

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
 Data File : BG017476.D
 Acq On : 5 Jun 2015 16:23
 Operator : TP/IZ
 Sample : G2501-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-5-6-2-15

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_G\METHODS\8270-BG060315.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.767	8	13	29	rVB	258044	379328	16.97%	3.250%
2	5.229	426	432	444	rBV	173425	298683	13.37%	2.559%
3	5.622	495	499	506	rBV2	15720	24022	1.07%	0.206%
4	6.850	703	708	718	rVB	104389	177568	7.95%	1.522%
5	7.174	756	763	775	rBV	383529	636707	28.49%	5.456%
6	7.268	775	779	785	rVB3	14503	23828	1.07%	0.204%
7	7.414	798	804	813	rVB	48657	73673	3.30%	0.631%
8	7.620	834	839	846	rBV2	100488	163237	7.30%	1.399%
9	7.943	887	894	900	rBV	377843	613276	27.44%	5.255%
10	8.789	1030	1038	1048	rBV	308595	511875	22.91%	4.386%
11	10.423	1307	1316	1321	rBV	112617	192113	8.60%	1.646%
12	11.615	1511	1519	1523	rBV	67150	95054	4.25%	0.814%
13	11.662	1523	1527	1529	rBV2	50755	75983	3.40%	0.651%
14	11.727	1534	1538	1546	rVB2	60912	95428	4.27%	0.818%
15	11.880	1557	1564	1568	rBV	413450	669898	29.98%	5.740%
16	11.921	1568	1571	1576	rVB	350531	504407	22.57%	4.322%
17	12.150	1600	1610	1618	rBV	215860	384954	17.23%	3.299%
18	12.908	1732	1739	1746	rBV	824787	1139157	50.98%	9.761%
19	13.149	1776	1780	1787	rVB	39460	57698	2.58%	0.494%
20	14.018	1923	1928	1933	rVB3	16030	25900	1.16%	0.222%
21	14.295	1967	1975	1985	rVB2	185321	283828	12.70%	2.432%
22	15.799	2223	2231	2238	rBV	804126	1122082	50.21%	9.615%
23	17.044	2438	2443	2450	rBV2	247632	356068	15.93%	3.051%
24	17.961	2593	2599	2607	rBV	133579	208971	9.35%	1.791%
25	19.242	2810	2817	2824	rBV	117742	184259	8.25%	1.579%
26	19.688	2887	2893	2898	rBV	1895150	2234726	100.00%	19.149%
27	20.487	3026	3029	3035	rBV7	16417	24403	1.09%	0.209%
28	20.957	3106	3109	3118	rBV6	22737	31371	1.40%	0.269%
29	21.157	3139	3143	3149	rBV	80165	96178	4.30%	0.824%
30	21.239	3152	3157	3165	rVB	398214	516308	23.10%	4.424%
31	23.484	3532	3539	3545	rVB	252008	469369	21.00%	4.022%

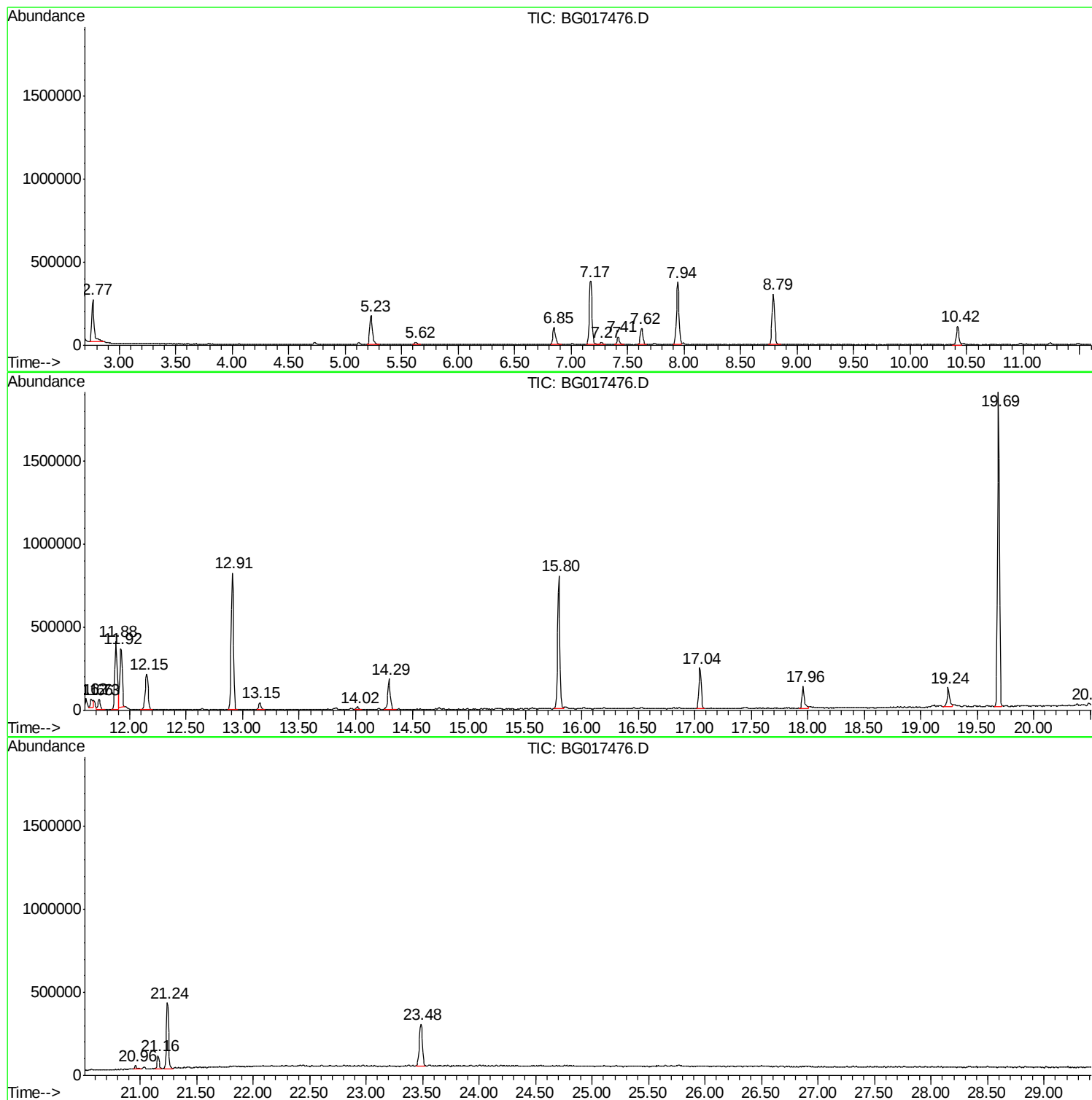
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Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
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Operator : TP/IZ
Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-5-6-2-15

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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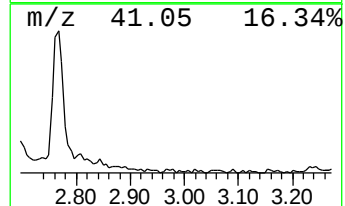
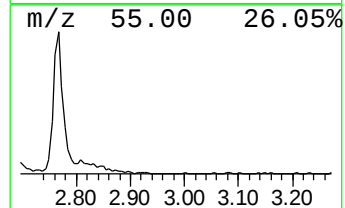
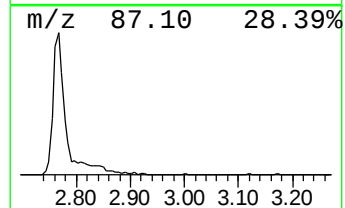
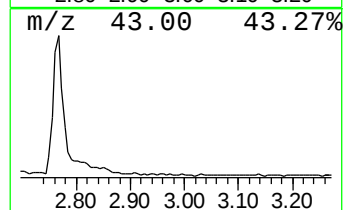
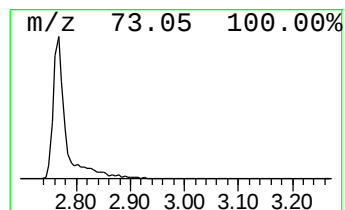
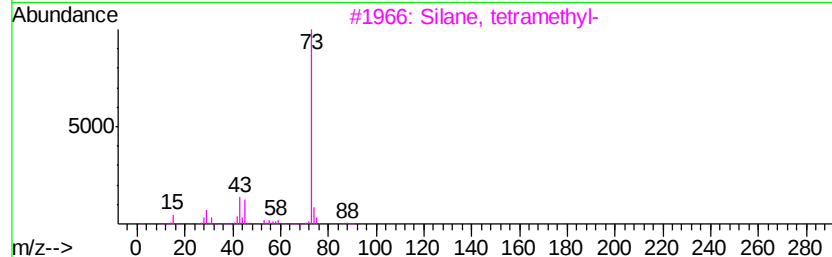
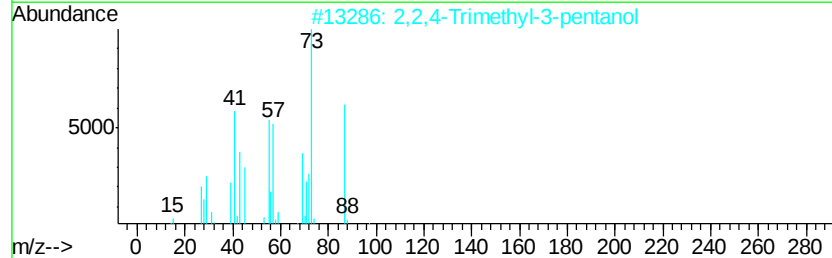
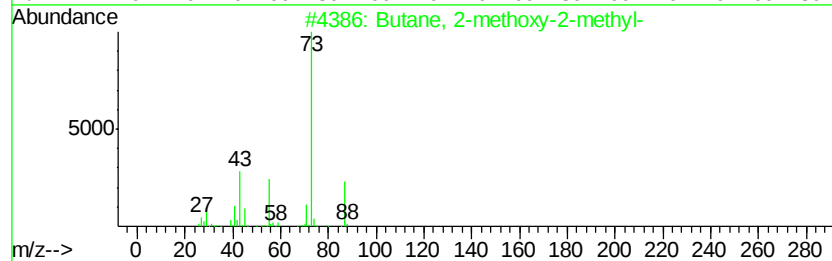
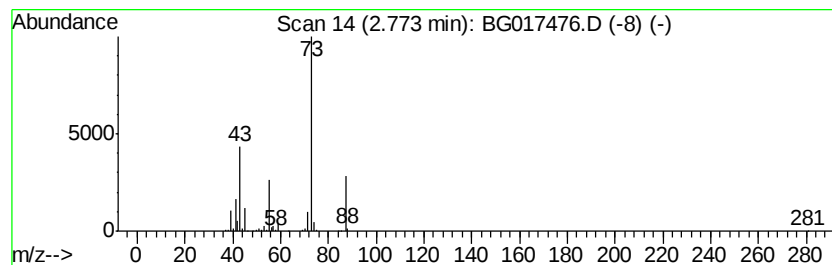
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.77	46.48 ng	379328	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	36
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	12
5		1-Propene-1-thiol	74	C3H6S	000925-89-3	9



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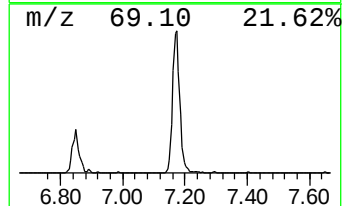
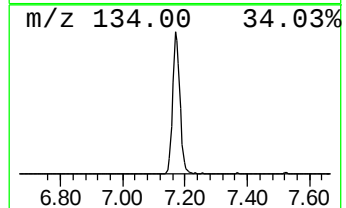
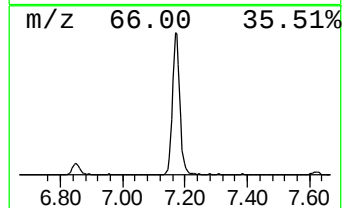
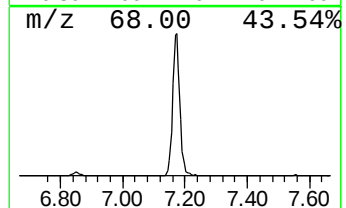
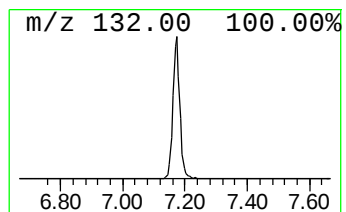
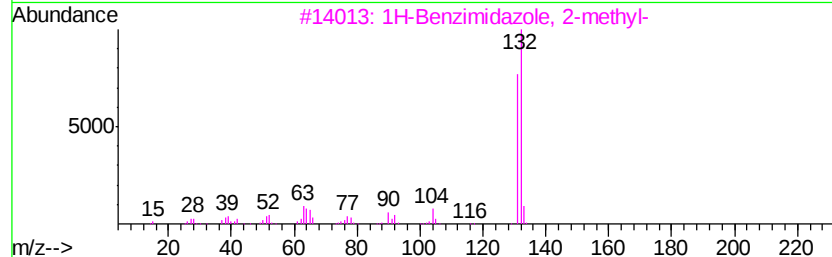
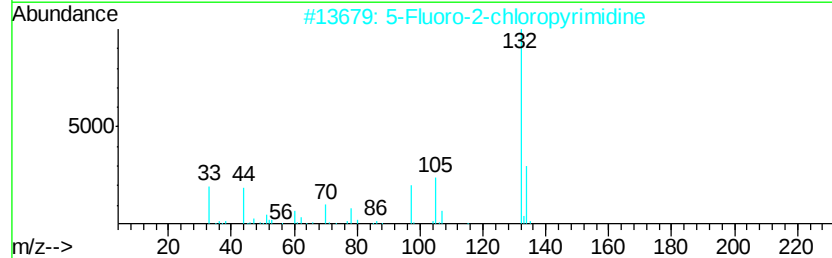
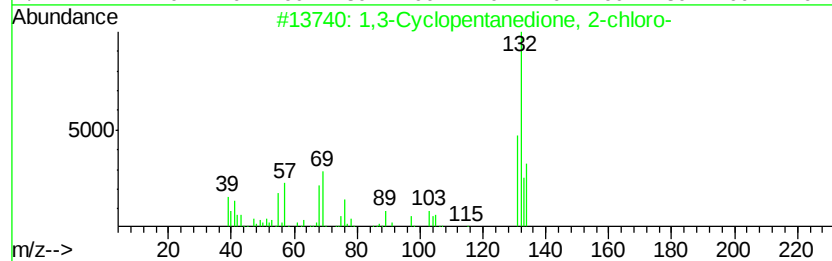
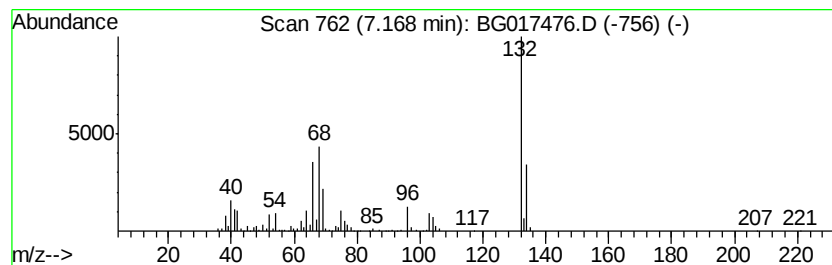
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown7.17 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	78.01 ng	636707	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	32
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	12



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Instrument :
BNA_G
ClientSampled :
MW-5-6-2-15

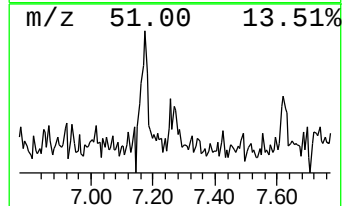
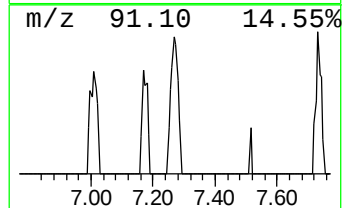
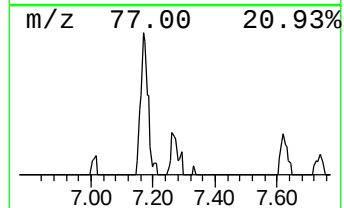
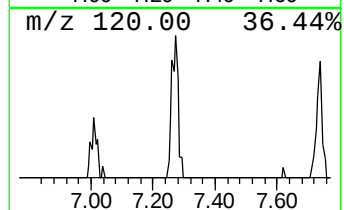
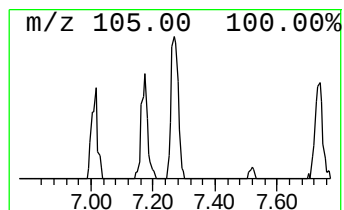
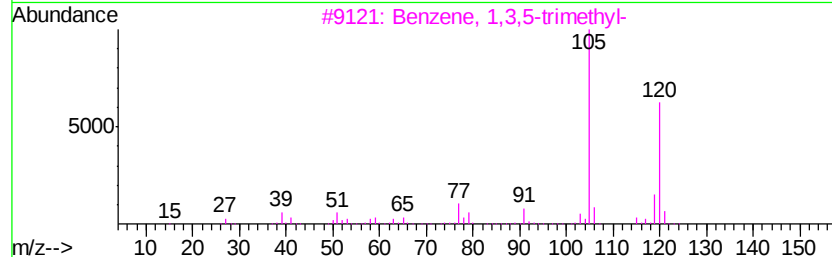
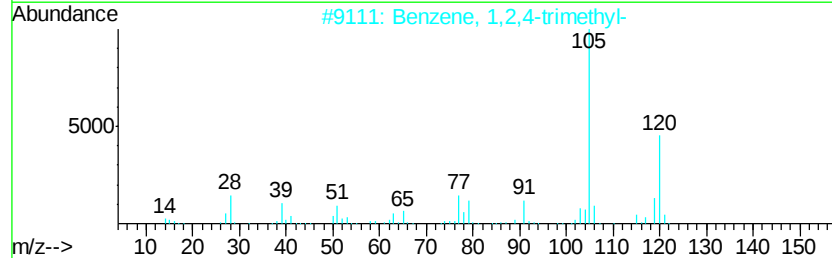
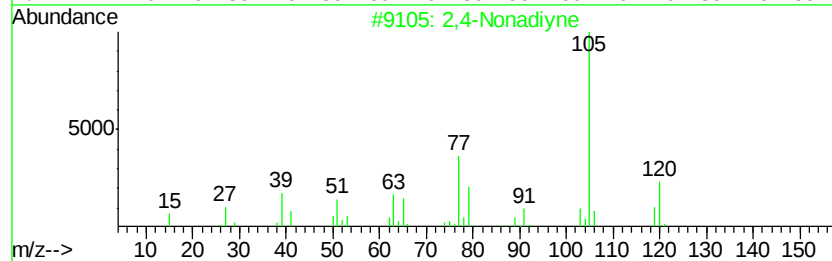
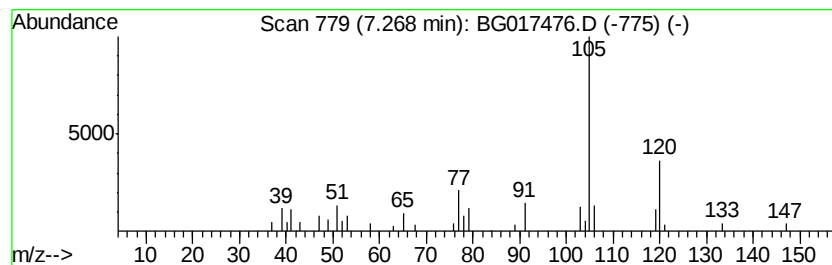
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 2,4-Nonadiyne Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.27	2.92 ng	23828	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Nonadiyne	120	C9H12	063621-15-8	81
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	81
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	81
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	76



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ALS Vial : 5 Sample Multiplier: 1

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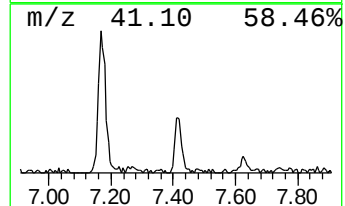
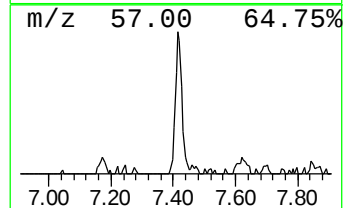
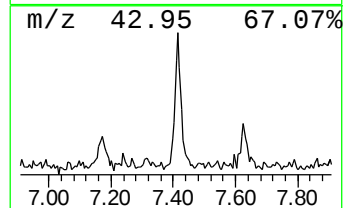
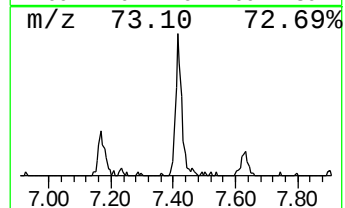
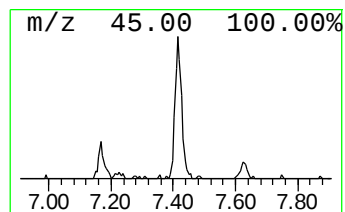
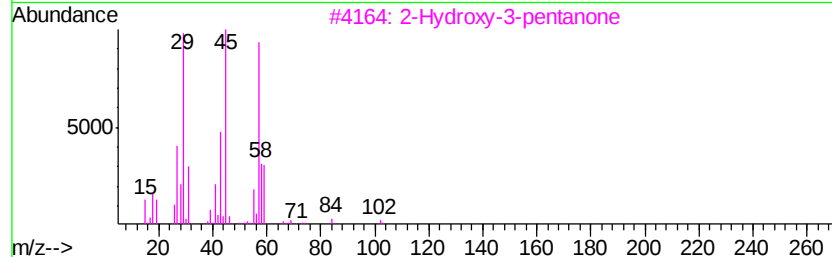
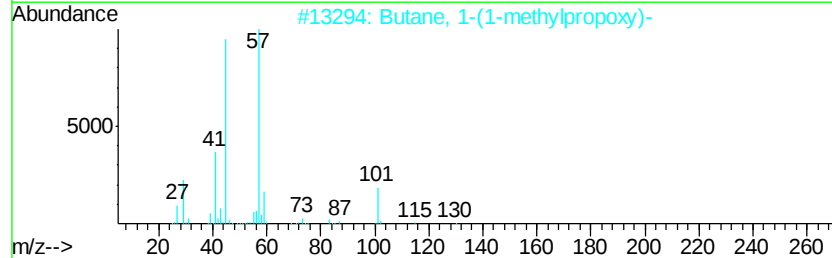
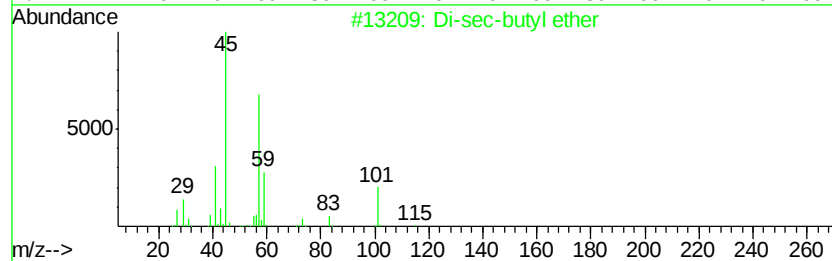
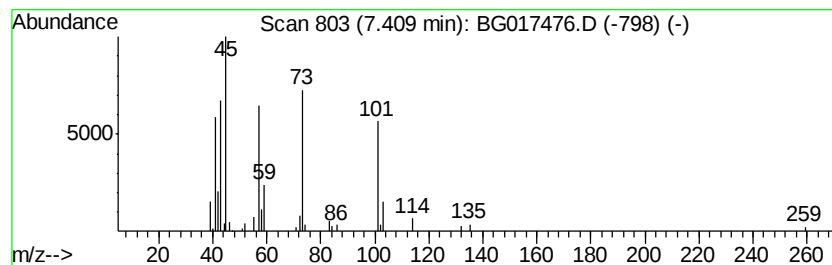
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Di-sec-butyl ether Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.41	9.03 ng	73673	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Di-sec-butyl ether	130	C8H18O	006863-58-7	53
2		Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50
3		2-Hydroxy-3-pentanone	102	C5H10O2	005704-20-1	43
4		3,4,5-Trimethyl-4-heptanol	158	C10H22O	060836-08-0	38
5		Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	36



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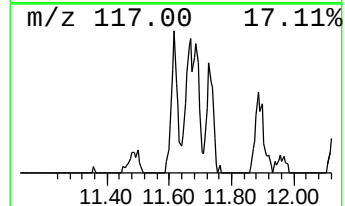
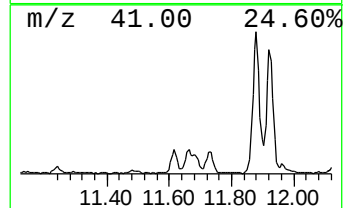
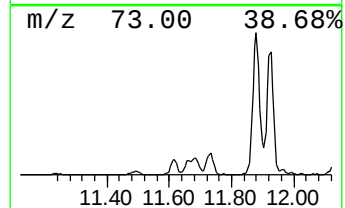
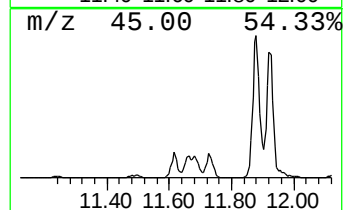
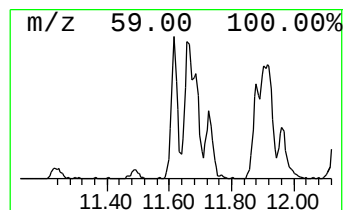
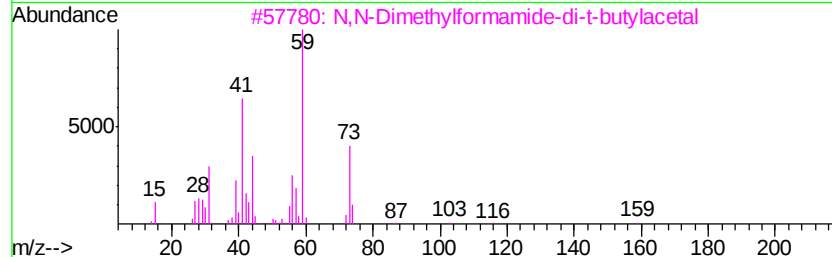
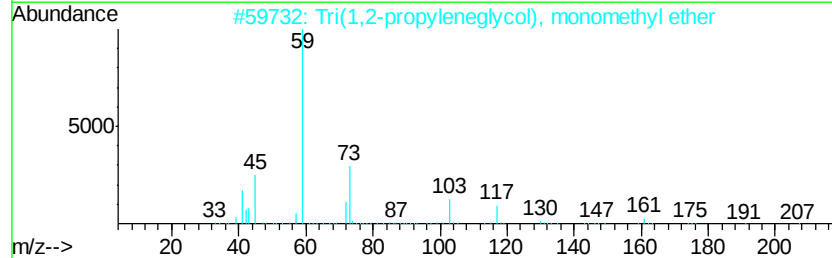
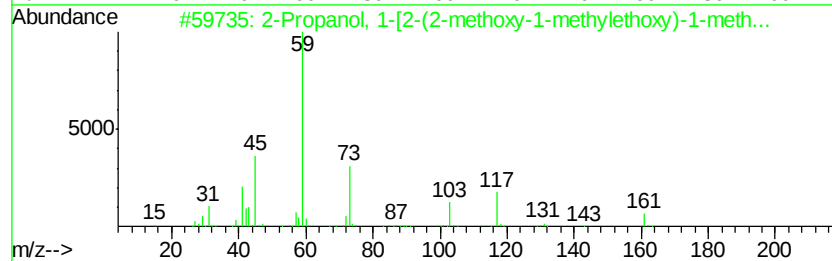
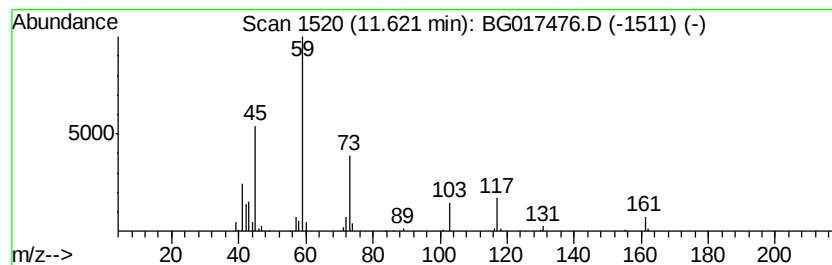
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.62	9.90 ng	95054	Naphthalene-d8	10.42		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	72	
2	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	50	
3	N,N-Dimethylformamide-di-t-butyl...	203	C11H25N02	036805-97-7	47	
4	2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	47	
5	Dipropylene glycol monomethyl ether	148	C7H16O3	034590-94-8	45	



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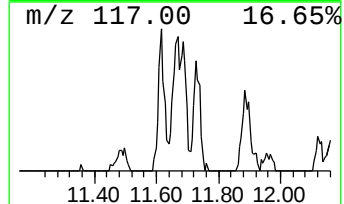
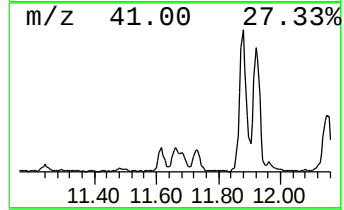
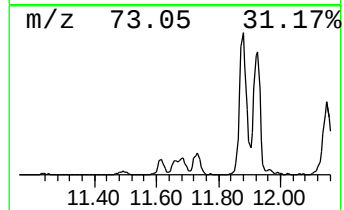
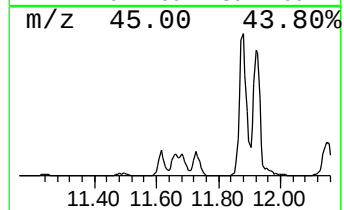
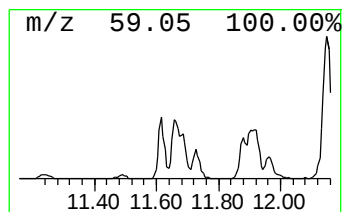
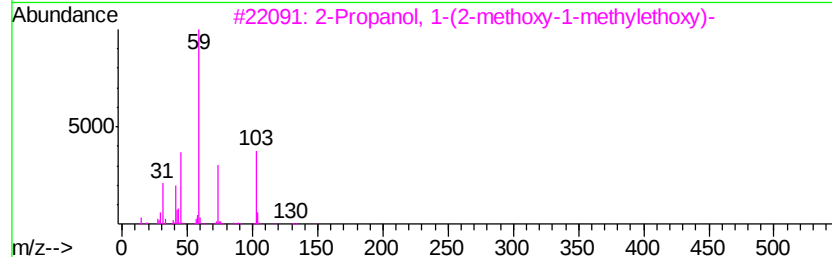
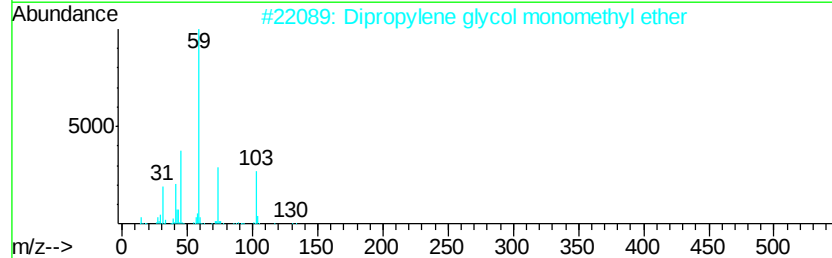
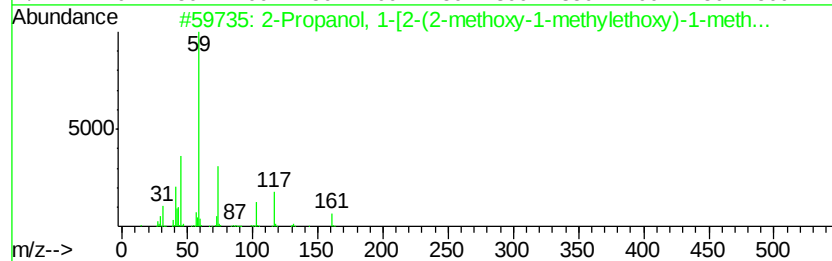
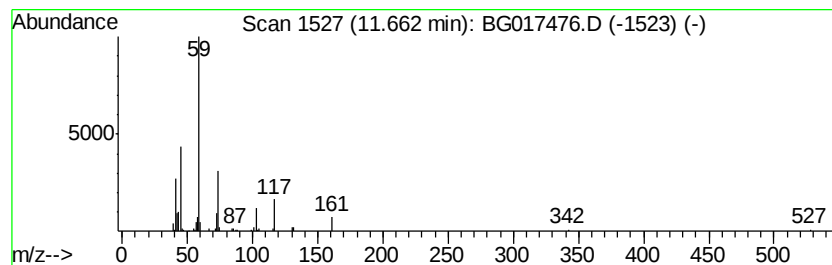
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Dipropylene glycol monometh... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	7.91 ng	75983	Napthalene-d8	10.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	90		
2	Dipropylene glycol monomethyl ether	148	C7H16O3	034590-94-8	72		
3	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72		
4	Tri(1,2-propyleneglycol), monome...	206	C10H22O4	1000262-26-6	64		
5	Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	50		



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
Acq On : 5 Jun 2015 16:23
Operator : TP/IZ
Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-5-6-2-15

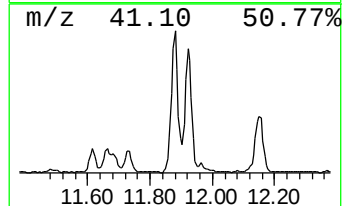
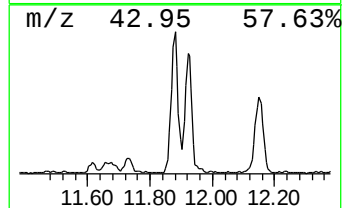
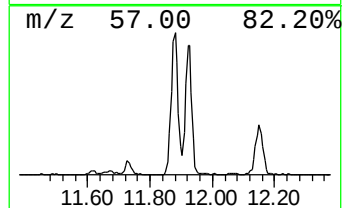
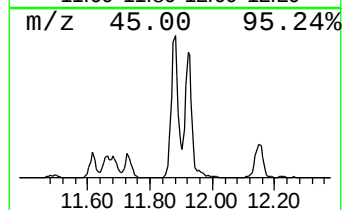
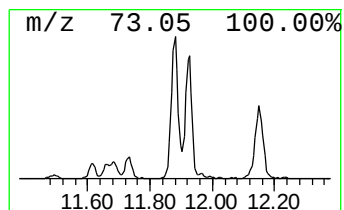
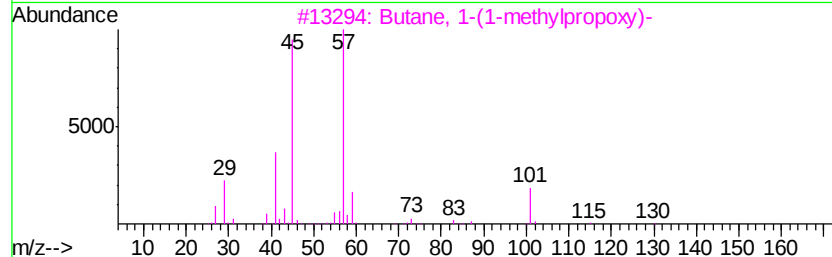
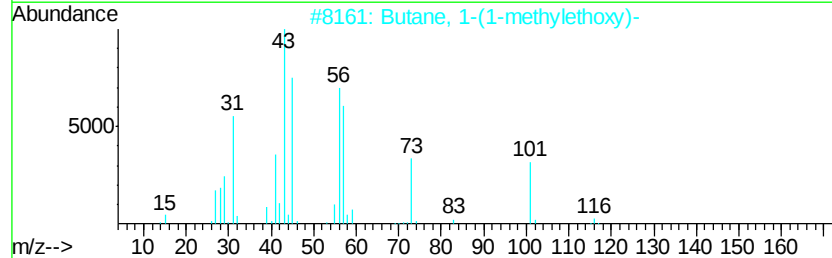
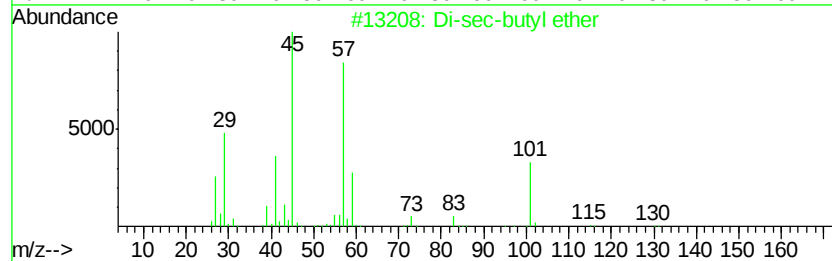
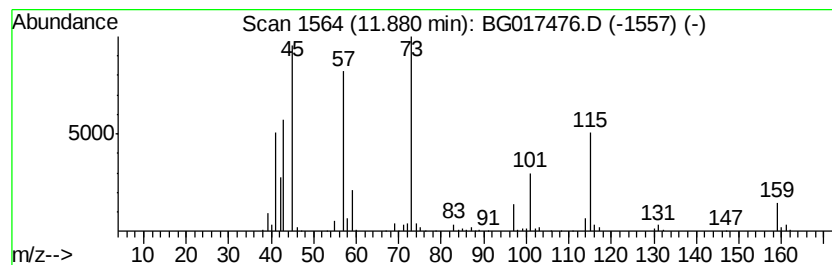
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown11.88 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.88	69.74 ng	669898	Napthalene-d8	10.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Di-sec-butyl ether	130	C8H18O	006863-58-7	47
2		Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	47
3		Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	43
4		Phloroglucitol	132	C6H12O3	002041-15-8	38
5		Butane, 1,1'-[ethylidenebis(oxy)...	174	C10H22O2	000871-22-7	37



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
Acq On : 5 Jun 2015 16:23
Operator : TP/IZ
Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-5-6-2-15

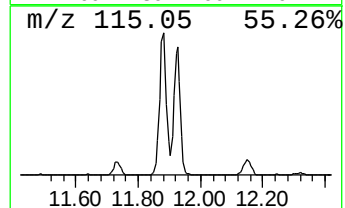
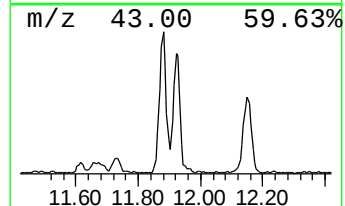
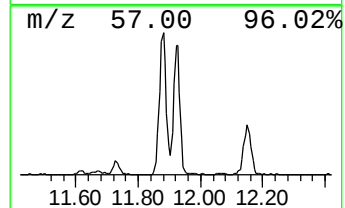
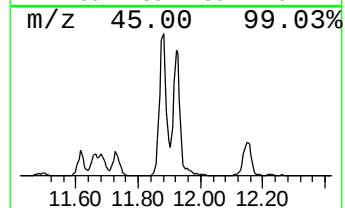
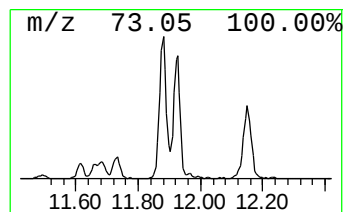
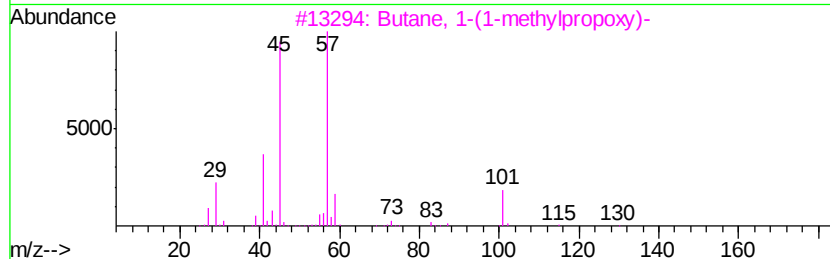
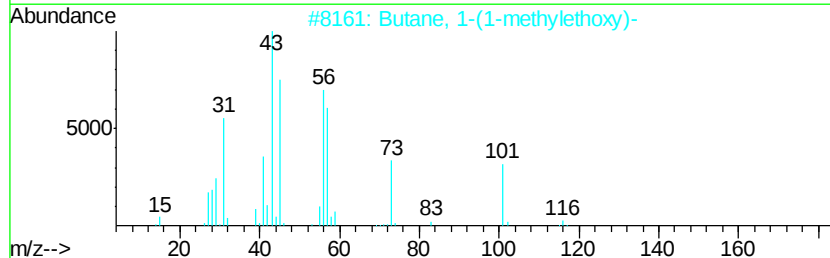
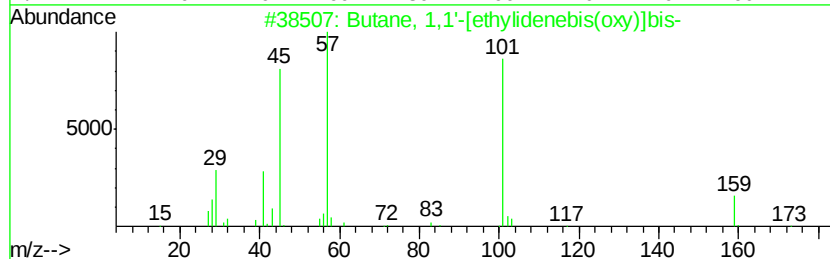
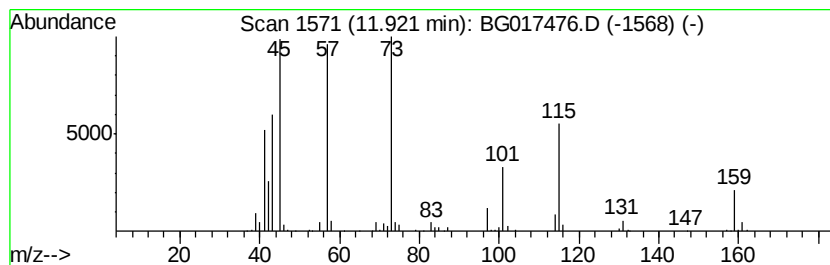
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown11.92 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	52.51 ng	504407	Napthalene-d8	10.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 1,1'-[ethylidenebis(oxy)]...	174	C10H22O2	000871-22-7	47
2		Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	43
3		Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	38
4		Di-sec-butyl ether	130	C8H18O	006863-58-7	38
5		4,6,9,11-Tetramethyl-1-oxa-4,6,9...	272	C12H24N4O3	1000139-36-0	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
Acq On : 5 Jun 2015 16:23
Operator : TP/IZ
Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

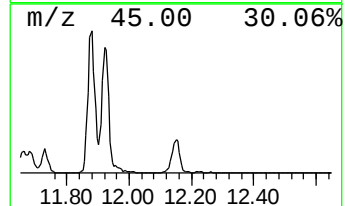
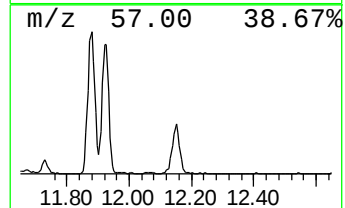
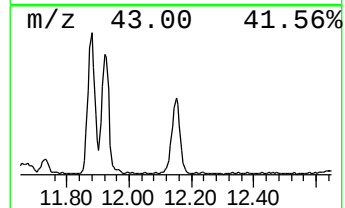
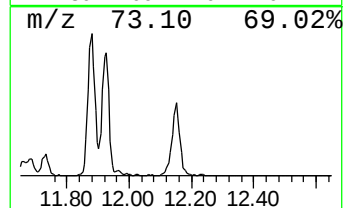
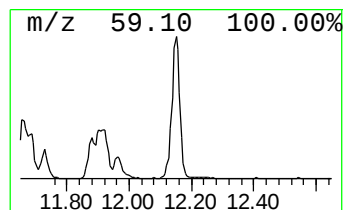
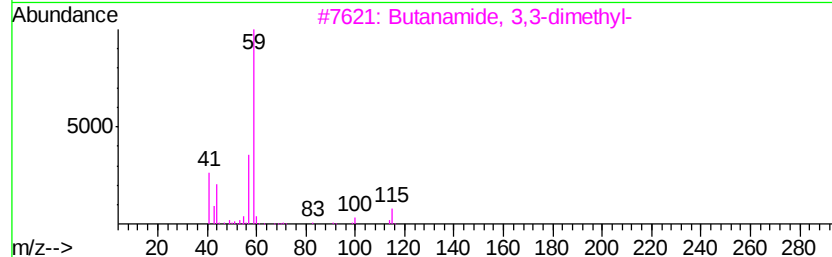
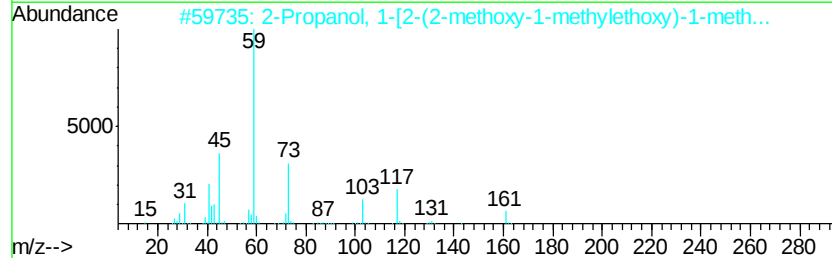
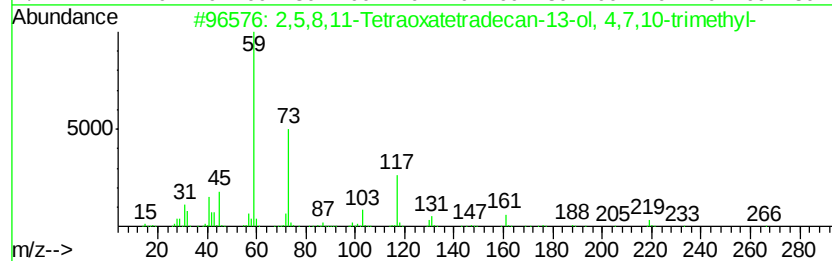
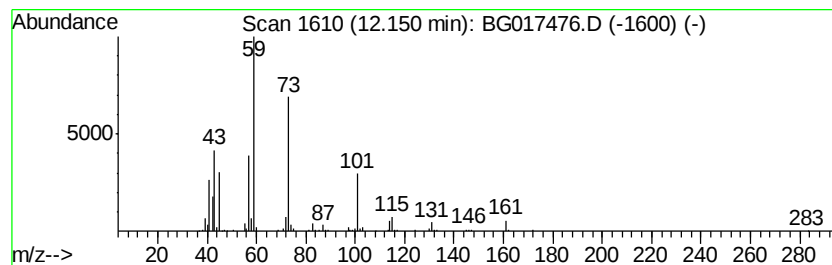
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ClientSampleId :
MW-5-6-2-15

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown12.15 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	40.08 ng	384954	Napthalene-d8	10.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,5,8,11-Tetraoxatetradecan-13-o...	264 C13H28O5	020324-34-9	40
2	2-Propanol, 1-[2-(2-methoxy-1-me...	206 C10H22O4	020324-33-8	39
3	Butanamide, 3,3-dimethyl-	115 C6H13NO	000926-04-5	37
4	2,4-Diethyl-6-methyl-1,3,5-trioxane	160 C8H16O3	117888-04-7	36
5	N-Acetyl-d-threo-0-methylthreonine	175 C7H13NO4	1000214-70-5	36



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
Acq On : 5 Jun 2015 16:23
Operator : TP/IZ
Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-5-6-2-15

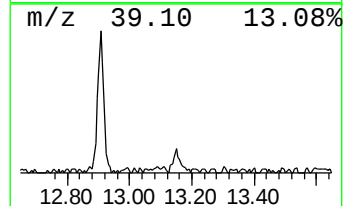
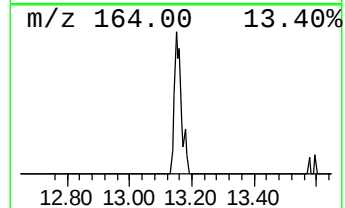
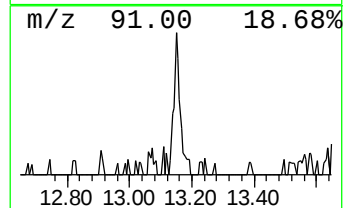
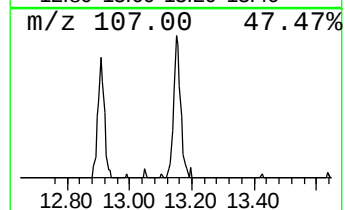
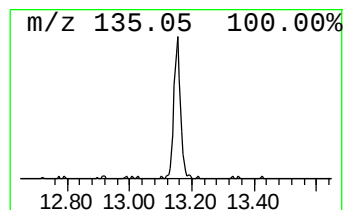
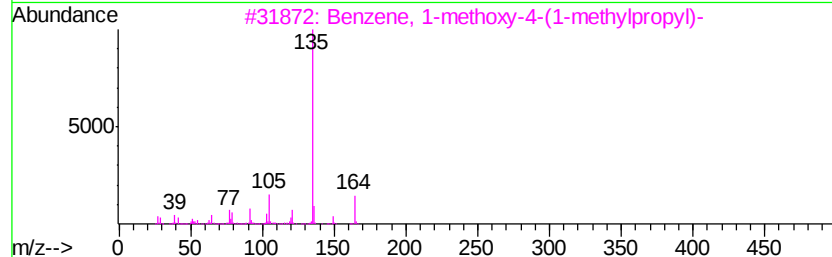
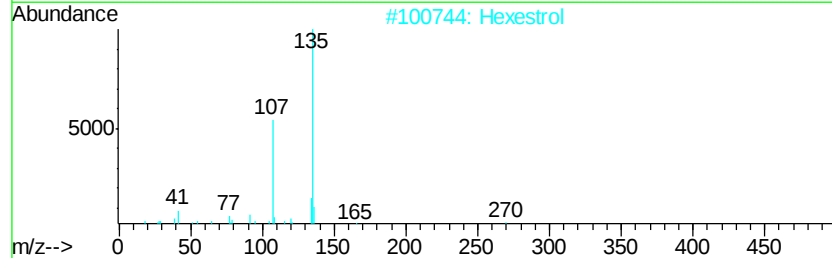
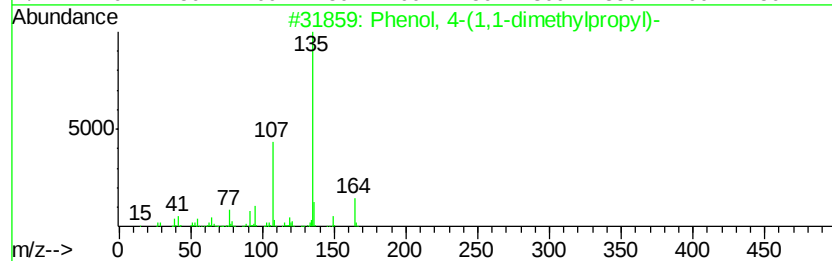
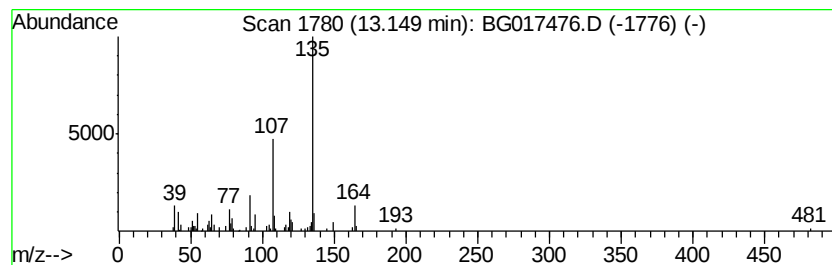
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenol, 4-(1,1-dimethylprop... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	4.07 ng	57698	Acenaphthene-d10	14.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	90
2		Hexestrol	270	C18H22O2	005635-50-7	59
3		Benzene, 1-methoxy-4-(1-methylpr...	164	C11H16O	004917-90-2	52
4		4-Isothiocyatomethyl-benzoic acid	193	C9H7NO2S	035009-14-4	50
5		Phenol, 5-methyl-2-(1-methylethy...	192	C12H16O2	000528-79-0	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
Acq On : 5 Jun 2015 16:23
Operator : TP/IZ
Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-5-6-2-15

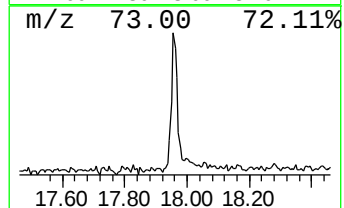
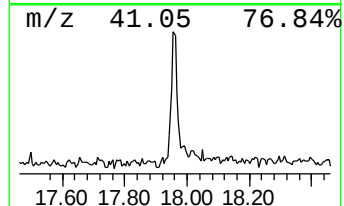
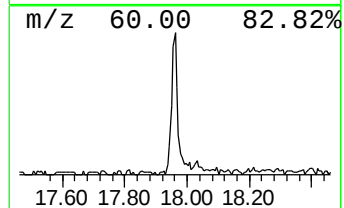
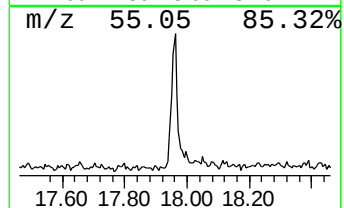
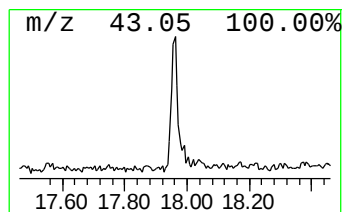
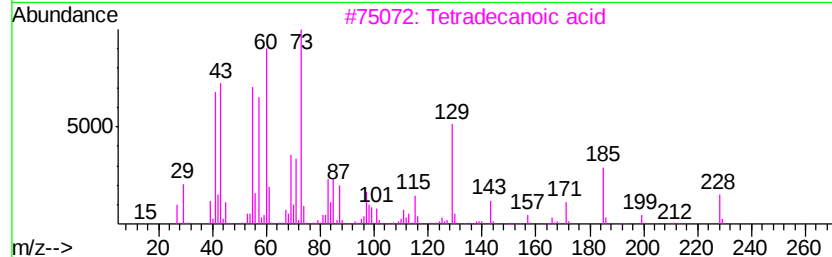
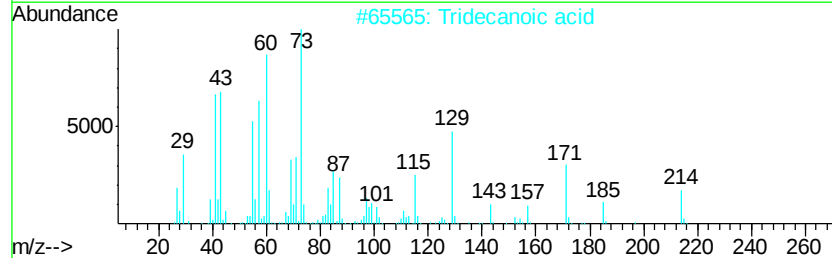
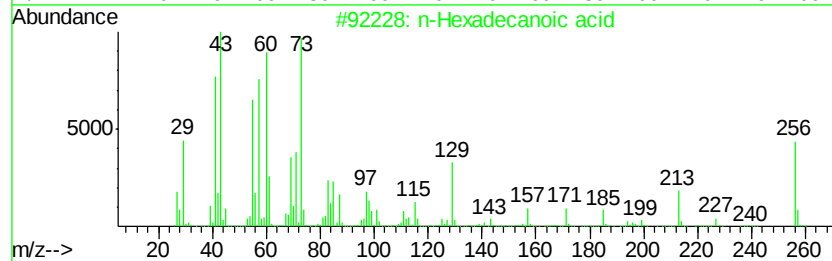
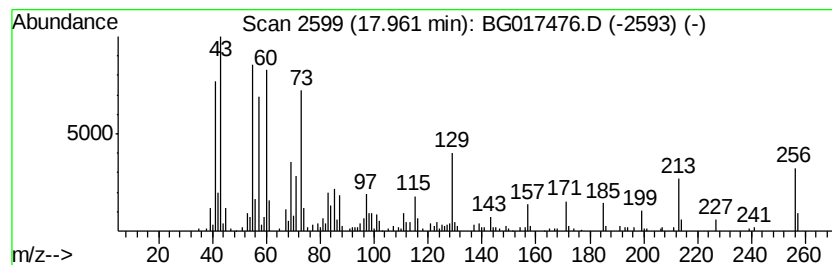
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	11.74 ng	208971	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92
2		Tridecanoic acid	214	C13H26O2	000638-53-9	64
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	58
4		n-Decanoic acid	172	C10H20O2	000334-48-5	50
5		Dodecanoic acid	200	C12H24O2	000143-07-7	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
Data File : BG017476.D
Acq On : 5 Jun 2015 16:23
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Sample : G2501-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-5-6-2-15

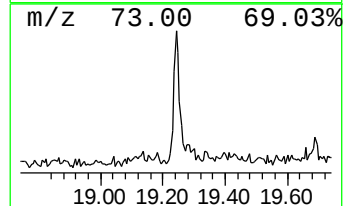
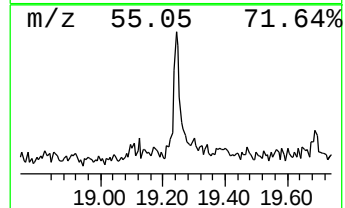
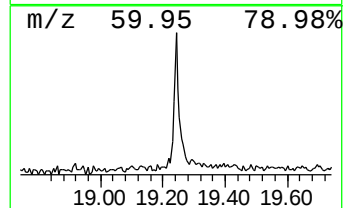
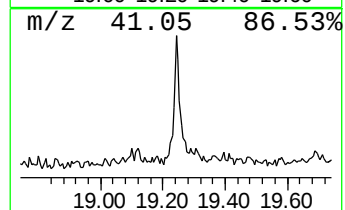
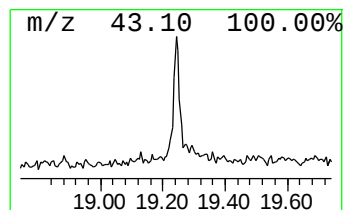
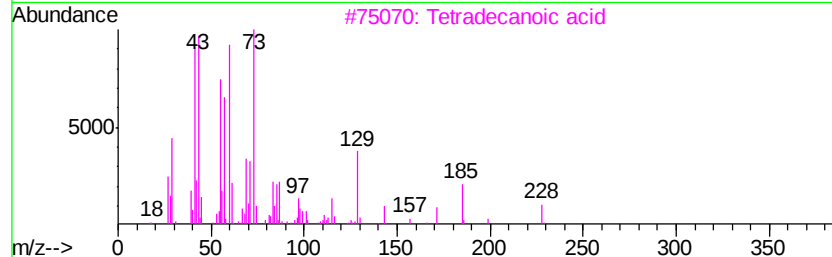
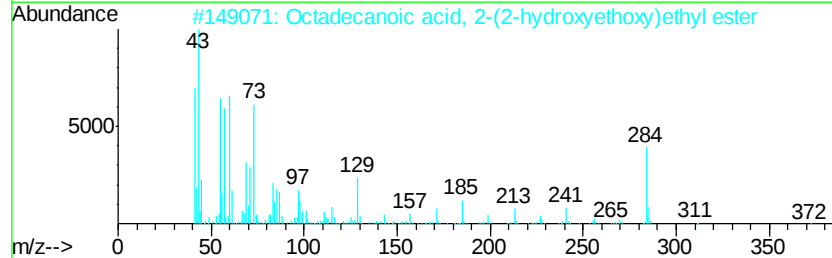
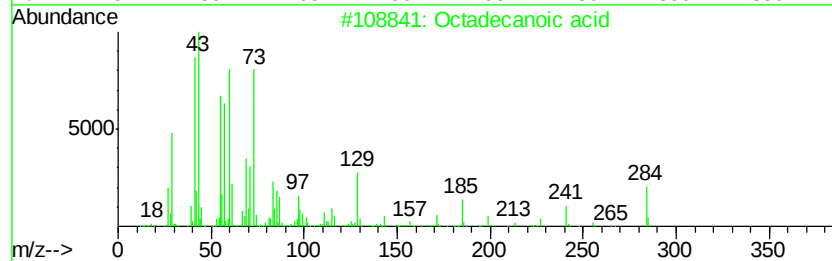
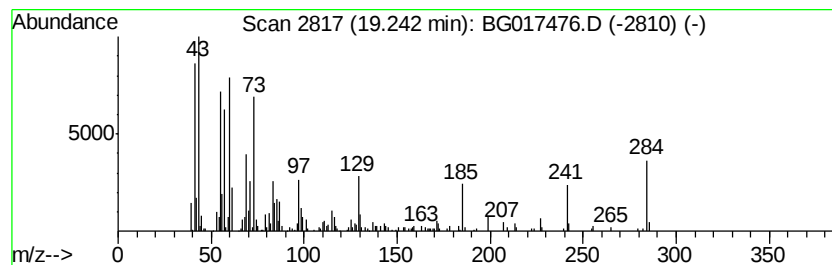
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	7.14 ng	184259	Chrysene-d12	21.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	93
2			Octadecanoic acid, 2-(2-hydroxyethyl) ester	372	C22H44O4	000106-11-6	74
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	52
4			n-Decanoic acid	172	C10H20O2	000334-48-5	50
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	49



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060515\
 Data File : BG017476.D
 Acq On : 5 Jun 2015 16:23
 Operator : TP/IZ
 Sample : G2501-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-5-6-2-15

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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...	2.77	46.5	ng	379328	1	7.62	163237	20.0
unknown7.17	7.17	78.0	ng	636707	1	7.62	163237	20.0
2,4-Nonadiyne	7.27	2.9	ng	23828	1	7.62	163237	20.0
Di-sec-butyl ether	7.41	9.0	ng	73673	1	7.62	163237	20.0
2-Propanol, 1-[2-...	11.62	9.9	ng	95054	2	10.42	192113	20.0
Dipropylene glyco...	11.66	7.9	ng	75983	2	10.42	192113	20.0
unknown11.73	11.73	9.9	ng	95428	2	10.42	192113	20.0
unknown11.88	11.88	69.7	ng	669898	2	10.42	192113	20.0
unknown11.92	11.92	52.5	ng	504407	2	10.42	192113	20.0
unknown12.15	12.15	40.1	ng	384954	2	10.42	192113	20.0
Phenol, 4-(1,1-di...	13.15	4.1	ng	57698	3	14.29	283828	20.0
n-Hexadecanoic acid	17.96	11.7	ng	208971	4	17.04	356068	20.0
Octadecanoic acid	19.24	7.1	ng	184259	5	21.24	516308	20.0