

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG012219\
 Data File : BG039253.D
 Acq On : 22 Jan 2019 16:26
 Operator : JU/SJ
 Sample : PB116551BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116551BL

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	5.088	335	340	348	rBV	11461	17714	1.33%	0.307%
2	5.558	413	420	431	rBV	216448	349874	26.23%	6.071%
3	7.168	686	694	704	rBV	240861	427342	32.04%	7.415%
4	7.532	749	756	768	rBV	225564	396773	29.75%	6.885%
5	8.008	829	837	843	rBB	53116	89608	6.72%	1.555%
6	8.326	884	891	899	rBV	183975	326551	24.48%	5.667%
7	9.184	1029	1037	1047	rBV	136082	249053	18.67%	4.322%
8	10.823	1308	1316	1325	rBV	64274	125823	9.43%	2.183%
9	13.267	1724	1732	1744	rBV	446634	702693	52.68%	12.194%
10	14.648	1960	1967	1978	rBV2	122000	202649	15.19%	3.516%
11	16.140	2214	2221	2237	rBV	314126	514676	38.59%	8.931%
12	17.392	2428	2434	2446	rBV	175191	279793	20.98%	4.855%
13	20.000	2871	2878	2890	rBV	992620	1333769	100.00%	23.144%
14	21.669	3154	3162	3173	rBV	196506	343774	25.77%	5.965%
15	23.890	3532	3540	3547	rVB6	6311	13653	1.02%	0.237%
16	24.924	3705	3716	3733	rVB2	129397	389087	29.17%	6.752%

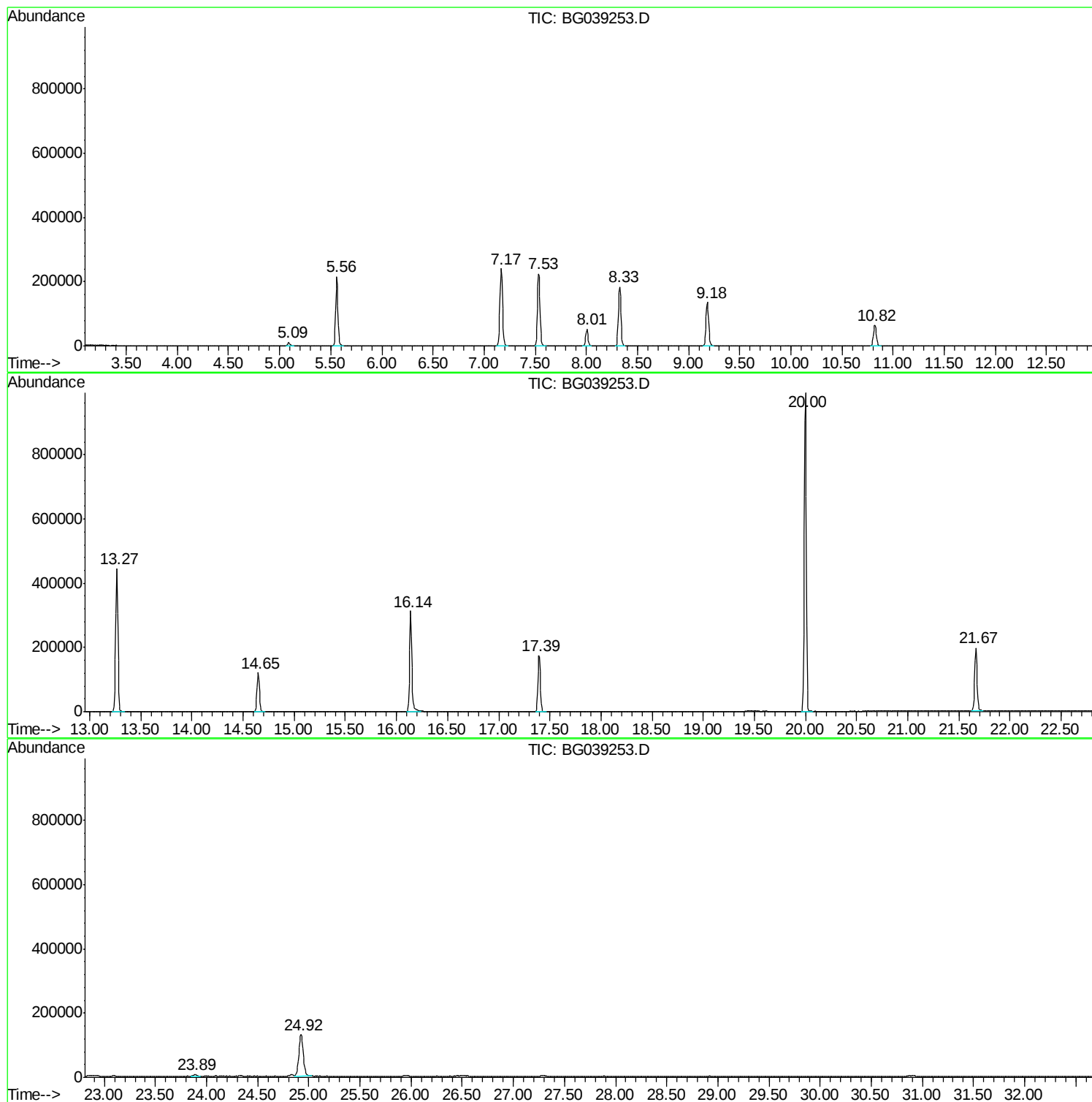
Sum of corrected areas: 5762832

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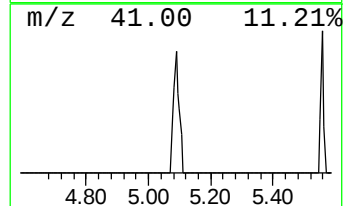
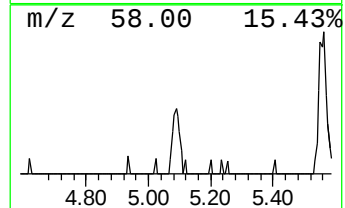
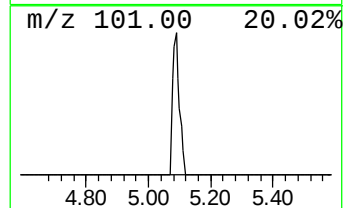
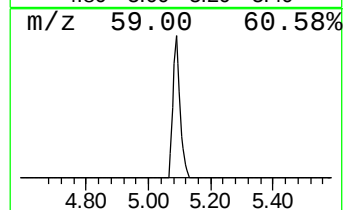
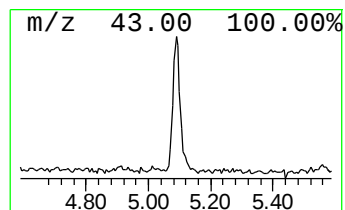
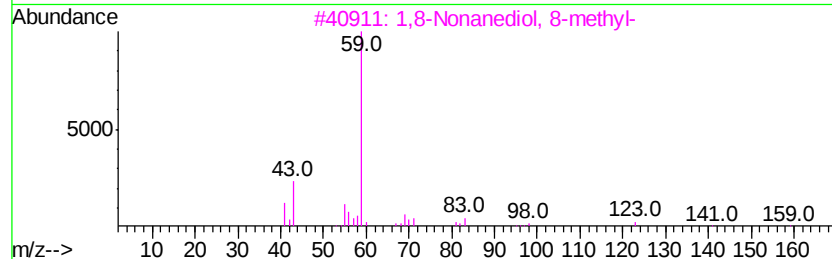
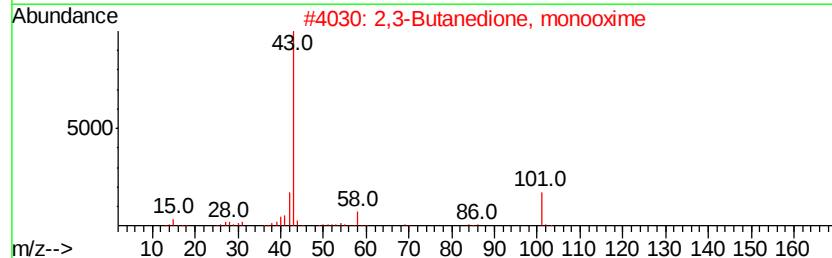
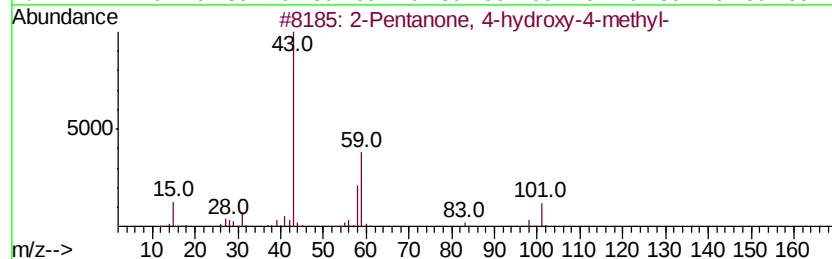
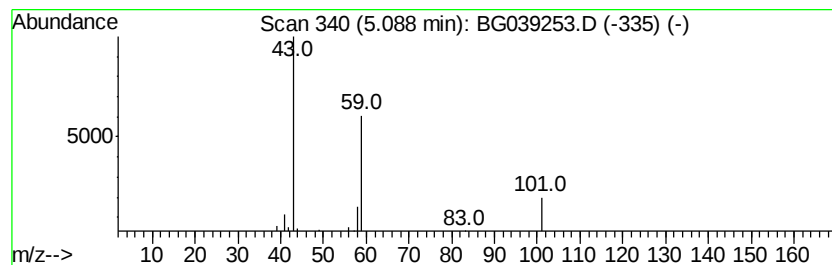
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	3.95 ng	17714	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	72
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	9



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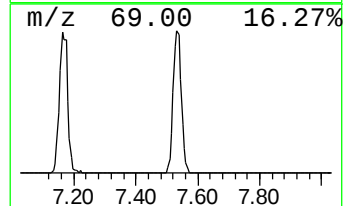
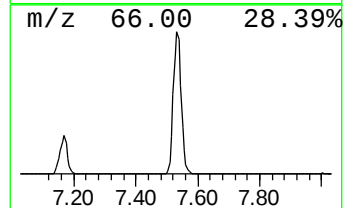
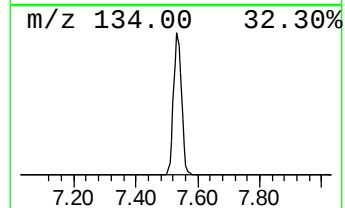
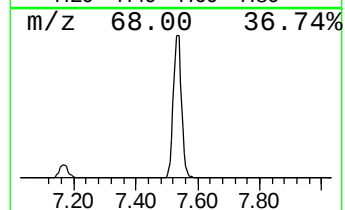
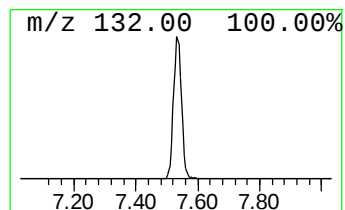
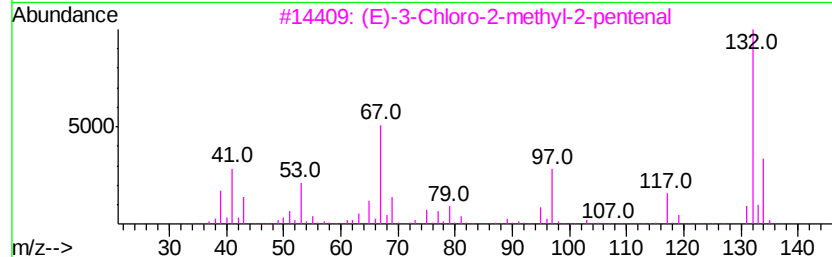
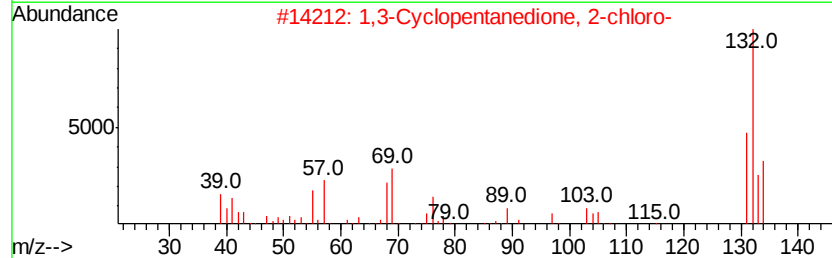
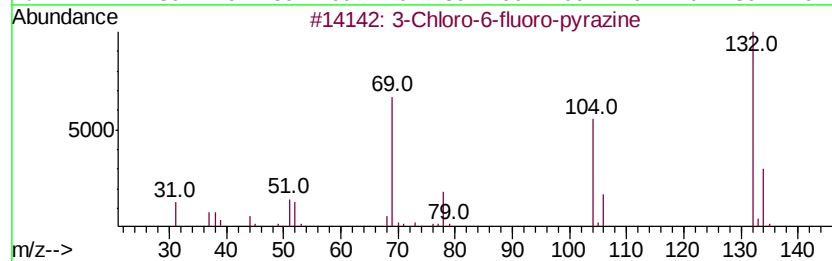
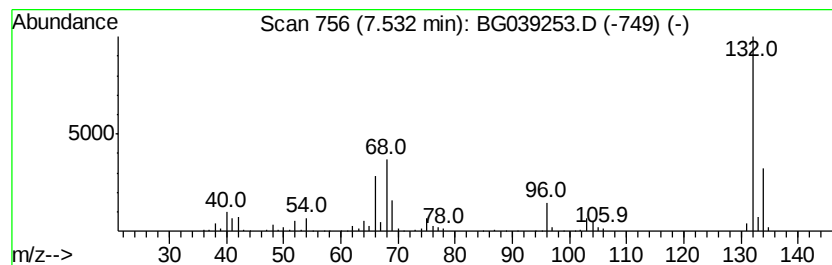
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Peak Number 2 unknown7.53 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.53	88.56 ng	396773	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		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	23
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	4.0	ng	17714	1	8.01	89608	20.0
unknown7.53	7.53	88.6	ng	396773	1	8.01	89608	20.0