

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012119\
 Data File : BG039231.D
 Acq On : 21 Jan 2019 13:23
 Operator : JU/SJ
 Sample : PB116511BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116511BL

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-BG010919.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	5.092	337	341	348	rVB2	13189	21156	1.47%	0.339%
2	5.562	414	421	436	rVB	249553	406801	28.20%	6.527%
3	7.172	688	695	705	rBV	266420	460308	31.91%	7.385%
4	7.536	750	757	768	rVB	249058	438293	30.38%	7.032%
5	8.006	829	837	844	rBV	51841	89632	6.21%	1.438%
6	8.329	884	892	901	rBB	196146	344701	23.90%	5.530%
7	9.187	1031	1038	1050	rBV	149153	274379	19.02%	4.402%
8	10.826	1308	1317	1326	rBV	66613	125616	8.71%	2.015%
9	13.264	1724	1732	1742	rBV	494148	780595	54.11%	12.524%
10	14.651	1961	1968	1978	rBB2	126629	208241	14.44%	3.341%
11	16.143	2215	2222	2238	rBV	435593	681753	47.26%	10.938%
12	17.401	2429	2436	2452	rBV2	192424	294395	20.41%	4.723%
13	20.004	2873	2879	2886	rBV	1124845	1442481	100.00%	23.143%
14	21.672	3156	3163	3170	rBV2	205265	343112	23.79%	5.505%
15	24.933	3708	3718	3734	rVB	111757	321474	22.29%	5.158%

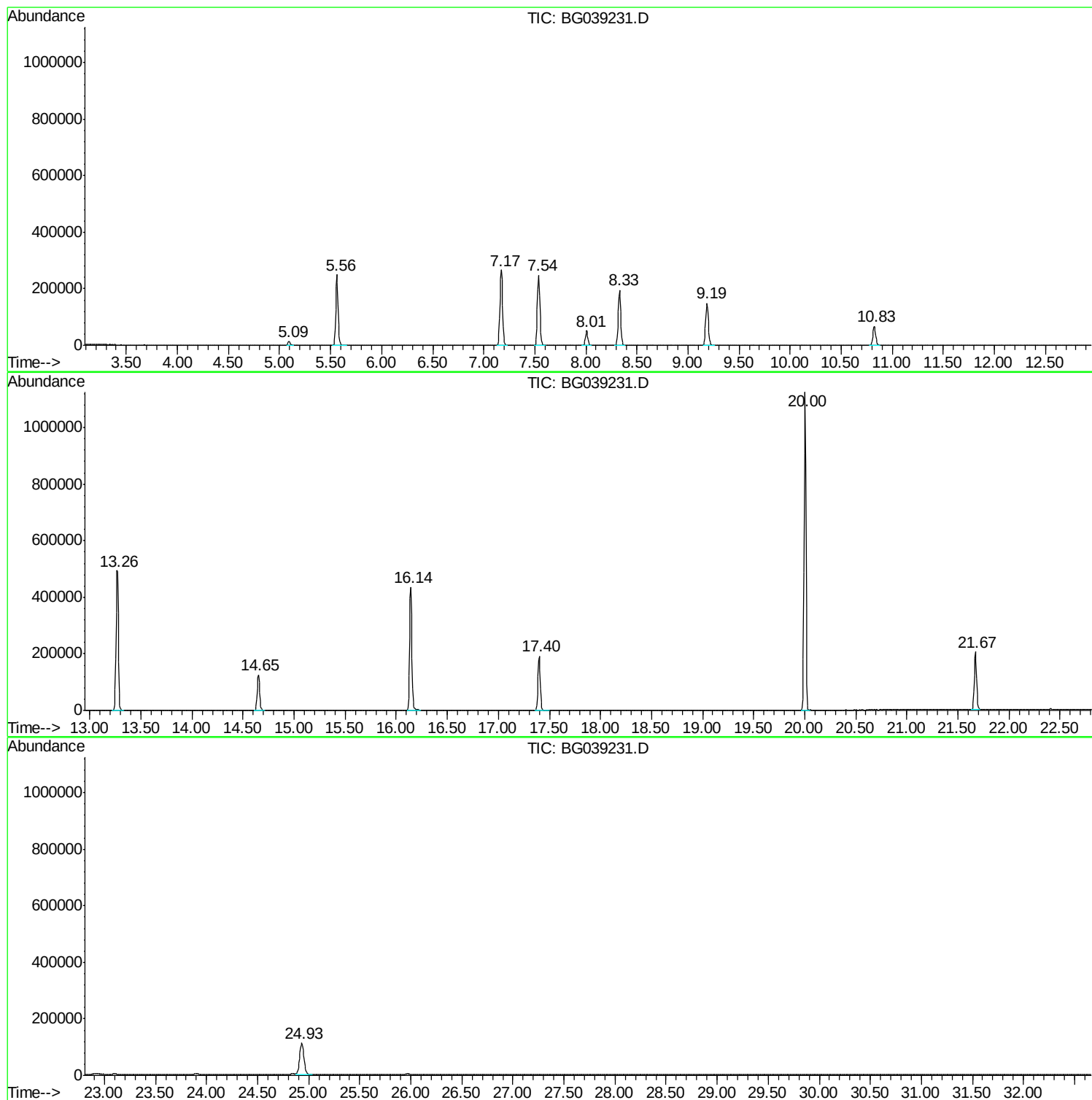
Sum of corrected areas: 6232937

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.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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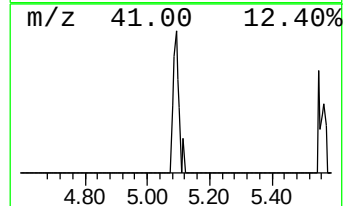
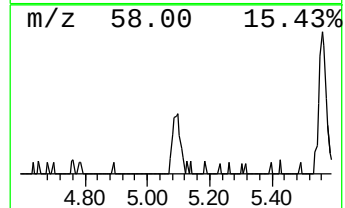
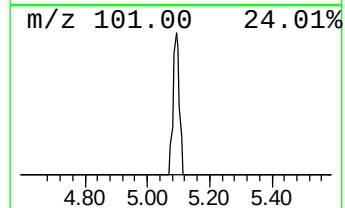
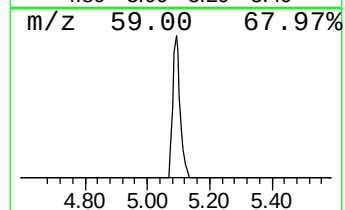
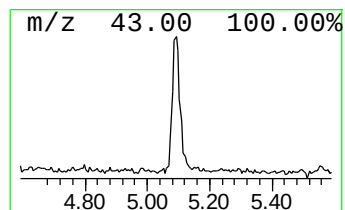
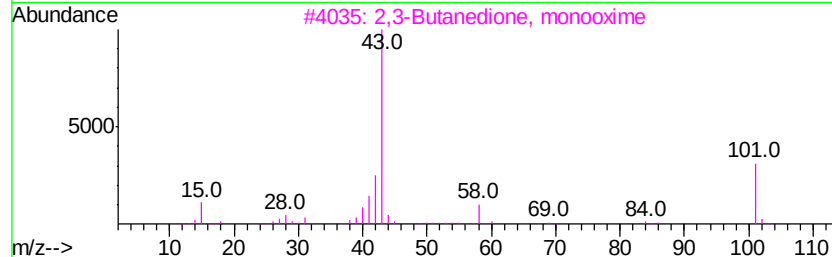
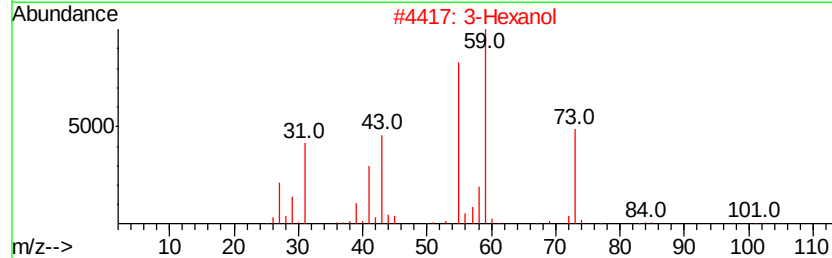
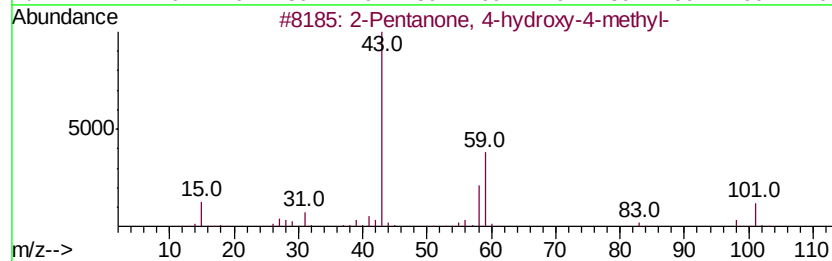
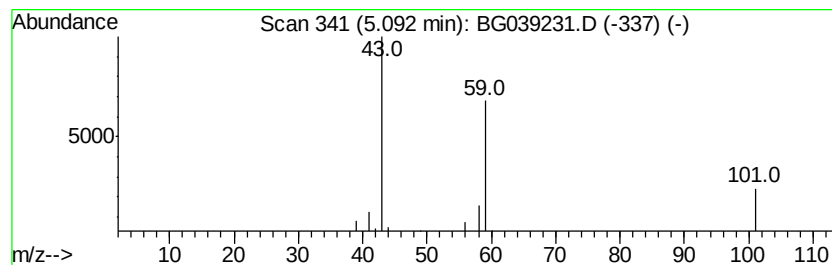
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	4.72 ng	21156	1,4-Dichlorobenzene-d4	8.01

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	102	C6H14O	000623-37-0	28
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9



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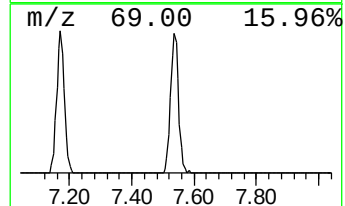
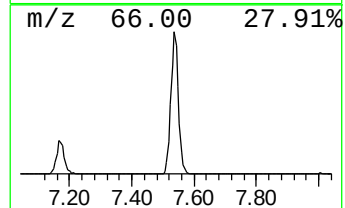
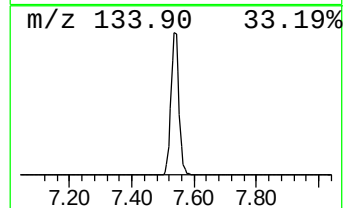
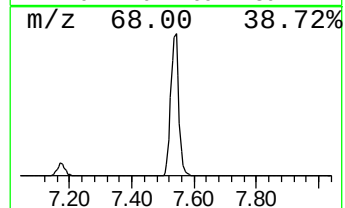
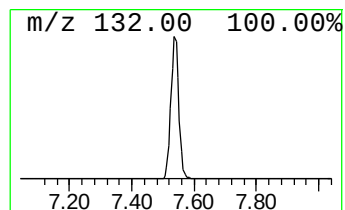
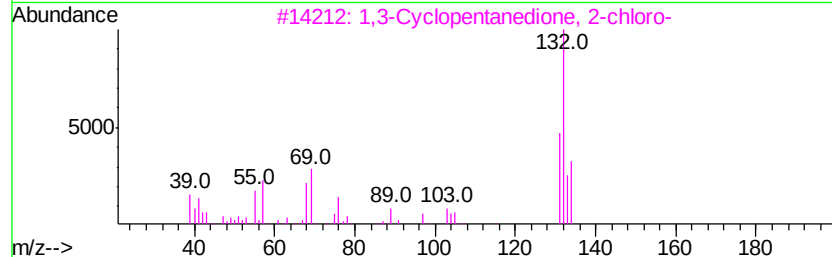
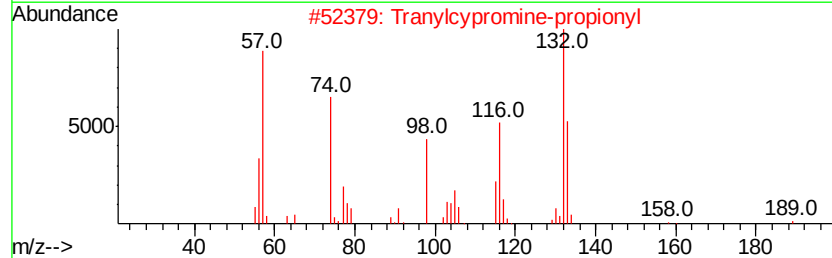
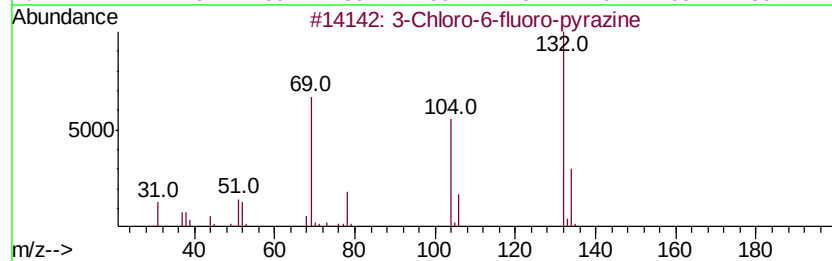
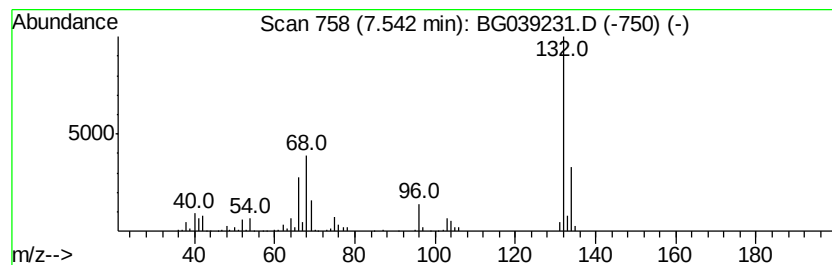
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Peak Number 2 unknown7.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	97.80 ng	438293	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	27
3		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	20
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	4.7	ng	21156	1	8.01	89632	20.0
unknown7.54	7.54	97.8	ng	438293	1	8.01	89632	20.0