

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
 Data File : BF086374.D
 Acq On : 23 Apr 2016 6:48
 Operator : UM/SJ
 Sample : H2634-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-11

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-BF042216.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	3.767	143	147	150	rBV	29156	54217	1.23%	0.148%
2	5.001	254	255	260	rBV	145253	192746	4.38%	0.526%
3	5.287	278	280	283	rBV	155063	159935	3.63%	0.436%
4	5.424	290	292	305	rVB	1362872	1532430	34.82%	4.179%
5	5.664	311	313	315	rBV	104689	106571	2.42%	0.291%
6	5.836	326	328	331	rBV2	47070	75225	1.71%	0.205%
7	6.362	371	374	379	rBV2	78863	126281	2.87%	0.344%
8	6.464	379	383	386	rBV	869971	1104385	25.09%	3.012%
9	6.522	386	388	393	rVV	293346	358212	8.14%	0.977%
10	6.602	393	395	398	rVV	2940689	2905501	66.02%	7.924%
11	6.659	399	400	405	rVV	142581	266885	6.06%	0.728%
12	6.796	409	412	414	rVV3	24257	53562	1.22%	0.146%
13	6.842	414	416	419	rVV	836000	956359	21.73%	2.608%
14	6.899	419	421	423	rVV	271323	290781	6.61%	0.793%
15	6.990	427	429	434	rVV2	2069102	3233370	73.47%	8.818%
16	7.093	436	438	442	rVV	141797	210222	4.78%	0.573%
17	7.162	442	444	449	rVB2	105535	165530	3.76%	0.451%
18	7.242	449	451	452	rBV	66541	82227	1.87%	0.224%
19	7.402	463	465	471	rVB2	1999623	2395035	54.42%	6.532%
20	7.493	471	473	474	rVB2	53910	73828	1.68%	0.201%
21	7.619	481	484	485	rBV	160820	227719	5.17%	0.621%
22	7.642	485	486	489	rVB2	99981	106536	2.42%	0.291%
23	7.767	495	497	498	rBV2	58624	89650	2.04%	0.244%
24	7.802	498	500	503	rVB2	153027	215818	4.90%	0.589%
25	7.859	503	505	507	rBV2	542300	672497	15.28%	1.834%
26	7.939	510	512	516	rVB4	40212	95141	2.16%	0.259%
27	8.065	521	523	525	rBV2	66226	143668	3.26%	0.392%
28	8.122	526	528	531	rBV	1260090	1447923	32.90%	3.949%
29	8.282	541	542	545	rBV2	41046	49200	1.12%	0.134%
30	8.407	548	553	558	rBV3	80807	168873	3.84%	0.461%
31	8.510	560	562	564	rVB	103506	93997	2.14%	0.256%
32	8.567	564	567	568	rBV2	36754	67170	1.53%	0.183%
33	8.590	568	569	571	rVV	92612	82736	1.88%	0.226%
34	8.670	574	576	579	rBV3	28532	45351	1.03%	0.124%

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35	8.933	597	599	604	rVB3	126839	179888	4.09%	0.491%
36	9.196	620	622	625	rBV	3024832	4401193	100.00%	12.003%
37	9.459	642	645	650	rBV5	21408	62051	1.41%	0.169%
38	9.596	655	657	661	rBV	50740	78630	1.79%	0.214%
39	9.813	675	676	679	rBV	66242	60483	1.37%	0.165%
40	9.870	679	681	684	rBV	1131203	1449720	32.94%	3.954%
41	10.168	703	707	708	rBV2	29142	59049	1.34%	0.161%
42	10.316	716	720	721	rBV2	100179	123883	2.81%	0.338%
43	10.396	724	727	733	rVB2	26082	82535	1.88%	0.225%
44	10.533	736	739	740	rBV	47985	72067	1.64%	0.197%
45	10.659	748	750	759	rBV	2229009	3014146	68.48%	8.220%
46	10.785	759	761	766	rVB	135731	191661	4.35%	0.523%
47	11.231	798	800	801	rBV	53271	48965	1.11%	0.134%
48	11.356	809	811	819	rBV	1430473	1657872	37.67%	4.521%
49	11.894	856	858	863	rBV2	176238	256194	5.82%	0.699%
50	12.682	924	927	930	rBV	126160	220640	5.01%	0.602%
51	12.774	933	935	936	rVV	52725	50610	1.15%	0.138%
52	12.945	947	950	960	rBV	3628495	4350291	98.84%	11.864%
53	13.985	1039	1041	1049	rBV	890131	1300874	29.56%	3.548%
54	15.402	1163	1165	1175	rVB	609562	1157523	26.30%	3.157%

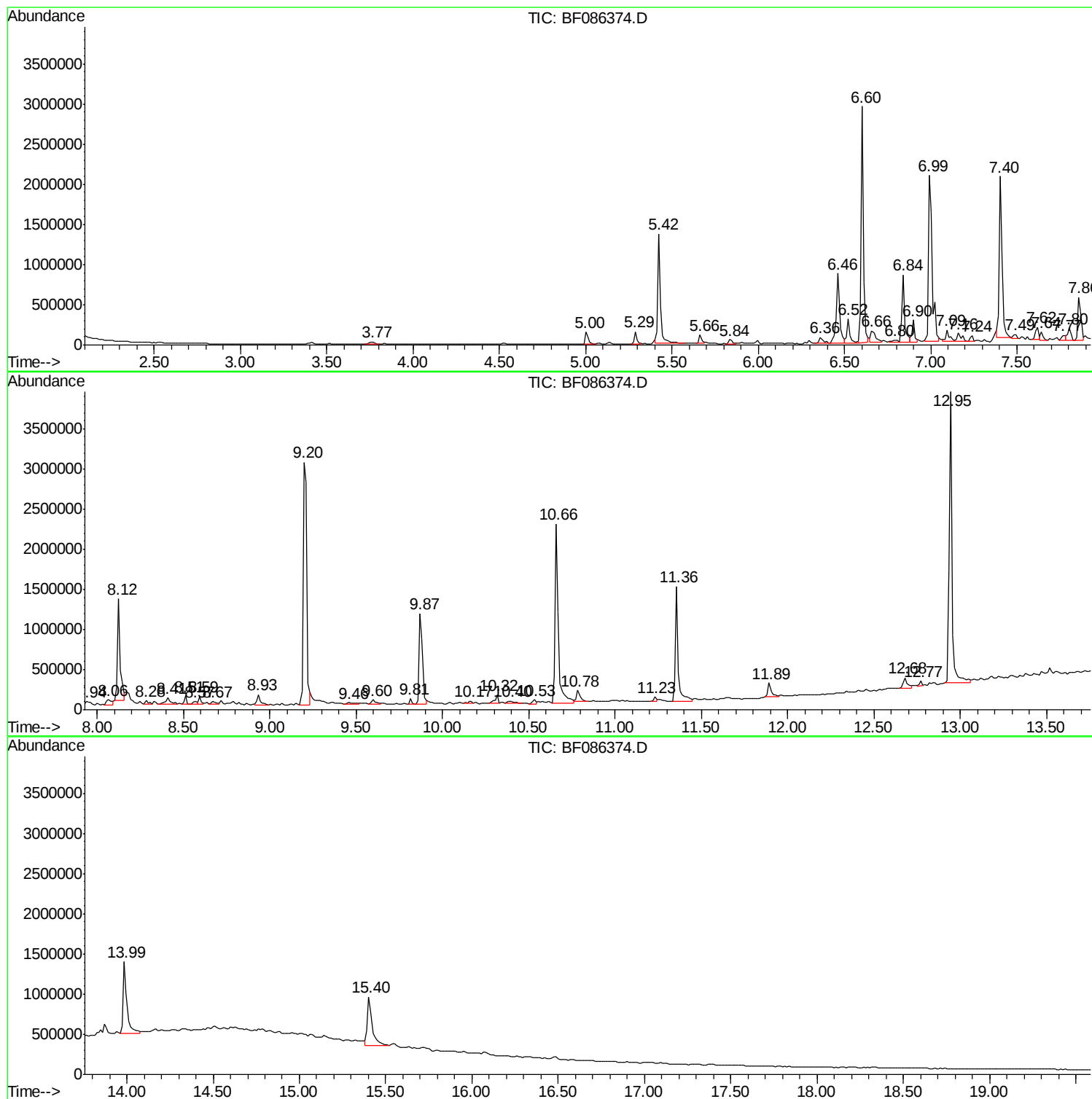
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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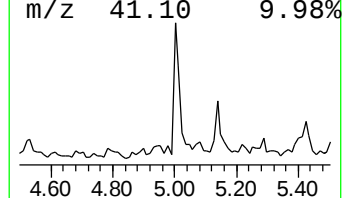
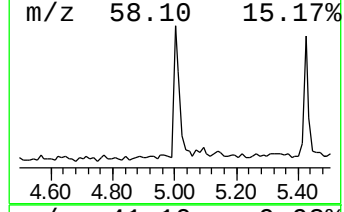
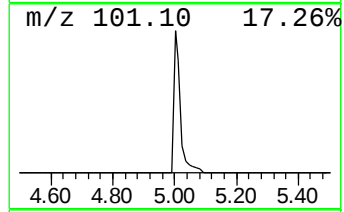
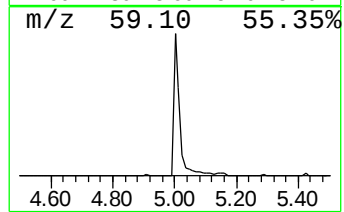
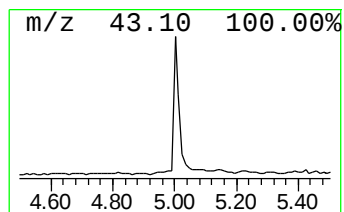
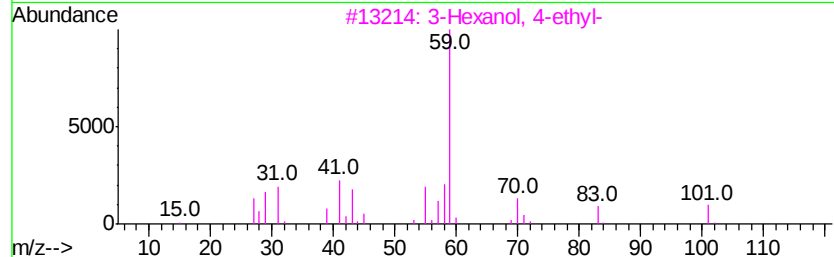
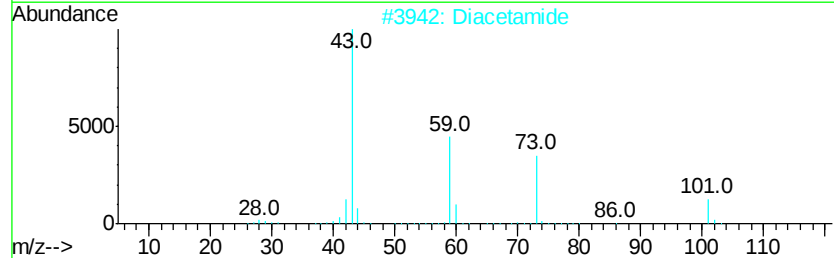
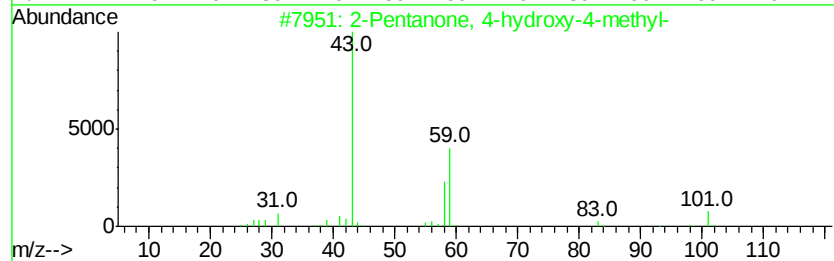
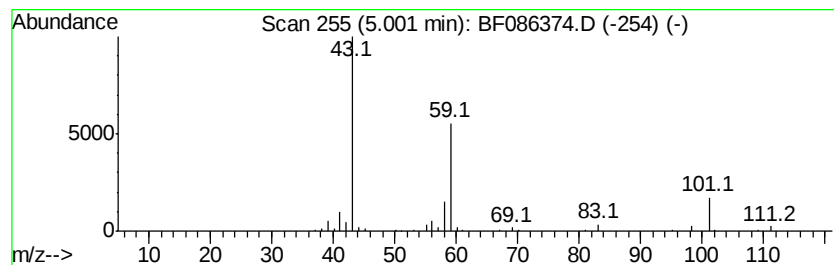
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	4.03 ng	192746	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Diacetamide	101	C4H7NO2	000625-77-4	36
3			3-Hexanol, 4-ethyl-	130	C8H18O	019780-44-0	28
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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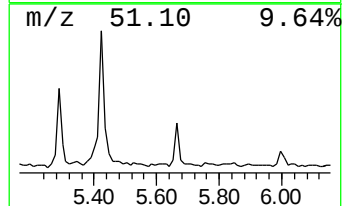
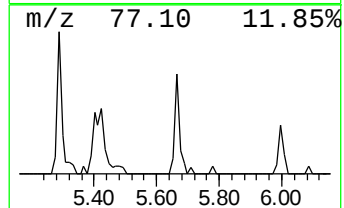
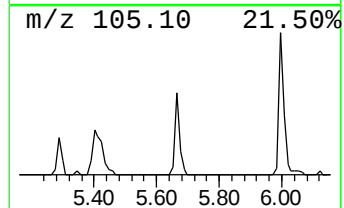
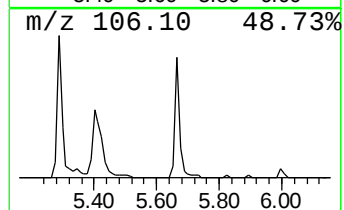
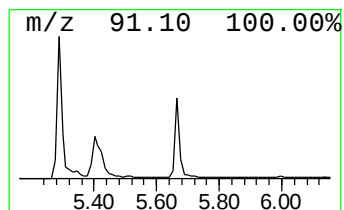
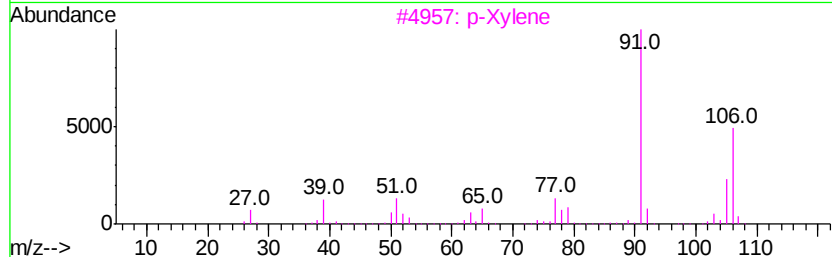
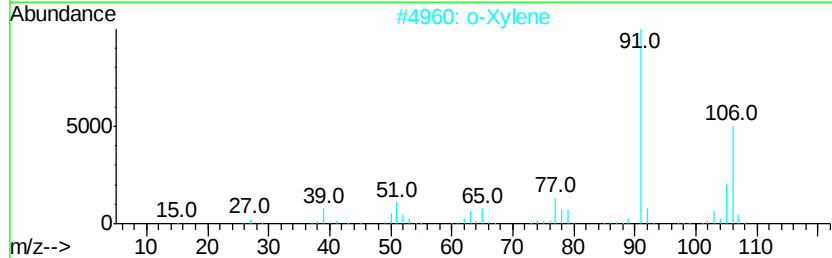
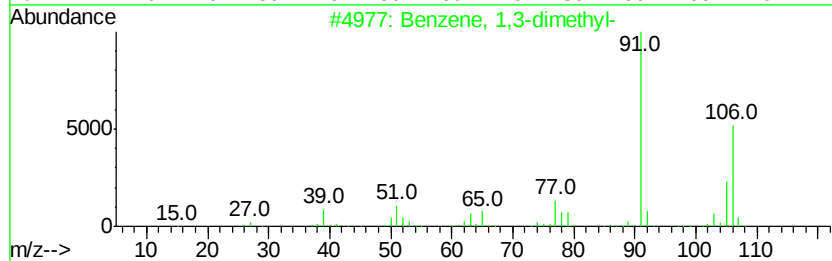
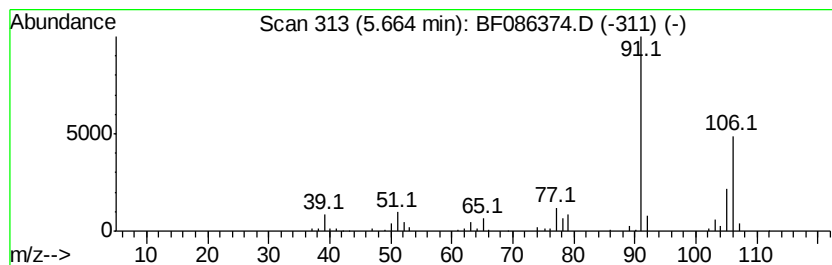
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.66	2.23 ng	106571	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2			o-Xylene	106	C8H10	000095-47-6	97
3			p-Xylene	106	C8H10	000106-42-3	97
4			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	90
5			Ethylbenzene	106	C8H10	000100-41-4	87



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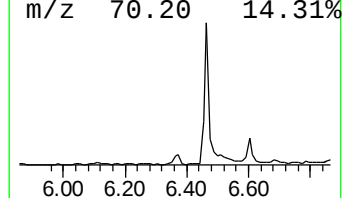
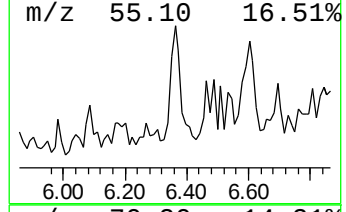
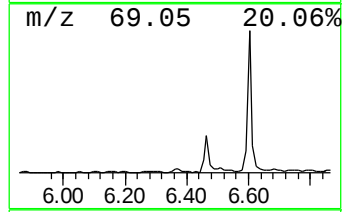
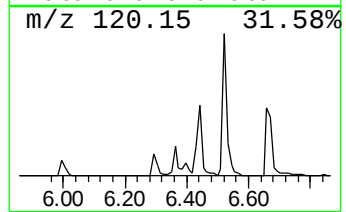
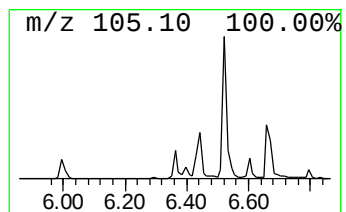
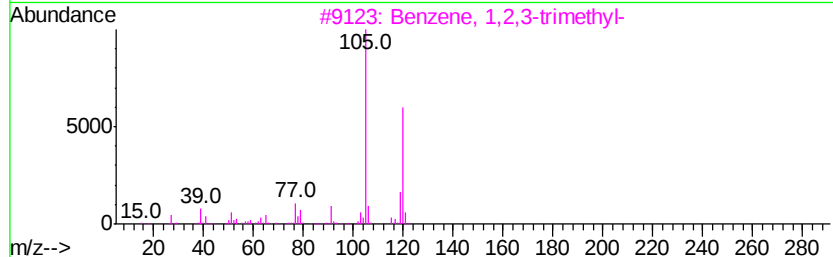
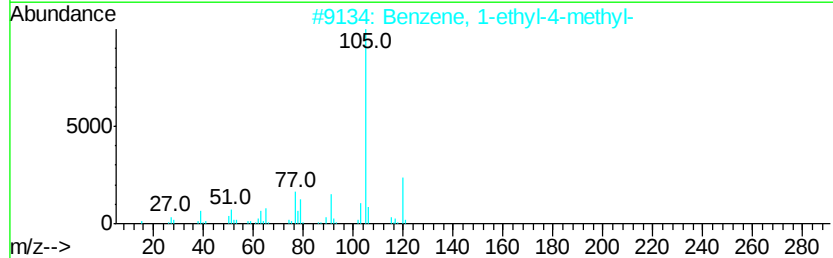
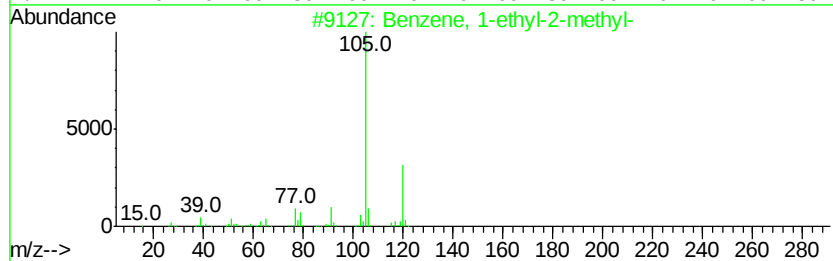
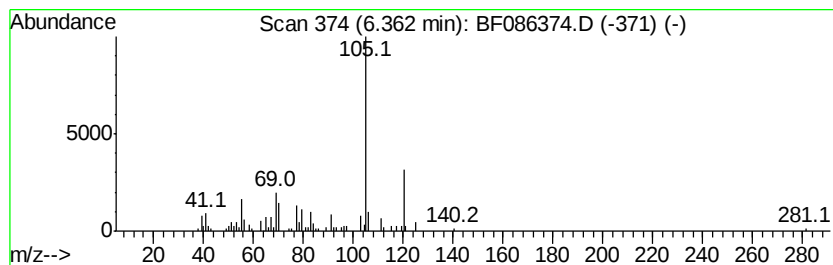
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.36	2.64 ng	126281	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	92
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	70
4		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	70
5		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	70



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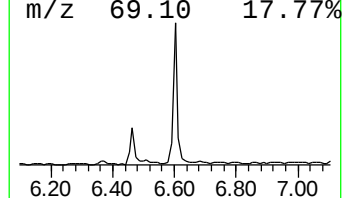
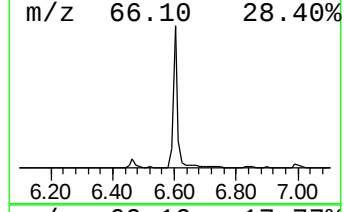
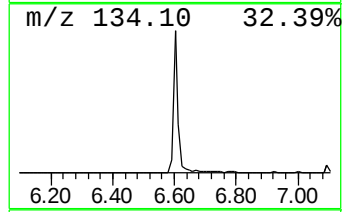
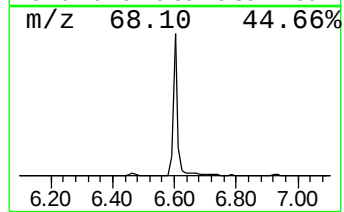
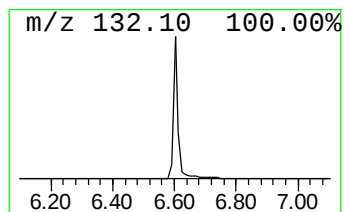
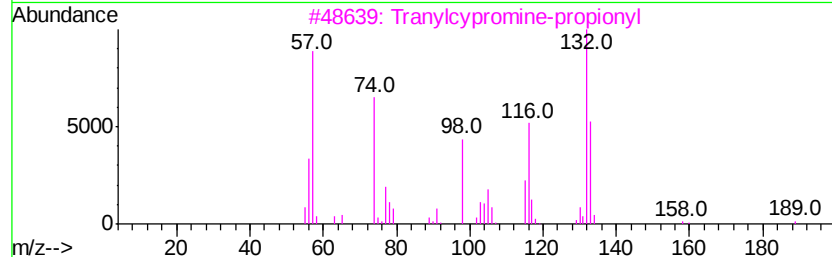
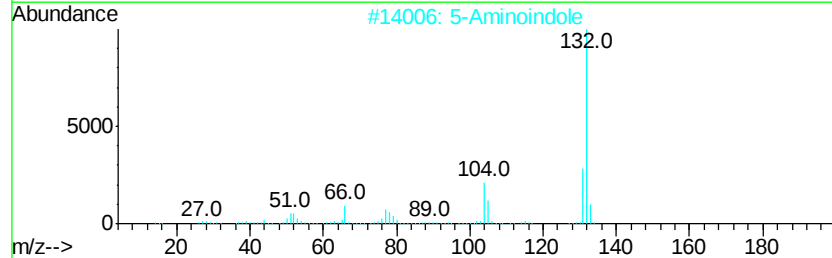
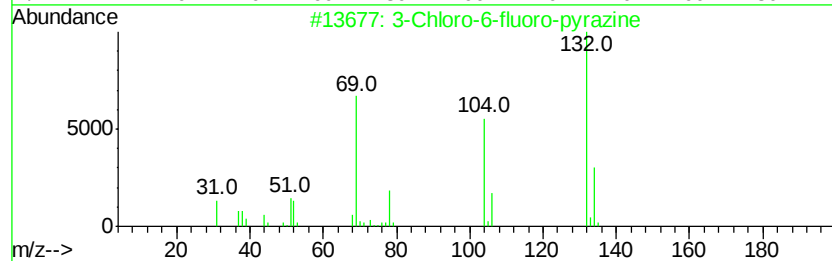
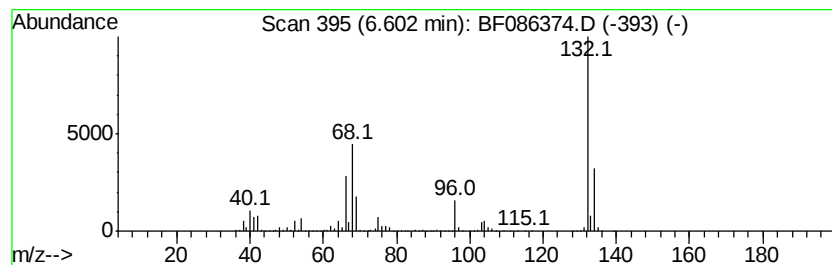
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 unknown6.60 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	60.76 ng	2905500	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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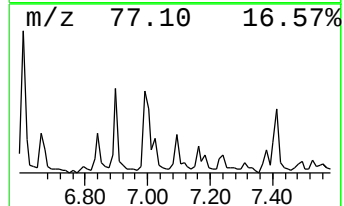
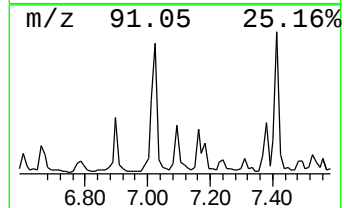
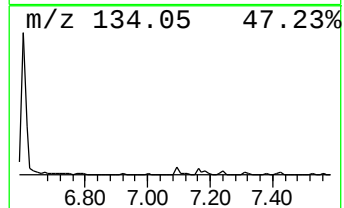
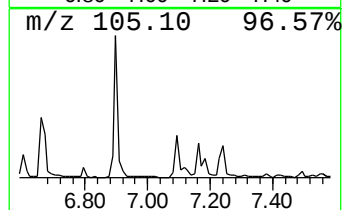
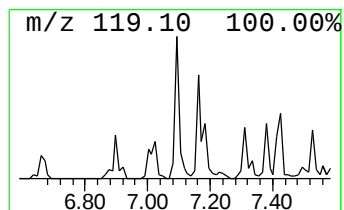
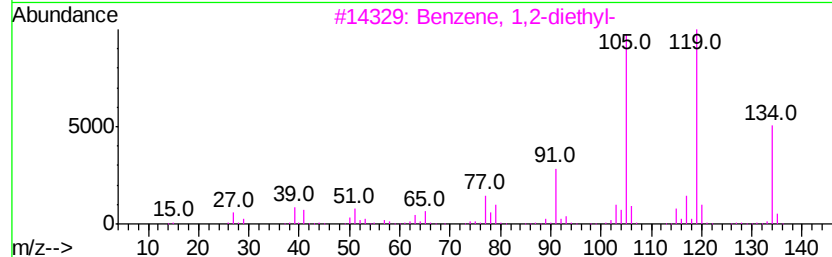
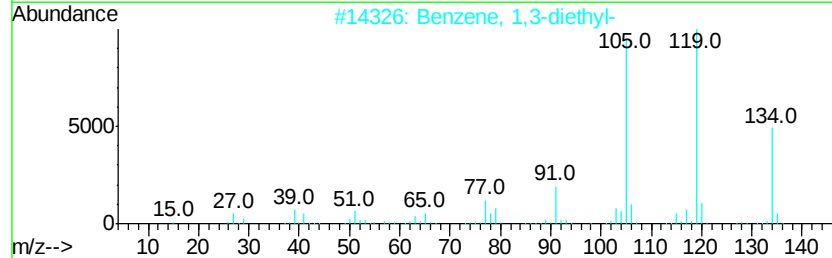
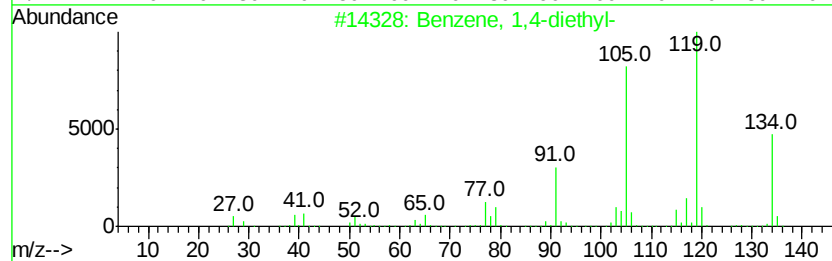
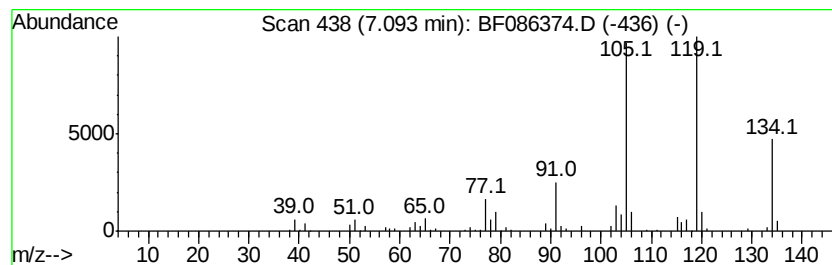
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1,4-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.09	4.40 ng	210222	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
3		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	91
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086374.D
Acq On : 23 Apr 2016 6:48
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Sample : H2634-08
Misc :
ALS Vial : 21 Sample Multiplier: 1

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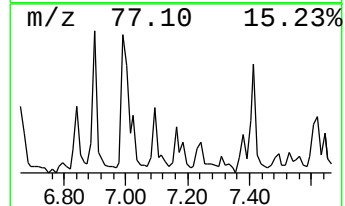
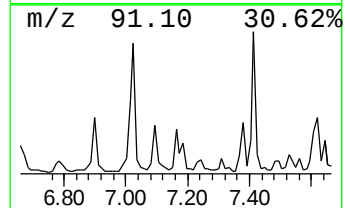
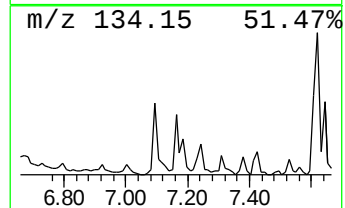
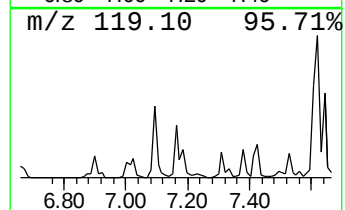
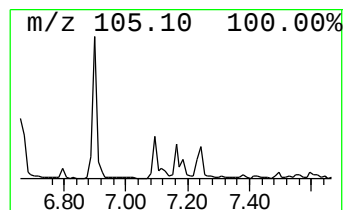
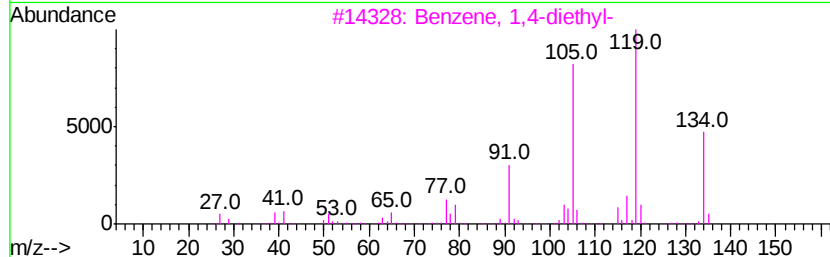
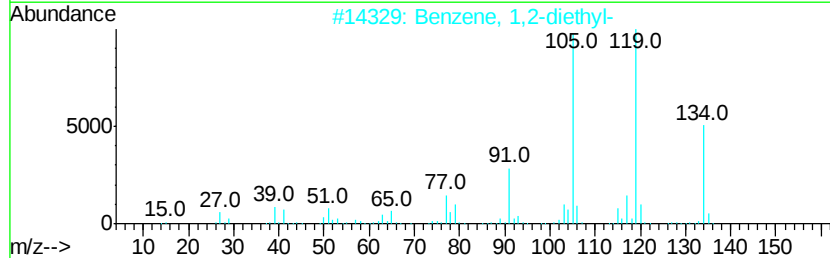
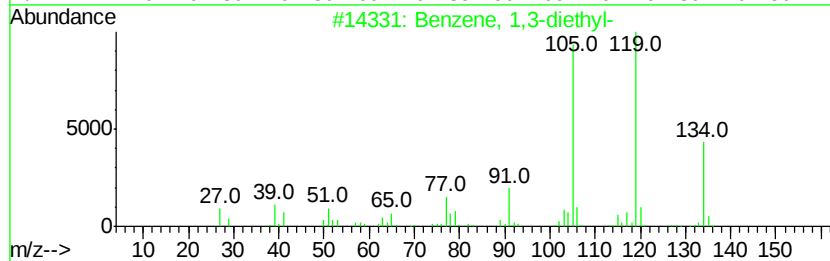
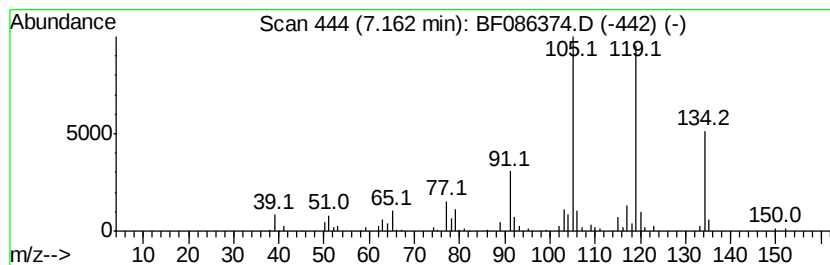
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,3-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.16	3.46 ng	165530	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
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Acq On : 23 Apr 2016 6:48
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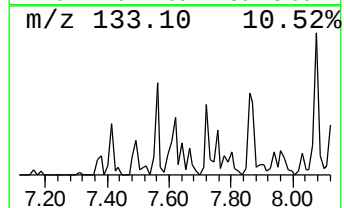
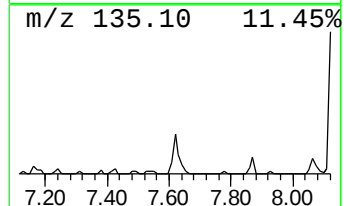
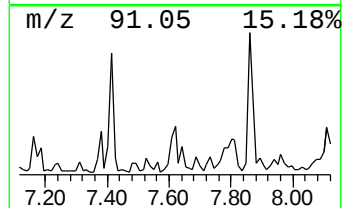
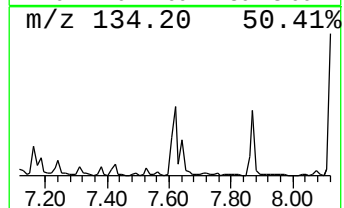
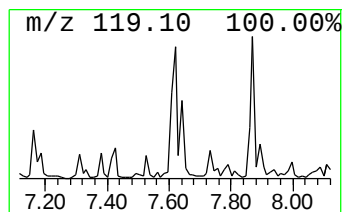
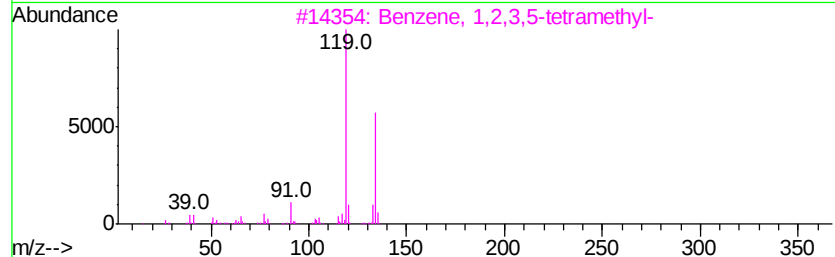
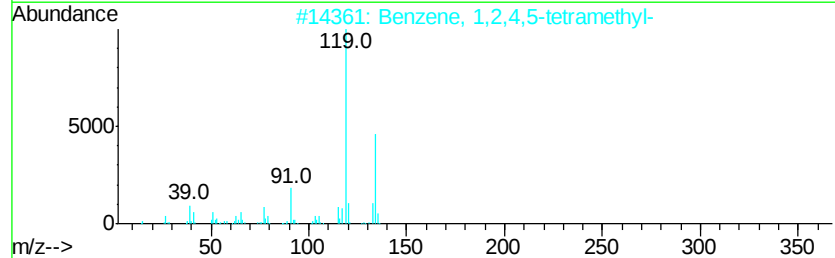
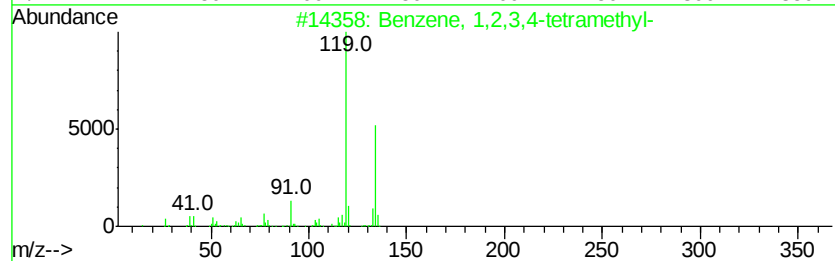
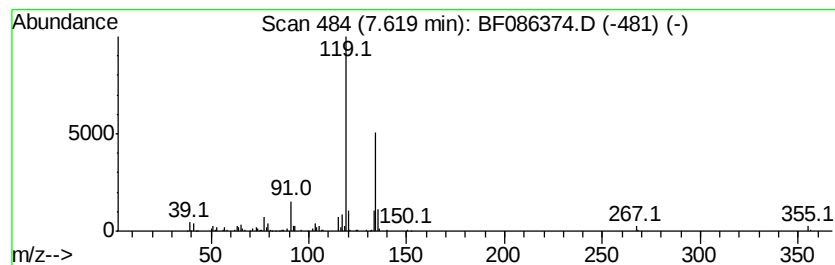
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.62	3.15 ng	227719	Naphthalene-d8	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



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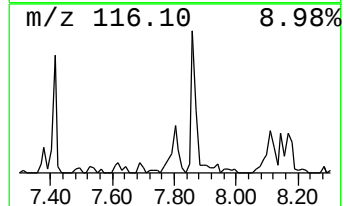
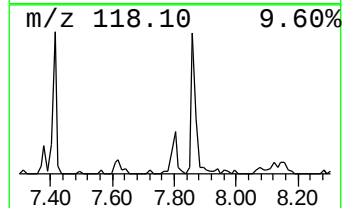
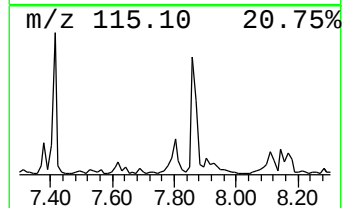
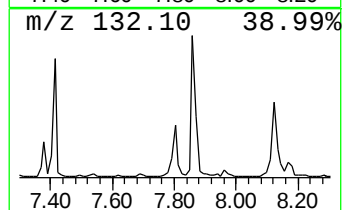
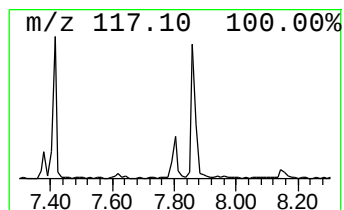
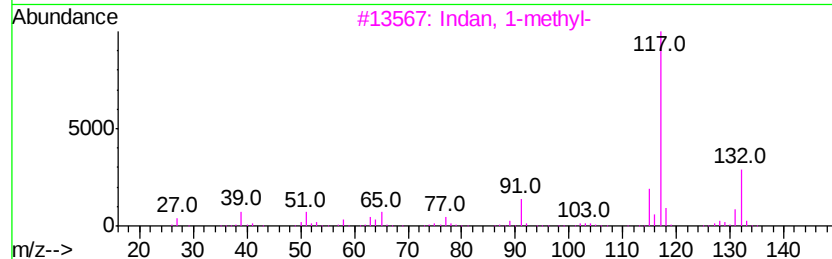
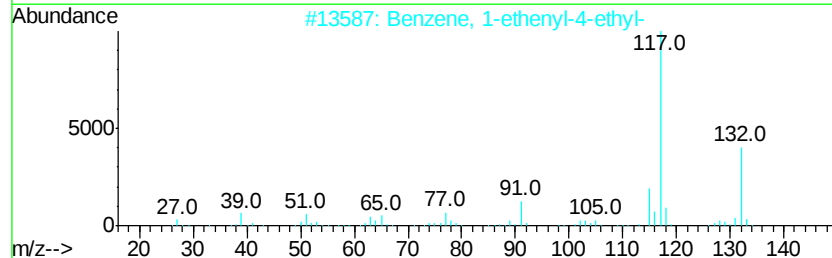
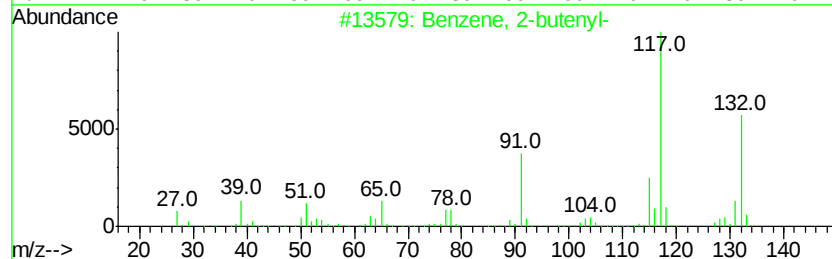
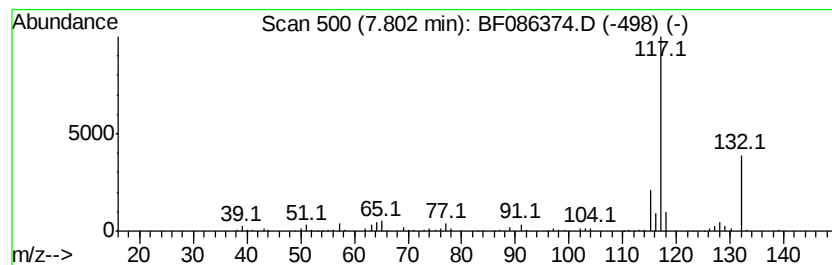
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 2-butenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	2.98 ng	215818	Naphthalene-d8	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
3		Indan, 1-methyl-	132	C10H12	000767-58-8	86
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83



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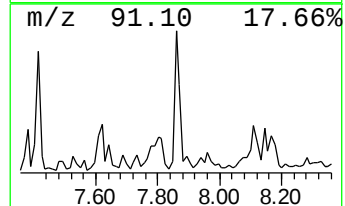
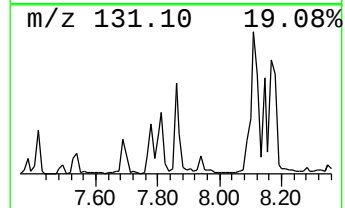
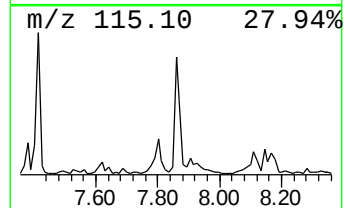
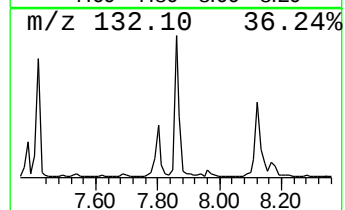
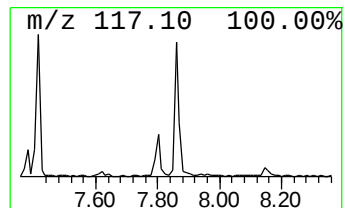
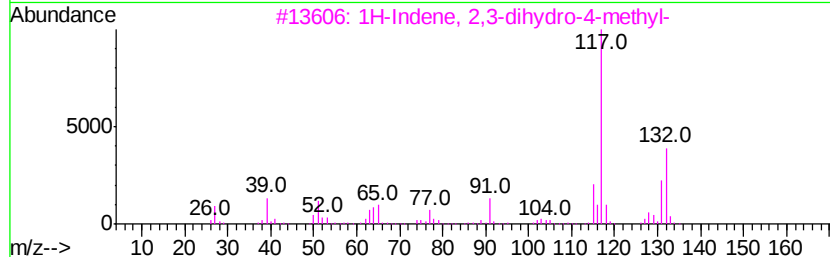
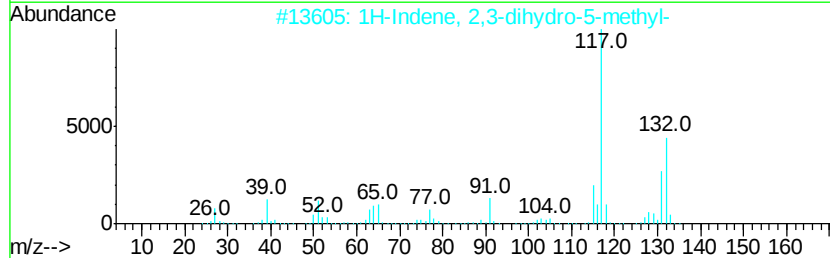
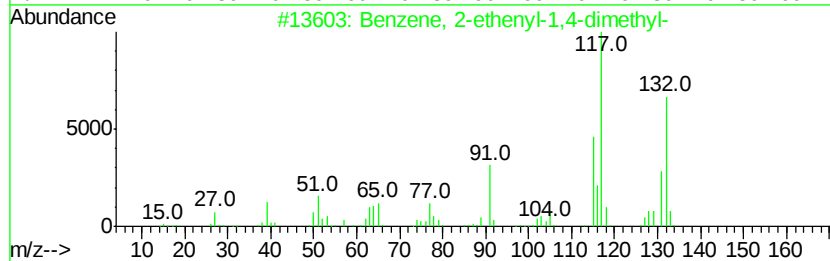
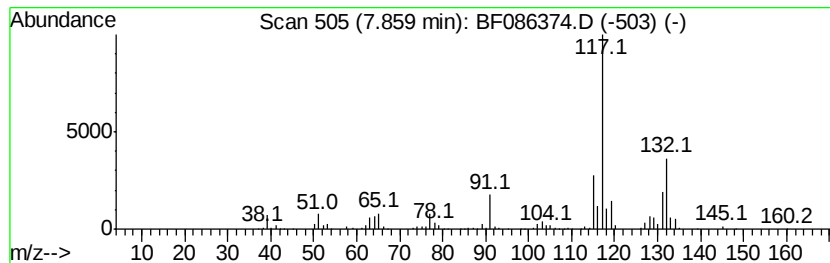
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	9.29 ng	672497	Naphthalene-d8	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	87



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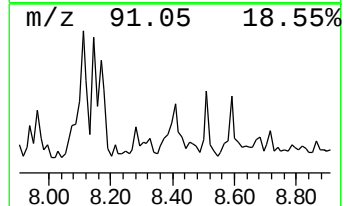
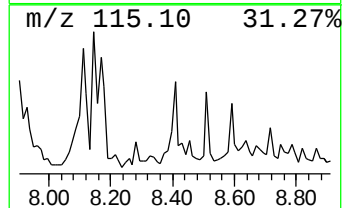
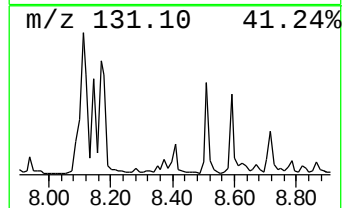
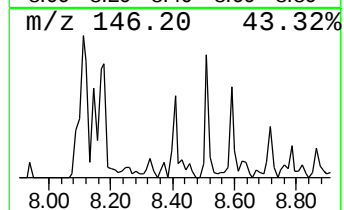
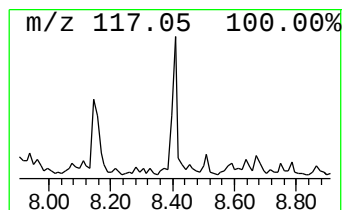
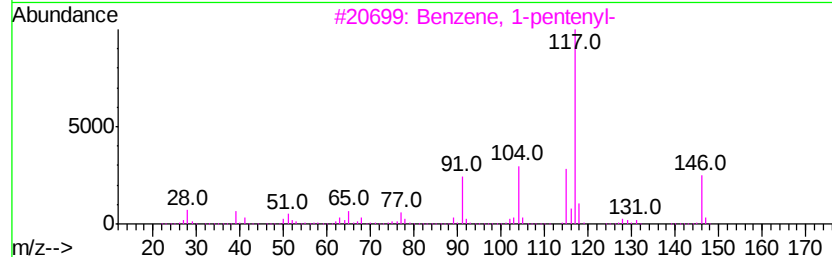
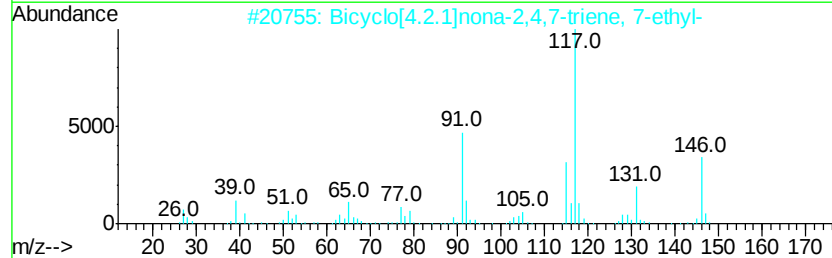
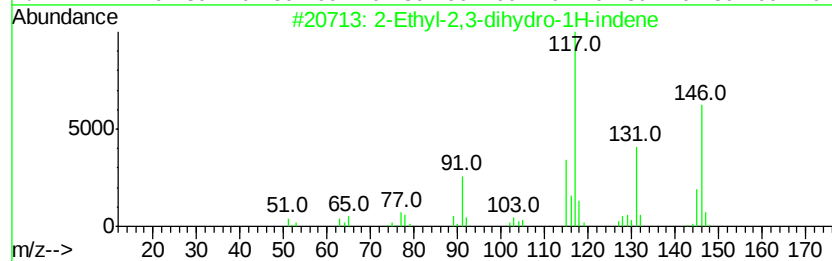
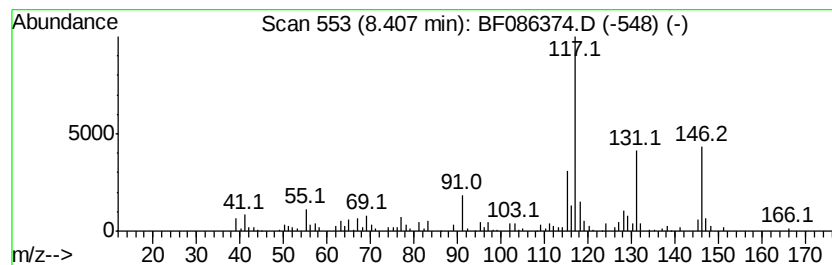
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 2-Ethyl-2,3-dihydro-1H-indene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.41	2.33 ng	168873	Napthalene-d8	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	91
2		Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	83
3		Benzene, 1-pentenyl-	146	C11H14	000826-18-6	64
4		trans-1-Phenyl-1-pentene	146	C11H14	016002-93-0	64
5		Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	60



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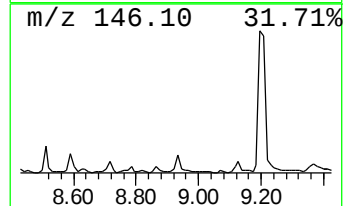
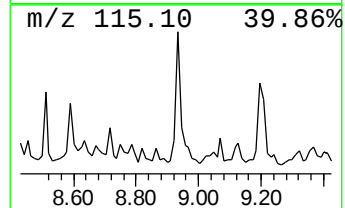
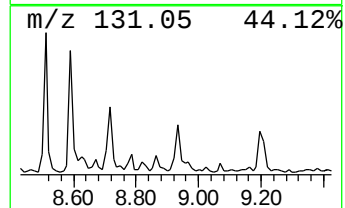
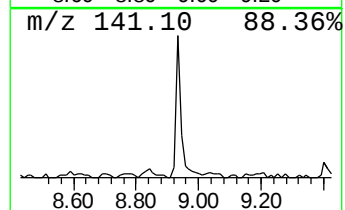
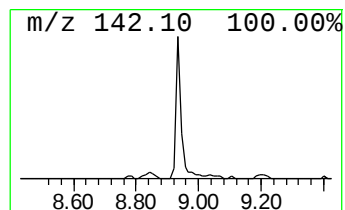
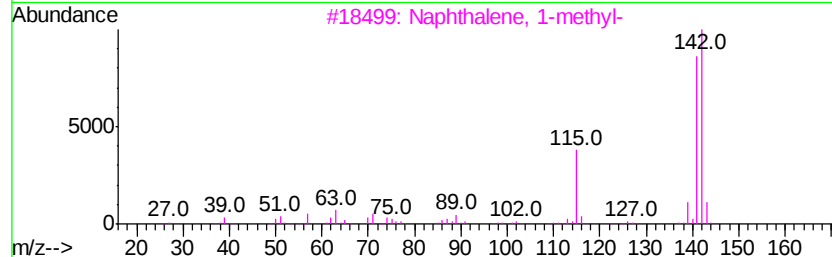
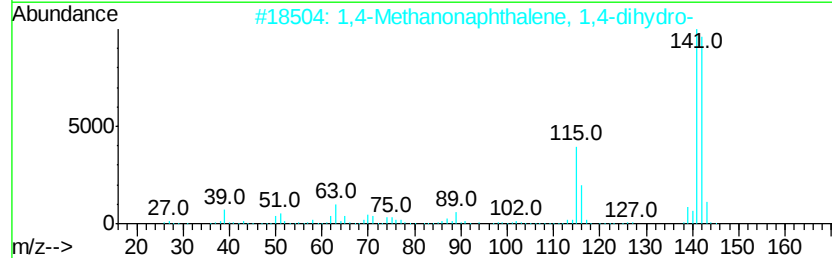
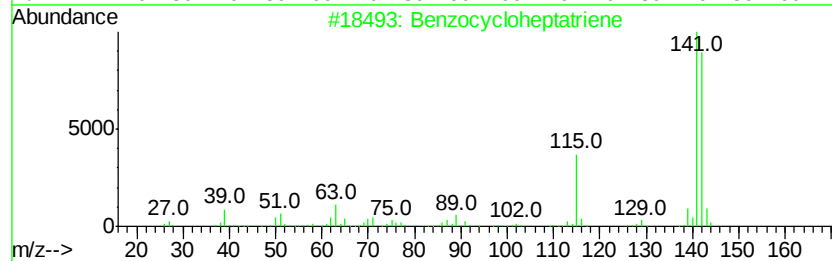
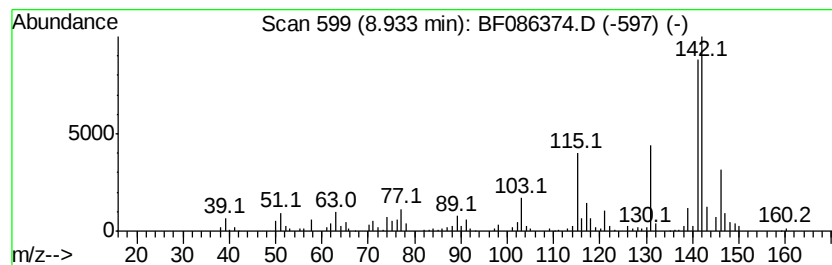
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzocycloheptatriene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.93	2.48 ng	179888	Naphthalene-d8	8.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzocycloheptatriene	142	C11H10	000264-09-5	70
2		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	70
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	70
4		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	70
5		2-Cyclopenten-1-ol, 1-phenyl-	160	C11H12O	056667-10-8	49



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ClientSampleId :
MW-11

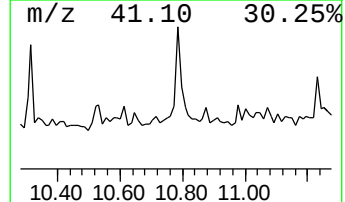
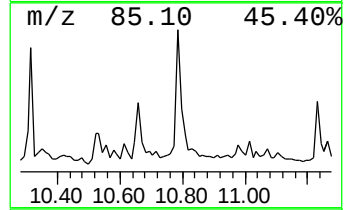
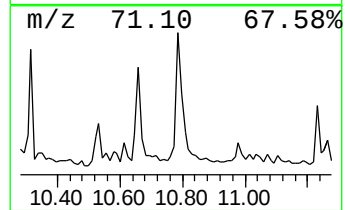
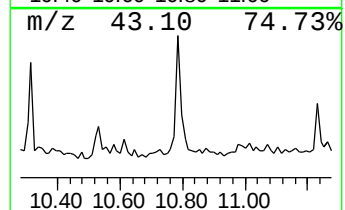
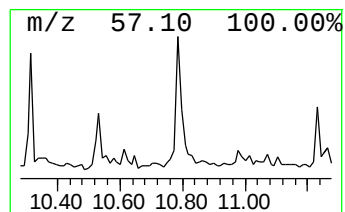
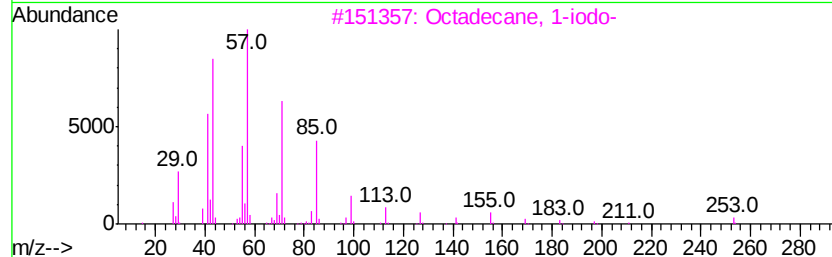
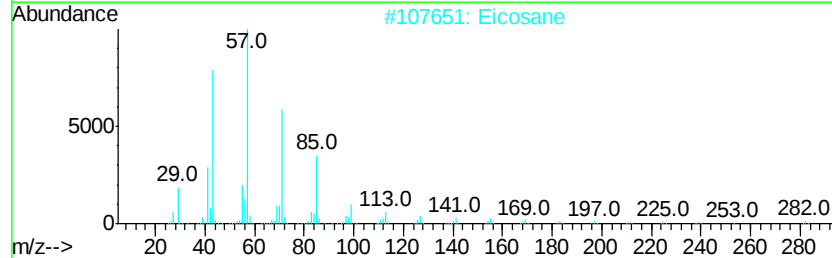
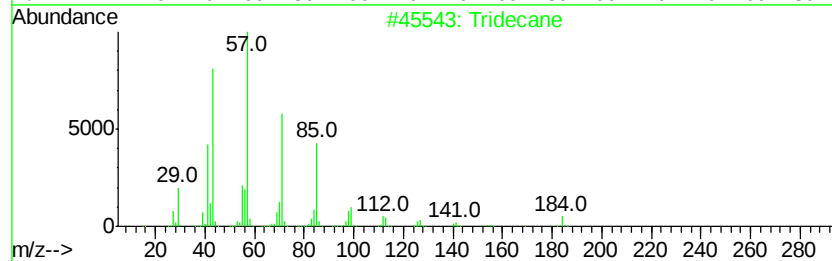
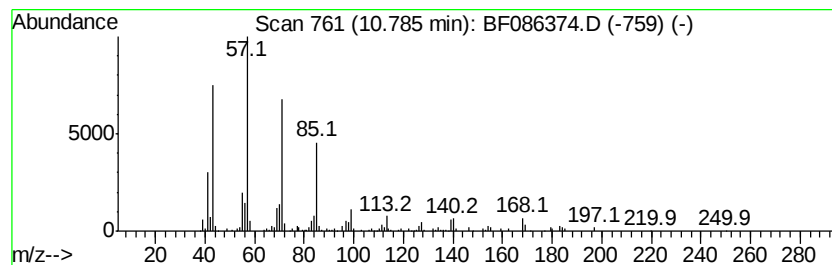
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tridecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	2.31 ng	191661	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	87
2		Eicosane	282	C20H42	000112-95-8	83
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	83
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
5		Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086374.D
Acq On : 23 Apr 2016 6:48
Operator : UM/SJ
Sample : H2634-08
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-11

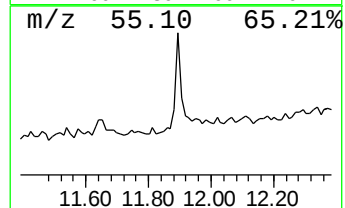
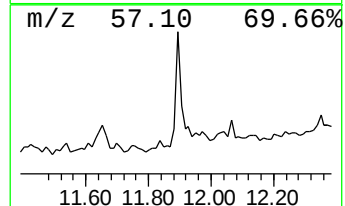
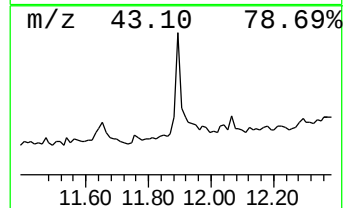
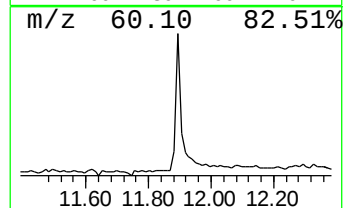
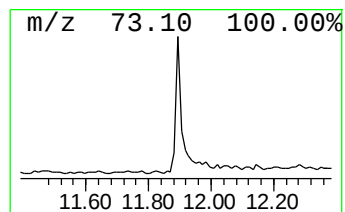
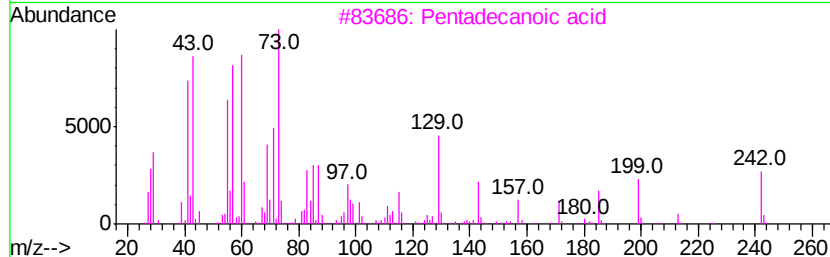
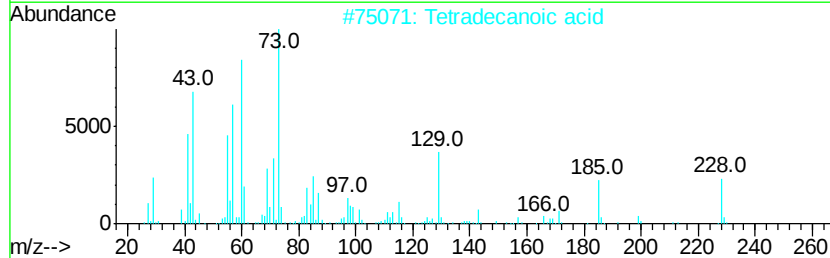
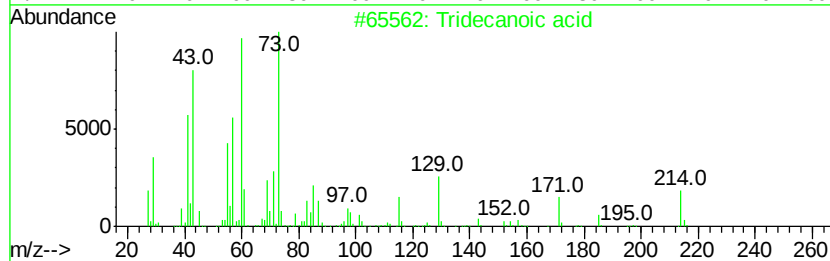
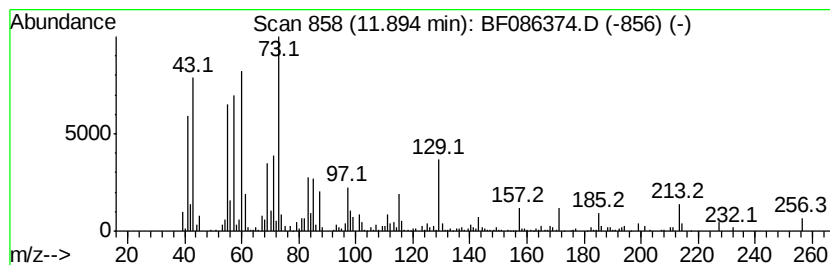
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Tridecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	3.09 ng	256194	Phenanthrene-d10	11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid	214	C13H26O2	000638-53-9	93
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	81
5			Undecane, 2-methyl-	170	C12H26	007045-71-8	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086374.D
Acq On : 23 Apr 2016 6:48
Operator : UM/SJ
Sample : H2634-08
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-11

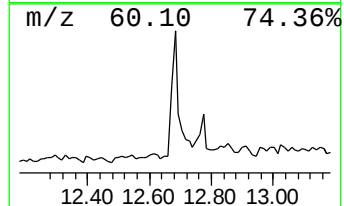
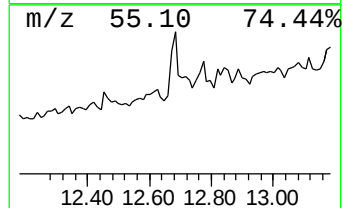
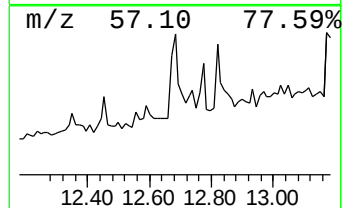
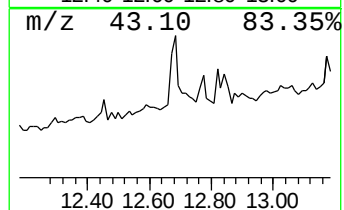
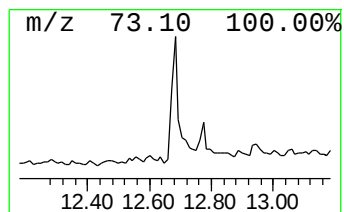
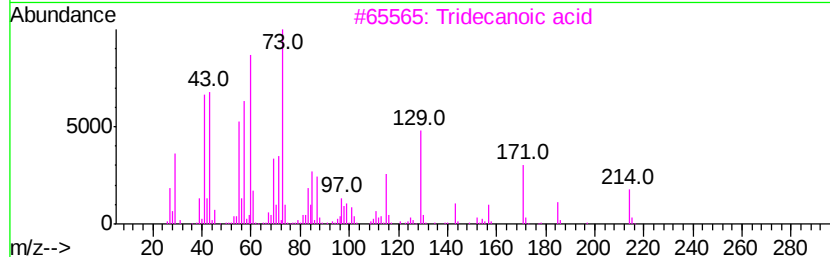
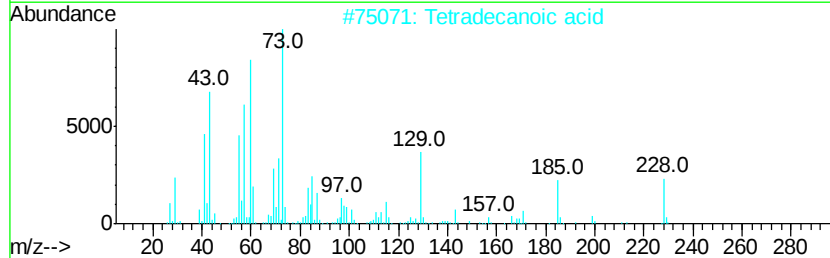
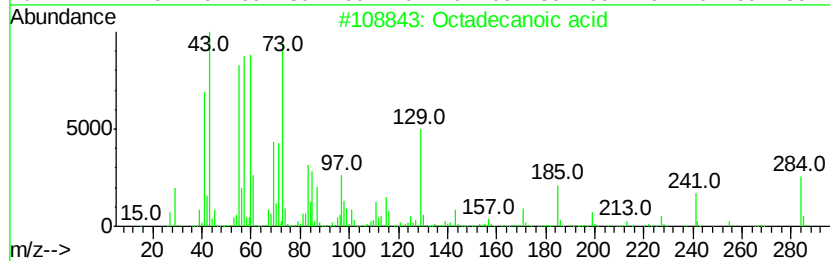
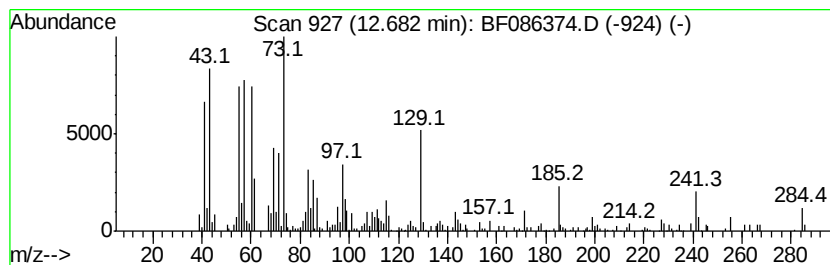
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	3.39 ng	220640	Chrysene-d12	13.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	95
3			Tridecanoic acid	214	C13H26O2	000638-53-9	64
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	62
5			Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
 Data File : BF086374.D
 Acq On : 23 Apr 2016 6:48
 Operator : UM/SJ
 Sample : H2634-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-11

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.00	4.0	ng	192746	1	6.84	956359	20.0
Benzene, 1,3-dime...	5.66	2.2	ng	106571	1	6.84	956359	20.0
Benzene, 1-ethyl-...	6.36	2.6	ng	126281	1	6.84	956359	20.0
unknown6.60	6.60	60.8	ng	2905500	1	6.84	956359	20.0
Benzene, 1,4-diet...	7.09	4.4	ng	210222	1	6.84	956359	20.0
Benzene, 1,3-diet...	7.16	3.5	ng	165530	1	6.84	956359	20.0
Benzene, 1,2,3,4-...	7.62	3.1	ng	227719	2	8.12	1447920	20.0
Benzene, 2-butenyl-	7.80	3.0	ng	215818	2	8.12	1447920	20.0
Benzene, 2-etheny...	7.86	9.3	ng	672497	2	8.12	1447920	20.0
2-Ethyl-2,3-dihyd...	8.41	2.3	ng	168873	2	8.12	1447920	20.0
Benzocycloheptatr...	8.93	2.5	ng	179888	2	8.12	1447920	20.0
Tridecane	10.78	2.3	ng	191661	4	11.36	1657870	20.0
Tridecanoic acid	11.89	3.1	ng	256194	4	11.36	1657870	20.0
Octadecanoic acid	12.68	3.4	ng	220640	5	13.99	1300870	20.0