

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
 Data File : BG026002.D
 Acq On : 28 Feb 2017 18:19
 Operator : SJ/MA
 Sample : I2006-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C19013051

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-BG021717.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	5.173	307	312	319	rBV	222766	326583	3.50%	0.727%
2	5.643	385	392	403	rVB	991884	1545412	16.55%	3.438%
3	7.230	654	662	673	rBV	636809	1083860	11.60%	2.411%
4	7.606	719	726	735	rBV	1649133	2812219	30.11%	6.256%
5	8.076	799	806	814	rBV	331156	543578	5.82%	1.209%
6	8.399	853	861	868	rBV	1383579	2442582	26.15%	5.434%
7	9.233	988	1003	1016	rBV	1180440	2246181	24.05%	4.997%
8	10.279	1175	1181	1188	rBV2	117565	214063	2.29%	0.476%
9	10.878	1271	1283	1288	rBV	499999	977785	10.47%	2.175%
10	10.931	1288	1292	1297	rVB4	60750	106725	1.14%	0.237%
11	10.990	1297	1302	1307	rBV2	58304	96853	1.04%	0.215%
12	11.760	1419	1433	1445	rBV2	283053	680246	7.28%	1.513%
13	11.977	1463	1470	1475	rBV	76386	132178	1.42%	0.294%
14	12.735	1588	1599	1609	rVB	215579	405185	4.34%	0.901%
15	13.311	1689	1697	1703	rBV	3732427	5708262	61.12%	12.699%
16	13.740	1763	1770	1777	rVB2	68277	136318	1.46%	0.303%
17	13.998	1808	1814	1820	rBV2	203517	367315	3.93%	0.817%
18	14.057	1820	1824	1831	rVB3	63968	146630	1.57%	0.326%
19	14.233	1848	1854	1858	rBV	113366	196290	2.10%	0.437%
20	14.268	1858	1860	1865	rVB2	85084	107347	1.15%	0.239%
21	14.680	1918	1930	1936	rBV2	909717	1519152	16.27%	3.380%
22	14.744	1936	1941	1947	rVV3	48755	116447	1.25%	0.259%
23	14.897	1961	1967	1973	rBV4	49285	118680	1.27%	0.264%
24	14.968	1974	1979	1986	rVV3	47232	118955	1.27%	0.265%
25	15.073	1989	1997	2004	rVB2	109278	219930	2.35%	0.489%
26	15.291	2028	2034	2040	rBV2	130013	224032	2.40%	0.498%
27	15.714	2100	2106	2111	rVB2	115749	193074	2.07%	0.430%
28	15.843	2118	2128	2132	rBV	165621	301615	3.23%	0.671%
29	16.154	2174	2181	2190	rBV	3284905	5078471	54.37%	11.298%
30	16.378	2216	2219	2227	rVB4	52356	112708	1.21%	0.251%
31	16.766	2280	2285	2298	rVB3	114580	236825	2.54%	0.527%
32	17.406	2388	2394	2403	rBV2	1272210	1961292	21.00%	4.363%
33	18.287	2538	2544	2549	rBV	172764	281482	3.01%	0.626%
34	19.550	2754	2759	2766	rBV	182298	259027	2.77%	0.576%

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Peak separation: 5

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35	20.003	2830	2836	2842	rBV	6581223	9339759	100.00%	20.778%
36	21.296	3052	3056	3066	rBV4	63458	122241	1.31%	0.272%
37	21.654	3110	3117	3127	rVB	1304315	2110373	22.60%	4.695%
38	22.453	3248	3253	3268	rVB	115797	245666	2.63%	0.547%
39	24.856	3651	3662	3677	rVB2	735427	2115737	22.65%	4.707%

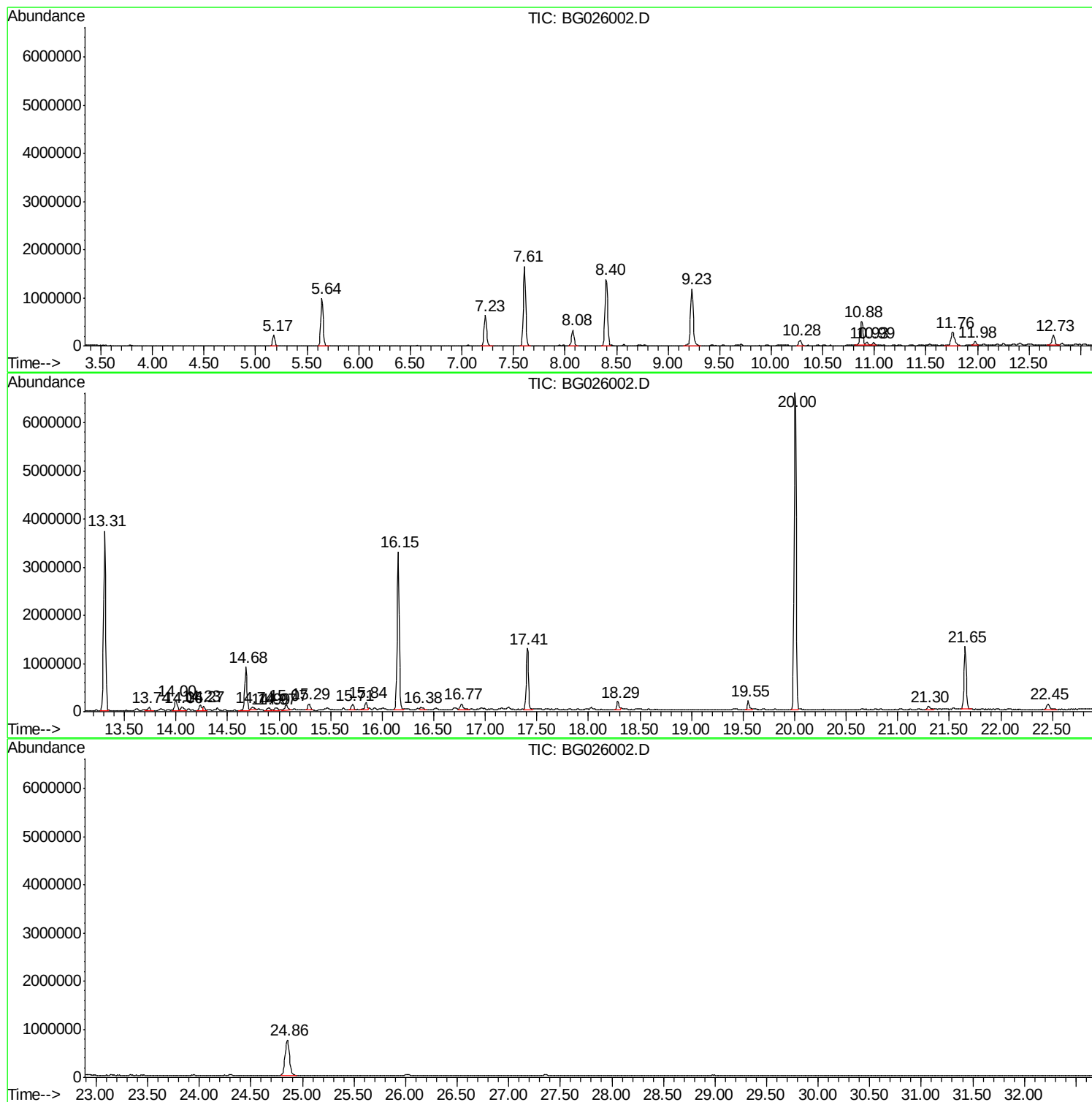
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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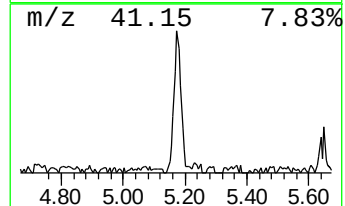
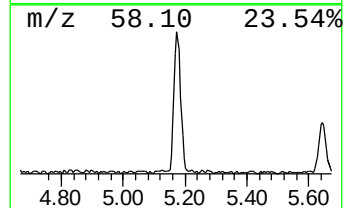
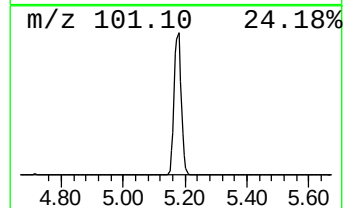
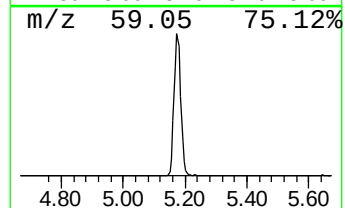
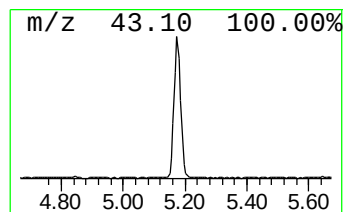
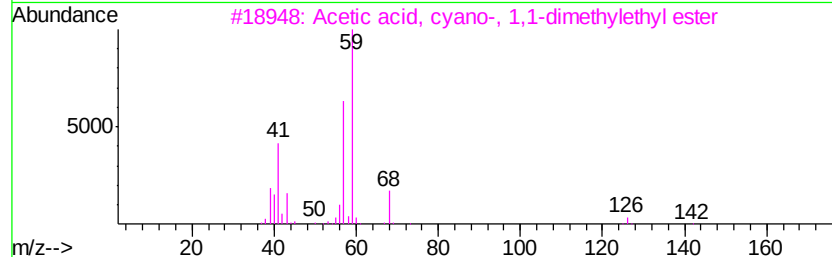
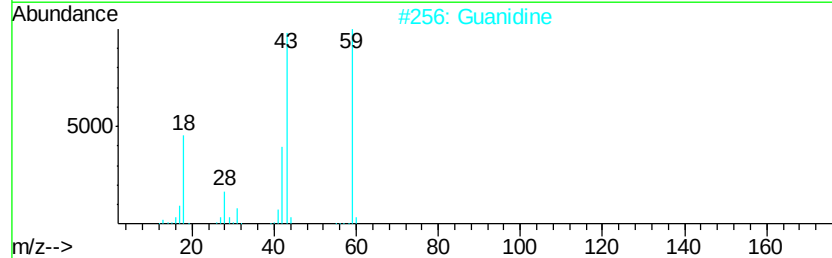
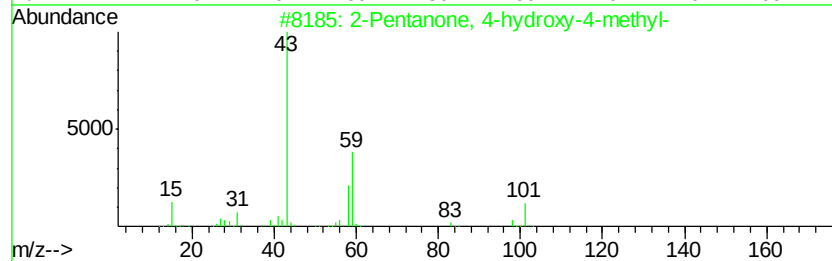
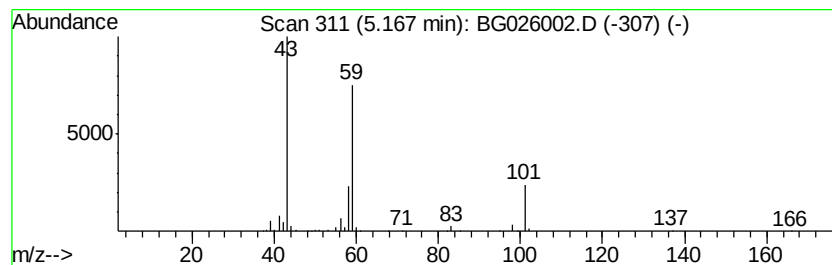
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	12.02 ng	326583	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Guanidine	59	CH5N3	000113-00-8	38
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	12
4		Acetone	58	C3H6O	000067-64-1	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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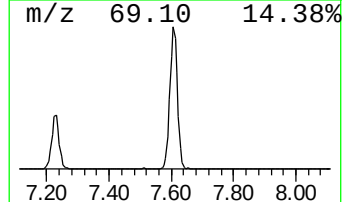
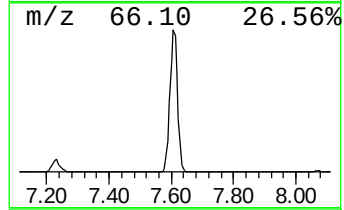
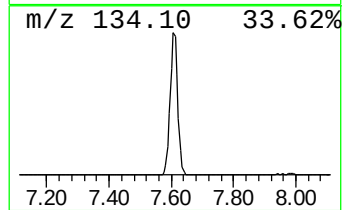
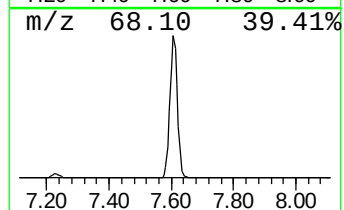
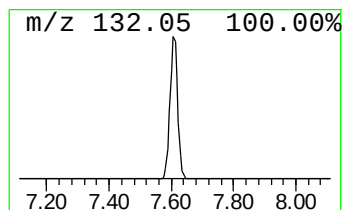
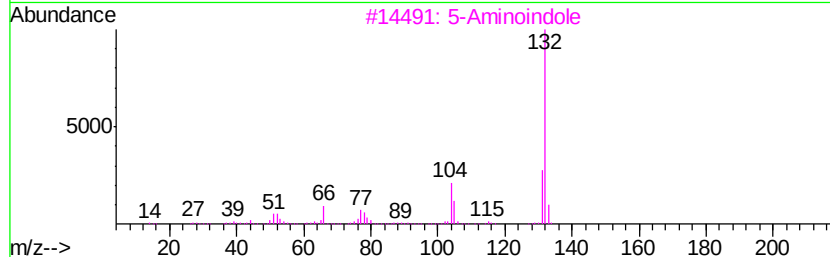
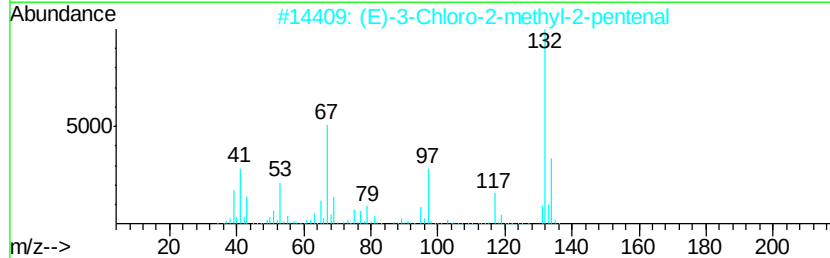
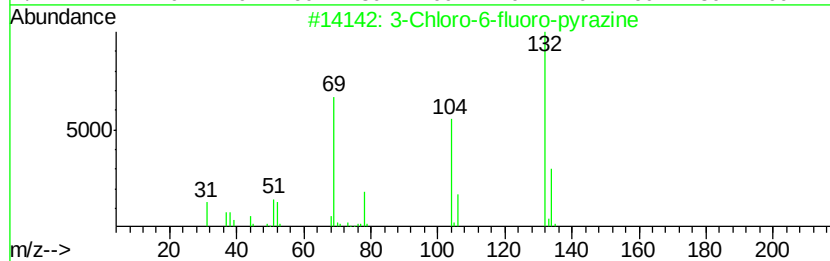
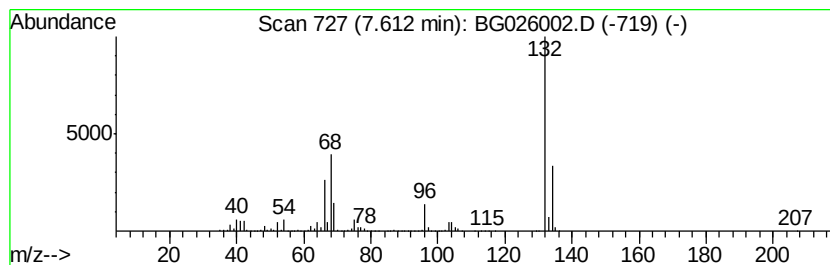
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	103.47 ng	2812220	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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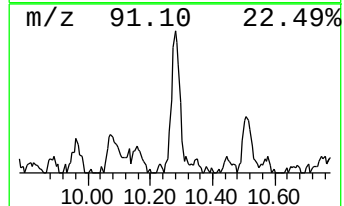
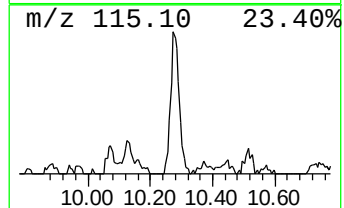
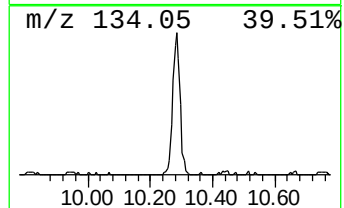
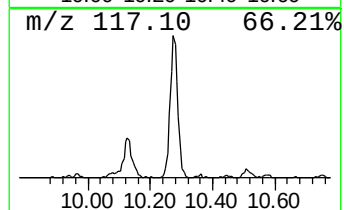
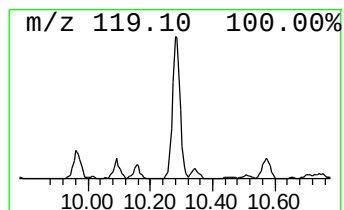
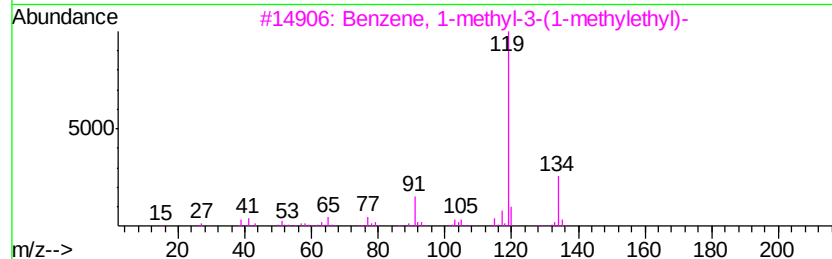
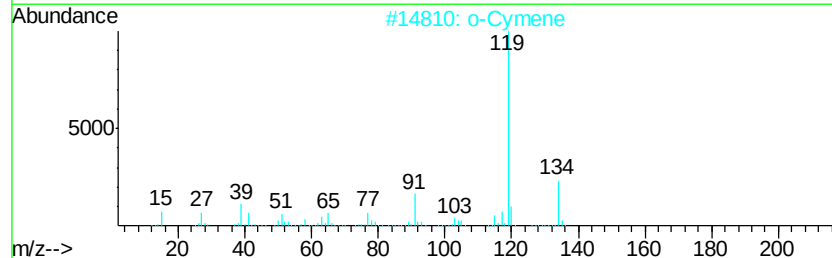
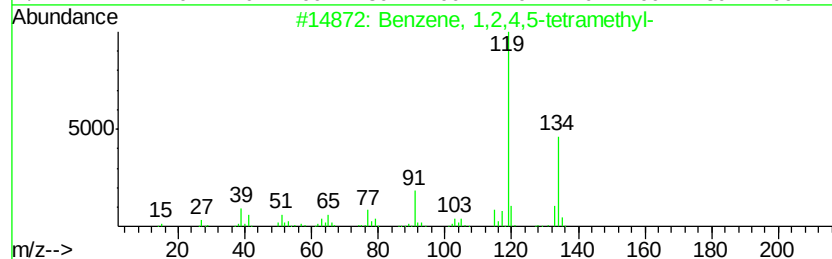
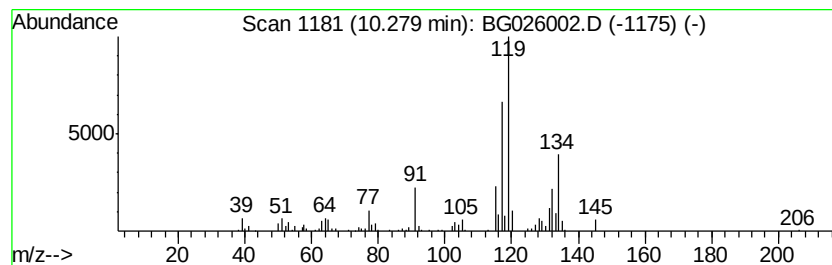
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.28	4.38 ng	214063	Naphthalene-d8	10.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2		o-Cymene	134	C10H14	000527-84-4	60
3		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	60
5		p-Cymene	134	C10H14	000099-87-6	60



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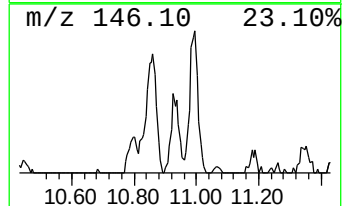
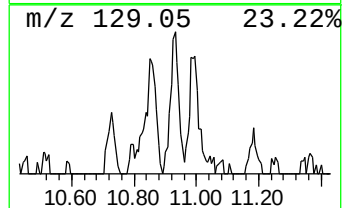
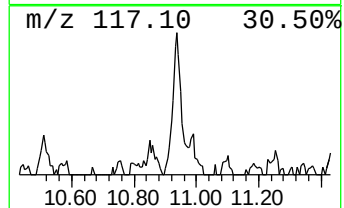
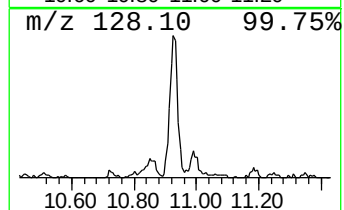
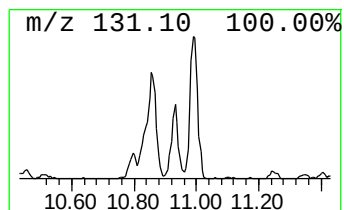
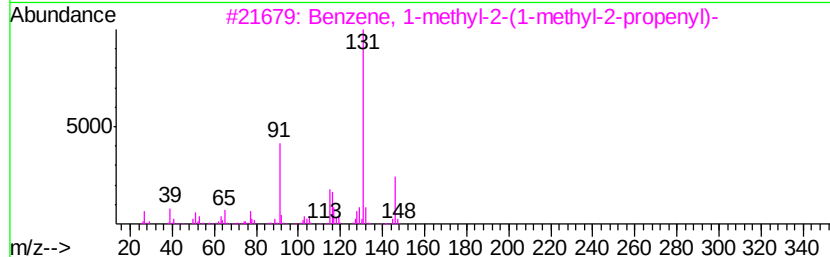
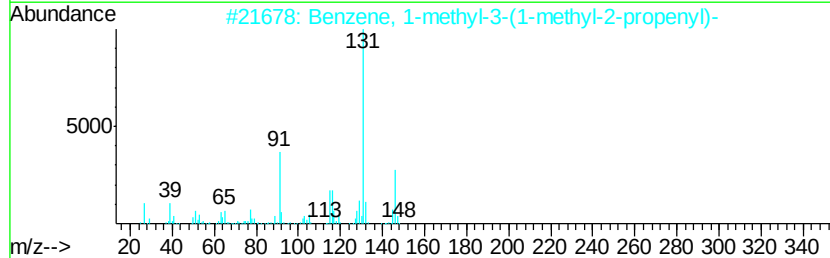
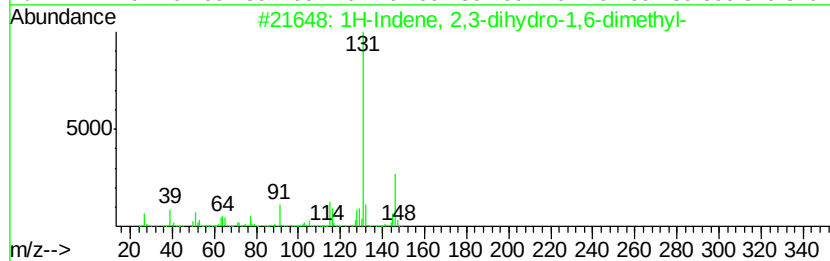
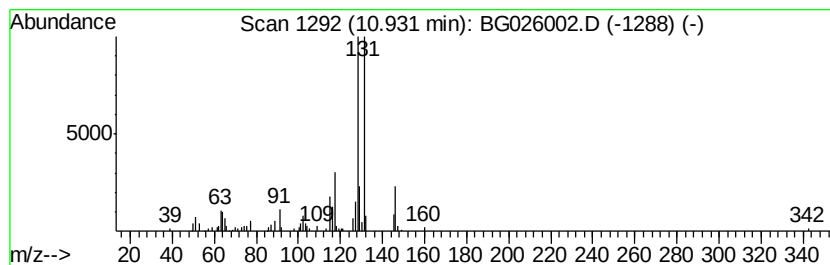
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown10.93 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.93	2.18 ng	106725	Naphthalene-d8	10.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	45
2		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	45
3		Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	43
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	43
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	43



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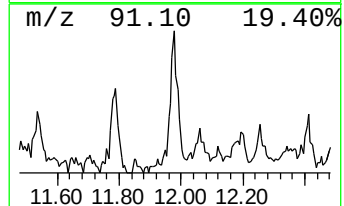
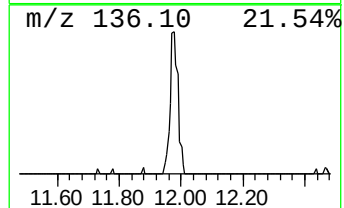
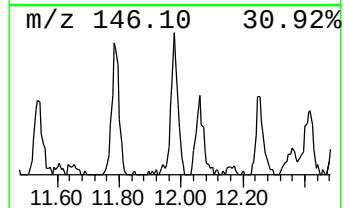
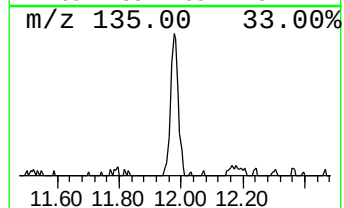
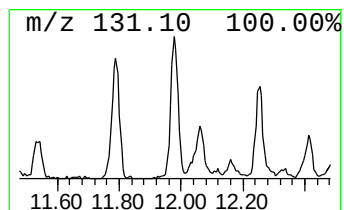
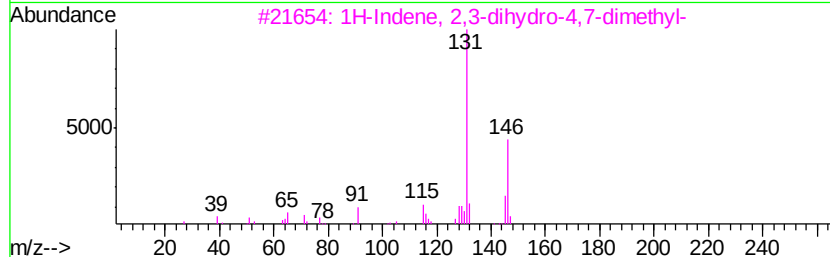
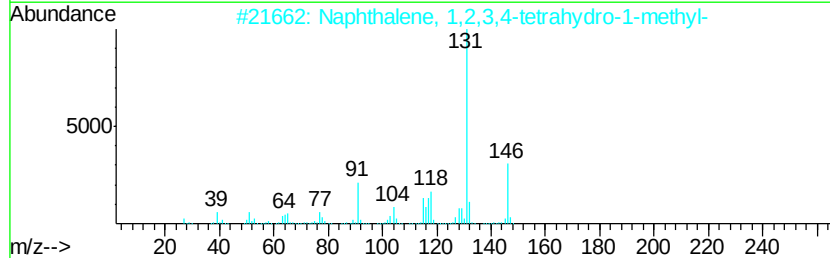
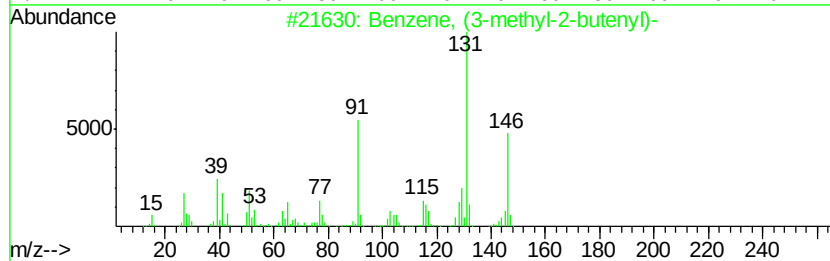
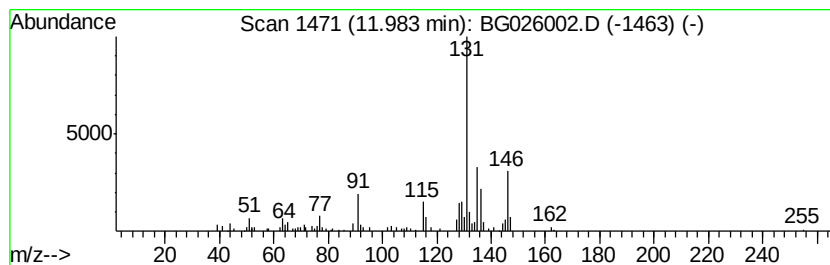
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Benzene, (3-methyl-2-butenyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	2.70 ng	132178	Naphthalene-d8	10.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	93
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	89
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	76
4		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	76
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
Operator : SJ/MA
Sample : I2006-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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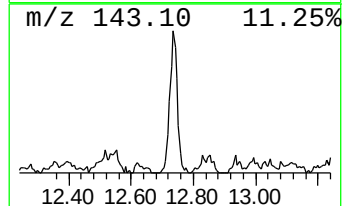
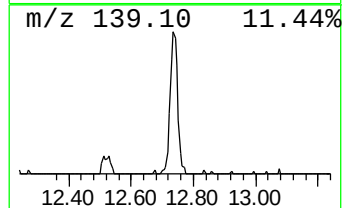
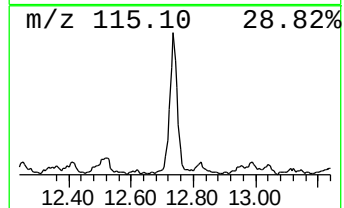
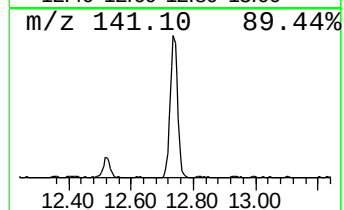
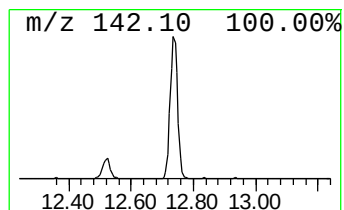
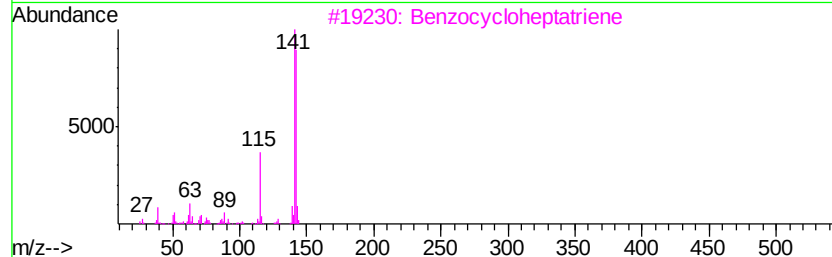
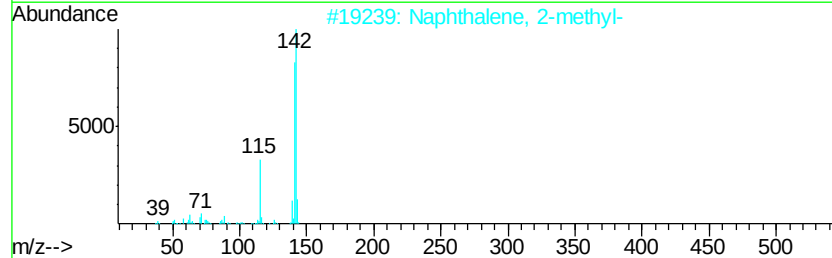
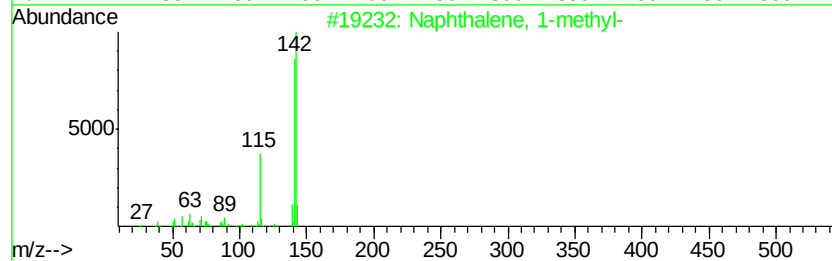
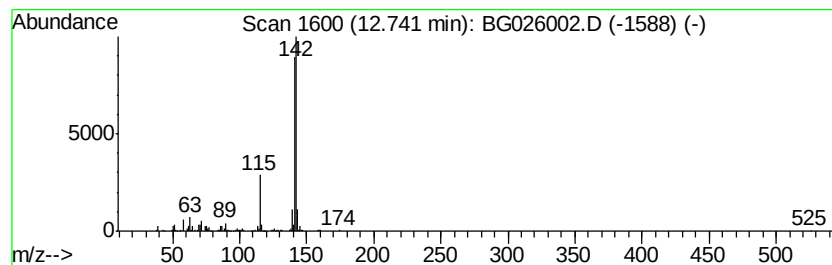
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	8.29 ng	405185	Naphthalene-d8	10.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1-Naphthaleneacetamide	185	C12H11NO	000086-86-2	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
Operator : SJ/MA
Sample : I2006-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

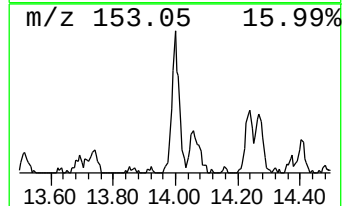
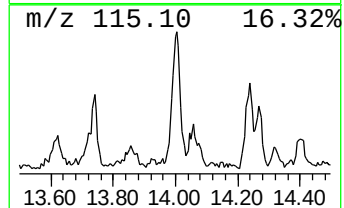
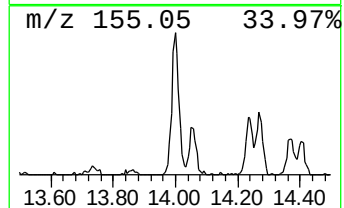
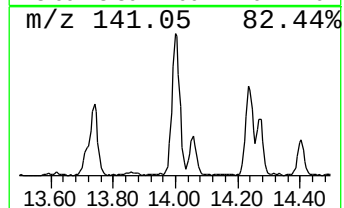
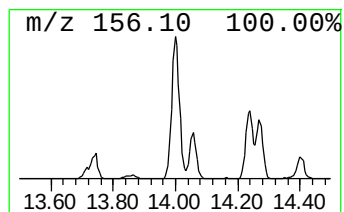
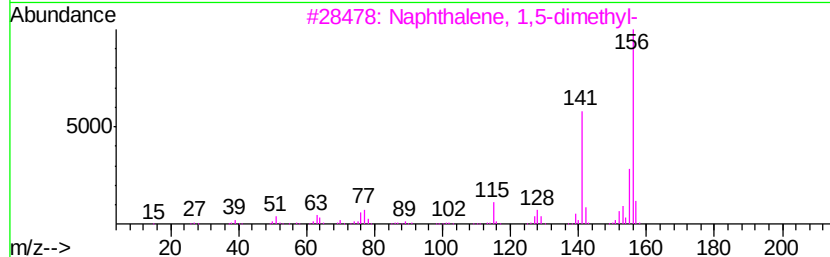
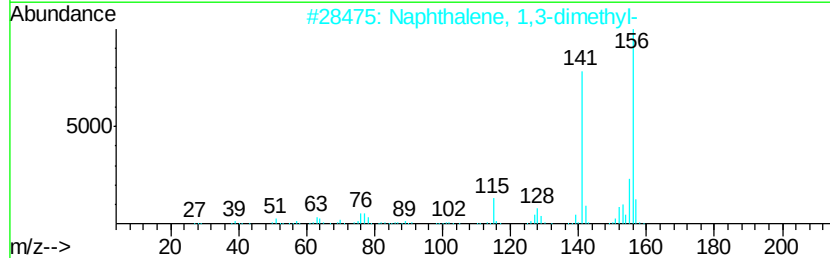
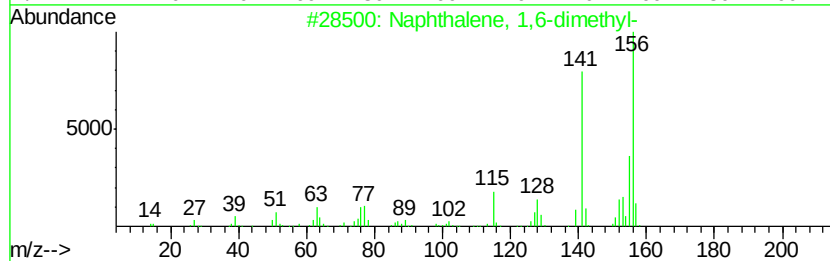
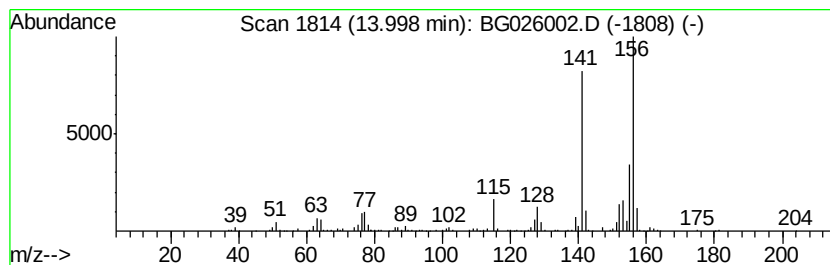
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 1,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.		
14.00	4.84 ng	367315	Acenaphthene-d10		14.68		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
Operator : SJ/MA
Sample : I2006-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

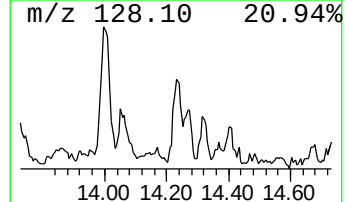
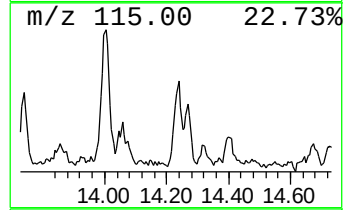
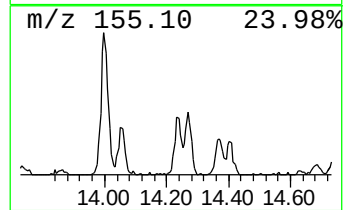
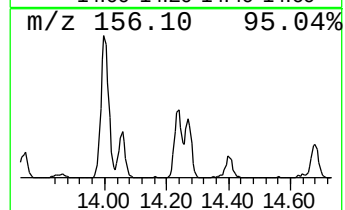
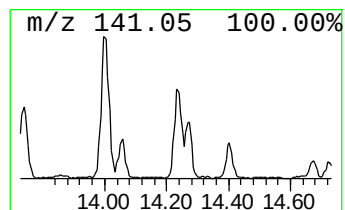
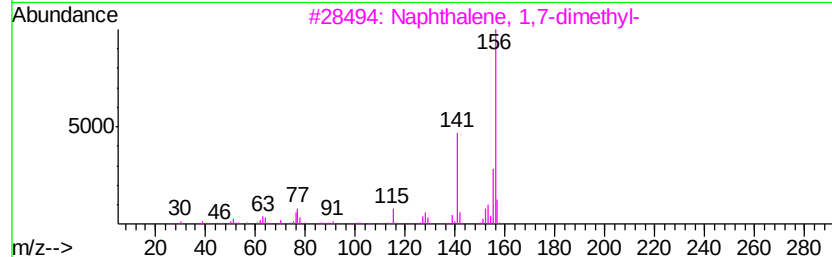
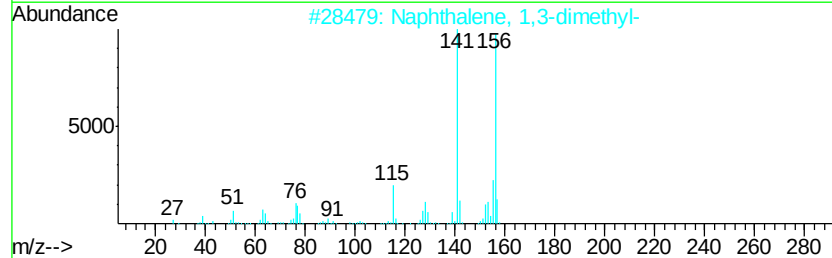
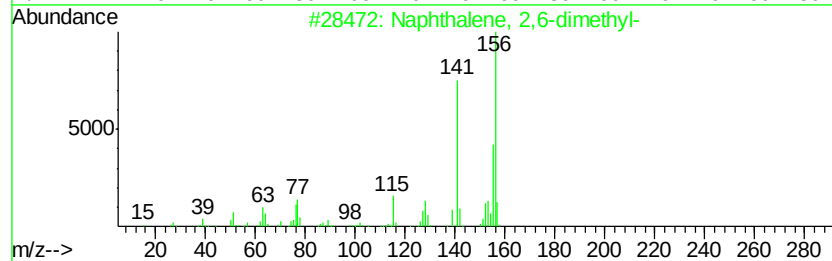
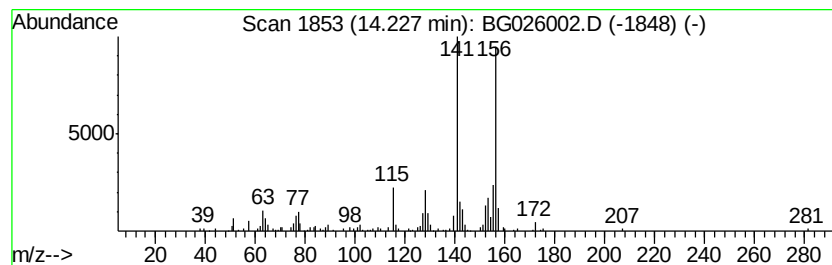
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.		
14.23	2.58 ng	196290	Acenaphthene-d10		14.68		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
Operator : SJ/MA
Sample : I2006-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

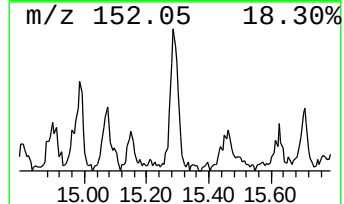
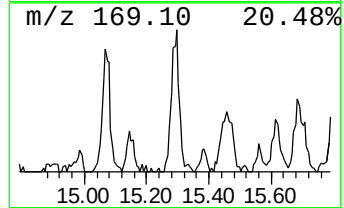
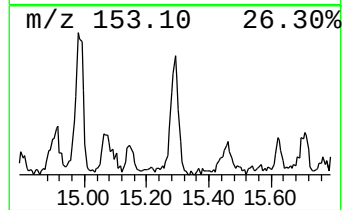
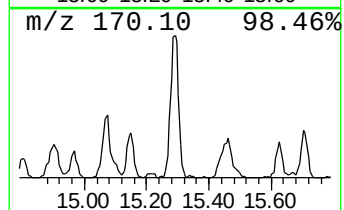
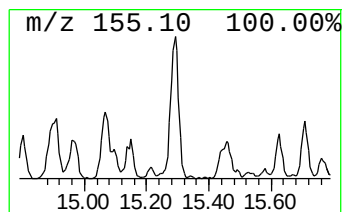
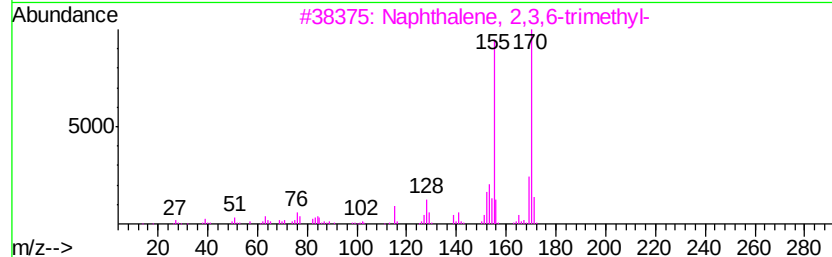
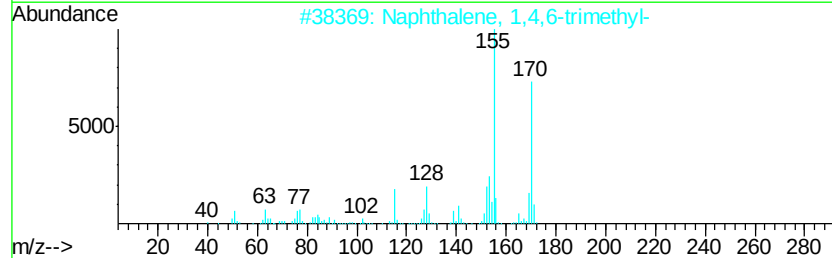
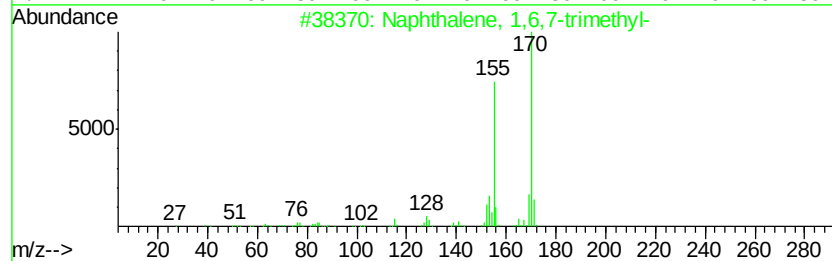
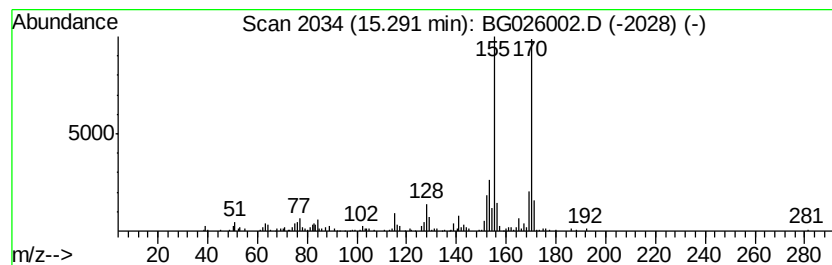
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 1,6,7-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.29	2.95 ng	224032	Acenaphthene-d10	14.68
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,6,7-trimethyl-	170 C13H14	002245-38-7 97
2		Naphthalene, 1,4,6-trimethyl-	170 C13H14	002131-42-2 97
3		Naphthalene, 2,3,6-trimethyl-	170 C13H14	000829-26-5 97
4		4,6,8-Trimethylazulene	170 C13H14	000941-81-1 91
5		Naphthalene, 1,4,5-trimethyl-	170 C13H14	002131-41-1 72



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
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Sample : I2006-01
Misc :
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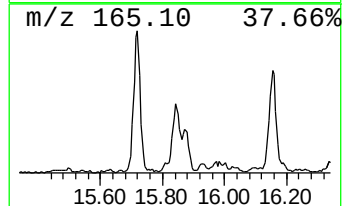
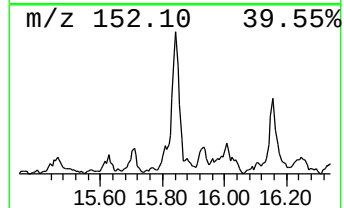
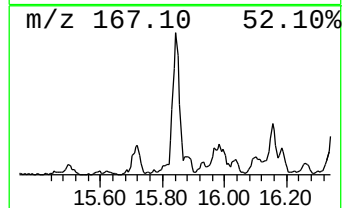
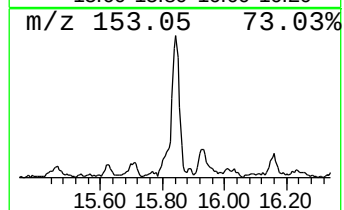
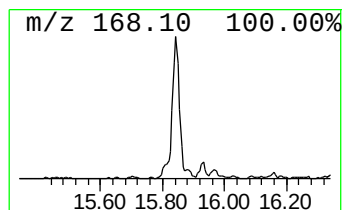
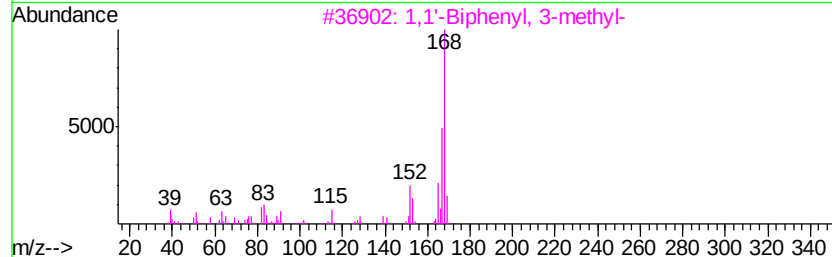
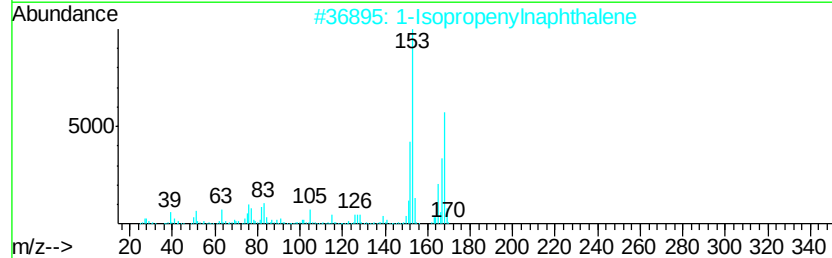
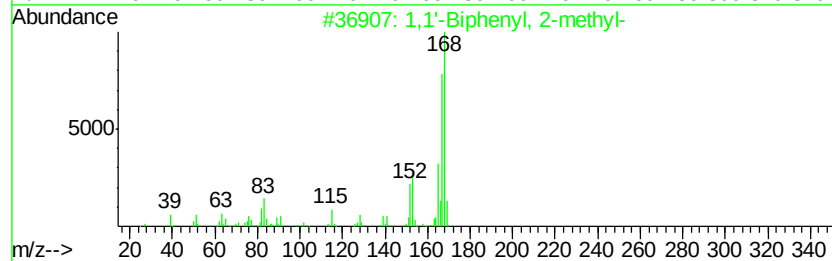
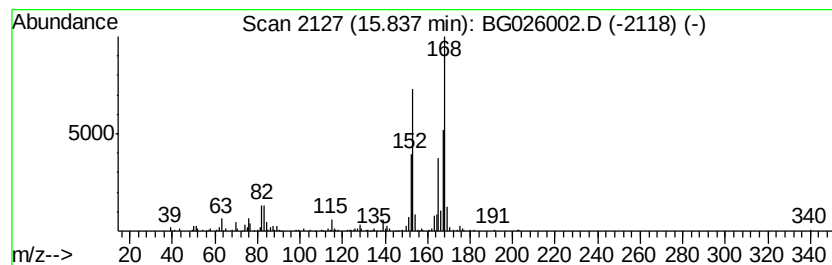
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,1'-Biphenyl, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.84	3.97 ng	301615	Acenaphthene-d10		14.68
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	81
2	1-Isopropenylnaphthalene	168	C13H12	001855-47-6	76
3	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	68
4	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	68
5	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
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Sample : I2006-01
Misc :
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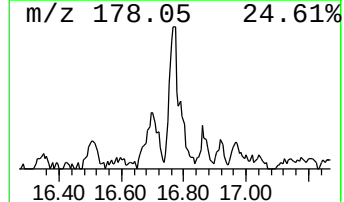
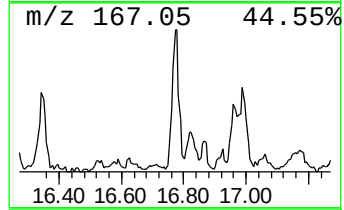
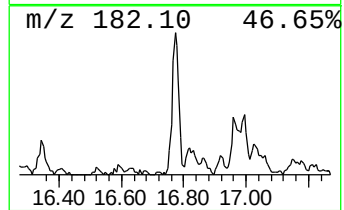
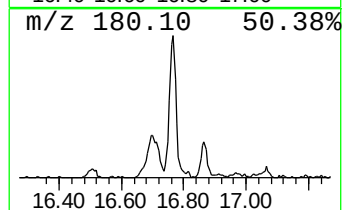
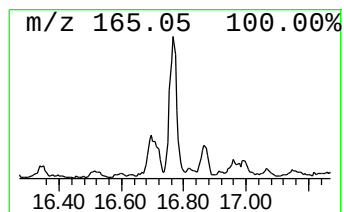
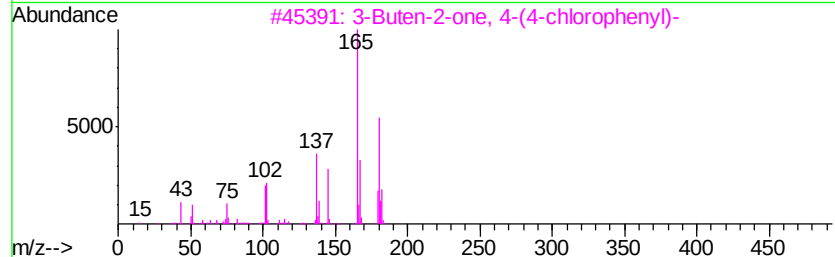
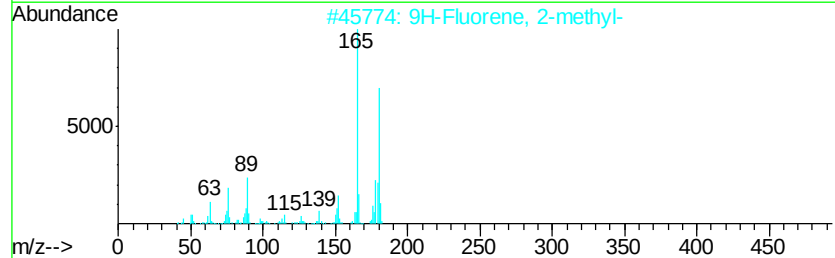
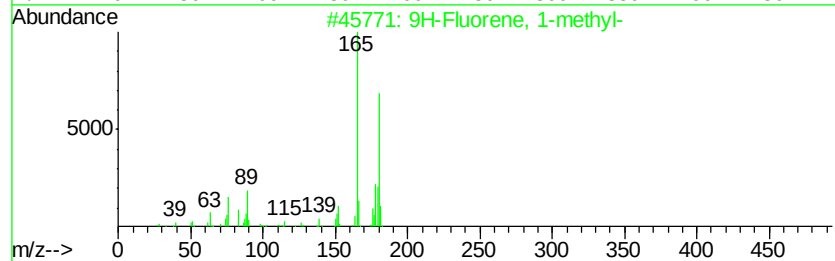
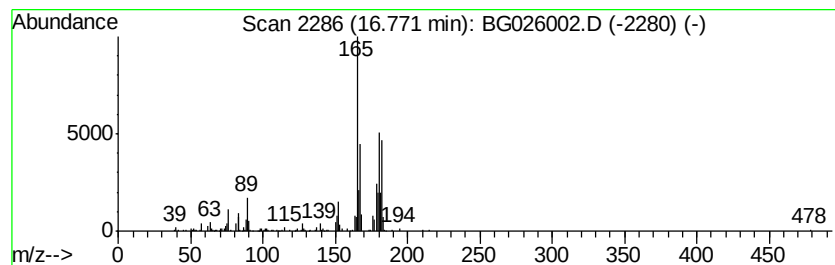
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 9H-Fluorene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.77	2.41 ng	236825	Phenanthrene-d10	17.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	91
2		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	83
3		3-Buten-2-one, 4-(4-chlorophenyl)-	180	C10H9ClO	003160-40-5	74
4		9H-Fluorene, 4-methyl-	180	C14H12	001556-99-6	70
5		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
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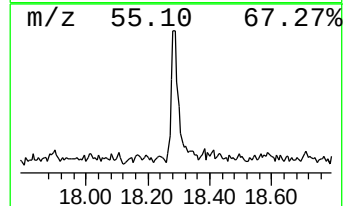
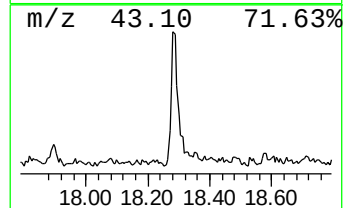
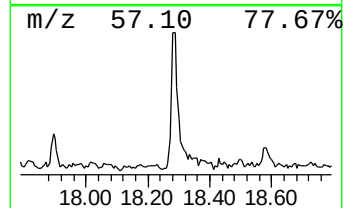
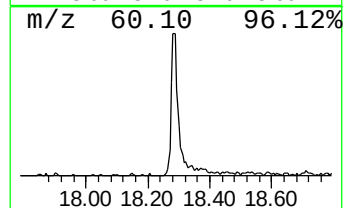
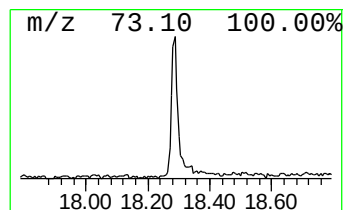
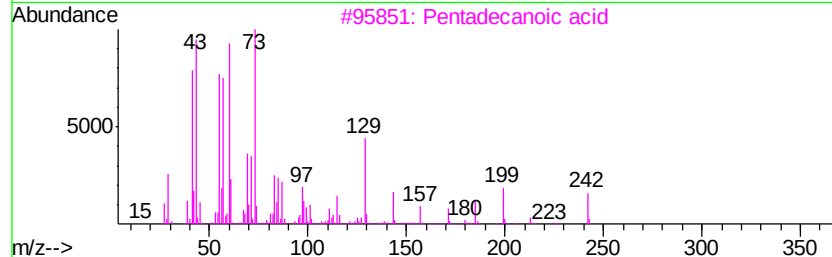
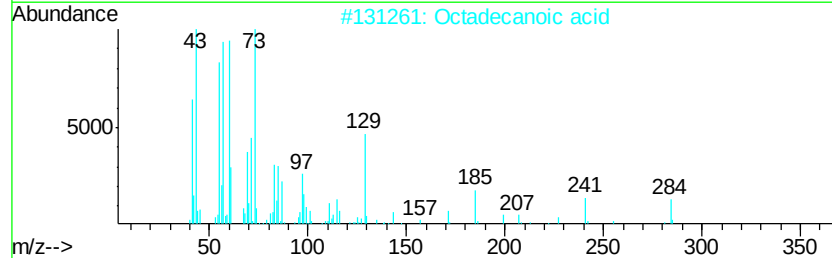
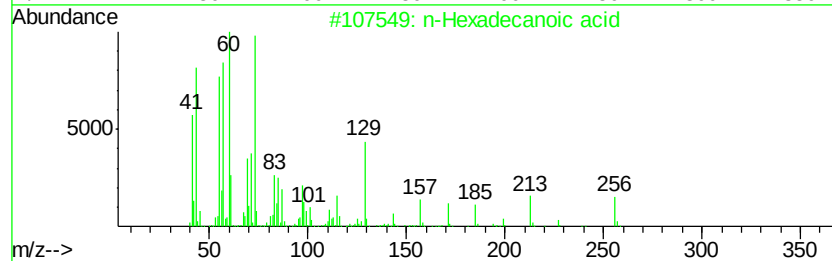
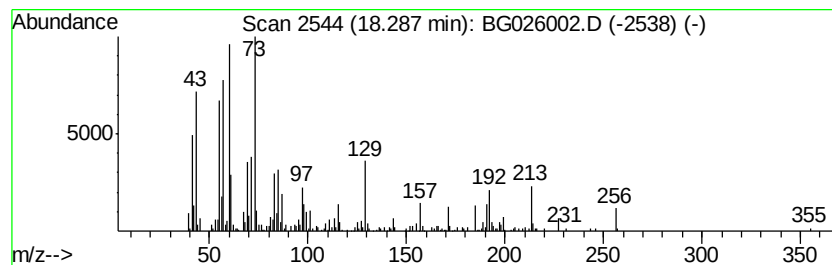
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.29	2.87 ng	281482	Phenanthrene-d10	17.41		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	72
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4		Tridecanoic acid	214	C13H26O2	000638-53-9	64
5		n-Decanoic acid	172	C10H20O2	000334-48-5	60



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
Operator : SJ/MA
Sample : I2006-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

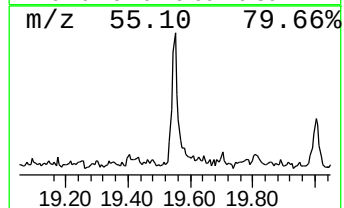
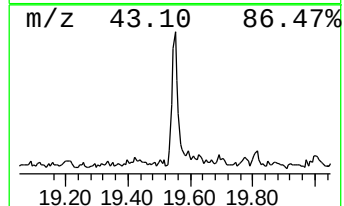
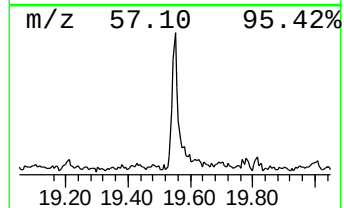
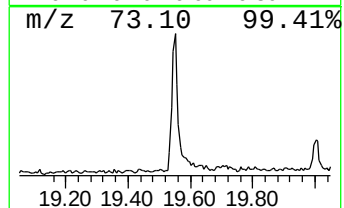
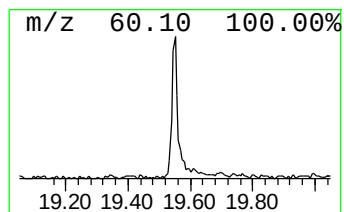
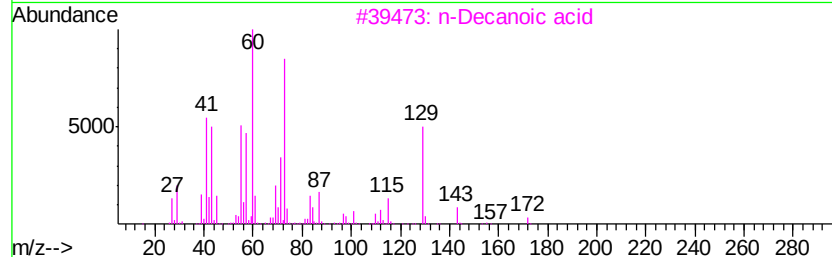
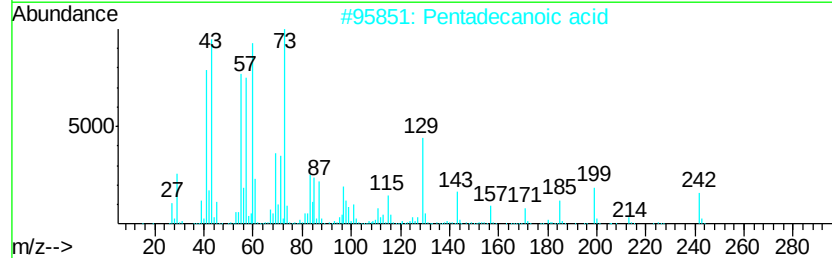
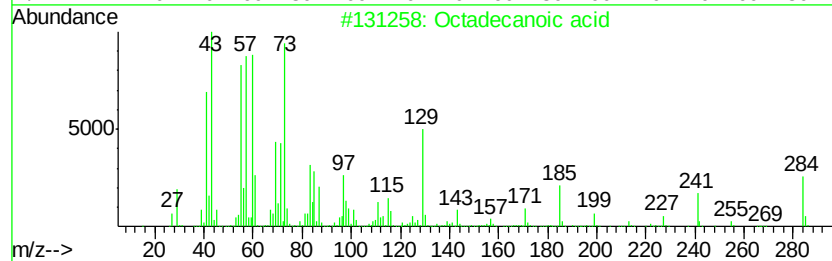
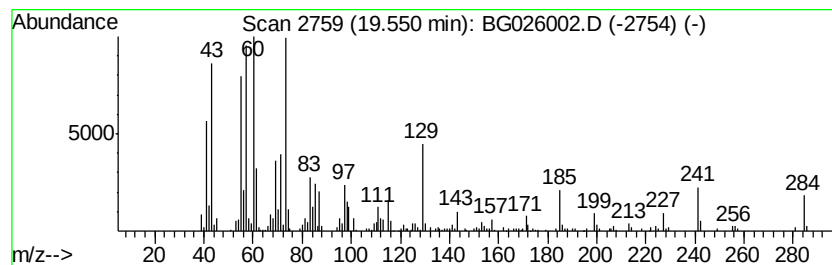
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
19.55	2.45 ng	259027	Chrysene-d12		21.65
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	96
3	n-Decanoic acid	172	C10H20O2	000334-48-5	96
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	83



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
Data File : BG026002.D
Acq On : 28 Feb 2017 18:19
Operator : SJ/MA
Sample : I2006-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

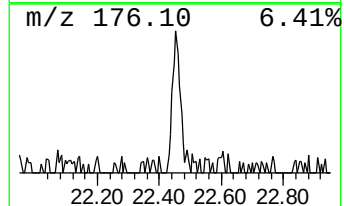
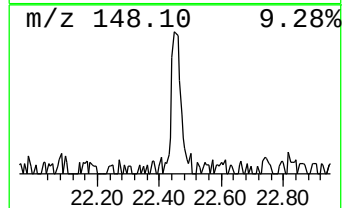
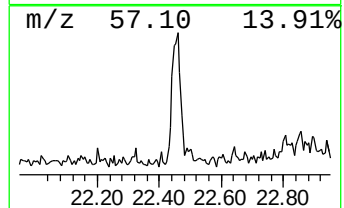
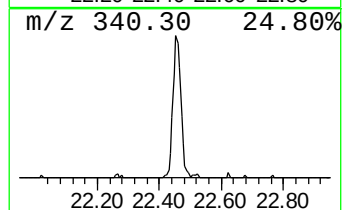
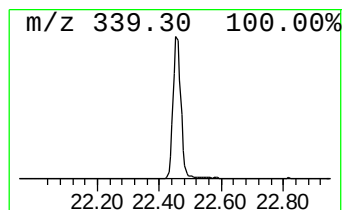
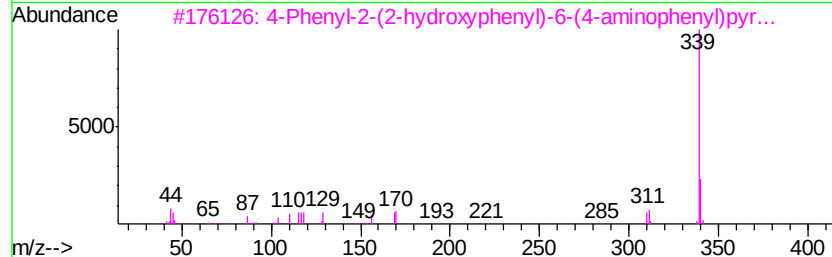
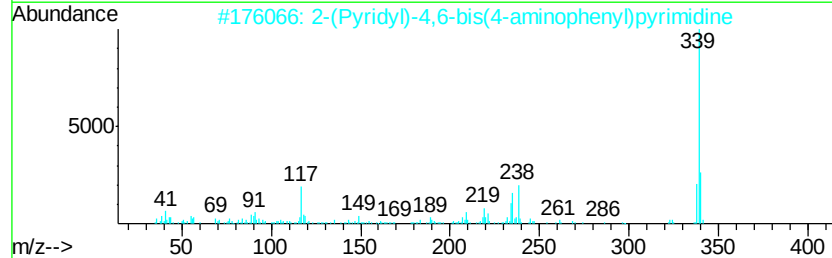
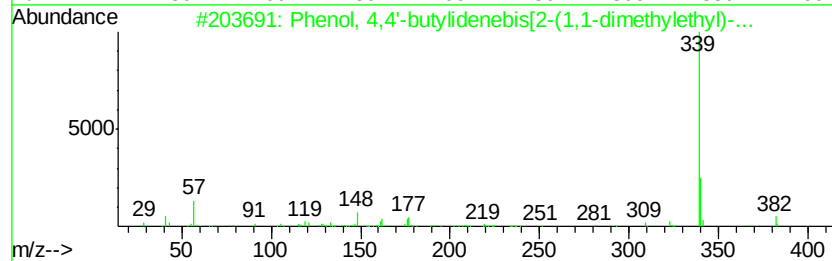
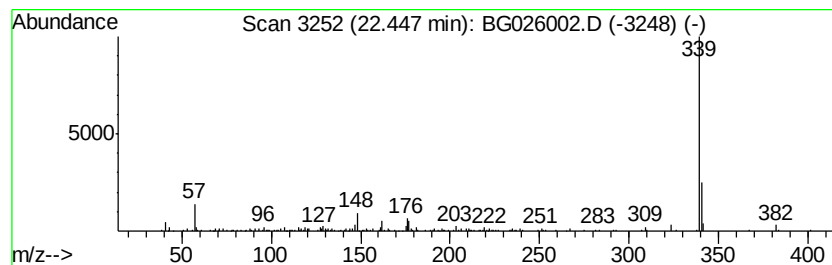
Instrument :
BNA_G
ClientSampleId :
C19013051

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Phenol, 4,4'-butylidenebis[... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
22.45	2.33 ng	245666	Chrysene-d12	21.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-butylidenebis[2-(1,...	382	C26H38O2	000085-60-9	95	
2	2-(Pyridyl)-4,6-bis(4-aminopheny...	339	C21H17N5	1000078-62-7	56	
3	4-Phenyl-2-(2-hydroxyphenyl)-6-(...	339	C22H17N3O	1000070-58-2	50	
4	2,4-Bis(4-aminophenyl)-6-(pyridy...	339	C21H17N5	136009-90-0	50	
5	Isoquinoline, 6,7-dimethoxy-1-me...	339	C20H21NO4	102012-79-3	50	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022817\
 Data File : BG026002.D
 Acq On : 28 Feb 2017 18:19
 Operator : SJ/MA
 Sample : I2006-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C19013051

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG021717.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
2-Pentanone, 4-hy...	5.17	12.0	ng	326583	1	8.08	543578	20.0
unknown7.61	7.61	103.5	ng	2812220	1	8.08	543578	20.0
Benzene, 1,2,4,5-...	10.28	4.4	ng	214063	2	10.88	977785	20.0
unknown10.93	10.93	2.2	ng	106725	2	10.88	977785	20.0
Benzene, (3-methy...	11.98	2.7	ng	132178	2	10.88	977785	20.0
Naphthalene, 1-me...	12.74	8.3	ng	405185	2	10.88	977785	20.0
Naphthalene, 1,6-...	14.00	4.8	ng	367315	3	14.68	1519150	20.0
Naphthalene, 2,6-...	14.23	2.6	ng	196290	3	14.68	1519150	20.0
Naphthalene, 1,6,...	15.29	3.0	ng	224032	3	14.68	1519150	20.0
1,1'-Biphenyl, 2-...	15.84	4.0	ng	301615	3	14.68	1519150	20.0
9H-Fluorene, 1-me...	16.77	2.4	ng	236825	4	17.41	1961290	20.0
n-Hexadecanoic acid	18.29	2.9	ng	281482	4	17.41	1961290	20.0
Octadecanoic acid	19.55	2.5	ng	259027	5	21.65	2110370	20.0
Phenol, 4,4'-buty...	22.45	2.3	ng	245666	5	21.65	2110370	20.0