

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
 Data File : BG025974.D
 Acq On : 27 Feb 2017 21:49
 Operator : SJ/MA
 Sample : I1978-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 UST-TWP

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	5.177	306	313	321	rVB	205603	316247	3.72%	0.726%
2	5.642	386	392	402	rBV	796981	1257785	14.79%	2.886%
3	7.228	654	662	671	rBV	538211	904120	10.63%	2.075%
4	7.610	718	727	735	rBV	1486663	2496412	29.36%	5.728%
5	8.074	799	806	815	rBV	308598	544208	6.40%	1.249%
6	8.403	854	862	869	rVB	1205647	2106228	24.77%	4.833%
7	9.232	995	1003	1012	rBV	1114777	1977279	23.26%	4.537%
8	10.877	1275	1283	1290	rBV	540910	988397	11.62%	2.268%
9	12.951	1630	1636	1637	rBV5	72622	118697	1.40%	0.272%
10	13.133	1662	1667	1671	rBV6	48315	88405	1.04%	0.203%
11	13.309	1690	1697	1702	rBV	3656941	5670877	66.70%	13.013%
12	13.356	1702	1705	1711	rVB2	247724	341946	4.02%	0.785%
13	13.515	1726	1732	1735	rBV8	44291	95679	1.13%	0.220%
14	13.568	1737	1741	1747	rVB5	85323	150500	1.77%	0.345%
15	13.638	1749	1753	1758	rVB6	68375	114428	1.35%	0.263%
16	13.803	1774	1781	1784	rBV4	104496	238687	2.81%	0.548%
17	14.126	1831	1836	1840	rBV	144581	215726	2.54%	0.495%
18	14.202	1845	1849	1859	rVB5	121029	296294	3.48%	0.680%
19	14.367	1874	1877	1889	rVB5	80665	214094	2.52%	0.491%
20	14.467	1889	1894	1895	rBV4	73845	105968	1.25%	0.243%
21	14.584	1910	1914	1919	rBV6	85875	218641	2.57%	0.502%
22	14.678	1924	1930	1935	rVV2	952416	1603704	18.86%	3.680%
23	14.725	1935	1938	1941	rVB	150426	194840	2.29%	0.447%
24	14.837	1953	1957	1963	rBV8	78628	134283	1.58%	0.308%
25	14.995	1979	1984	1990	rBV4	213282	385134	4.53%	0.884%
26	15.084	1997	1999	2004	rVB	100657	125131	1.47%	0.287%
27	15.454	2057	2062	2070	rVB6	354860	673072	7.92%	1.544%
28	15.665	2095	2098	2102	rBV2	164750	219997	2.59%	0.505%
29	15.718	2103	2107	2114	rBV3	278746	550588	6.48%	1.263%
30	16.006	2152	2156	2162	rVB5	161665	250267	2.94%	0.574%
31	16.159	2176	2182	2186	rBV	3051016	4740391	55.75%	10.878%
32	17.017	2324	2328	2333	rBV4	145233	319854	3.76%	0.734%
33	17.346	2381	2384	2388	rBV3	104832	157510	1.85%	0.361%
34	17.410	2389	2395	2408	rVB2	1350101	2244231	26.39%	5.150%

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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

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

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35	18.297	2543	2546	2550	rVB	213560	243676	2.87%	0.559%
36	19.337	2721	2723	2728	rVB2	115979	151215	1.78%	0.347%
37	19.555	2757	2760	2766	rVB3	182349	242711	2.85%	0.557%
38	20.007	2831	2837	2842	rVB	6476238	8502499	100.00%	19.510%
39	21.652	3111	3117	3124	rVB	1363355	2181154	25.65%	5.005%
40	24.854	3653	3662	3673	rVB2	768845	2198325	25.86%	5.044%

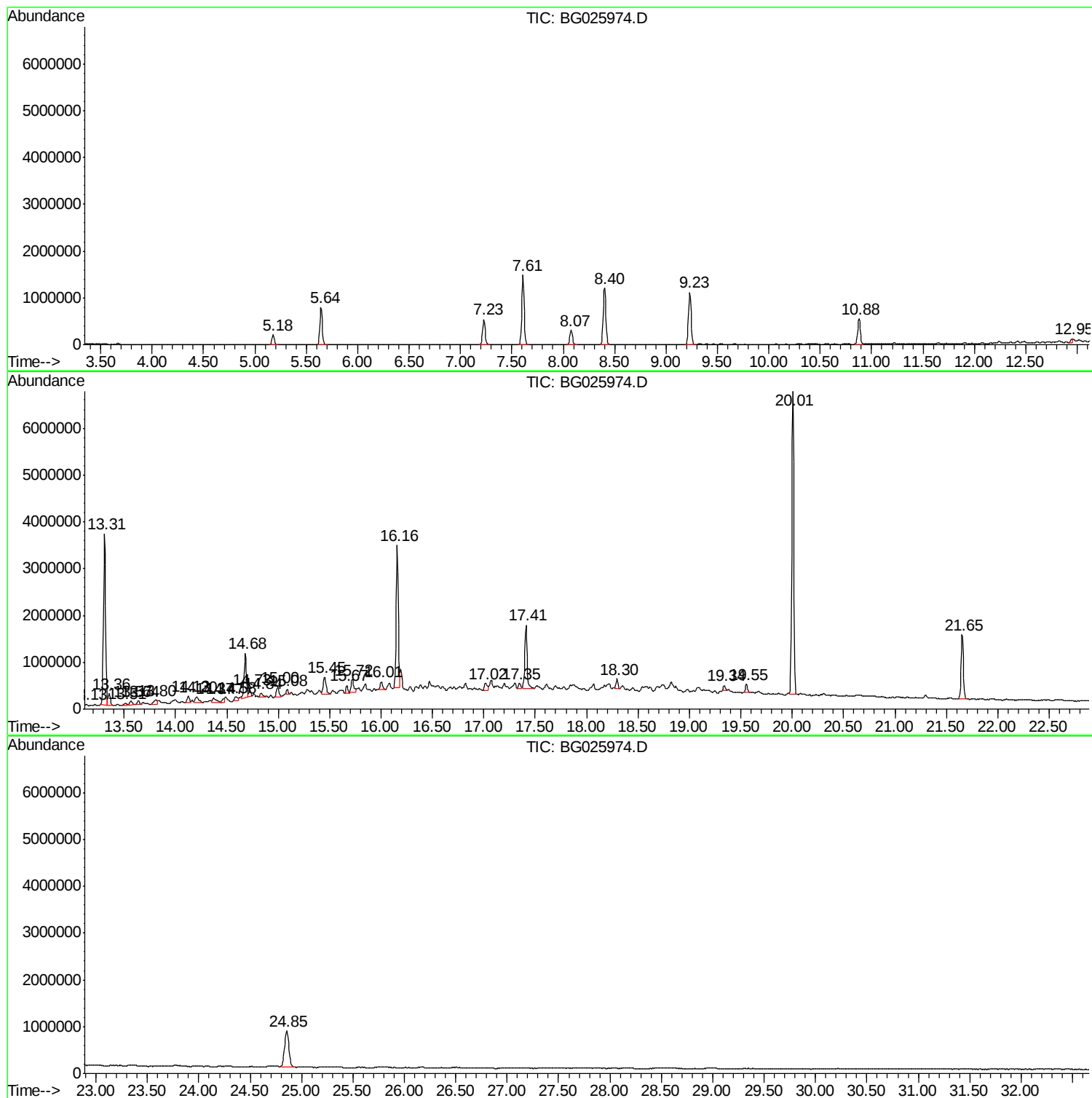
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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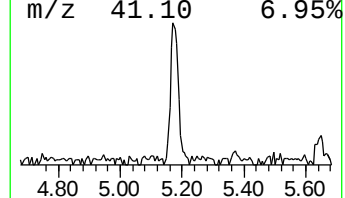
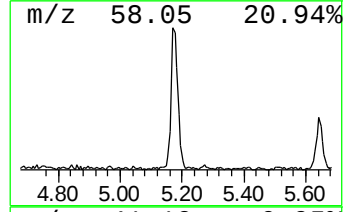
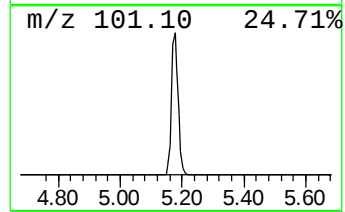
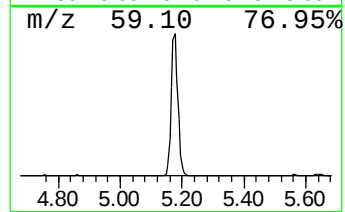
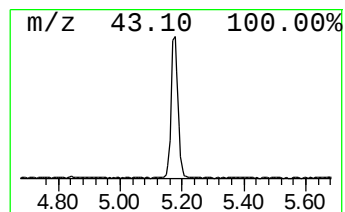
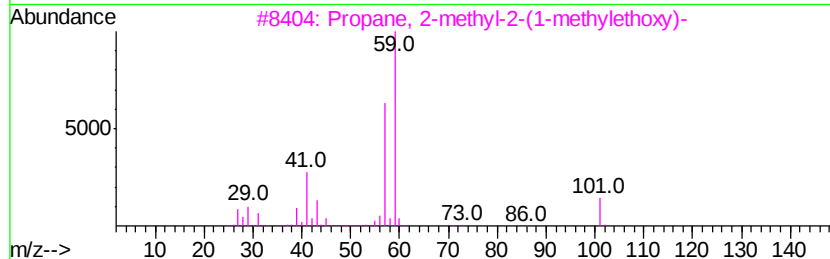
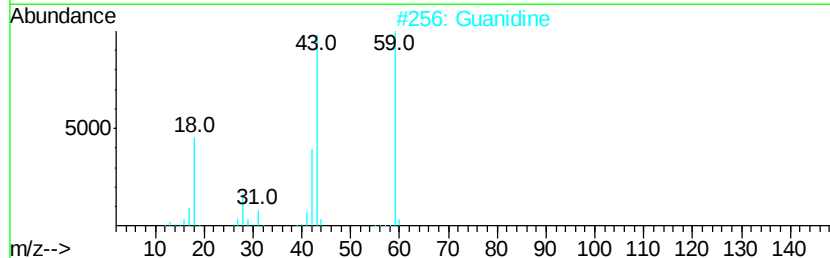
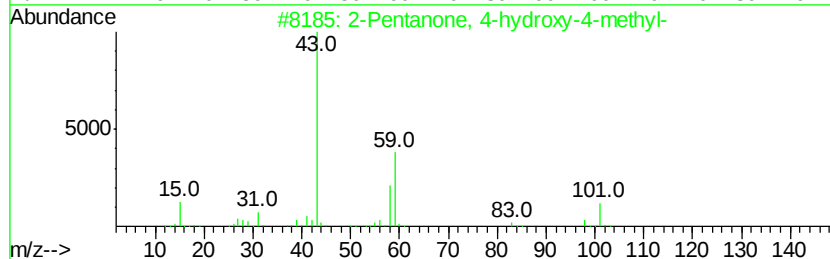
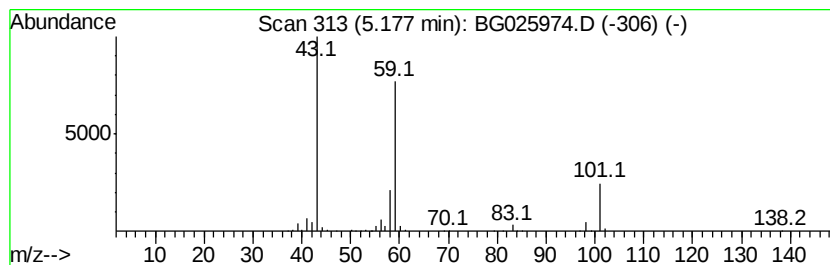
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.18	11.62 ng	316247	1,4-Dichlorobenzene-d4	8.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Guanidine	59	CH5N3	000113-00-8	47
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	28
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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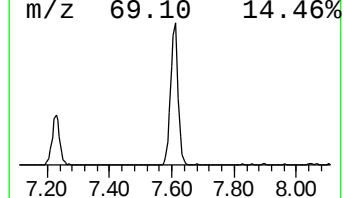
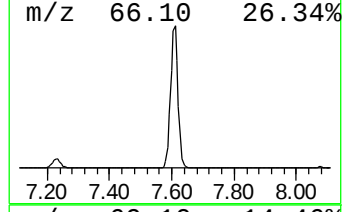
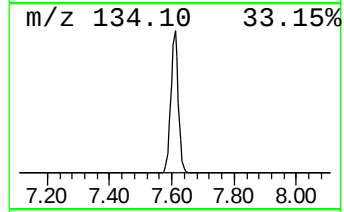
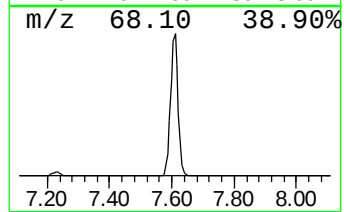
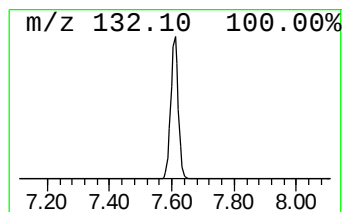
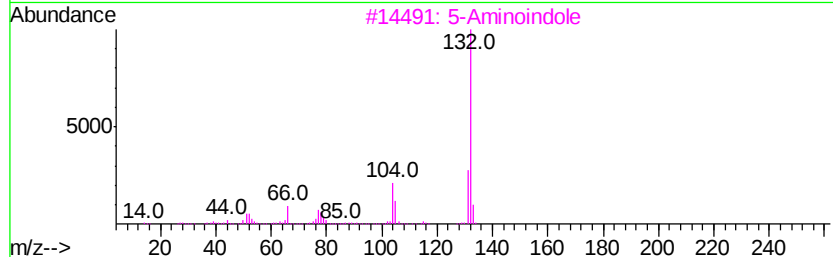
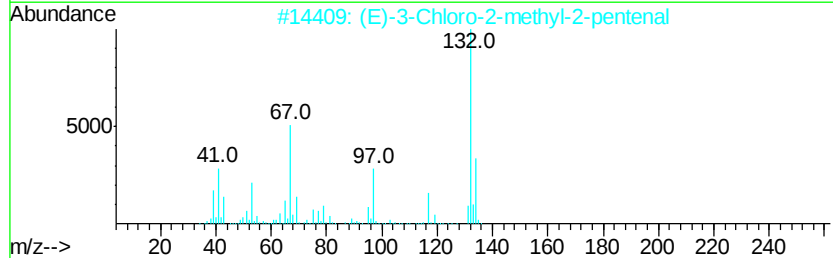
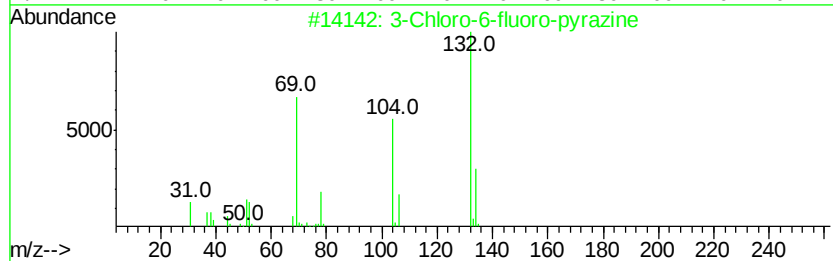
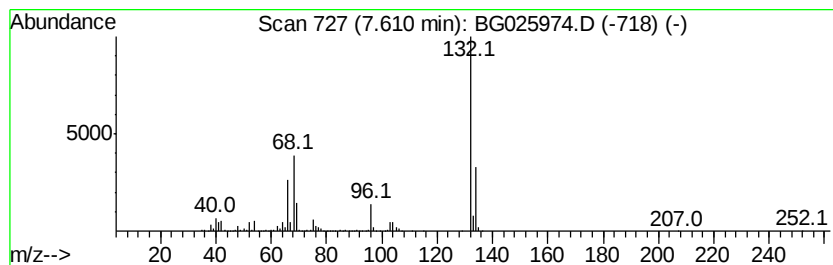
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	91.74 ng	2496410	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		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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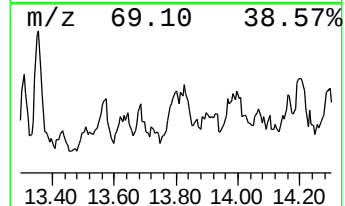
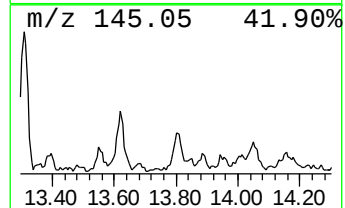
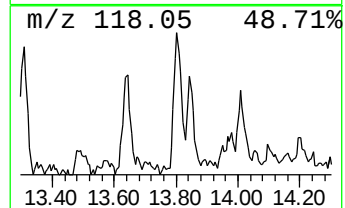
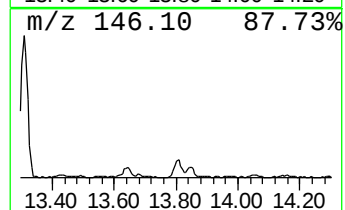
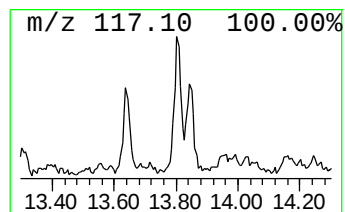
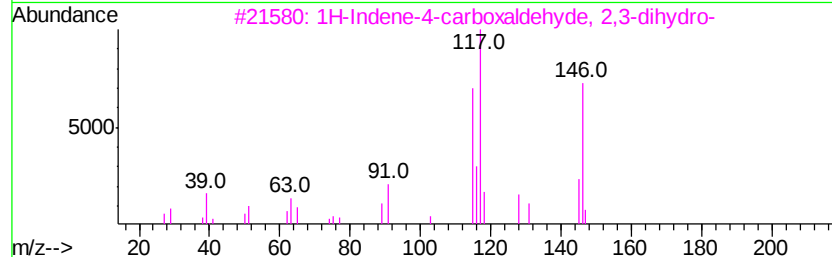
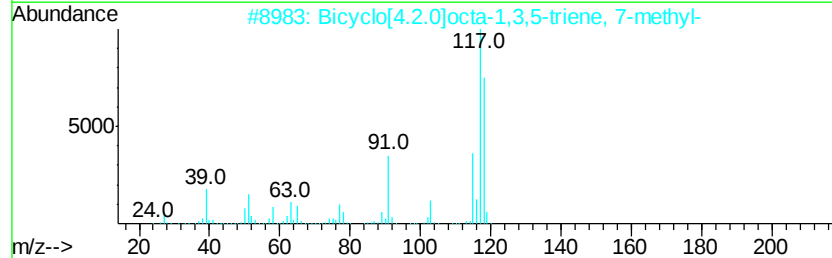
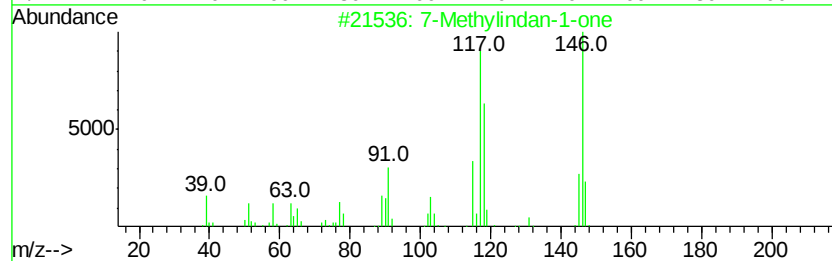
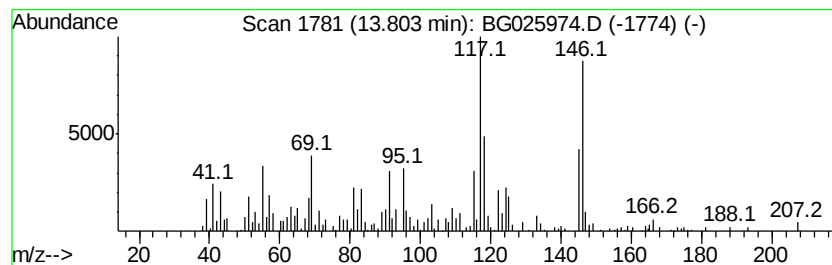
Instrument :
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ClientSampled :
UST-TWP

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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 4 7-Methylindan-1-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	2.98 ng	238687	Acenaphthene-d10	14.68
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	7-Methylindan-1-one	146 C10H10O	039627-61-7	83
2	Bicyclo[4.2.0]octa-1,3,5-triene,...	118 C9H10	055337-80-9	51
3	1H-Indene-4-carboxaldehyde, 2,3-...	146 C10H10O	051932-70-8	49
4	1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3	46
5	Benzeneacetaldehyde, .alpha.-eth...	146 C10H10O	004411-89-6	46



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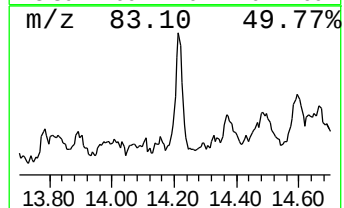
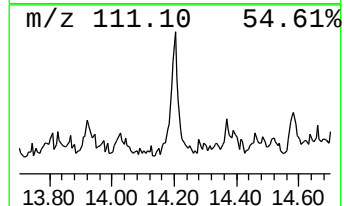
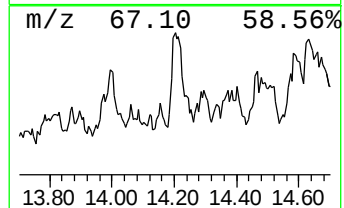
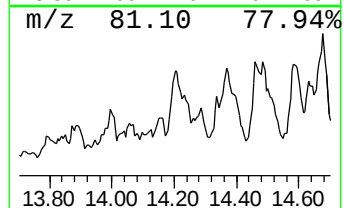
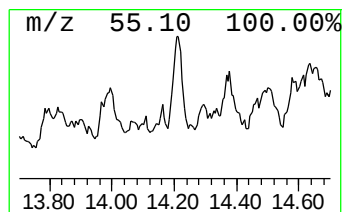
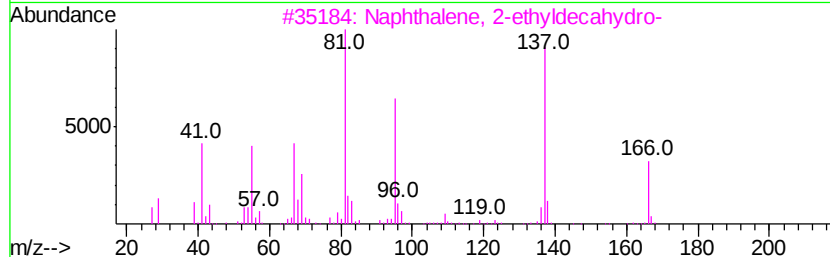
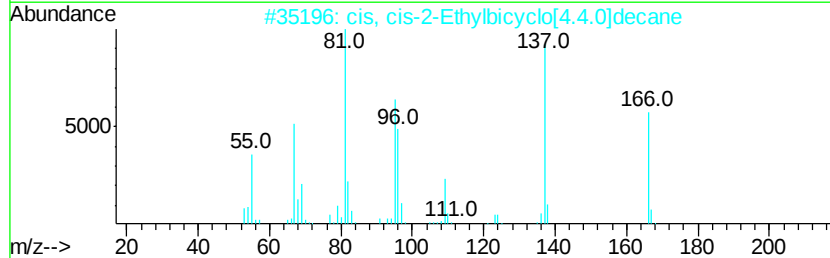
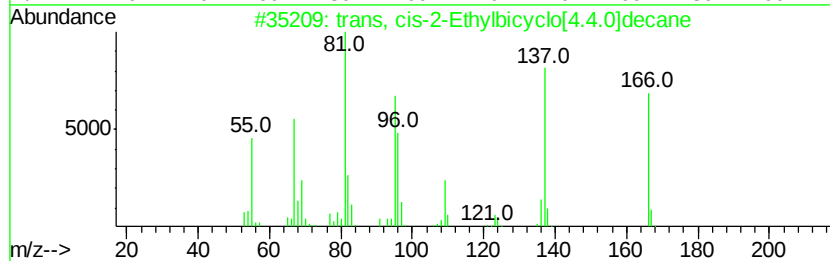
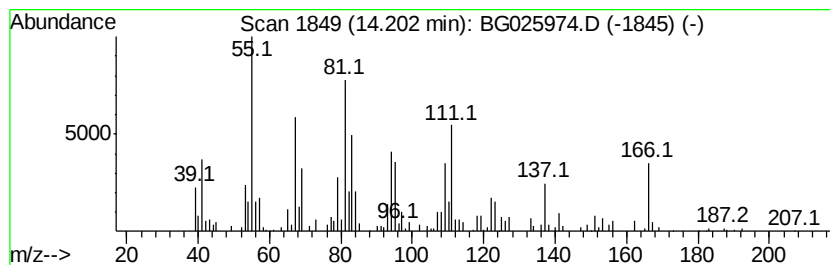
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 trans, cis-2-Ethylbicyclo[4... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.20	3.70 ng	296294	Acenaphthene-d10	14.68	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans, cis-2-Ethylbicyclo[4.4.0]...	166	C12H22	066660-39-7	55
2	cis, cis-2-Ethylbicyclo[4.4.0]de...	166	C12H22	066660-40-0	49
3	Naphthalene, 2-ethyldecahydro-	166	C12H22	001618-23-1	43
4	4-Cyclohexylidene-n-butanol	154	C10H18O	004441-58-1	38
5	1H-Imidazole-1-propanoic acid,	156	C6H8N2O3	051103-59-4	30



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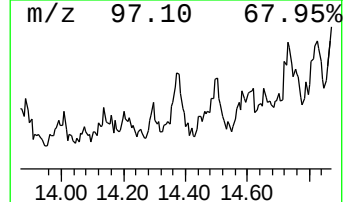
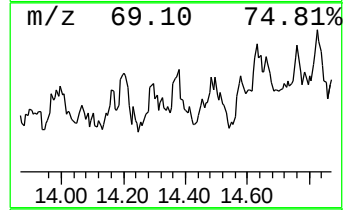
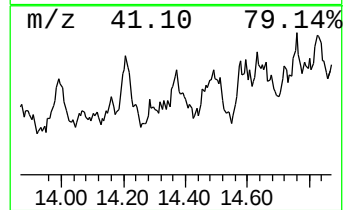
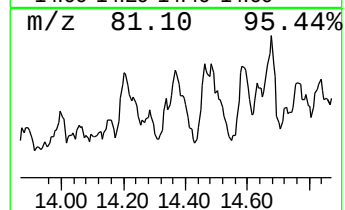
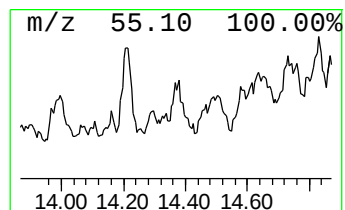
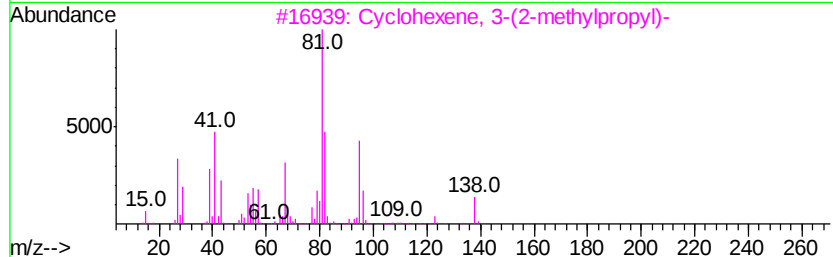
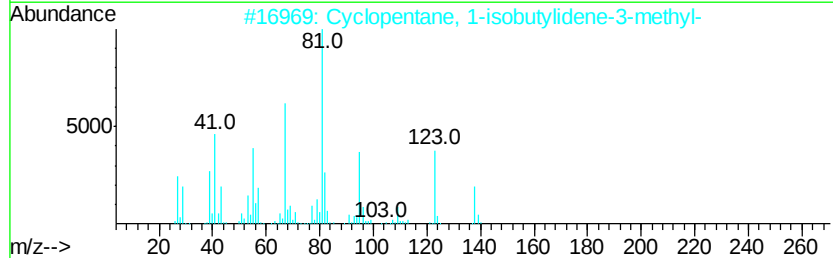
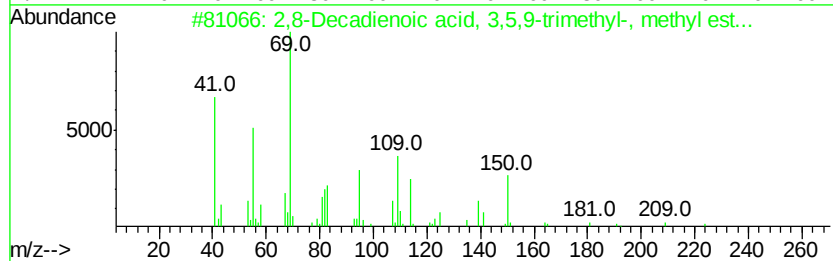
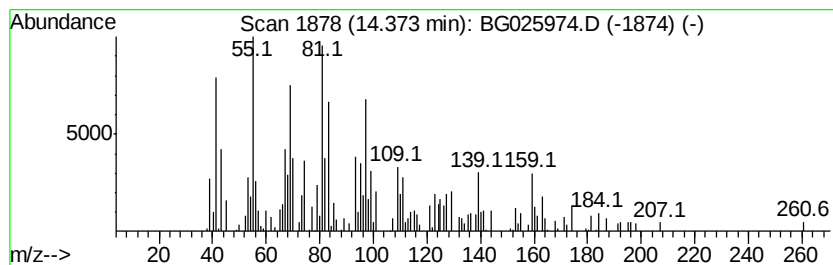
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TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown14.37 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.37	2.67 ng	214094	Acenaphthene-d10	14.68	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,8-Decadienoic acid, 3,5,9-trim...	224	C14H24O2	055283-32-4	38
2	Cyclopentane, 1-isobutylidene-3-...	138	C10H18	1000150-62-1	25
3	Cyclohexene, 3-(2-methylpropyl)-	138	C10H18	004104-56-7	25
4	Cyclohexadecane	224	C16H32	000295-65-8	25
5	Pentadec-7-ene, 7-bromomethyl-	302	C16H31Br	1000259-58-5	25



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
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Operator : SJ/MA
Sample : I1978-02
Misc :
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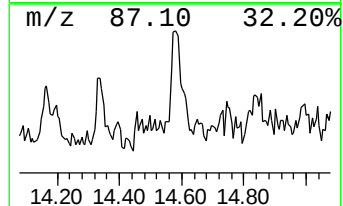
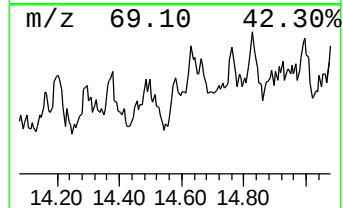
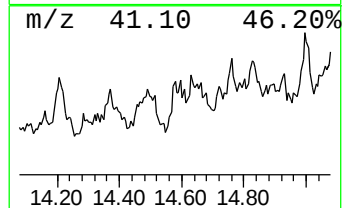
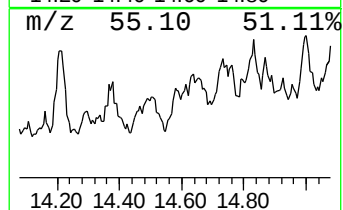
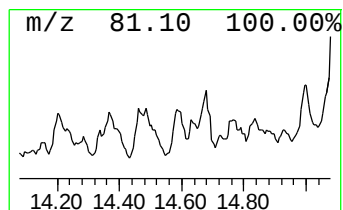
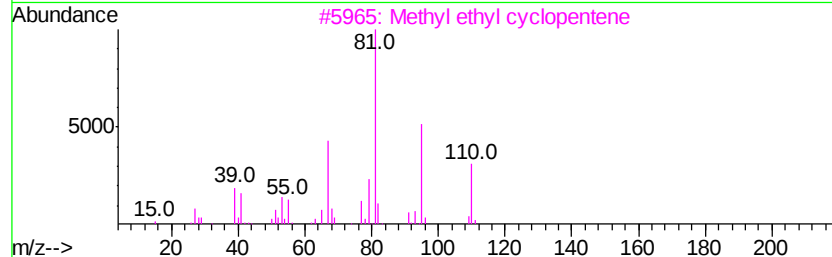
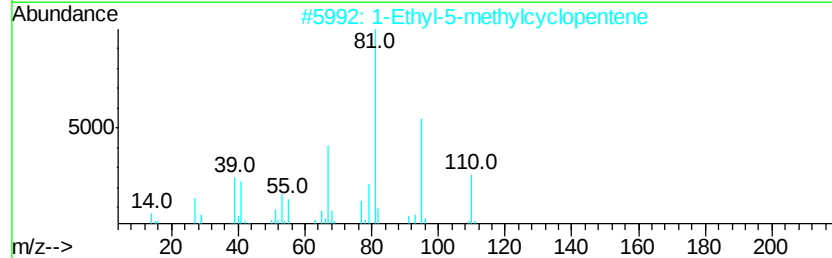
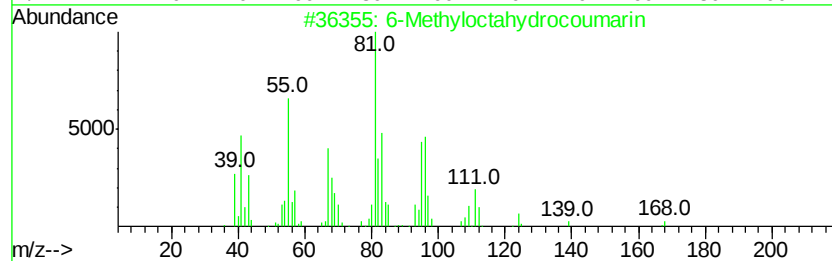
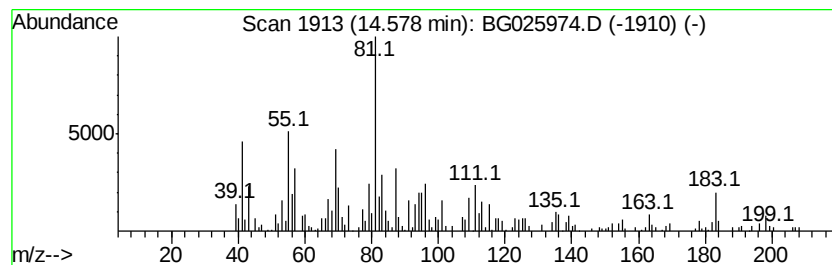
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 6-Methyloctahydrocoumarin Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.58	2.73 ng	218641	Acenaphthene-d10	14.68
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	6-Methyloctahydrocoumarin	168 C10H16O2	080648-29-9	59
2	1-Ethyl-5-methylcyclopentene	110 C8H14	097797-57-4	35
3	Methyl ethyl cyclopentene	110 C8H14	019780-56-4	35
4	Ethanone, 1-(3-methylenecyclopent...	124 C8H12O	054829-98-0	27
5	Bicyclo[3.1.0]hexan-2-one, 5-(1-...	138 C9H14O	000513-20-2	25



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
Data File : BG025974.D
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ClientSampleId :
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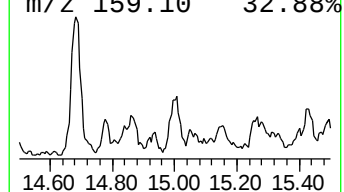
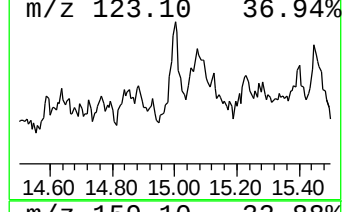
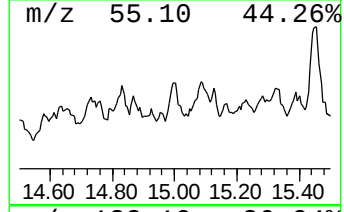
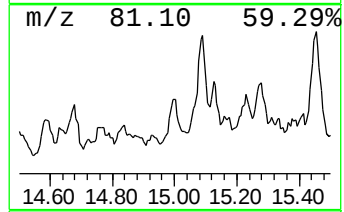
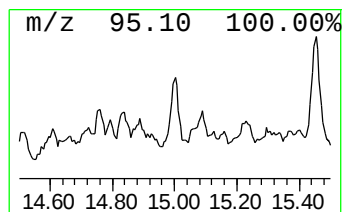
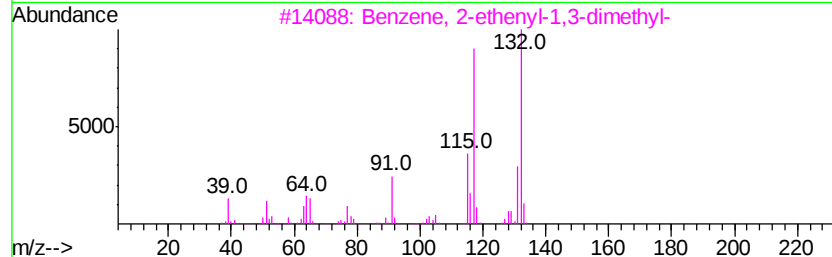
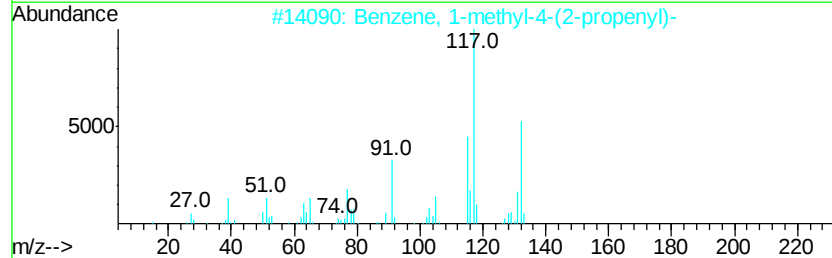
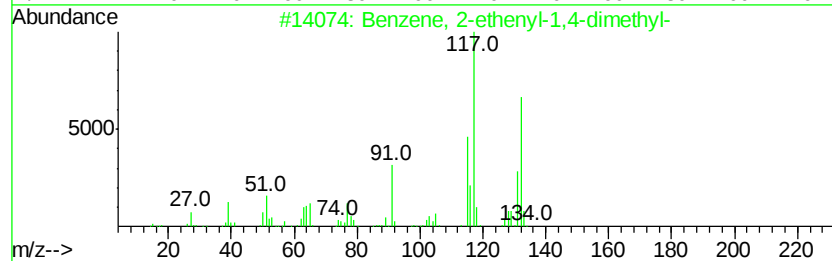
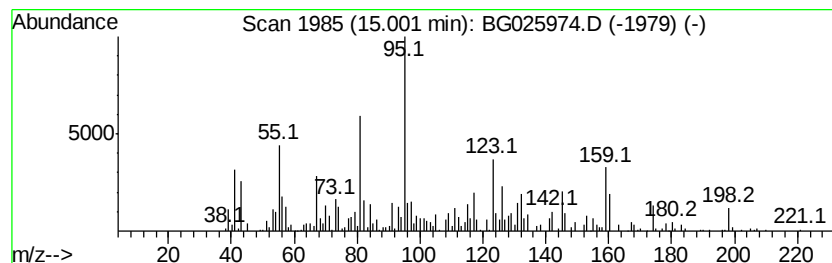
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown15.00 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	4.80 ng	385134	Acenaphthene-d10	14.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	35
2			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	35
3			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	35
4			1,3-Naphthalenediol	160	C10H8O2	000132-86-5	25
5			1H-Inden-1-one, 2,3-dihydro-4,7-...	160	C11H12O	005037-60-5	25



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
Data File : BG025974.D
Acq On : 27 Feb 2017 21:49
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Sample : I1978-02
Misc :
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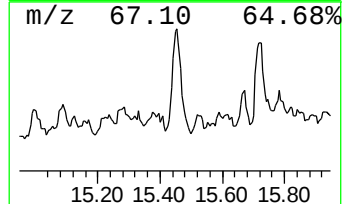
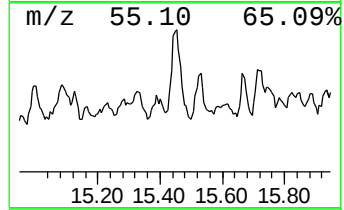
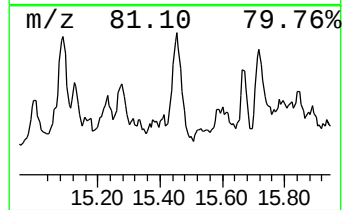
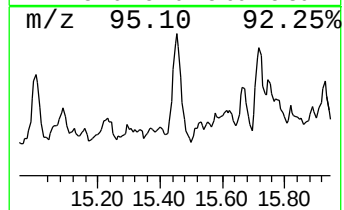
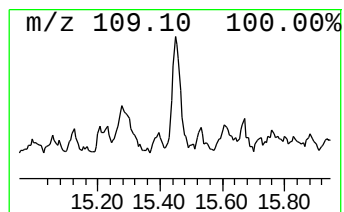
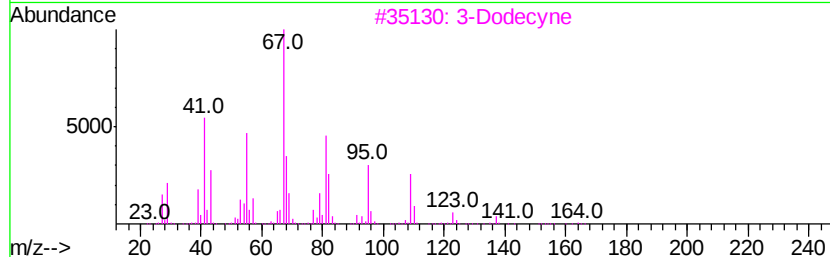
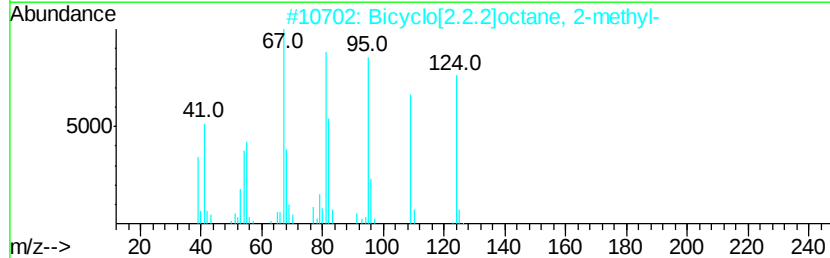
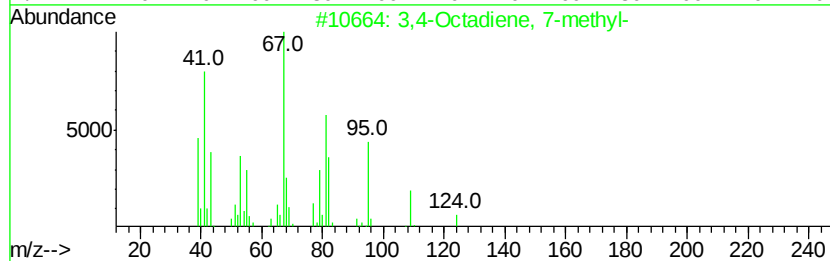
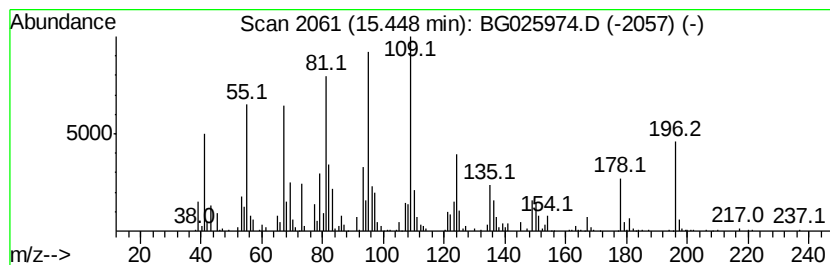
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 3,4-Octadiene, 7-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	8.39 ng	673072	Acenaphthene-d10	14.68
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	3,4-Octadiene, 7-methyl-	124 C9H16	037050-05-8	86
2	Bicyclo[2.2.2]octane, 2-methyl-	124 C9H16	000766-53-0	53
3	3-Dodecyne	166 C12H22	006790-27-8	38
4	Cyclohexene, 3,3,5-trimethyl-	124 C9H16	000503-45-7	30
5	Cyclohexane, (1-methylethylidene)-	124 C9H16	005749-72-4	30



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
Data File : BG025974.D
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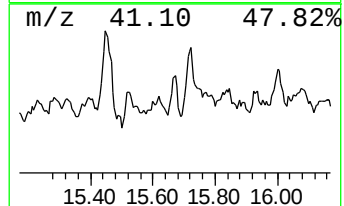
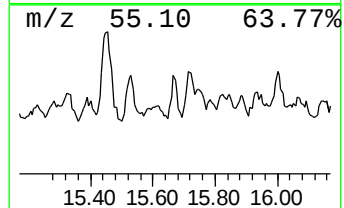
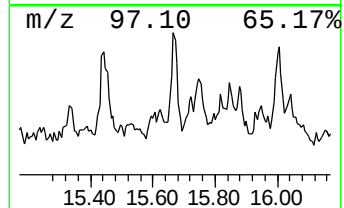
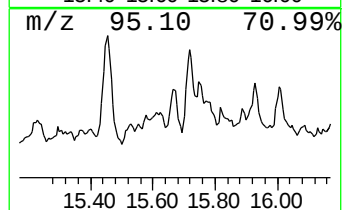
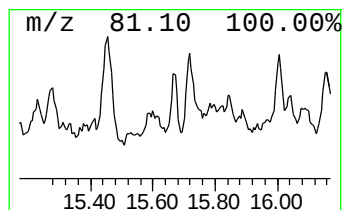
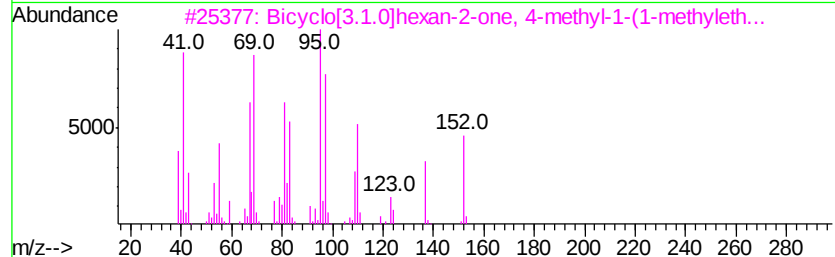
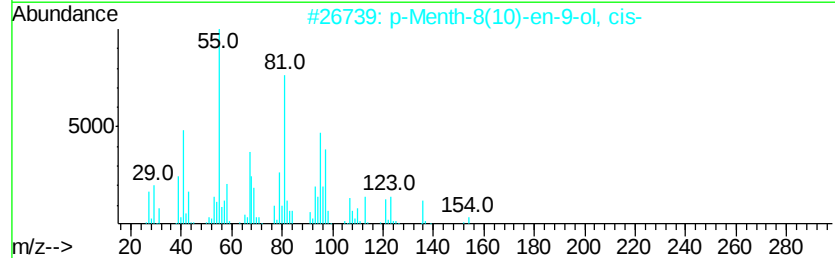
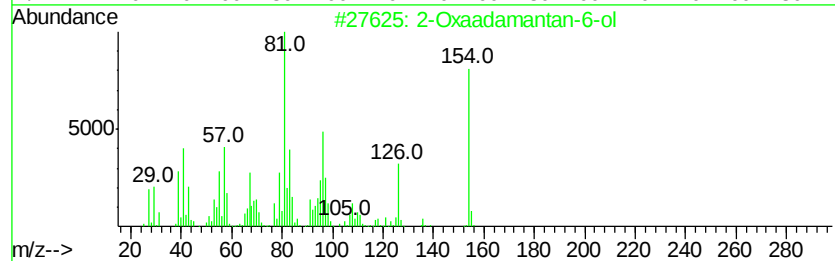
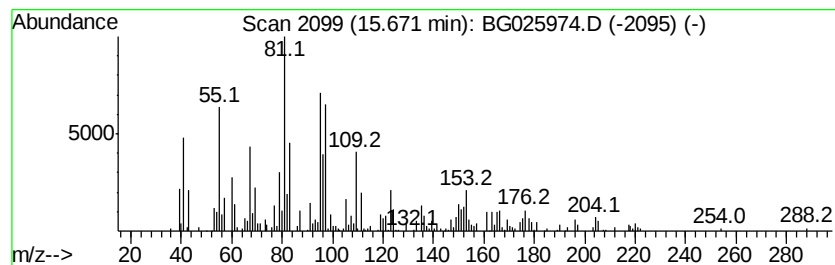
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-Oxaadamantan-6-ol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.67	2.74 ng	219997	Acenaphthene-d10	14.68	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Oxaadamantan-6-ol	154	C9H14O2	058512-55-3	53
2	p-Menth-8(10)-en-9-ol, cis-	154	C10H18O	015714-13-3	46
3	Bicyclo[3.1.0]hexan-2-one, 4-met...	152	C10H16O	002506-61-8	46
4	Cyclohexene, 3-pentyl-	152	C11H20	015232-92-5	45
5	Cyclopropane, 1-(1-methylethyl)-...	152	C11H20	056259-17-7	38



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
Data File : BG025974.D
Acq On : 27 Feb 2017 21:49
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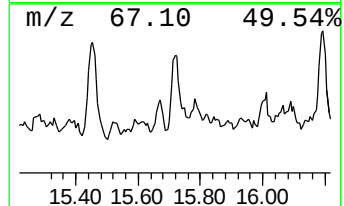
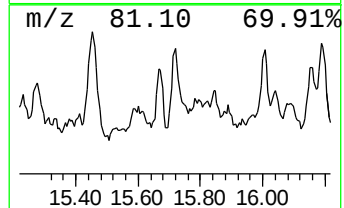
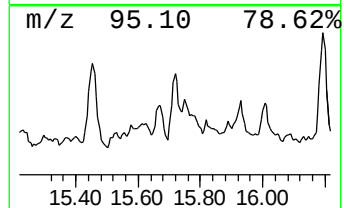
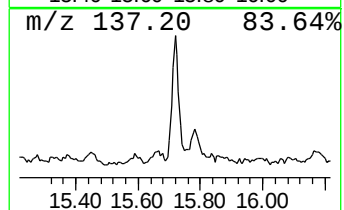
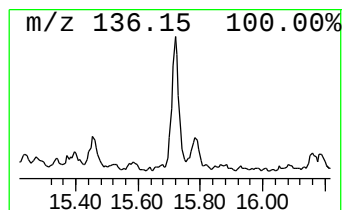
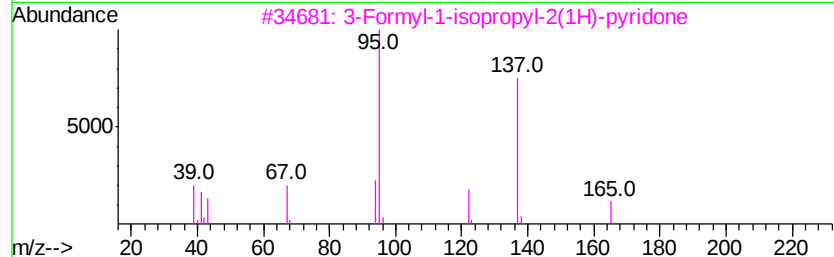
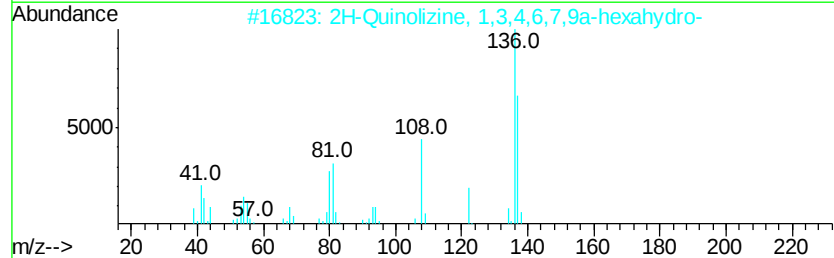
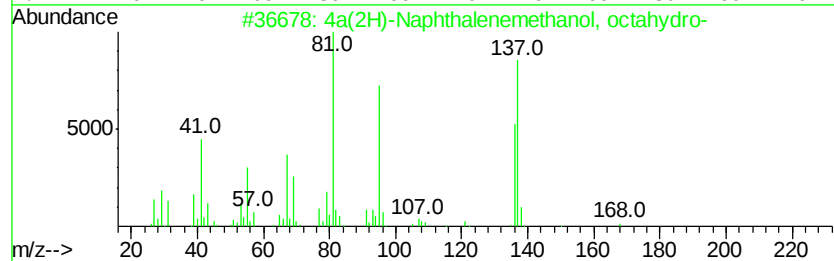
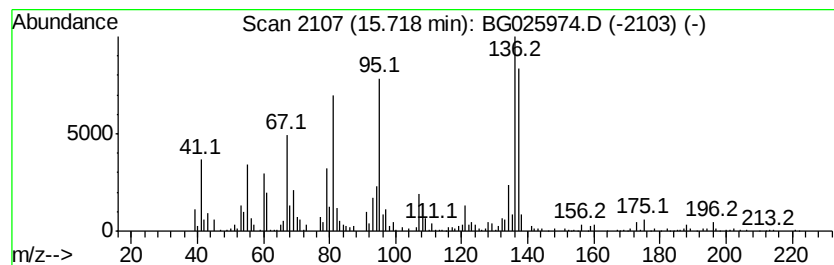
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown15.72 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	6.87 ng	550588	Acenaphthene-d10	14.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4a(2H)-Naphthalenemethanol, octa...	168	C11H20O	099992-19-5	43
2			2H-Quinolizine, 1,3,4,6,7,9a-hex...	137	C9H15N	001004-90-6	38
3			3-Formyl-1-isopropyl-2(1H)-pyridone	165	C9H11NO2	1000306-46-8	38
4			Cyclopenta[c]pyrazol-3-amine, 2,...	137	C7H11N3	1000351-26-4	38
5			Phenol, 3-(dimethylamino)-	137	C8H11NO	000099-07-0	38



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
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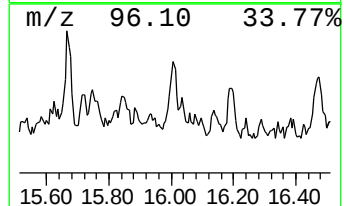
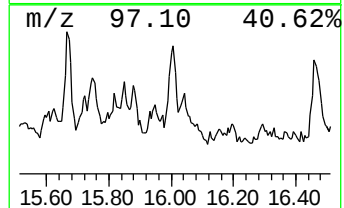
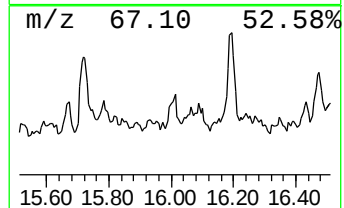
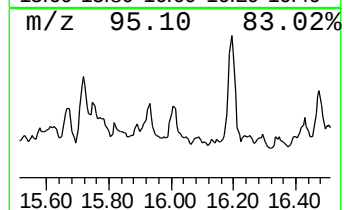
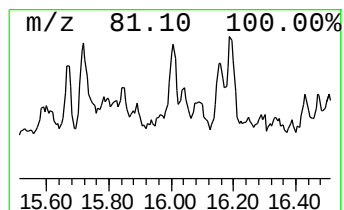
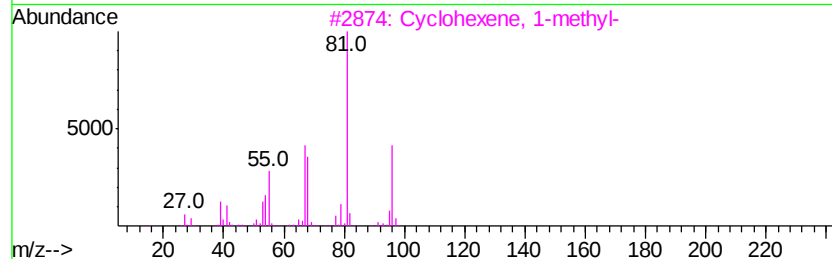
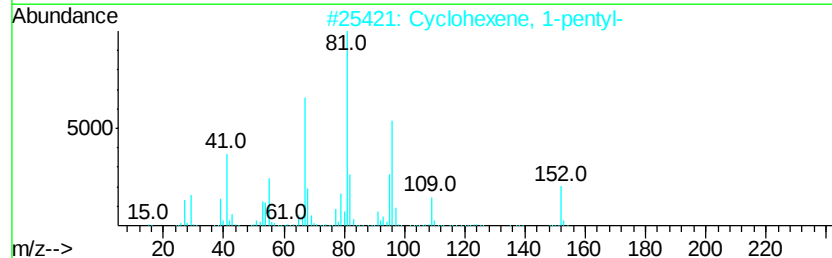
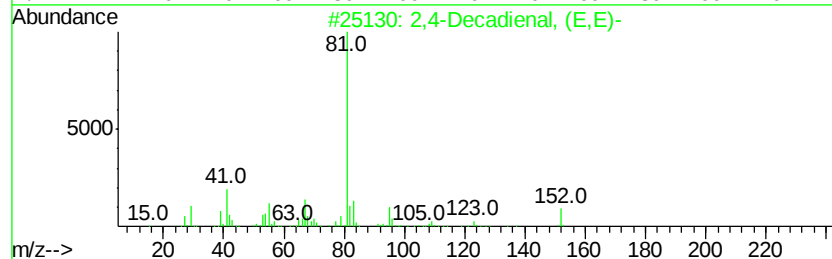
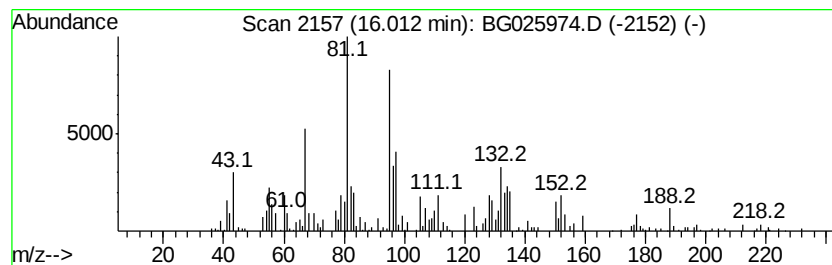
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown16.01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.01	3.12 ng	250267	Acenaphthene-d10	14.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Decadienal, (E,E)-	152	C10H16O	025152-84-5	38
2		Cyclohexene, 1-pentyl-	152	C11H20	015232-85-6	35
3		Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	30
4		2,4-Heptadiene, (E,E)-	96	C7H12	002384-94-3	27
5		Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	27



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
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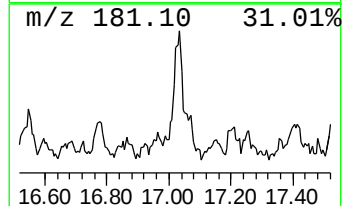
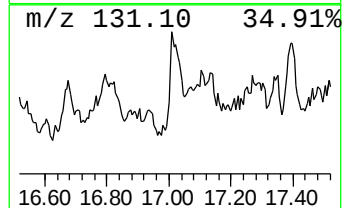
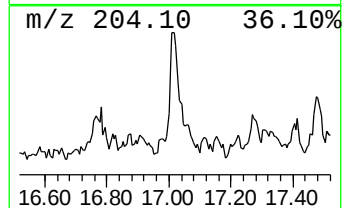
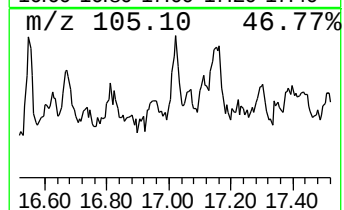
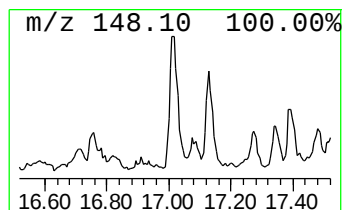
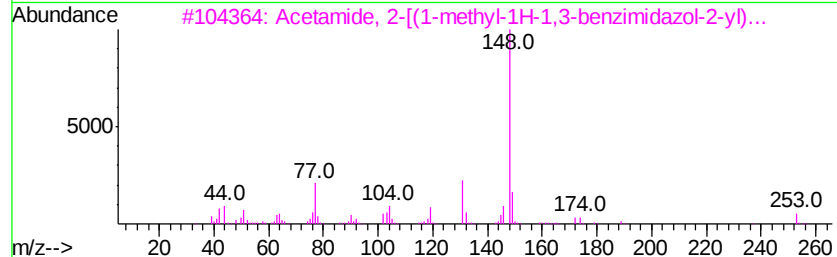
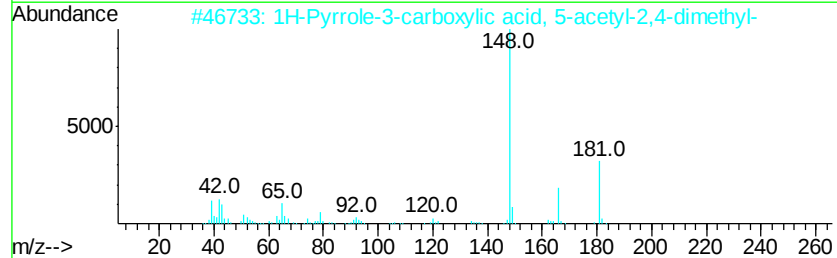
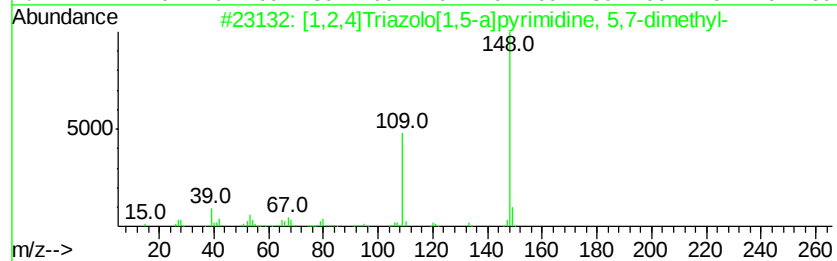
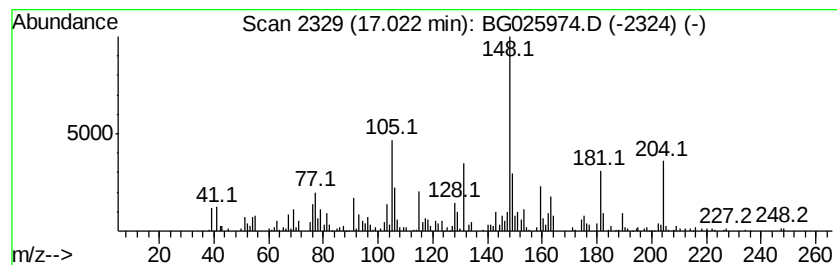
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown17.02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.02	2.85 ng	319854	Phenanthrene-d10	17.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	[1,2,4]Triazolo[1,5-a]pyrimidine...	148	C7H8N4	007681-99-4	38
2		1H-Pyrrole-3-carboxylic acid, 5-...	181	C9H11N03	1000337-47-2	38
3		Acetamide, 2-[(1-methyl-1H-1,3-b...	253	C10H11N3O3S	1000320-08-0	37
4		Oxalic acid, monoamide, N-(2-eth...	361	C22H35N03	1000309-36-5	35
5		[1,2,4]-Triazolo[2,3-a]pyrimidin...	248	C12H16N4O2	300361-61-9	35



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
Data File : BG025974.D
Acq On : 27 Feb 2017 21:49
Operator : SJ/MA
Sample : I1978-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
UST-TWP

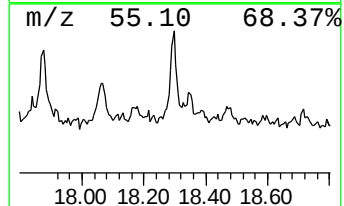
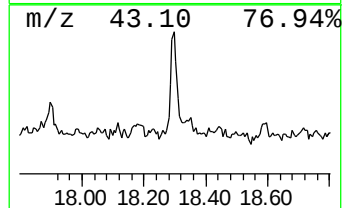
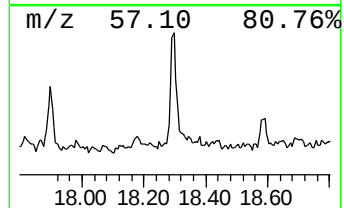
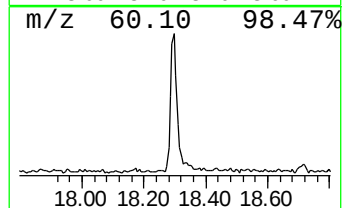
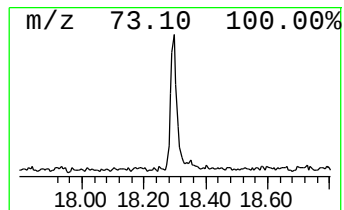
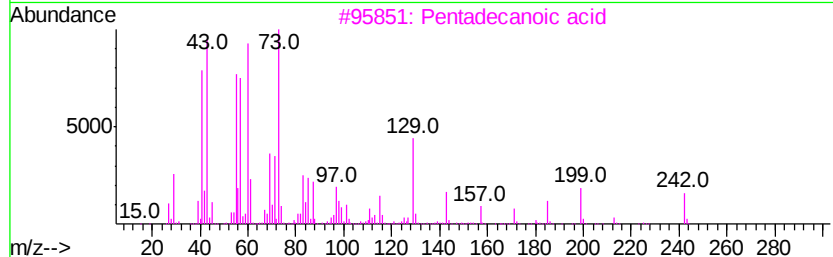
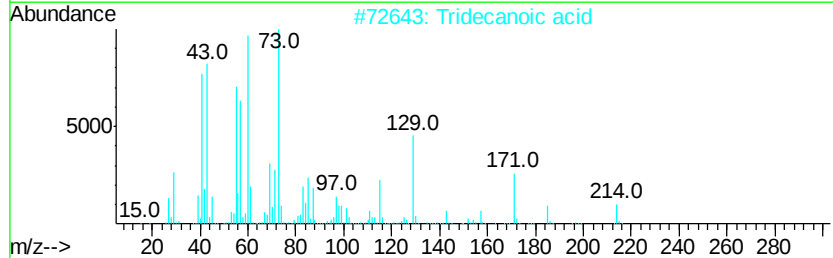
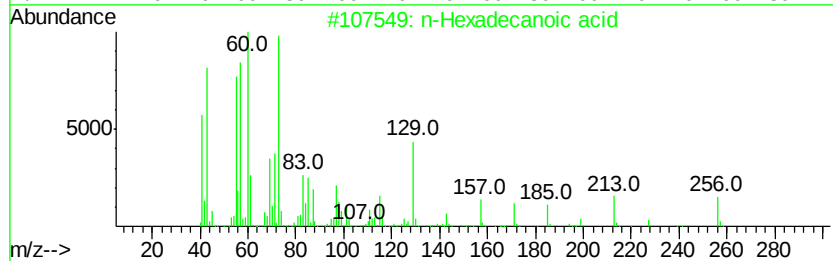
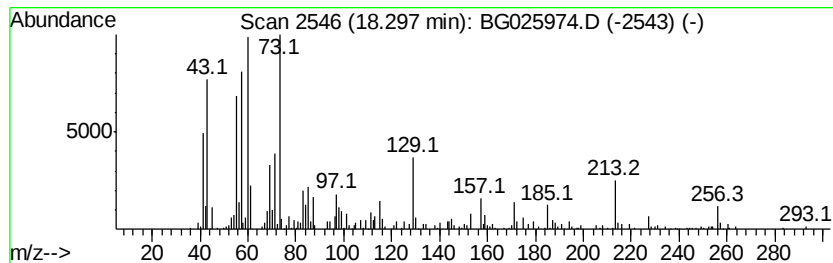
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.30	2.17 ng	243676	Phenanthrene-d10	17.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	94
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
Data File : BG025974.D
Acq On : 27 Feb 2017 21:49
Operator : SJ/MA
Sample : I1978-02
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
UST-TWP

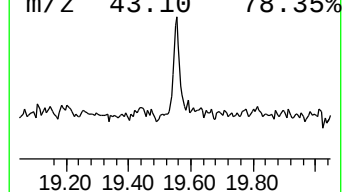
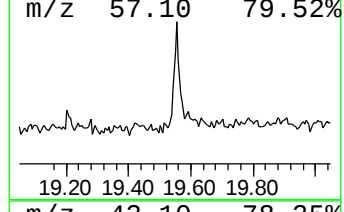
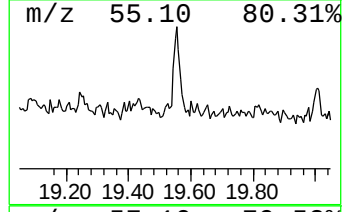
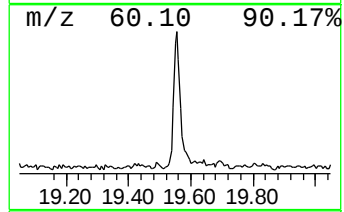
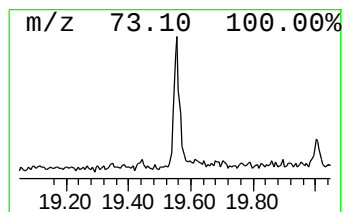
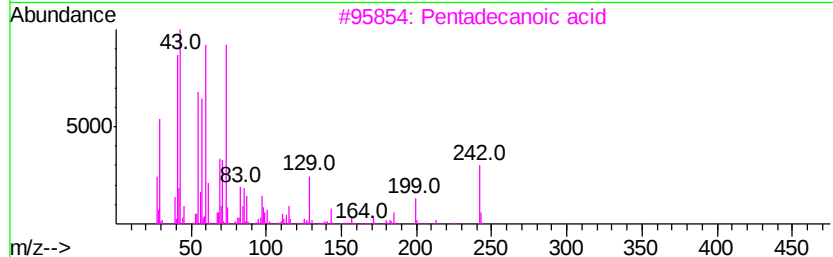
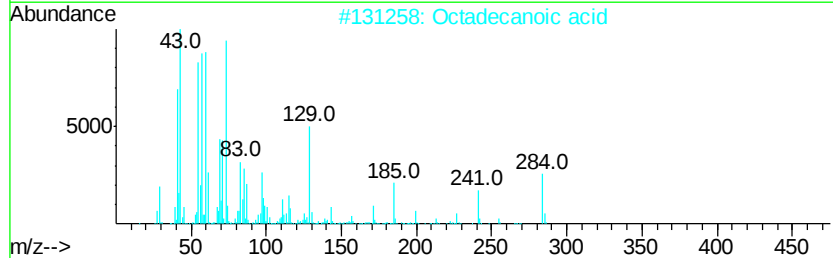
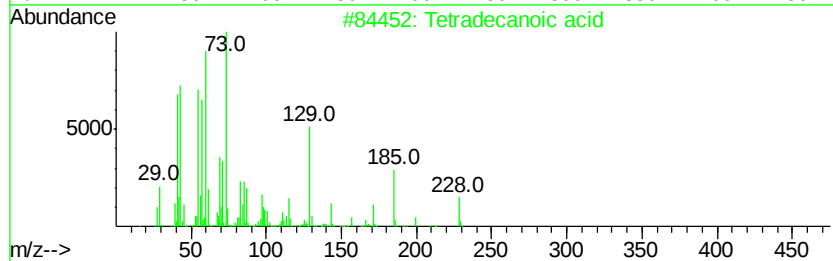
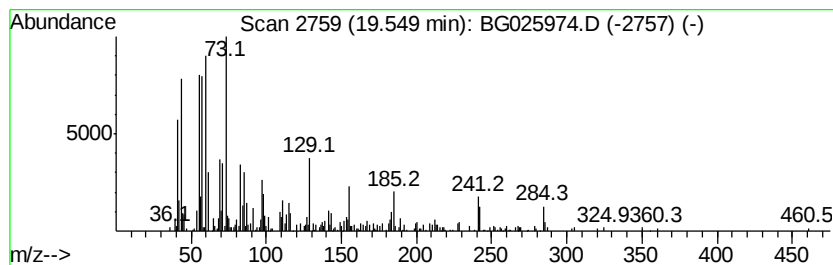
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetradecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.55	2.23 ng	242711	Chrysene-d12	21.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanoic acid	228	C14H28O2	000544-63-8	92
2		Octadecanoic acid	284	C18H36O2	000057-11-4	90
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	80
5		Tridecanoic acid	214	C13H26O2	000638-53-9	76



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022717\
 Data File : BG025974.D
 Acq On : 27 Feb 2017 21:49
 Operator : SJ/MA
 Sample : I1978-02
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 UST-TWP

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...	5.18	11.6 ng		316247	1	8.08	544208	20.0
unknown7.61	7.61	91.7 ng		2496410	1	8.08	544208	20.0
7-Methylindan-1-one	13.80	3.0 ng		238687	3	14.68	1603700	20.0
trans, cis-2-Ethy...	14.20	3.7 ng		296294	3	14.68	1603700	20.0
unknown14.37	14.37	2.7 ng		214094	3	14.68	1603700	20.0
6-Methyloctahydro...	14.58	2.7 ng		218641	3	14.68	1603700	20.0
unknown15.00	15.00	4.8 ng		385134	3	14.68	1603700	20.0
3,4-Octadiene, 7-...	15.45	8.4 ng		673072	3	14.68	1603700	20.0
2-Oxaadamantan-6-ol	15.67	2.7 ng		219997	3	14.68	1603700	20.0
unknown15.72	15.72	6.9 ng		550588	3	14.68	1603700	20.0
unknown16.01	16.01	3.1 ng		250267	3	14.68	1603700	20.0
unknown17.02	17.02	2.9 ng		319854	4	17.41	2244230	20.0
n-Hexadecanoic acid	18.30	2.2 ng		243676	4	17.41	2244230	20.0
Tetradecanoic acid	19.55	2.2 ng		242711	5	21.66	2181150	20.0