

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082418\
 Data File : BG036401.D
 Acq On : 24 Aug 2018 6:38
 Operator : JU/SJ
 Sample : PB112276BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112276BL

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG082318.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	5.171	314	320	328	rBV	39117	56954	1.59%	0.350%
2	5.635	391	399	410	rBV	599016	937819	26.20%	5.770%
3	7.222	661	669	676	rBV	603485	1032994	28.86%	6.356%
4	7.598	725	733	740	rBV	571707	959608	26.81%	5.904%
5	8.068	806	813	820	rVB	186217	303843	8.49%	1.869%
6	8.391	860	868	876	rVB	624315	1071684	29.94%	6.594%
7	9.225	1000	1010	1018	rBV	515961	897984	25.09%	5.525%
8	10.876	1283	1291	1299	rVB	244469	435510	12.17%	2.679%
9	13.320	1698	1707	1715	rBV	1417189	2209426	61.73%	13.594%
10	14.689	1933	1940	1951	rBV2	380293	622519	17.39%	3.830%
11	16.176	2186	2193	2204	rBV	753224	1153164	32.22%	7.095%
12	17.433	2401	2407	2417	rVB2	579815	851465	23.79%	5.239%
13	20.042	2845	2851	2860	rBV	2843377	3579279	100.00%	22.022%
14	21.699	3126	3133	3140	rBV2	675062	1099671	30.72%	6.766%
15	24.919	3670	3681	3692	rBV2	370707	1041557	29.10%	6.408%

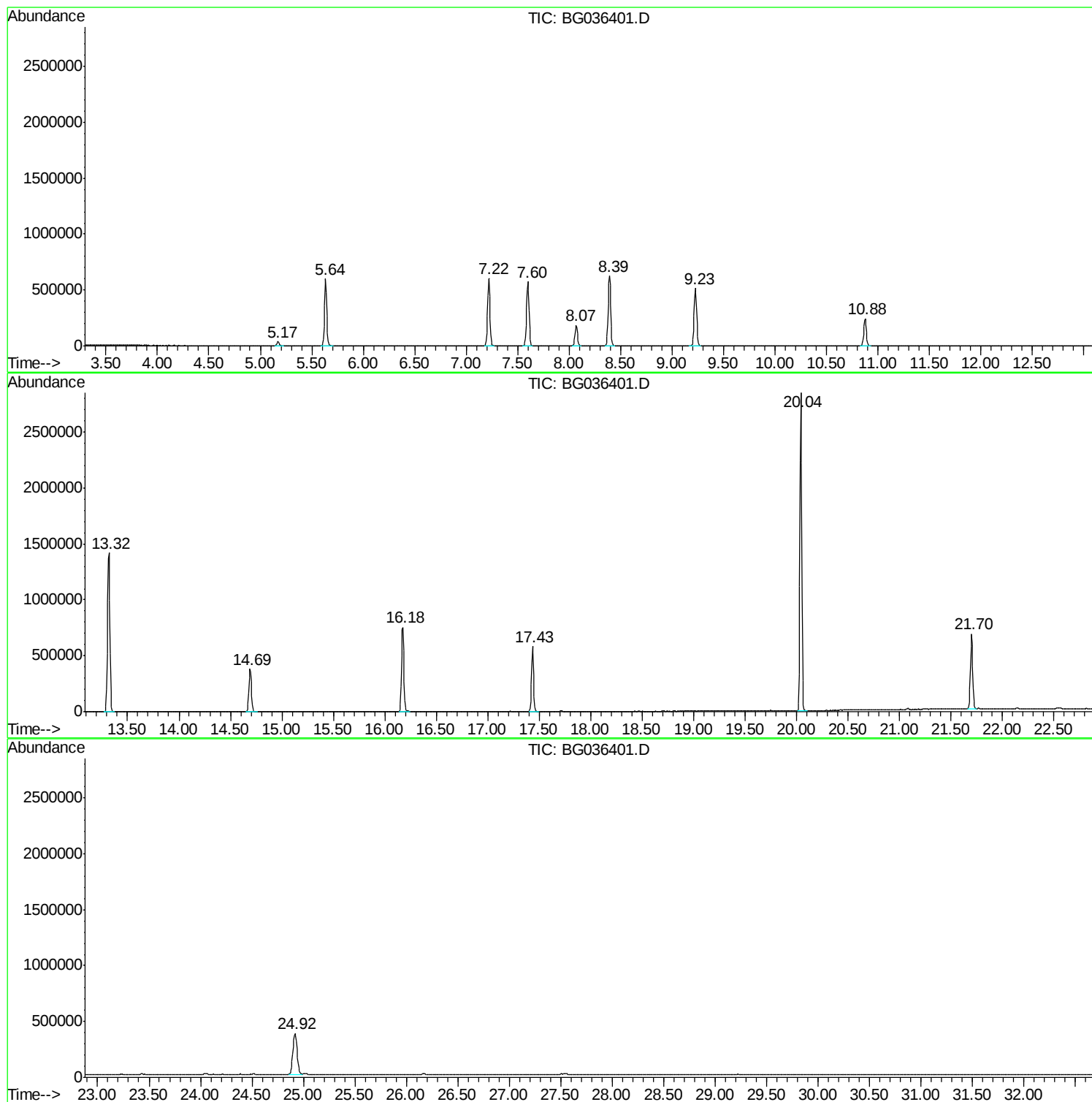
Sum of corrected areas: 16253477

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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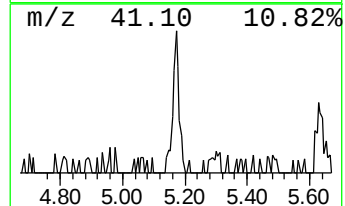
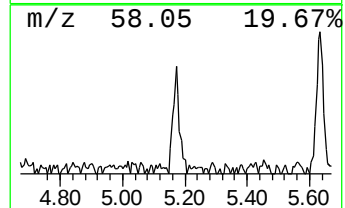
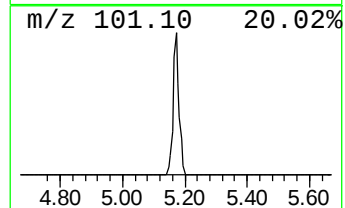
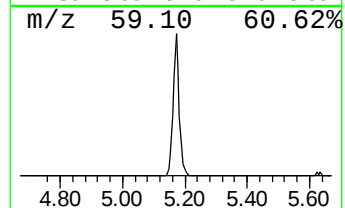
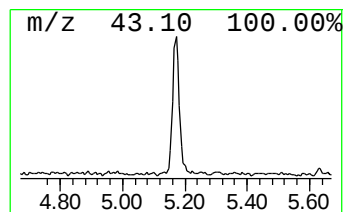
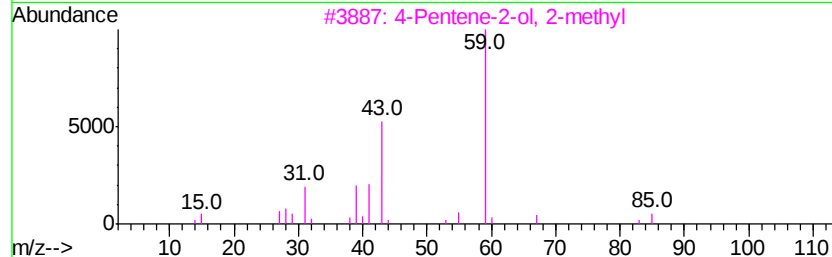
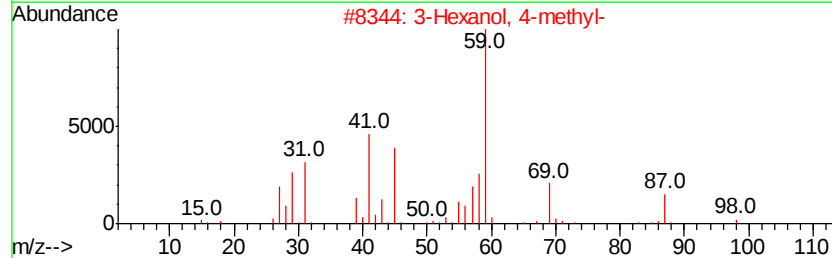
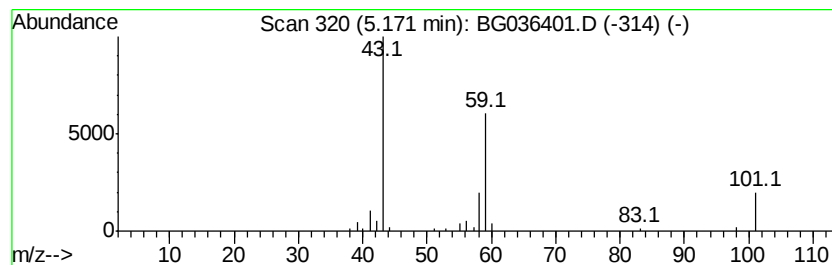
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.17	3.75 ng	56954	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			4-Pentene-2-ol, 2-methyl	100	C6H12O	000624-97-5	17
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5			Acetone	58	C3H6O	000067-64-1	9



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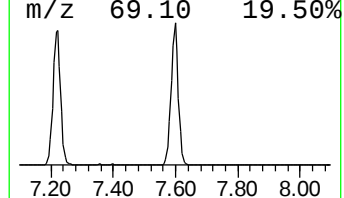
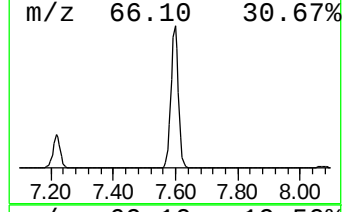
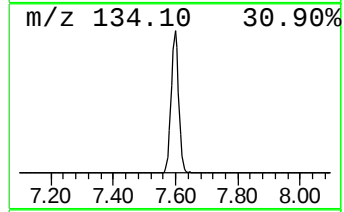
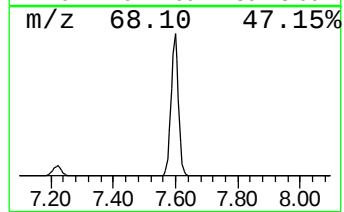
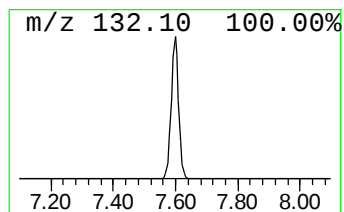
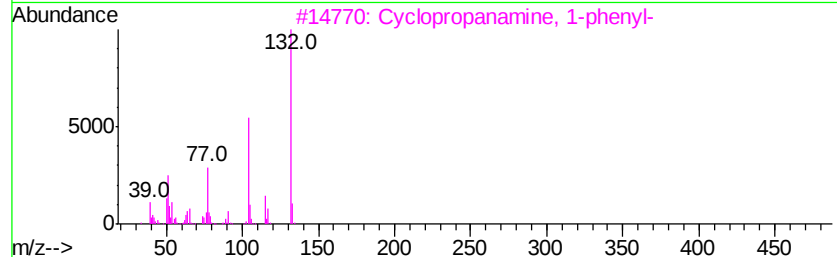
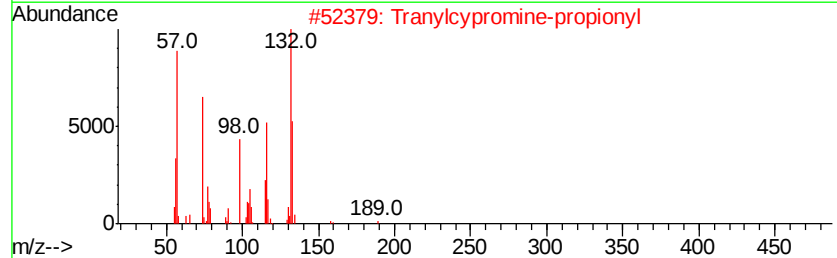
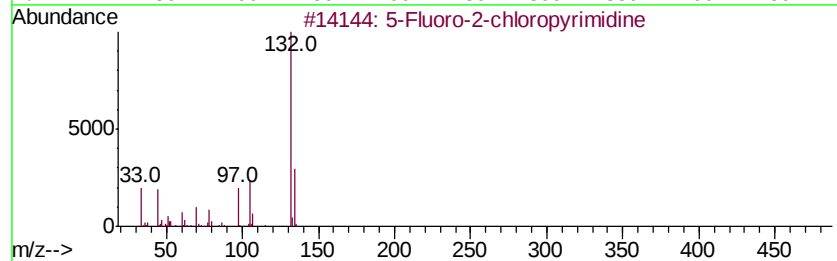
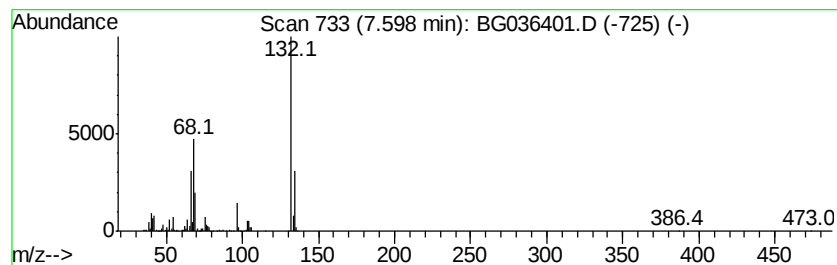
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Peak Number 2 unknown7.60 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.60	63.16 ng	959608	1,4-Dichlorobenzene-d4	8.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2			Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	10
3			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.17	3.8	ng	56954	1	8.07	303843	20.0
unknown7.60	7.60	63.2	ng	959608	1	8.07	303843	20.0