

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG121417\
 Data File : BG031571.D
 Acq On : 14 Dec 2017 22:56
 Operator : SJ/JU
 Sample : I6884-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-11-20171211

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_G\METHODS\8270-BG121117.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	4.772	247	252	265	rVB	116717	213856	3.44%	0.917%
2	5.171	315	320	335	rVB	57966	95244	1.53%	0.408%
3	5.659	397	403	418	rBV	415048	744998	11.97%	3.195%
4	7.269	670	677	691	rBV	247337	474620	7.62%	2.035%
5	7.633	731	739	758	rBV	809688	1485002	23.86%	6.369%
6	8.097	811	818	830	rBV	233526	409268	6.57%	1.755%
7	8.420	865	873	886	rBV	887016	1609043	25.85%	6.901%
8	9.266	1009	1017	1036	rBV	591469	1174216	18.86%	5.036%
9	10.911	1286	1297	1318	rBV	253901	574716	9.23%	2.465%
10	13.344	1703	1711	1723	rBV	2060646	3351139	53.83%	14.372%
11	14.725	1937	1946	1966	rBV2	485931	868613	13.95%	3.725%
12	16.211	2193	2199	2220	rBV	1102713	1827308	29.35%	7.837%
13	17.468	2406	2413	2430	rBV2	753891	1241513	19.94%	5.324%
14	20.060	2847	2854	2871	rBV	4564202	6224923	100.00%	26.696%
15	21.740	3132	3140	3154	rBV	908904	1633769	26.25%	7.007%
16	25.018	3687	3698	3717	rBV2	467298	1389390	22.32%	5.959%

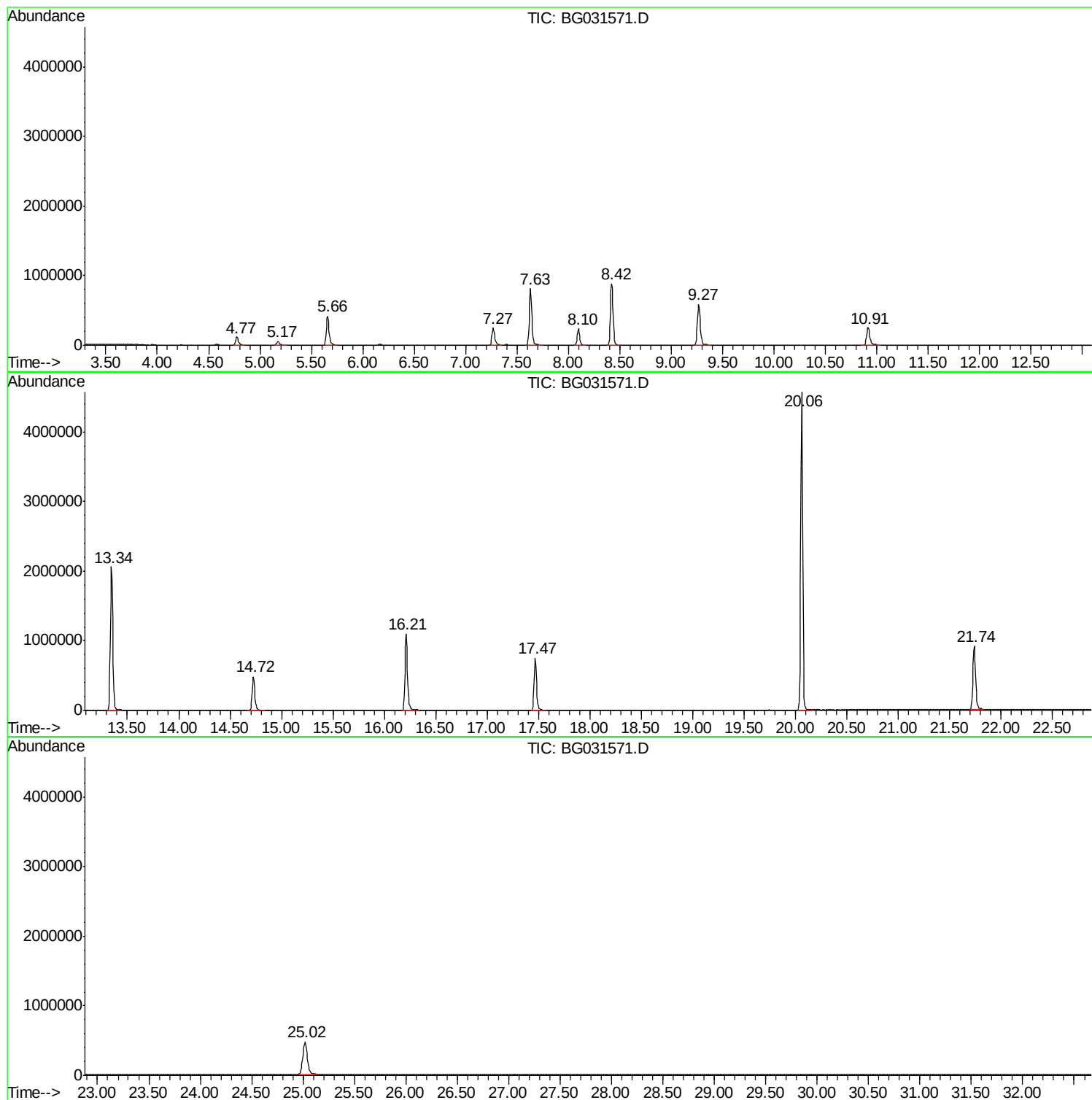
Sum of corrected areas: 23317618

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



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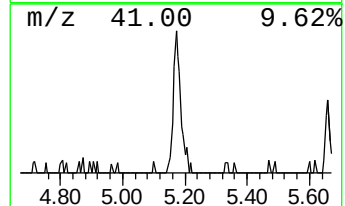
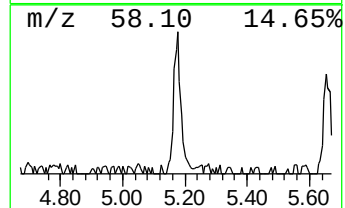
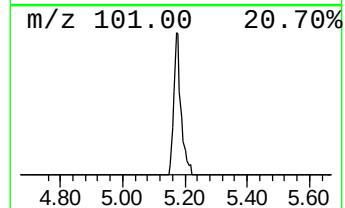
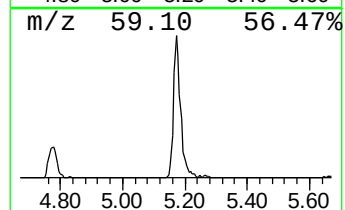
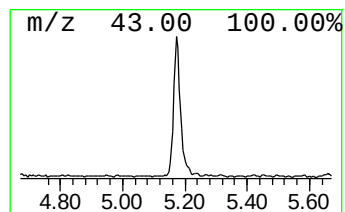
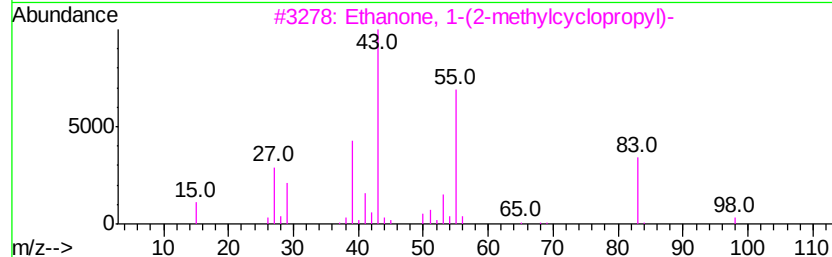
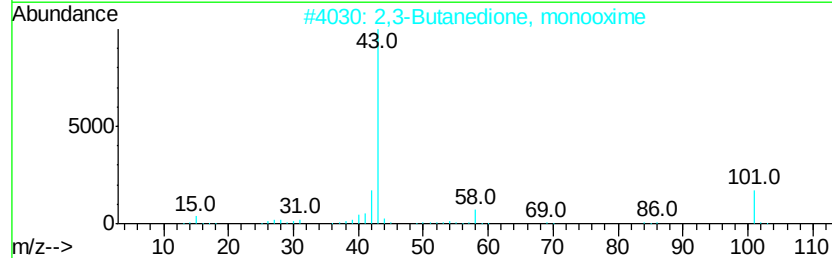
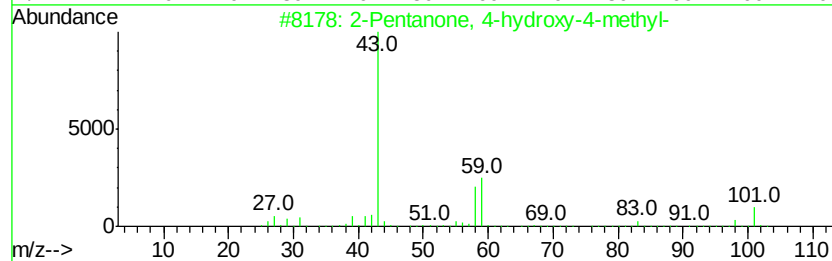
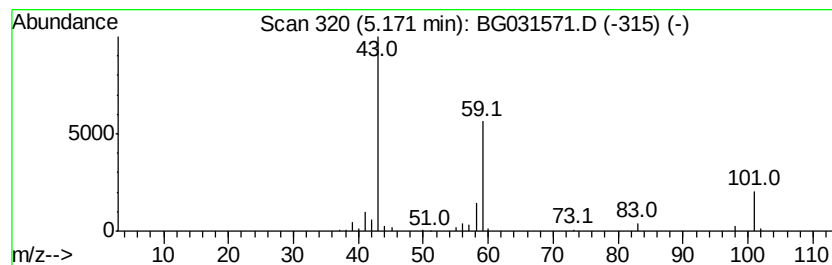
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	4.65 ng	95244	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3		Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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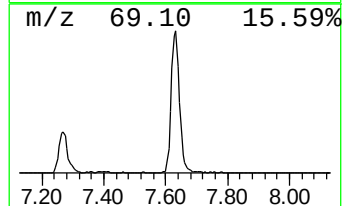
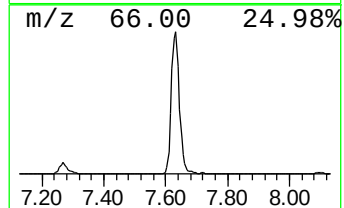
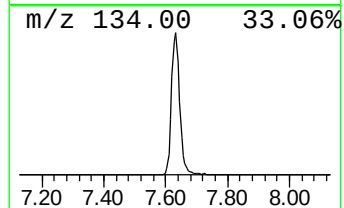
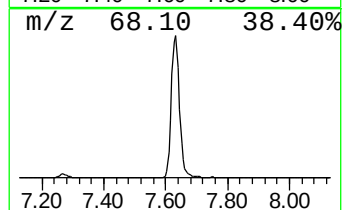
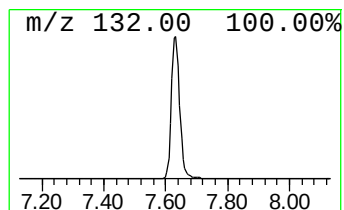
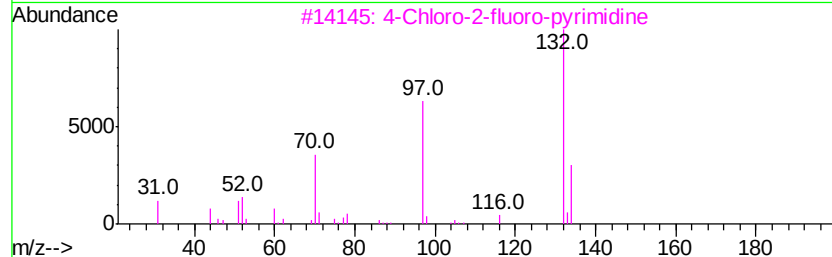
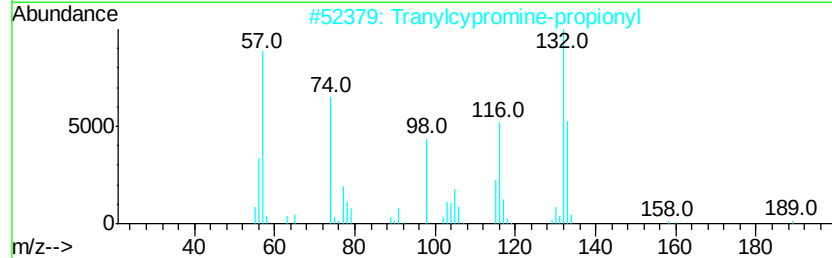
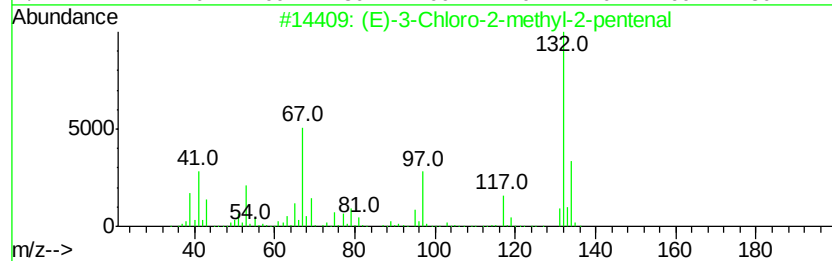
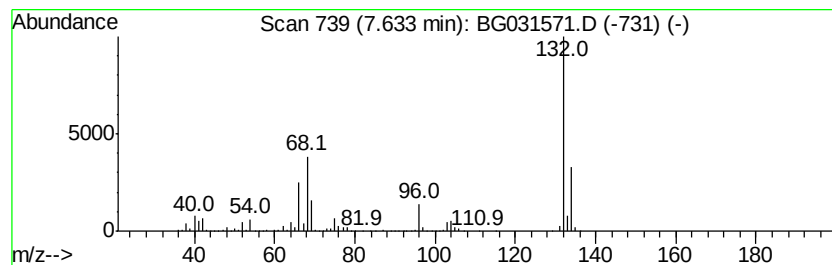
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Peak Number 3 unknown7.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	72.57 ng	1485000	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	4.7	ng	95244	1	8.10	409268	20.0
unknown7.63	7.63	72.6	ng	1485000	1	8.10	409268	20.0