

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050818\
 Data File : BF105174.D
 Acq On : 9 May 2018 5:10
 Operator : JU/SJ
 Sample : PB108327BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB108327BL

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-BF050818.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	5.028	494	500	509	rBV	201653	308553	4.71%	0.590%
2	5.410	559	565	568	rBV	4938034	5684360	86.74%	10.863%
3	6.422	730	737	740	rBV	4798681	5555951	84.79%	10.617%
4	6.563	755	761	764	rBV	5072702	5515898	84.17%	10.541%
5	6.792	796	800	803	rBV	1494787	1267737	19.35%	2.423%
6	6.945	822	826	830	rVB	4254595	4371010	66.70%	8.353%
7	7.357	890	896	899	rBV	3385509	3628750	55.38%	6.935%
8	8.069	1012	1017	1021	rBV	1835591	1692716	25.83%	3.235%
9	9.151	1195	1201	1204	rBV	6023252	6320951	96.46%	12.079%
10	9.827	1310	1316	1319	rBV	1867705	1853818	28.29%	3.543%
11	10.616	1443	1450	1453	rBV	3999743	4072785	62.15%	7.783%
12	11.310	1562	1568	1571	rBV	2344169	1979367	30.21%	3.783%
13	12.910	1834	1840	1843	rBV	5936041	6552964	100.00%	12.523%
14	13.986	2019	2023	2027	rBV	1850646	1900862	29.01%	3.633%
15	15.521	2279	2284	2300	rBV	1278101	1622949	24.77%	3.101%

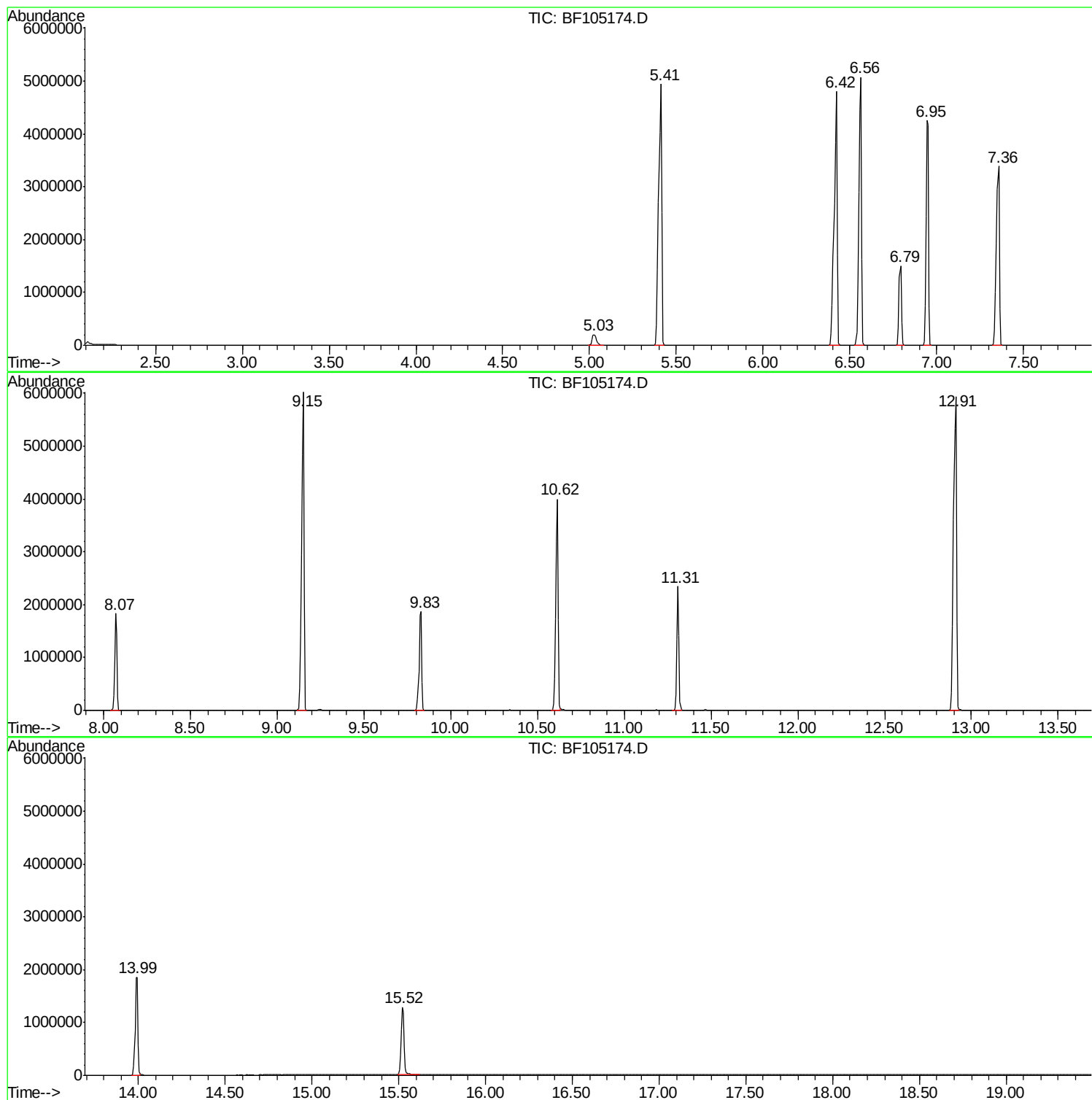
Sum of corrected areas: 52328671

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050818.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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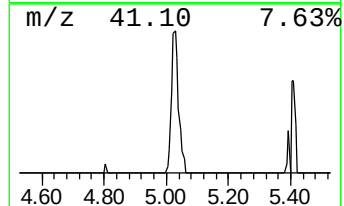
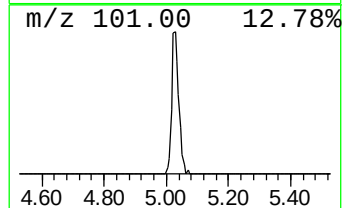
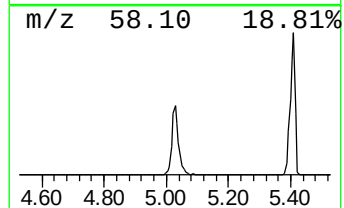
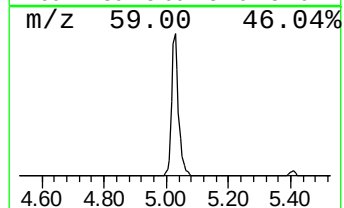
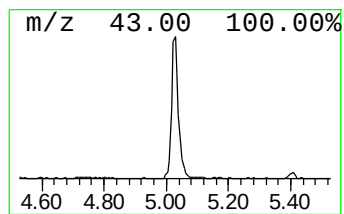
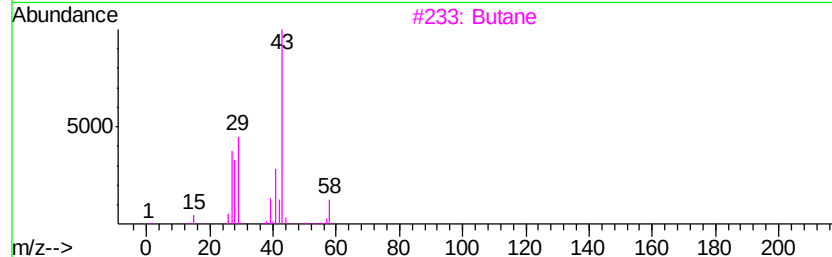
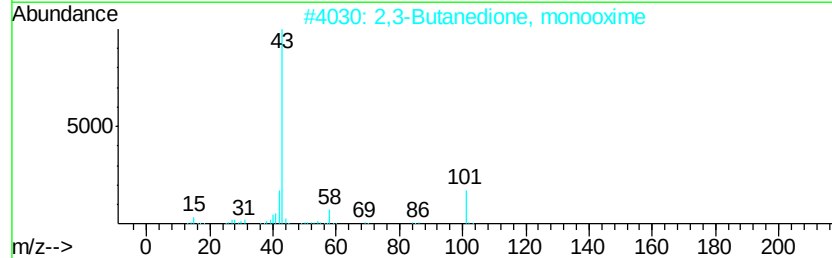
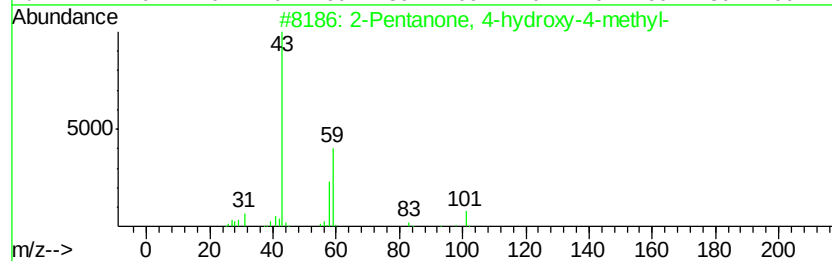
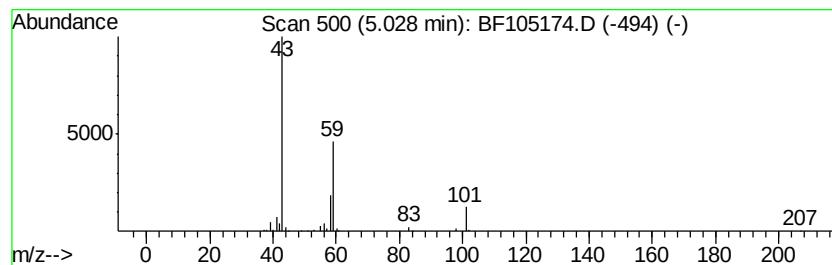
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	4.87 ng	308553	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
3		Butane	58	C4H10	000106-97-8	9
4		Diacetamide	101	C4H7NO2	000625-77-4	9
5		Acetone	58	C3H6O	000067-64-1	9



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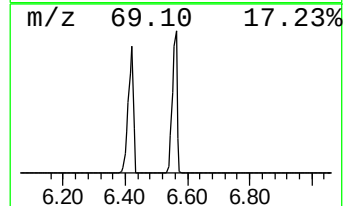
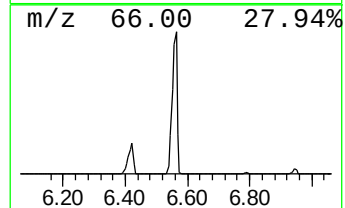
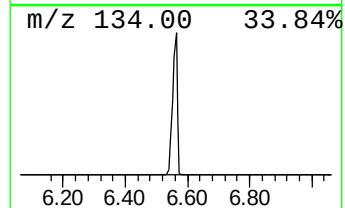
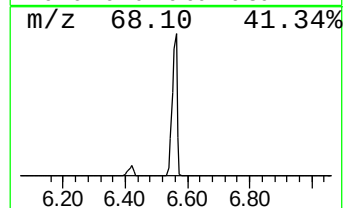
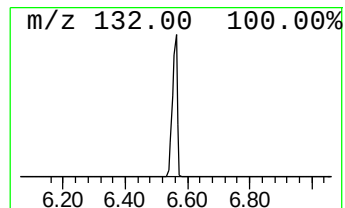
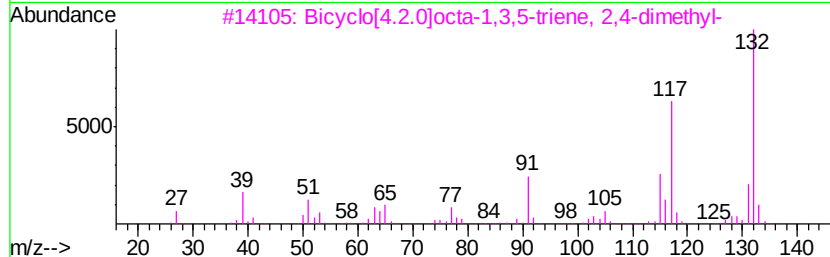
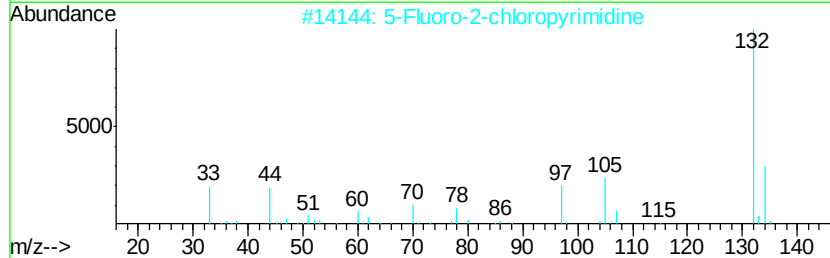
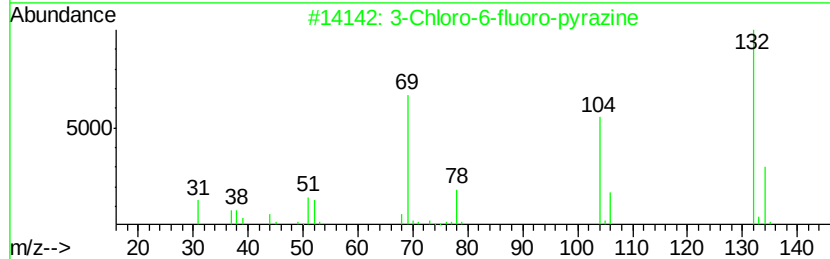
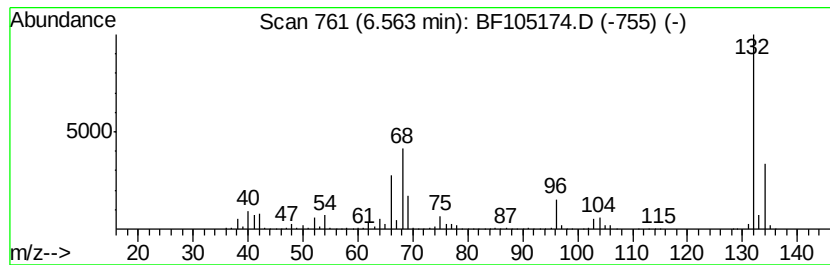
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Peak Number 2 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	87.02 ng	5515900	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
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...	5.03	4.9	ng	308553	1	6.79	1267740	20.0
unknown6.56	6.56	87.0	ng	5515900	1	6.79	1267740	20.0