

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
 Data File : BF078833.D
 Acq On : 30 Apr 2015 2:47
 Operator : TP/IZ
 Sample : G2003-10
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PZ-2(45-47)

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-BF042815.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	1.438	20	22	28	rVB	129773	207271	2.36%	0.552%
2	1.541	28	31	34	rVB	7397005	8797033	100.00%	23.410%
3	1.678	40	43	46	rVB	157235	260501	2.96%	0.693%
4	1.781	51	52	57	rBV	315395	466410	5.30%	1.241%
5	1.873	57	60	62	rVV	723381	1064318	12.10%	2.832%
6	1.918	62	64	92	rVB	2148591	3855890	43.83%	10.261%
7	5.153	345	347	349	rBV	92824	123031	1.40%	0.327%
8	5.187	349	350	355	rVB	383686	447013	5.08%	1.190%
9	5.679	390	393	396	rBV	2366856	2324732	26.43%	6.186%
10	6.924	500	502	505	rBV	2249949	2346716	26.68%	6.245%
11	7.096	514	517	519	rBV	1945151	2490027	28.31%	6.626%
12	7.382	540	542	545	rBV	515912	459945	5.23%	1.224%
13	7.565	556	558	561	rBV	1378421	1747607	19.87%	4.651%
14	8.067	600	602	605	rBV	1300113	1385306	15.75%	3.687%
15	8.959	677	680	682	rBV	729352	633316	7.20%	1.685%
16	10.308	795	798	800	rBV	3099880	2756855	31.34%	7.336%
17	10.811	839	842	844	rBV	117014	106712	1.21%	0.284%
18	11.142	868	871	873	rBV	564263	675505	7.68%	1.798%
19	12.034	947	949	952	rBV	189766	248898	2.83%	0.662%
20	12.114	954	956	959	rBV2	1594479	1806628	20.54%	4.808%
21	12.982	1030	1032	1034	rBV	708727	708556	8.05%	1.886%
22	14.982	1204	1207	1210	rBV	2908419	2901030	32.98%	7.720%
23	16.148	1306	1309	1318	rBV	115561	183528	2.09%	0.488%
24	16.274	1318	1320	1323	rBV	655483	785376	8.93%	2.090%
25	17.988	1467	1470	1473	rBV	611995	795397	9.04%	2.117%

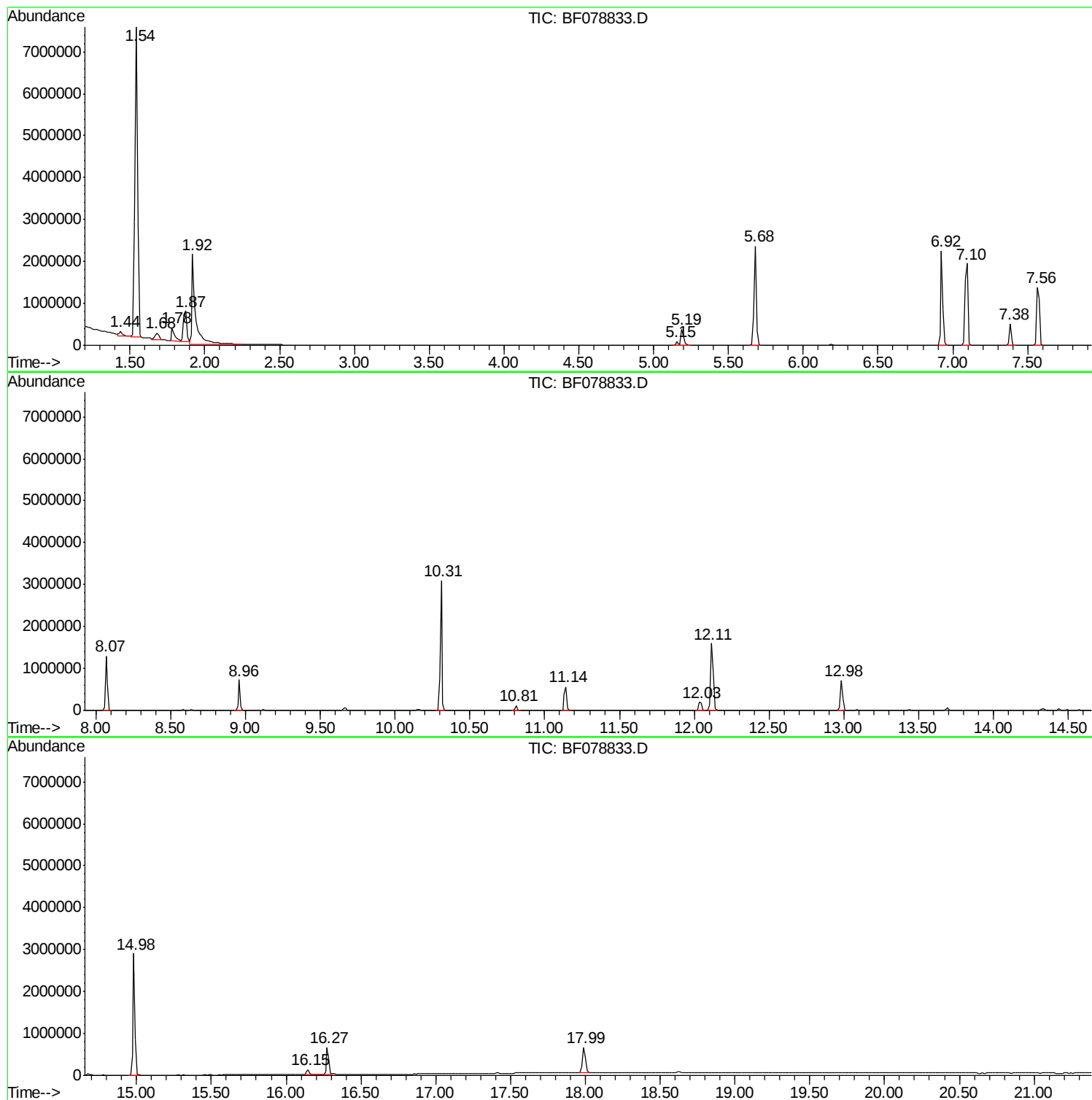
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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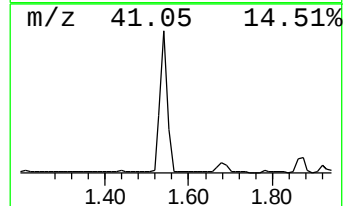
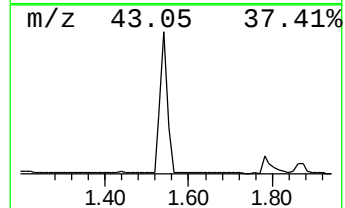
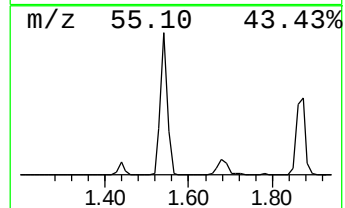
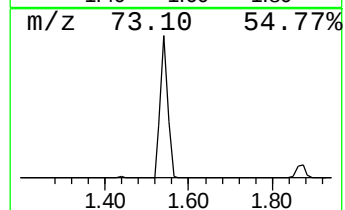
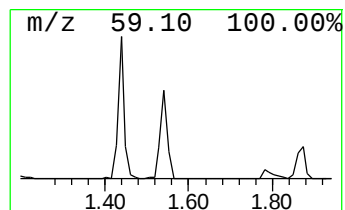
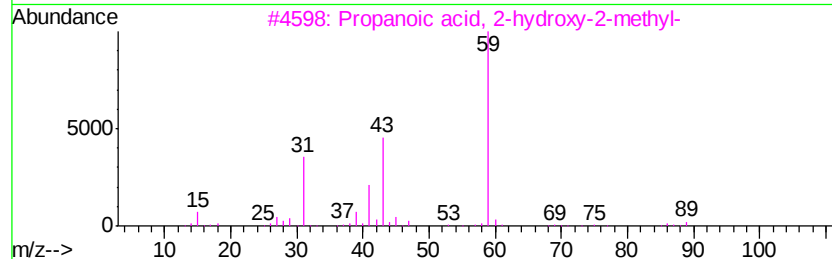
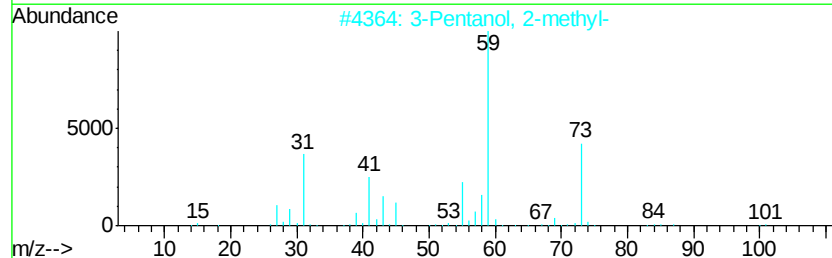
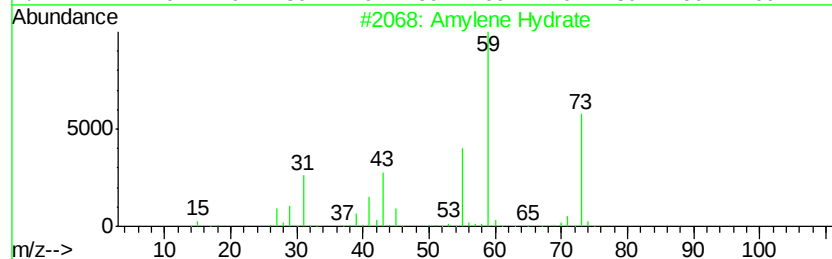
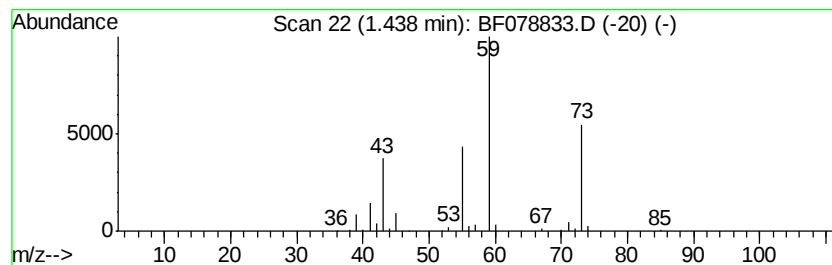
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Amylene Hydrate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.44	9.01 ng	207271	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Amylene Hydrate	88	C5H12O	000075-85-4	83
2		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	56
3		Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	42
4		Silane, trimethyl-	74	C3H10Si	000993-07-7	39
5		3-Hexanol	102	C6H14O	000623-37-0	39



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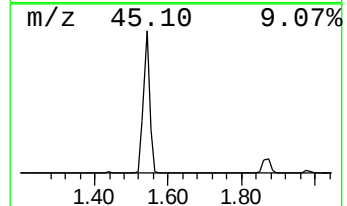
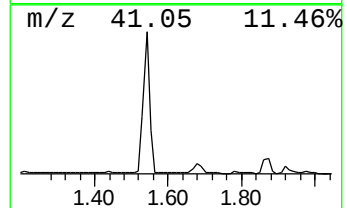
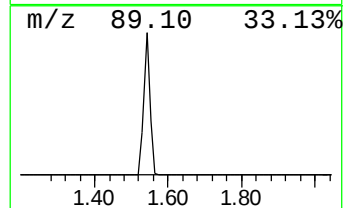
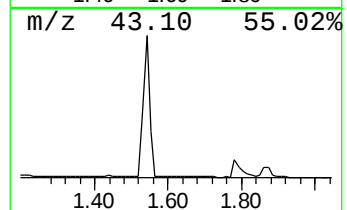
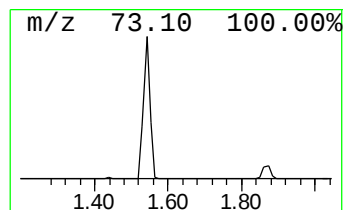
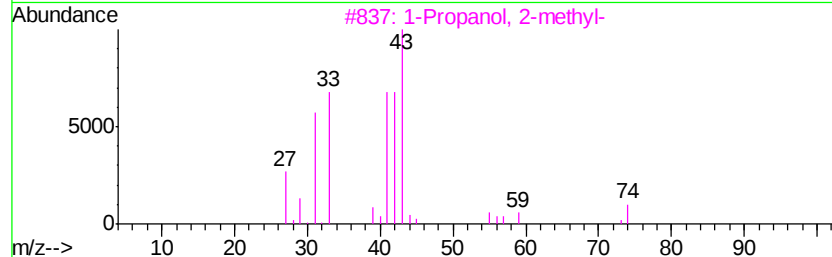
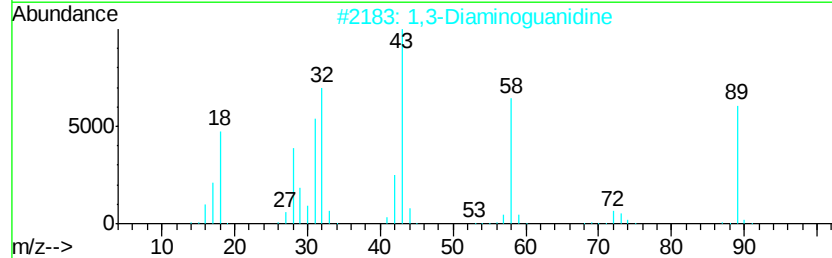
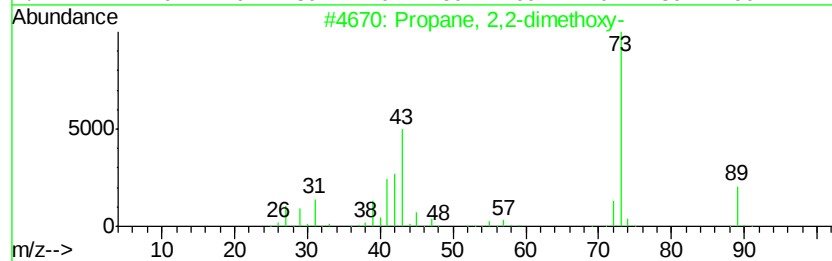
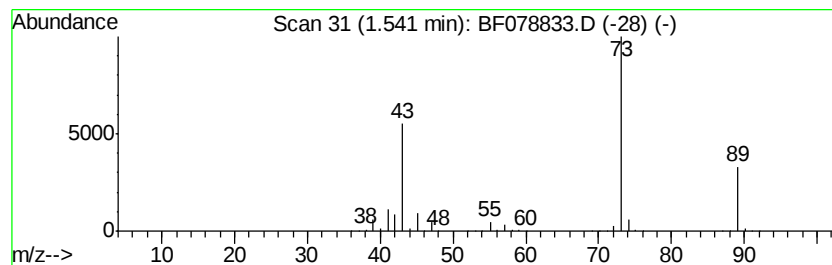
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.54	382.53 ng	8797030	1,4-Dichlorobenzene-d4	7.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	56
2			1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
3			1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9



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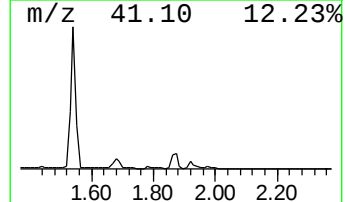
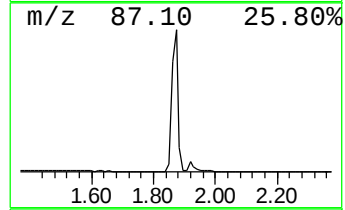
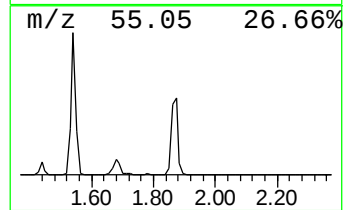
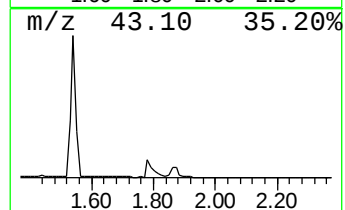
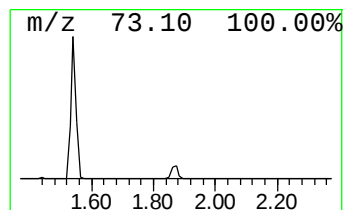
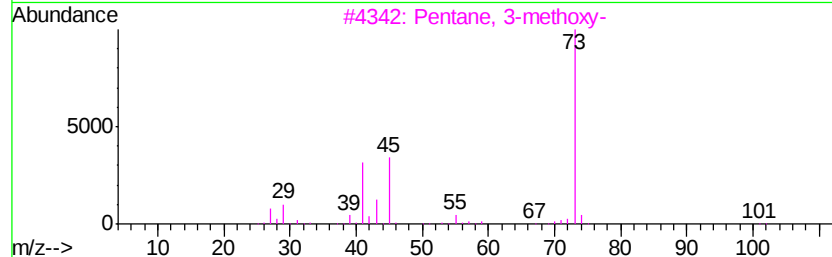
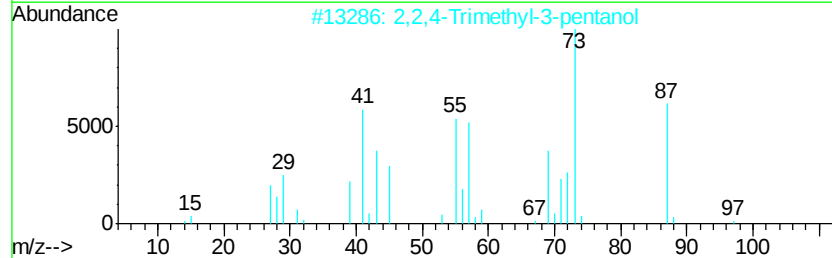
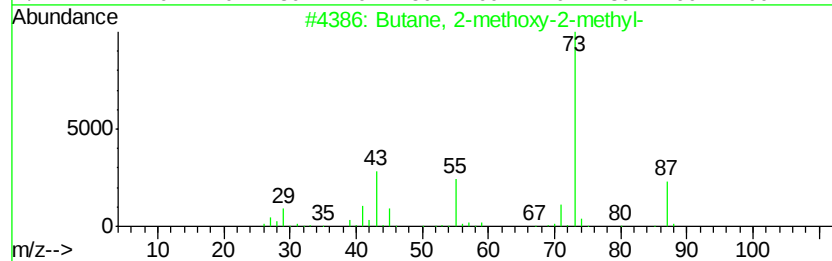
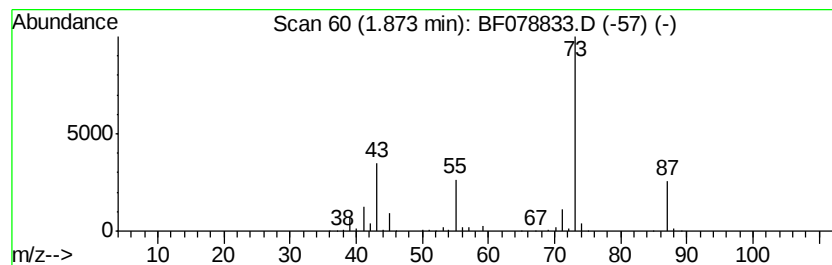
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.87	46.28 ng	1064320	1,4-Dichlorobenzene-d4	7.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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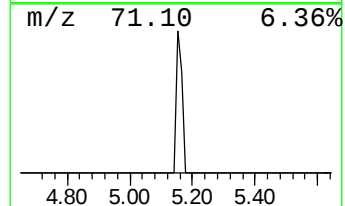
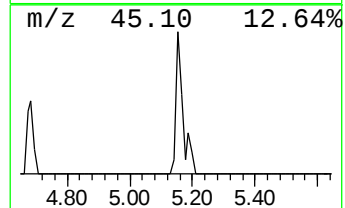
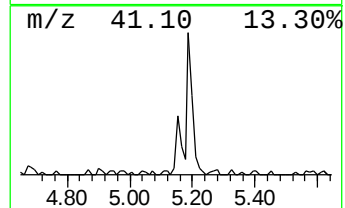
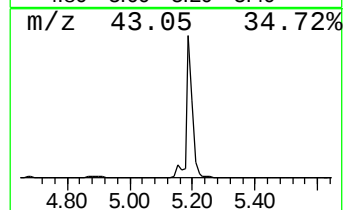
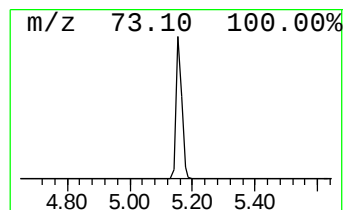
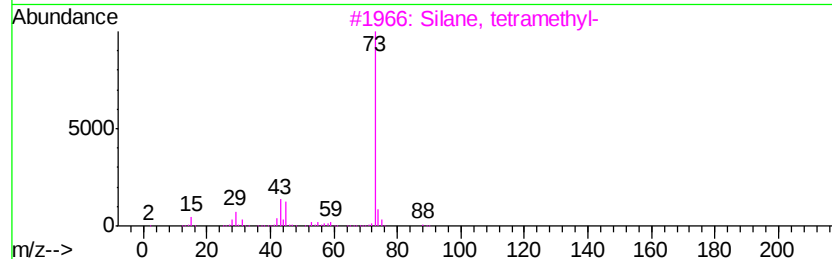
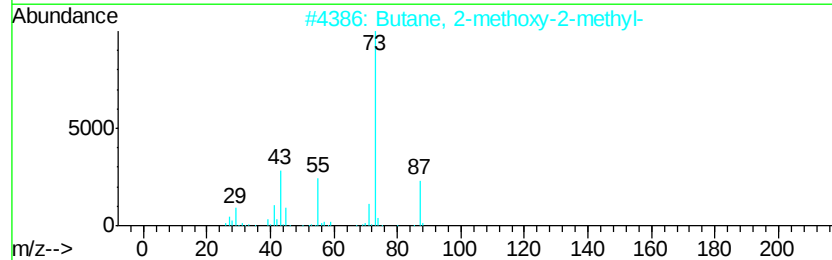
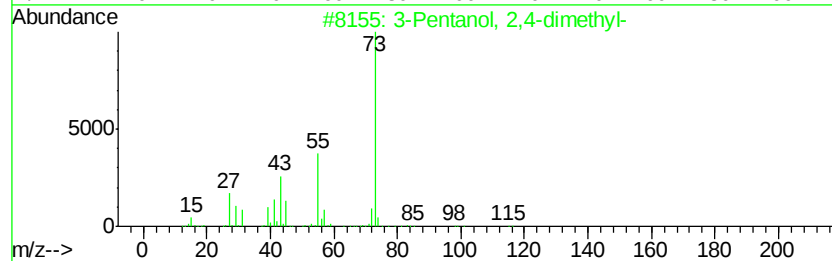
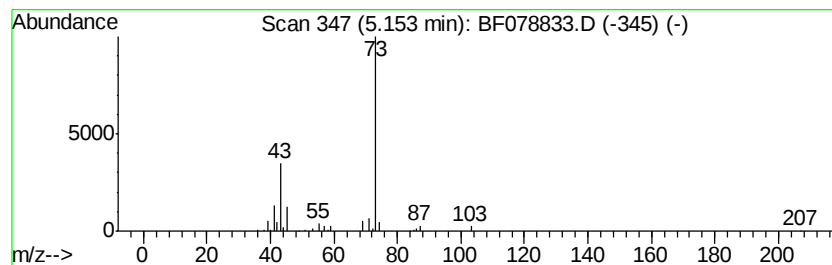
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown5.15 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	5.35 ng	123031	1,4-Dichlorobenzene-d4	7.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	36		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	36		
3	Silane, tetramethyl-	88	C4H12Si	000075-76-3	28		
4	Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9		
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9		



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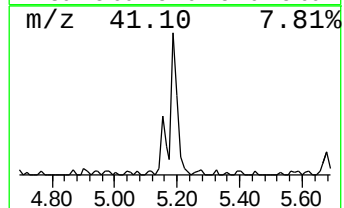
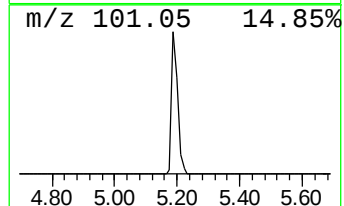
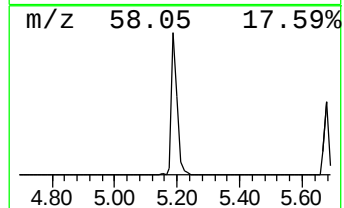
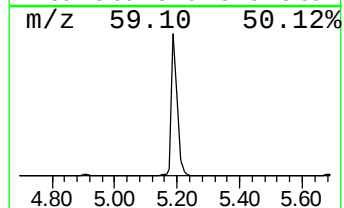
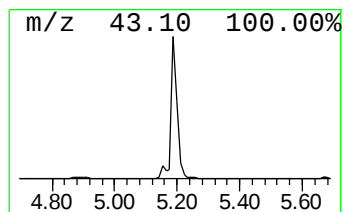
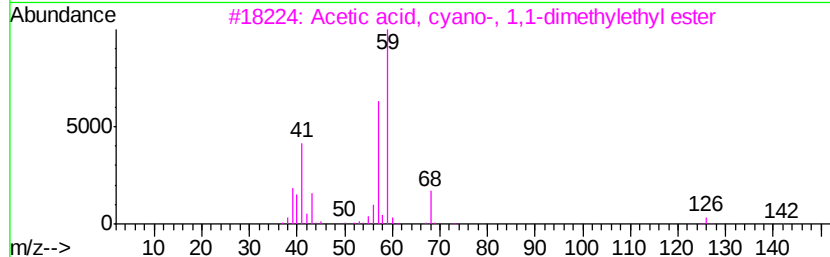
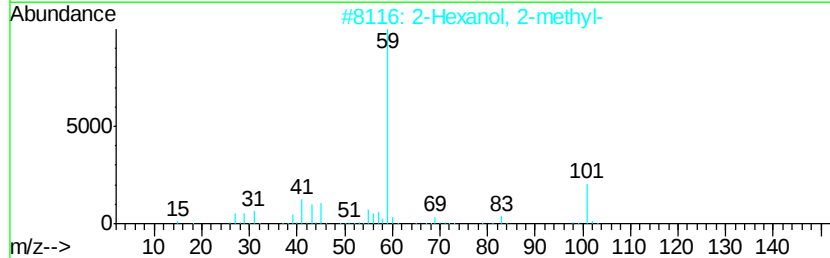
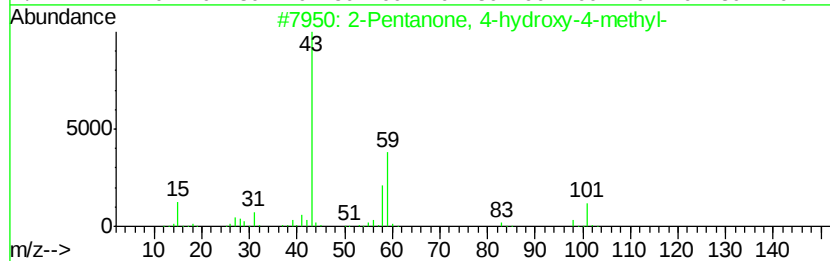
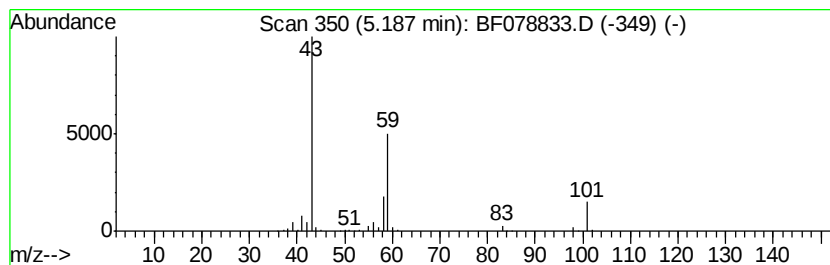
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	19.44 ng	447013	1,4-Dichlorobenzene-d4	7.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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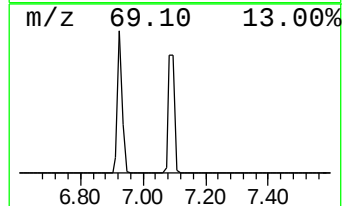
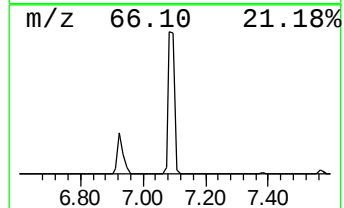
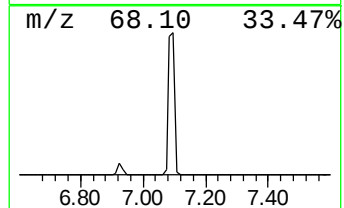
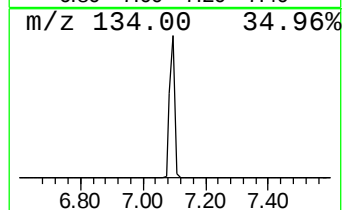
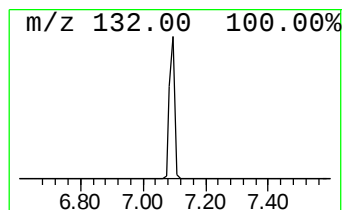
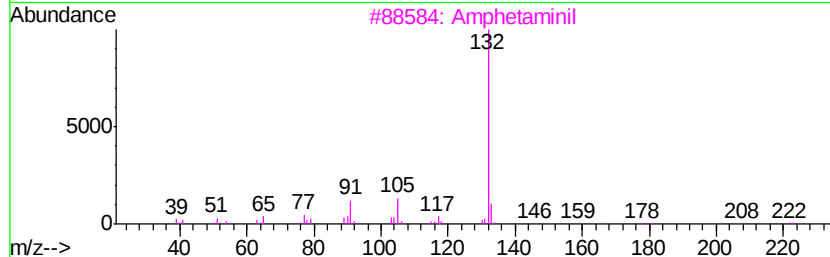
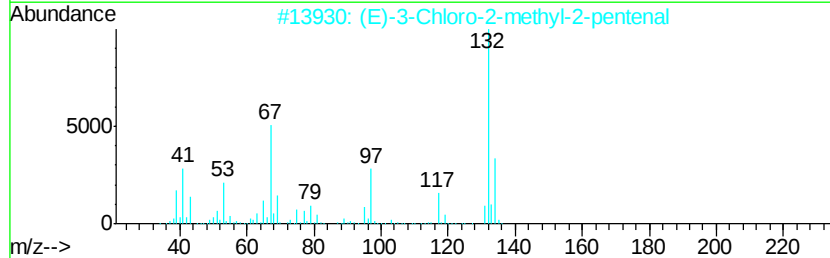
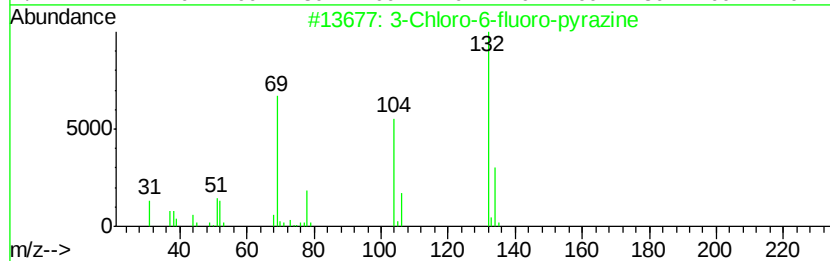
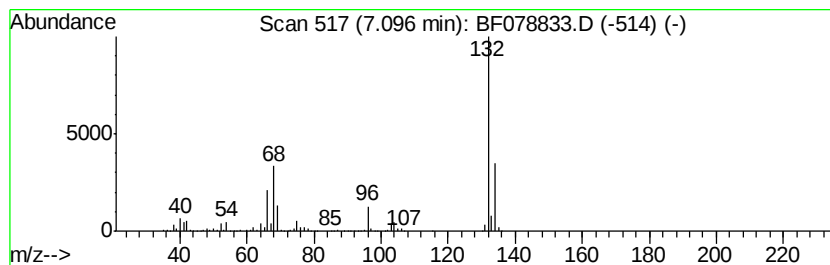
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown7.10 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.10	108.28 ng	2490030	1,4-Dichlorobenzene-d4	7.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
Data File : BF078833.D
Acq On : 30 Apr 2015 2:47
Operator : TP/IZ
Sample : G2003-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PZ-2(45-47)

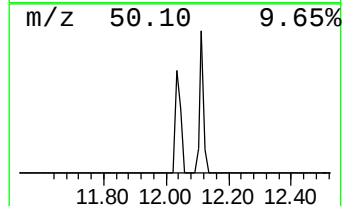
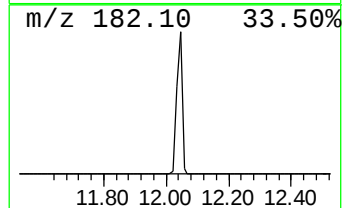
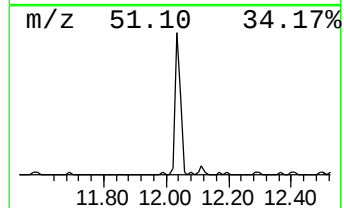
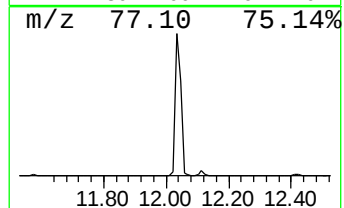
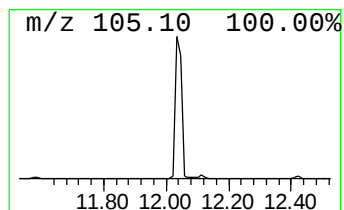
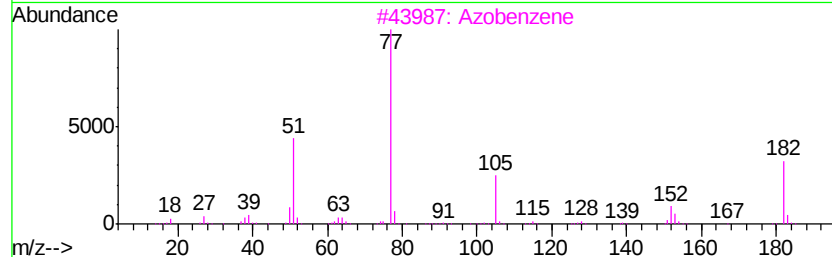
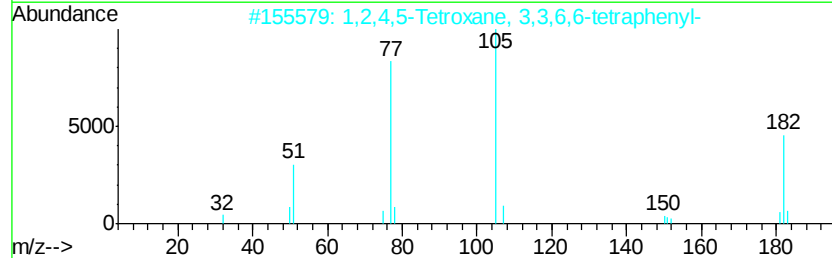
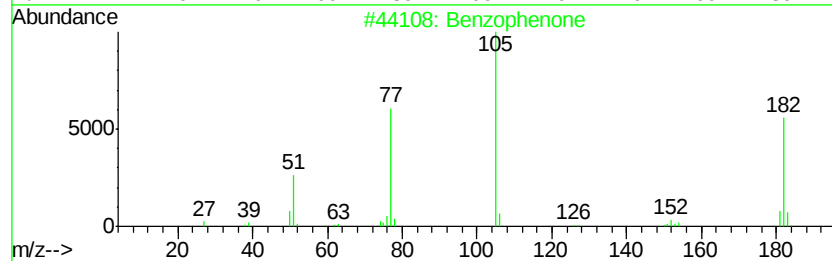
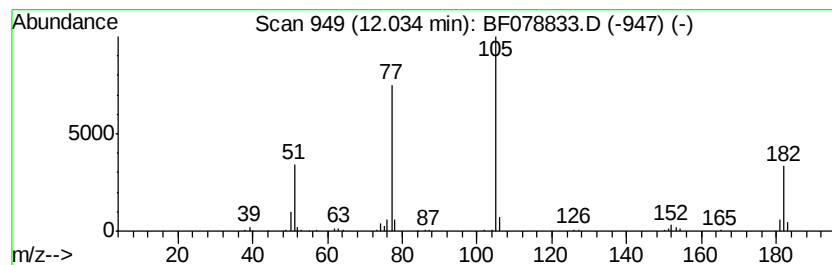
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	7.37 ng	248898	Acenaphthene-d10	11.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	78
3		Azobenzene	182	C12H10N2	000103-33-3	72
4		Benzenebutanoic acid, .gamma.-ox...	192	C11H12O3	025333-24-8	64
5		N-(1-Cyanovinyl)benzamide	172	C10H8N2O	1000186-19-1	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042915\
Data File : BF078833.D
Acq On : 30 Apr 2015 2:47
Operator : TP/IZ
Sample : G2003-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PZ-2(45-47)

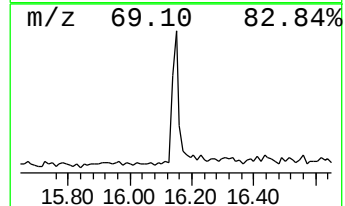
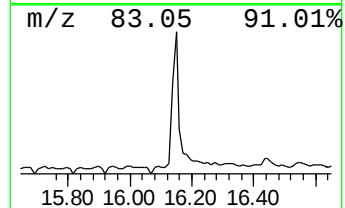
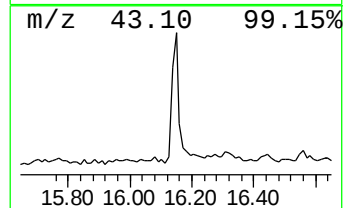
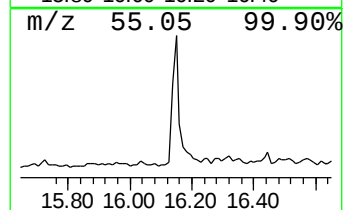
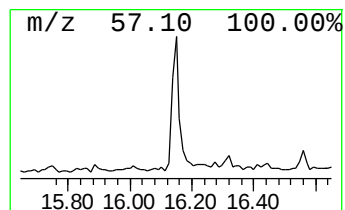
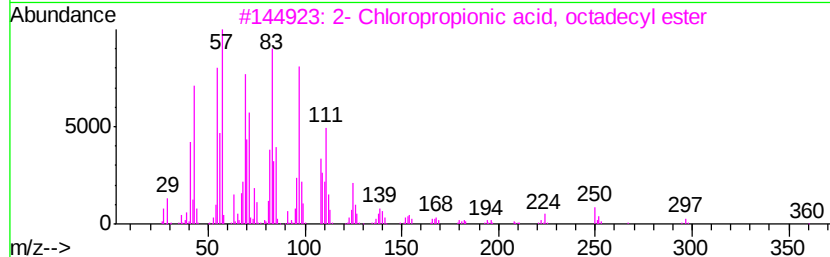
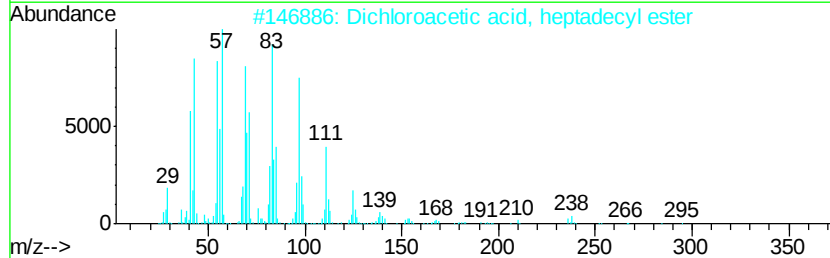
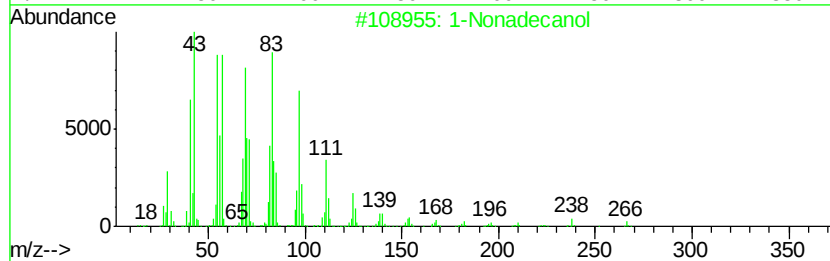
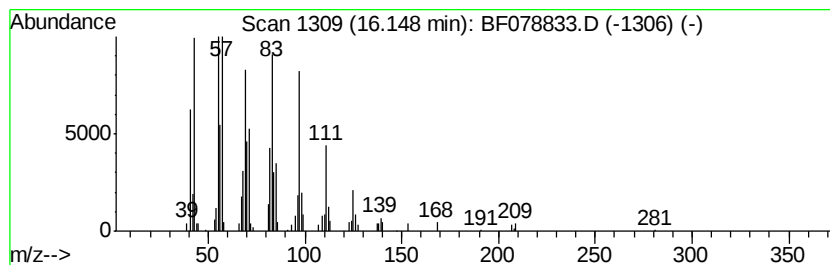
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Nonadecanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	4.67 ng	183528	Chrysene-d12	16.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecanol	284	C19H40O	001454-84-8	95
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
3		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	93
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042915\
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Acq On : 30 Apr 2015 2:47
Operator : TP/IZ
Sample : G2003-10
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PZ-2(45-47)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042815.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
Amylene Hydrate	1.44	9.0	ng	207271	1	7.38	459945	20.0
Propane, 2,2-dime...	1.54	382.5	ng	8797030	1	7.38	459945	20.0
Butane, 2-methoxy...	1.87	46.3	ng	1064320	1	7.38	459945	20.0
unknown5.15	5.15	5.3	ng	123031	1	7.38	459945	20.0
2-Pentanone, 4-hy...	5.19	19.4	ng	447013	1	7.38	459945	20.0
unknown7.10	7.10	108.3	ng	2490030	1	7.38	459945	20.0
Benzophenone	12.03	7.4	ng	248898	3	11.14	675505	20.0
1-Nonadecanol	16.15	4.7	ng	183528	5	16.27	785376	20.0