

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011817\
 Data File : BF092353.D
 Acq On : 18 Jan 2017 16:10
 Operator : UM/SJ
 Sample : PB96092BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96092BL

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-BF122116.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	4.656	248	251	264	rBV	1056049	2029992	35.12%	4.694%
2	5.136	290	293	295	rBV	3988947	4676308	80.91%	10.814%
3	6.210	384	387	389	rBV	4167387	4383169	75.84%	10.136%
4	6.313	394	396	399	rVB	4619059	4245337	73.45%	9.817%
5	6.542	414	416	419	rBV	1054173	915671	15.84%	2.117%
6	6.702	427	430	432	rBV	2710929	3505610	60.65%	8.106%
7	7.113	463	466	469	rBV	2486887	2649049	45.83%	6.126%
8	7.833	526	529	531	rBV	1029400	1207243	20.89%	2.792%
9	8.907	620	623	626	rBV	5102614	5779658	100.00%	13.365%
10	9.570	679	681	684	rBV	1433440	1680768	29.08%	3.887%
11	10.370	748	751	753	rBV2	2448646	3240805	56.07%	7.494%
12	11.056	808	811	813	rBV	1359630	1619586	28.02%	3.745%
13	12.645	947	950	952	rBV	3648186	5116688	88.53%	11.832%
14	13.674	1038	1040	1043	rBV	830158	1114734	19.29%	2.578%
15	15.034	1156	1159	1172	rVB	805828	1080240	18.69%	2.498%

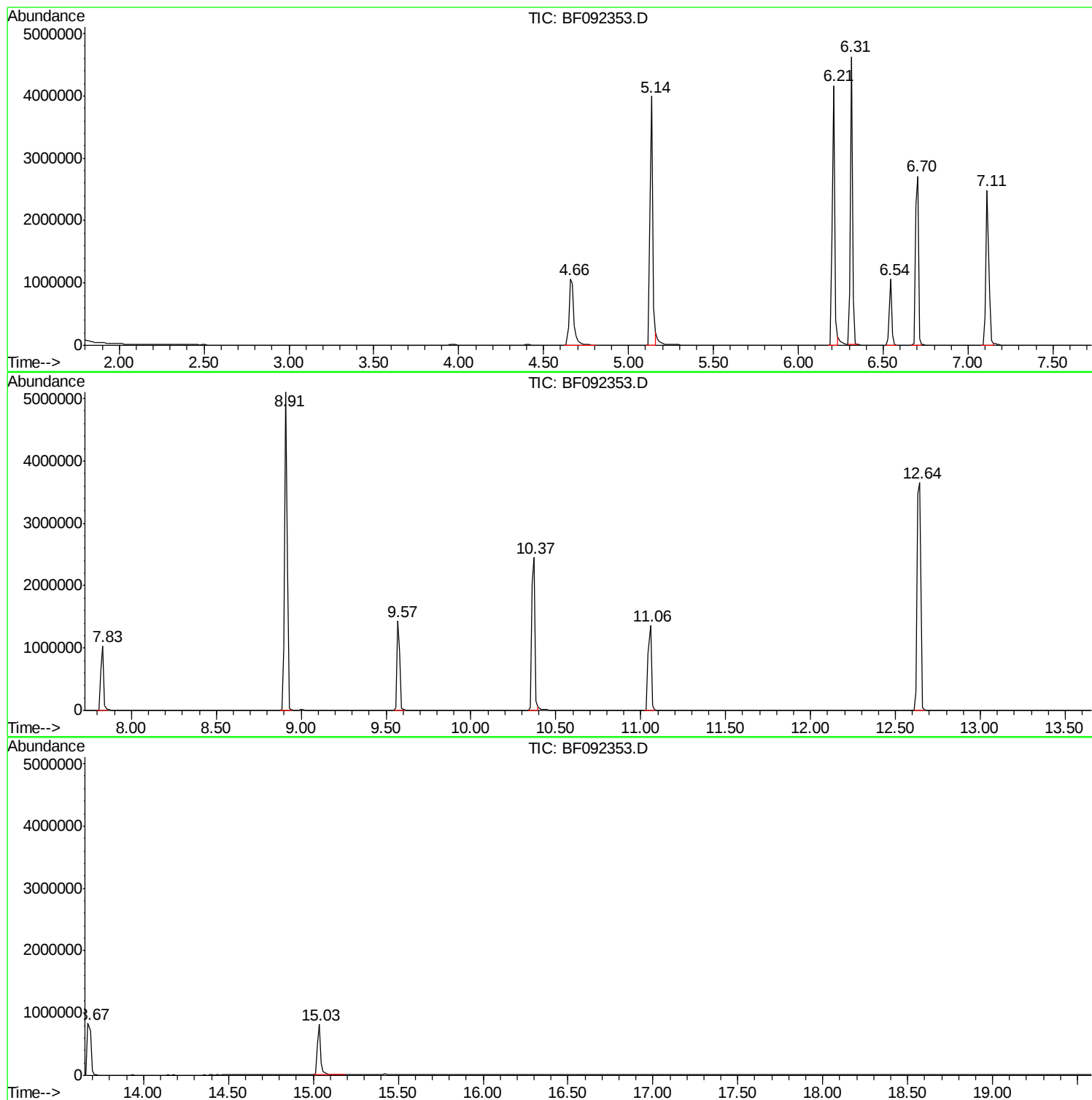
Sum of corrected areas: 43244858

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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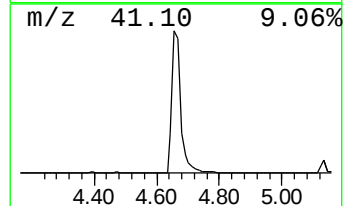
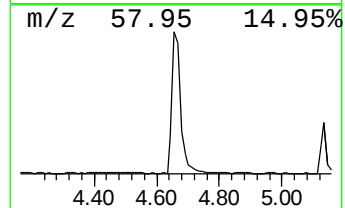
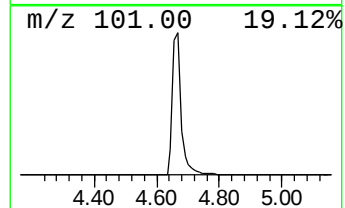
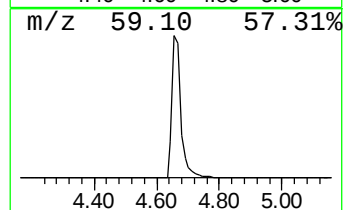
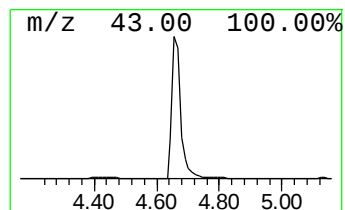
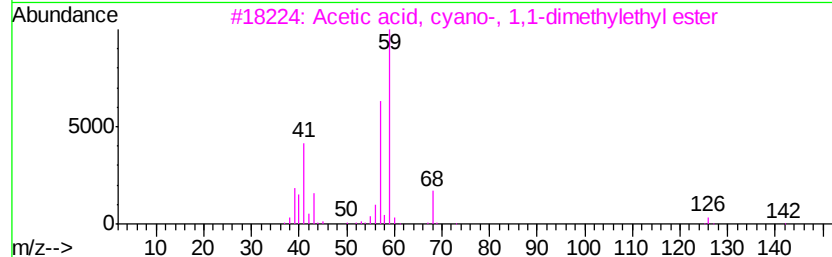
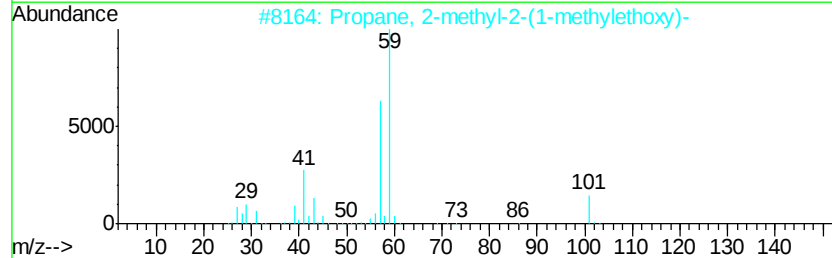
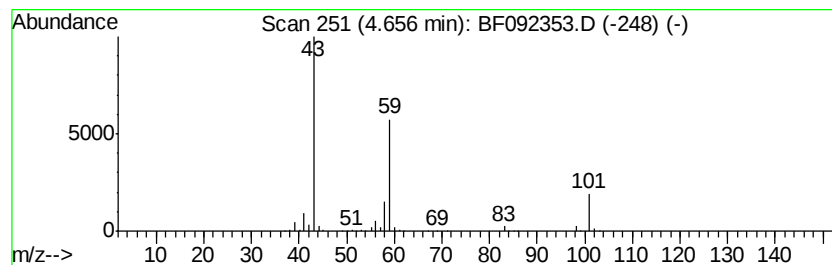
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.66	44.34 ng	2029990	1,4-Dichlorobenzene-d4	6.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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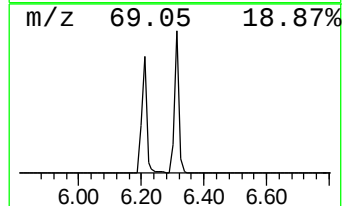
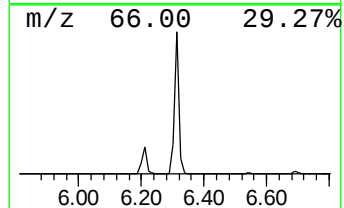
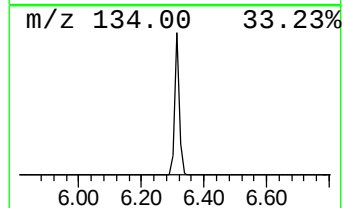
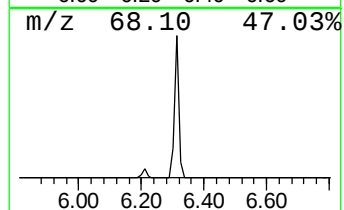
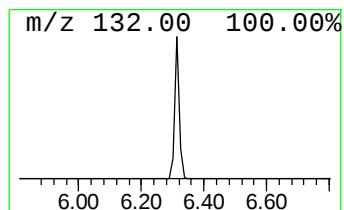
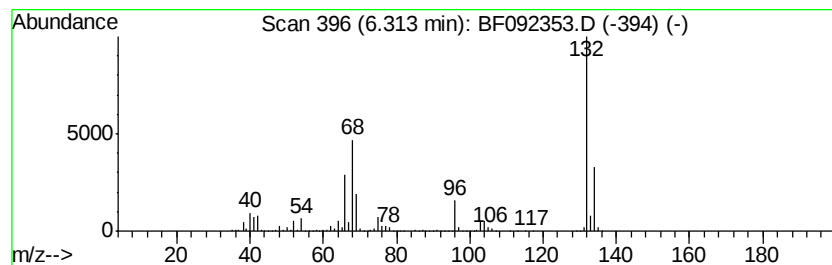
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Peak Number 2 unknown6.31 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	92.73 ng	4245340	1,4-Dichlorobenzene-d4	6.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.66	44.3	ng	2029990	1	6.54	915671	20.0
unknown6.31	6.31	92.7	ng	4245340	1	6.54	915671	20.0