

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
 Data File : BG049872.D
 Acq On : 17 Aug 2021 10:55
 Operator : CG/JU
 Sample : PB138424BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB138424BL

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-BG080321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG049872.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.097	344	350	358	rBV	39074	64835	2.97%	0.732%
2	5.579	425	432	443	rBV	447862	742825	34.05%	8.384%
3	7.194	699	707	719	rBV	489286	838211	38.42%	9.460%
4	8.029	841	849	857	rVB	87708	155228	7.12%	1.752%
5	9.197	1040	1048	1057	rBV	317122	561181	25.72%	6.334%
6	10.848	1322	1329	1336	rBV	122551	225700	10.35%	2.547%
7	13.298	1739	1746	1754	rBV	834828	1304531	59.80%	14.723%
8	14.678	1975	1981	1989	rBV	203348	333235	15.28%	3.761%
9	16.176	2228	2236	2248	rBV	738472	1192972	54.69%	13.464%
10	17.433	2443	2450	2462	rVB	278234	431270	19.77%	4.867%
11	20.042	2888	2894	2900	rBV	1664672	2181513	100.00%	24.621%
12	21.716	3172	3179	3186	rVB	246920	418931	19.20%	4.728%
13	24.988	3727	3736	3747	rVB2	138853	410007	18.79%	4.627%

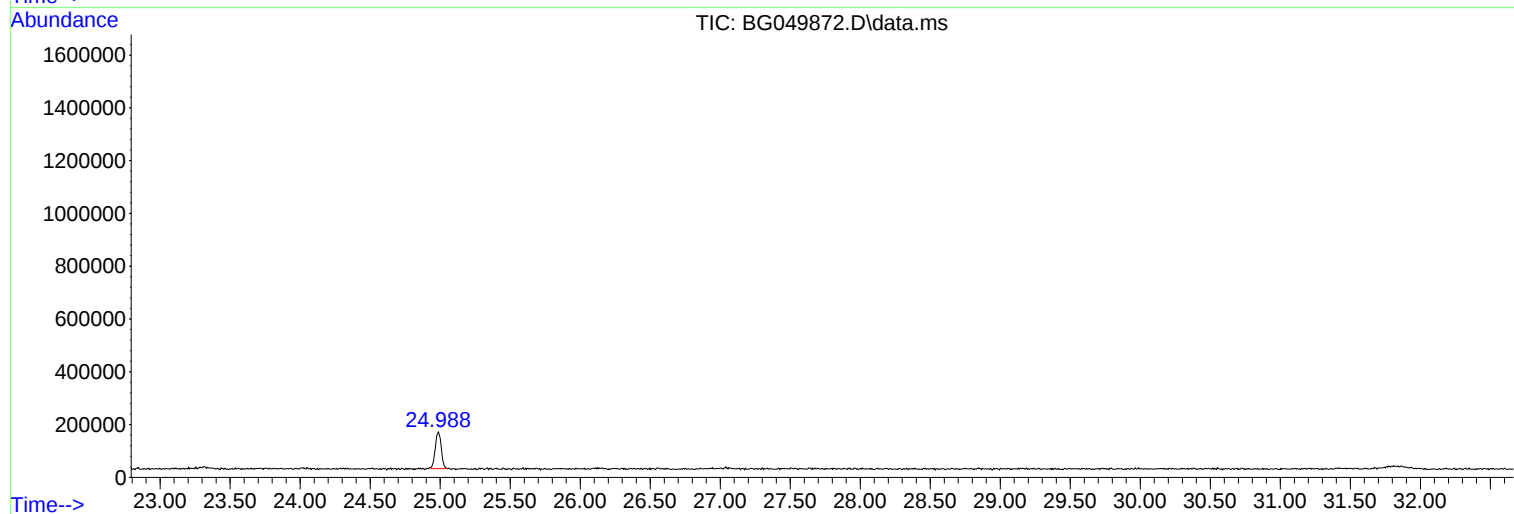
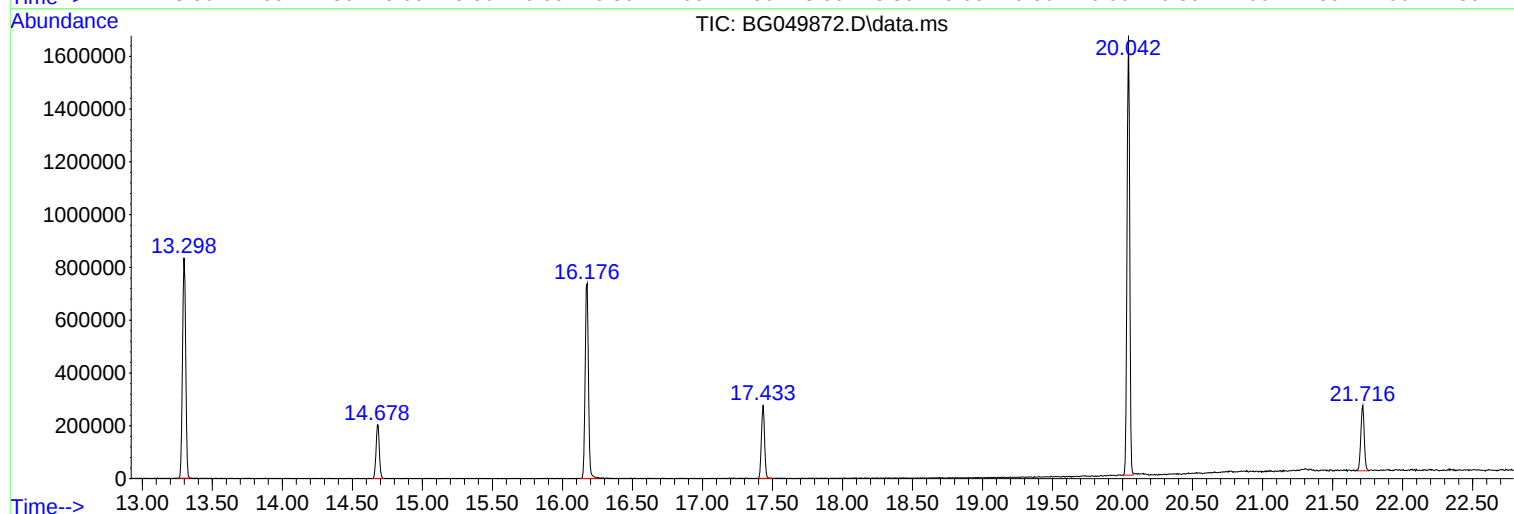
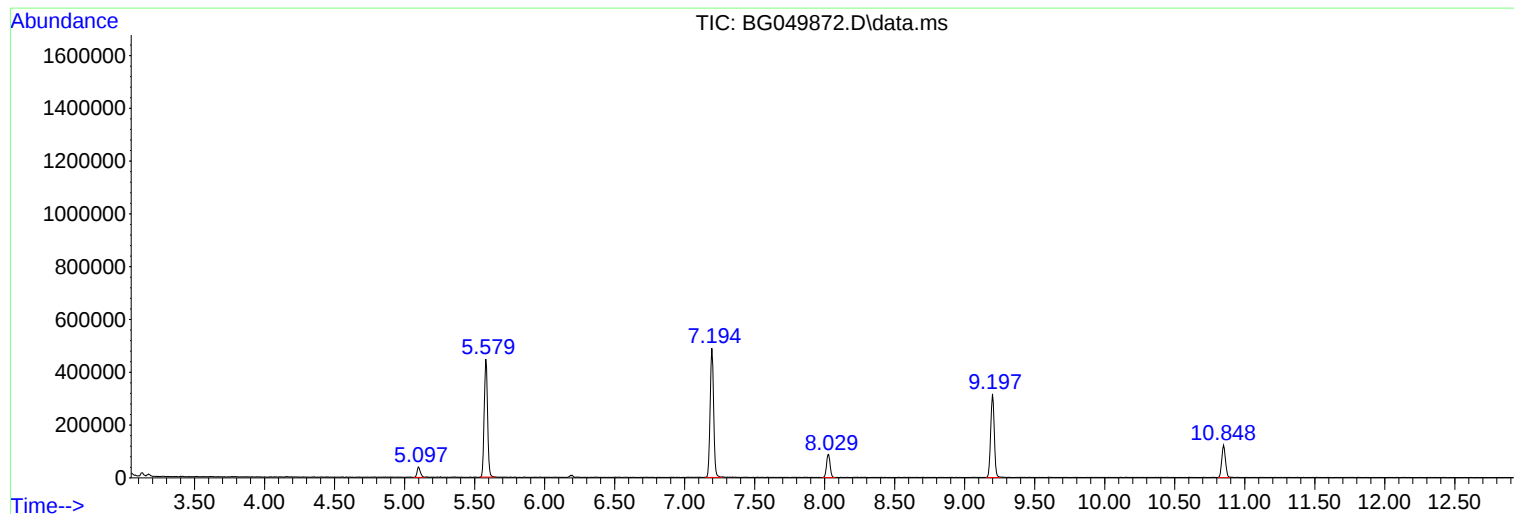
Sum of corrected areas: 8860439

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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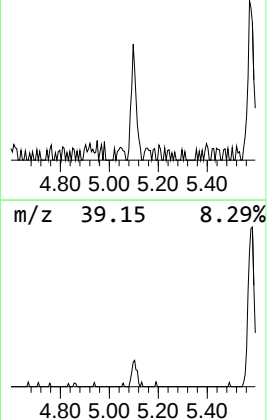
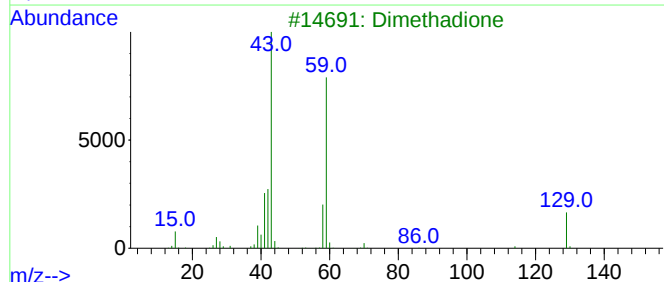
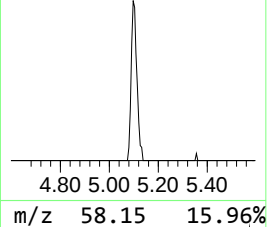
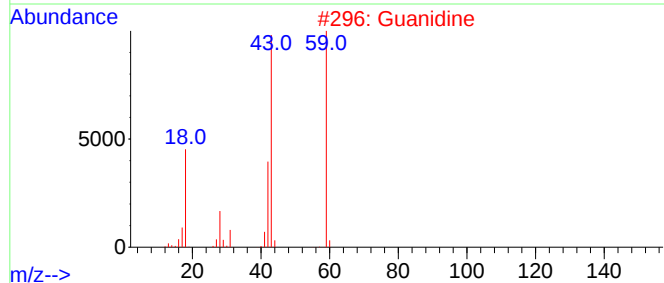
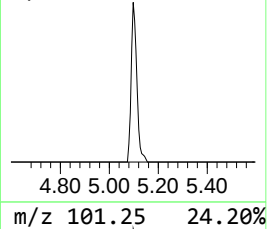
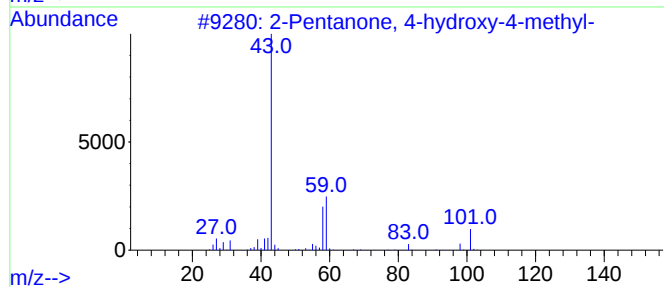
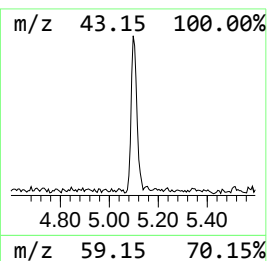
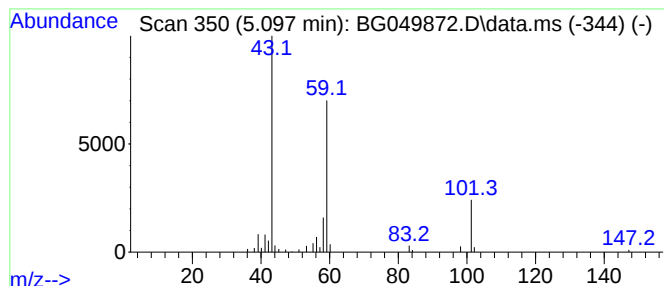
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.097	8.35 ng	64835	1,4-Dichlorobenzene-d4	8.028

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Guanidine	59	CH5N3	000113-00-8	47
3		Dimethadione	129	C5H7NO3	000695-53-4	33
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32
5		n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.097	8.3	ng	64835	1	8.028	155228	20.0