

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
 Data File : BF084861.D
 Acq On : 5 Feb 2016 15:56
 Operator : SJ/IZ
 Sample : H1347-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM-196-242-COMPOSITE

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-BF020116.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.201	8	10	20	rVB	38724	71957	1.76%	0.163%
2	2.910	69	72	81	rBV2	12822	76667	1.87%	0.174%
3	4.807	236	238	250	rVB	826673	1191889	29.07%	2.705%
4	5.253	275	277	280	rBV	3849427	3842699	93.74%	8.723%
5	6.236	362	363	367	rBV	130143	144811	3.53%	0.329%
6	6.316	367	370	372	rBV	3627659	4035876	98.45%	9.161%
7	6.430	377	380	382	rVV	4408849	4066361	99.19%	9.230%
8	6.659	398	400	402	rBV	1350150	1113104	27.15%	2.527%
9	6.819	411	414	416	rBV	3069249	3124615	76.22%	7.093%
10	7.230	447	450	452	rBV	2550616	2506871	61.15%	5.690%
11	7.356	455	461	463	rBV	75821	113062	2.76%	0.257%
12	7.470	466	471	472	rBV3	32017	51945	1.27%	0.118%
13	7.710	488	492	495	rBV3	69503	138835	3.39%	0.315%
14	7.870	503	506	507	rBV	52681	59953	1.46%	0.136%
15	7.950	510	513	515	rVV	1449091	1607897	39.22%	3.650%
16	8.019	518	519	523	rVV	28945	47573	1.16%	0.108%
17	8.247	535	539	543	rBV4	33033	96143	2.35%	0.218%
18	8.304	543	544	546	rVV	66382	91483	2.23%	0.208%
19	8.384	549	551	555	rVB	116235	160100	3.91%	0.363%
20	8.556	564	566	568	rVB	187719	192955	4.71%	0.438%
21	8.647	572	574	576	rVB3	40366	52942	1.29%	0.120%
22	8.876	588	594	596	rBV5	65656	214267	5.23%	0.486%
23	8.922	596	598	599	rVV2	52551	80674	1.97%	0.183%
24	8.979	602	603	604	rVV	120039	98984	2.41%	0.225%
25	9.025	604	607	610	rVV	4035140	3987756	97.28%	9.052%
26	9.116	612	615	617	rVV	269157	287197	7.01%	0.652%
27	9.173	617	620	623	rVB4	52249	111138	2.71%	0.252%
28	9.367	635	637	640	rBV3	36332	73826	1.80%	0.168%
29	9.425	640	642	644	rBV	1512645	1345402	32.82%	3.054%
30	9.527	649	651	653	rVB3	39209	62462	1.52%	0.142%
31	9.642	659	661	663	rVB	255088	220519	5.38%	0.501%
32	9.699	663	666	668	rVB	1396008	1592624	38.85%	3.615%
33	9.859	678	680	683	rBV2	85243	149030	3.64%	0.338%
34	9.927	683	686	688	rVV	126894	187352	4.57%	0.425%

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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	9.962	688	689	691 rVB	59138	52264	1.27%	0.119%
36	10.076	697	699	701 rBV	142310	134572	3.28%	0.305%
37	10.145	701	705	706 rVB	159130	178193	4.35%	0.404%
38	10.362	721	724	727 rVB	66120	88729	2.16%	0.201%
39	10.488	732	735	737 rVV2	2170981	2945532	71.85%	6.686%
40	10.613	744	746	748 rBV	194818	232514	5.67%	0.528%
41	10.648	748	749	753 rVB2	41493	79041	1.93%	0.179%
42	10.705	753	754	758 rBV3	25543	45768	1.12%	0.104%
43	10.808	761	763	765 rBV2	31080	50498	1.23%	0.115%
44	11.059	782	785	786 rBV	156203	139480	3.40%	0.317%
45	11.093	786	788	792 rVV2	37434	66356	1.62%	0.151%
46	11.173	792	795	797 rVV	1812754	1591068	38.81%	3.612%
47	11.230	797	800	805 rVB4	22351	54212	1.32%	0.123%
48	11.413	813	816	820 rBV3	23647	56709	1.38%	0.129%
49	11.482	820	822	824 rBV	105890	91490	2.23%	0.208%
50	11.722	841	843	844 rBV	41529	42817	1.04%	0.097%
51	11.893	854	858	859 rBV	55518	67224	1.64%	0.153%
52	11.973	862	865	870 rBV	85448	140728	3.43%	0.319%
53	12.762	931	934	936 rBV	4332125	4099434	100.00%	9.305%
54	13.231	973	975	978 rBV	50062	62969	1.54%	0.143%
55	13.414	989	991	992 rBV	35013	48322	1.18%	0.110%
56	13.665	1011	1013	1016 rBV2	106837	172309	4.20%	0.391%
57	13.722	1016	1018	1020 rVB2	40008	51421	1.25%	0.117%
58	13.802	1022	1025	1027 rVB	1531417	1404592	34.26%	3.188%
59	14.316	1066	1070	1074 rBV3	25432	48254	1.18%	0.110%
60	15.162	1142	1144	1147 rBV	607089	910978	22.22%	2.068%

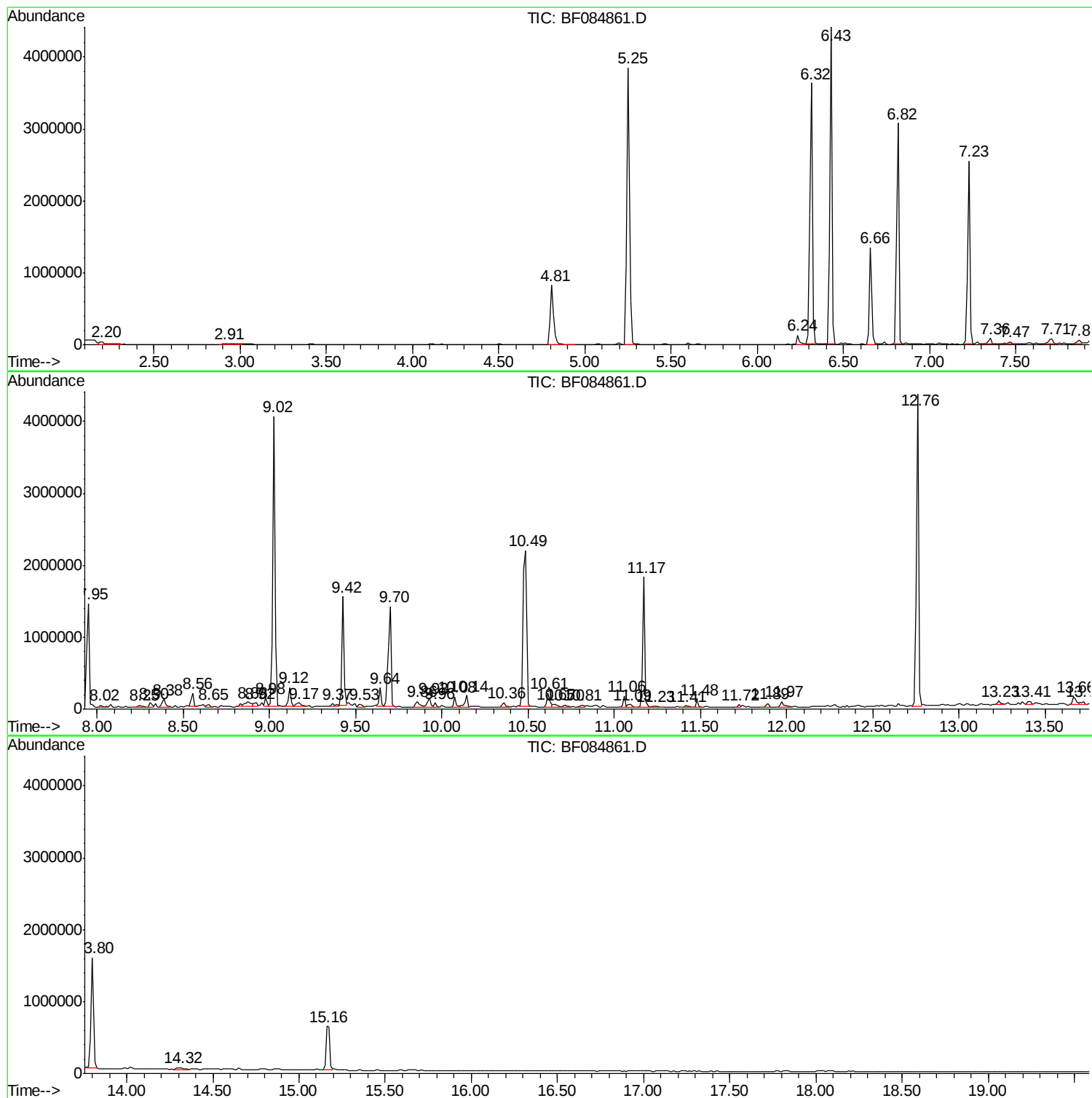
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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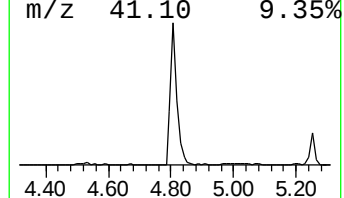
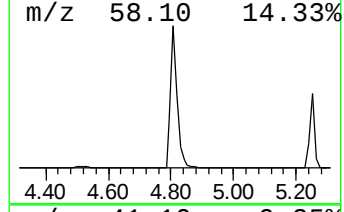
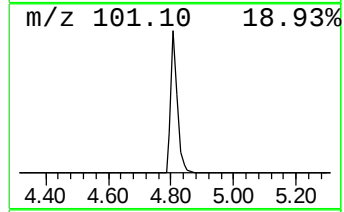
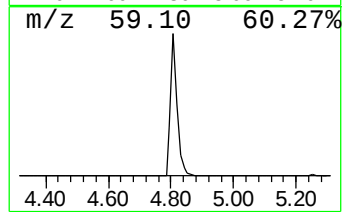
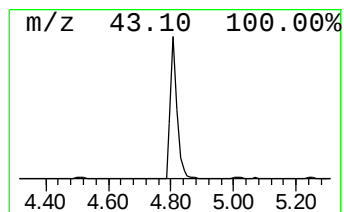
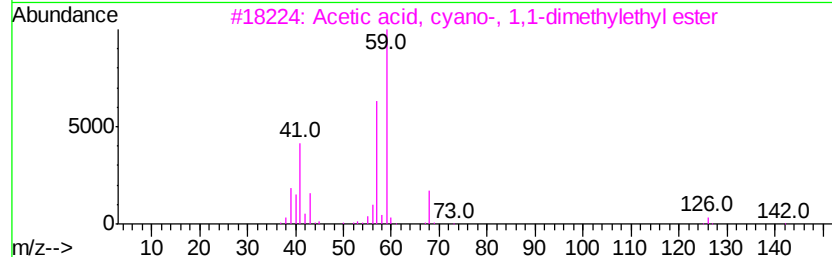
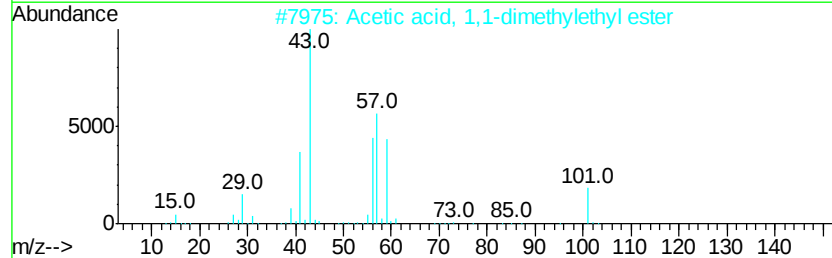
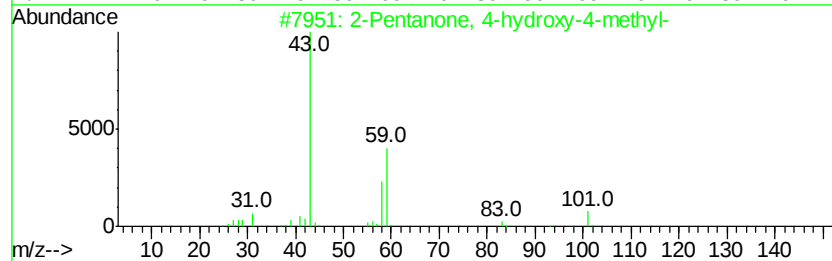
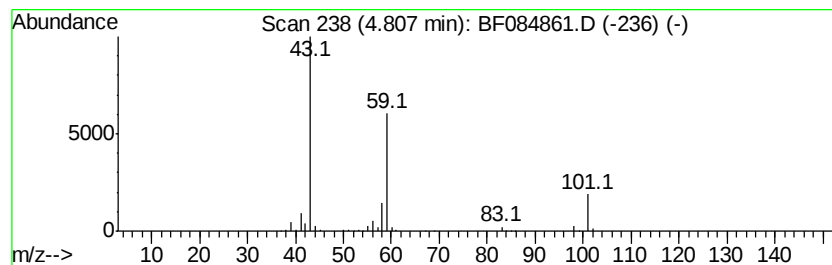
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.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.
4.81	21.42 ng	1191890	1,4-Dichlorobenzene-d4	6.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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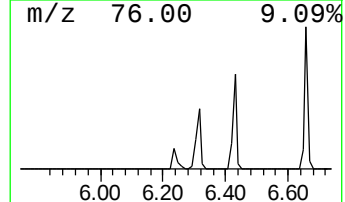
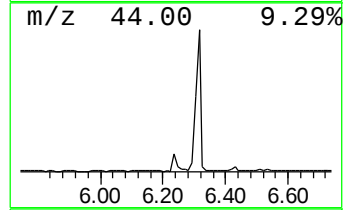
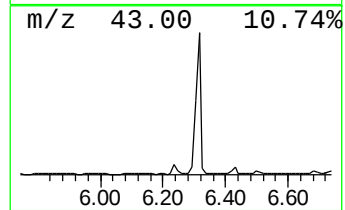
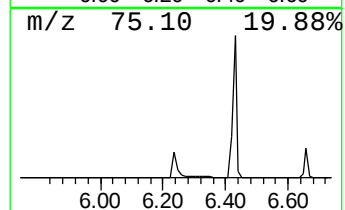
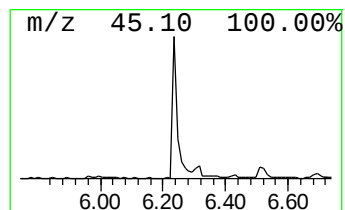
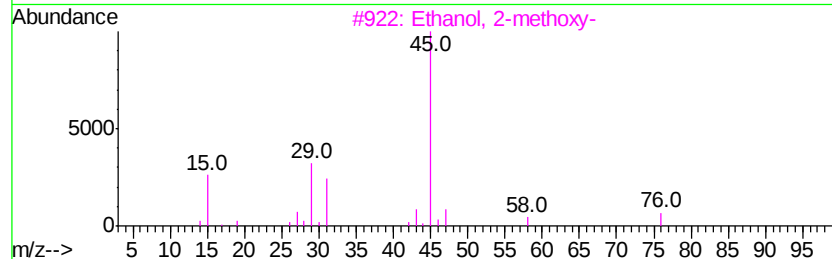
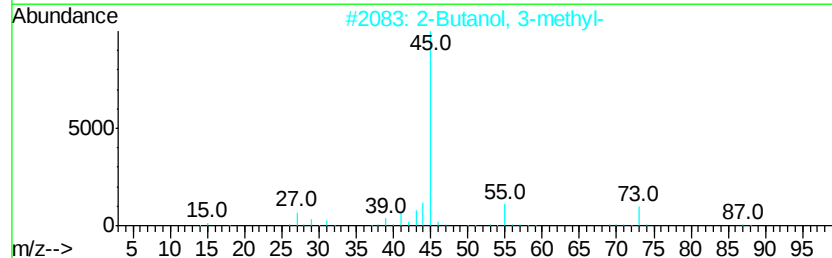
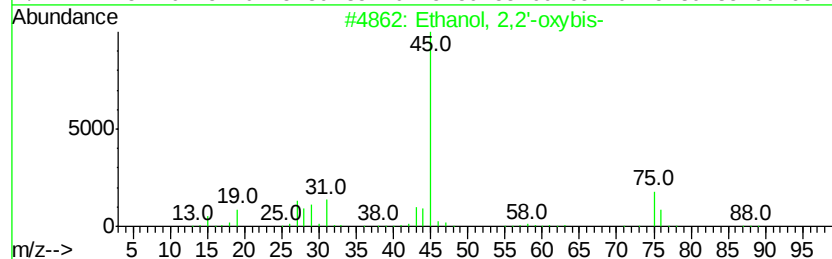
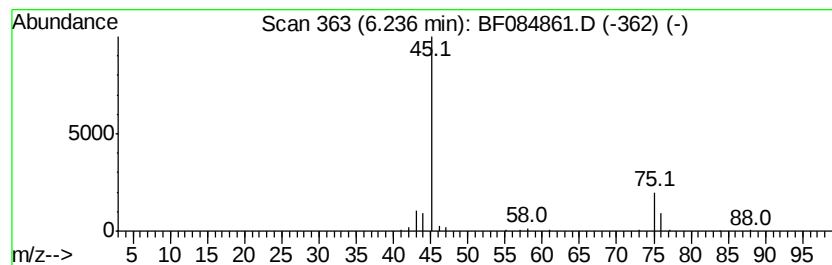
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2,2'-oxybis- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.24	2.60 ng	144811	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2,2'-oxybis-	106	C4H10O3	000111-46-6	83
2		2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	42
3		Ethanol, 2-methoxy-	76	C3H8O2	000109-86-4	9
4		2-Hydroxyethylhydrazine	76	C2H8N2O	000109-84-2	9
5		2-Propanol, 1,3-dimethoxy-	120	C5H12O3	000623-69-8	9



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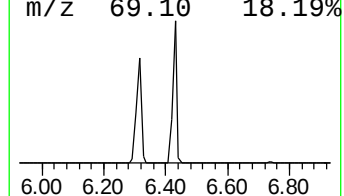
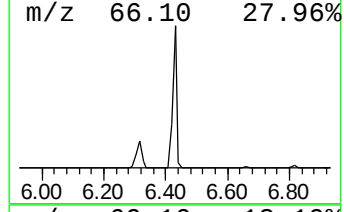
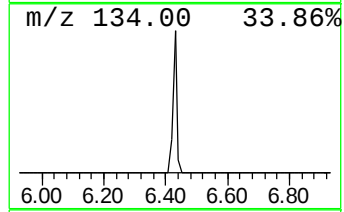
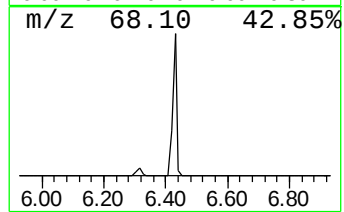
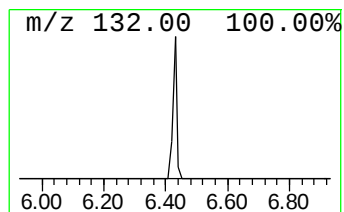
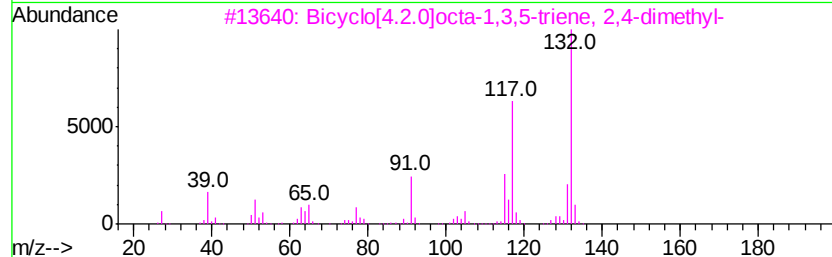
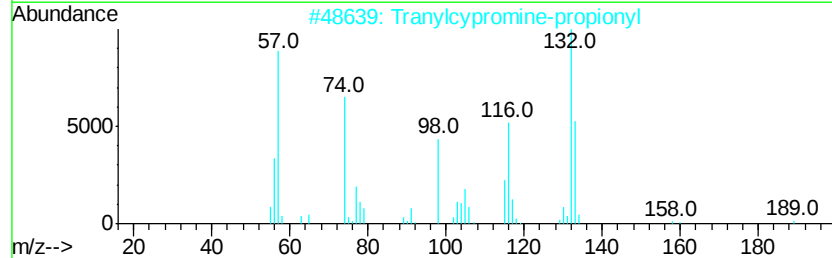
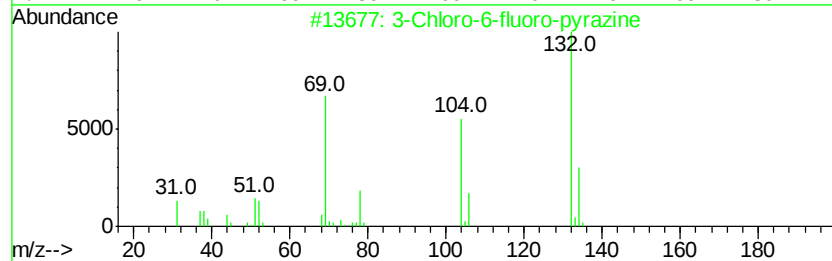
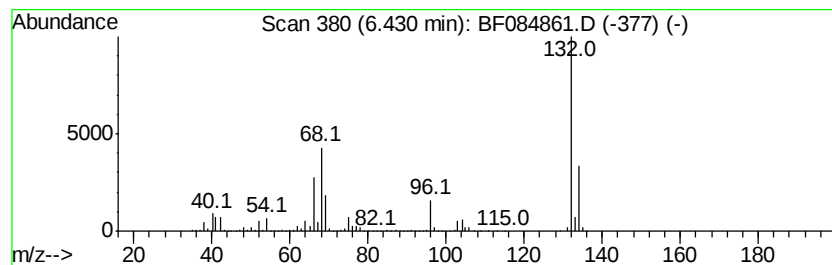
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.43 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	73.06 ng	4066360	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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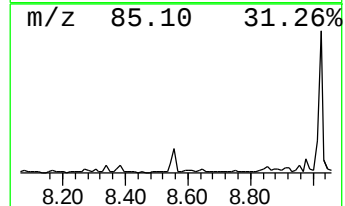
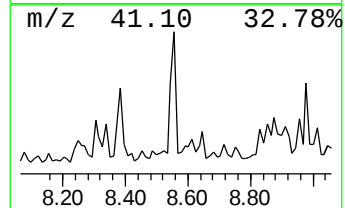
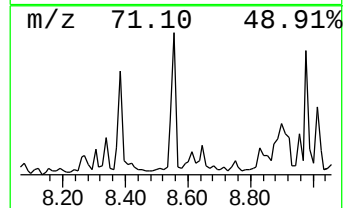
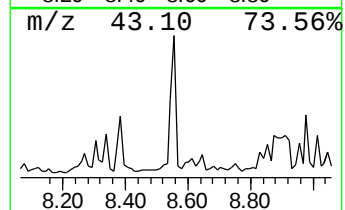
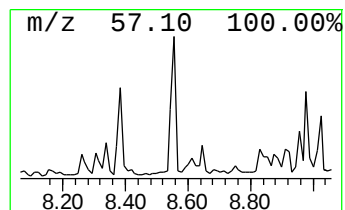
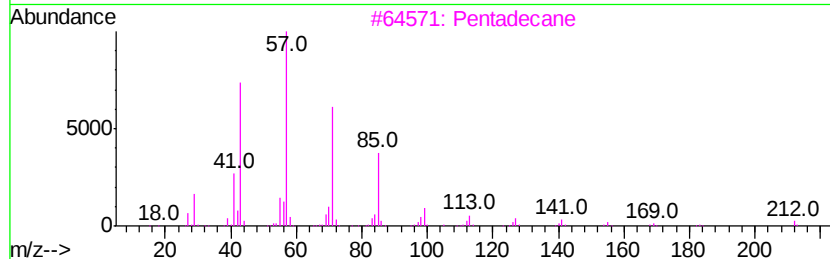
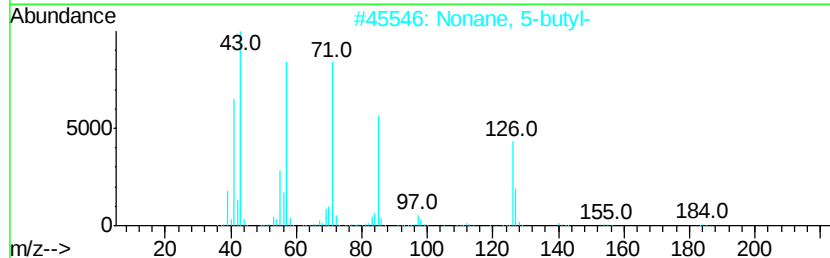
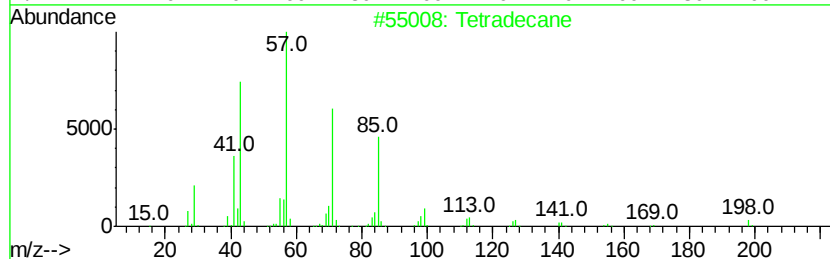
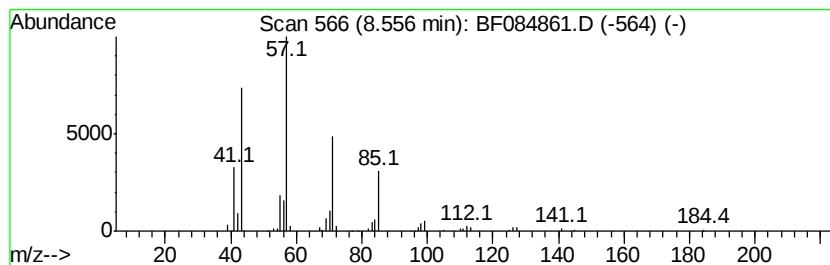
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Tetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	2.40 ng	192955	Napthalene-d8	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	90
2		Nonane, 5-butyl-	184	C13H28	017312-63-9	86
3		Pentadecane	212	C15H32	000629-62-9	86
4		Tridecane	184	C13H28	000629-50-5	86
5		Hexadecane	226	C16H34	000544-76-3	86



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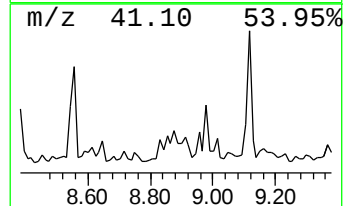
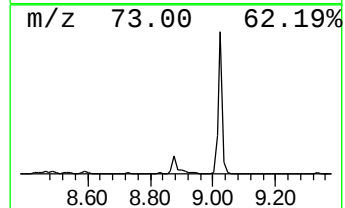
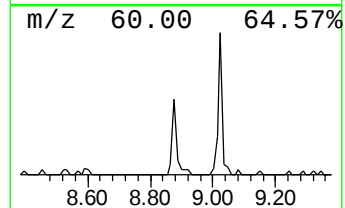
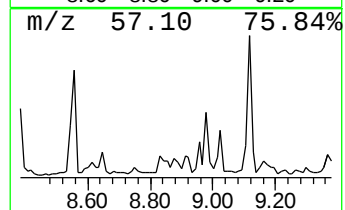
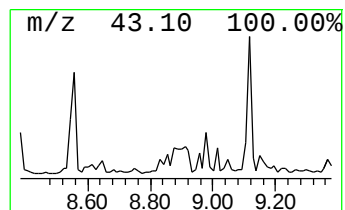
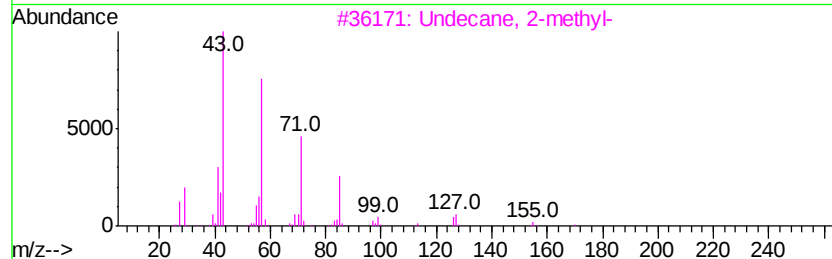
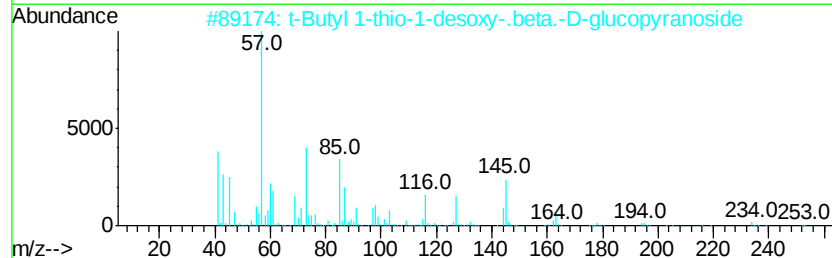
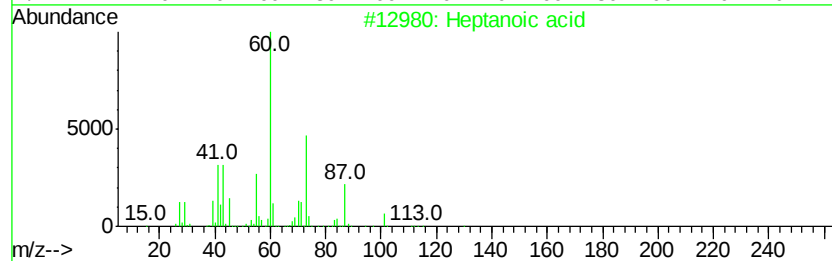
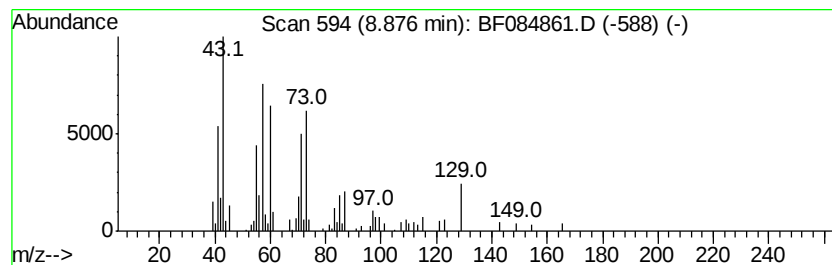
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown8.88 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.88	2.69 ng	214267	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptanoic acid	130	C7H14O2	000111-14-8	35
2			t-Butyl 1-thio-1-desoxy-.beta.-D...	252	C10H20O5S	1000126-98-1	35
3			Undecane, 2-methyl-	170	C12H26	007045-71-8	27
4			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	27
5			Hexadecane	226	C16H34	000544-76-3	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
Data File : BF084861.D
Acq On : 5 Feb 2016 15:56
Operator : SJ/IZ
Sample : H1347-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM-196-242-COMPOSITE

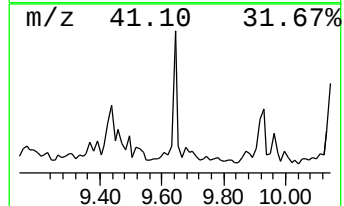
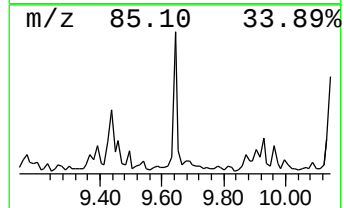
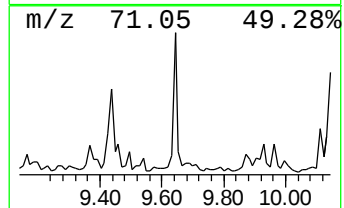
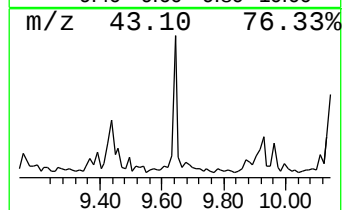
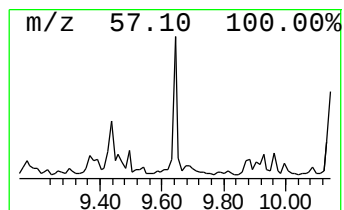
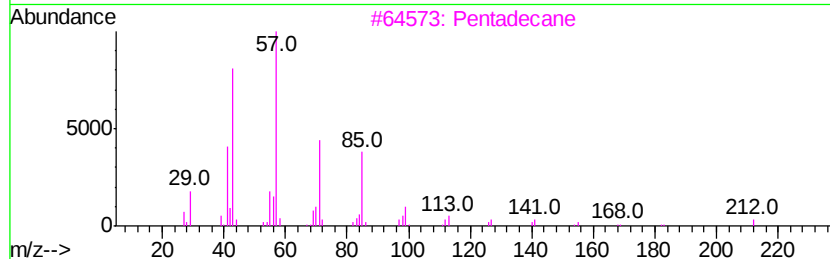
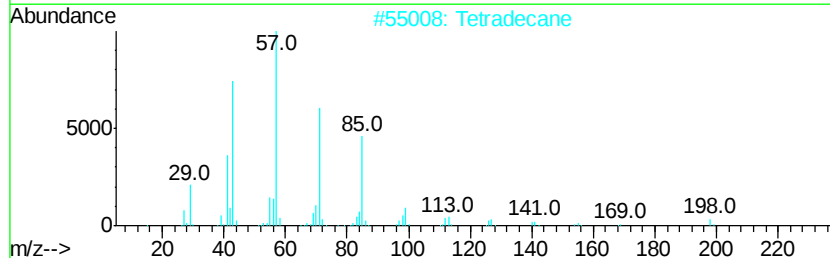
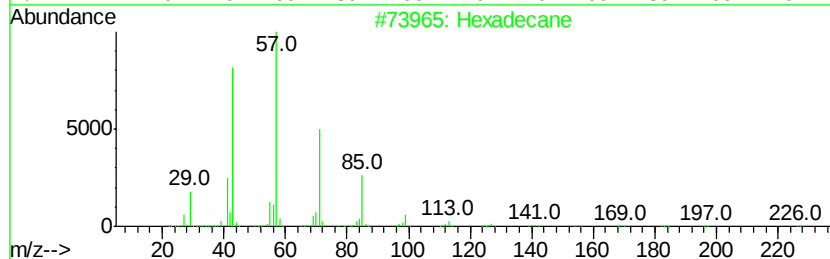
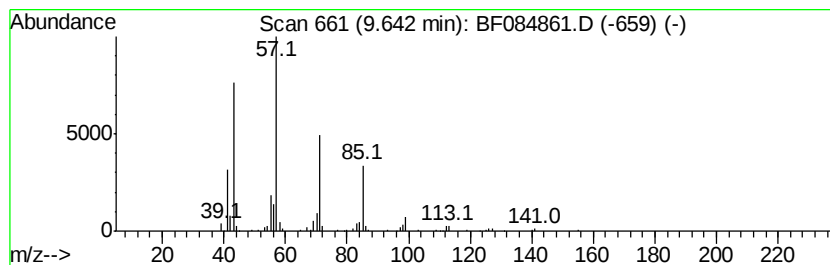
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.64	2.77 ng	220519	Acenaphthene-d10	9.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetradecane		198	C14H30	000629-59-4	90
3	Pentadecane		212	C15H32	000629-62-9	80
4	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	59
5	Heptane, 3-ethyl-5-methyl-		142	C10H22	052896-90-9	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
Data File : BF084861.D
Acq On : 5 Feb 2016 15:56
Operator : SJ/IZ
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DRUM-196-242-COMPOSITE

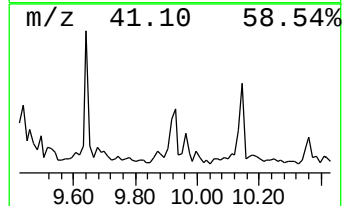
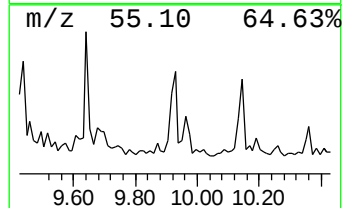
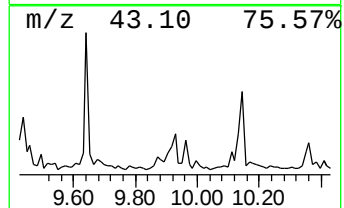
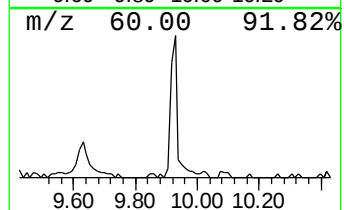
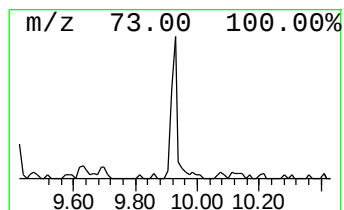
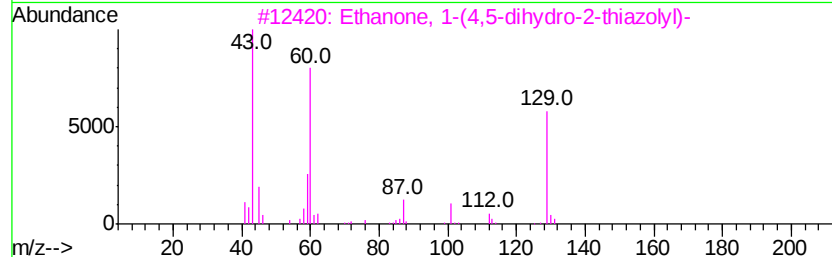
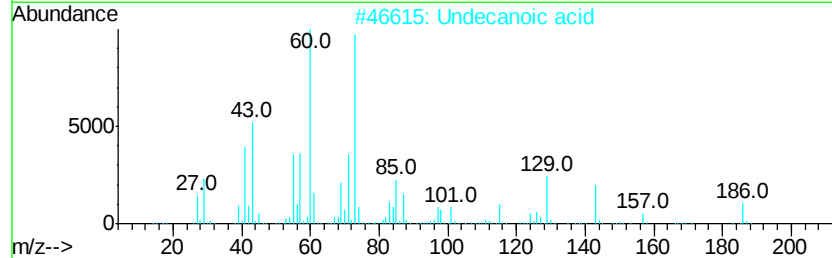
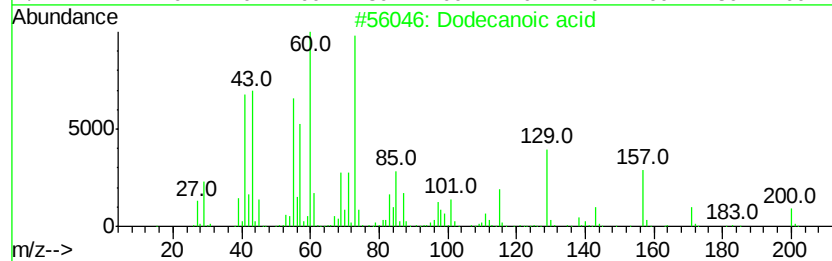
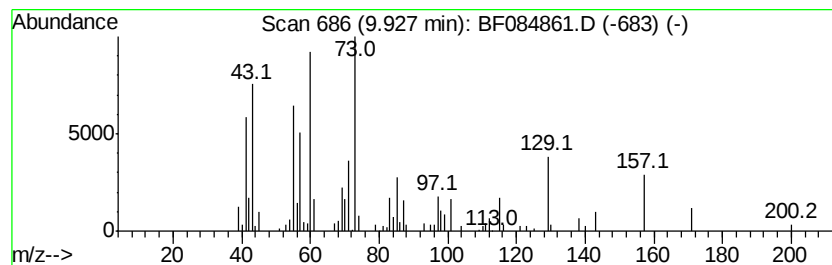
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Dodecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.93	2.35 ng	187352	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanoic acid	200	C12H24O2	000143-07-7	93
2		Undecanoic acid	186	C11H22O2	000112-37-8	53
3		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	43
4		Octanoic Acid	144	C8H16O2	000124-07-2	43
5		n-Decanoic acid	172	C10H20O2	000334-48-5	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
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Operator : SJ/IZ
Sample : H1347-01
Misc :
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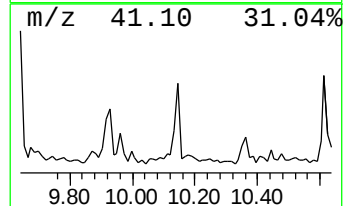
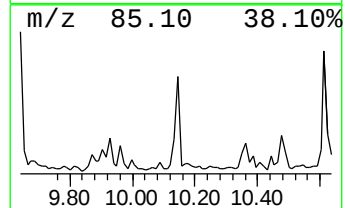
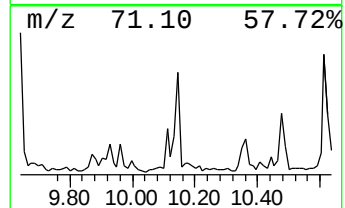
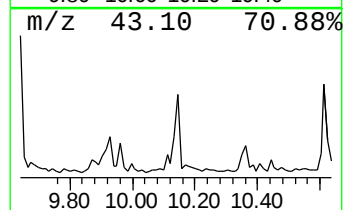
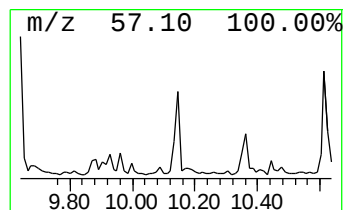
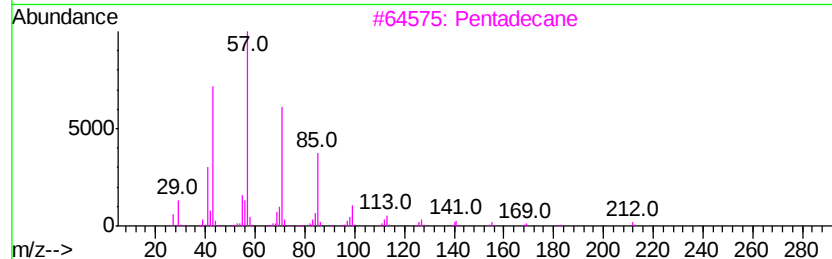
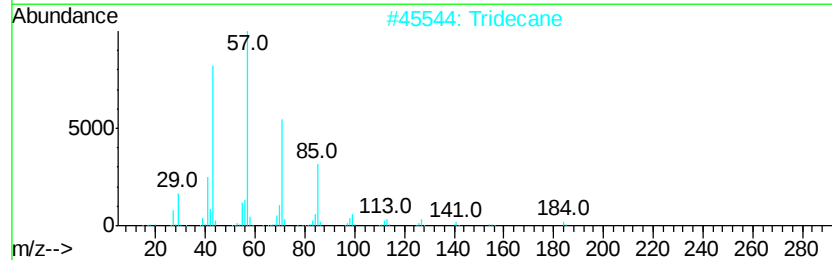
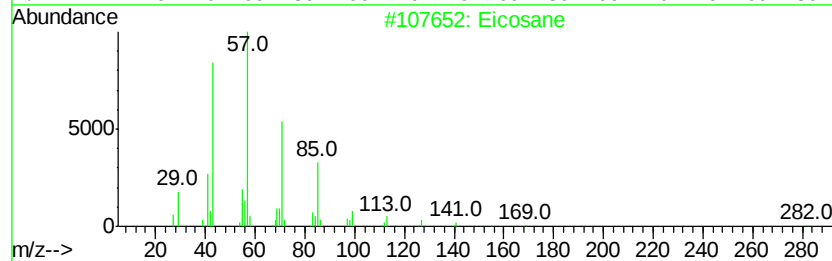
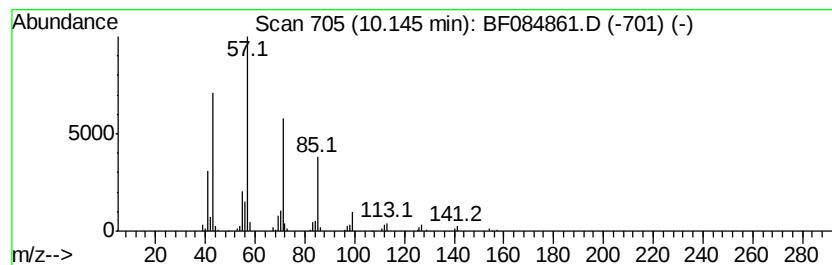
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.24 ng	178193	Acenaphthene-d10	9.70
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 91
2	Tridecane		184 C13H28	000629-50-5 90
3	Pentadecane		212 C15H32	000629-62-9 90
4	Hexadecane		226 C16H34	000544-76-3 83
5	Decane, 2,3,5-trimethyl-		184 C13H28	062238-11-3 78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
Data File : BF084861.D
Acq On : 5 Feb 2016 15:56
Operator : SJ/IZ
Sample : H1347-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DRUM-196-242-COMPOSITE

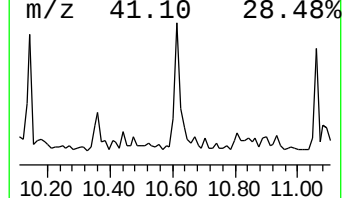
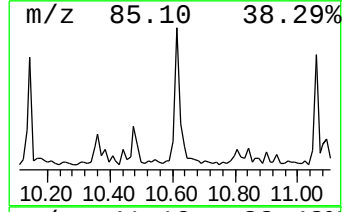
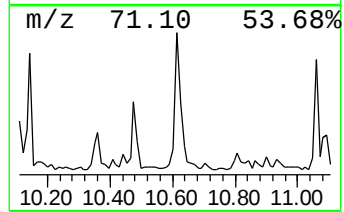
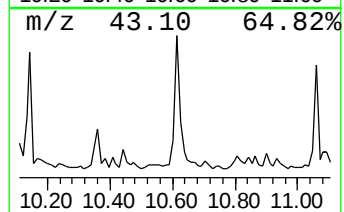
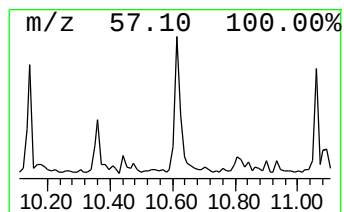
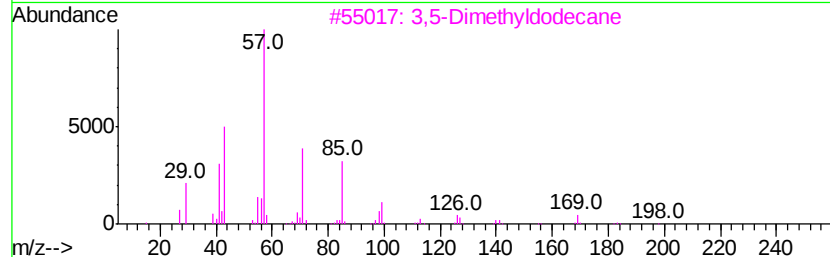
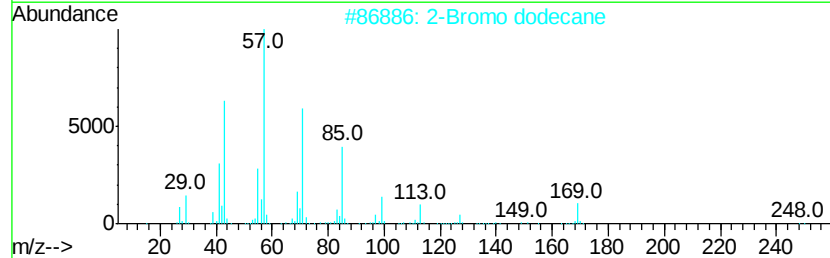
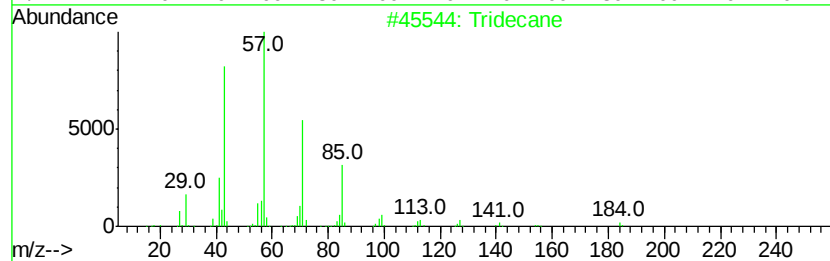
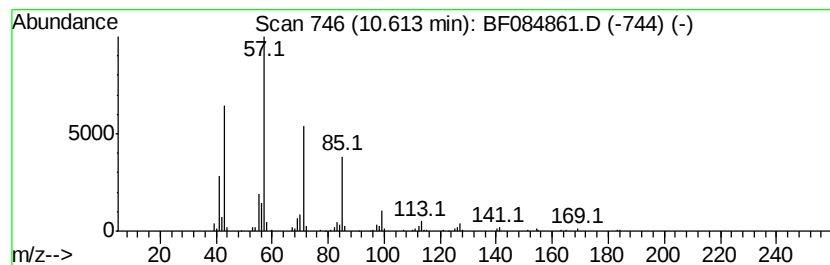
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Tridecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.61	2.92 ng	232514	Phenanthrene-d10	11.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	90
2		2-Bromo dodecane	248	C12H25Br	013187-99-0	90
3		3,5-Dimethyldodecane	198	C14H30	107770-99-0	83
4		Hexadecane	226	C16H34	000544-76-3	78
5		Heptacosane	380	C27H56	000593-49-7	68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
Data File : BF084861.D
Acq On : 5 Feb 2016 15:56
Operator : SJ/IZ
Sample : H1347-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DRUM-196-242-COMPOSITE

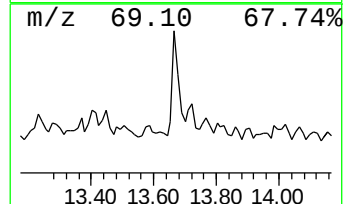
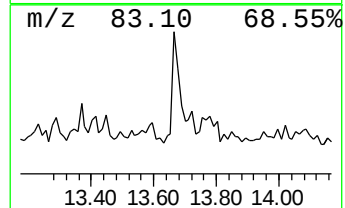
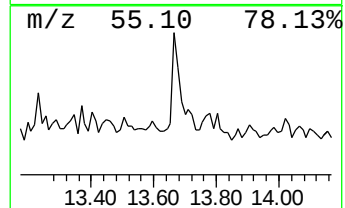
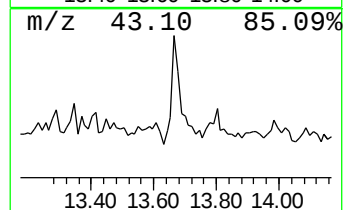
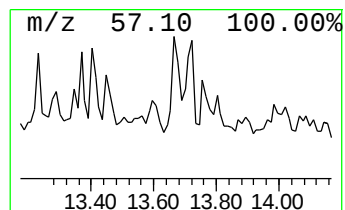
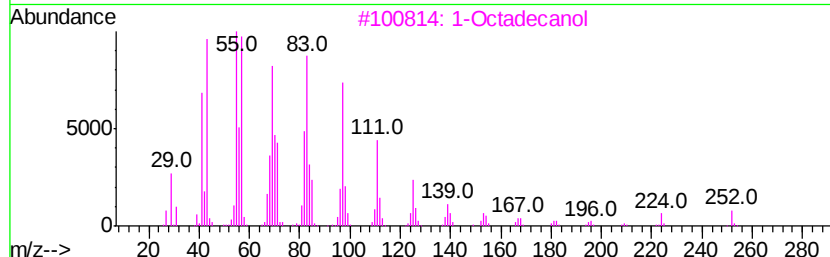
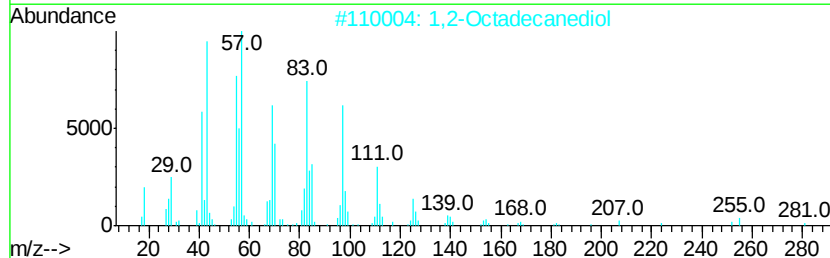
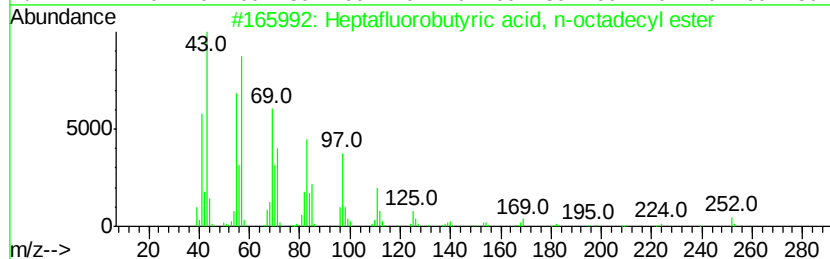
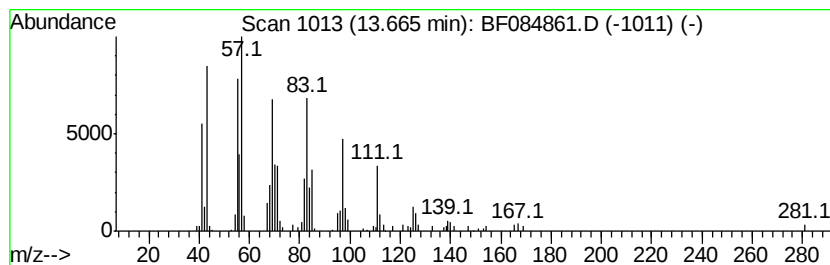
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptafluorobutyric acid, n-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	2.45 ng	172309	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91
2		1,2-Octadecanediol	286	C18H38O2	020294-76-2	91
3		1-Octadecanol	270	C18H38O	000112-92-5	91
4		Pentafluoropropionic acid, heptafluorobutyl ester	402	C20H35F5O2	1000283-04-2	91
5		Trifluoroacetic acid, n-octadecyl ester	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020516\
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Sample : H1347-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.81	21.4	ng	1191890	1	6.66	1113100	20.0
Ethanol, 2,2'-oxy...	6.24	2.6	ng	144811	1	6.66	1113100	20.0
unknown6.43	6.43	73.1	ng	4066360	1	6.66	1113100	20.0
Tetradecane	8.56	2.4	ng	192955	2	7.95	1607900	20.0
unknown8.88	8.88	2.7	ng	214267	3	9.70	1592620	20.0
Hexadecane	9.64	2.8	ng	220519	3	9.70	1592620	20.0
Dodecanoic acid	9.93	2.4	ng	187352	3	9.70	1592620	20.0
Eicosane	10.14	2.2	ng	178193	3	9.70	1592620	20.0
Tridecane	10.61	2.9	ng	232514	4	11.17	1591070	20.0
Heptafluorobutyri...	13.67	2.5	ng	172309	5	13.80	1404590	20.0