

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
 Data File : BF088939.D
 Acq On : 13 Jul 2016 00:55
 Operator : UM/SJ
 Sample : H3935-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C3-04

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-BF063016.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	1.667	6	7	11	rVB	44197	64427	3.45%	0.413%
2	1.792	16	18	24	rVV	463568	597589	32.00%	3.828%
3	1.884	24	26	40	rVB	43177	91694	4.91%	0.587%
4	2.867	109	112	118	rBV	25420	52777	2.83%	0.338%
5	4.856	284	286	291	rBV	147431	217798	11.66%	1.395%
6	5.301	322	325	328	rVB	1423963	1531679	82.02%	9.812%
7	6.353	414	417	420	rVB	1229301	1676085	89.75%	10.737%
8	6.479	425	428	430	rBV	1738495	1867475	100.00%	11.963%
9	6.707	446	448	449	rBV	340380	359446	19.25%	2.303%
10	6.856	459	461	464	rVB	982470	1265686	67.78%	8.108%
11	7.267	494	497	500	rBV	829655	1025330	54.90%	6.568%
12	7.747	536	539	541	rBV2	30933	40702	2.18%	0.261%
13	7.987	558	560	563	rVB	538994	463279	24.81%	2.968%
14	9.073	652	655	657	rBV	1726543	1812041	97.03%	11.608%
15	9.462	687	689	691	rBV	313521	246695	13.21%	1.580%
16	9.736	711	713	716	rVB	486950	429137	22.98%	2.749%
17	10.525	780	782	784	rBV	919760	831410	44.52%	5.326%
18	10.650	792	793	798	rVB	21904	36215	1.94%	0.232%
19	10.971	819	821	828	rVB	61500	101032	5.41%	0.647%
20	11.096	831	832	834	rBV	32684	45805	2.45%	0.293%
21	11.211	840	842	845	rVB	353151	330564	17.70%	2.118%
22	11.291	845	849	853	rVB2	12641	25836	1.38%	0.166%
23	11.451	861	863	868	rVB2	14337	25954	1.39%	0.166%
24	11.759	888	890	892	rBV	38846	30436	1.63%	0.195%
25	11.816	892	895	900	rVB	30926	43890	2.35%	0.281%
26	12.799	978	981	983	rBV	1270044	1135336	60.80%	7.273%
27	13.016	994	1000	1004	rBV3	12867	24838	1.33%	0.159%
28	13.119	1007	1009	1017	rBV	56277	100952	5.41%	0.647%
29	13.657	1053	1056	1057	rBV	49072	44464	2.38%	0.285%
30	13.702	1057	1060	1064	rBV	111529	119532	6.40%	0.766%
31	13.839	1069	1072	1074	rBV	413007	419244	22.45%	2.686%
32	13.977	1083	1084	1086	rVB	49761	36374	1.95%	0.233%
33	14.022	1086	1088	1091	rBV	50335	88147	4.72%	0.565%
34	14.331	1113	1115	1119	rBV2	30848	60384	3.23%	0.387%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.834	1157	1159	1162	rVB	14786	28303	1.52%	0.181%
36	14.925	1165	1167	1170	rBV2	21578	33495	1.79%	0.215%
37	15.211	1189	1192	1195	rBV	282052	306046	16.39%	1.961%

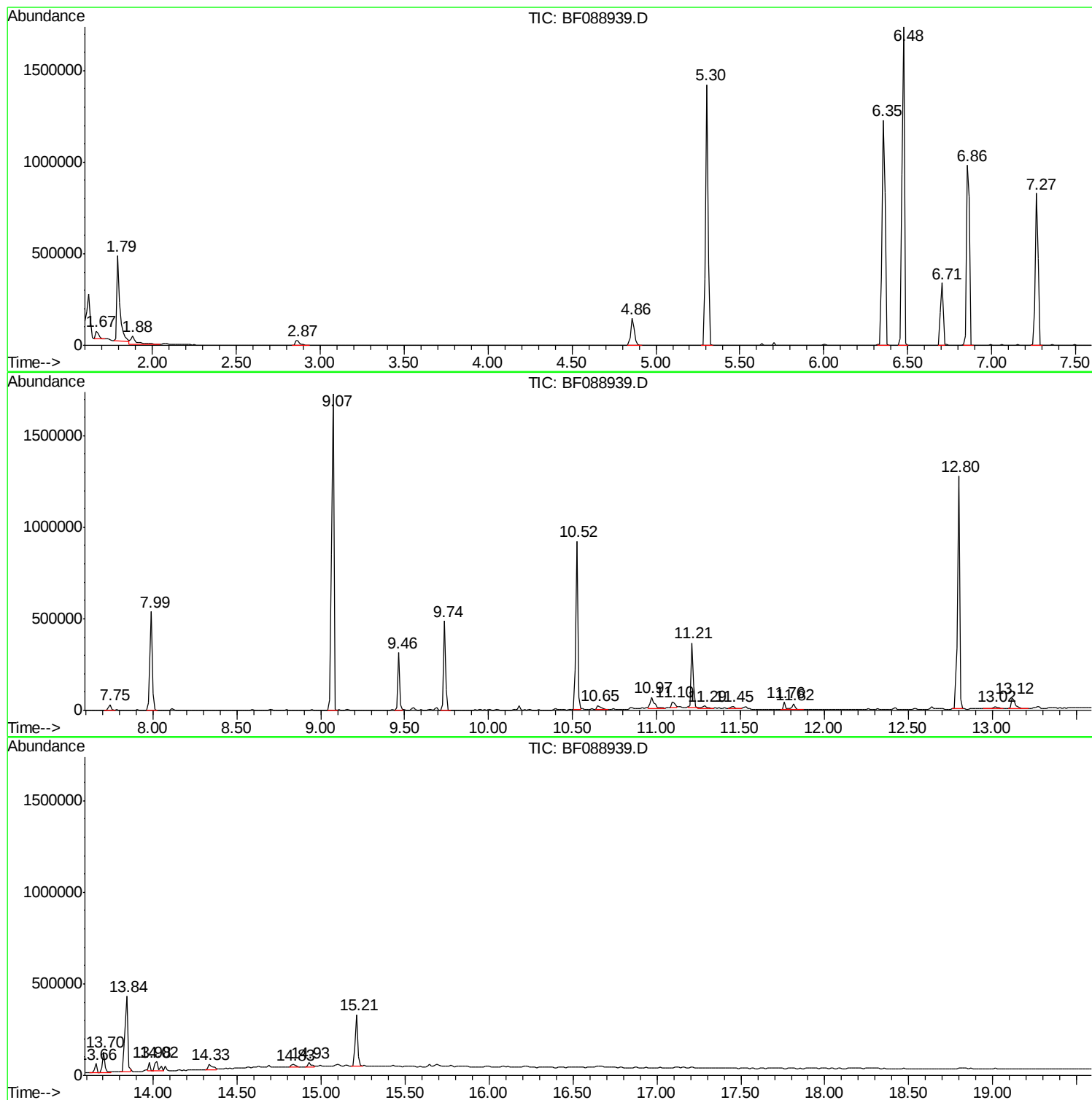
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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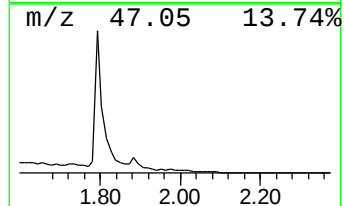
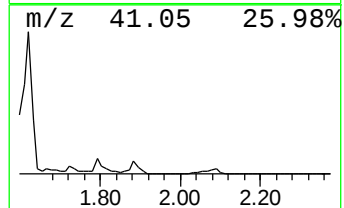
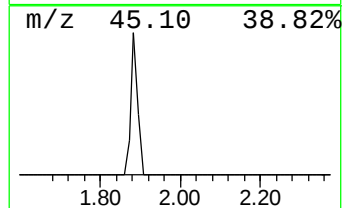
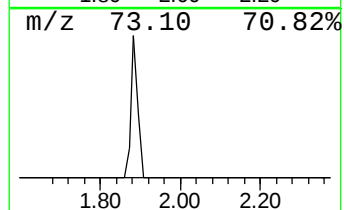
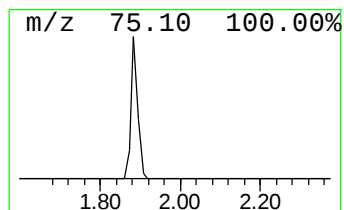
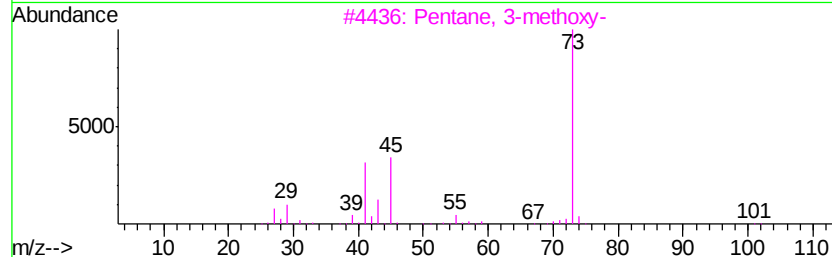
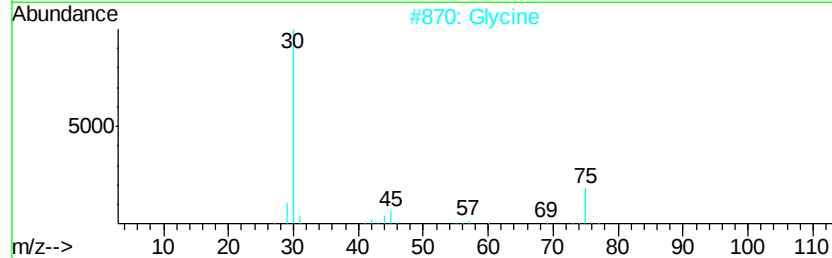
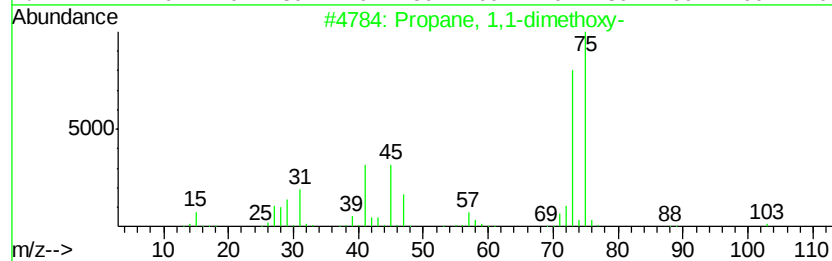
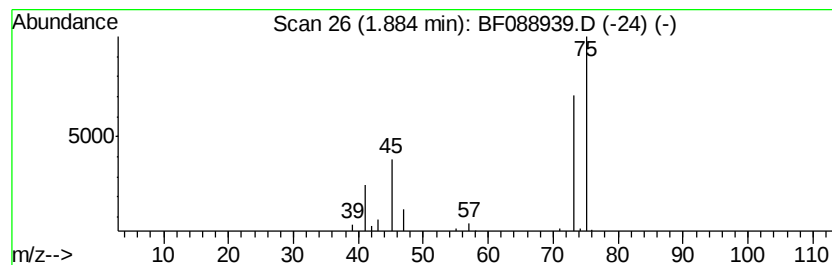
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.88	5.10 ng	91694	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	72
2		Glycine	75	C2H5NO2	000056-40-6	9
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9
4		1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9
5		Methane, nitroso-	45	CH3NO	000865-40-7	4



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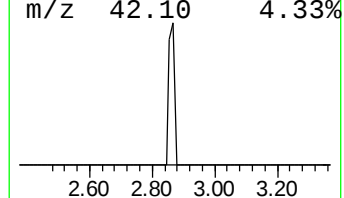
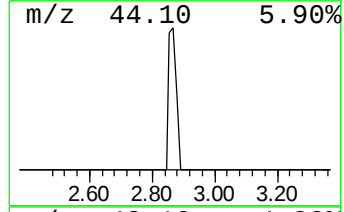
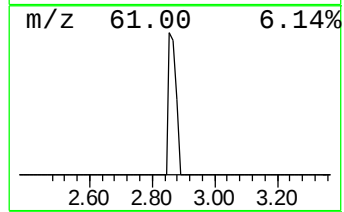
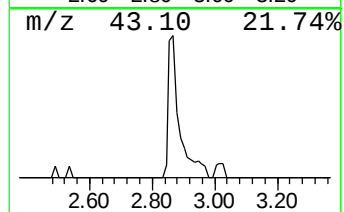
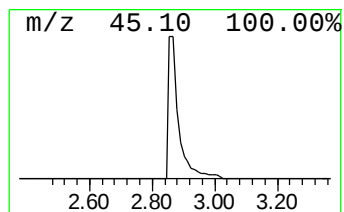
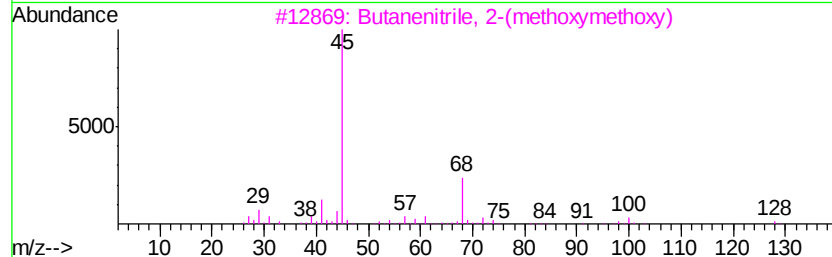
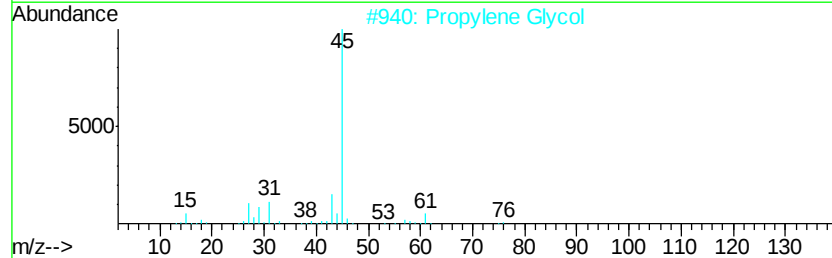
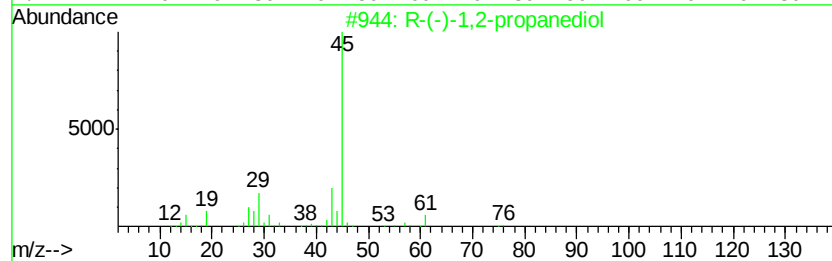
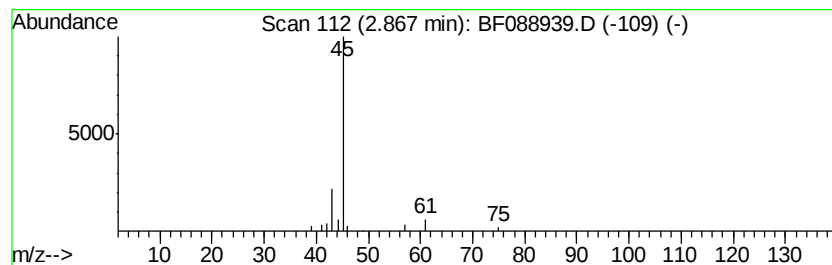
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 R-(-)-1,2-propanediol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	2.94 ng	52777	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	64
2		Propylene Glycol	76	C3H8O2	000057-55-6	9
3		Butanenitrile, 2-(methoxymethoxy)	129	C6H11NO2	1000196-86-5	9
4		Methoxyacetic acid, pentyl ester	160	C8H16O3	168920-35-2	9
5		Isopropyl Alcohol	60	C3H8O	000067-63-0	9



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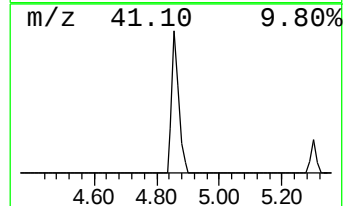
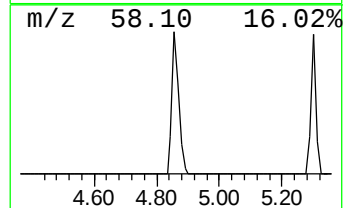
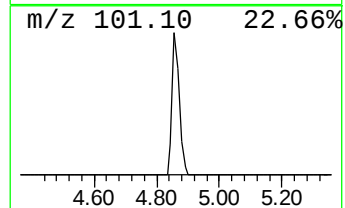
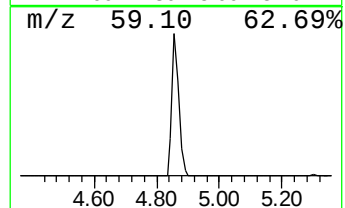
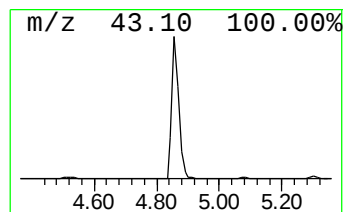
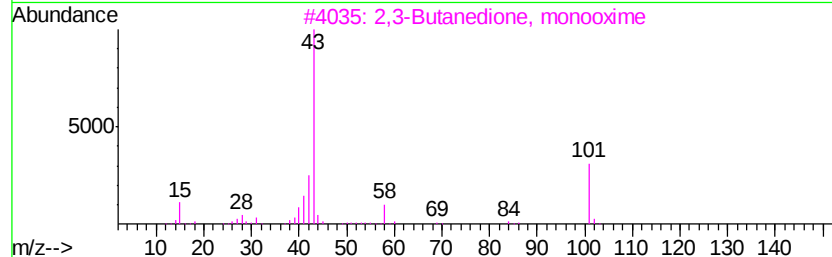
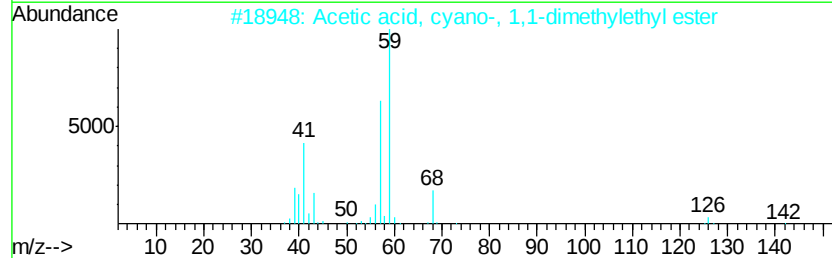
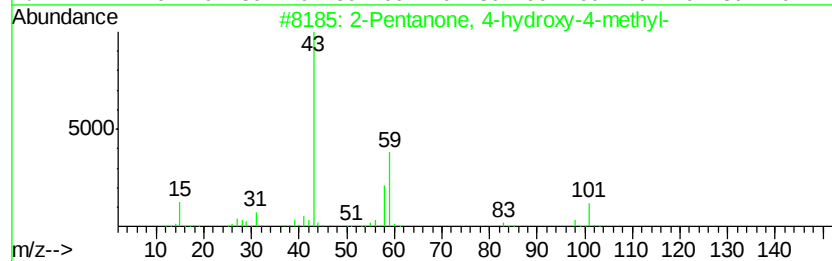
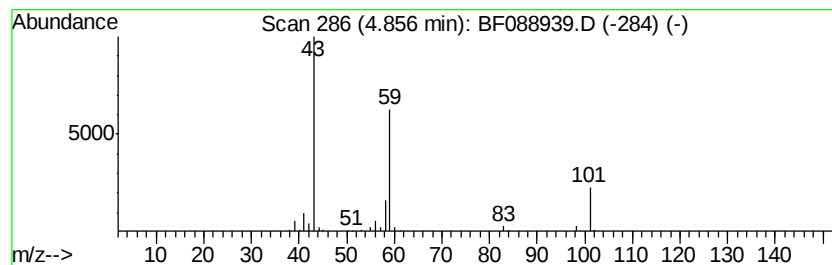
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	12.12 ng	217798	1,4-Dichlorobenzene-d4	6.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	17
4	Morpholine, 4-methyl-	101	C5H11NO		000109-02-4	9
5	Butane, 1-ethoxy-	102	C6H14O		000628-81-9	9



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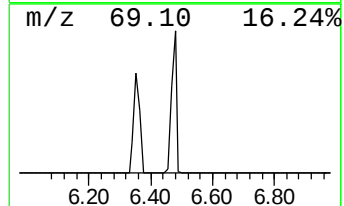
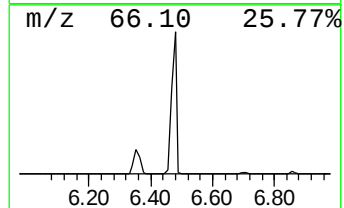
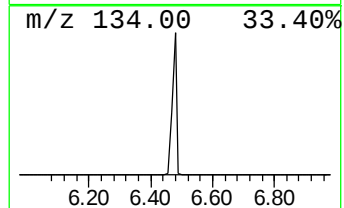
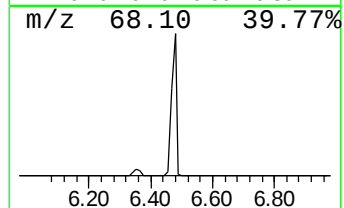
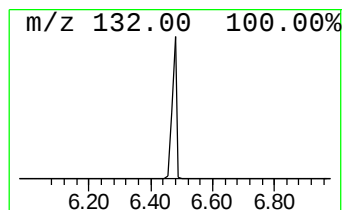
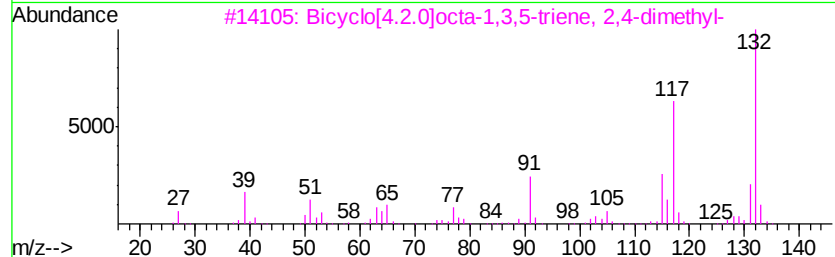
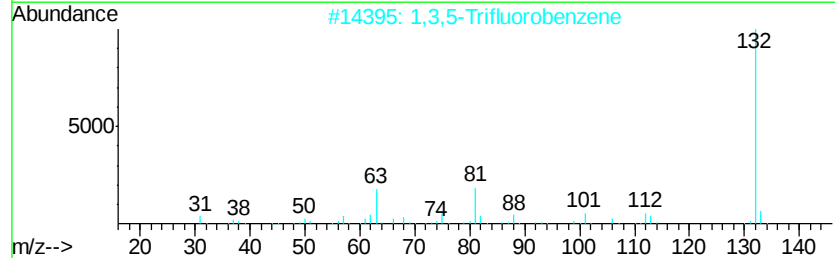
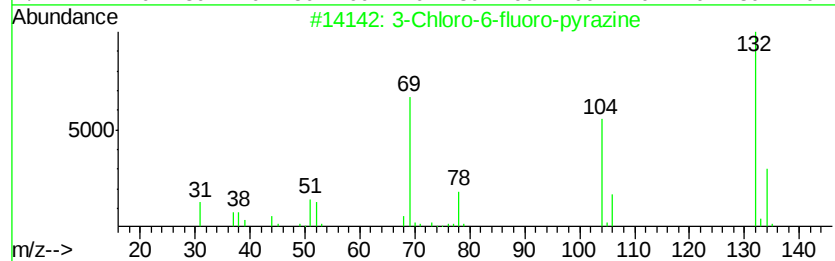
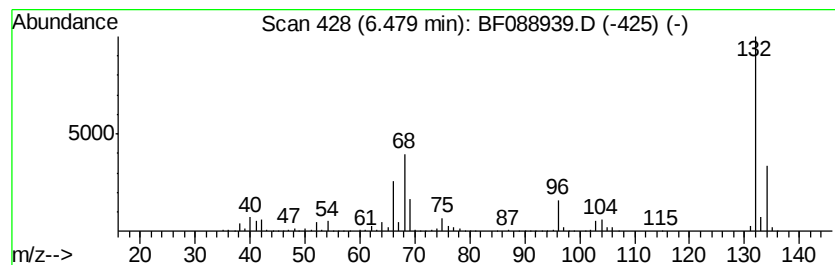
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	103.91 ng	1867480	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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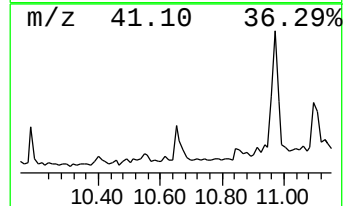
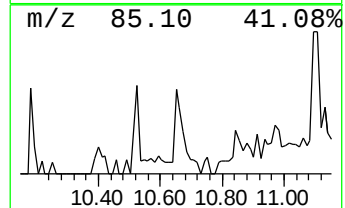
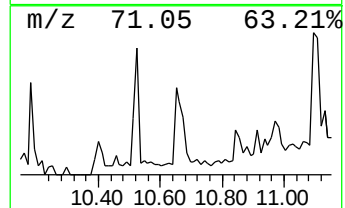
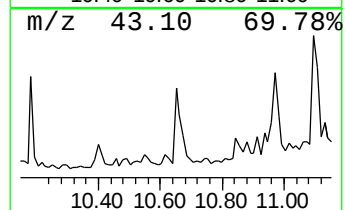
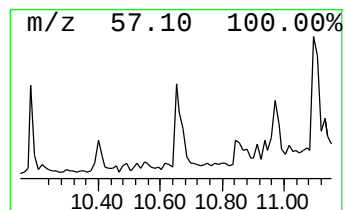
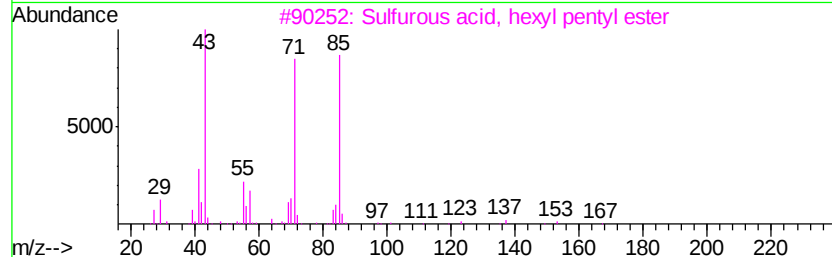
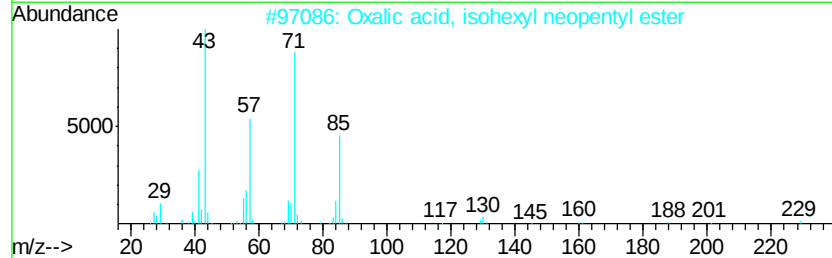
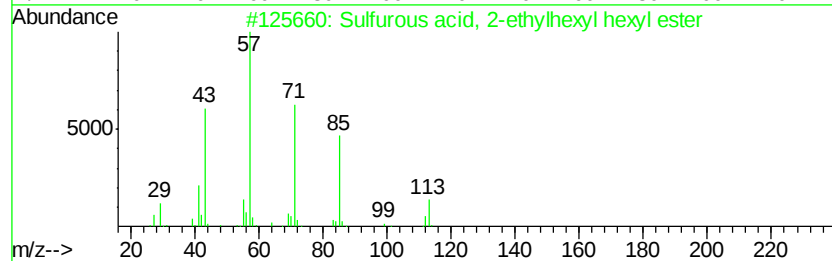
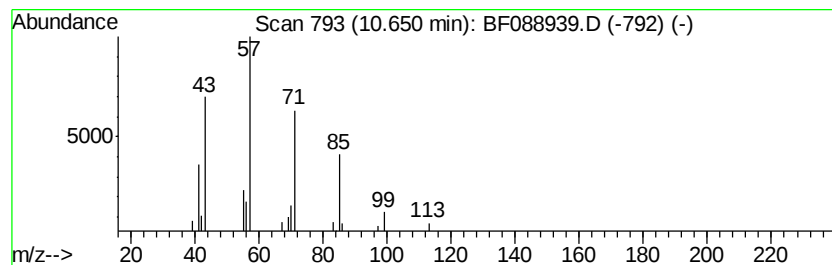
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Peak Number 8 Sulfurous acid, 2-ethylhexy... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	2.19 ng	36215	Phenanthrene-d10	11.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
2		Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	64
3		Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	59
4		Tetradecane	198	C14H30	000629-59-4	56
5		Tridecane	184	C13H28	000629-50-5	56



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Sample : H3935-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

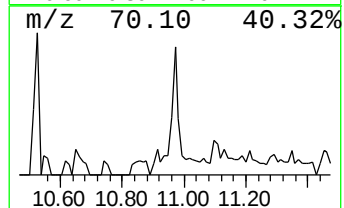
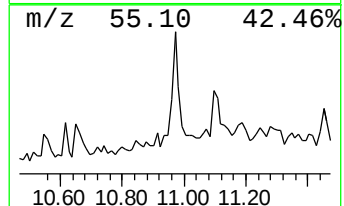
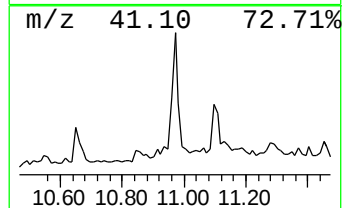
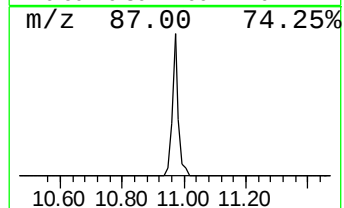
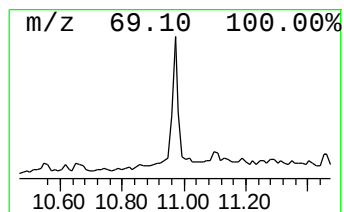
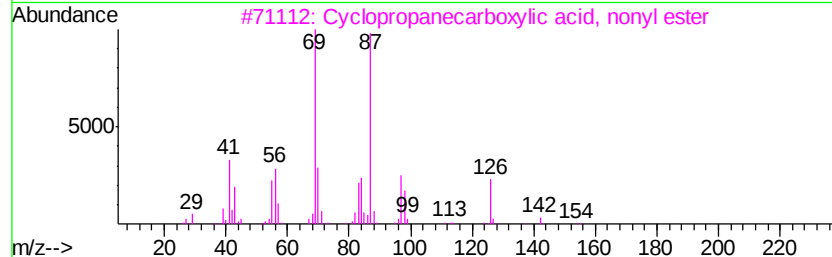
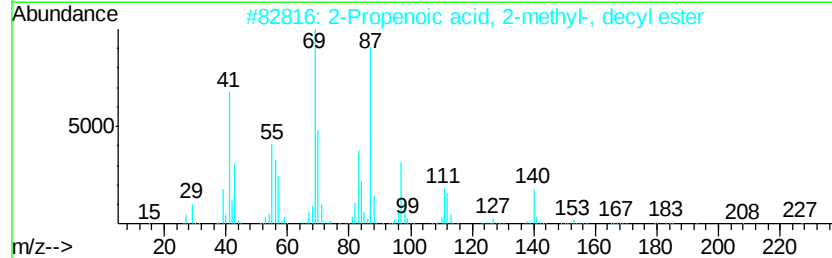
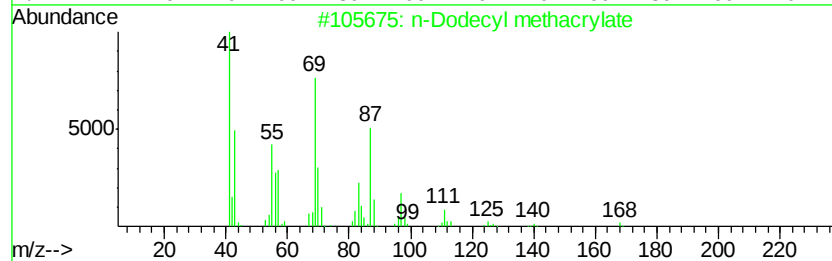
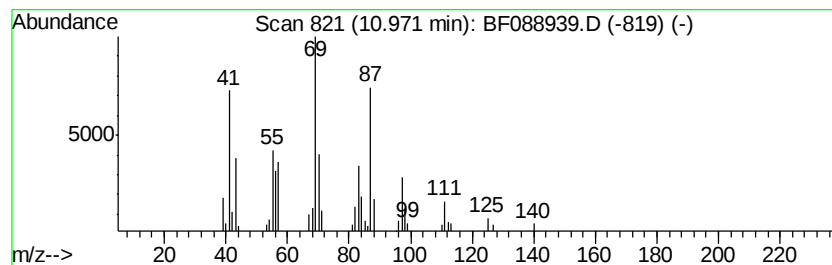
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Dodecyl methacrylate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.97	6.11 ng	101032	Phenanthrene-d10	11.21	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	91
2	2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	91
3	Cyclopropanecarboxylic acid, non...	212	C13H24O2	060128-06-5	72
4	Cyclopentane, ethyl-	98	C7H14	001640-89-7	38
5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088939.D
Acq On : 13 Jul 2016 00:55
Operator : UM/SJ
Sample : H3935-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

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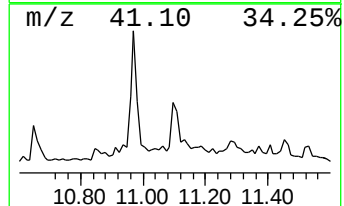
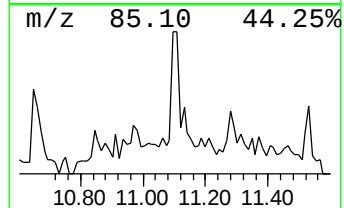
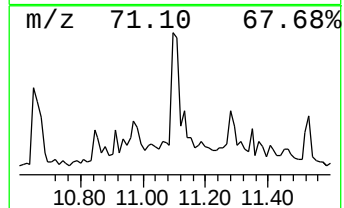
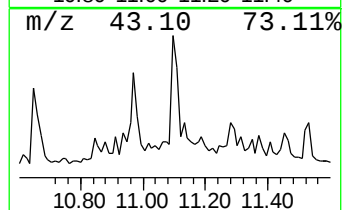
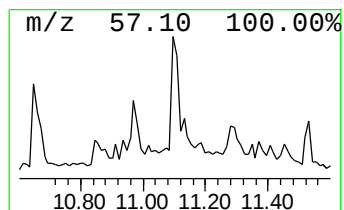
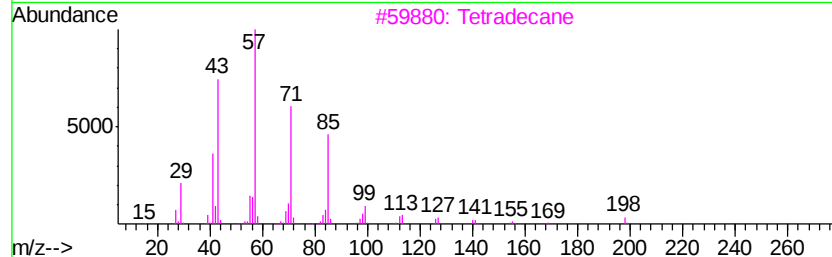
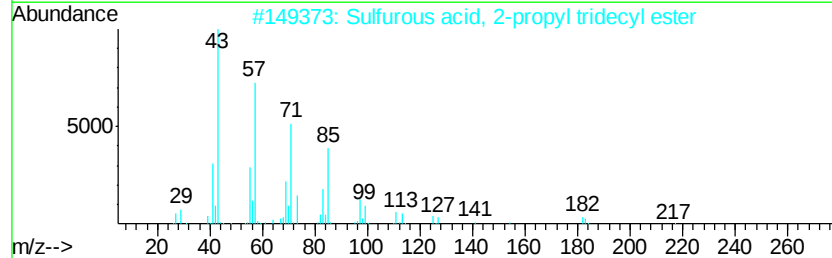
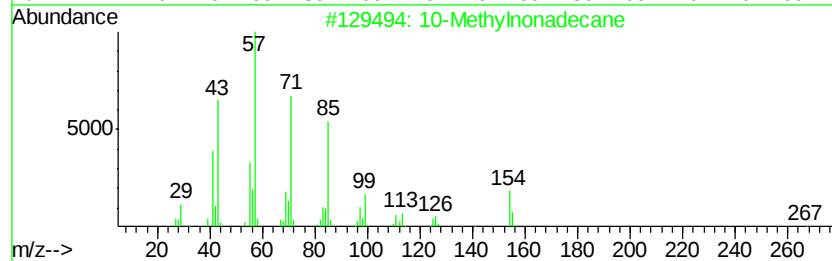
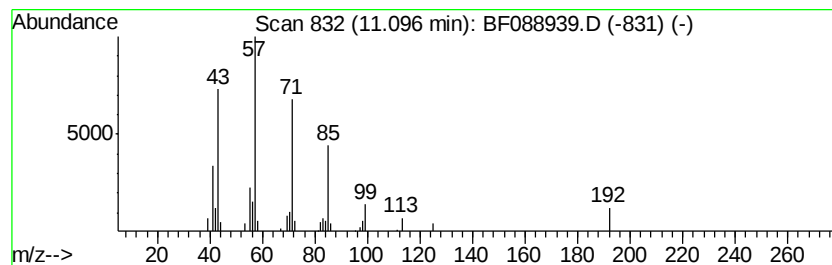
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 10-Methylnonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.77 ng	45805	Phenanthrene-d10	11.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Methylnonadecane	282	C20H42	056862-62-5	86
2		Sulfurous acid, 2-propyl tridecy...	306	C16H34O3S	1000309-12-4	86
3		Tetradecane	198	C14H30	000629-59-4	80
4		Tridecane	184	C13H28	000629-50-5	72
5		Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
Data File : BF088939.D
Acq On : 13 Jul 2016 00:55
Operator : UM/SJ
Sample : H3935-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

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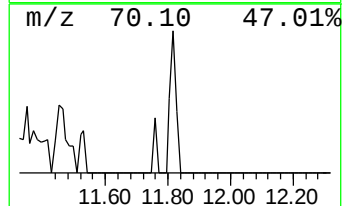
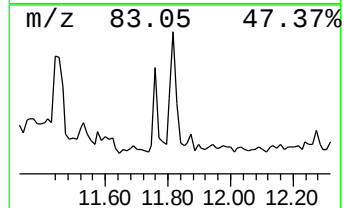
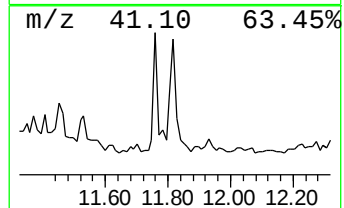
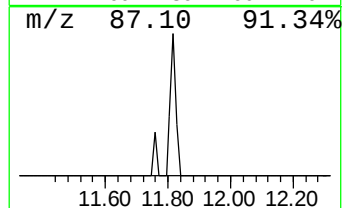
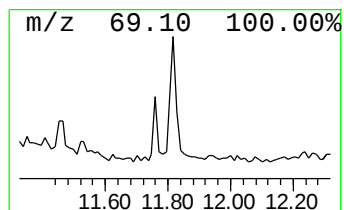
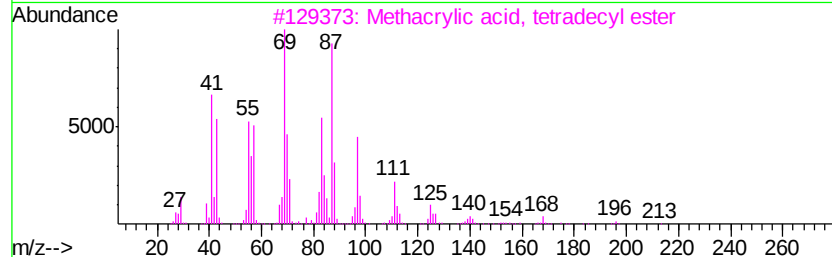
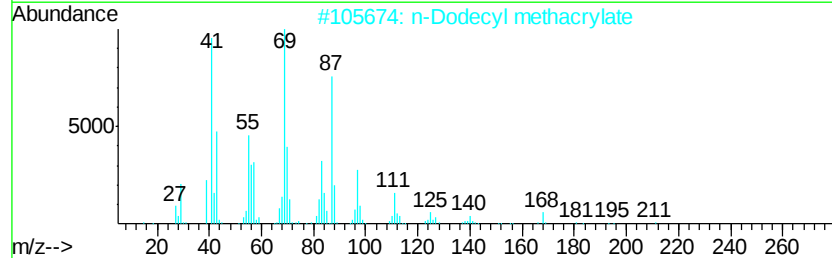
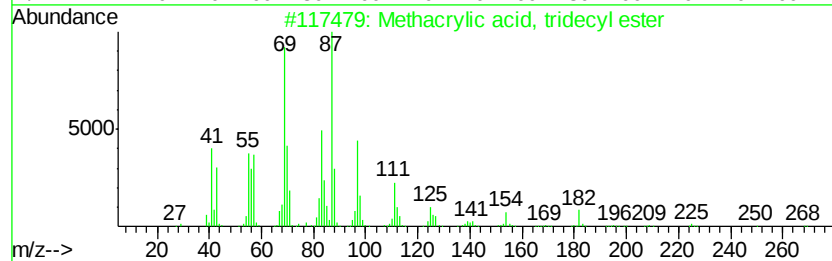
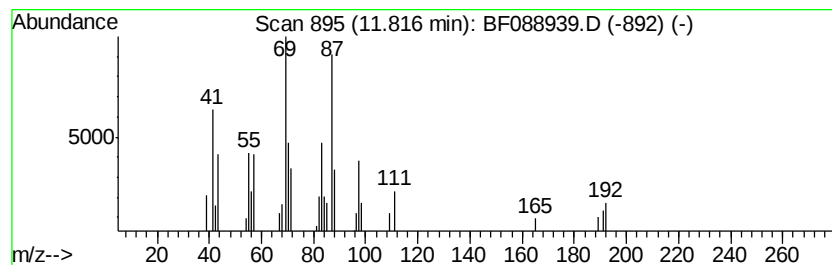
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methacrylic acid, tridecyl ... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.82	2.66 ng	43890	Phenanthrene-d10	11.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methacrylic acid, tridecyl ester	268	C17H32O2	1000340-28-9	91
2		n-Dodecyl methacrylate	254	C16H30O2	000142-90-5	90
3		Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	78
4		2-Propenoic acid, 2-methyl-, dec...	226	C14H26O2	003179-47-3	78
5		5-Hexadecanol	242	C16H34O	021078-87-5	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088939.D
Acq On : 13 Jul 2016 00:55
Operator : UM/SJ
Sample : H3935-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

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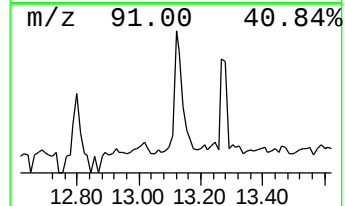
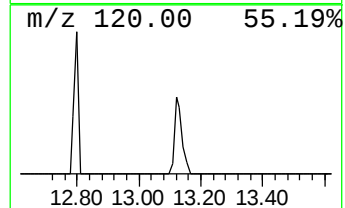
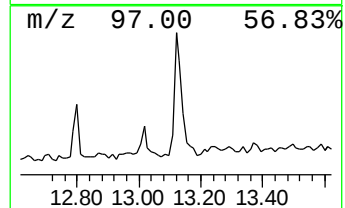
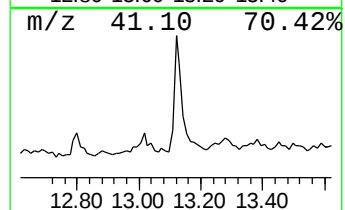
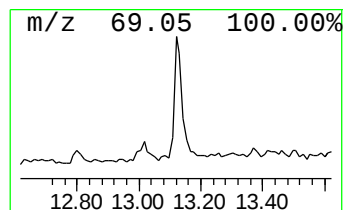
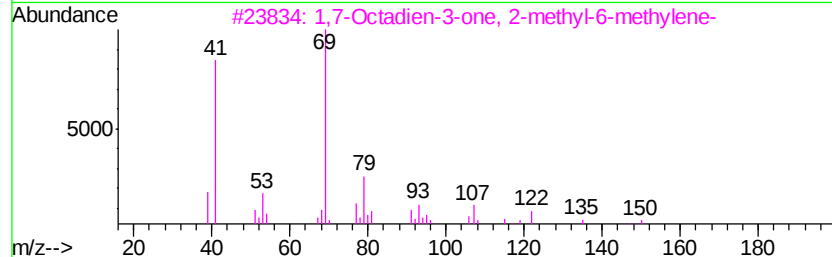
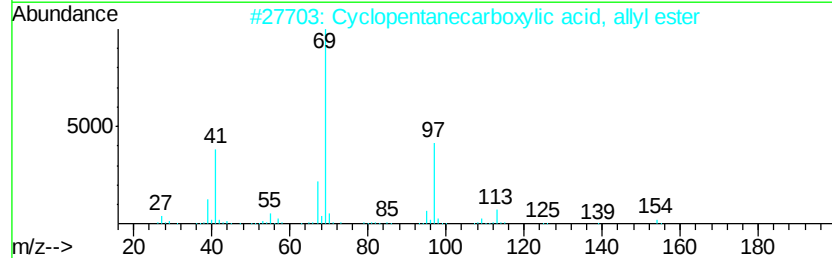
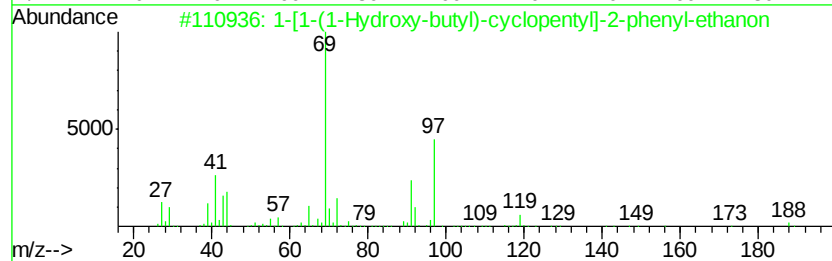
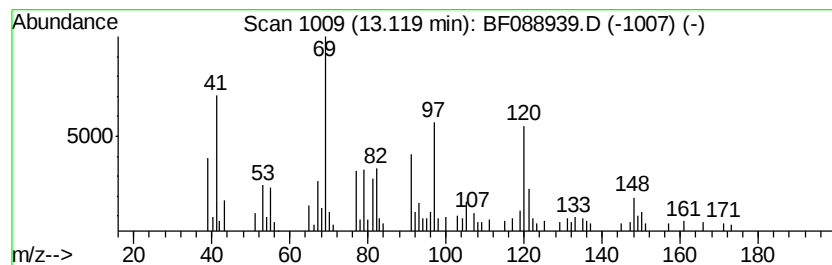
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown13.12 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	4.82 ng	100952	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-[1-(1-Hydroxy-butyl)-cyclopent...	260	C17H24O2	097234-38-3	43
2		Cyclopentanecarboxylic acid, all...	154	C9H14O2	178905-33-4	38
3		1,7-Octadien-3-one, 2-methyl-6-m...	150	C10H14O	041702-60-7	38
4		1-(10,10-Dimethyl-3,3-dioxo-3-th...	311	C16H25N03S	1000210-42-6	38
5		Pentane, 1,5-dibromo-	228	C5H10Br2	000111-24-0	35



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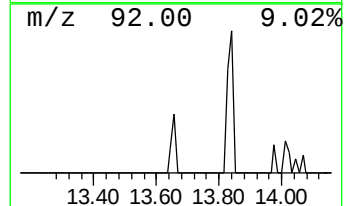
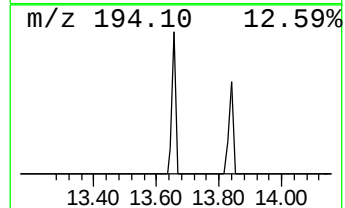
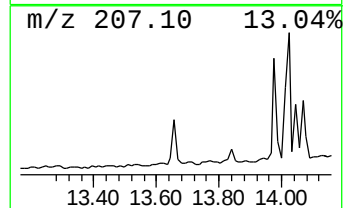
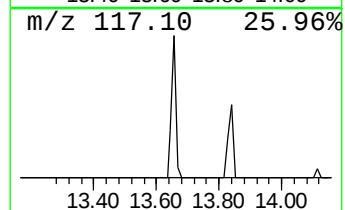
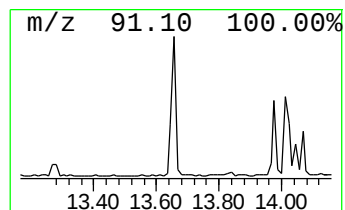
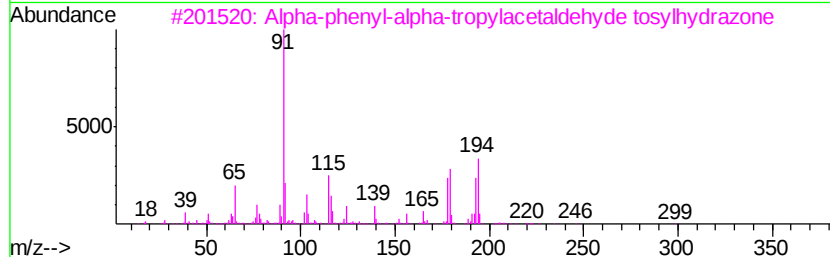
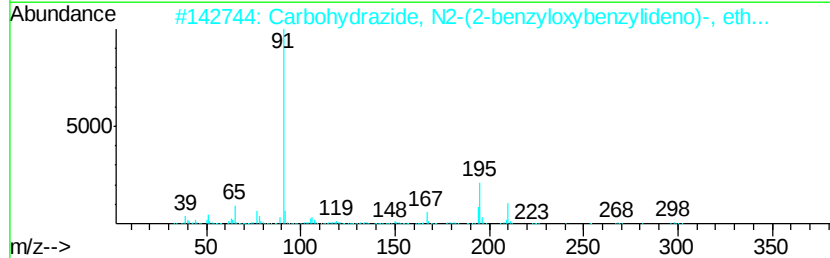
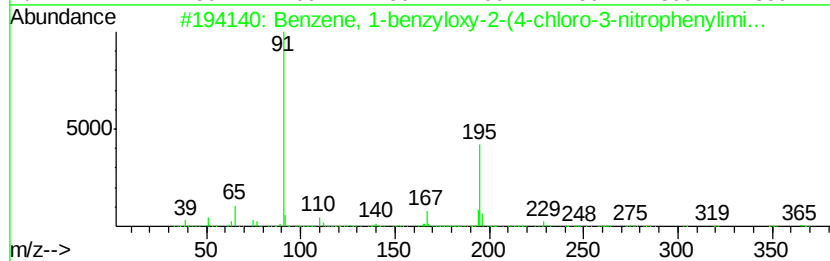
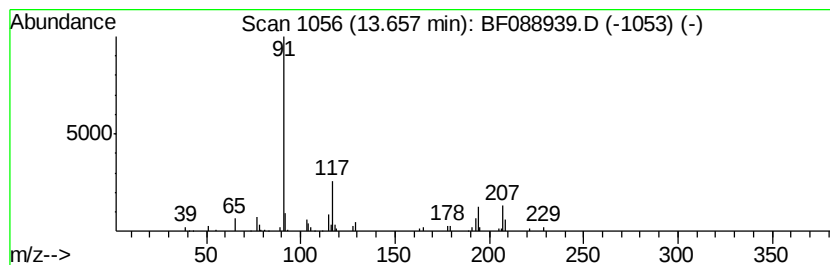
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown13.66 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	2.12 ng	44464	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-benzvloxy-2-(4-chloro...	366	C20H15ClN2O3	303761-04-8	12
2		Carbohydrazide, N2-(2-benzvloxyb...	298	C17H18N2O3	1000263-89-4	12
3		Alpha-phenyl-alpha-tropylacetalde...	378	C22H22N2O2S	022532-16-7	12
4		4-Imidazolidinone, 5-(phenylmeth...	206	C10H10N2OS	006330-09-2	12
5		(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	12



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
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Acq On : 13 Jul 2016 00:55
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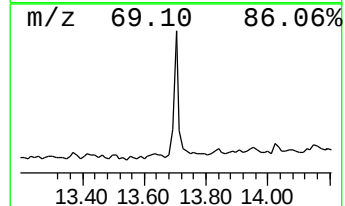
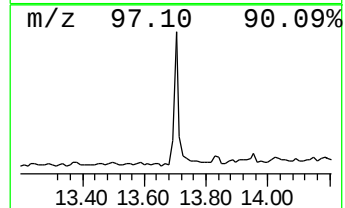
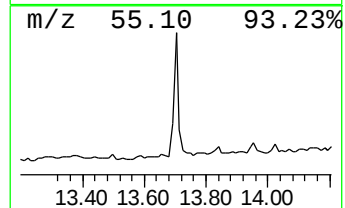
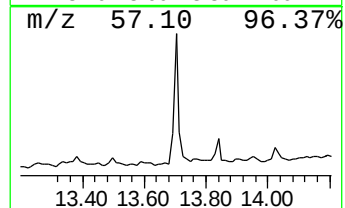
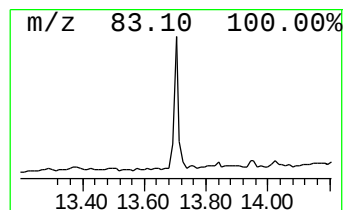
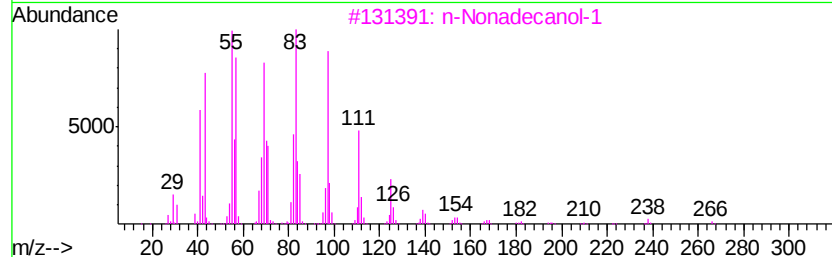
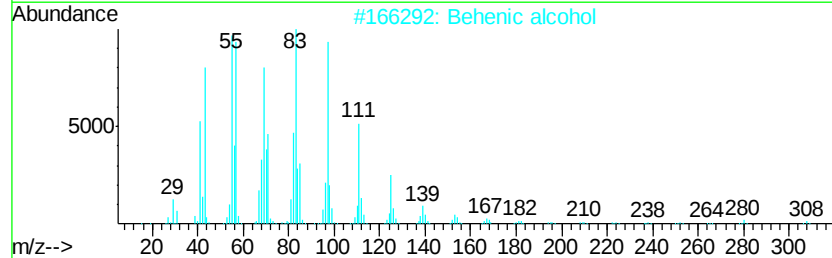
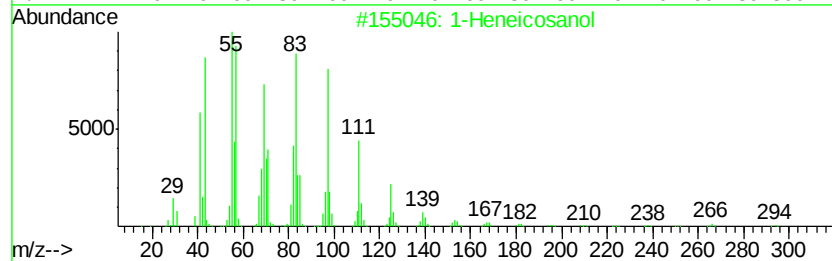
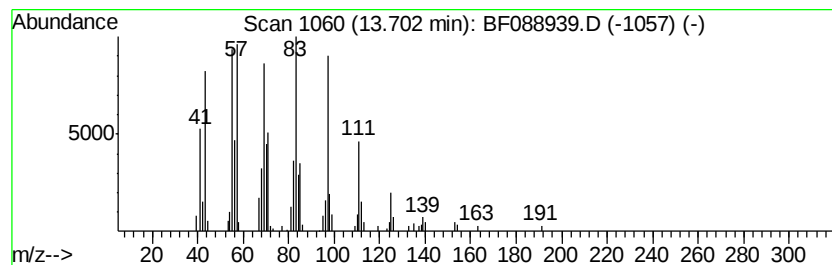
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Heneicosanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.70	5.70 ng	119532	Chrysene-d12		13.84
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Behenic alcohol	326	C22H46O	000661-19-8	95
3	n-Nonadecanol-1	284	C19H40O	001454-84-8	95
4	1-Nonadecene	266	C19H38	018435-45-5	94
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071216\
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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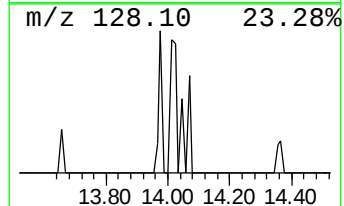
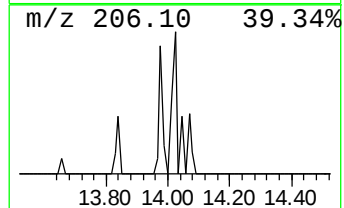
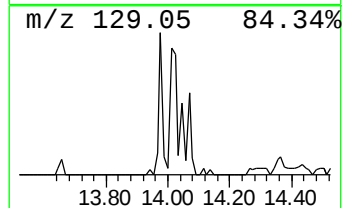
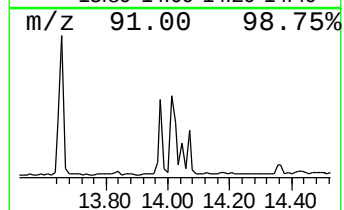
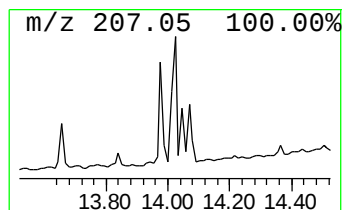
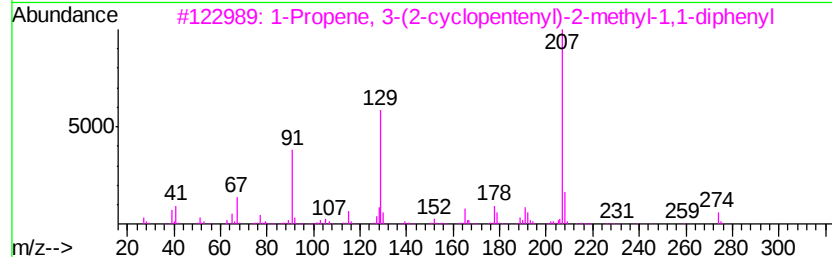
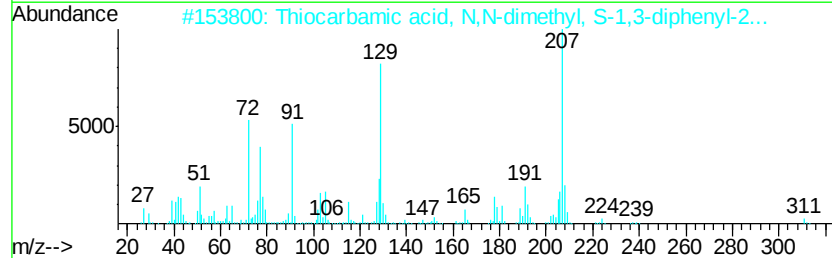
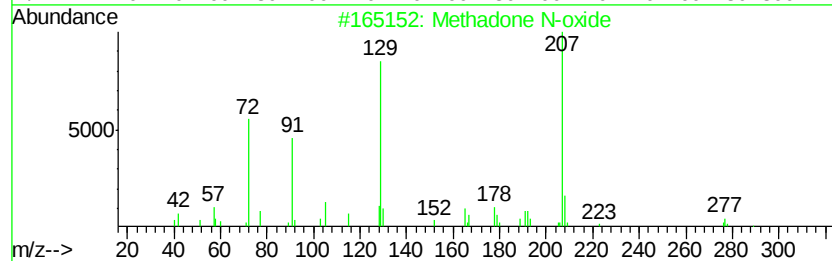
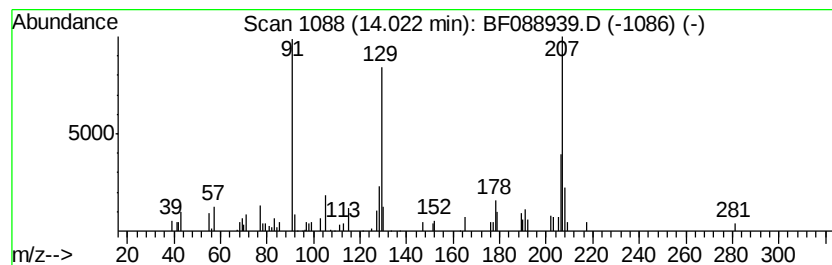
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Methadone N-oxide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	4.21 ng	88147	Chrysene-d12	13.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Methadone N-oxide	325	C21H27NO2	1000120-80-7	64
2		Thiocarbamic acid, N,N-dimethyl,...	311	C19H21NOS	1000192-89-2	59
3		1-Propene, 3-(2-cyclopentenvl)-2...	274	C21H22	1000154-23-3	59
4		5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	50
5		Cyclopentene-1-carboxylic acid, ...	332	C23H24O2	1000159-40-6	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088939.D
Acq On : 13 Jul 2016 00:55
Operator : UM/SJ
Sample : H3935-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C3-04

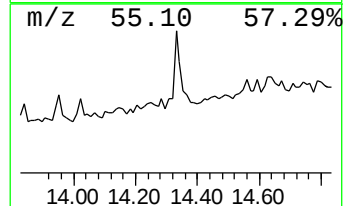
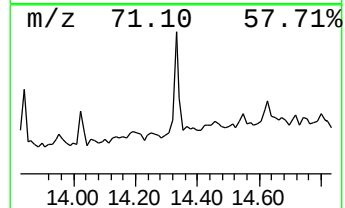
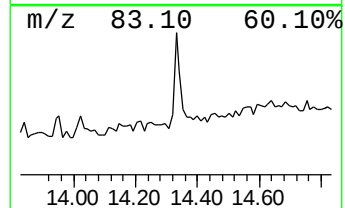
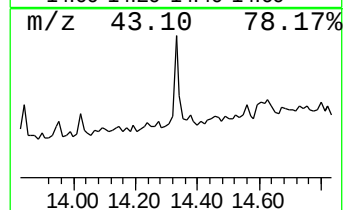
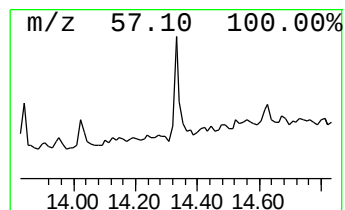
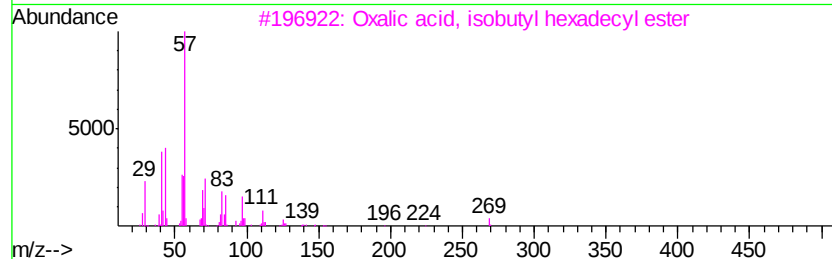
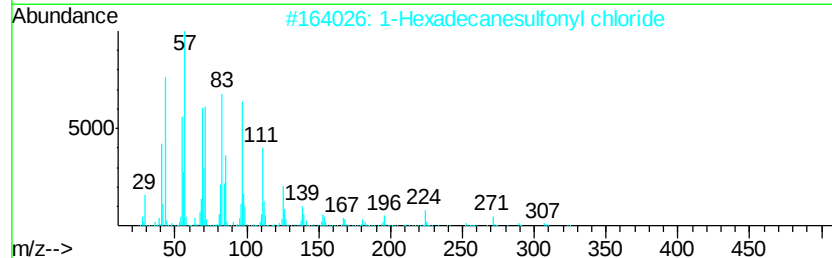
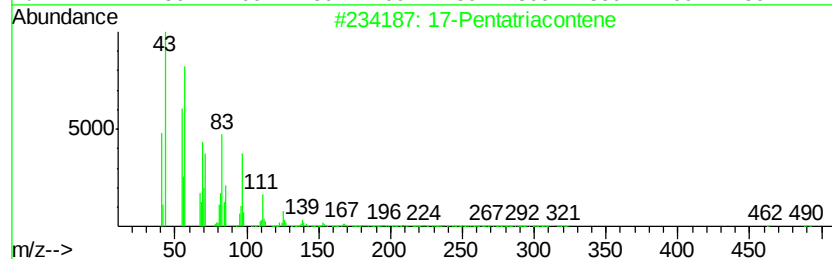
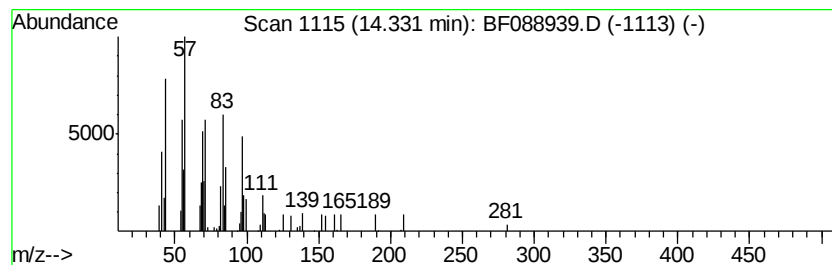
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 17-Pentatriacontene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	2.88 ng	60384	Chrysene-d12	13.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			17-Pentatriacontene	491	C35H70	006971-40-0	83
2			1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	72
3			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	58
4			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	52
5			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
Data File : BF088939.D
Acq On : 13 Jul 2016 00:55
Operator : UM/SJ
Sample : H3935-18
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C3-04

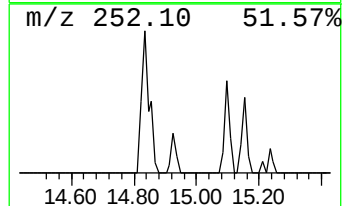
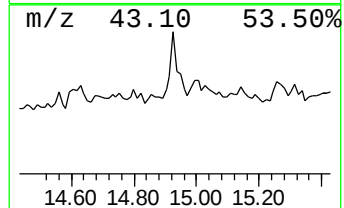
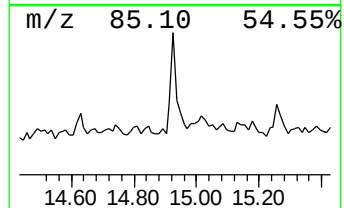
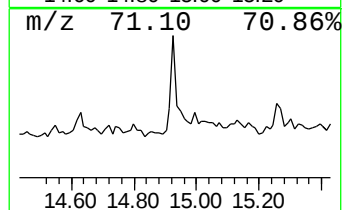
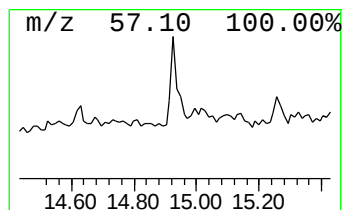
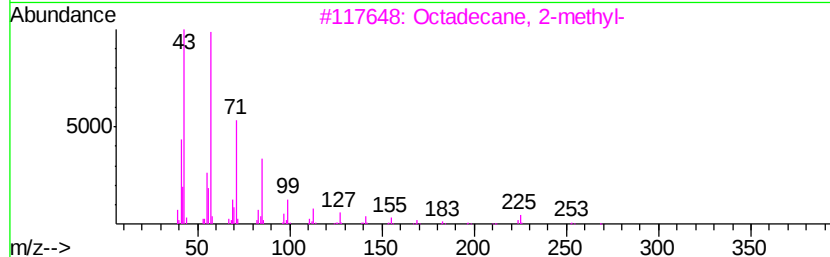
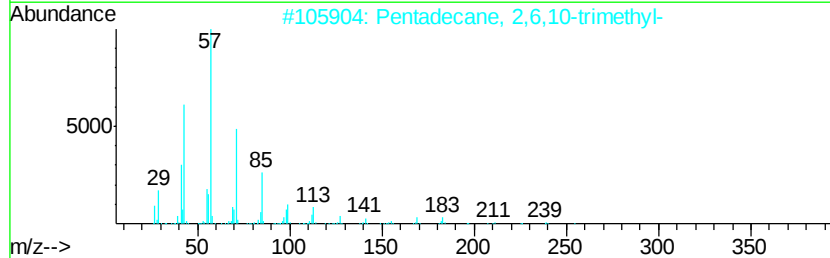
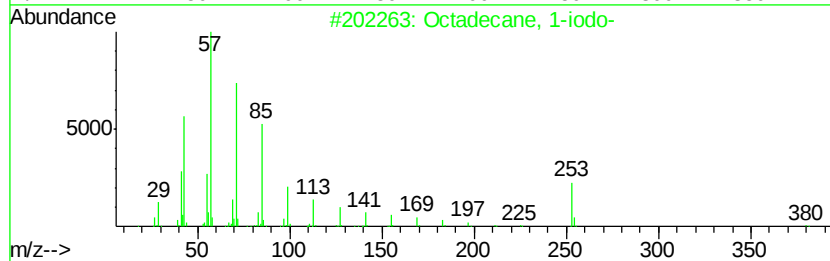
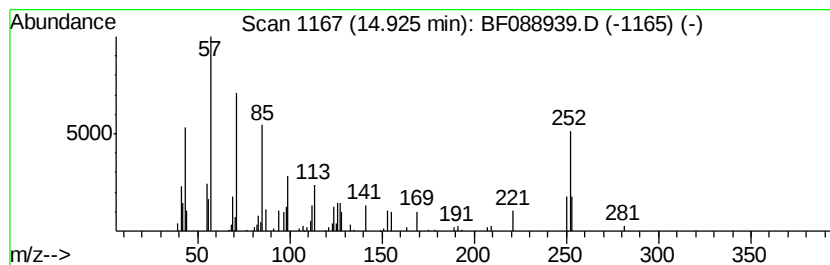
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Octadecane, 1-iodo- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	2.19 ng	33495	Perylene-d12	15.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	50
2		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	43
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	43
4		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	43
5		Tetradecane, 2-methyl-	212	C15H32	001560-95-8	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071216\
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 Sample : H3935-18
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C3-04

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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
Propane, 1,1-dime...	1.88	5.1 ng		91694	1	6.71	359446	20.0
R-(-)-1,2-propane...	2.87	2.9 ng		52777	1	6.71	359446	20.0
2-Pentanone, 4-hy...	4.86	12.1 ng		217798	1	6.71	359446	20.0
unknown6.48	6.48	103.9 ng		1867480	1	6.71	359446	20.0
Sulfurous acid, 2...	10.65	2.2 ng		36215	4	11.21	330564	20.0
n-Dodecyl methacr...	10.97	6.1 ng		101032	4	11.21	330564	20.0
10-Methylnonadecane	11.10	2.8 ng		45805	4	11.21	330564	20.0
Methacrylic acid,...	11.82	2.7 ng		43890	4	11.21	330564	20.0
unknown13.12	13.12	4.8 ng		100952	5	13.84	419244	20.0
unknown13.66	13.66	2.1 ng		44464	5	13.84	419244	20.0
1-Heneicosanol	13.70	5.7 ng		119532	5	13.84	419244	20.0
Methadone N-oxide	14.02	4.2 ng		88147	5	13.84	419244	20.0
17-Pentatriacontene	14.33	2.9 ng		60384	5	13.84	419244	20.0
Octadecane, 1-iodo-	14.93	2.2 ng		33495	6	15.21	306046	20.0