

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
 Data File : BF077385.D
 Acq On : 20 Feb 2015 16:49
 Operator : TP/IZ
 Sample : G1325-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-GW-01

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-BF021915.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	0.990	6	9	12	rVB2	278153	502710	13.13%	1.962%
2	1.641	62	66	69	rBV	413787	657082	17.16%	2.564%
3	1.813	79	81	84	rVB	377019	467165	12.20%	1.823%
4	1.893	86	88	103	rVB2	119612	215702	5.63%	0.842%
5	5.127	369	371	377	rVB	201923	296743	7.75%	1.158%
6	5.619	411	414	416	rBV	765439	770142	20.11%	3.005%
7	6.876	521	524	526	rBV	365460	475456	12.42%	1.855%
8	7.036	535	538	540	rBV	1710744	1530601	39.98%	5.973%
9	7.333	561	564	566	rBV	428590	498354	13.02%	1.945%
10	7.516	577	580	582	rBV	2409622	2137067	55.82%	8.340%
11	8.019	621	624	626	rBV	1859750	1633378	42.66%	6.374%
12	8.899	699	701	704	rBV	563163	706982	18.46%	2.759%
13	9.493	748	753	761	rBV6	19794	58698	1.53%	0.229%
14	9.608	761	763	766	rBV	95796	90369	2.36%	0.353%
15	10.248	817	819	822	rBV	2695544	3638952	95.04%	14.201%
16	10.751	862	863	866	rBV	97122	104630	2.73%	0.408%
17	10.808	866	868	871	rVV	44688	95487	2.49%	0.373%
18	10.911	874	877	879	rVB	52183	59180	1.55%	0.231%
19	11.082	889	892	894	rBV	927651	882684	23.05%	3.445%
20	11.116	894	895	897	rVB	299090	301472	7.87%	1.176%
21	11.231	901	905	906	rBV4	24498	41364	1.08%	0.161%
22	11.265	906	908	911	rVB3	28840	49913	1.30%	0.195%
23	11.425	920	922	926	rVV	212712	376753	9.84%	1.470%
24	11.482	926	927	933	rVV3	48477	98407	2.57%	0.384%
25	11.608	933	938	942	rVV	75953	111270	2.91%	0.434%
26	11.756	949	951	954	rVB3	41149	63131	1.65%	0.246%
27	11.825	954	957	962	rBV2	43337	119747	3.13%	0.467%
28	11.905	962	964	965	rBV	30711	40746	1.06%	0.159%
29	11.985	968	971	973	rBV	252195	269930	7.05%	1.053%
30	12.065	975	978	980	rBV2	1077117	1530590	39.98%	5.973%
31	12.317	997	1000	1002	rBV4	28404	47276	1.23%	0.184%
32	12.442	1008	1011	1013	rVB2	34280	58005	1.51%	0.226%
33	12.511	1013	1017	1019	rBV	46749	89957	2.35%	0.351%
34	12.557	1019	1021	1026	rVV4	24718	84883	2.22%	0.331%

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Integration Parameters: rteint.p
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

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
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35	12.877	1046	1049	1051	rBV	31641	49277	1.29%	0.192%
36	12.922	1051	1053	1055	rVV	803500	901719	23.55%	3.519%
37	13.162	1071	1074	1081	rVB	106612	207771	5.43%	0.811%
38	13.551	1105	1108	1110	rBV2	29364	57003	1.49%	0.222%
39	13.642	1113	1116	1118	rVV	179698	221793	5.79%	0.866%
40	13.688	1118	1120	1126	rVB3	41284	87147	2.28%	0.340%
41	14.625	1200	1202	1208	rVB5	65962	113208	2.96%	0.442%
42	14.922	1226	1228	1231	rVB	3174551	3828784	100.00%	14.942%
43	16.100	1328	1331	1339	rVV	113396	215305	5.62%	0.840%
44	16.225	1339	1342	1344	rVV	749176	955244	24.95%	3.728%
45	17.986	1493	1496	1499	rBV	710098	883009	23.06%	3.446%

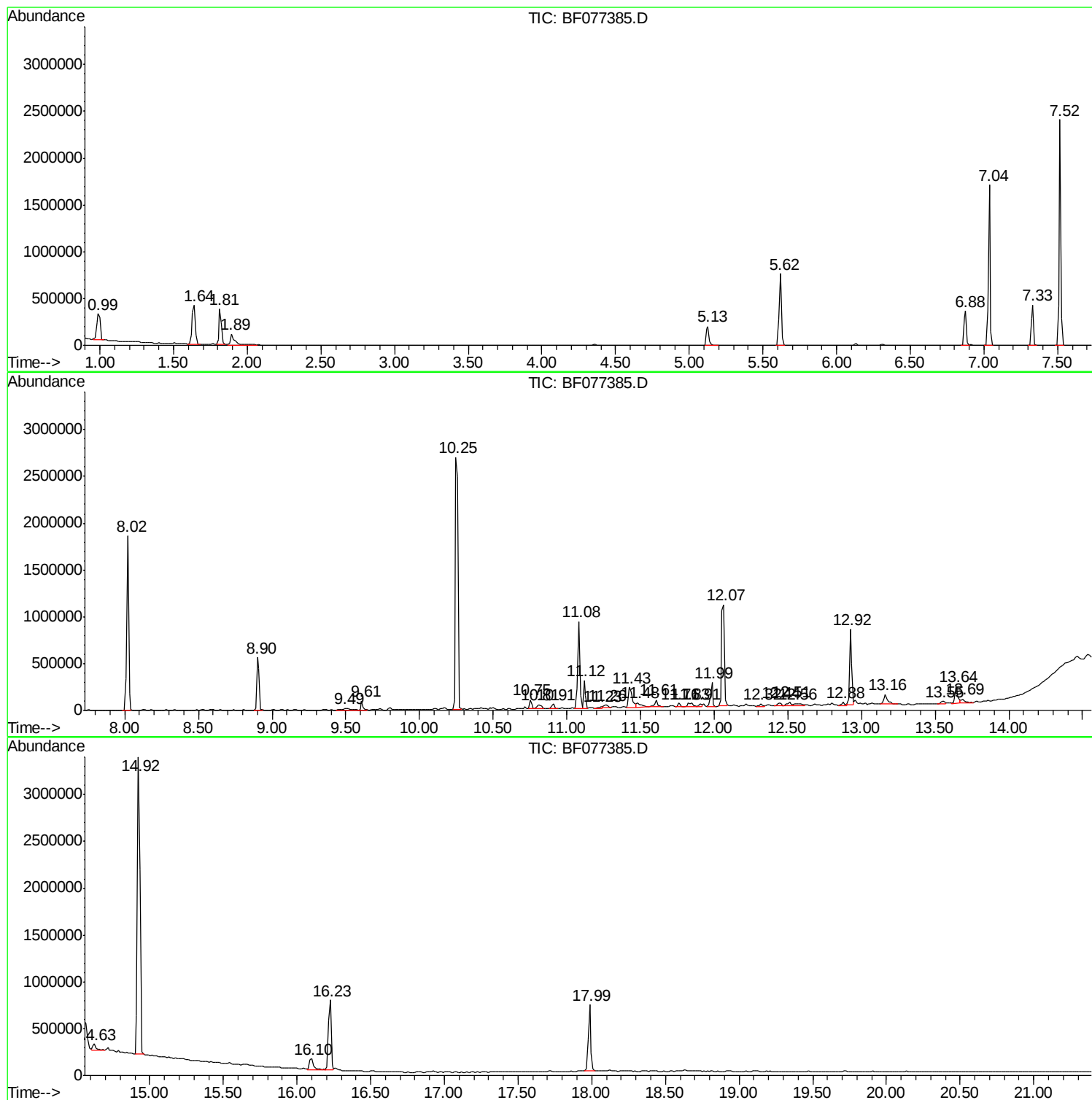
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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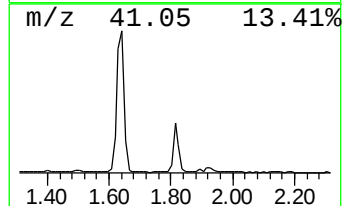
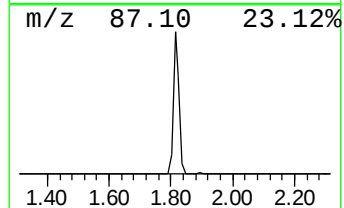
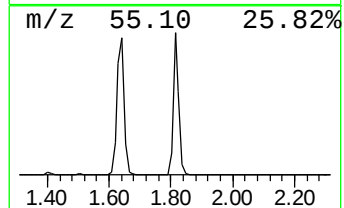
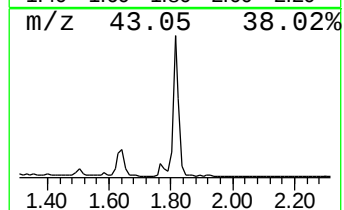
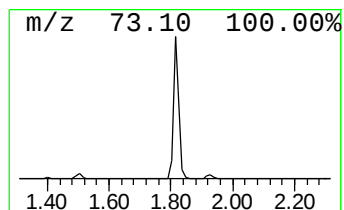
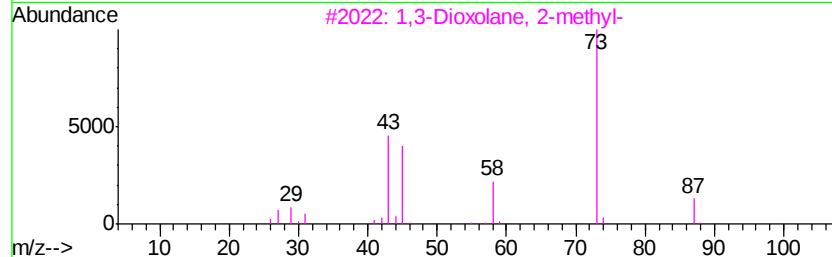
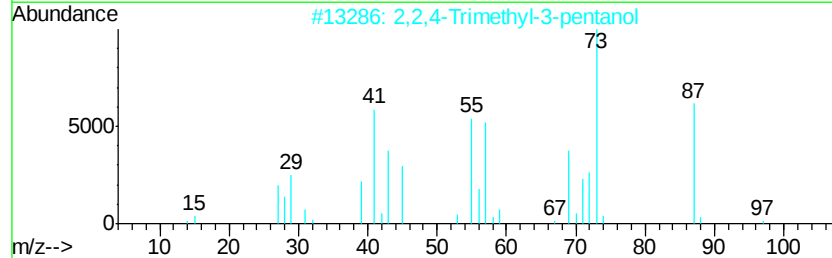
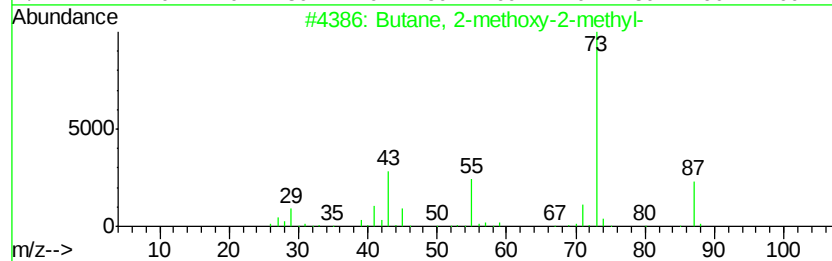
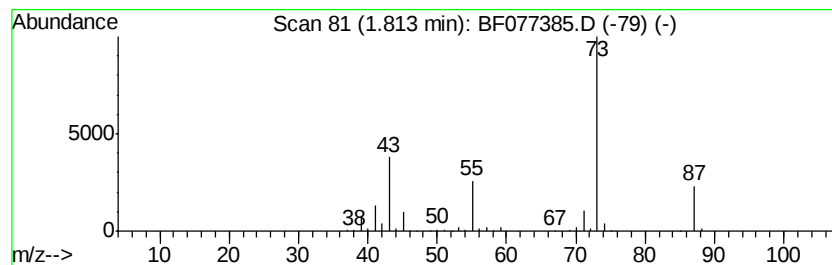
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.81	18.75 ng	467165	1,4-Dichlorobenzene-d4	7.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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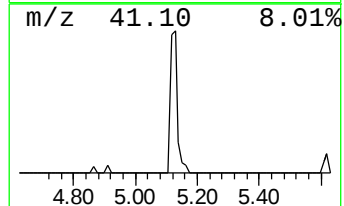
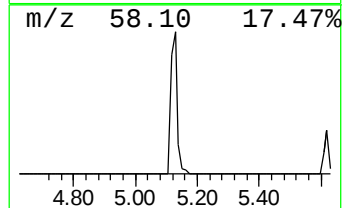
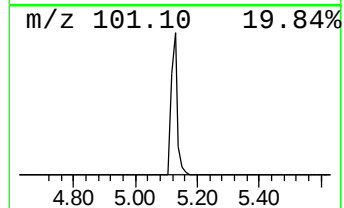
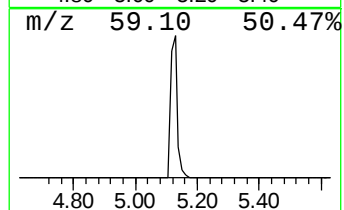
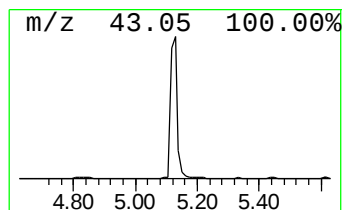
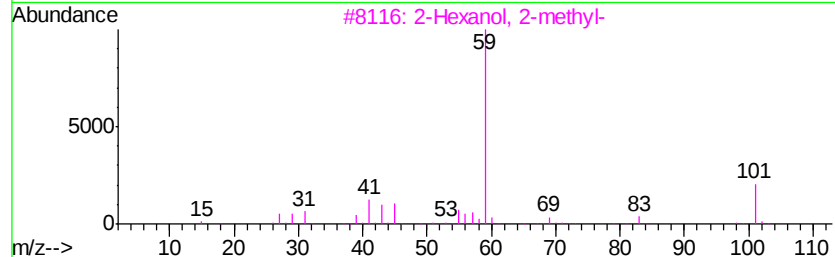
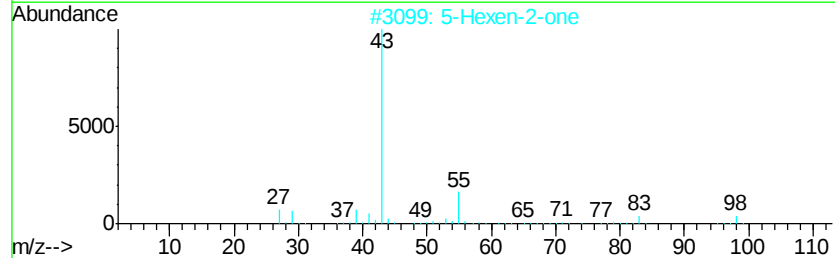
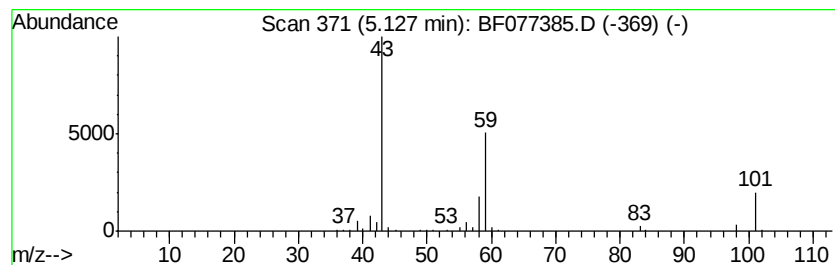
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	11.91 ng	296743	1,4-Dichlorobenzene-d4	7.33

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	5-Hexen-2-one	98	C6H10O	000109-49-9	9	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9	
4	Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9	
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9	



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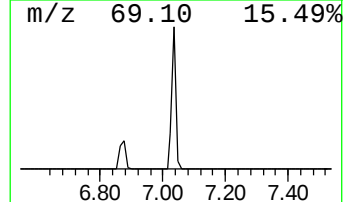
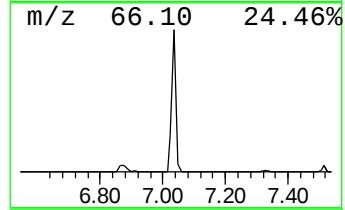
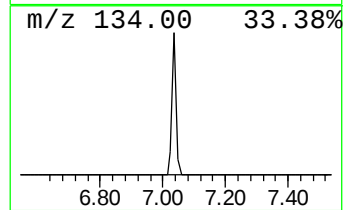
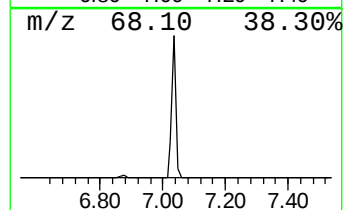
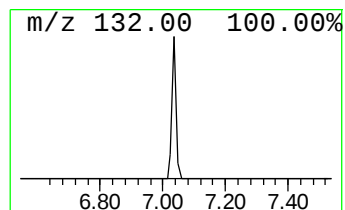
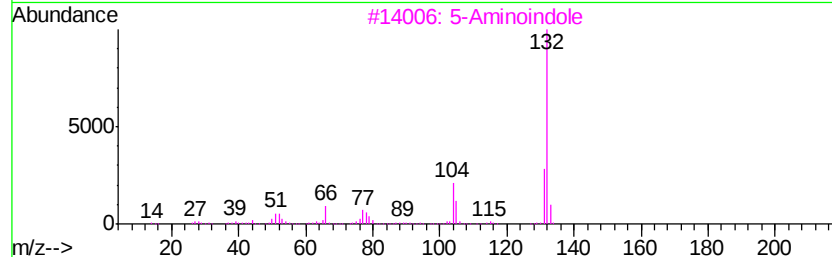
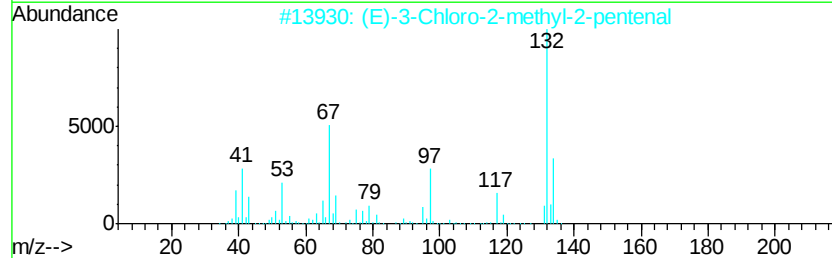
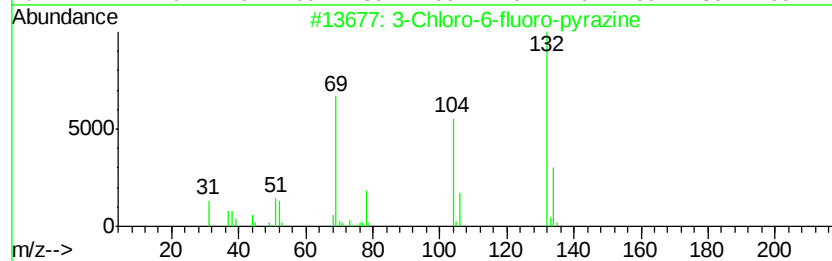
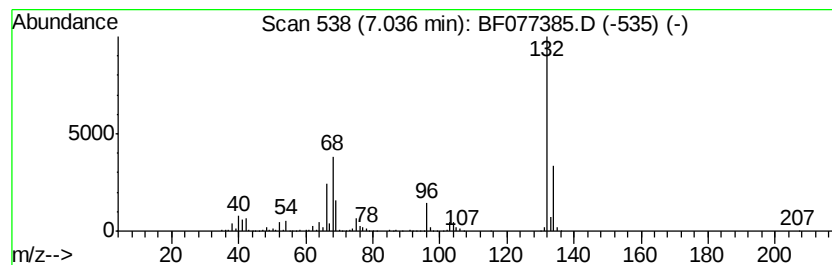
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown7.04 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	61.43 ng	1530600	1,4-Dichlorobenzene-d4	7.33

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3			5-Aminoindole	132	C8H8N2	005192-03-0	12
4			Amphetaminil	250	C17H18N2	017590-01-1	10
5			2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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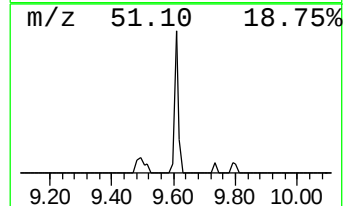
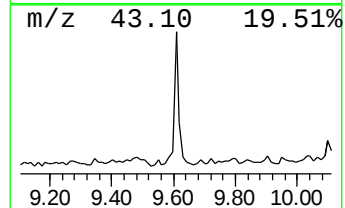
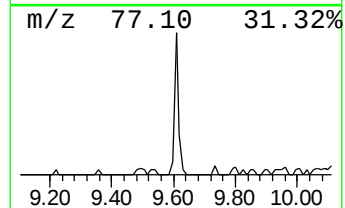
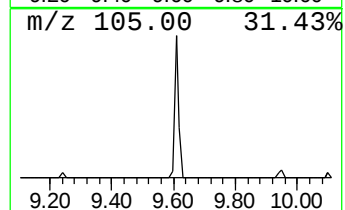
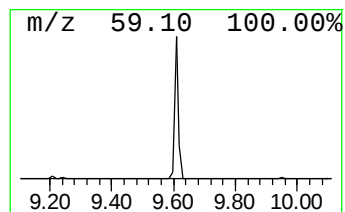
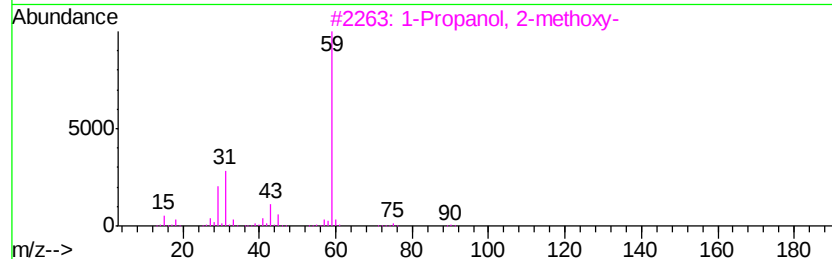
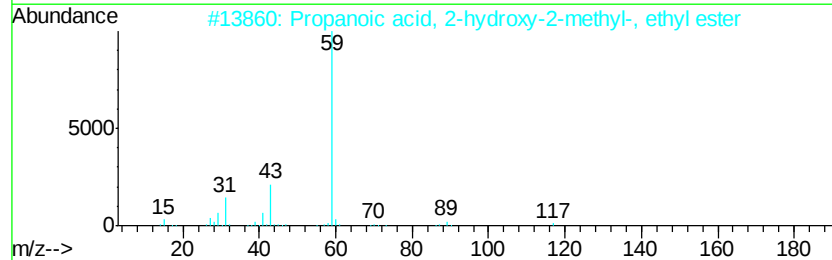
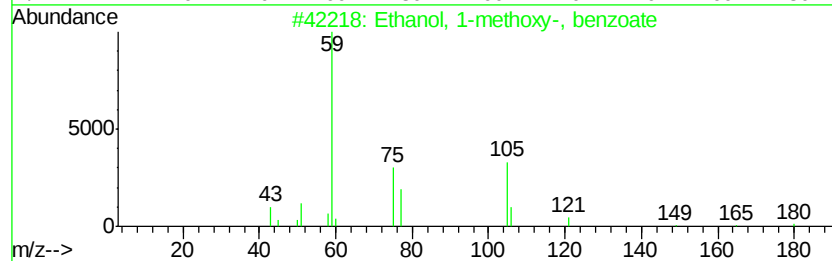
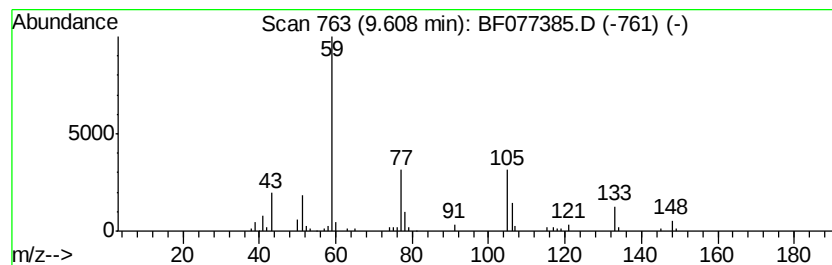
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 unknown9.61 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.61	2.56 ng	90369	Napthalene-d8	8.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 1-methoxy-, benzoate	180	C10H12O3	051835-44-0	45
2		Propanoic acid, 2-hydroxy-2-meth...	132	C6H12O3	000080-55-7	9
3		1-Propanol, 2-methoxy-	90	C4H10O2	001589-47-5	9
4		Ethanone, 2-amino-1-phenyl-	135	C8H9NO	000613-89-8	9
5		Ethane, 1,1-dimethoxy-	90	C4H10O2	000534-15-6	9



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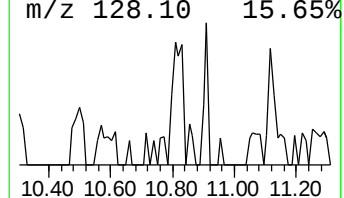
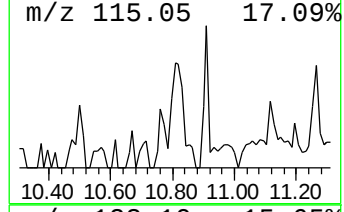
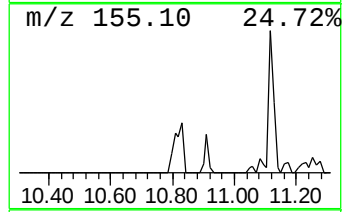
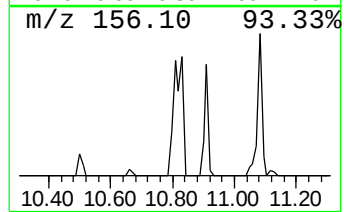
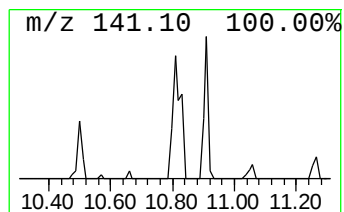
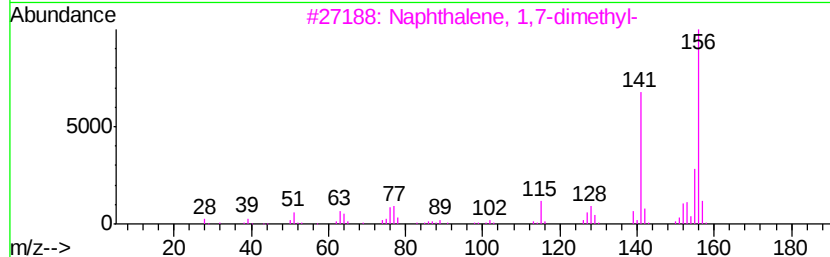
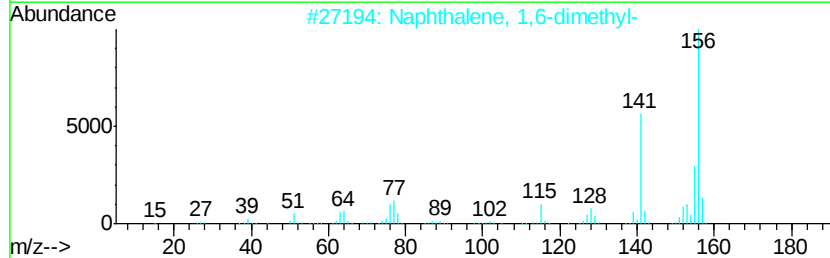
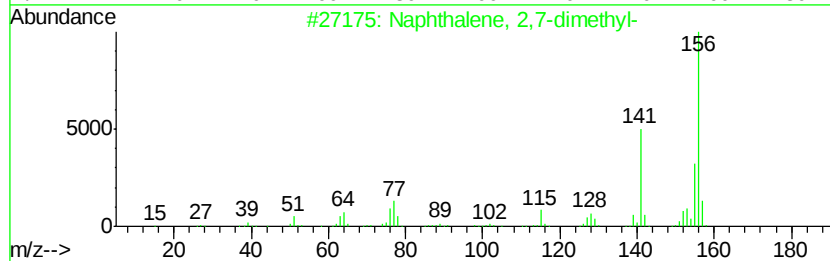
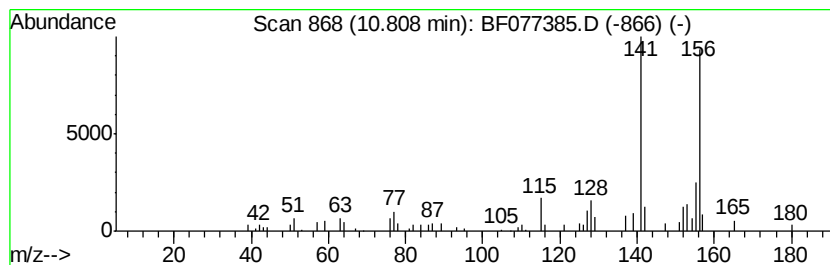
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 2,7-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.81	2.16 ng	95487	Acenaphthene-d10		11.08
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
5	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
Data File : BF077385.D
Acq On : 20 Feb 2015 16:49
Operator : TP/IZ
Sample : G1325-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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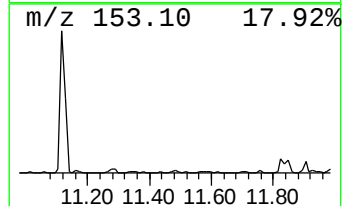
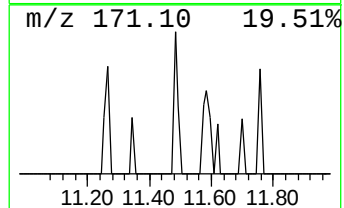
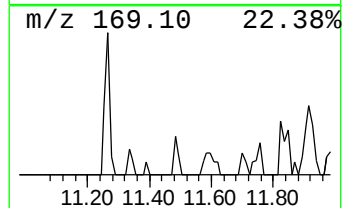
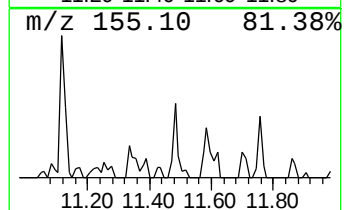
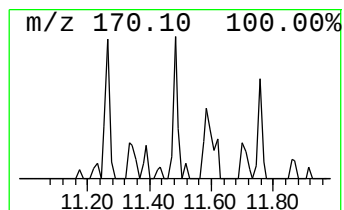
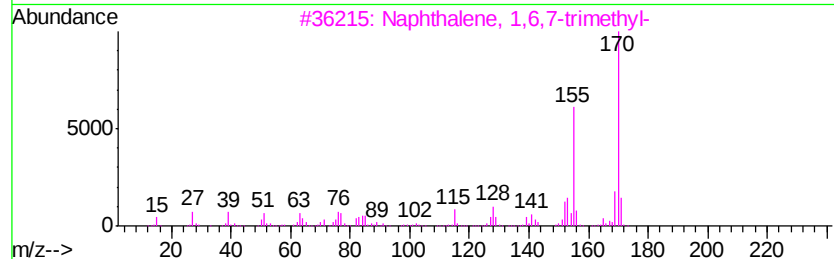
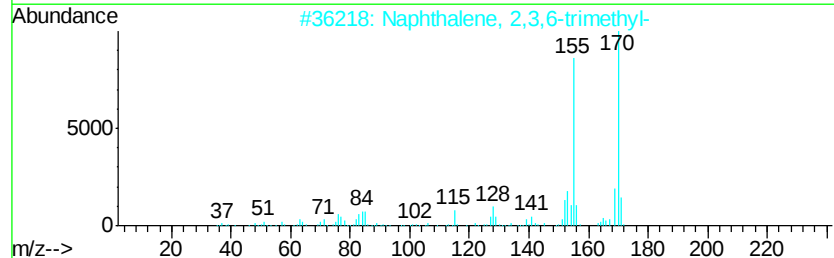
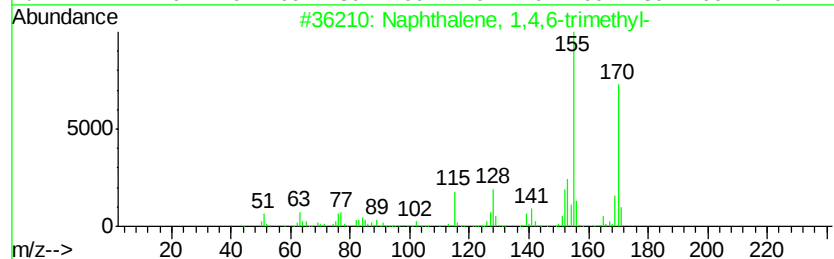
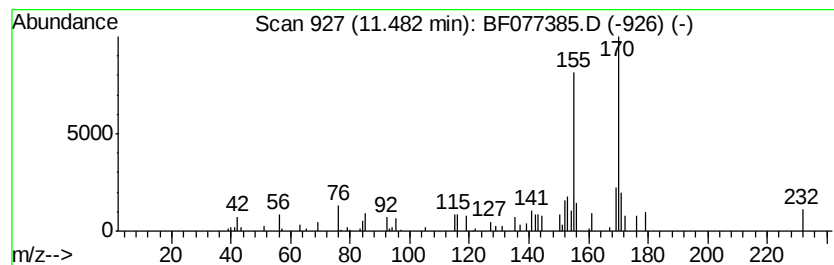
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,4,6-trimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	2.23 ng	98407	Acenaphthene-d10	11.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
2		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	90
4		3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	87
5		1H,3H-Thieno[3,4-c]thiophene, 4,...	170	C8H10S2	015441-54-0	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
Data File : BF077385.D
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Misc :
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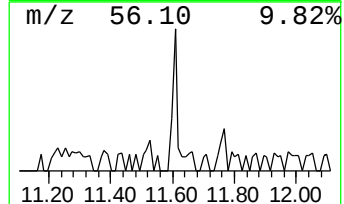
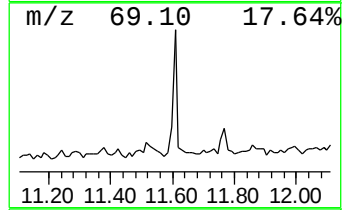
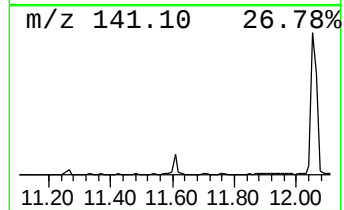
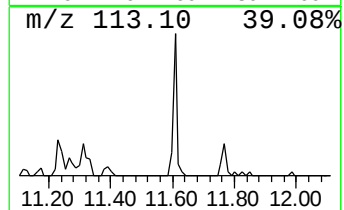
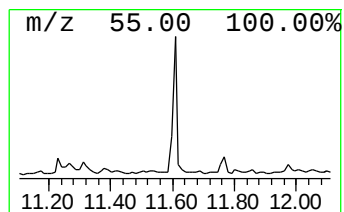
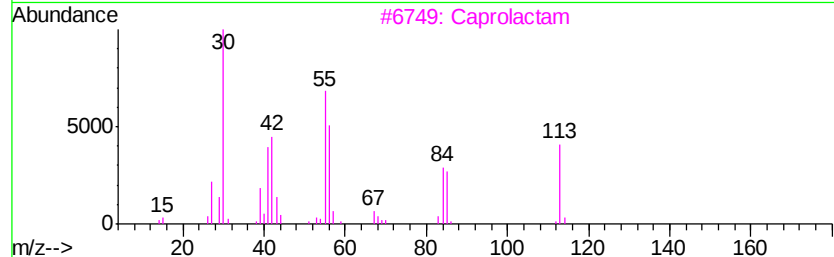
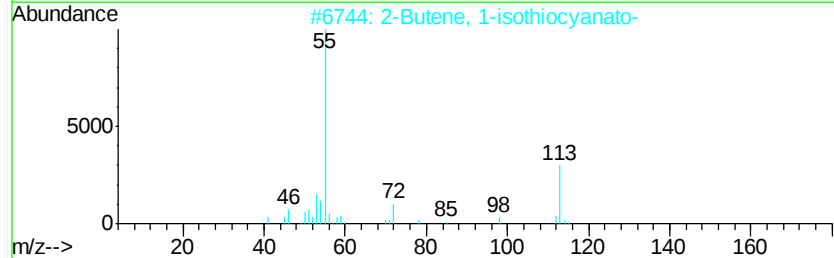
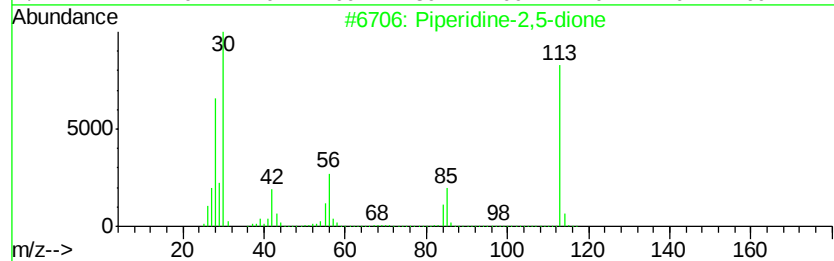
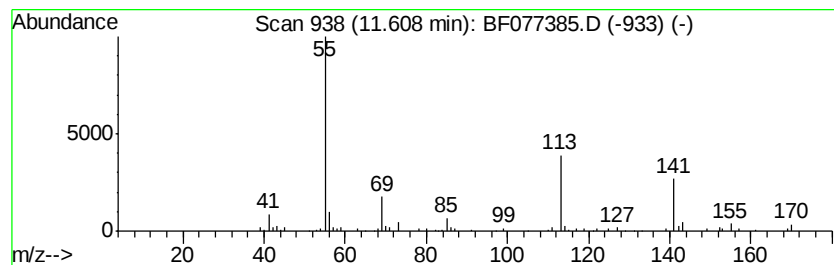
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown11.61 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.61	2.52 ng	111270	Acenaphthene-d10	11.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Piperidine-2,5-dione	113	C5H7N02	052065-78-8	35
2			2-Butene, 1-isothiocyanato-	113	C5H7NS	002253-93-2	9
3			Caprolactam	113	C6H11NO	000105-60-2	9
4			2-Propenoic acid, 2-methylpropyl...	128	C7H12O2	000106-63-8	9
5			2-Ethyl-3-hydroxy-2-methyl-succi...	204	C9H16O5	1000192-42-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
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Sample : G1325-01
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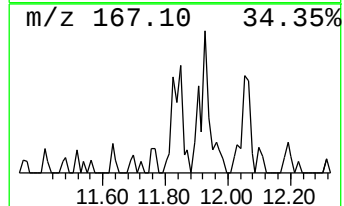
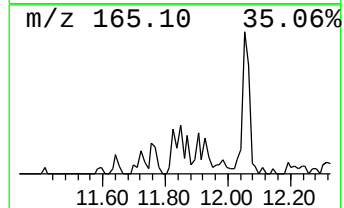
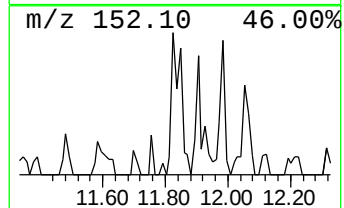
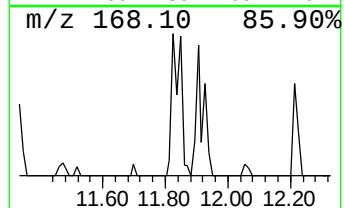
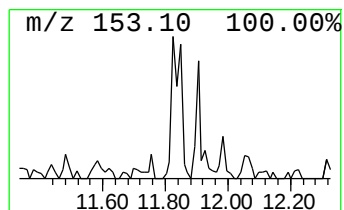
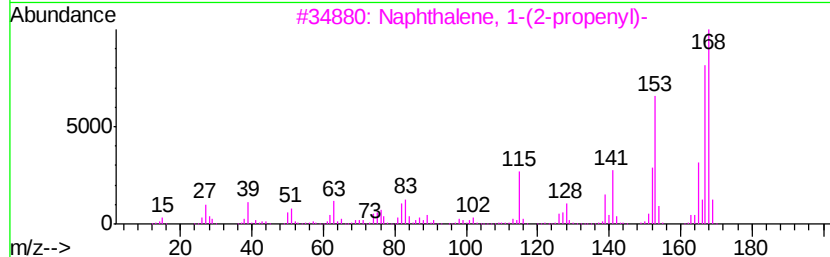
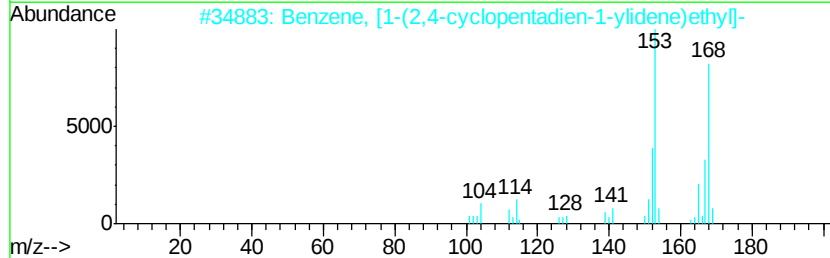
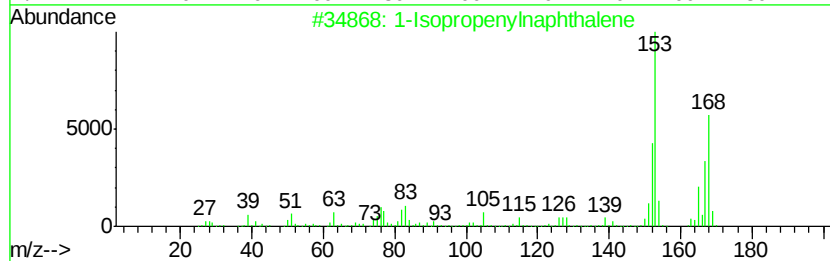
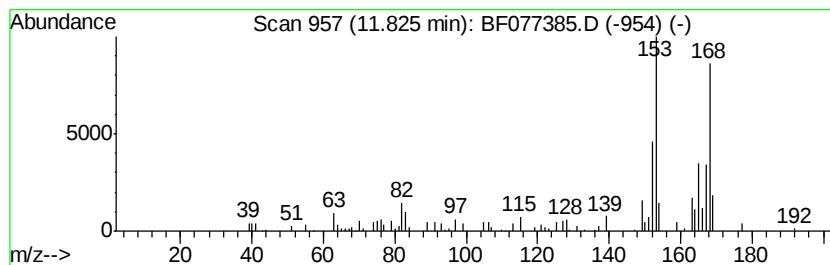
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Isopropenylnaphthalene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.		
11.83	2.71 ng	119747	Acenaphthene-d10		11.08		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Isopropenylnaphthalene	168	C13H12	001855-47-6	87
2			Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	74
3			Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	64
4			1-Methoxy-2-methyl-4-(methylthio...	168	C9H12OS	050390-78-8	47
5			1,1,4,4-Tetramethyl-1,4-disilacy...	168	C8H16Si2	020627-53-6	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
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Acq On : 20 Feb 2015 16:49
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Sample : G1325-01
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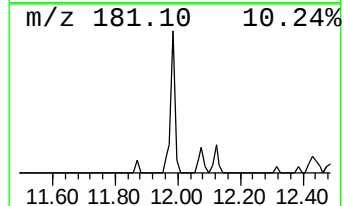
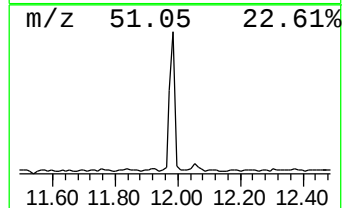
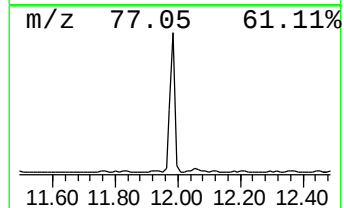
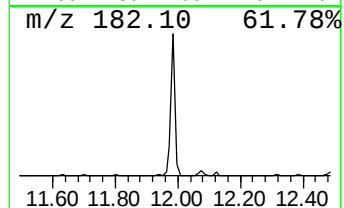
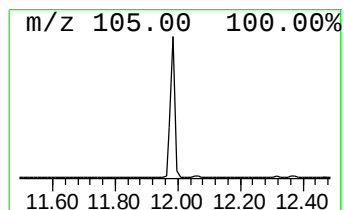
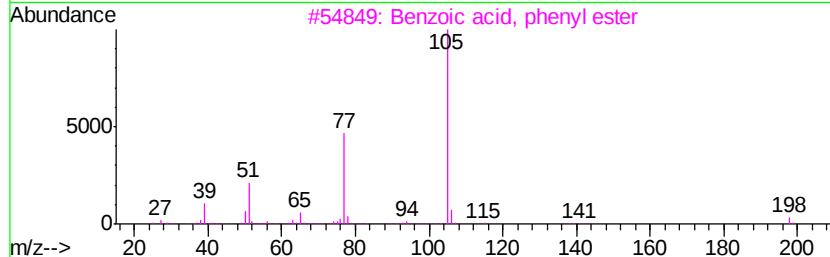
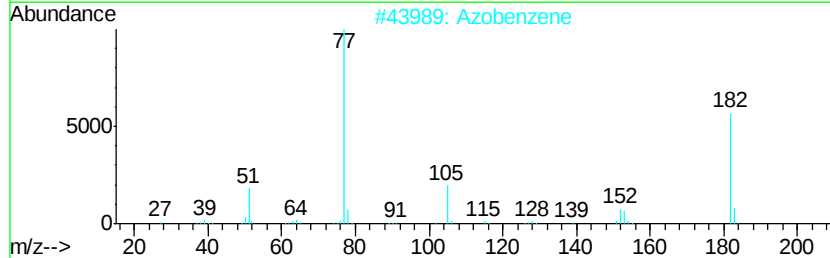
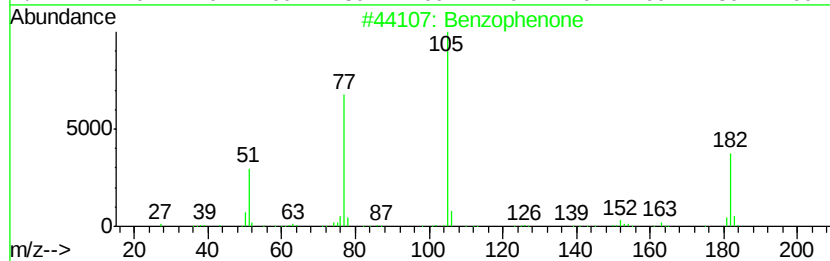
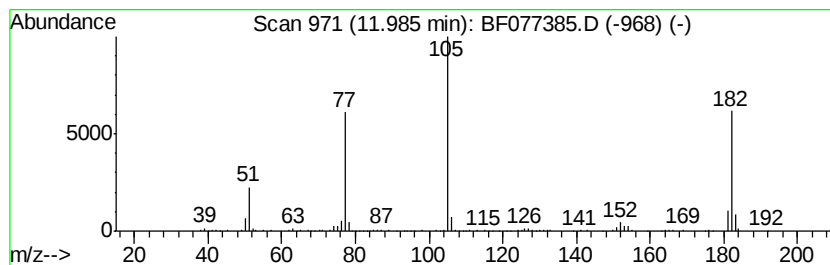
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	6.12 ng	269930	Acenaphthene-d10	11.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzophenone	182 C13H10O	000119-61-9 96
2		Azobenzene	182 C12H10N2	000103-33-3 59
3		Benzoic acid, phenyl ester	198 C13H10O2	000093-99-2 47
4		Benzenebutanoic acid, .gamma.-oxo-	178 C10H10O3	002051-95-8 47
5		Benzoyl bromide	184 C7H5BrO	000618-32-6 47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
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Acq On : 20 Feb 2015 16:49
Operator : TP/IZ
Sample : G1325-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

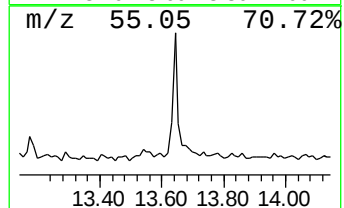
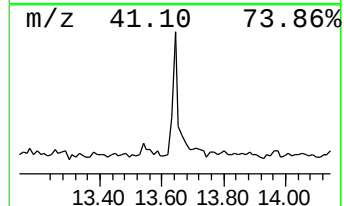
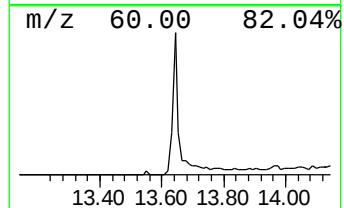
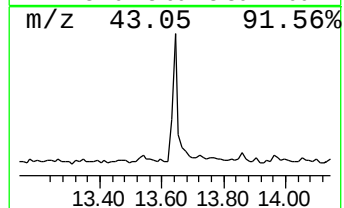
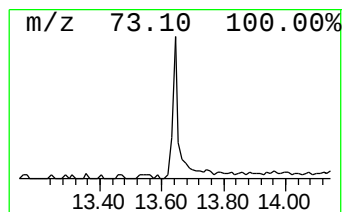
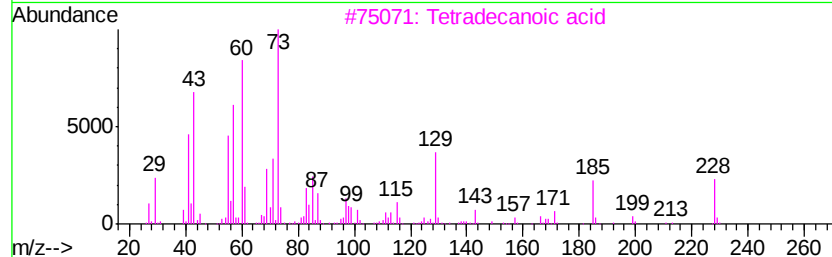
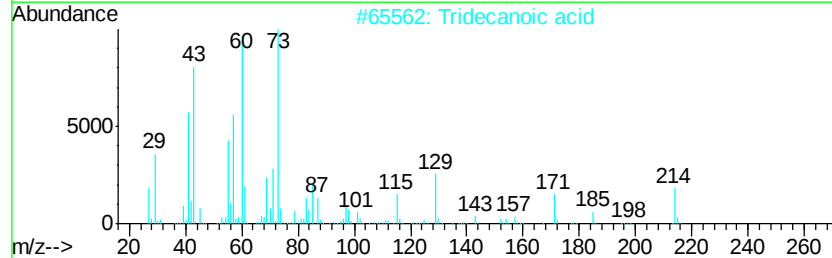
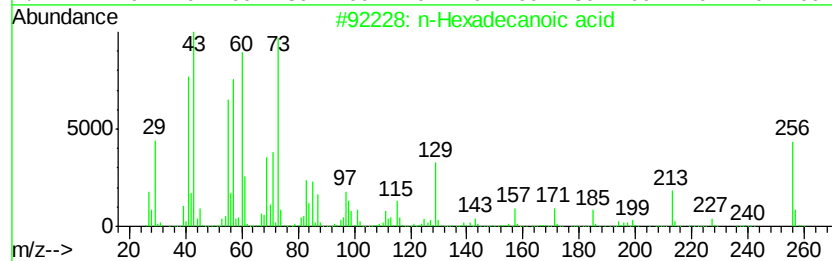
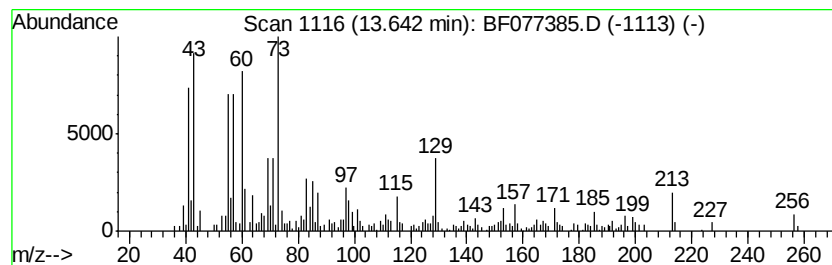
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.64	4.92 ng	221793	Phenanthrene-d10		12.92	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4		Dodecanoic acid	200	C12H24O2	000143-07-7	64
5		n-Decanoic acid	172	C10H20O2	000334-48-5	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
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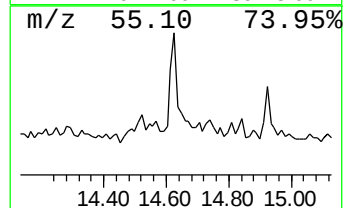
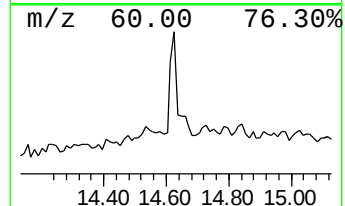
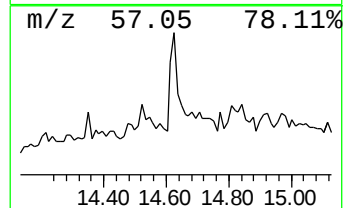
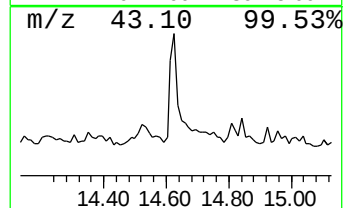
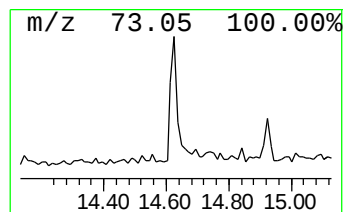
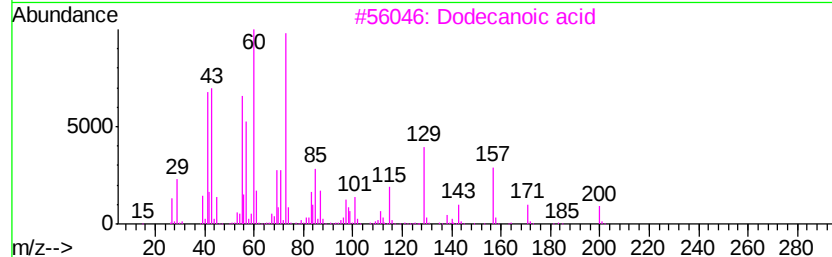
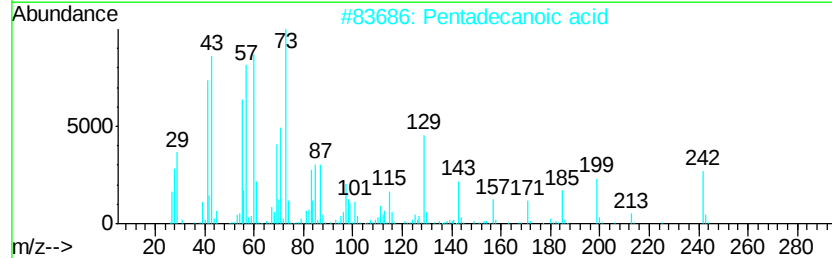
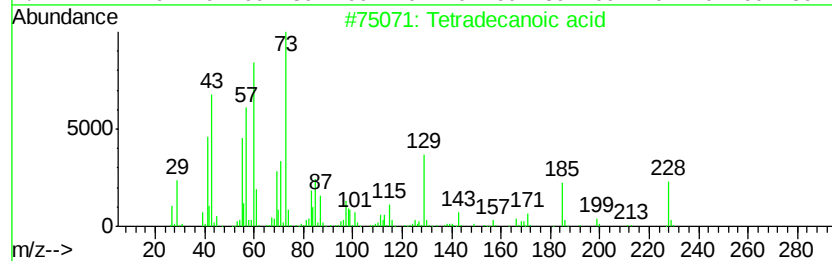
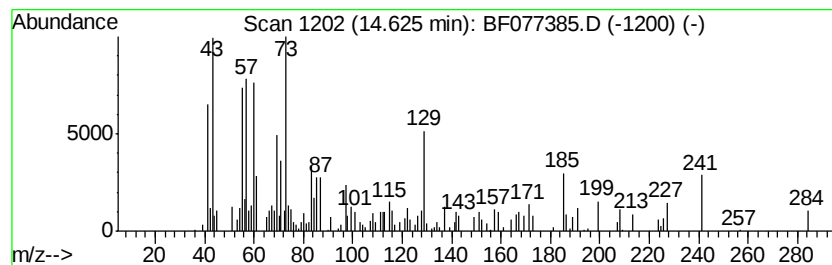
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tetradecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.63	2.37 ng	113208	Chrysene-d12		16.23
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	52
3	Dodecanoic acid	200	C12H24O2	000143-07-7	47
4	n-Decanoic acid	172	C10H20O2	000334-48-5	47
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
Data File : BF077385.D
Acq On : 20 Feb 2015 16:49
Operator : TP/IZ
Sample : G1325-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

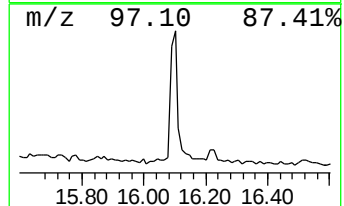
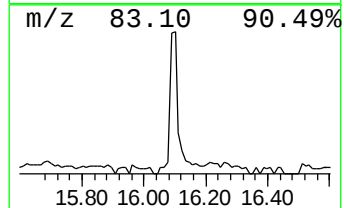
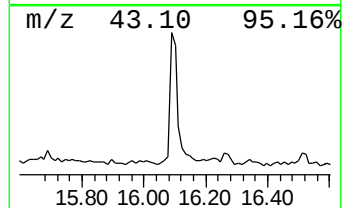
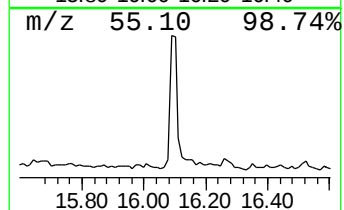
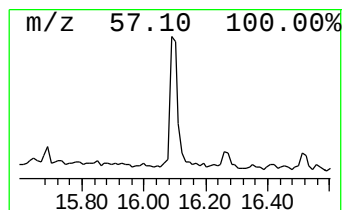
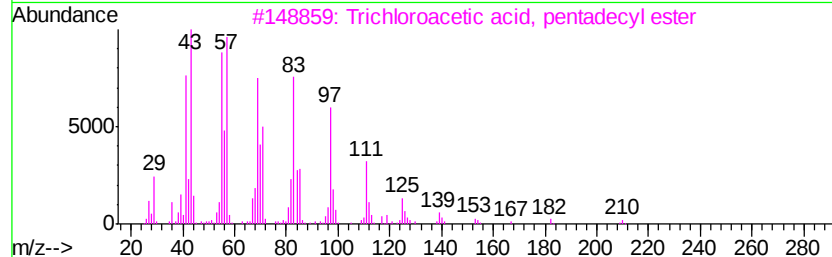
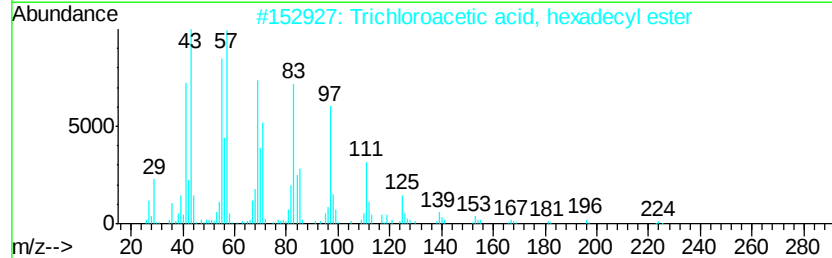
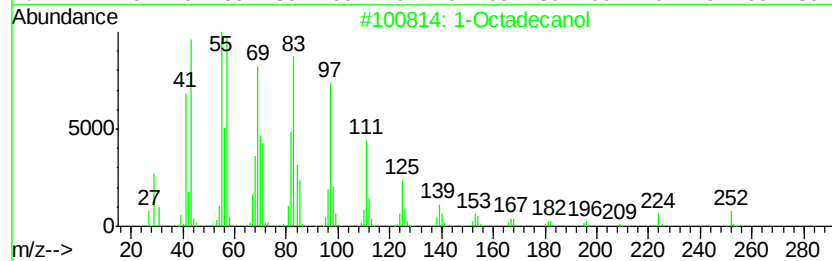
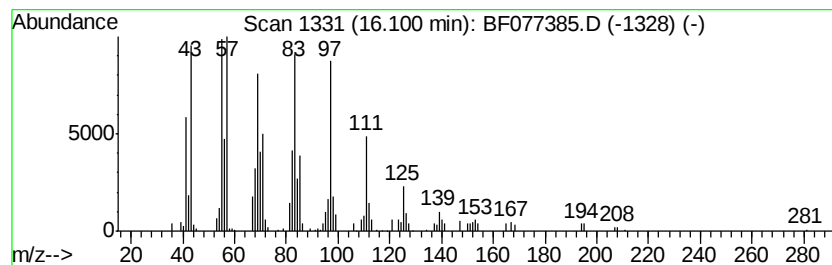
Instrument :
BNA_F
ClientSampleId :
WC-GW-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Octadecanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	4.51 ng	215305	Chrysene-d12	16.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Octadecanol		270 C18H38O	000112-92-5 95
2	Trichloroacetic acid, hexadecyl ...		386 C18H33Cl3O2	074339-54-1 91
3	Trichloroacetic acid, pentadecyl...		372 C17H31Cl3O2	074339-53-0 91
4	Trifluoroacetic acid, n-octadecy...		366 C20H37F3O2	079392-43-1 91
5	1-Heptadecanol		256 C17H36O	001454-85-9 90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022015\
Data File : BF077385.D
Acq On : 20 Feb 2015 16:49
Operator : TP/IZ
Sample : G1325-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-GW-01

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.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
Butane, 2-methoxy...	1.81	18.8	ng	467165	1	7.33	498354	20.0
2-Pentanone, 4-hy...	5.13	11.9	ng	296743	1	7.33	498354	20.0
unknown7.04	7.04	61.4	ng	1530600	1	7.33	498354	20.0
unknown9.61	9.61	2.6	ng	90369	2	8.90	706982	20.0
Naphthalene, 2,7-...	10.81	2.2	ng	95487	3	11.08	882684	20.0
Naphthalene, 1,4,...	11.48	2.2	ng	98407	3	11.08	882684	20.0
unknown11.61	11.61	2.5	ng	111270	3	11.08	882684	20.0
1-Isopropenyl naph...	11.83	2.7	ng	119747	3	11.08	882684	20.0
Benzophenone	11.99	6.1	ng	269930	3	11.08	882684	20.0
n-Hexadecanoic acid	13.64	4.9	ng	221793	4	12.92	901719	20.0
Tetradecanoic acid	14.63	2.4	ng	113208	5	16.23	955244	20.0
1-Octadecanol	16.10	4.5	ng	215305	5	16.23	955244	20.0