

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080615\
 Data File : BF080912.D
 Acq On : 7 Aug 2015 4:13
 Operator : TP/UM
 Sample : PB84869BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84869BL

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080615.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	4.956	249	251	254	rBV	124813	124860	15.62%	2.268%
2	5.379	286	288	291	rBV	179564	184893	23.13%	3.359%
3	6.430	377	380	382	rBV	127502	175176	21.92%	3.182%
4	6.567	390	392	394	rBV	218045	188981	23.65%	3.433%
5	6.807	411	413	415	rBV	563653	483510	60.50%	8.783%
6	6.956	424	426	428	rVB	139872	137339	17.18%	2.495%
7	7.367	459	462	464	rBB	138532	118122	14.78%	2.146%
8	8.087	523	525	528	rBV	555211	619504	77.52%	11.253%
9	9.162	617	619	622	rBV	190006	204435	25.58%	3.714%
10	9.848	676	679	681	rBV	700174	743055	92.97%	13.498%
11	10.625	745	747	749	rBV	137055	117677	14.72%	2.138%
12	11.322	806	808	810	rBV	955815	799205	100.00%	14.518%
13	12.899	944	946	949	rBB	258793	230696	28.87%	4.191%
14	13.951	1035	1038	1040	rBV	821019	758380	94.89%	13.776%
15	15.357	1158	1161	1164	rBV	512129	619157	77.47%	11.247%

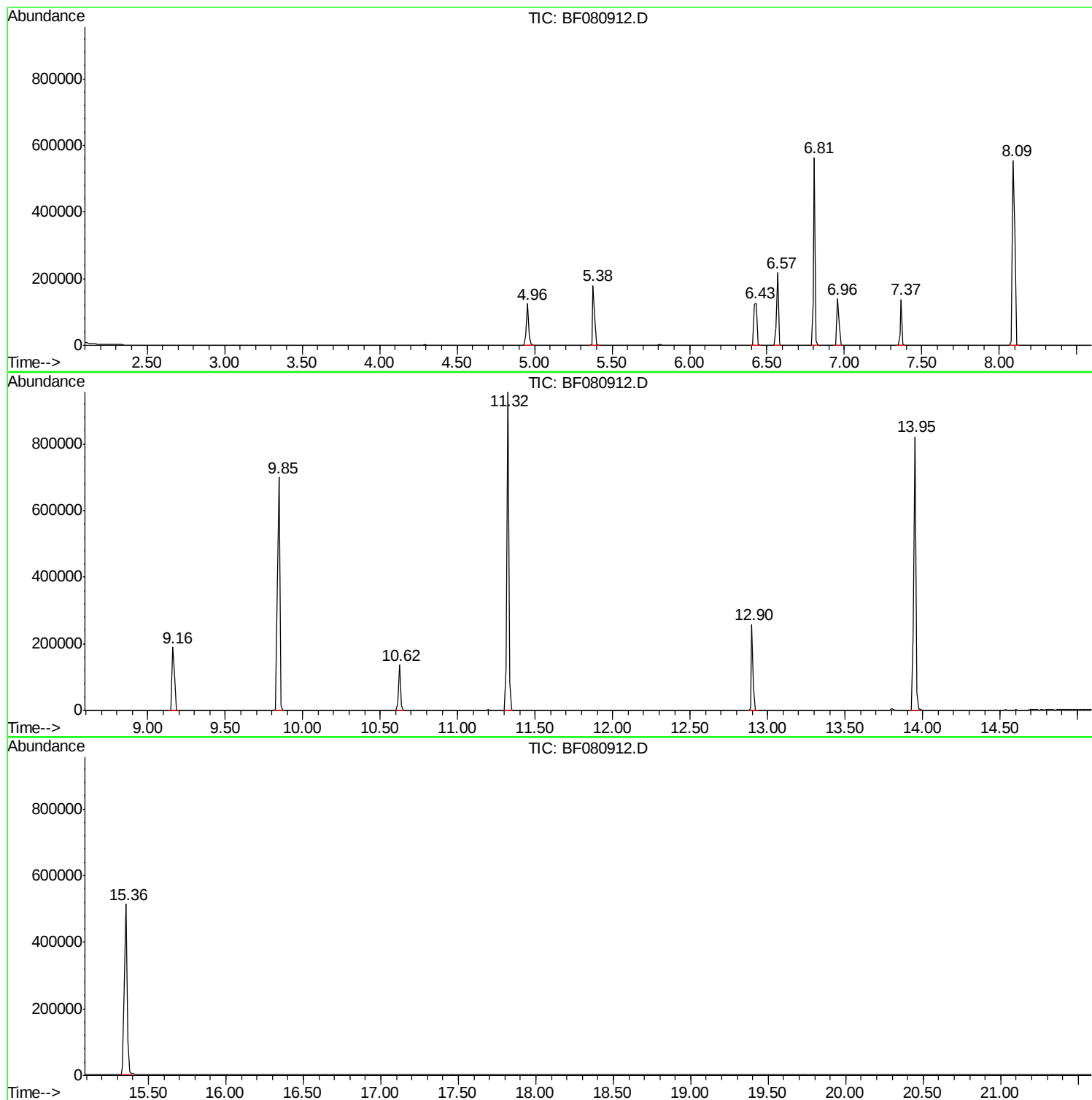
Sum of corrected areas: 5504990

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080615.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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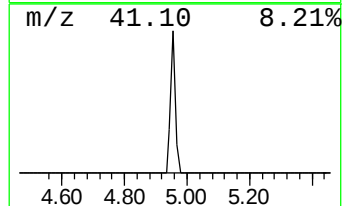
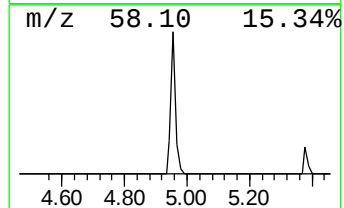
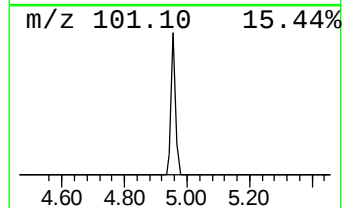
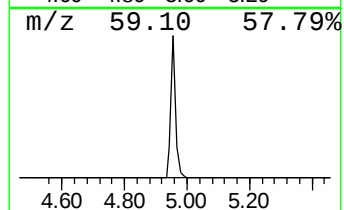
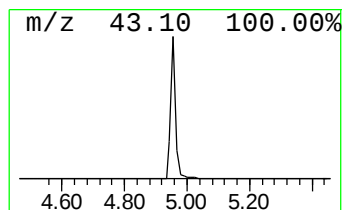
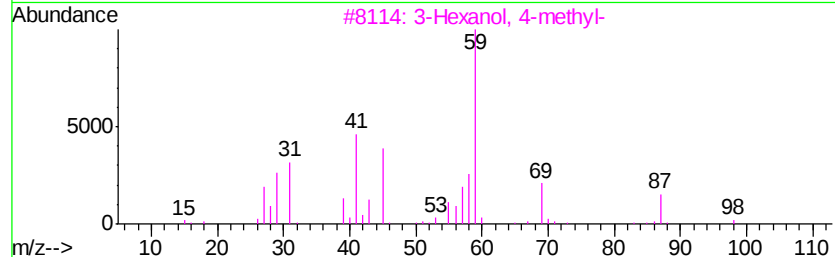
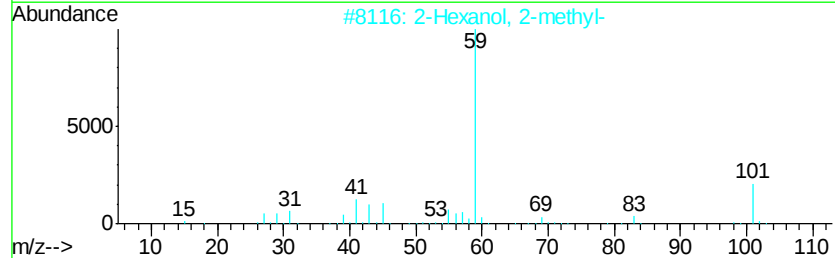
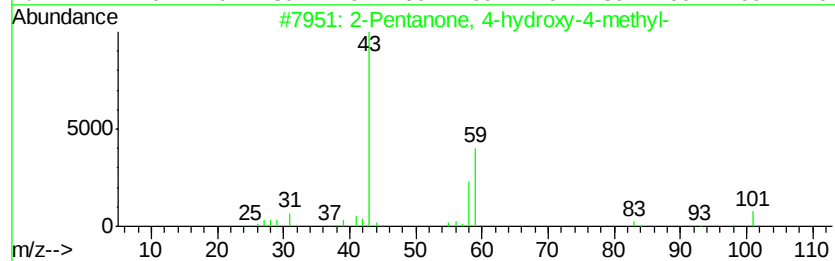
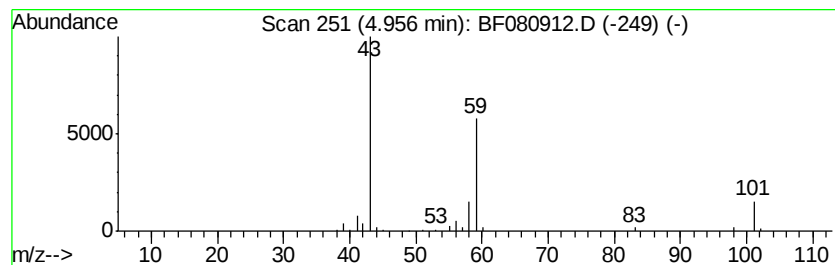
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	5.16 ng	124860	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	17



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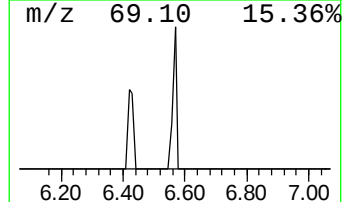
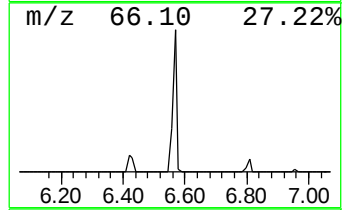
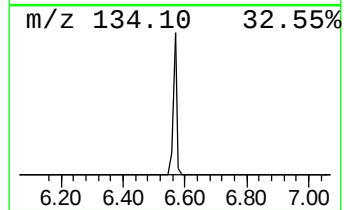
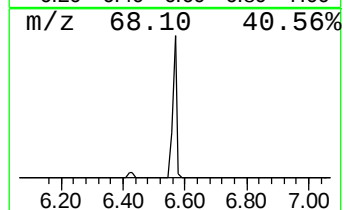
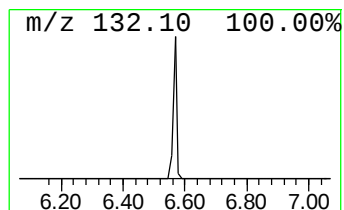
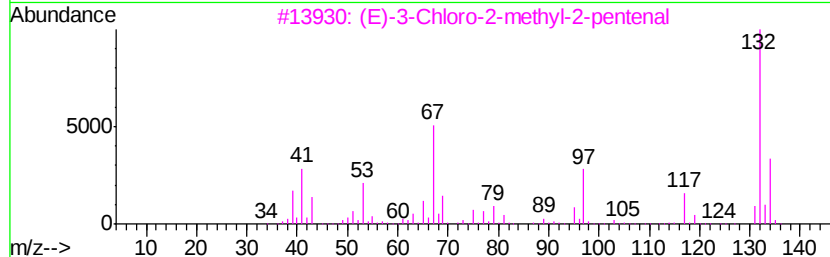
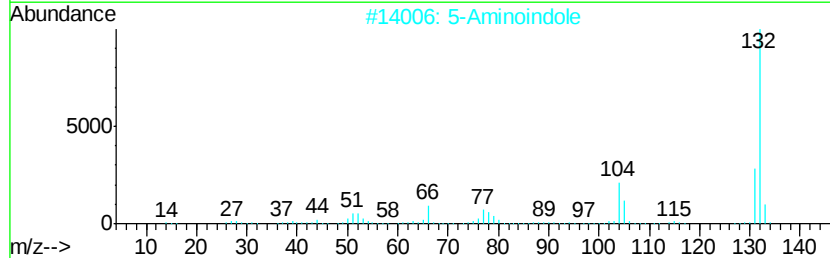
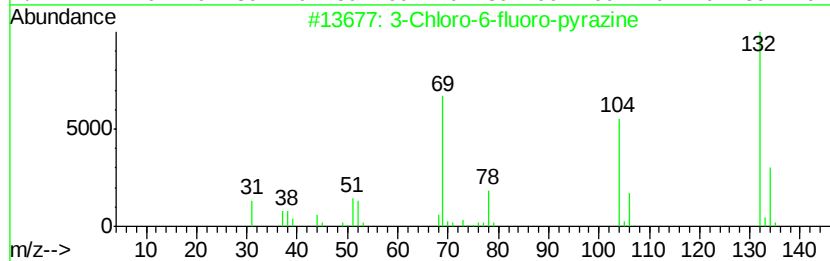
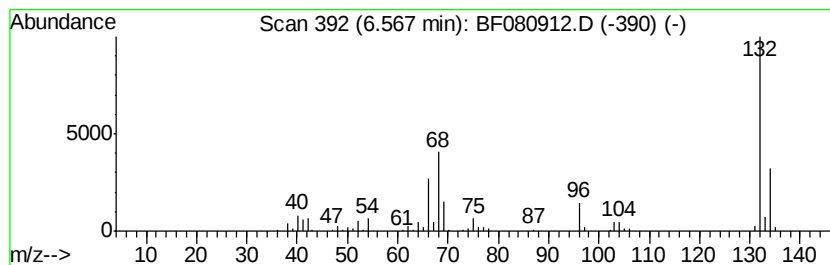
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Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	7.82 ng	188981	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	5.2	ng	124860	1	6.81	483510	20.0
unknown6.57	6.57	7.8	ng	188981	1	6.81	483510	20.0