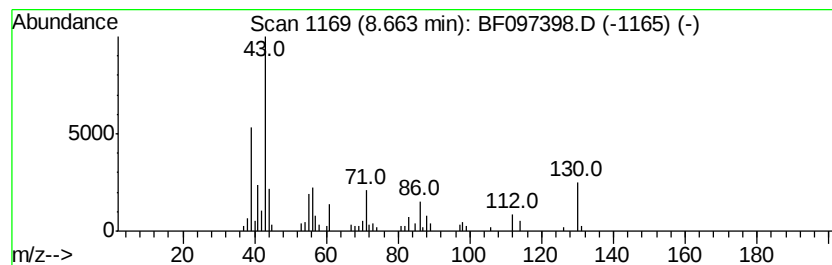
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown8.66 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.66	2.63 ng	137341	Acenaphthene-d10	9.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N-Acetyl-L-methioninamide	190	C7H14N2O2S	1000130-47-4	12
2		2-Propene-1,1-diol, 2-methyl-, d...	172	C8H12O4	010476-95-6	12
3		Piperazine, 1,4-dimethyl-	114	C6H14N2	000106-58-1	10
4		Propane, 2-(ethenyl)-	86	C5H10	000926-65-8	10
5		2-Butanone, 3-methyl-	86	C5H10	000563-80-4	9



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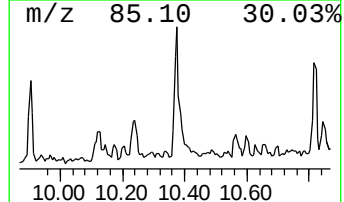
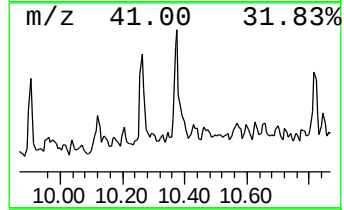
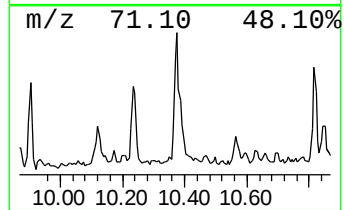
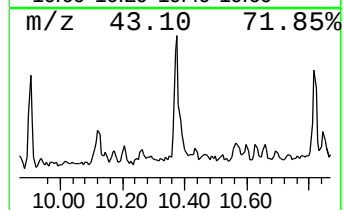
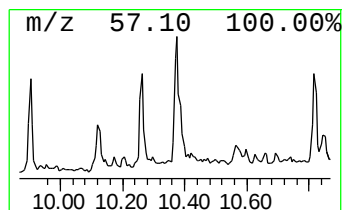
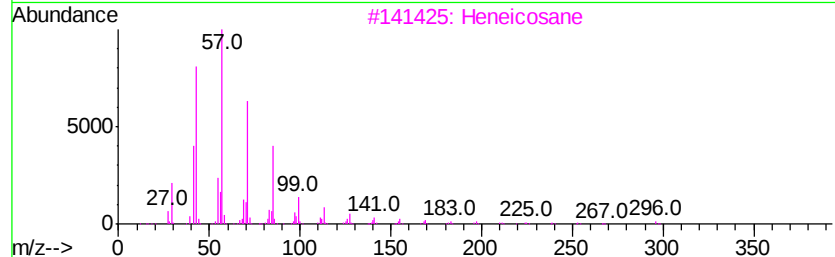
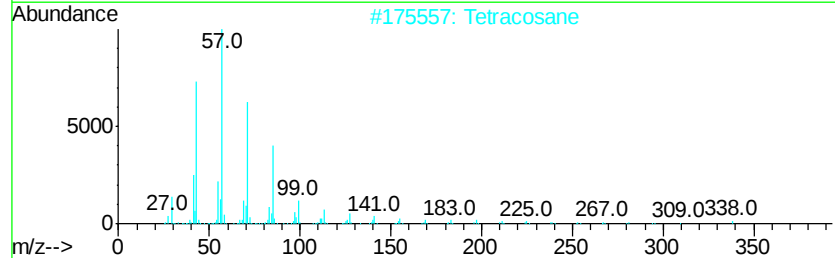
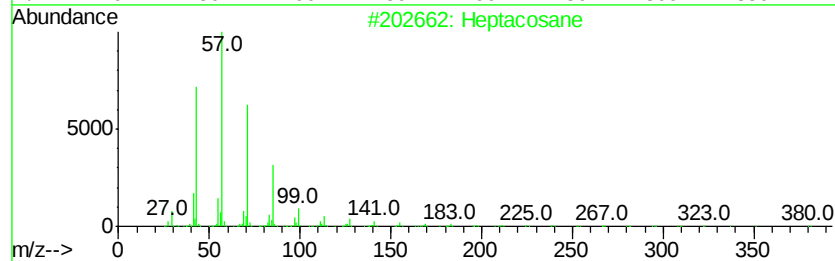
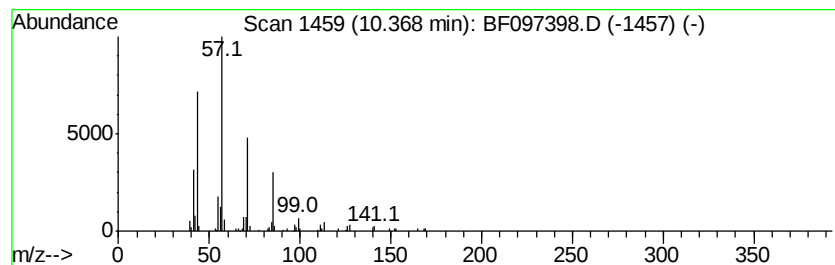
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	2.80 ng	148147	Phenanthrene-d10	10.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Tetracosane	338	C24H50	000646-31-1	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Eicosane	282	C20H42	000112-95-8	87
5		Eicosane, 3-methyl-	296	C21H44	006418-46-8	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
Sample : I4592-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
15-50-RT3460-RT4215-RT3451

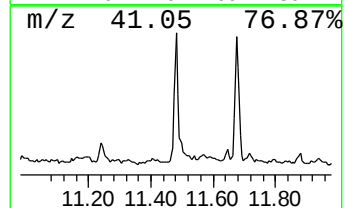
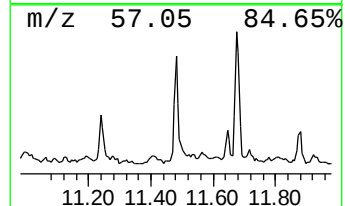
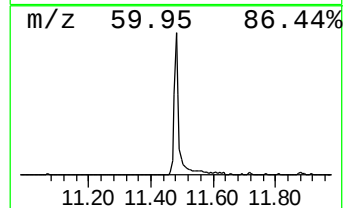
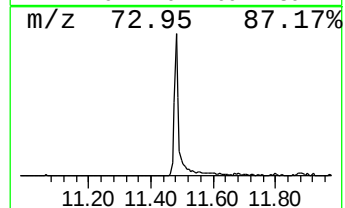
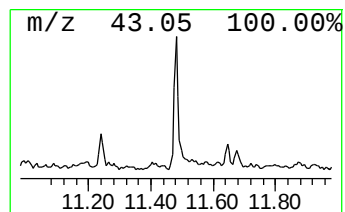
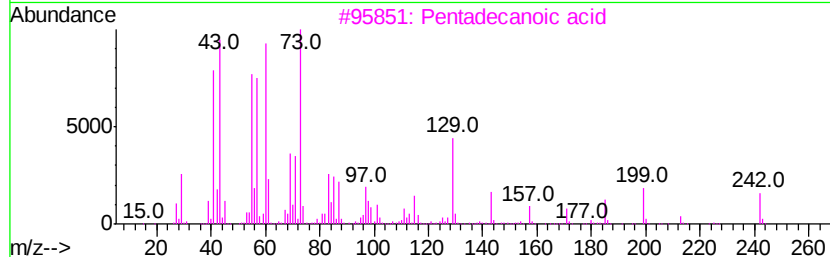
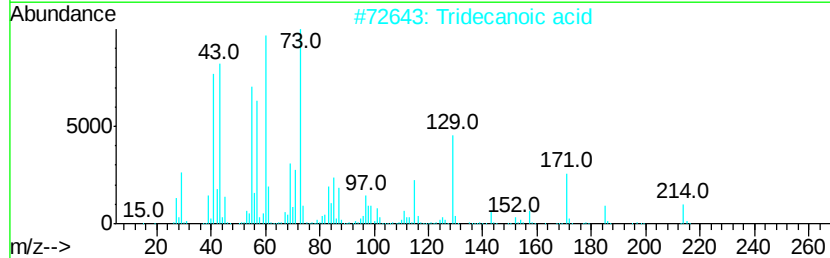
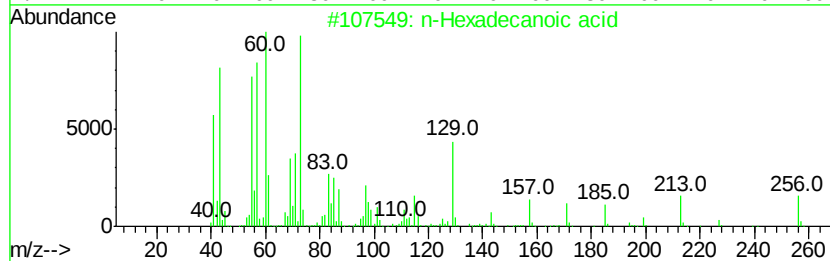
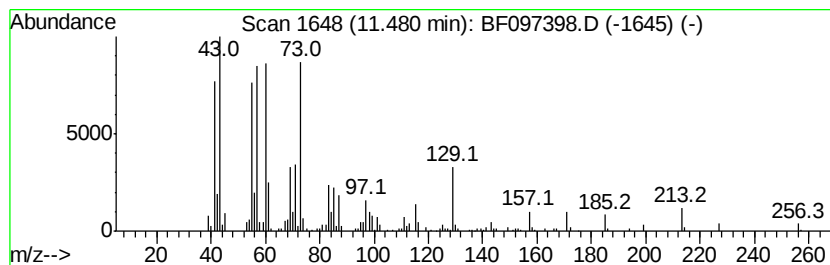
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	6.29 ng	333116	Phenanthrene-d10	10.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2			Tridecanoic acid	214	C13H26O2	000638-53-9	91
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4			Undecanoic acid	186	C11H22O2	000112-37-8	72
5			Dodecanoic acid	200	C12H24O2	000143-07-7	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
Sample : I4592-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
15-50-RT3460-RT4215-RT3451

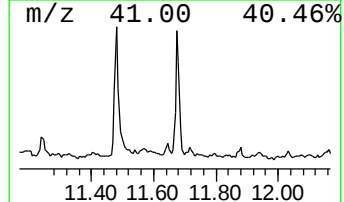
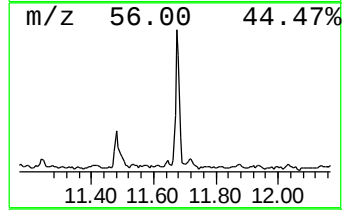
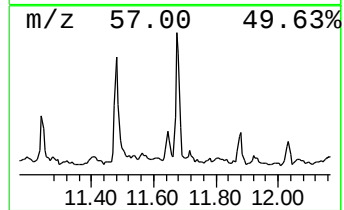
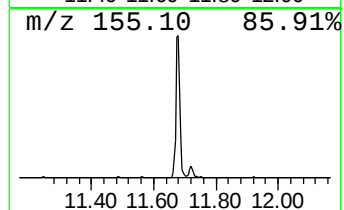
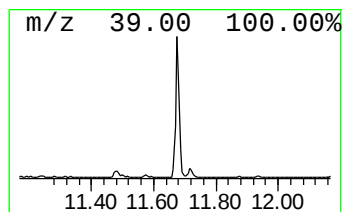
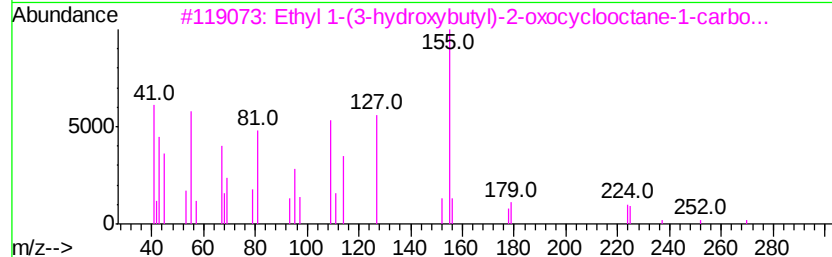
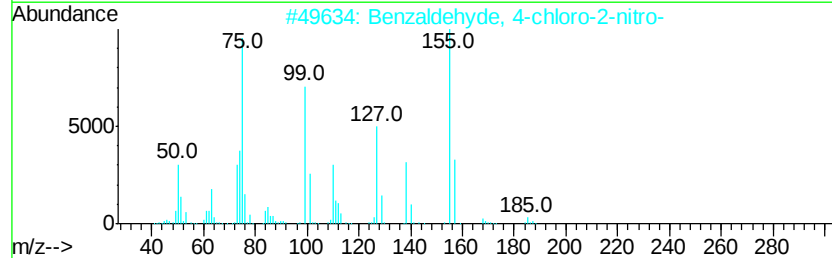
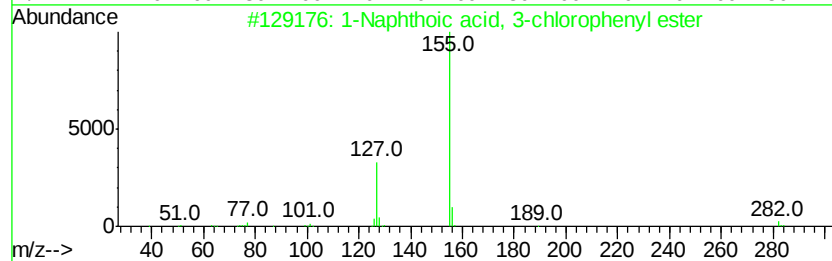
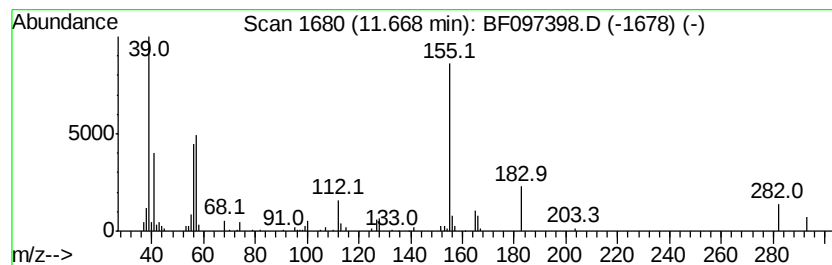
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown11.67 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.67	5.80 ng	306999	Phenanthrene-d10	10.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthoic acid, 3-chlorophenyl...	282	C17H11ClO2	1000325-54-1	38
2		Benzaldehyde, 4-chloro-2-nitro-	185	C7H4ClNO3	005551-11-1	35
3		Ethyl 1-(3-hydroxybutyl)-2-oxocyclo...	270	C15H26O4	110288-16-9	27
4		1,3-Benzenediol, 0-(2,6-difluoro...	404	C24H14F2O4	1000345-52-7	25
5		1,2-Benzenediol, 0-(1-naphthoyl)...	382	C25H18O4	1000330-48-2	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
Sample : I4592-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

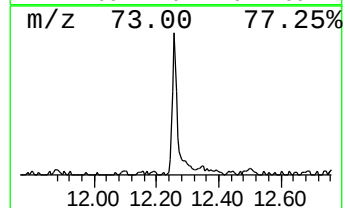
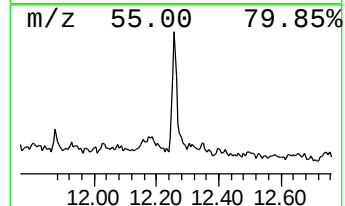
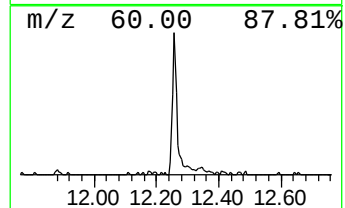
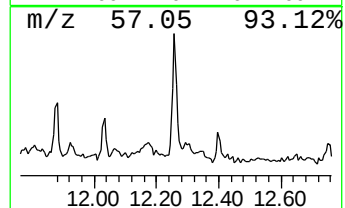
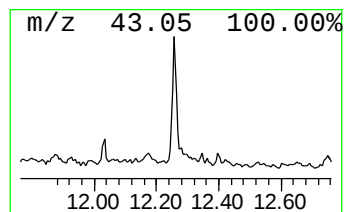
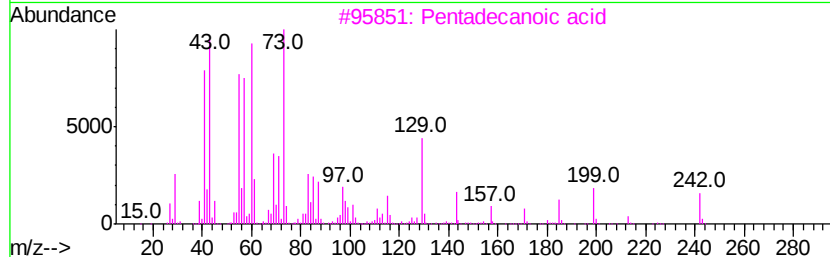
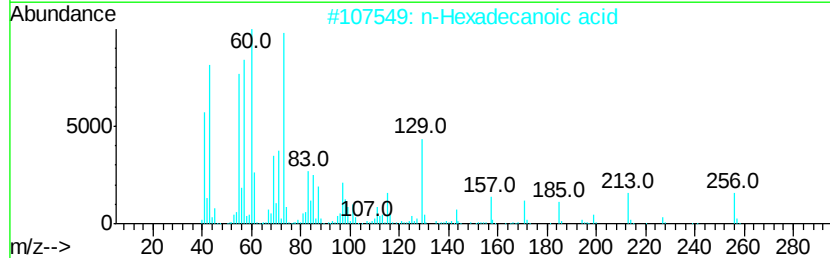
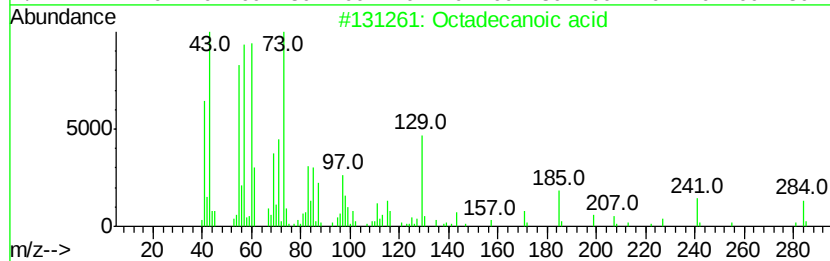
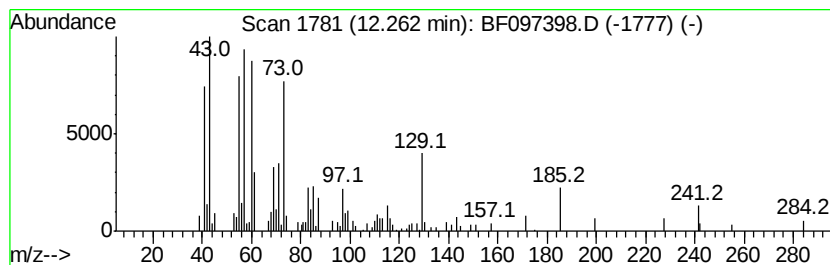
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15-50-RT3460-RT4215-RT3451

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.26	7.12 ng	236315	Chrysene-d12	13.54	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	94
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	60



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
Sample : I4592-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
15-50-RT3460-RT4215-RT3451

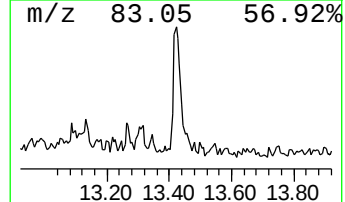
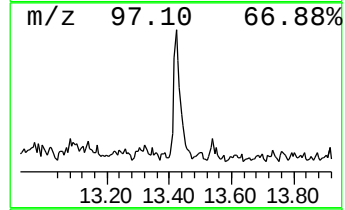
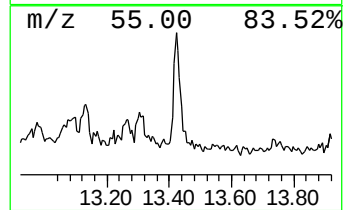
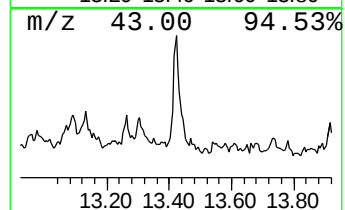
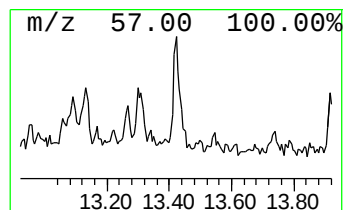
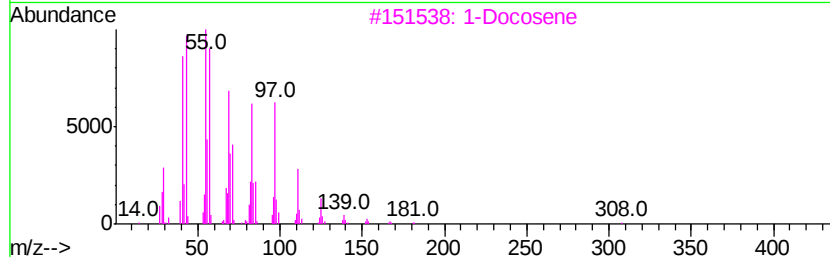
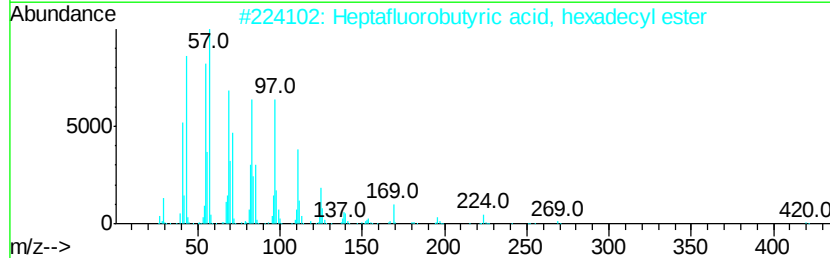
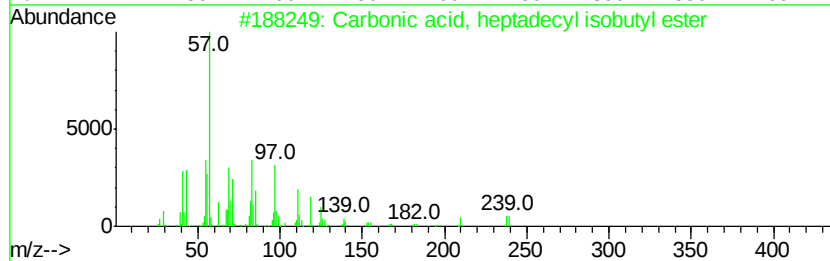
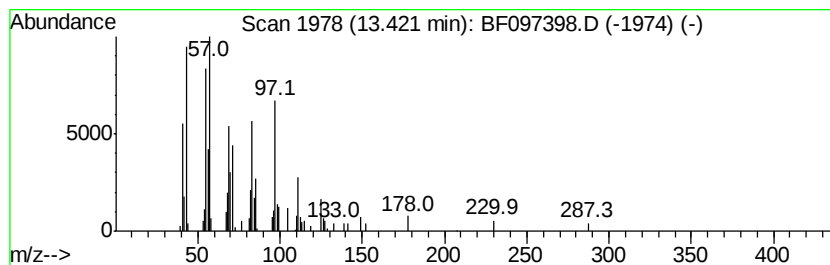
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Carbonic acid, heptadecyl i... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	3.46 ng	115023	Chrysene-d12	13.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, heptadecyl isobut...	356	C22H44O3	1000314-61-4	91
2			Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	90
3			1-Docosene	308	C22H44	001599-67-3	90
4			Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	83
5			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
Sample : I4592-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
15-50-RT3460-RT4215-RT3451

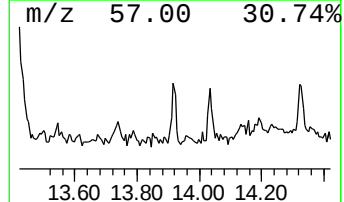
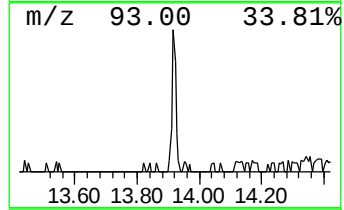
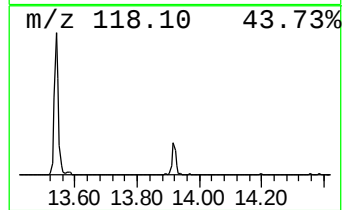
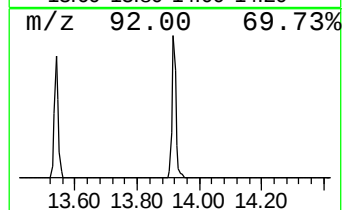
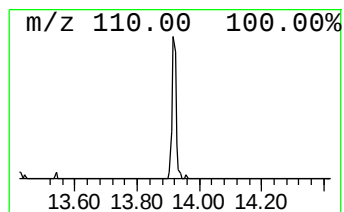
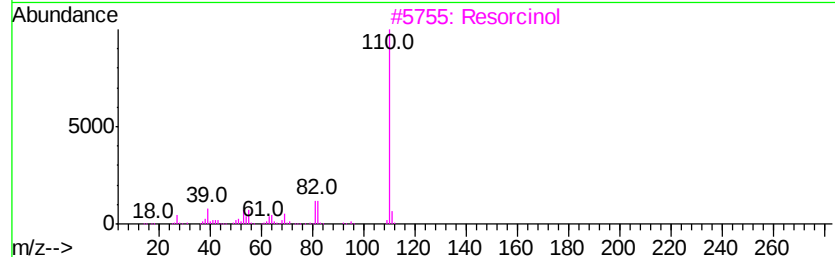
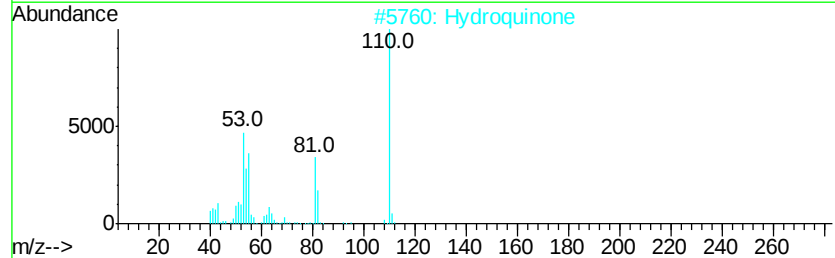
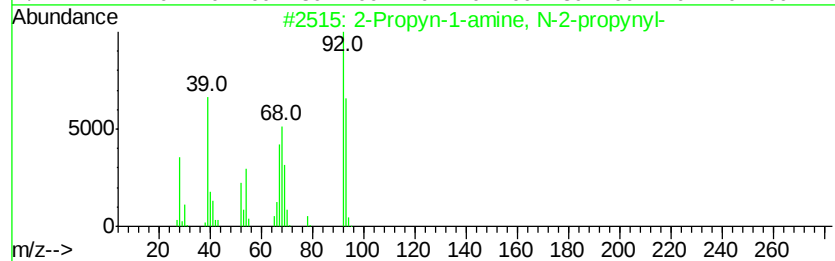
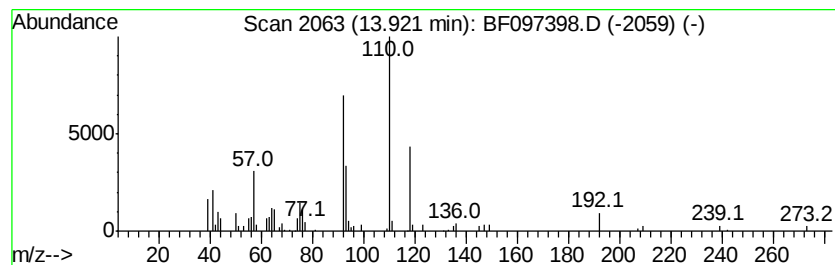
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown13.92 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	2.16 ng	71887	Chrysene-d12	13.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propyn-1-amine, N-2-propynyl-	93	C6H7N	006921-28-4	30
2			Hydroquinone	110	C6H6O2	000123-31-9	18
3			Resorcinol	110	C6H6O2	000108-46-3	14
4			4,5-Pyrimidinediamine	110	C4H6N4	013754-19-3	14
5			1H-Pyrrole-2-carbonitrile	92	C5H4N2	009513-94-4	12



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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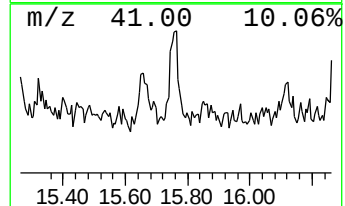
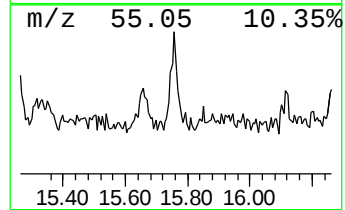
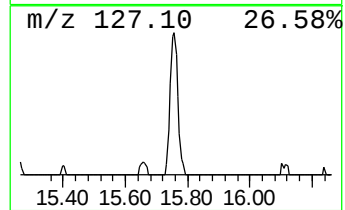
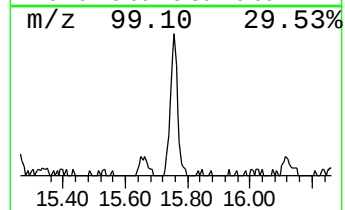
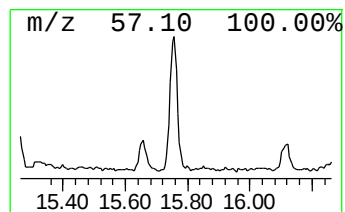
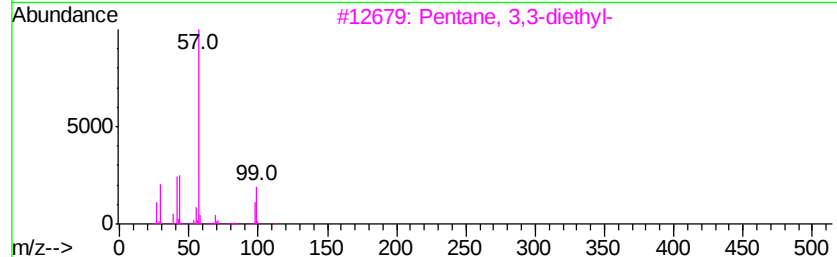
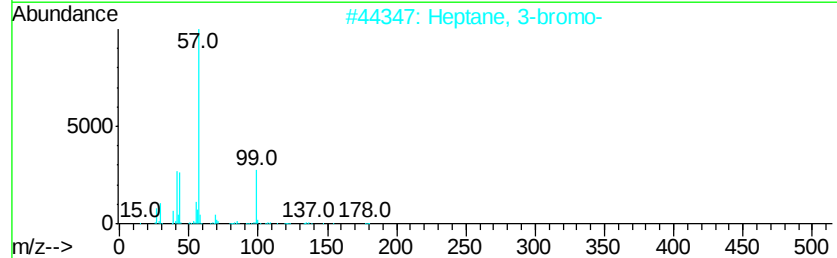
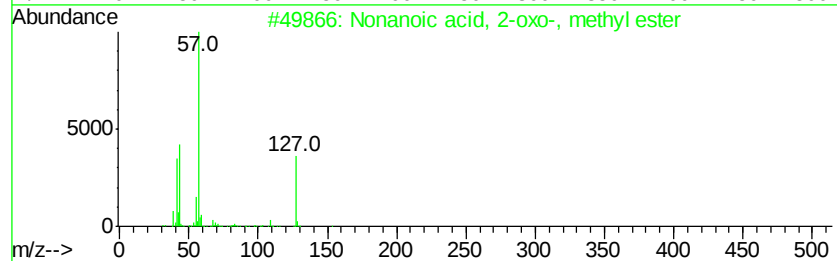
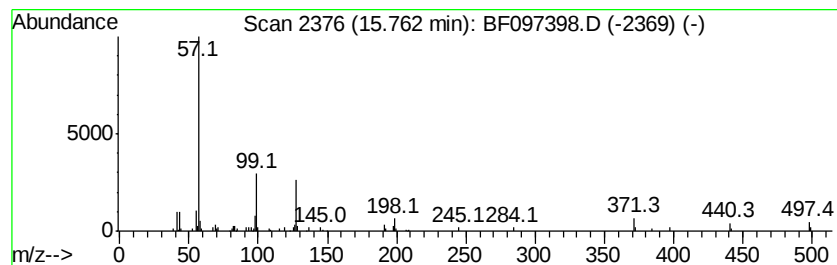
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.76 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	4.45 ng	132637	Perylene-d12	14.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanoic acid, 2-oxo-, methyl ester	186	C10H18O3	056275-54-8	35
2			Heptane, 3-bromo-	178	C7H15Br	001974-05-6	25
3			Pentane, 3,3-diethyl-	128	C9H20	001067-20-5	25
4			2,2,4,4-Tetramethyloctane	170	C12H26	062183-79-3	25
5			N-2-Ethylhexanoylimidazole	194	C11H18N2O	062615-83-2	17



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
Data File : BF097398.D
Acq On : 5 Aug 2017 15:41
Operator : SJ/JU
Sample : I4592-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
15-50-RT3460-RT4215-RT3451

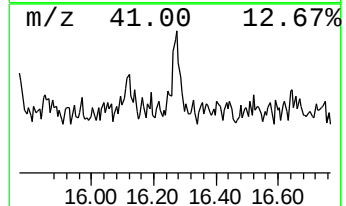
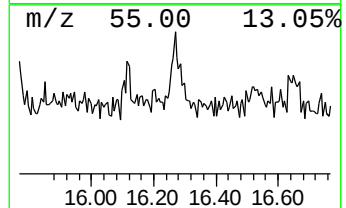
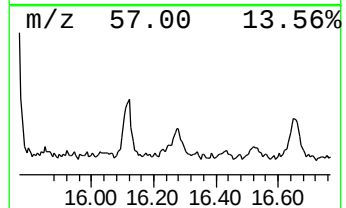
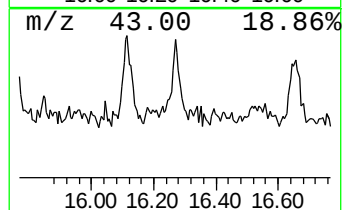
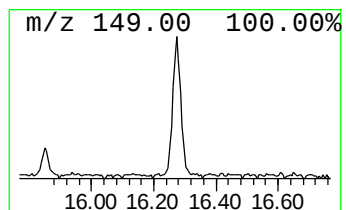
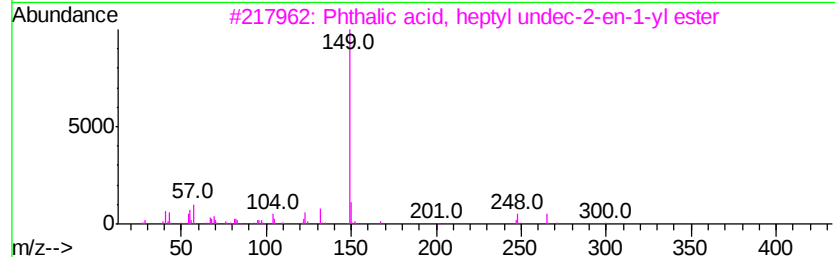
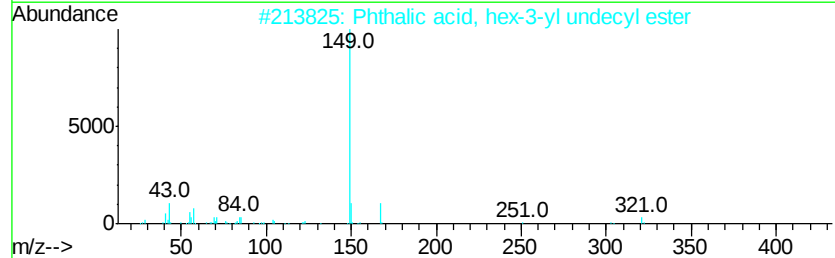
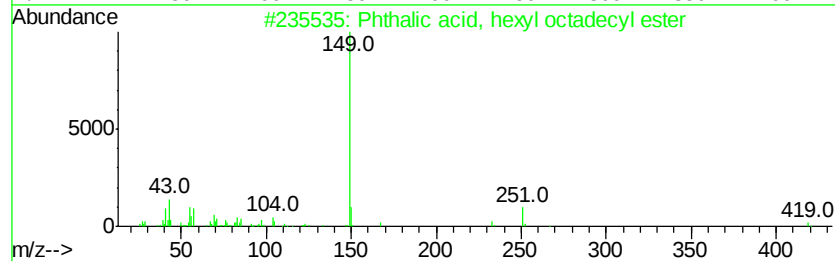
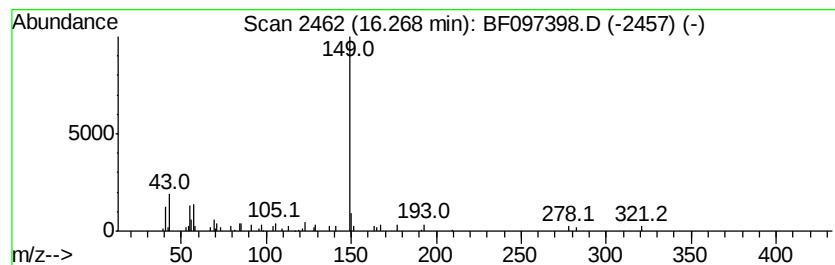
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phthalic acid, hexyl octade... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.27	2.27 ng	67833	Perylene-d12	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, hexyl octadecyl e...	502	C32H54O4	1000309-06-2	72
2		Phthalic acid, hex-3-yl undecyl ...	404	C25H40O4	1000356-96-2	72
3		Phthalic acid, heptyl undec-2-en...	416	C26H40O4	1000315-41-8	64
4		Phthalic acid, ethyl pentadecyl ...	404	C25H40O4	1000308-93-3	64
5		Phthalic acid, heptyl 4-methylpe...	348	C21H32O4	1000356-87-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080517\
 Data File : BF097398.D
 Acq On : 5 Aug 2017 15:41
 Operator : SJ/JU
 Sample : I4592-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 15-50-RT3460-RT4215-RT3451

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072517.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...	4.55	18.9	ng	714006	1	6.42	755117	20.0
unknown6.20	6.20	70.3	ng	2654890	1	6.42	755117	20.0
unknown7.90	7.90	25.4	ng	1397080	2	7.70	1098960	20.0
unknown8.66	8.66	2.6	ng	137341	3	9.45	1044900	20.0
Heptacosane	10.37	2.8	ng	148147	4	10.92	1059370	20.0
n-Hexadecanoic acid	11.48	6.3	ng	333116	4	10.92	1059370	20.0
unknown11.67	11.67	5.8	ng	306999	4	10.92	1059370	20.0
Octadecanoic acid	12.26	7.1	ng	236315	5	13.54	664133	20.0
Carbonic acid, he...	13.42	3.5	ng	115023	5	13.54	664133	20.0
unknown13.92	13.92	2.2	ng	71887	5	13.54	664133	20.0
unknown15.76	15.76	4.5	ng	132637	6	14.88	596711	20.0
Phthalic acid, he...	16.27	2.3	ng	67833	6	14.88	596711	20.0