

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112717\
 Data File : BF100917.D
 Acq On : 28 Nov 2017 3:17
 Operator : SJ/JU
 Sample : I6584-06
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 112117-TIDO-F-D-S

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	5.069	504	507	517	rVB	100941	99433	4.85%	0.874%
2	5.469	572	575	583	rBV	515283	455460	22.21%	4.003%
3	6.475	743	746	755	rVB	297729	305539	14.90%	2.686%
4	6.622	767	771	775	rBV	1023674	920821	44.90%	8.093%
5	6.857	807	811	814	rBV	458424	376093	18.34%	3.306%
6	7.016	833	838	841	rBV	1443072	1339346	65.31%	11.772%
7	7.422	901	907	910	rBV	1106129	1075819	52.46%	9.456%
8	8.139	1024	1029	1032	rBV	575219	490084	23.90%	4.308%
9	9.216	1206	1212	1215	rBV	2287893	2050880	100.00%	18.026%
10	9.892	1322	1327	1331	rBV2	603432	554529	27.04%	4.874%
11	10.680	1457	1461	1472	rVB	795713	732857	35.73%	6.441%
12	11.380	1576	1580	1584	rBV	617897	512179	24.97%	4.502%
13	12.968	1845	1850	1853	rBV	1795021	1719068	83.82%	15.110%
14	13.857	1998	2001	2009	rBV2	23117	33031	1.61%	0.290%
15	14.021	2025	2029	2036	rVB	398905	375513	18.31%	3.301%
16	15.486	2273	2278	2288	rBV	265249	336677	16.42%	2.959%

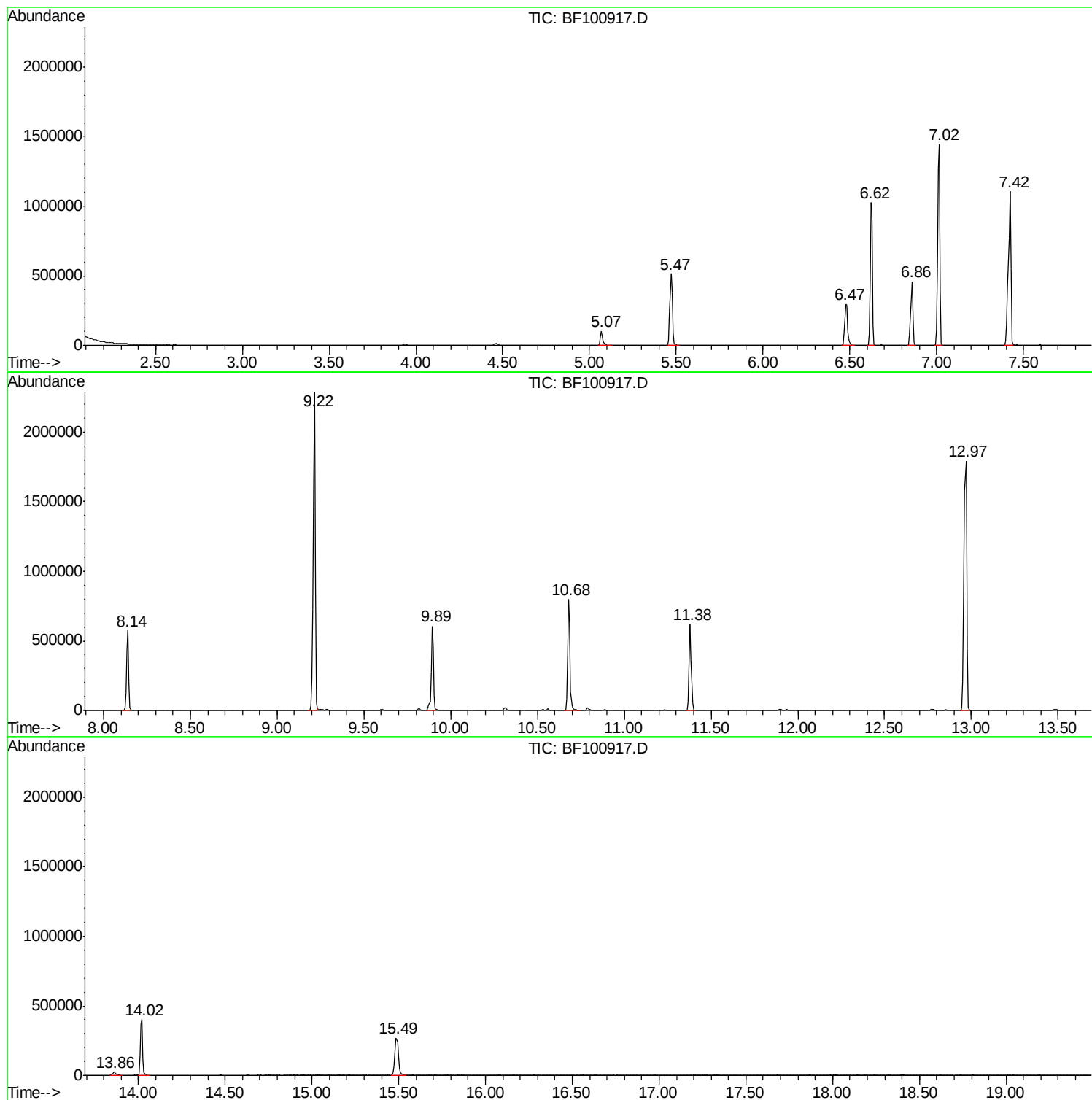
Sum of corrected areas: 11377329

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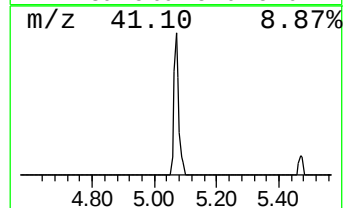
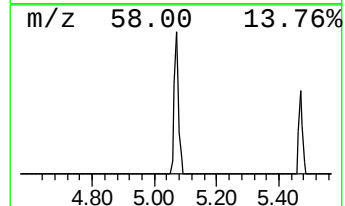
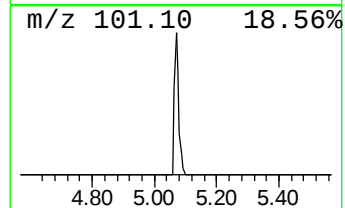
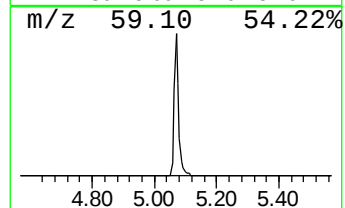
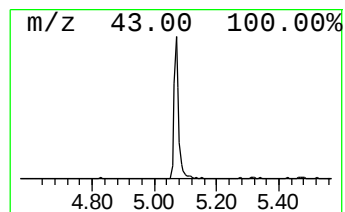
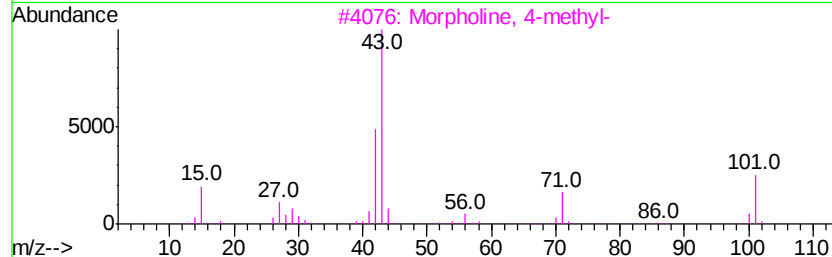
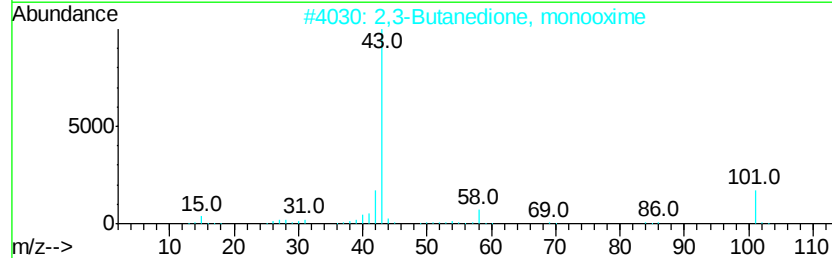
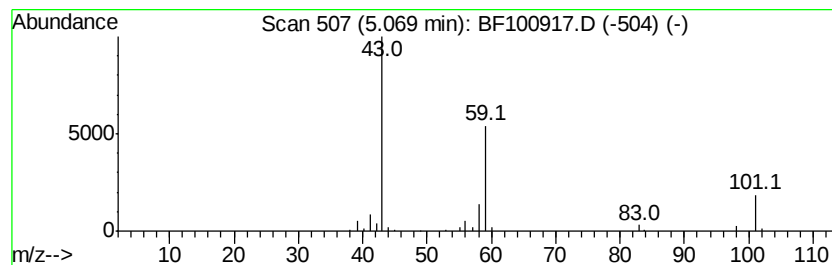
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	5.29 ng	99433	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9



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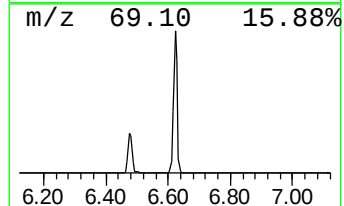
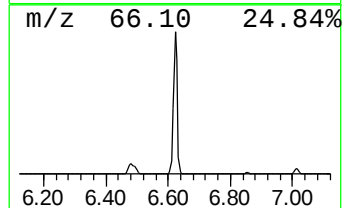
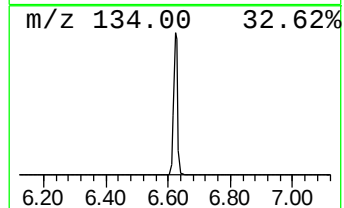
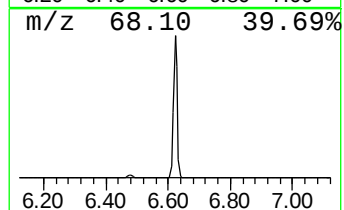
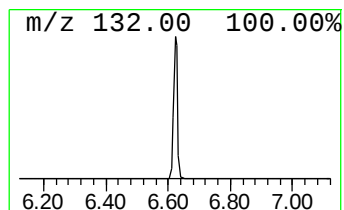
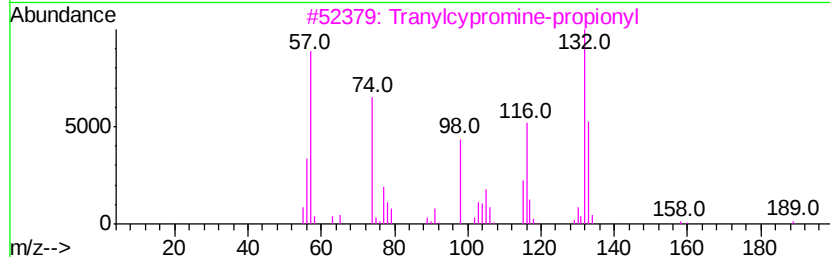
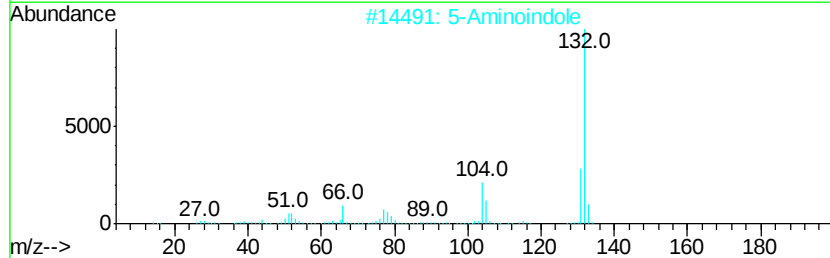
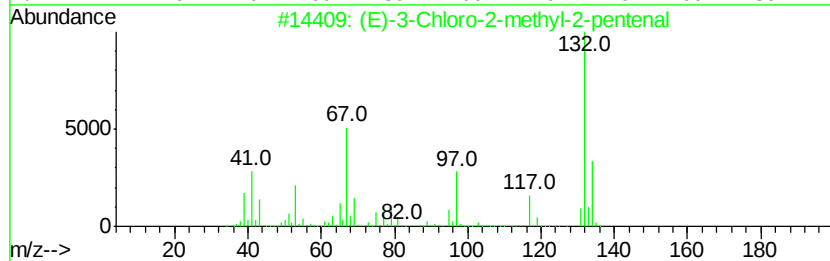
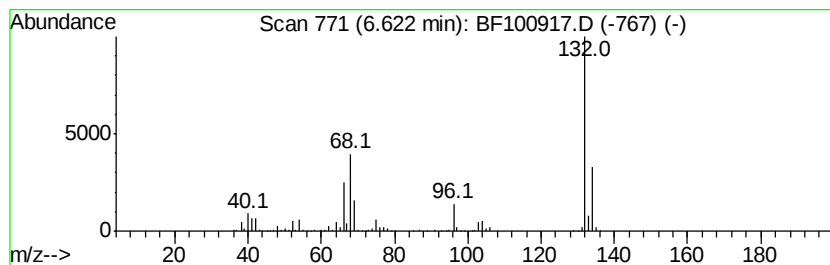
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Peak Number 2 unknown6.62 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	48.97 ng	920821	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	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...	5.07	5.3	ng	99433	1	6.86	376093	20.0
unknown6.62	6.62	49.0	ng	920821	1	6.86	376093	20.0