

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061816\
 Data File : BF088159.D
 Acq On : 19 Jun 2016 3:47
 Operator : UM/SJ
 Sample : H3628-03
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-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_F\METHODS\8270-BF060716.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.667	6	7	11	rBV	18550	26030	1.21%	0.154%
2	1.781	15	17	46	rVB	327564	545204	25.29%	3.231%
3	4.833	282	284	290	rBV	34145	51893	2.41%	0.308%
4	5.278	321	323	336	rBV	1078759	1011683	46.93%	5.995%
5	6.341	413	416	424	rBV	538811	683671	31.72%	4.051%
6	6.456	424	426	428	rBV	1807640	1650144	76.55%	9.778%
7	6.684	444	446	448	rBV	556380	473424	21.96%	2.805%
8	6.844	457	460	462	rVB	1554121	1690897	78.44%	10.020%
9	7.256	493	496	498	rBV	1257617	1267499	58.80%	7.511%
10	7.976	556	559	561	rBV	426495	569356	26.41%	3.374%
11	8.547	606	609	615	rBV4	10618	26910	1.25%	0.159%
12	8.719	621	624	627	rVB	63890	80514	3.74%	0.477%
13	8.833	631	634	636	rBV	20715	31342	1.45%	0.186%
14	8.970	643	646	650	rVB3	30996	53686	2.49%	0.318%
15	9.050	650	653	655	rBV	2396600	2155591	100.00%	12.774%
16	9.096	655	657	659	rVV	66883	61578	2.86%	0.365%
17	9.450	683	688	692	rVB	38839	65222	3.03%	0.386%
18	9.724	709	712	714	rBV	473779	557468	25.86%	3.303%
19	9.930	727	730	731	rBV	19179	25292	1.17%	0.150%
20	9.987	731	735	736	rVV4	13936	36848	1.71%	0.218%
21	10.022	736	738	740	rVV	46778	76876	3.57%	0.456%
22	10.079	742	743	747	rVB4	16364	26164	1.21%	0.155%
23	10.376	767	769	771	rBV	47273	66193	3.07%	0.392%
24	10.513	778	781	783	rBV	862478	1059284	49.14%	6.277%
25	10.650	790	793	795	rVB	104604	114992	5.33%	0.681%
26	10.696	795	797	799	rVV	149784	178894	8.30%	1.060%
27	10.730	799	800	802	rVV2	195319	237814	11.03%	1.409%
28	10.765	802	803	807	rVV	175790	253911	11.78%	1.505%
29	10.833	807	809	811	rVV2	72951	141223	6.55%	0.837%
30	10.879	811	813	815	rVV2	124384	161681	7.50%	0.958%
31	10.913	815	816	819	rVV	157298	211665	9.82%	1.254%
32	10.959	819	820	822	rVB	72841	52344	2.43%	0.310%
33	11.073	828	830	834	rVB3	18121	26972	1.25%	0.160%
34	11.199	839	841	844	rVB	578517	496885	23.05%	2.944%

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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	11.748	887	889	894	rBV2	21545	36767	1.71%	0.218%
36	11.976	907	909	911	rVB	18260	24725	1.15%	0.147%
37	12.582	959	962	964	rBV	16361	26673	1.24%	0.158%
38	12.788	977	980	982	rBV	1475191	1703317	79.02%	10.094%
39	13.359	1028	1030	1032	rBV	64870	68528	3.18%	0.406%
40	13.828	1069	1071	1073	rBV	501592	459803	21.33%	2.725%
41	13.873	1073	1075	1077	rVB2	15780	24223	1.12%	0.144%
42	15.211	1189	1192	1198	rVB	310055	362169	16.80%	2.146%

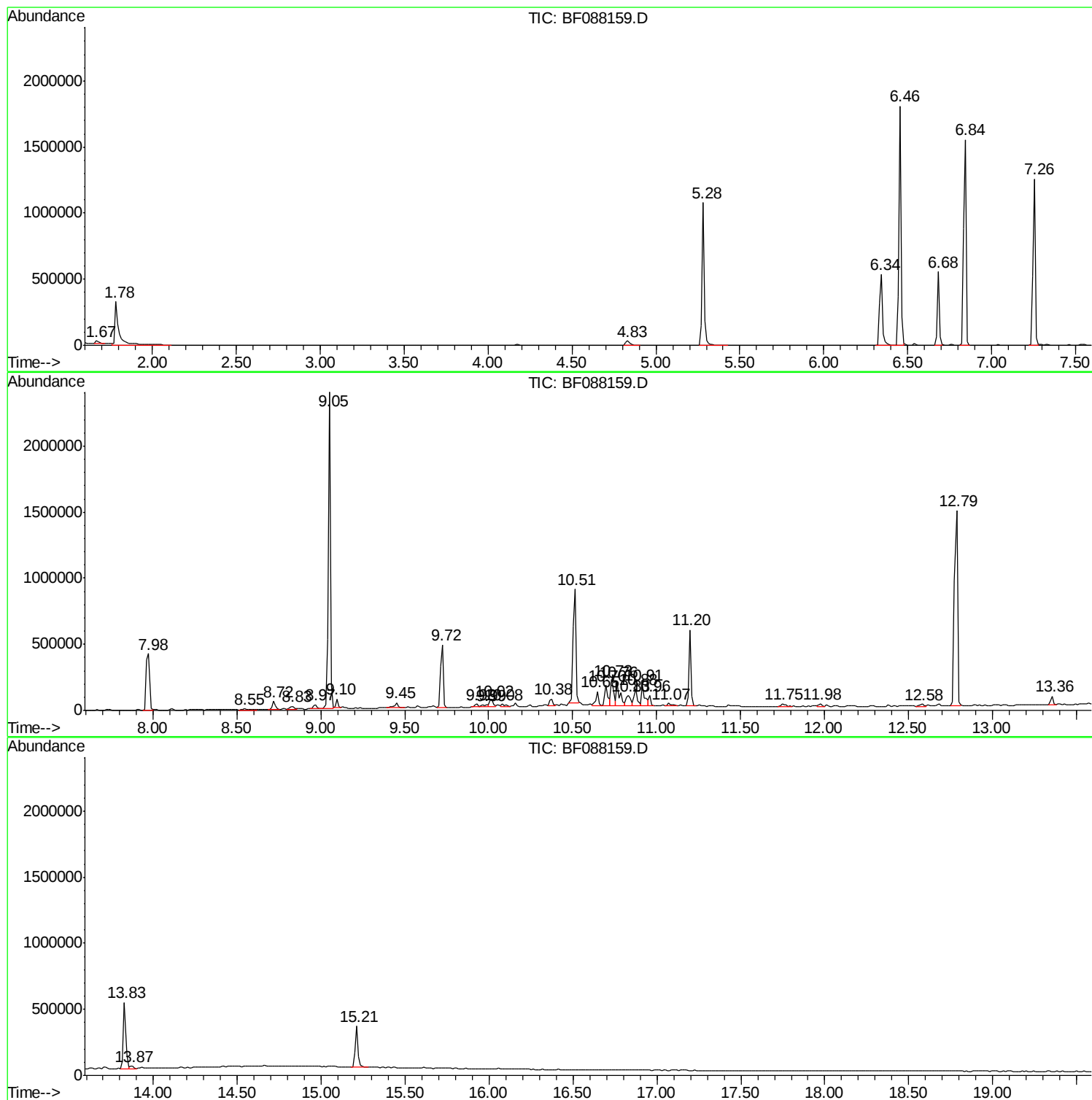
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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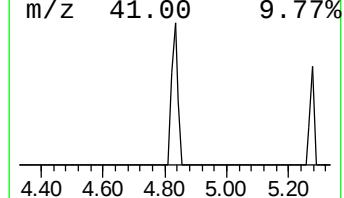
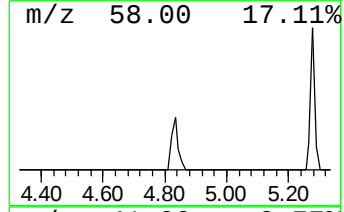
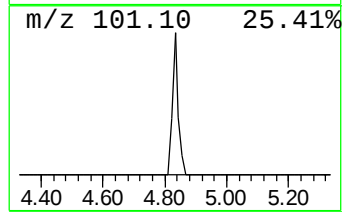
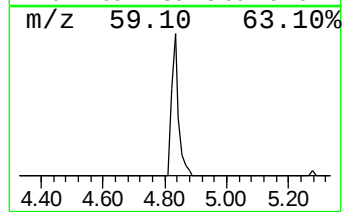
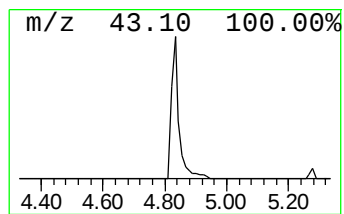
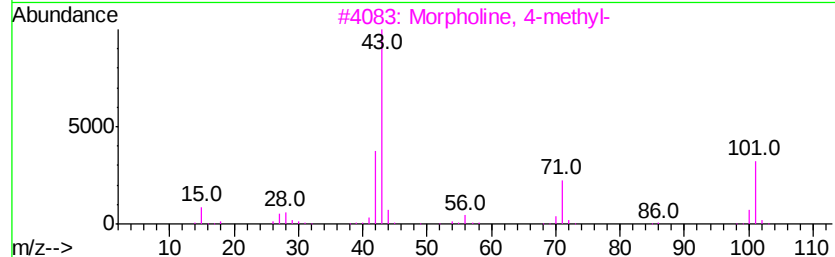
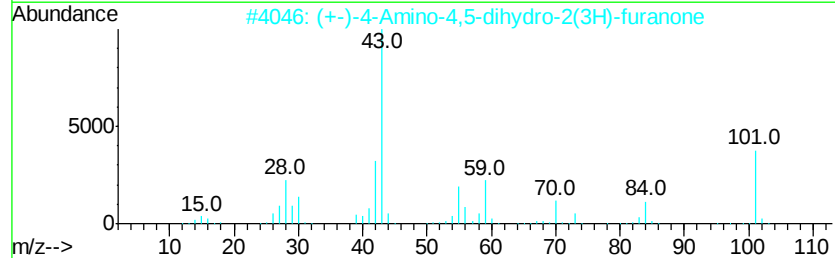
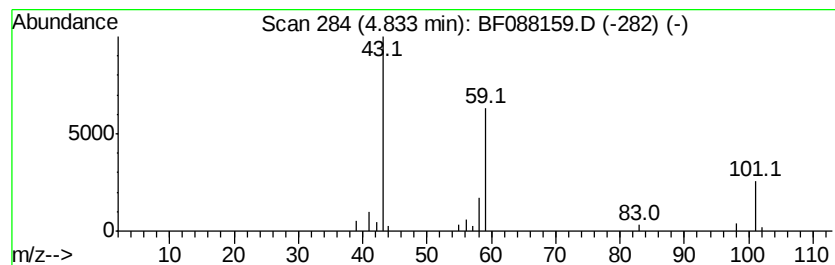
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	2.19 ng	51893	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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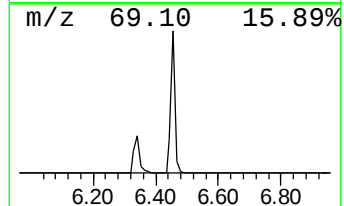
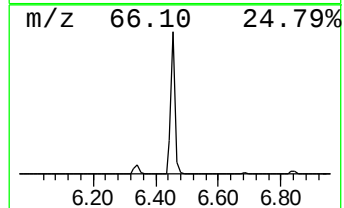
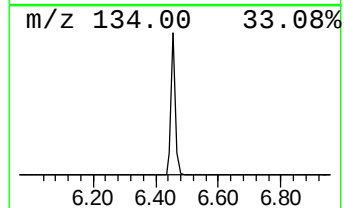
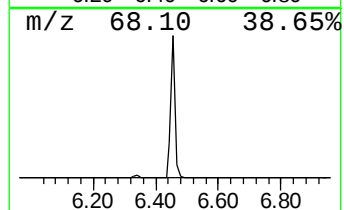
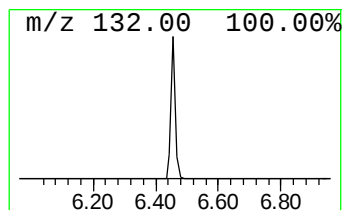
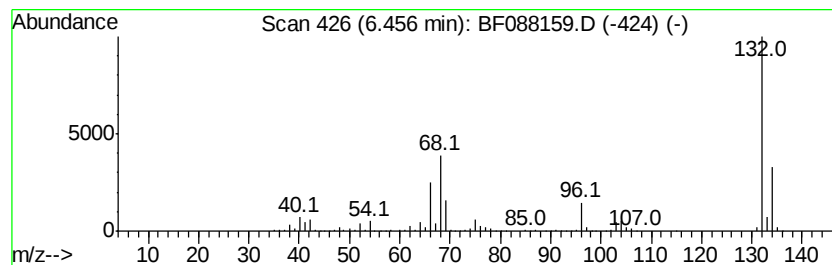
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.46 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	69.71 ng	1650140	1,4-Dichlorobenzene-d4	6.68

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	12
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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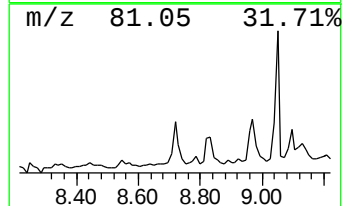
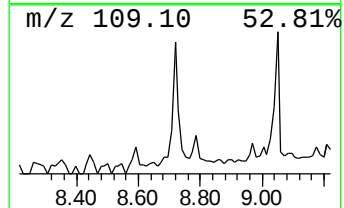
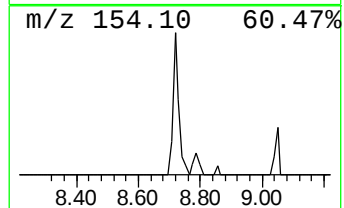
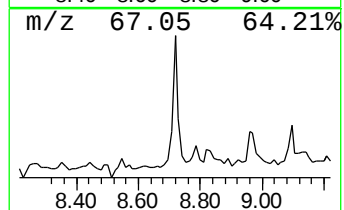
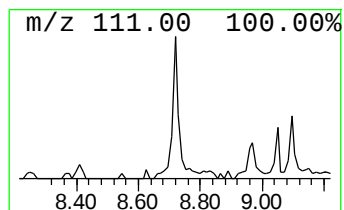
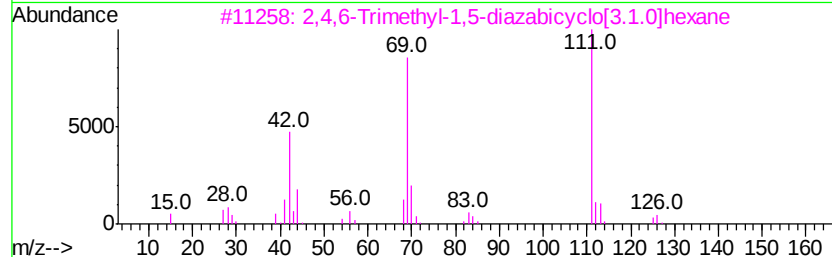
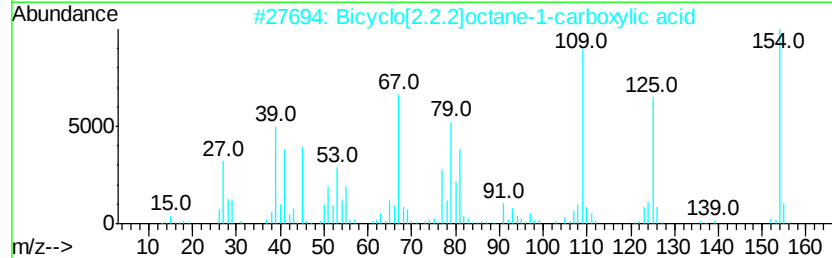
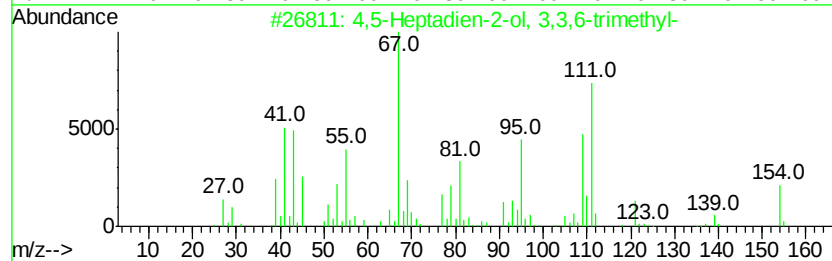
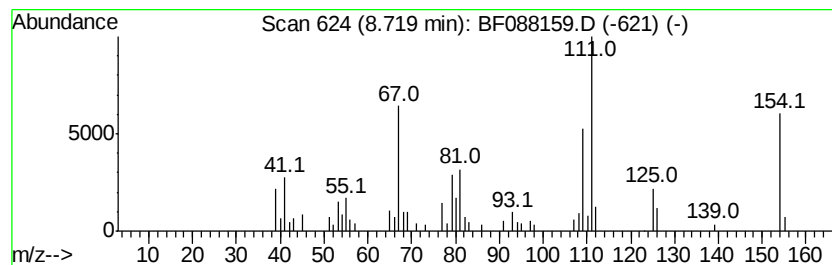
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Peak Number 5 4,5-Heptadien-2-ol, 3,3,6-t... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	2.83 ng	80514	Naphthalene-d8	7.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4,5-Heptadien-2-ol, 3,3,6-trimet...	154	C10H18O	060432-69-1	59
2		Bicyclo[2.2.2]octane-1-carboxyli...	154	C9H14O2	000699-55-8	45
3		2,4,6-Trimethyl-1,5-diazabicyclo...	126	C7H14N2	1000283-21-2	25
4		Phenol, 4-amino-2-nitro-	154	C6H6N2O3	000119-34-6	20
5		2-(Cyanoimino)oxazolidine	111	C4H5N3O	050411-06-8	14



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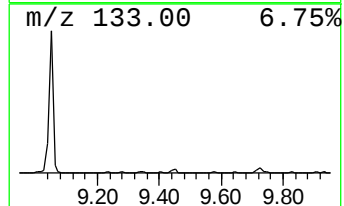
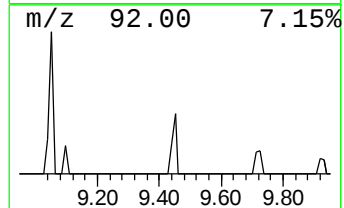
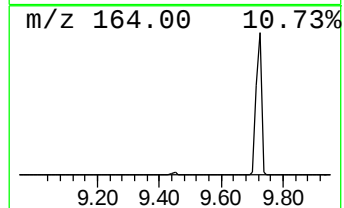
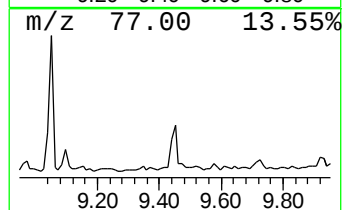
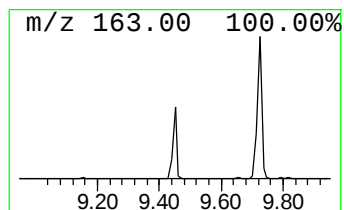
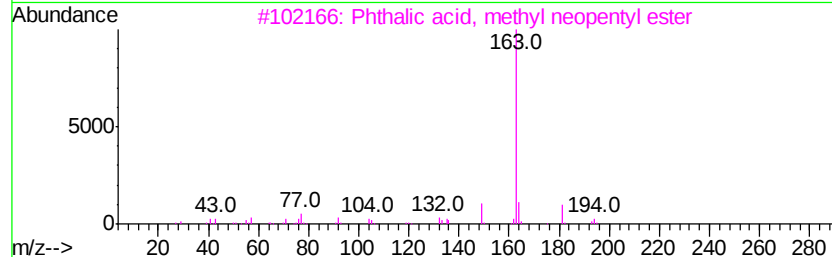
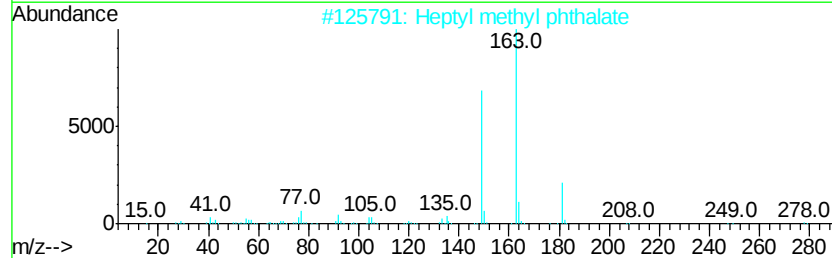
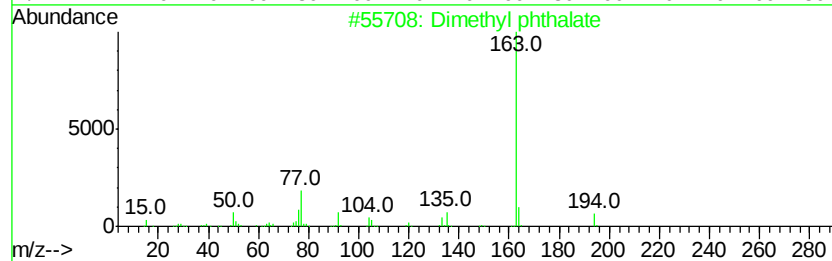
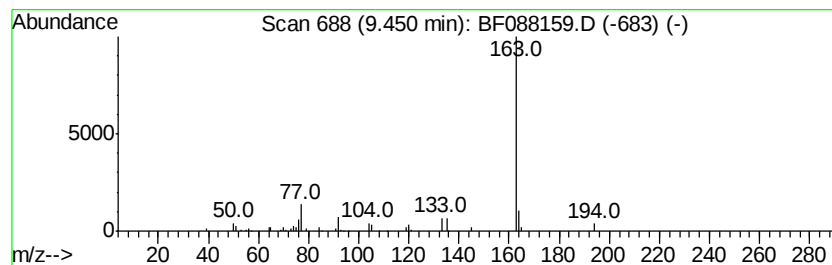
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 Dimethyl phthalate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.45	2.34 ng	65222	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dimethyl phthalate	194	C10H10O4	000131-11-3	91
2		Heptyl methyl phthalate	278	C16H22O4	1000373-88-4	83
3		Phthalic acid, methyl neopentyl ...	250	C14H18O4	1000315-53-9	83
4		Methyl 4-methylpentan-2-yl phtha...	264	C15H20O4	1000373-82-3	83
5		Methyl 4-methylpentyl phthalate	264	C15H20O4	1000373-82-6	83



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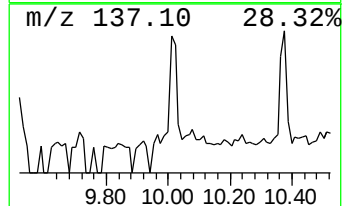
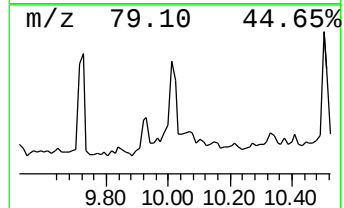
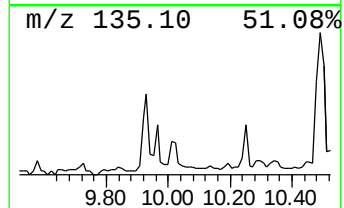
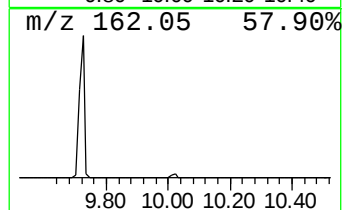
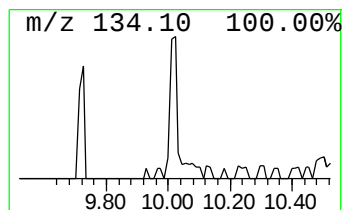
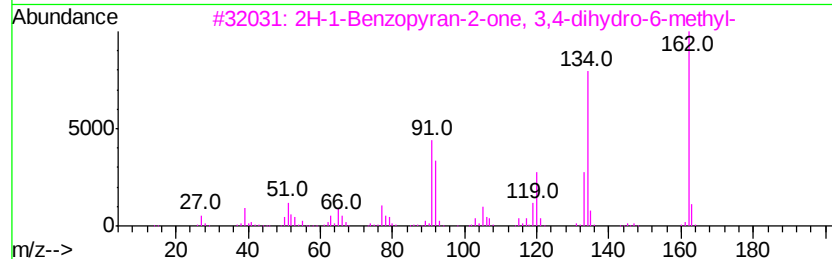
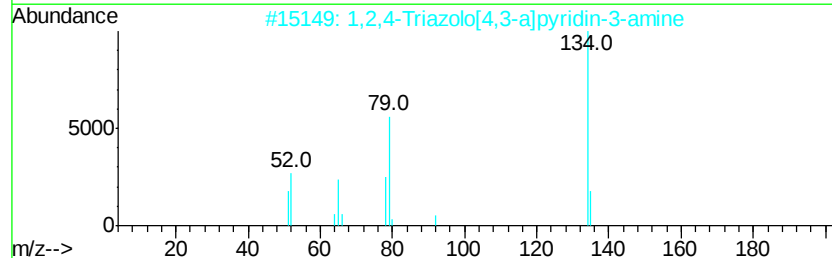
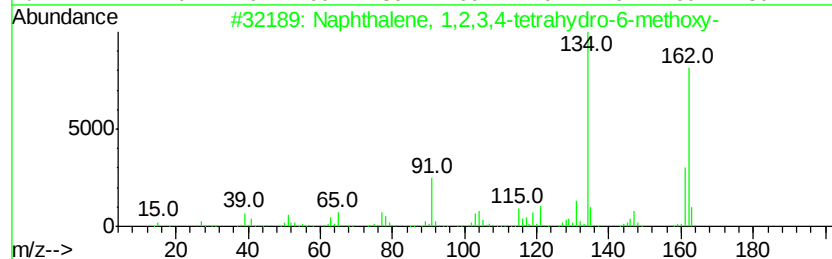
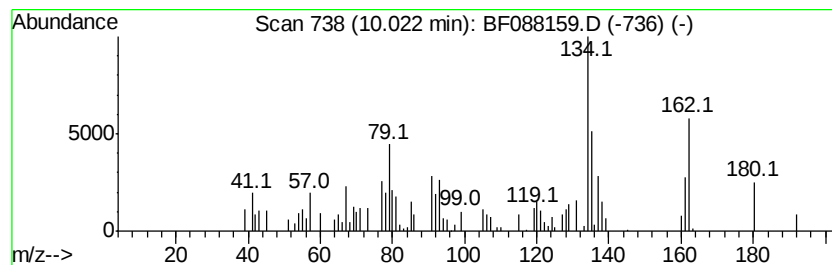
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Peak Number 7 unknown10.02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.02	2.76 ng	76876	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	162	C11H14O	001730-48-9	38
2		1,2,4-Triazolo[4,3-a]pyridin-3-a...	134	C6H6N4	000767-62-4	38
3		2H-1-Benzopyran-2-one, 3,4-dihyd...	162	C10H10O2	000092-47-7	35
4		Tricyclo[4.3.1.0(3,8)]decan-2-ol	152	C10H16O	1000197-64-5	32
5		1-Adamantyl fluoroformate	198	C11H15FO2	062087-82-5	27



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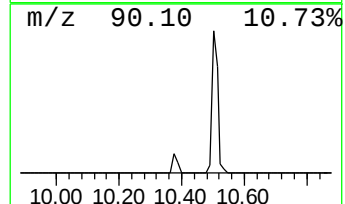
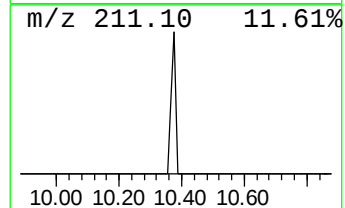
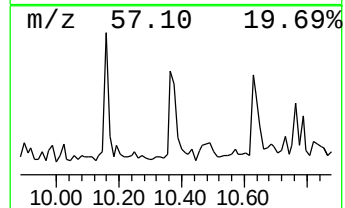
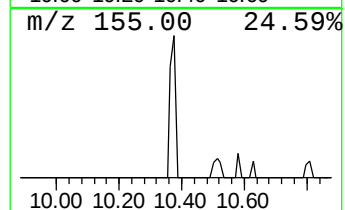
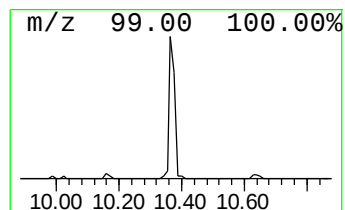
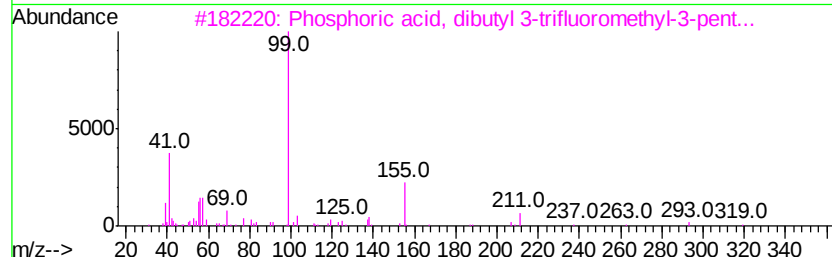
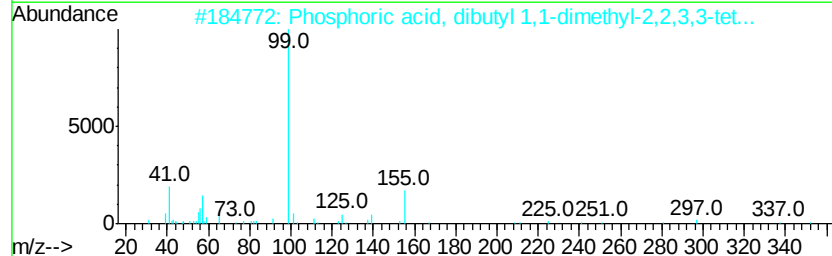
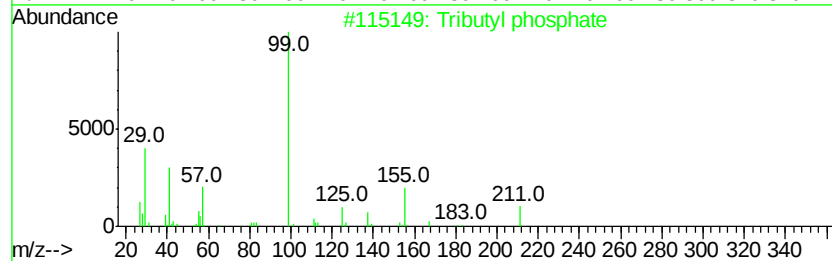
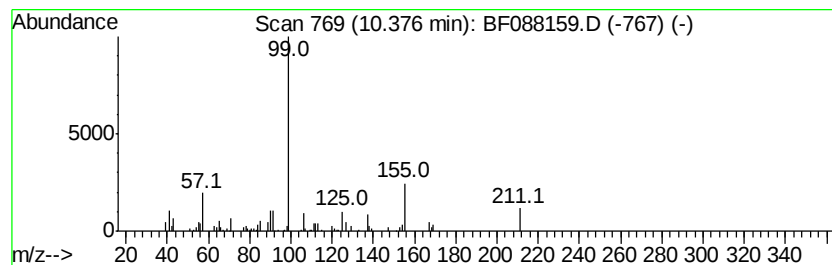
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tributyl phosphate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.38	2.37 ng	66193	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tributyl phosphate	266	C12H27O4P	000126-73-8	80
2		Phosphoric acid, dibutyl 1,1-dim...	352	C13H25F4O4P	1000298-93-9	42
3		Phosphoric acid, dibutyl 3-trifl...	348	C14H28F3O4P	1000298-93-8	36
4		Butyl diethyl phosphate	210	C8H19O4P	1000298-58-0	33
5		Diallyl maleate	196	C10H12O4	000999-21-3	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061816\
Data File : BF088159.D
Acq On : 19 Jun 2016 3:47
Operator : UM/SJ
Sample : H3628-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

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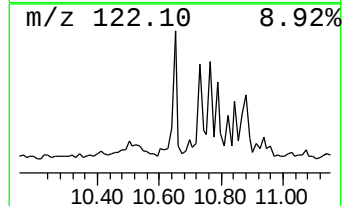
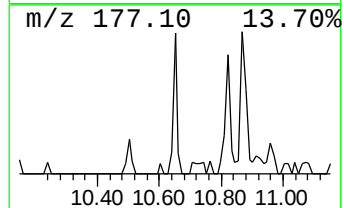
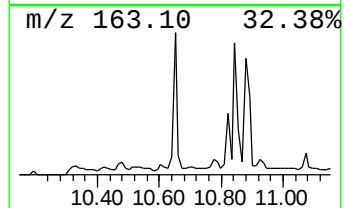
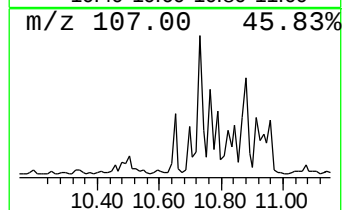
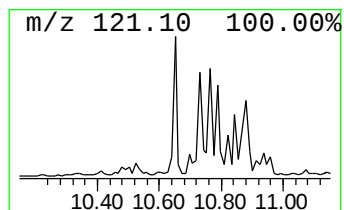
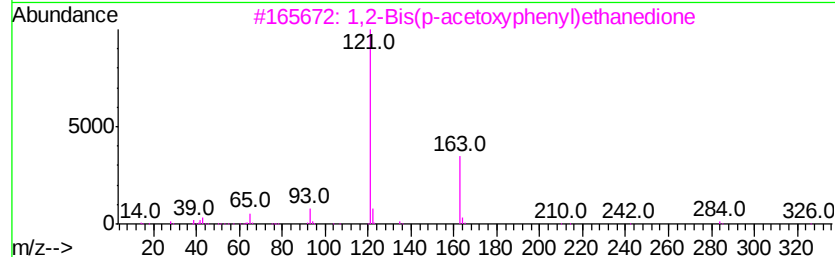
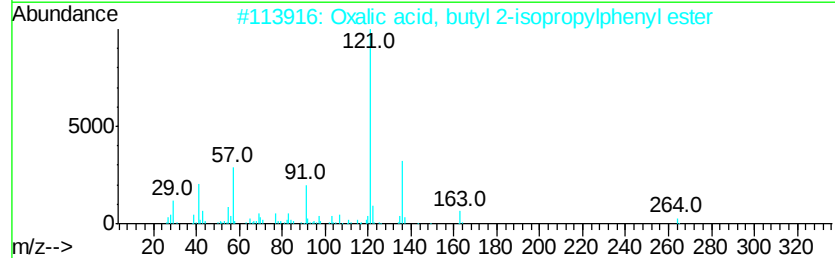
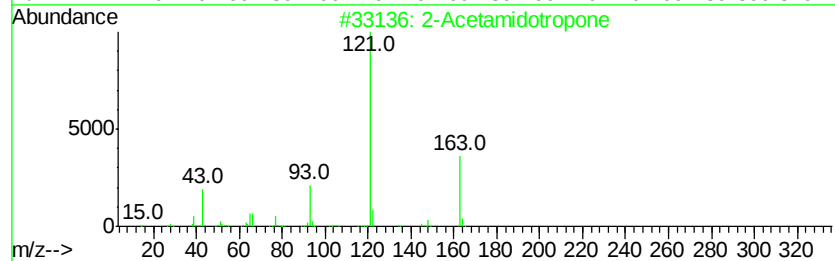
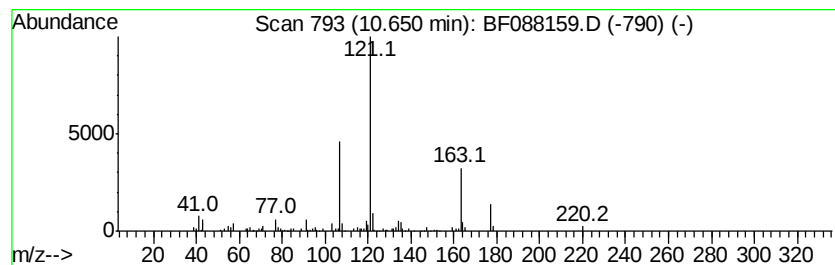
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown10.65 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	4.63 ng	114992	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2		2-Acetamidotropone	163	C9H9NO2	006422-12-4	47
2			Oxalic acid, butyl 2-isopropylph...	264	C15H20O4	1000309-56-3	47
3			1,2-Bis(p-acetoxyphehyl)ethanedione	326	C18H14O6	093655-79-9	47
4			Oxalic acid, 2-isopropylphenyl p...	250	C14H18O4	1000309-56-1	47
5			2-Propanone, 1-(4-methoxyphenyl)-	164	C10H12O2	000122-84-9	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061816\
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Misc :
ALS Vial : 38 Sample Multiplier: 1

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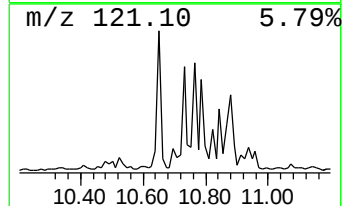
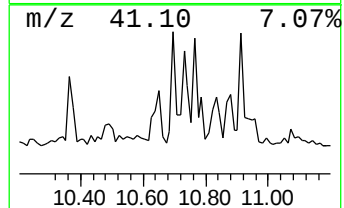
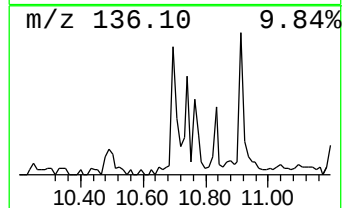
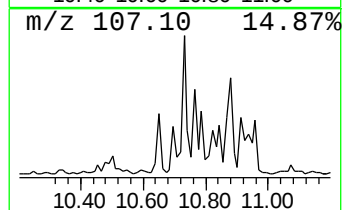
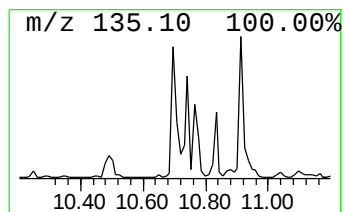
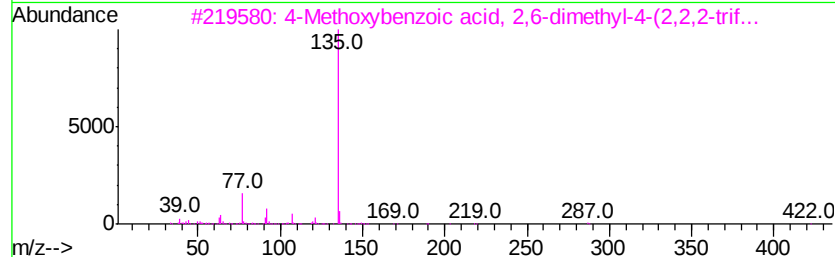
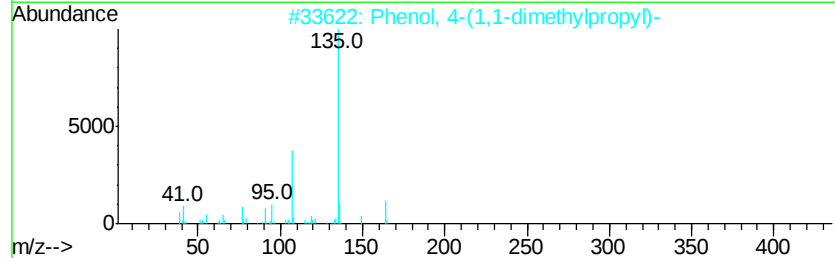
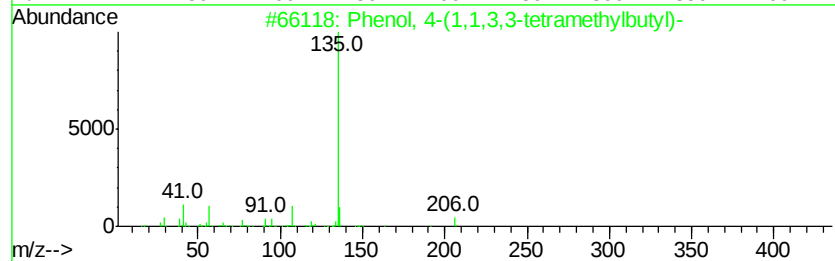
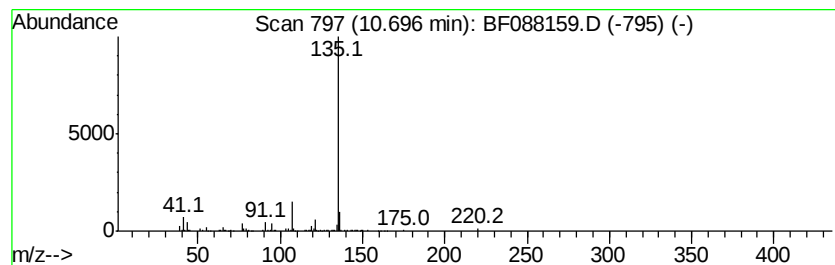
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	7.20 ng	178894	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2		Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3		4-Methoxybenzoic acid, 2,6-dimet...	422	C19H16F6O4	1000304-50-0	64
4		4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	64
5		1-Adamantanecarboxylic acid, 3,4...	324	C17H18Cl2O2	1000307-66-4	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061816\
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Acq On : 19 Jun 2016 3:47
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Misc :
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Instrument :
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ClientSampleId :
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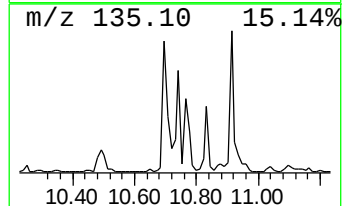
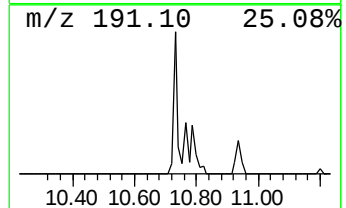
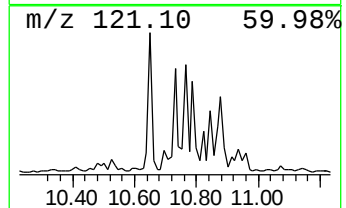
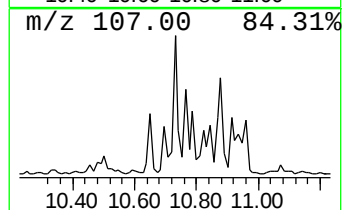
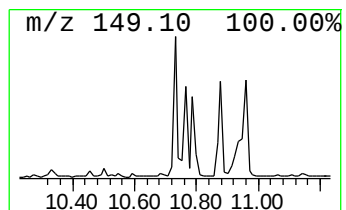
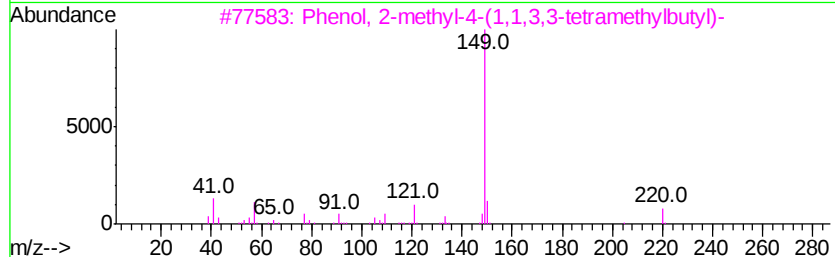
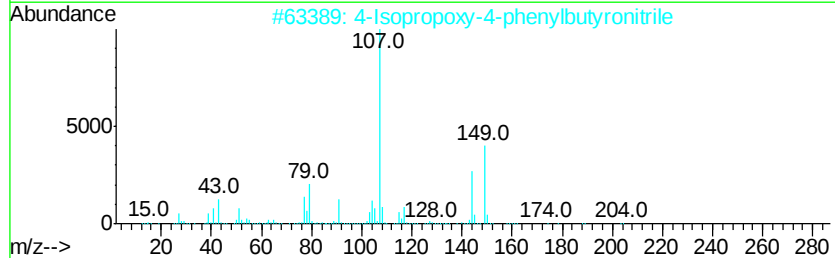
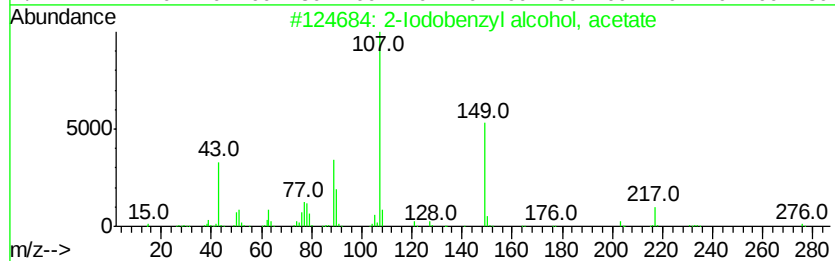
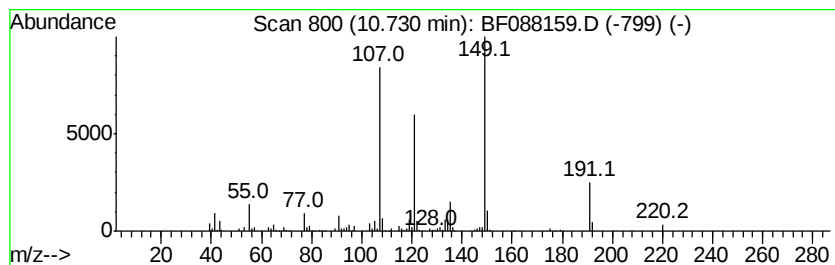
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 2-Iodobenzyl alcohol, acetate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.73	9.57 ng	237814	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Iodobenzyl alcohol, acetate	276	C9H9IO2	1000364-47-1	59
2		4-Isopropoxy-4-phenylbutyronitrile	203	C13H17NO	1000370-35-3	45
3		Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	38
4		Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	35
5		Phenol, 2-(1,1-dimethylethyl)-3-...	164	C11H16O	013037-79-1	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061816\
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Acq On : 19 Jun 2016 3:47
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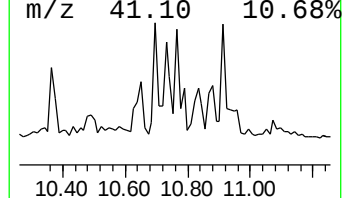
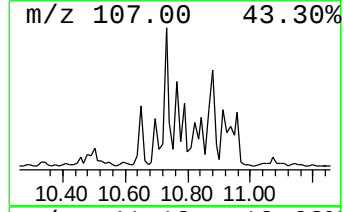
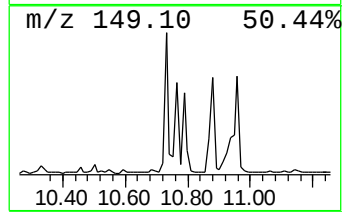
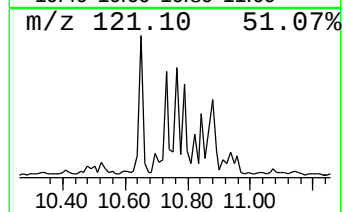
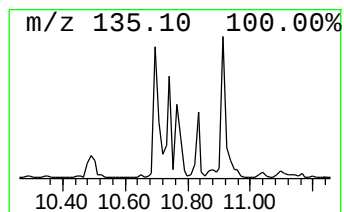
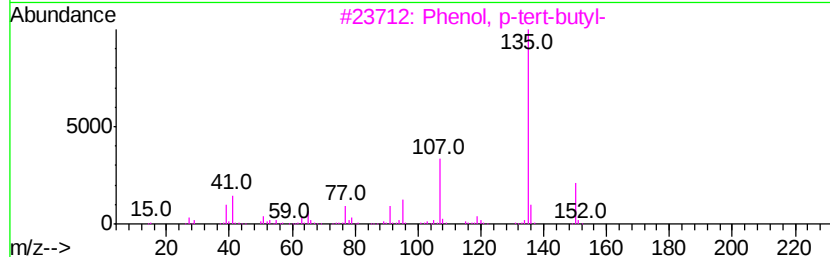
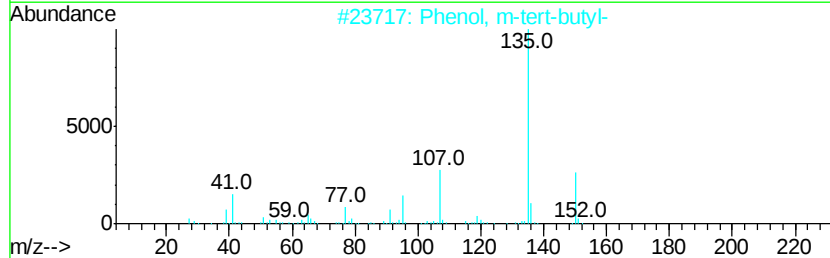
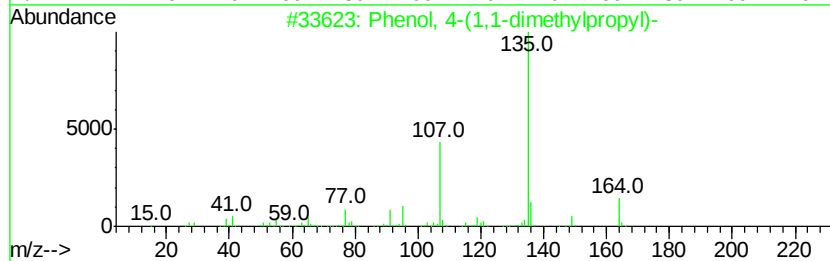
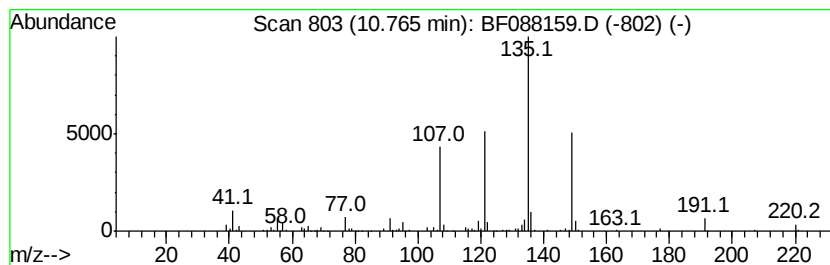
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Phenol, 4-(1,1-dimethylprop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	10.22 ng	253911	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50
2		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	50
3		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	50
4		((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	50
5		Hexestrol, 0-acetyl-	312	C20H24O3	1000374-87-8	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061816\
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Misc :
ALS Vial : 38 Sample Multiplier: 1

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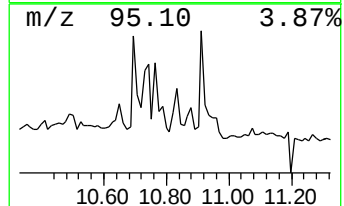
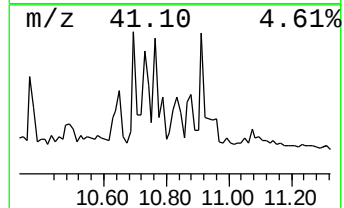
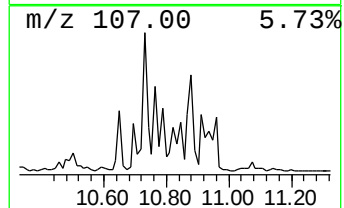
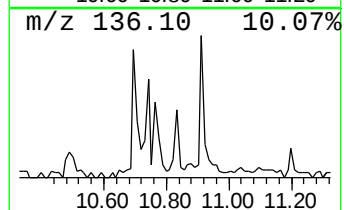
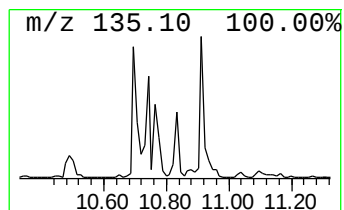
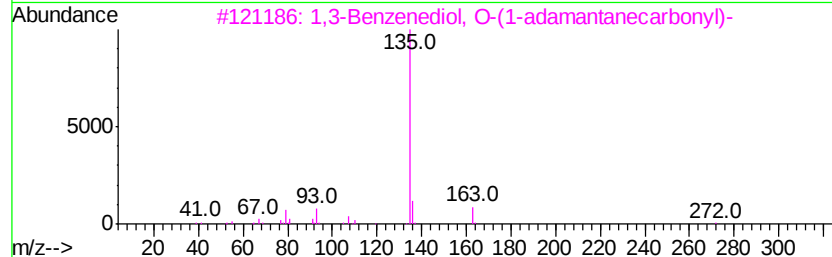
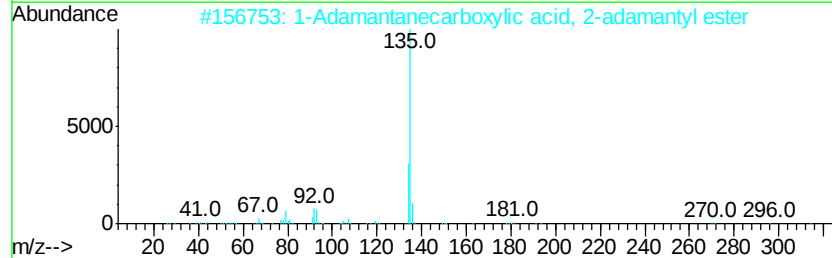
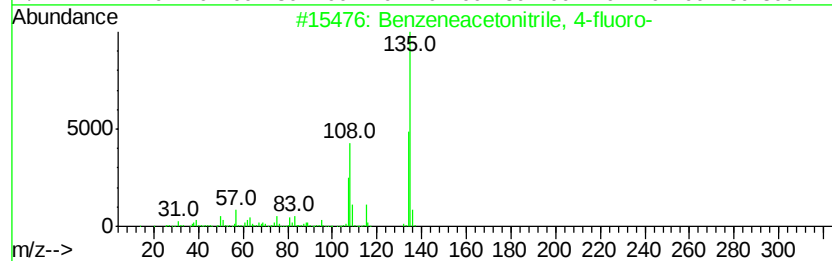
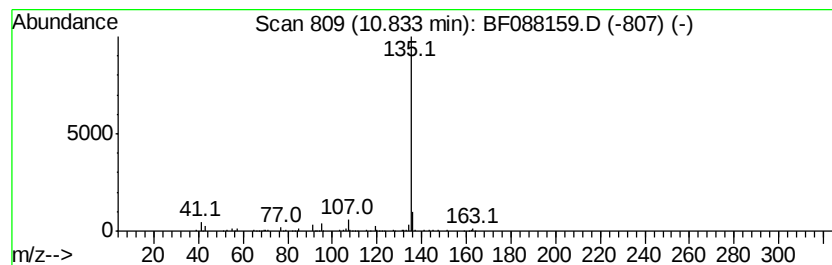
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzeneacetonitrile, 4-fluoro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.83	5.68 ng	141223	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetonitrile, 4-fluoro-	135	C8H6FN	000459-22-3	56
2			1-Adamantanecarboxylic acid, 2-a...	314	C21H30O2	1000282-94-3	56
3			1,3-Benzenediol, O-(1-adamantane...	272	C17H20O3	1000345-57-3	50
4			1-Adamantanecarboxylic acid, 4-c...	281	C18H19NO2	1000307-66-3	50
5			p-Anisic acid, 3,4-dichloropheny...	296	C14H10Cl2O3	1000307-63-8	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061816\
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ALS Vial : 38 Sample Multiplier: 1

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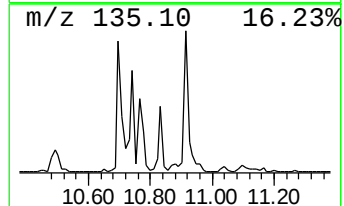
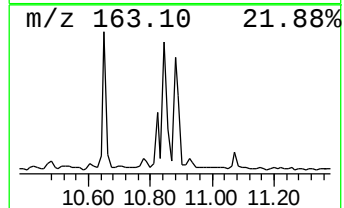
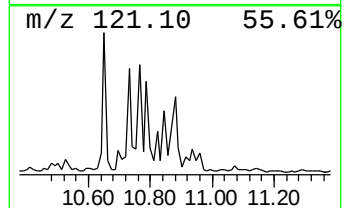
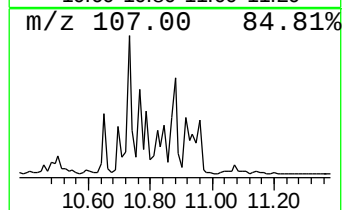
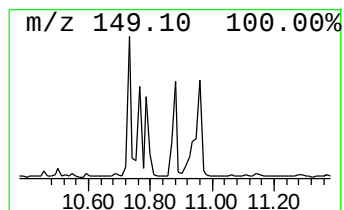
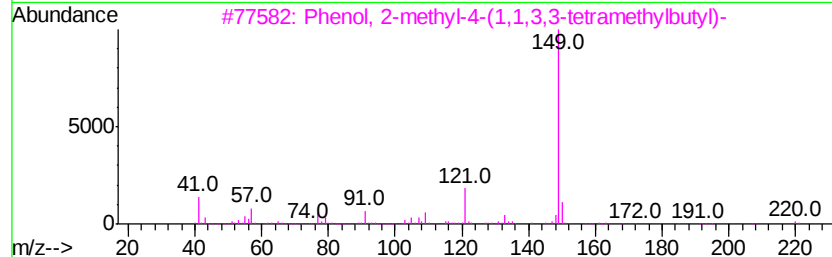
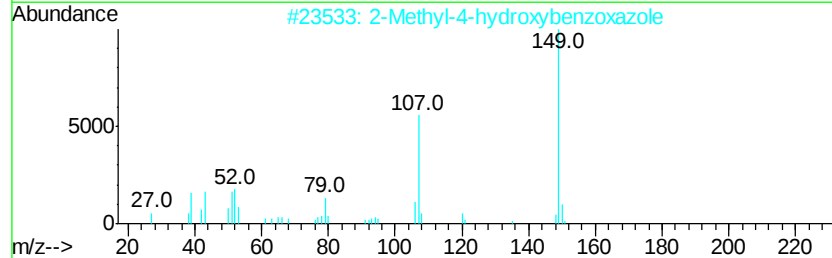
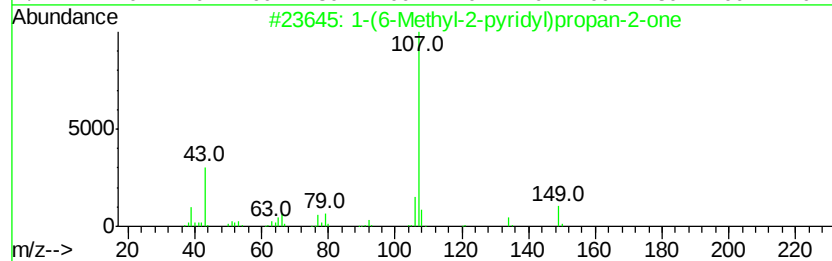
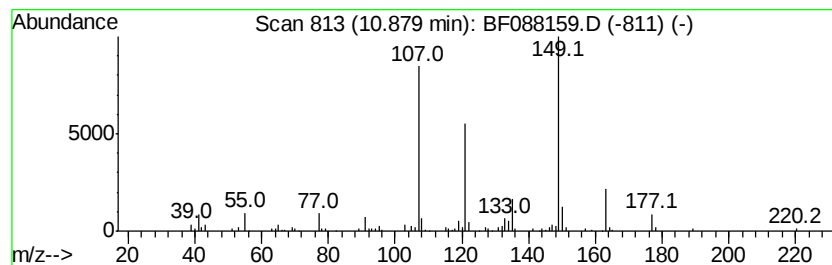
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-(6-Methyl-2-pyridyl)propa... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	6.51 ng	161681	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(6-Methyl-2-pyridyl)propan-2-one	149	C9H11NO	065702-08-1	59
2			2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	59
3			Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	43
4			Benzene, 1-ethoxy-4-ethyl-	150	C10H14O	001585-06-4	38
5			Phenol, 2-(1,1-dimethylethyl)-3-...	164	C11H16O	013037-79-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061816\
Data File : BF088159.D
Acq On : 19 Jun 2016 3:47
Operator : UM/SJ
Sample : H3628-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-15

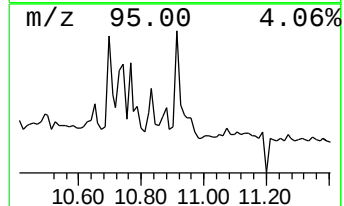
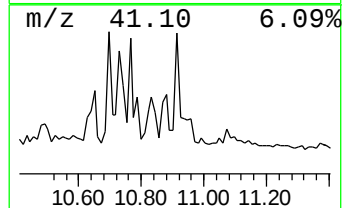
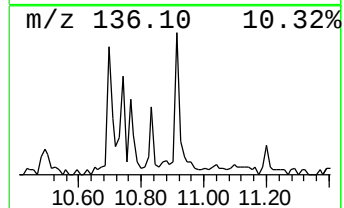
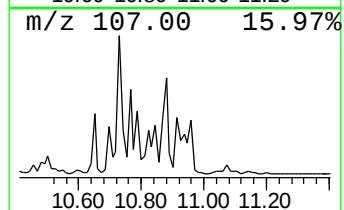
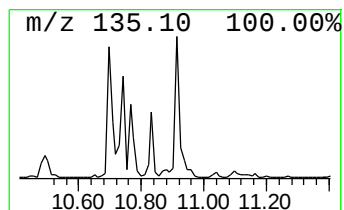
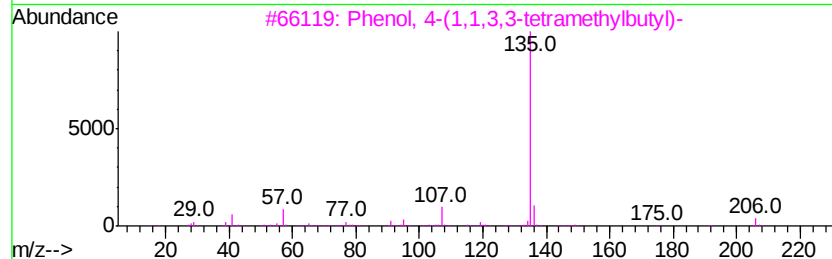
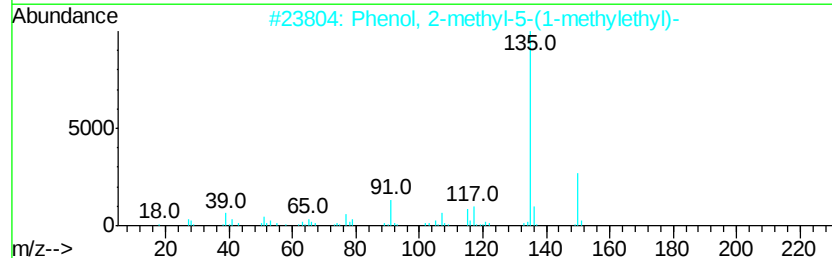
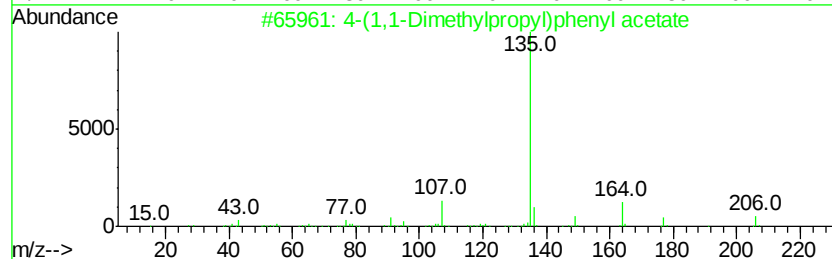
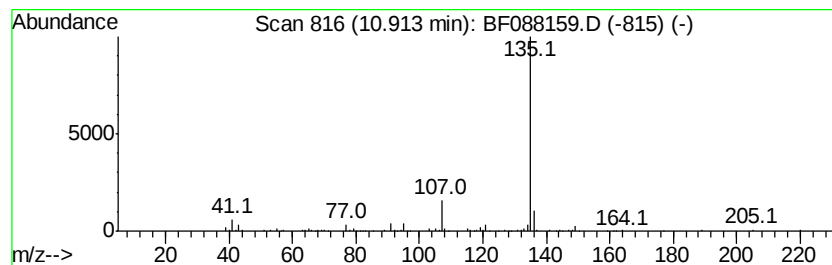
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 4-(1,1-Dimethylpropyl)pheny... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	8.52 ng	211665	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	83
2		Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	78
3		Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
4		Butane, 1,3-dibromo-	214	C4H8Br2	000107-80-2	78
5		4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78



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Operator : UM/SJ
Sample : H3628-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-15

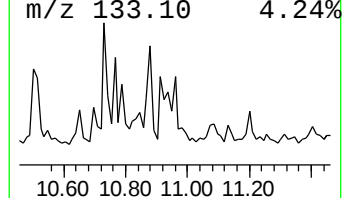
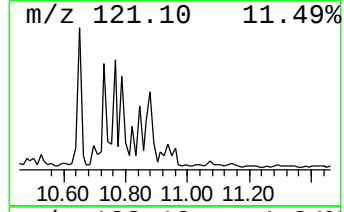
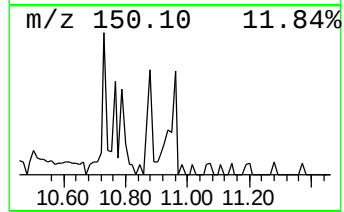
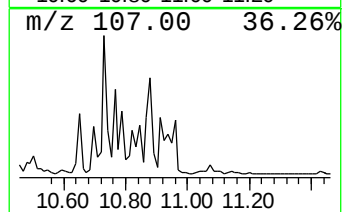
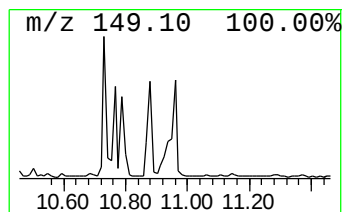
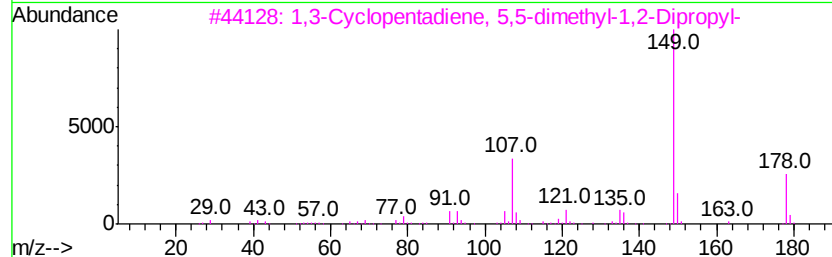
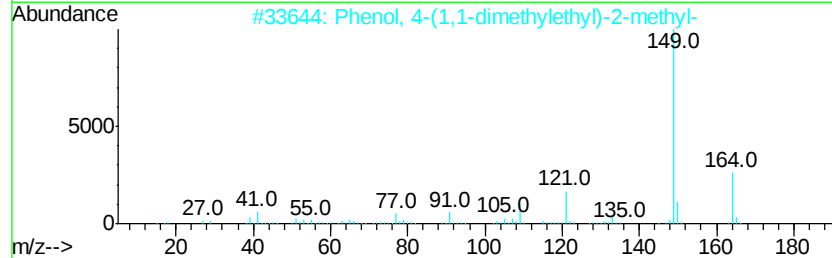
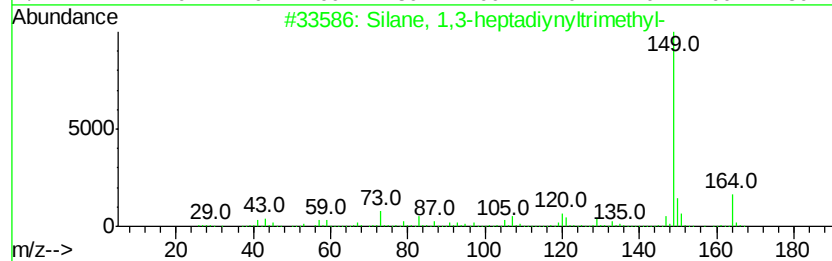
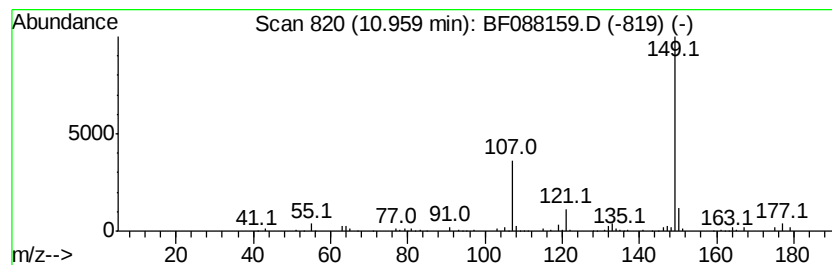
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Silane, 1,3-heptadiynyltrim... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.96	2.11 ng	52344	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, 1,3-heptadiynyltrimethyl-	164	C10H16Si	019024-47-6	59
2			Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	59
3			1,3-Cyclopentadiene, 5,5-dimethv...	178	C13H22	1000163-88-0	56
4			2-t-Butyl-5-[hydroxy-(2,4,6-trim...	306	C18H26O4	092572-63-9	45
5			Hexestrol dimethyl ether	298	C20H26O2	000130-78-9	42



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ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
W-15

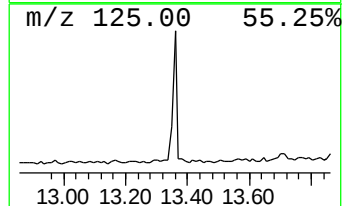
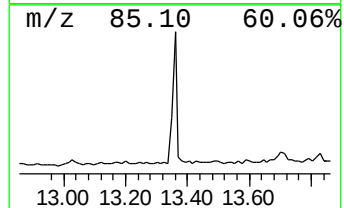
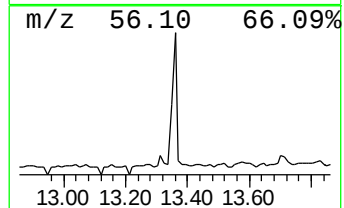
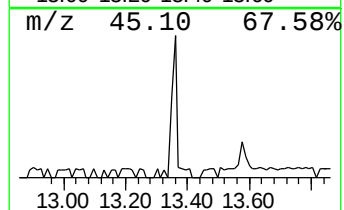
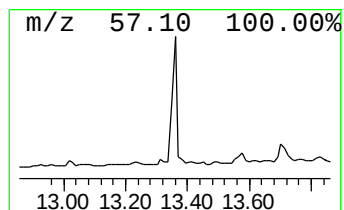
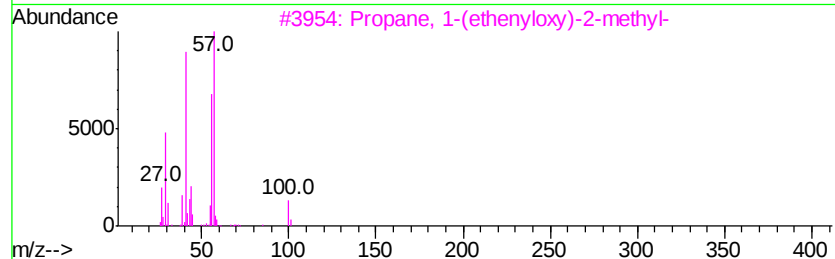
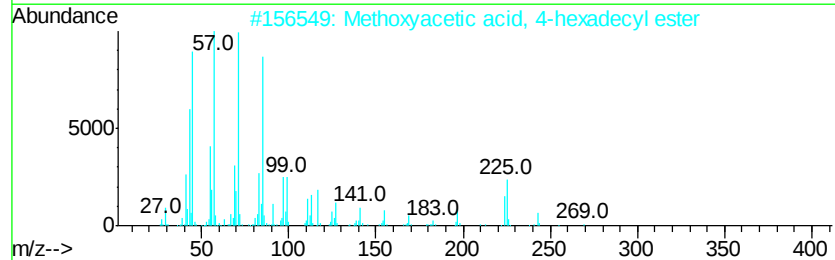
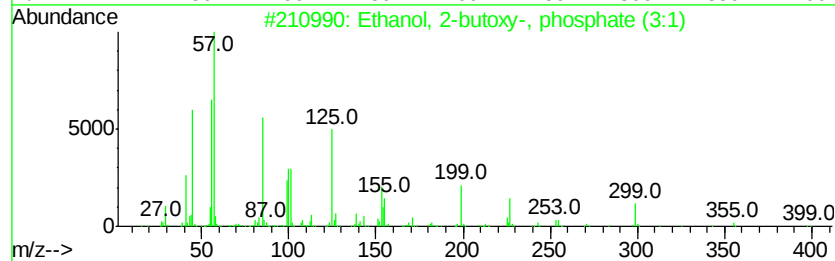
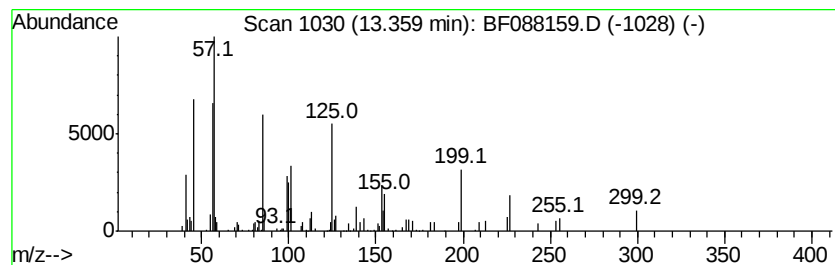
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Ethanol, 2-butoxy-, phospho... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	2.98 ng	68528	Chrysene-d12	13.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-butoxy-, phosphate (3:1)	398	C18H39O7P	000078-51-3	91
2		Methoxyacetic acid, 4-hexadecyl ...	314	C19H38O3	1000282-99-0	10
3		Propane, 1-(ethenyloxy)-2-methyl-	100	C6H12O	000109-53-5	10
4		Butanoic acid, 2-methyl-, 2-meth...	158	C9H18O2	002445-67-2	10
5		Carbonic acid, isobutyl isohexyl...	202	C11H22O3	1000314-60-3	10



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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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...	4.83	2.2	ng	51893	1	6.68	473424	20.0
unknown6.46	6.46	69.7	ng	1650140	1	6.68	473424	20.0
4,5-Heptadien-2-o...	8.72	2.8	ng	80514	2	7.98	569356	20.0
Dimethyl phthalate	9.45	2.3	ng	65222	3	9.72	557468	20.0
unknown10.02	10.02	2.8	ng	76876	3	9.72	557468	20.0
Tributyl phosphate	10.38	2.4	ng	66193	3	9.72	557468	20.0
unknown10.65	10.65	4.6	ng	114992	4	11.20	496885	20.0
Phenol, 4-(1,1,3,...	10.70	7.2	ng	178894	4	11.20	496885	20.0
2-Iodobenzyl alco...	10.73	9.6	ng	237814	4	11.20	496885	20.0
Phenol, 4-(1,1-di...	10.76	10.2	ng	253911	4	11.20	496885	20.0
Benzeneacetoneitri...	10.83	5.7	ng	141223	4	11.20	496885	20.0
1-(6-Methyl-2-pyr...	10.88	6.5	ng	161681	4	11.20	496885	20.0
4-(1,1-Dimethylpr...	10.91	8.5	ng	211665	4	11.20	496885	20.0
Silane, 1,3-hepta...	10.96	2.1	ng	52344	4	11.20	496885	20.0
Ethanol, 2-butoxy...	13.36	3.0	ng	68528	5	13.83	459803	20.0