

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030317\
 Data File : BF093364.D
 Acq On : 3 Mar 2017 17:49
 Operator : SJ/MA
 Sample : PB97311BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB97311BL

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-BF013017.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	248	251	266	rVB	420635	805444	15.37%	1.950%
2	5.390	286	289	291	rBV	3005814	4338693	82.77%	10.504%
3	6.442	378	381	383	rBV	3617073	4695442	89.57%	11.367%
4	6.556	388	391	393	rBV	4321484	4130976	78.81%	10.001%
5	6.784	409	411	413	rBV	1172987	1051012	20.05%	2.544%
6	6.944	422	425	427	rBV	2943719	3538725	67.51%	8.567%
7	7.356	458	461	464	rBV	2592794	2855707	54.48%	6.913%
8	8.076	521	524	526	rBV	1297757	1514162	28.89%	3.666%
9	9.150	614	618	620	rBV	4744334	5241917	100.00%	12.690%
10	9.825	674	677	679	rBV	1503289	1610861	30.73%	3.900%
11	10.613	743	746	749	rBV	2464699	2585698	49.33%	6.260%
12	11.311	804	807	809	rBV	1172273	1602513	30.57%	3.880%
13	12.888	942	945	948	rBV	3618293	4644302	88.60%	11.243%
14	13.939	1035	1037	1040	rBV	1261999	1492913	28.48%	3.614%
15	15.357	1158	1161	1167	rVB	819088	1144030	21.82%	2.770%
16	16.214	1234	1236	1240	rBV	28941	54372	1.04%	0.132%

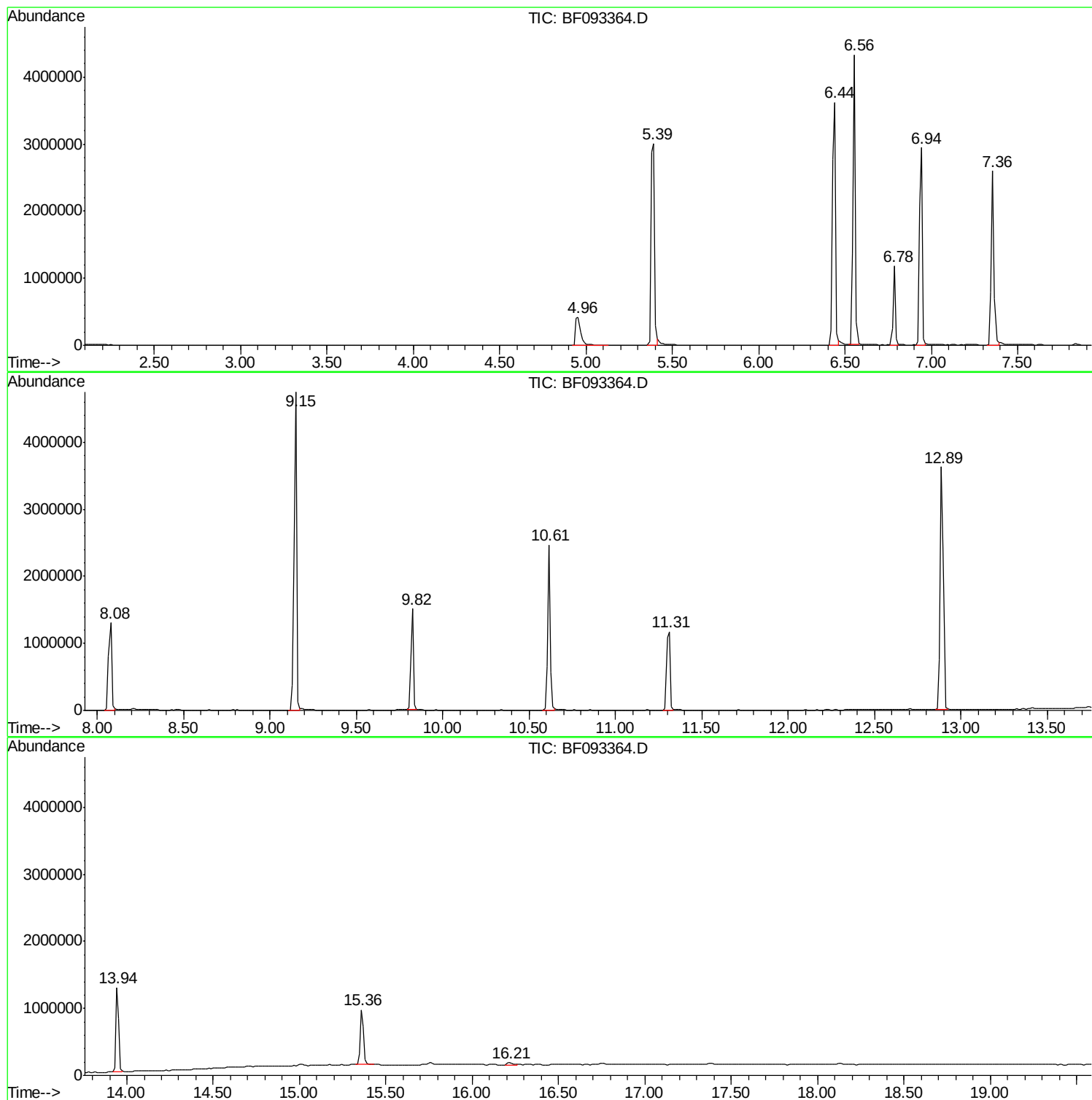
Sum of corrected areas: 41306767

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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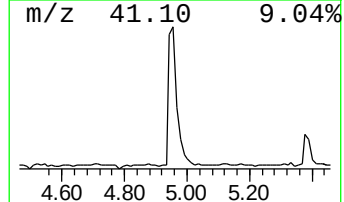
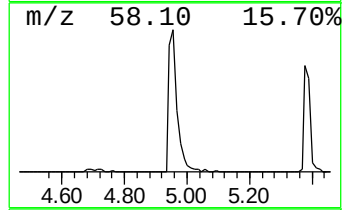
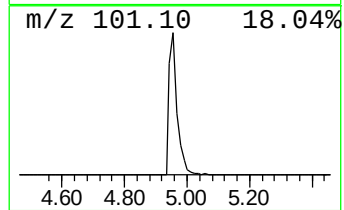
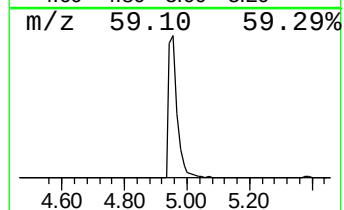
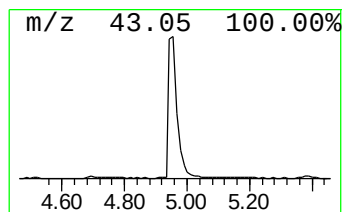
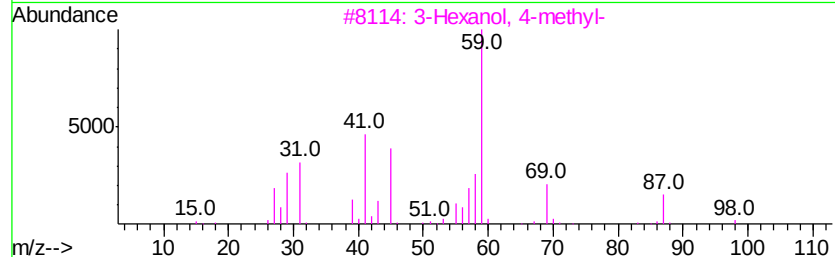
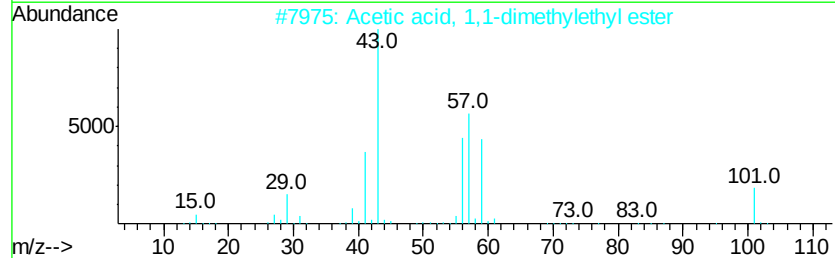
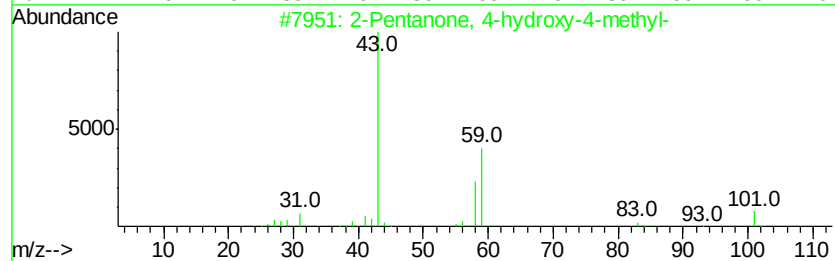
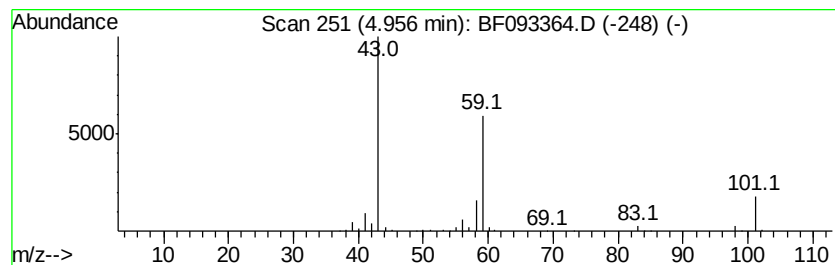
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	15.33 ng	805444	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
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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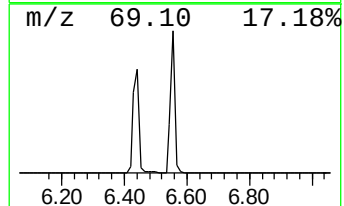
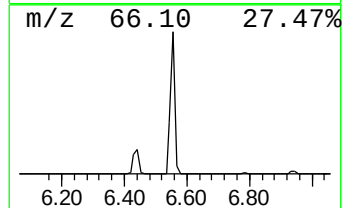
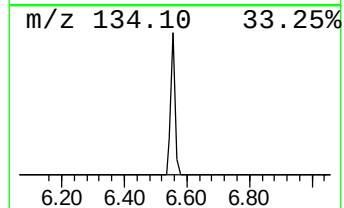
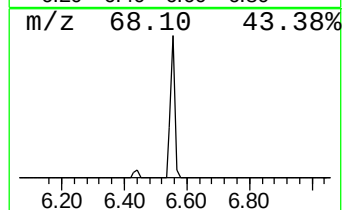
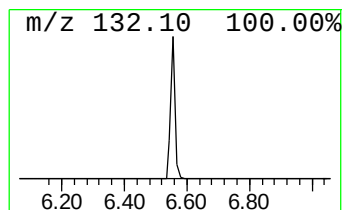
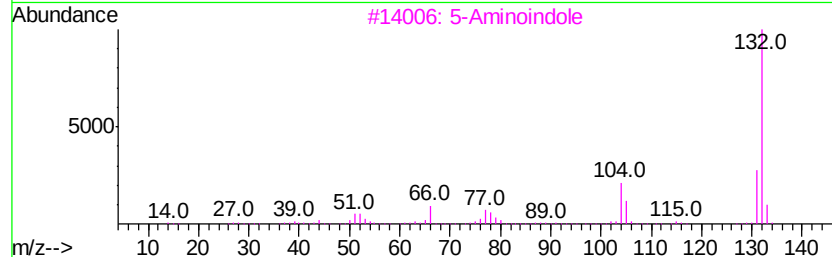
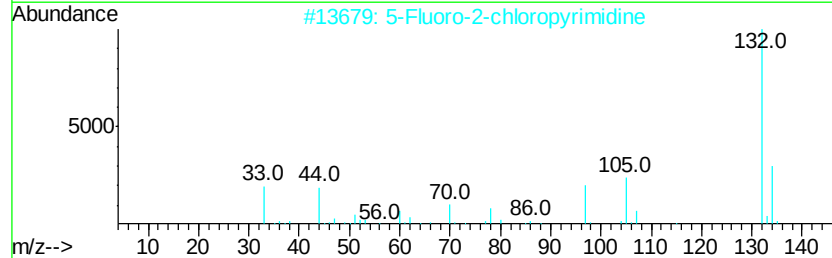
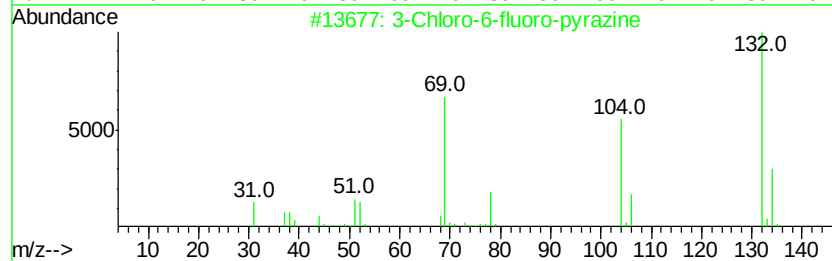
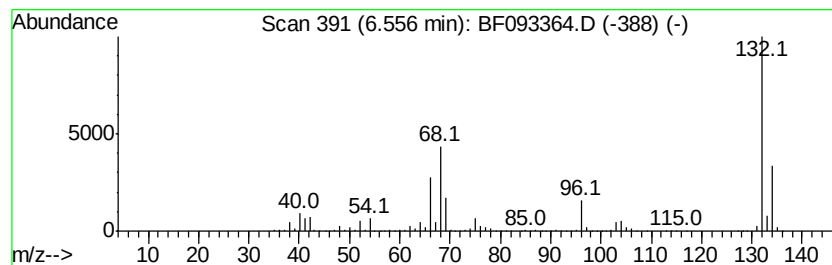
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	78.61 ng	4130980	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	15.3	ng	805444	1	6.78	1051010	20.0
unknown6.56	6.56	78.6	ng	4130980	1	6.78	1051010	20.0