

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112717\
 Data File : BF100916.D
 Acq On : 28 Nov 2017 2:50
 Operator : SJ/JU
 Sample : I6584-05
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

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

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	514	rBV	104491	101791	4.82%	0.894%
2	5.469	572	575	586	rBV	577627	511006	24.18%	4.490%
3	6.481	743	747	755	rVB	329240	327355	15.49%	2.876%
4	6.622	767	771	775	rBV	1066741	993449	47.01%	8.728%
5	6.857	807	811	814	rBV	415361	340918	16.13%	2.995%
6	7.016	833	838	841	rBV	1507055	1359522	64.33%	11.945%
7	7.422	901	907	910	rBV	1124196	1136880	53.80%	9.989%
8	8.139	1024	1029	1032	rBV	524971	439040	20.78%	3.857%
9	9.216	1206	1212	1215	rBV	2373524	2113265	100.00%	18.567%
10	9.892	1321	1327	1335	rVB2	558984	497176	23.53%	4.368%
11	10.680	1457	1461	1470	rVB	839678	770705	36.47%	6.771%
12	11.380	1576	1580	1583	rBV	569998	444465	21.03%	3.905%
13	12.969	1845	1850	1853	rBV	1909439	1734763	82.09%	15.242%
14	14.016	2025	2028	2034	rVB	336634	319607	15.12%	2.808%
15	15.492	2273	2279	2288	rBV	222593	291807	13.81%	2.564%

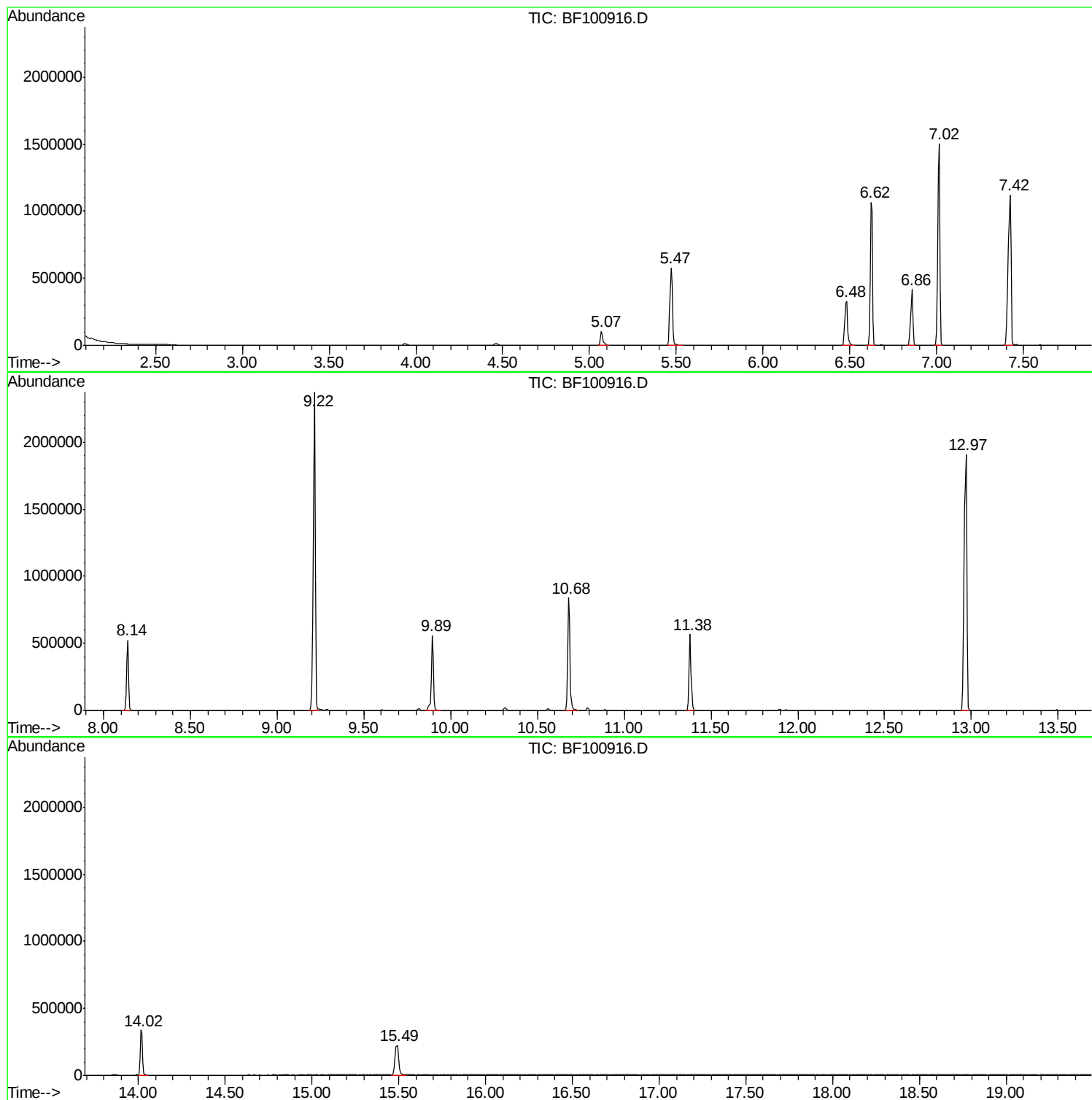
Sum of corrected areas: 11381749

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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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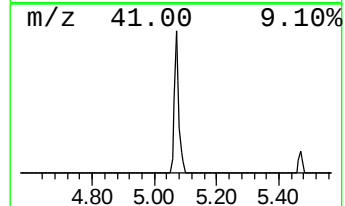
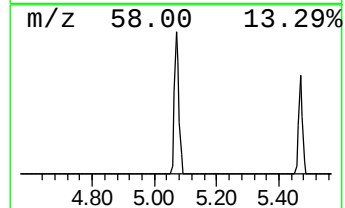
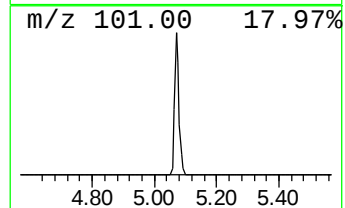
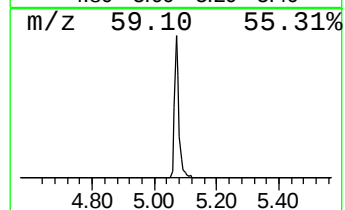
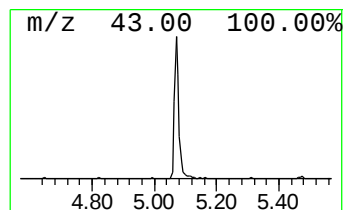
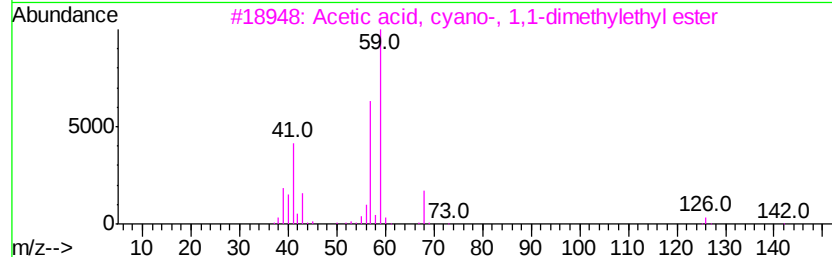
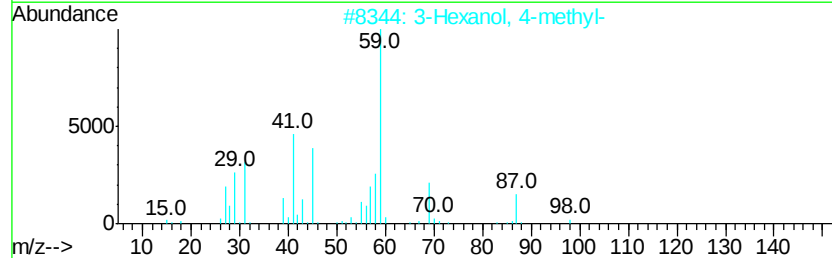
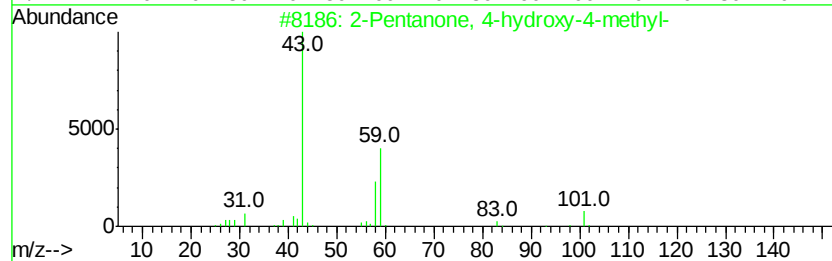
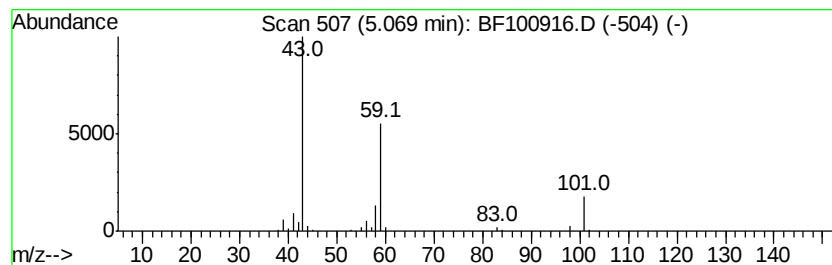
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	5.97 ng	101791	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	56
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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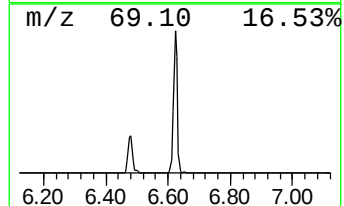
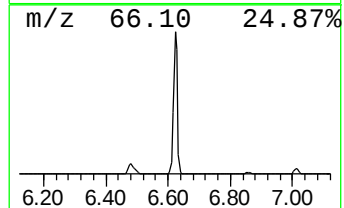
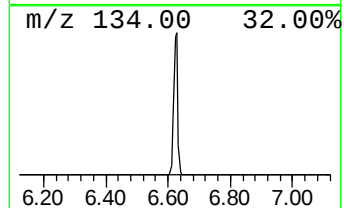
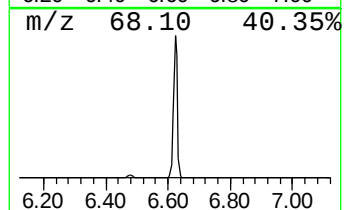
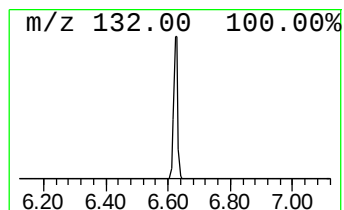
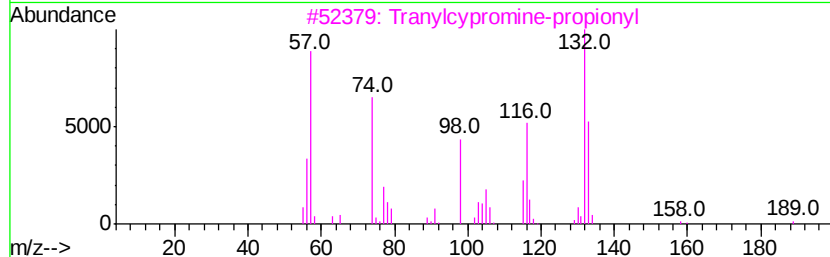
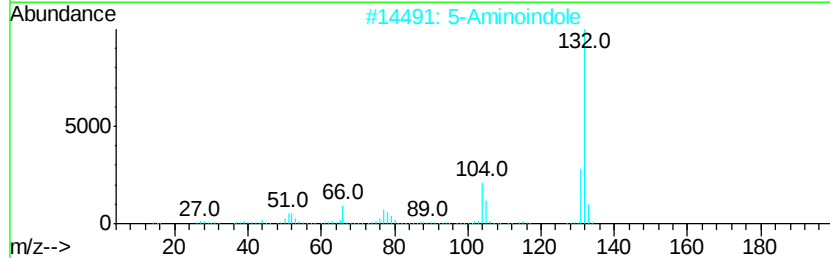
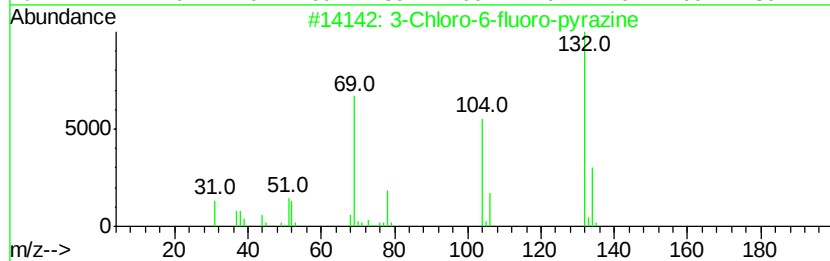
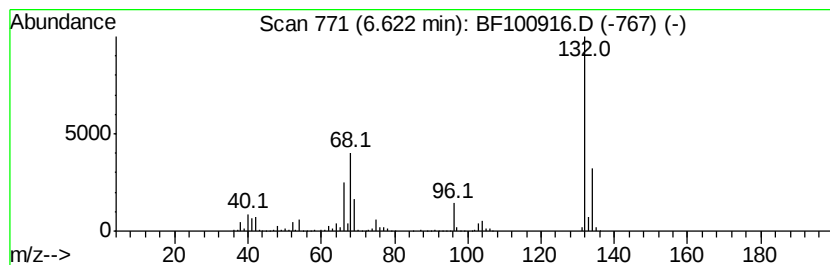
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Peak Number 2 unknown6.62 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	58.28 ng	993449	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Aminoindole	132	C8H8N2	005192-03-0	12	
3	Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10	
4	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9	
5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	6.0	ng	101791	1	6.86	340918	20.0
unknown6.62	6.62	58.3	ng	993449	1	6.86	340918	20.0