

Data Path : Z:\HPCHEM1\BNA F\DATA\BF030817\
 Data File : BF093439.D
 Acq On : 8 Mar 2017 20:12
 Operator : SJ/MA
 Sample : I2137-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 38D-9-11

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-BF030717.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	1.955	4	6	10	rVB	88074	120311	2.08%	0.248%
2	3.590	146	149	161	rBV	36230	93459	1.61%	0.192%
3	4.916	263	265	277	rBV	884554	1456506	25.16%	2.999%
4	5.361	302	304	326	rBV	3973207	5625077	97.16%	11.581%
5	5.761	335	339	347	rBV	61292	93818	1.62%	0.193%
6	6.413	393	396	404	rBV	4240756	5123890	88.51%	10.549%
7	6.527	404	406	409	rBV	3686610	4989077	86.18%	10.272%
8	6.756	424	426	429	rBV	736732	996131	17.21%	2.051%
9	6.916	438	440	443	rBV	3976227	3852453	66.54%	7.932%
10	7.339	474	477	479	rBV	3446564	3598201	62.15%	7.408%
11	8.047	537	539	542	rBV	1093084	1311274	22.65%	2.700%
12	9.133	630	634	636	rBV	4401810	5789295	100.00%	11.919%
13	9.522	666	668	671	rBV	817846	940409	16.24%	1.936%
14	9.739	684	687	690	rBV2	50755	128728	2.22%	0.265%
15	9.796	690	692	695	rVB	1138457	1441875	24.91%	2.969%
16	10.230	728	730	732	rVB	148678	116146	2.01%	0.239%
17	10.448	746	749	752	rBV	45051	76376	1.32%	0.157%
18	10.596	759	762	768	rBV	2821760	3007021	51.94%	6.191%
19	10.699	770	771	779	rVB	145895	206483	3.57%	0.425%
20	11.145	809	810	816	rVB	93711	156969	2.71%	0.323%
21	11.282	820	822	825	rBV2	1204030	1498363	25.88%	3.085%
22	12.871	958	961	964	rBV	4164159	4978712	86.00%	10.250%
23	13.762	1037	1039	1046	rBV	138532	250247	4.32%	0.515%
24	13.922	1052	1053	1056	rVB	1087182	1386388	23.95%	2.854%
25	14.745	1123	1125	1127	rBV	53777	57942	1.00%	0.119%
26	15.340	1174	1177	1186	rBV	911786	1276026	22.04%	2.627%

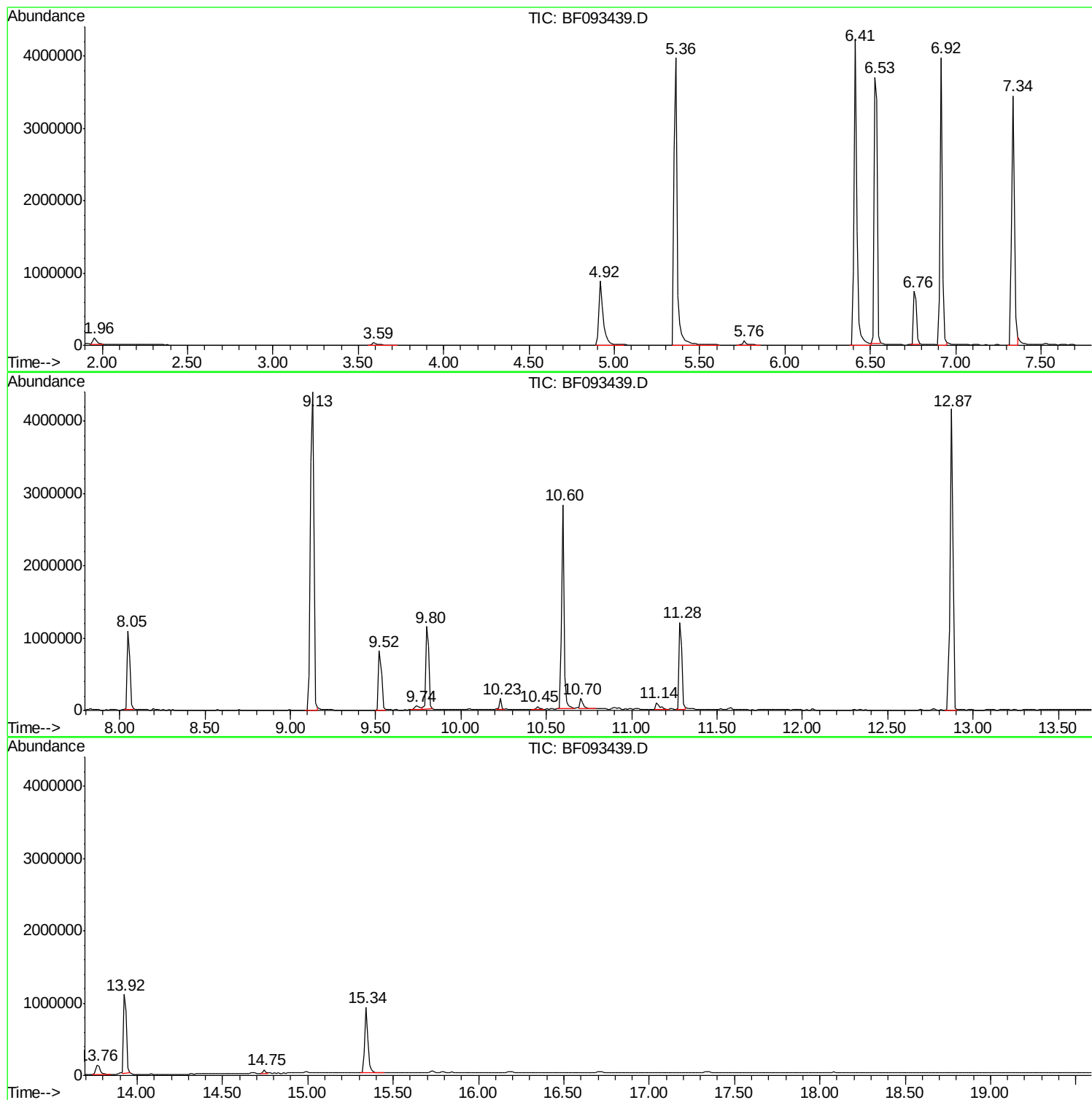
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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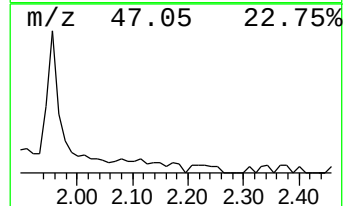
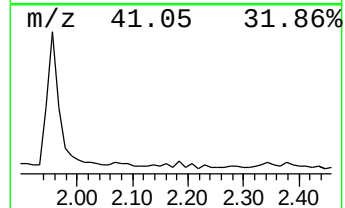
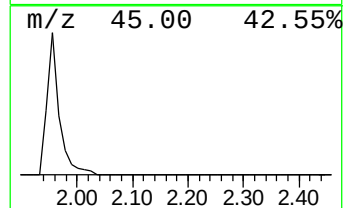
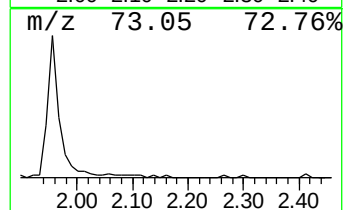
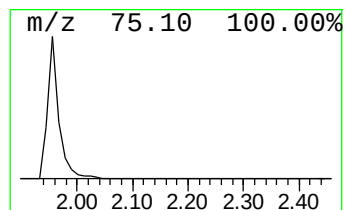
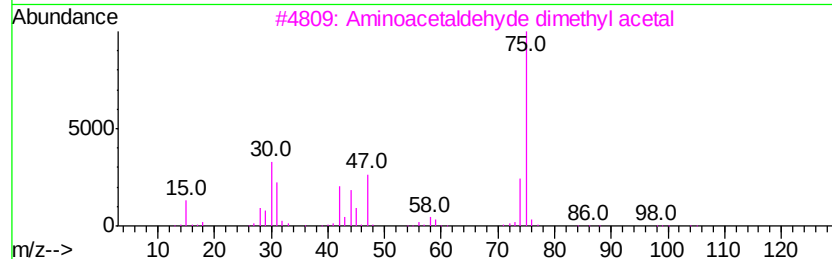
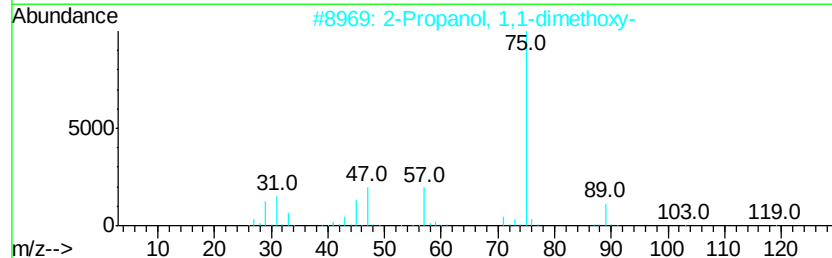
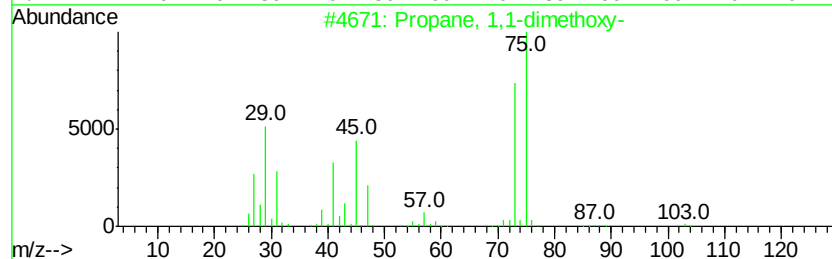
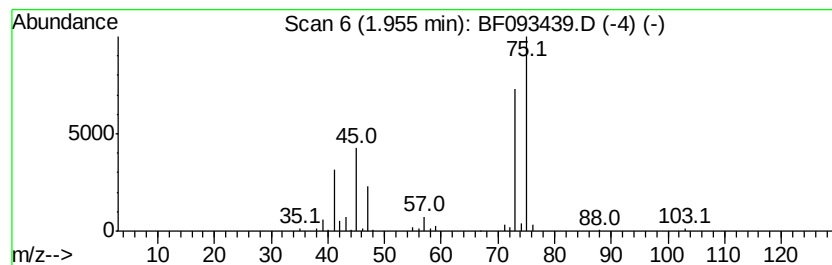
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	2.42 ng	120311	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2		2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	42
3		Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	33
4		Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	33
5		Methane, trimethoxy-	106	C4H10O3	000149-73-5	33



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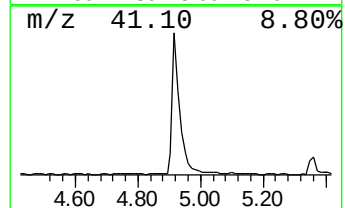
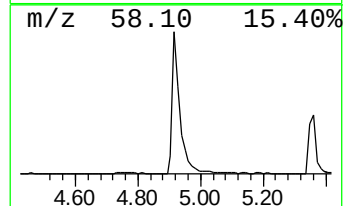
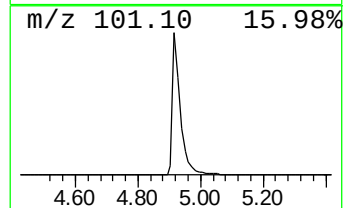
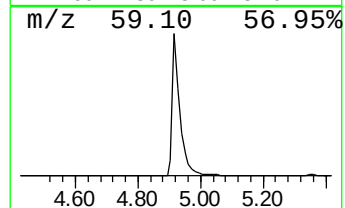
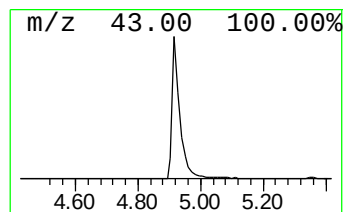
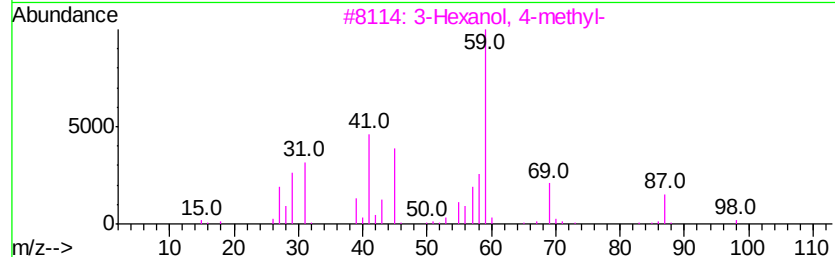
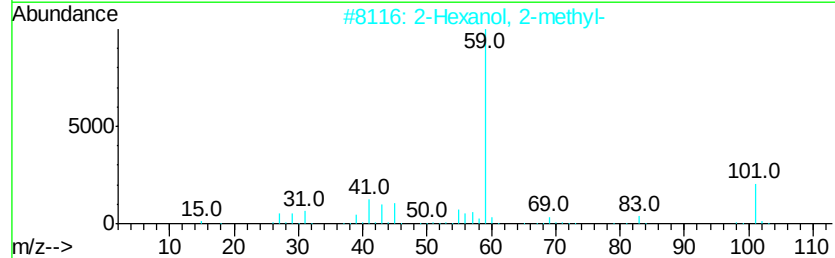
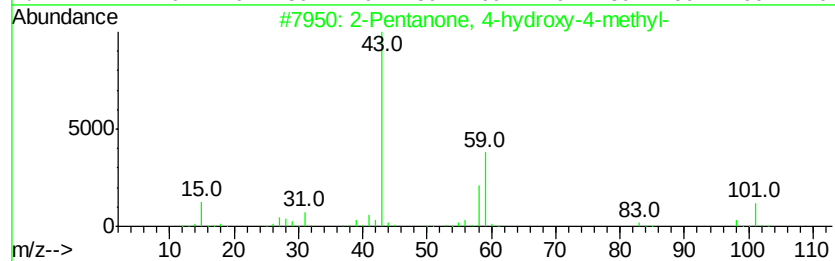
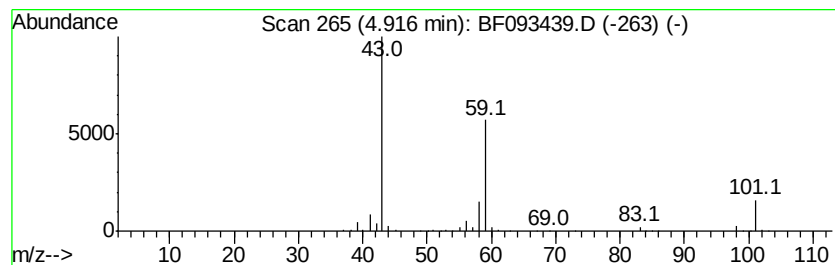
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	29.24 ng	1456510	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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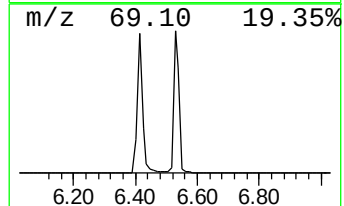
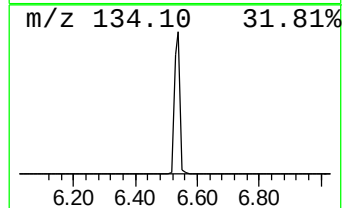
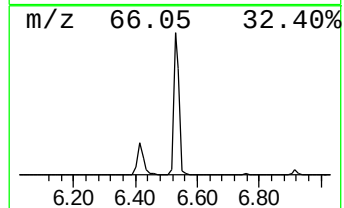
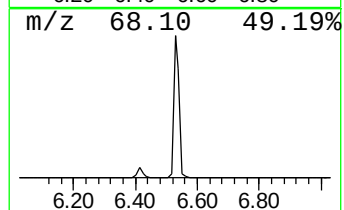
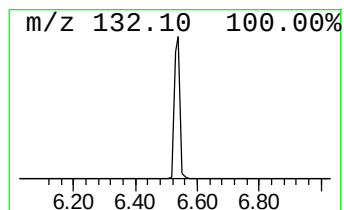
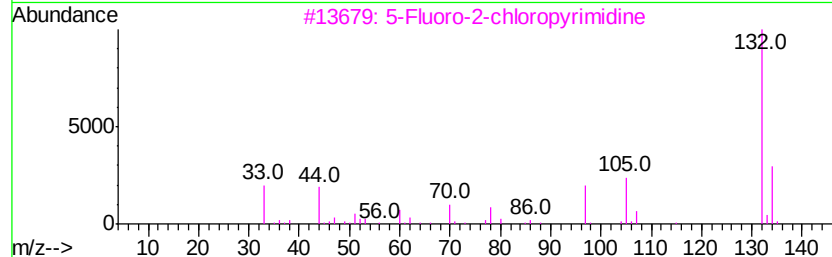
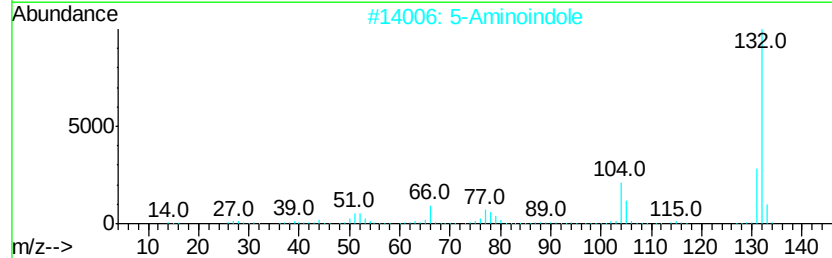
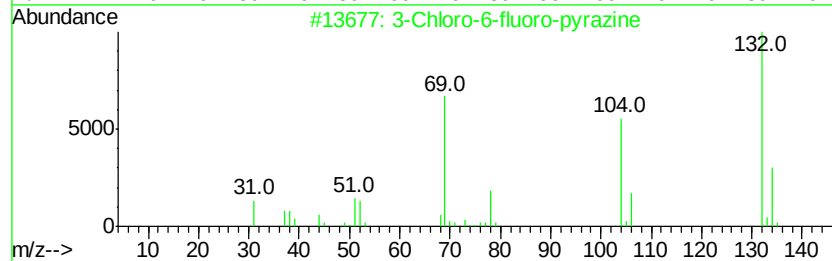
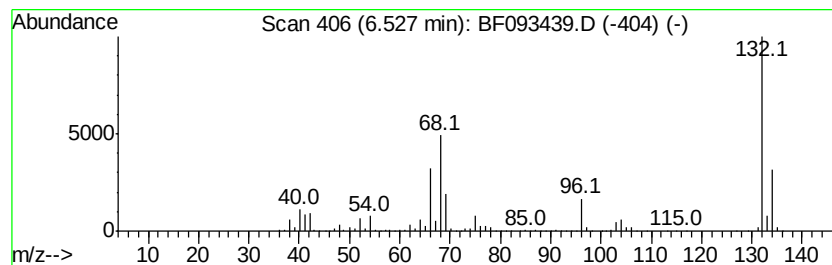
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	100.17 ng	4989080	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5	Aminoindole	132	C8H8N2	005192-03-0	27
3	5	Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranvlcypromine-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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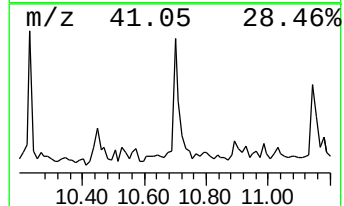
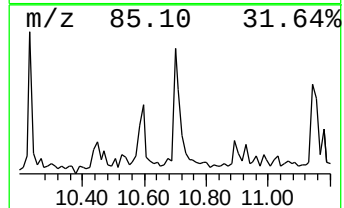
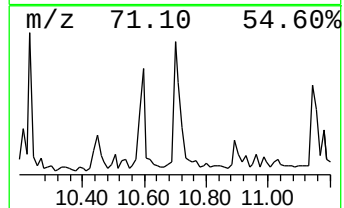
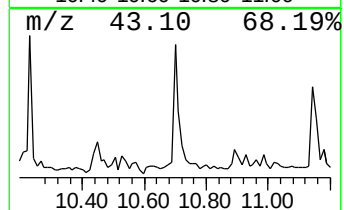
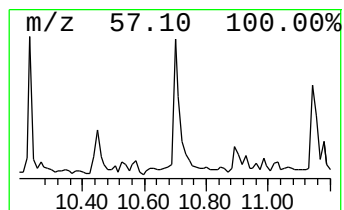
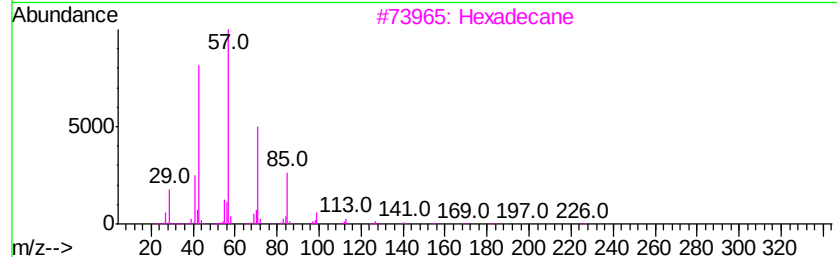
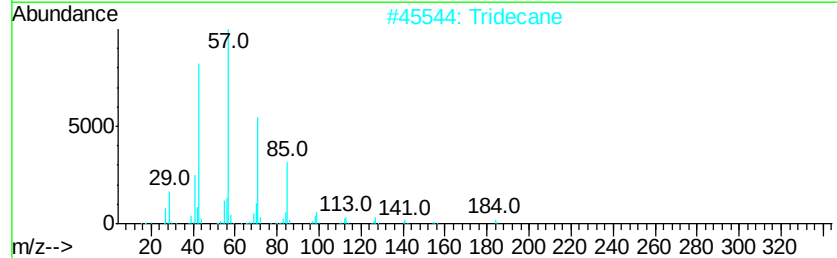
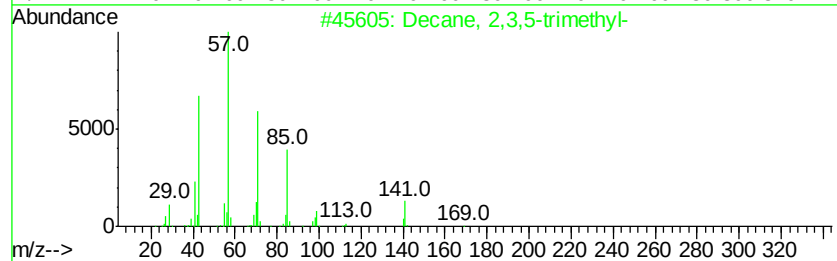
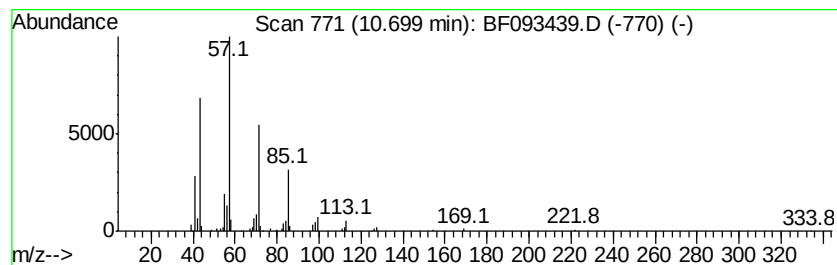
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Decane, 2,3,5-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	2.76 ng	206483	Phenanthrene-d10	11.28
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 2,3,5-trimethyl-	184 C13H28	062238-11-3	86
2	Tridecane	184 C13H28	000629-50-5	86
3	Hexadecane	226 C16H34	000544-76-3	86
4	Pentadecane, 7-methyl-	226 C16H34	006165-40-8	83
5	Pentadecane	212 C15H32	000629-62-9	78



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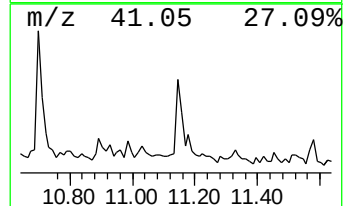
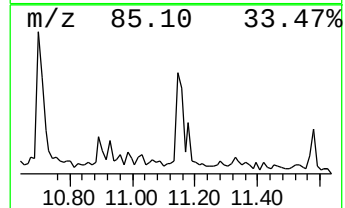
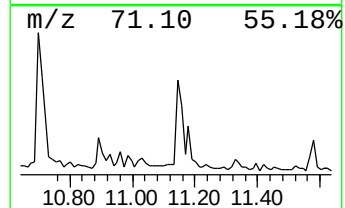
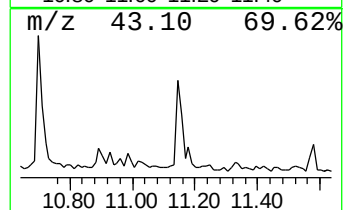
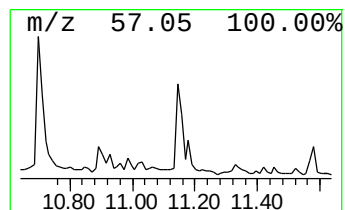
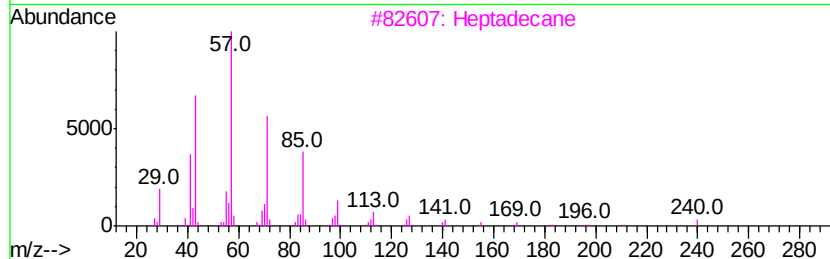
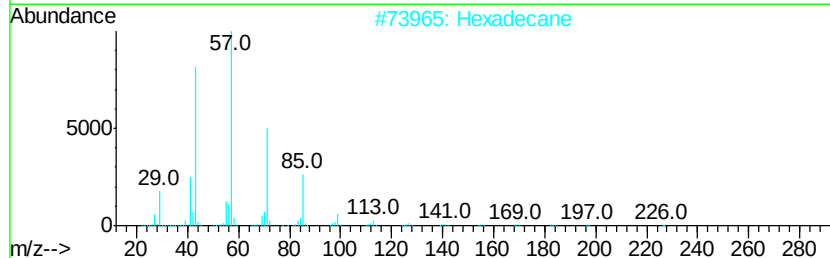
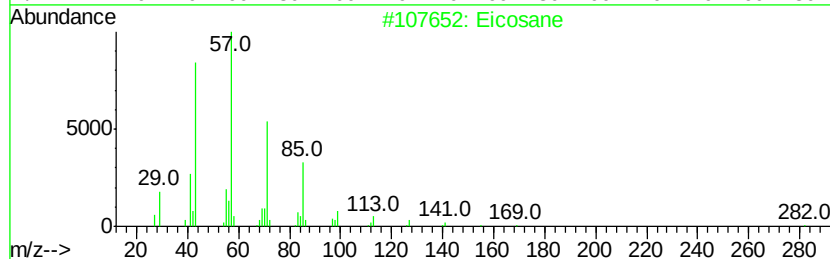
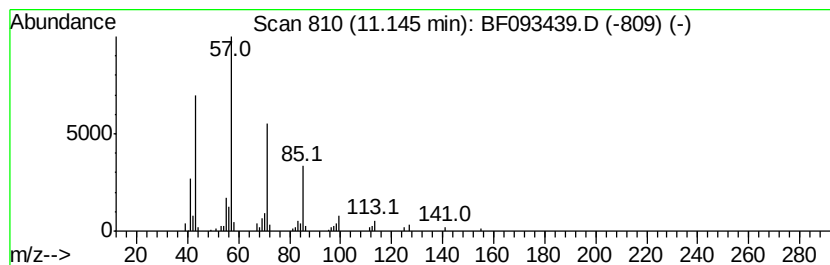
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.14	2.10 ng	156969	Phenanthrene-d10		11.28		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			Hexadecane	226	C16H34	000544-76-3	91
3			Heptadecane	240	C17H36	000629-78-7	90
4			Tridecane	184	C13H28	000629-50-5	90
5			Pentadecane	212	C15H32	000629-62-9	86



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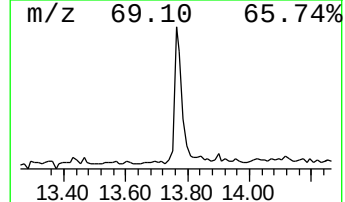
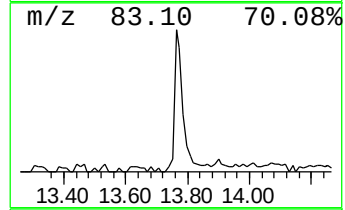
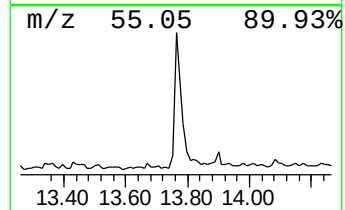
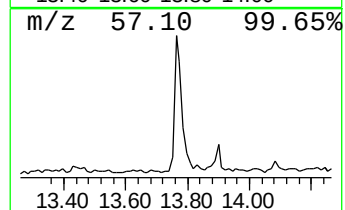
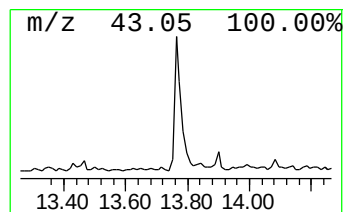
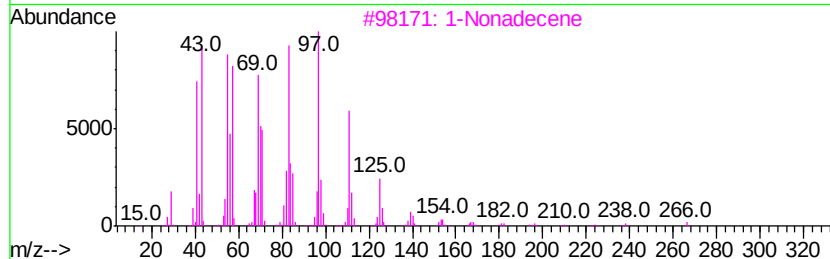
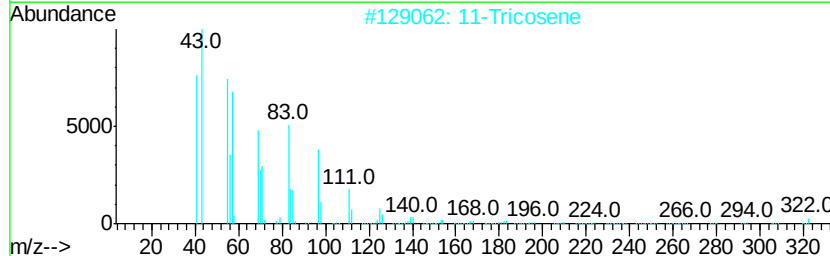
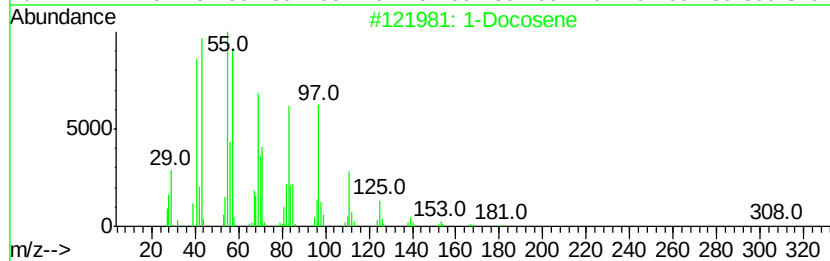
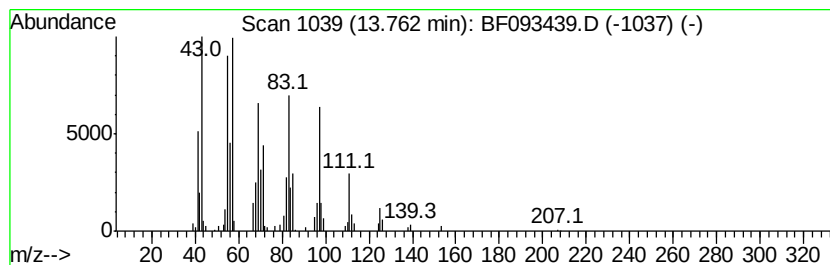
Instrument :
BNA_F
ClientSampleId :
38D-9-11

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	3.61 ng	250247	Chrysene-d12	13.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Docosene		308 C22H44	001599-67-3 94
2	11-Tricosene		322 C23H46	052078-56-5 91
3	1-Nonadecene		266 C19H38	018435-45-5 91
4	3-Eicosene, (E)-		280 C20H40	074685-33-9 91
5	Dichloroacetic acid, tetradecyl ...		324 C16H30Cl2O2	083005-02-1 90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030817\
Data File : BF093439.D
Acq On : 8 Mar 2017 20:12
Operator : SJ/MA
Sample : I2137-04
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
38D-9-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030717.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
Propane, 1,1-dime...	1.96	2.4	ng	120311	1	6.77	996131	20.0
2-Pentanone, 4-hy...	4.92	29.2	ng	1456510	1	6.77	996131	20.0
unknown6.53	6.53	100.2	ng	4989080	1	6.77	996131	20.0
Decane, 2,3,5-tri...	10.70	2.8	ng	206483	4	11.28	1498360	20.0
Eicosane	11.14	2.1	ng	156969	4	11.28	1498360	20.0
1-Docosene	13.76	3.6	ng	250247	5	13.92	1386390	20.0