

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112217\
 Data File : BF100856.D
 Acq On : 23 Nov 2017 6:42
 Operator : SJ/JU
 Sample : PB104490BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB104490BL

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-BF110817.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	5.092	506	511	522	rBV	280245	376982	15.97%	2.513%
2	5.487	573	578	581	rBV	1429511	1274745	54.01%	8.498%
3	6.486	743	748	751	rBV	1273696	1284867	54.44%	8.566%
4	6.628	768	772	775	rBV	1508604	1342742	56.89%	8.952%
5	6.857	807	811	815	rBB	539647	431471	18.28%	2.876%
6	7.016	833	838	841	rBV	1741732	1557972	66.01%	10.386%
7	7.422	901	907	910	rBV	1113112	1231798	52.19%	8.212%
8	8.139	1024	1029	1033	rBV	650892	565857	23.97%	3.772%
9	9.216	1206	1212	1215	rBV	2245530	2360291	100.00%	15.735%
10	9.898	1323	1328	1331	rBV	707407	613356	25.99%	4.089%
11	10.686	1457	1462	1465	rBV	940682	852230	36.11%	5.681%
12	11.386	1576	1581	1584	rBV	610862	566827	24.02%	3.779%
13	12.968	1846	1850	1854	rBV	1799279	1729723	73.28%	11.531%
14	14.021	2026	2029	2033	rBV	469113	387382	16.41%	2.583%
15	15.492	2274	2279	2286	rBV	328132	423913	17.96%	2.826%

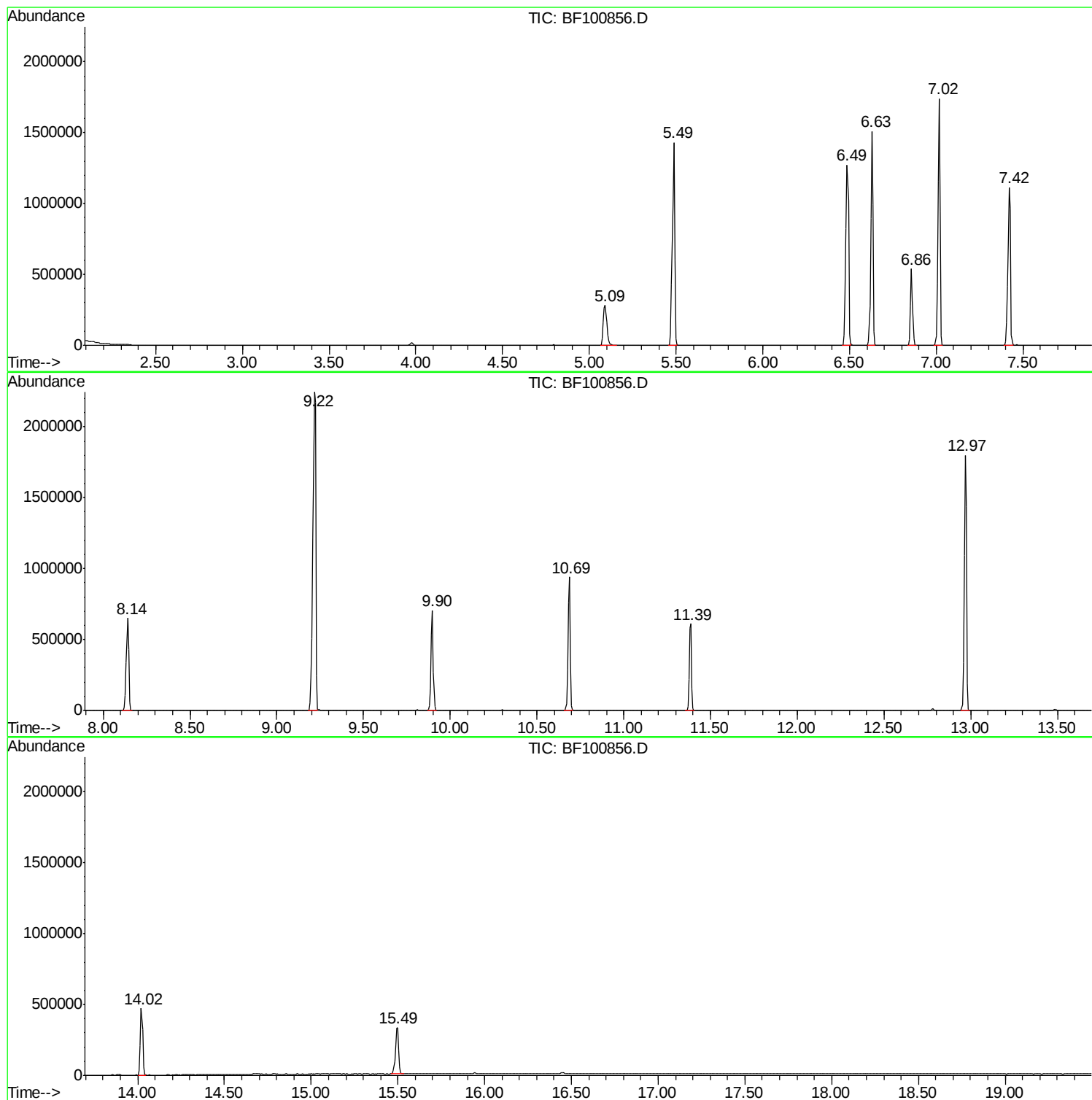
Sum of corrected areas: 15000156

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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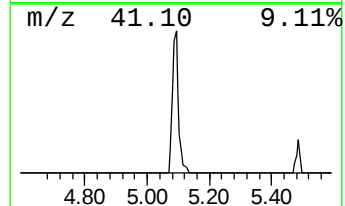
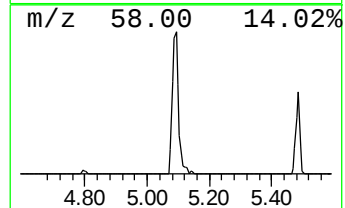
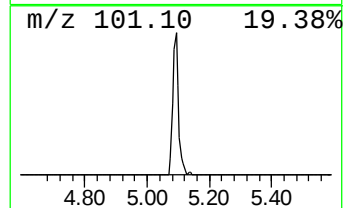
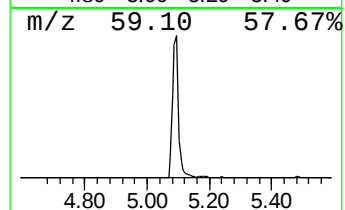
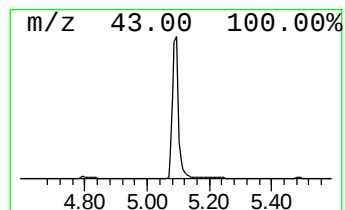
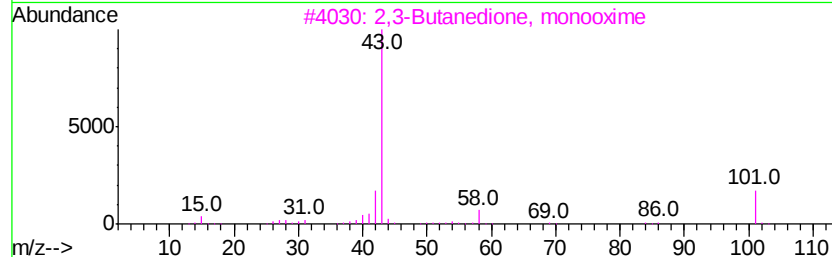
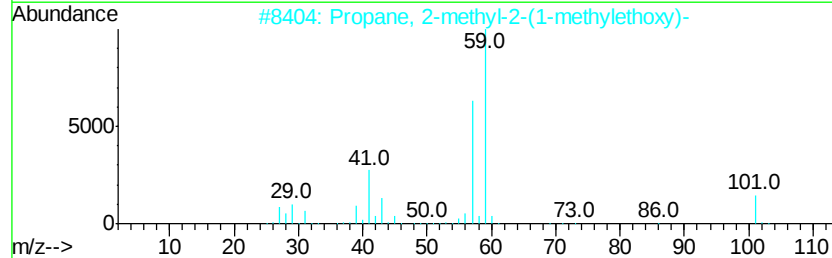
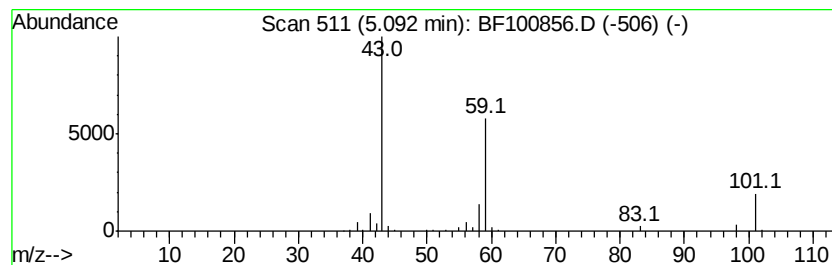
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	17.47 ng	376982	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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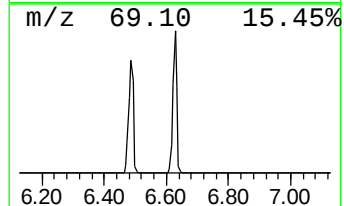
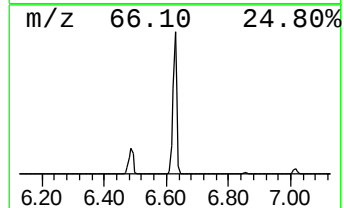
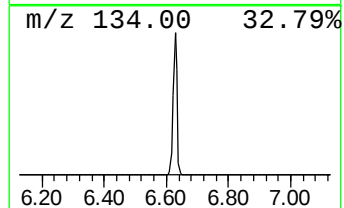
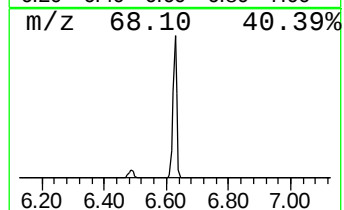
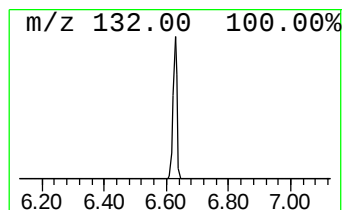
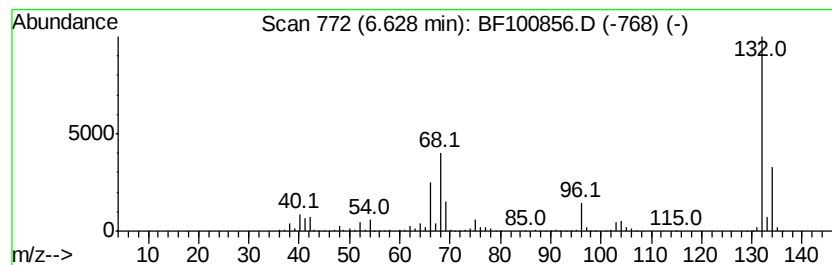
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Peak Number 2 unknown6.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	62.24 ng	1342740	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	17.5	ng	376982	1	6.86	431471	20.0
unknown6.63	6.63	62.2	ng	1342740	1	6.86	431471	20.0