

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031315\
 Data File : BF077741.D
 Acq On : 13 Mar 2015 15:53
 Operator : TP/IZ
 Sample : G1517-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW2

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-BF031115.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	1.447	46	49	52	rVB	259623	432433	24.54%	3.483%
2	1.618	61	64	67	rBV	238103	335124	19.02%	2.700%
3	1.710	71	72	77	rVB2	20021	25341	1.44%	0.204%
4	1.790	77	79	87	rBV	59728	85663	4.86%	0.690%
5	4.933	352	354	359	rBV	36936	42915	2.44%	0.346%
6	5.459	397	400	403	rBV	459143	536195	30.42%	4.319%
7	6.739	509	512	518	rBV	225966	338071	19.18%	2.723%
8	6.876	522	524	527	rBV	1115850	1055065	59.87%	8.499%
9	7.173	547	550	552	rBV	219549	287293	16.30%	2.314%
10	7.356	563	566	568	rBV	1210312	1109282	62.94%	8.936%
11	7.859	608	610	613	rBV	823505	777478	44.11%	6.263%
12	8.739	685	687	690	rBV	323509	395958	22.47%	3.190%
13	9.448	747	749	752	rBV	30978	36059	2.05%	0.290%
14	10.088	803	805	808	rBV	1969211	1762393	100.00%	14.197%
15	10.591	847	849	852	rBV	35324	38533	2.19%	0.310%
16	10.911	875	877	879	rBV	535166	492188	27.93%	3.965%
17	10.945	879	880	883	rVB	61568	82562	4.68%	0.665%
18	11.448	920	924	927	rBV2	16168	23166	1.31%	0.187%
19	11.814	953	956	960	rVB	126261	116657	6.62%	0.940%
20	11.894	960	963	965	rBV	815098	1003753	56.95%	8.086%
21	12.751	1036	1038	1040	rBV	498142	509374	28.90%	4.103%
22	14.740	1210	1212	1215	rBV	1388079	1581168	89.72%	12.737%
23	15.940	1315	1317	1323	rBV	84422	134241	7.62%	1.081%
24	16.043	1323	1326	1329	rBV	551167	583144	33.09%	4.698%
25	17.849	1481	1484	1487	rBV	445243	568517	32.26%	4.580%
26	18.512	1539	1542	1549	rVB7	8604	27895	1.58%	0.225%
27	18.957	1576	1581	1586	rVB2	10947	33355	1.89%	0.269%

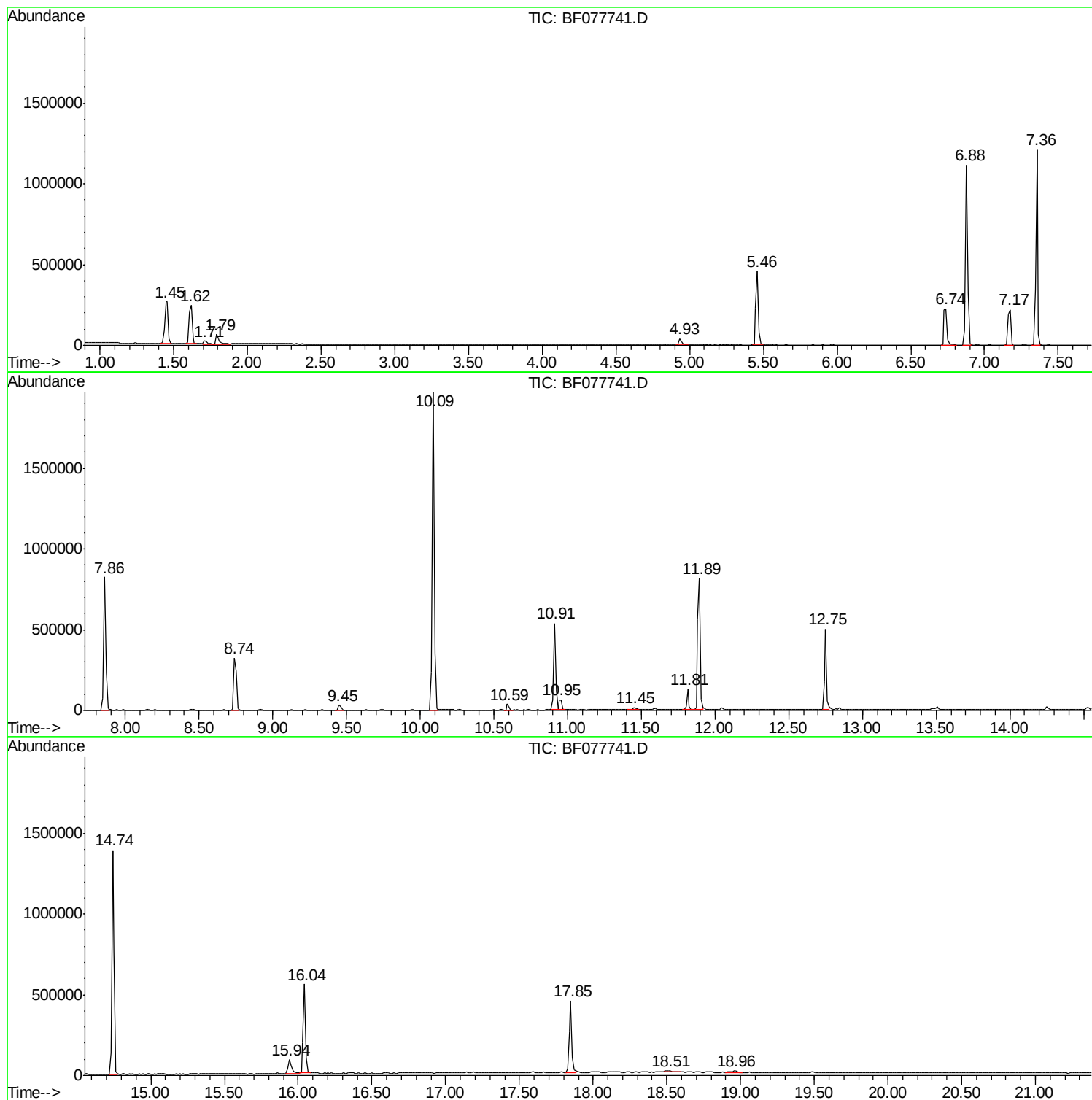
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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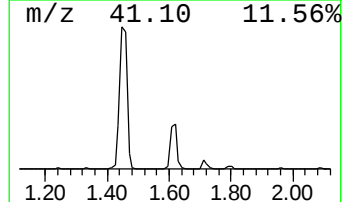
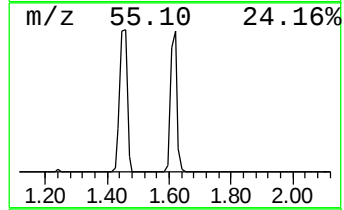
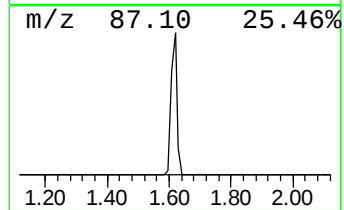
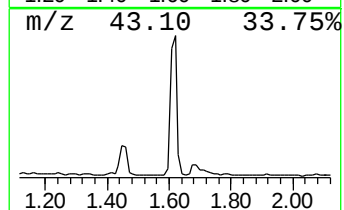
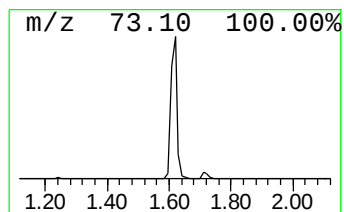
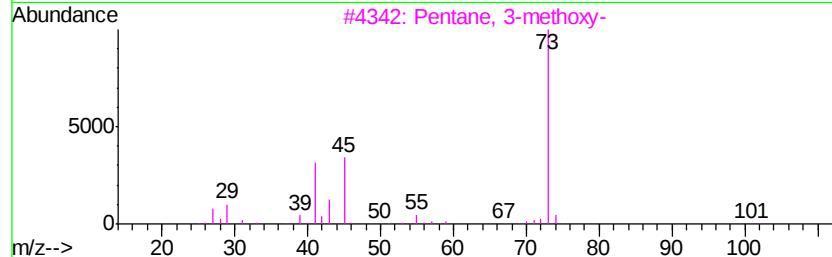
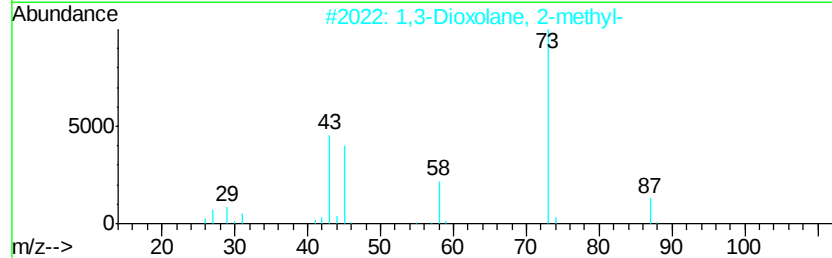
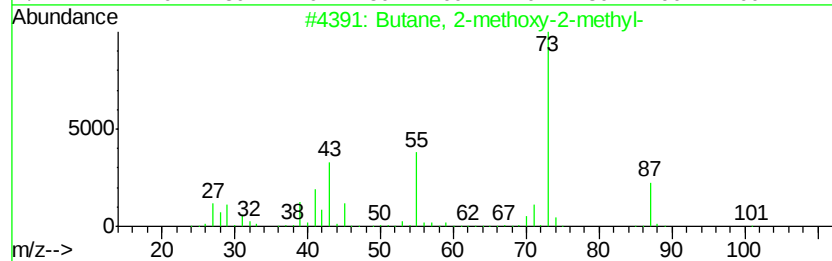
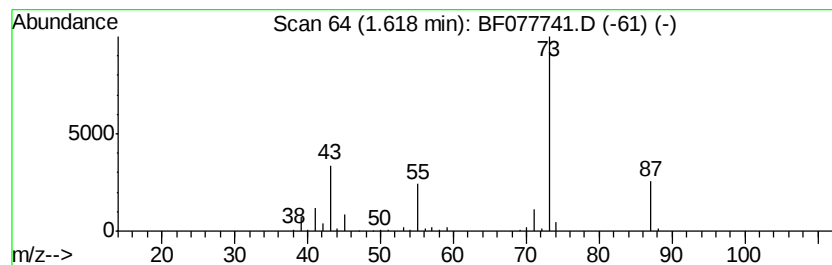
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.62	23.33 ng	335124	1,4-Dichlorobenzene-d4	7.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10



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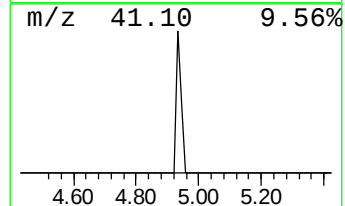
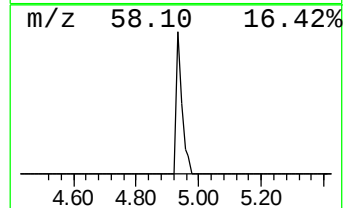
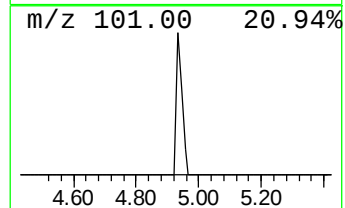
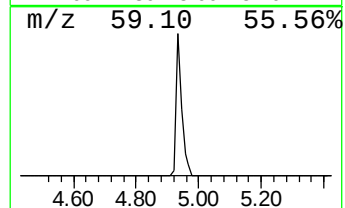
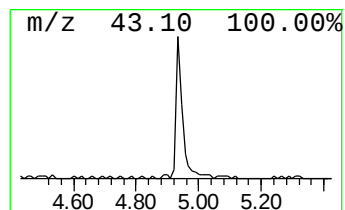
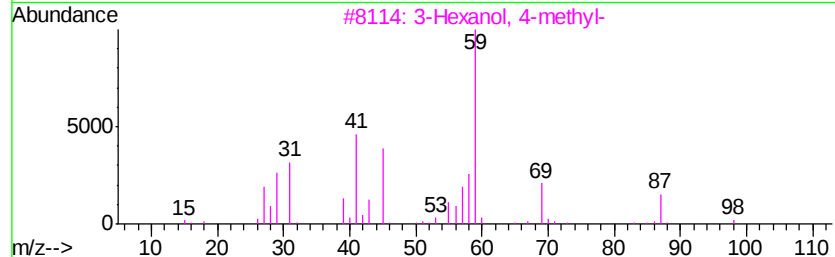
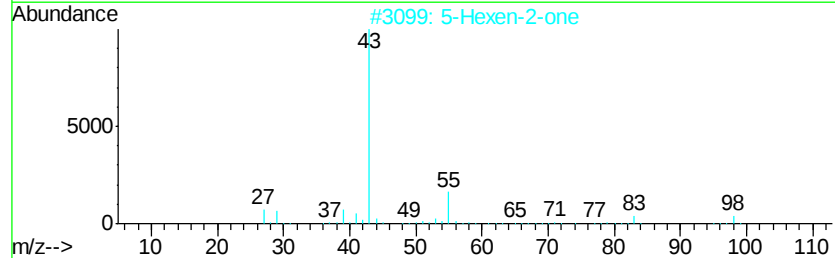
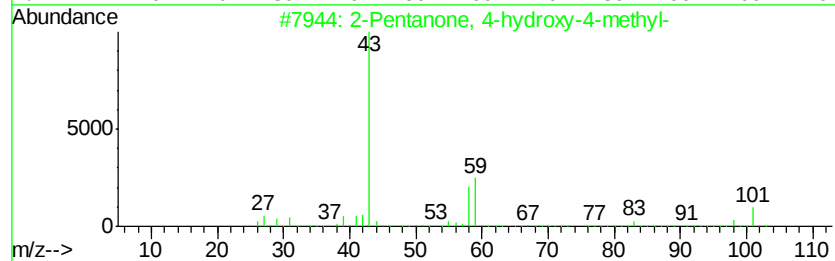
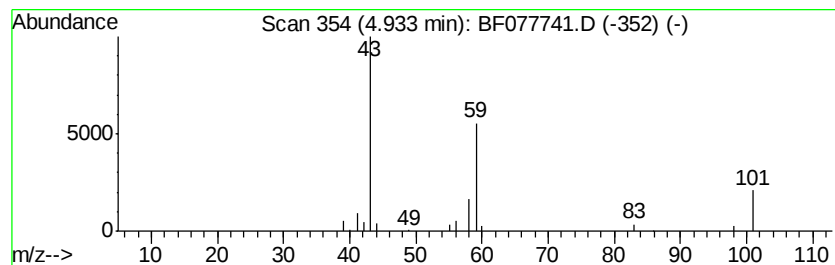
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	2.99 ng	42915	1,4-Dichlorobenzene-d4	7.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9
5		3-Hexanol	102	C6H14O	000623-37-0	9



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Instrument :
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ClientSampleId :
MW2

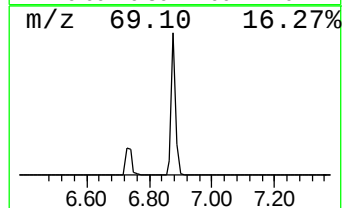
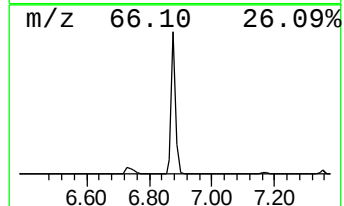
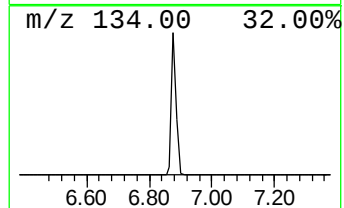
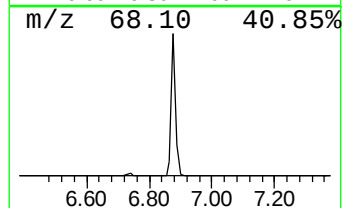
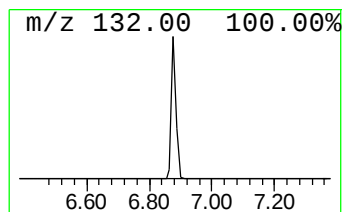
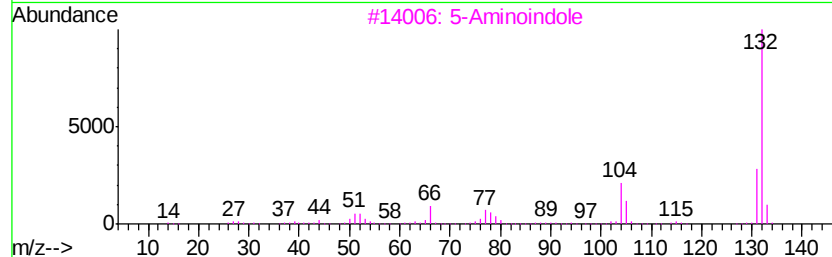
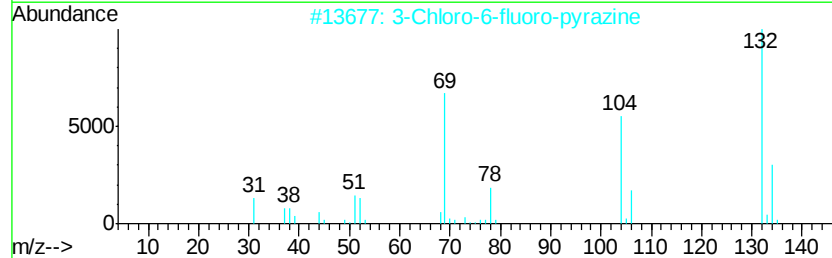
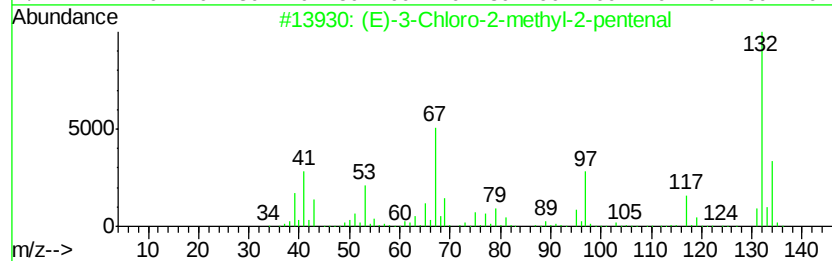
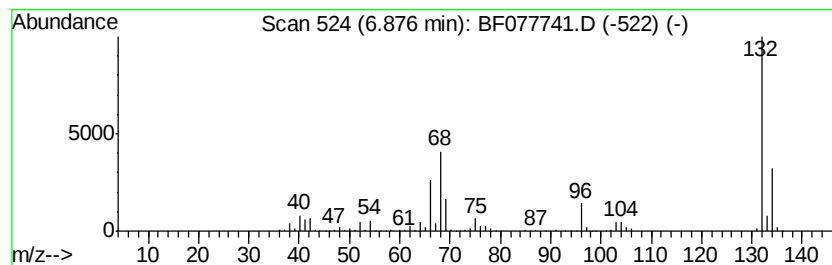
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.88 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	73.45 ng	1055070	1,4-Dichlorobenzene-d4	7.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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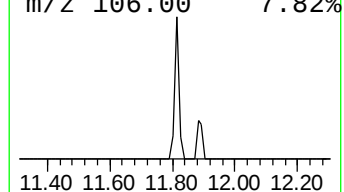
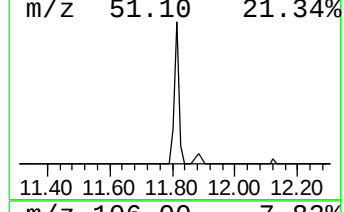
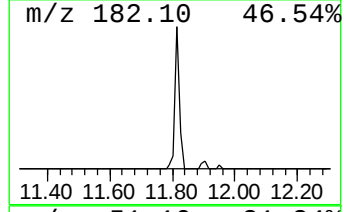
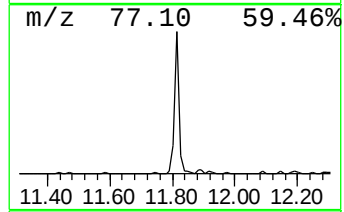
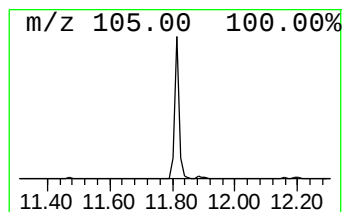
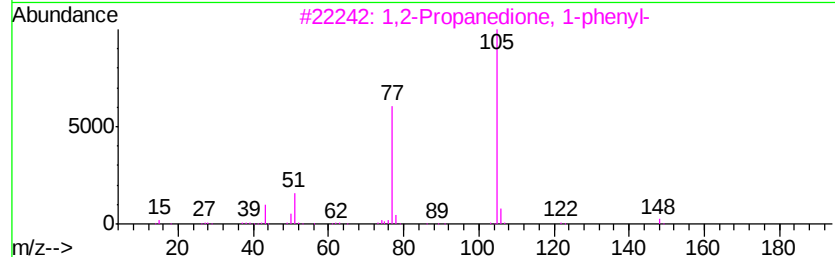
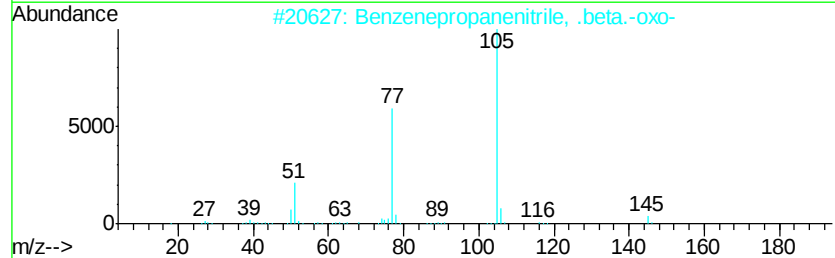
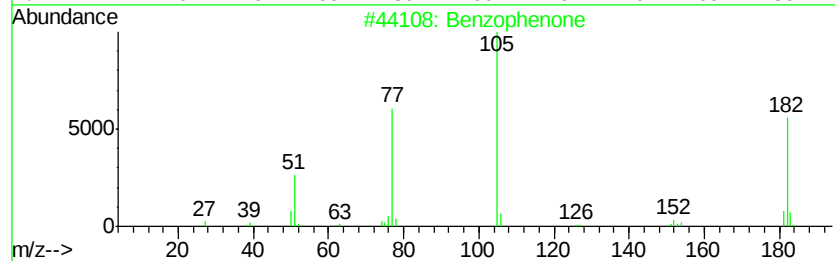
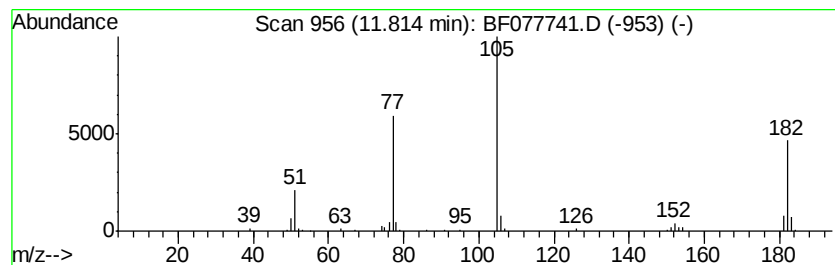
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.81	4.74 ng	116657	Acenaphthene-d10		10.91	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	96
2		Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	47
3		1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	47
4		5-[N-Benzoylaminomethyl]tetrazole	203	C9H9N5O	1000227-36-3	47
5		1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47



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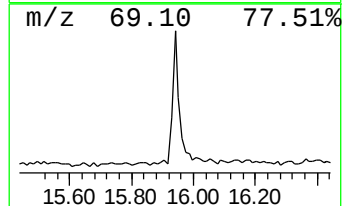
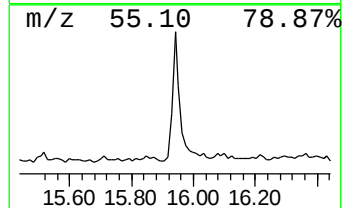
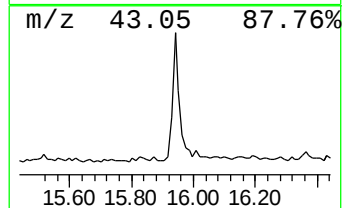
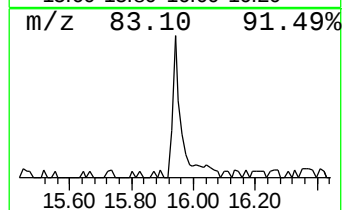
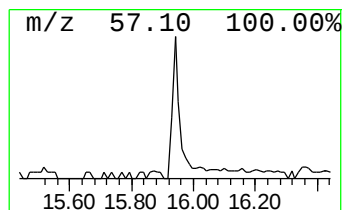
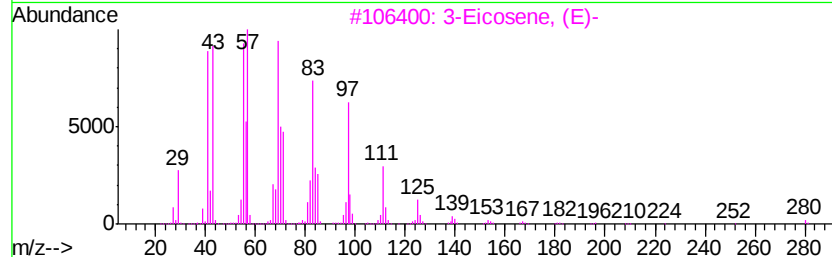
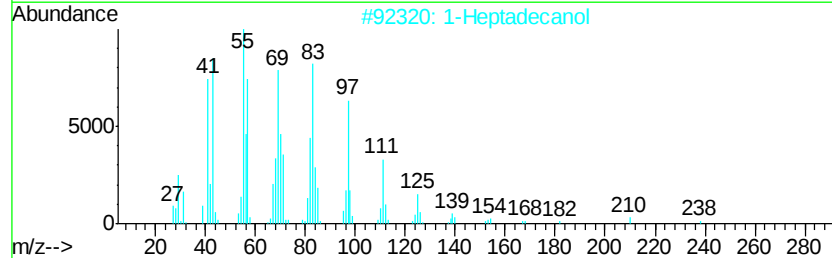
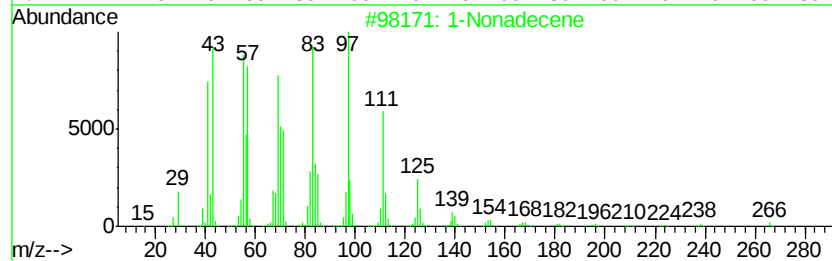
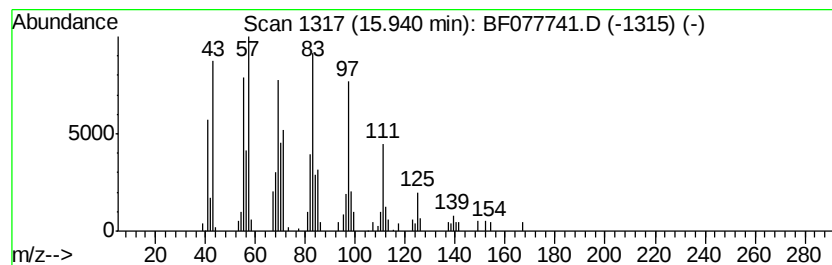
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1-Nonadecene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.94	4.60 ng	134241	Chrysene-d12	16.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene		266	C19H38	018435-45-5	91
2	1-Heptadecanol		256	C17H36O	001454-85-9	91
3	3-Eicosene, (E)-		280	C20H40	074685-33-9	90
4	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	90
5	9-Eicosene, (E)-		280	C20H40	074685-29-3	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.62	23.3	ng	335124	1	7.17	287293	20.0
2-Pentanone, 4-hy...	4.93	3.0	ng	42915	1	7.17	287293	20.0
unknown6.88	6.88	73.5	ng	1055070	1	7.17	287293	20.0
Benzophenone	11.81	4.7	ng	116657	3	10.91	492188	20.0
1-Nonadecene	15.94	4.6	ng	134241	5	16.04	583144	20.0