

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020617\
 Data File : BF092800.D
 Acq On : 6 Feb 2017 17:20
 Operator : SJ/MA
 Sample : PB96743BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96743BL

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-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	5.104	261	264	273	rBV	986212	1372560	31.39%	3.859%
2	5.527	298	301	304	rBV	3264906	3355644	76.75%	9.433%
3	6.556	388	391	393	rBV	3393153	3550470	81.20%	9.981%
4	6.682	399	402	405	rBV	3304097	3350559	76.63%	9.419%
5	6.910	420	422	425	rBV	1085993	905680	20.71%	2.546%
6	7.070	433	436	438	rBV	3100389	2875922	65.78%	8.085%
7	7.482	469	472	474	rBV	1696580	2392800	54.73%	6.727%
8	8.190	532	534	537	rBV	901505	1166599	26.68%	3.280%
9	9.276	625	629	631	rBV	3526985	4372278	100.00%	12.291%
10	9.859	674	680	685	rBV	47594	106182	2.43%	0.299%
11	9.951	685	688	690	rVB	1554863	1415924	32.38%	3.980%
12	10.739	754	757	760	rBV	2098688	2333449	53.37%	6.560%
13	11.436	815	818	820	rBV	1617222	1435181	32.82%	4.035%
14	13.025	954	957	960	rBV	3856659	4108559	93.97%	11.550%
15	14.077	1046	1049	1051	rBV	1608893	1491581	34.11%	4.193%
16	15.528	1173	1176	1179	rBV	1074462	1338291	30.61%	3.762%

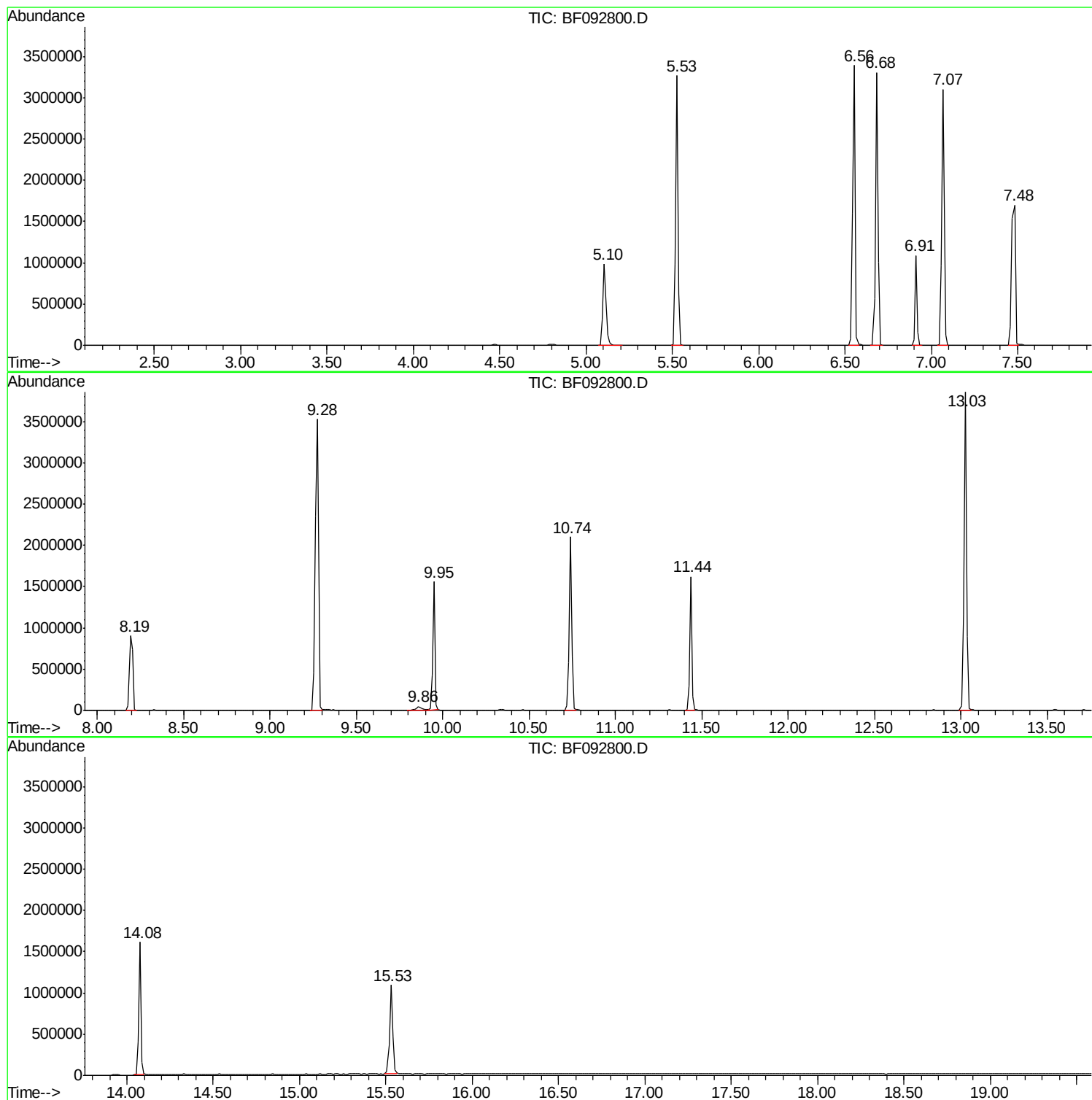
Sum of corrected areas: 35571679

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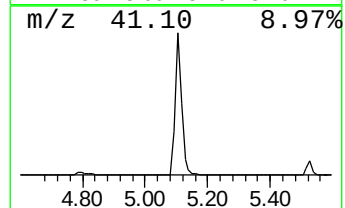
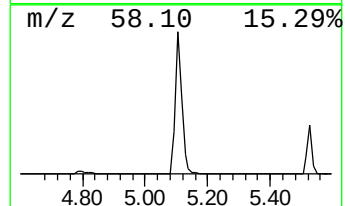
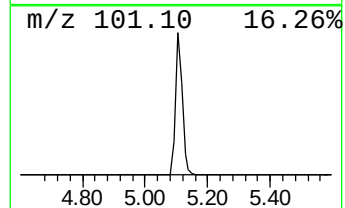
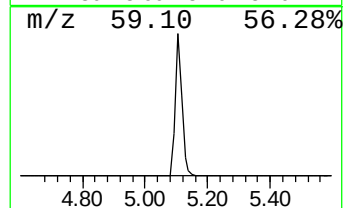
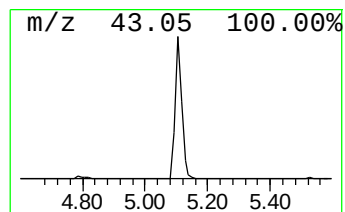
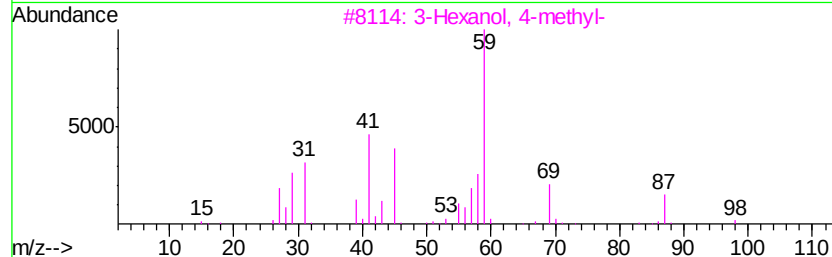
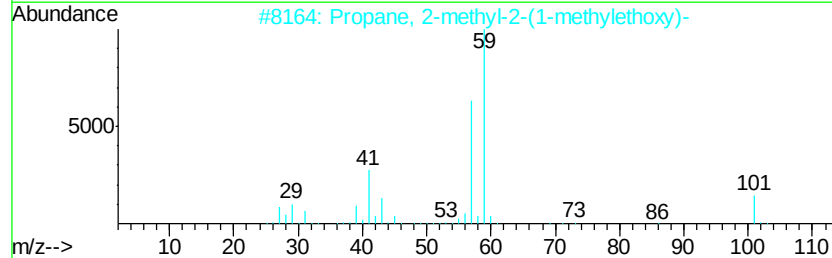
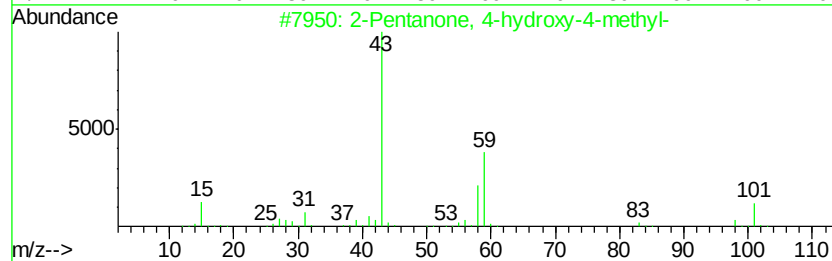
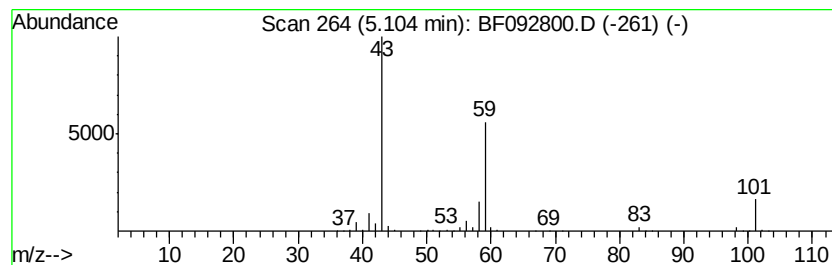
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	30.31 ng	1372560	1,4-Dichlorobenzene-d4	6.91

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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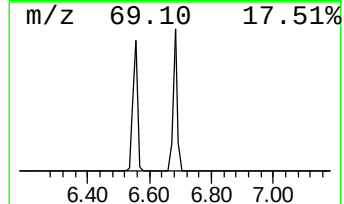
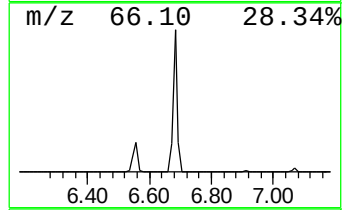
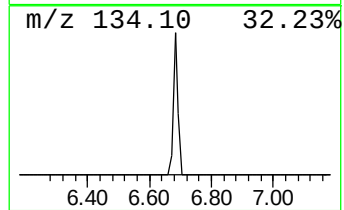
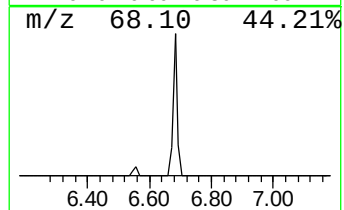
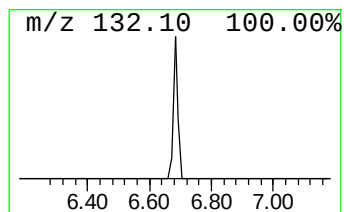
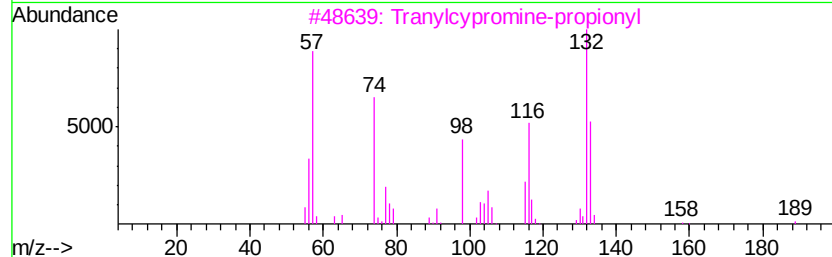
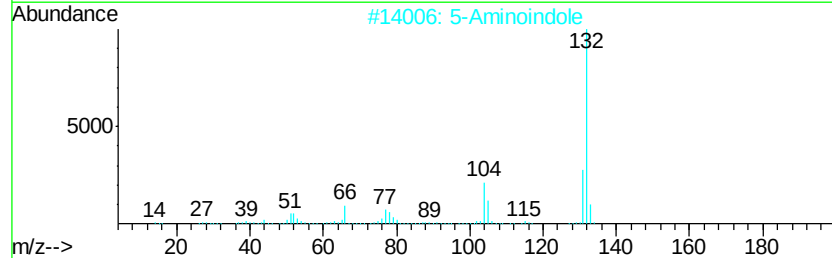
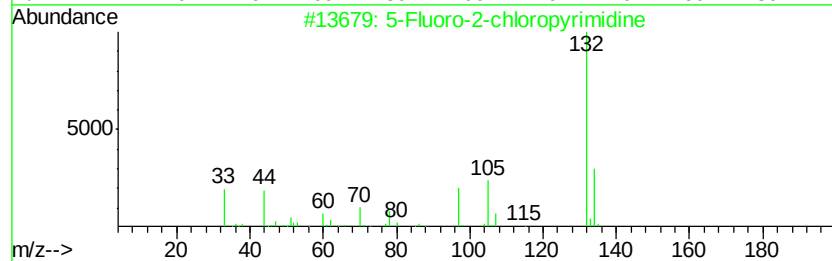
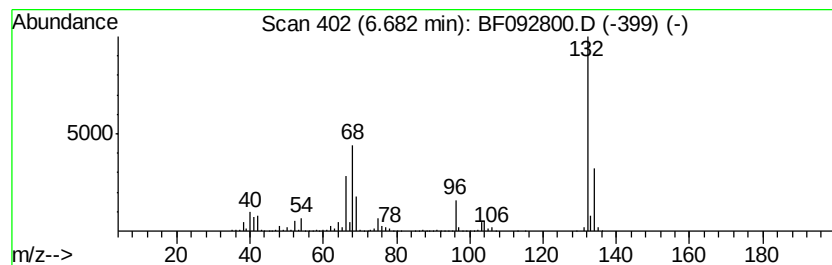
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Peak Number 2 unknown6.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.68	73.99 ng	3350560	1,4-Dichlorobenzene-d4	6.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	30.3	ng	1372560	1	6.91	905680	20.0
unknown6.68	6.68	74.0	ng	3350560	1	6.91	905680	20.0