

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033121\
 Data File : BP005118.D
 Acq On : 31 Mar 2021 10:09
 Operator : CG/JU
 Sample : PB135372BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB135372BL

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_P\Methods\8270E-BP032321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005118.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.852	296	300	307	rBV	24833	34201	1.74%	0.422%
2	5.305	371	377	389	rBV	547841	799670	40.59%	9.868%
3	6.887	639	646	657	rBV	515615	812991	41.26%	10.033%
4	7.710	776	786	795	rBV	97799	153992	7.82%	1.900%
5	8.869	976	983	1002	rBV	319098	520856	26.44%	6.428%
6	10.499	1252	1260	1269	rBV	121541	203752	10.34%	2.514%
7	12.981	1672	1682	1694	rBV	840056	1213566	61.59%	14.976%
8	14.363	1910	1917	1925	rBV	193531	285902	14.51%	3.528%
9	15.863	2165	2172	2189	rBV	686620	944179	47.92%	11.652%
10	17.122	2380	2386	2393	rBV	248390	355108	18.02%	4.382%
11	19.698	2818	2824	2837	rBV	1681658	1970275	100.00%	24.315%
12	21.216	3077	3082	3092	rBV	307335	412465	20.93%	5.090%
13	23.504	3464	3471	3480	rVB2	199466	396312	20.11%	4.891%

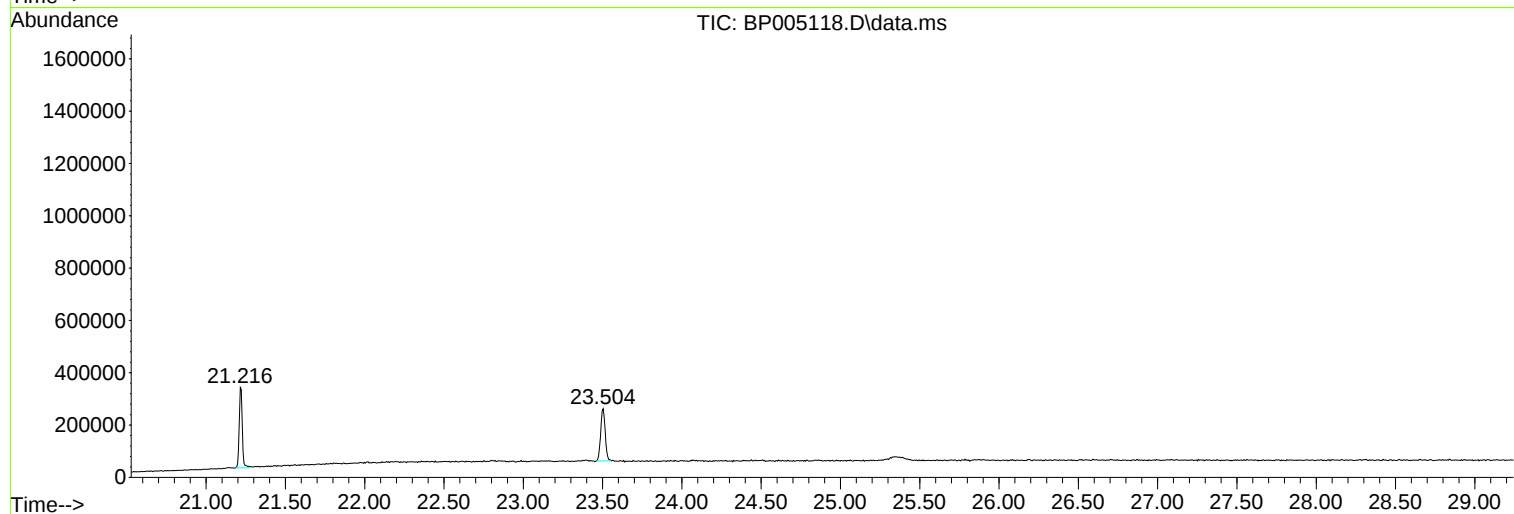
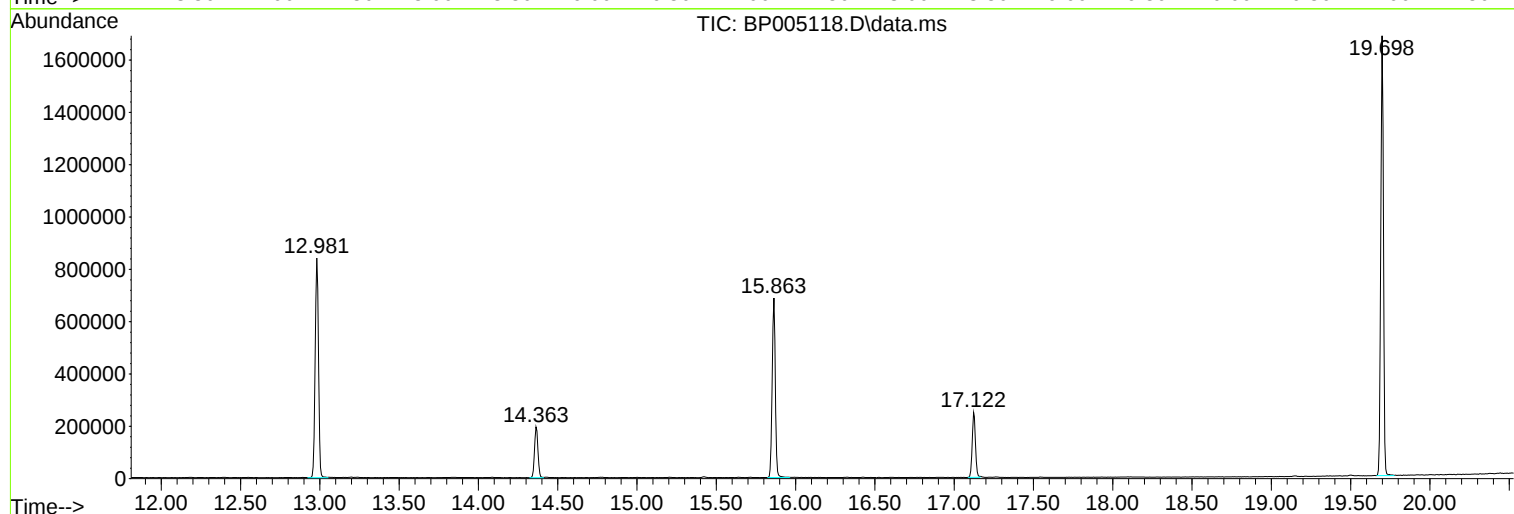
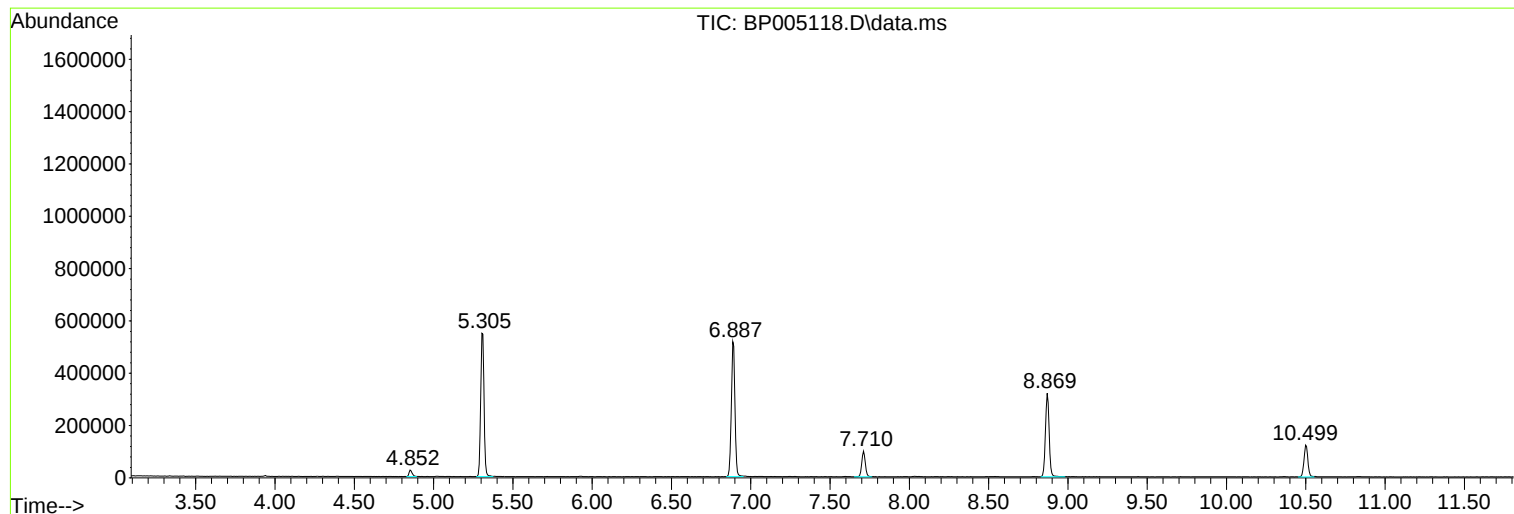
Sum of corrected areas: 8103269

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.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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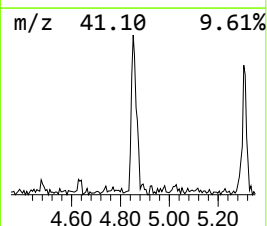
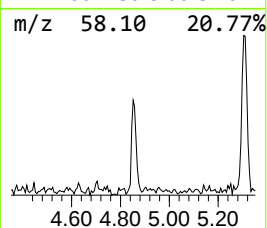
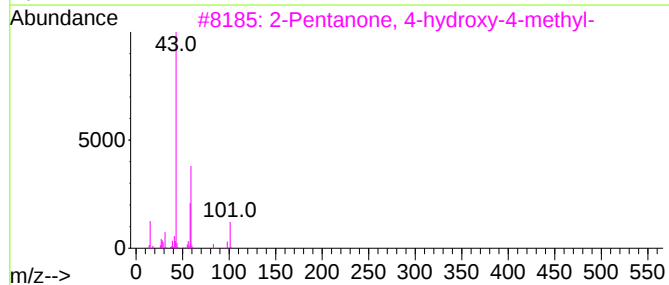
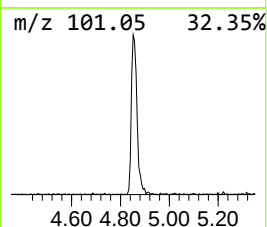
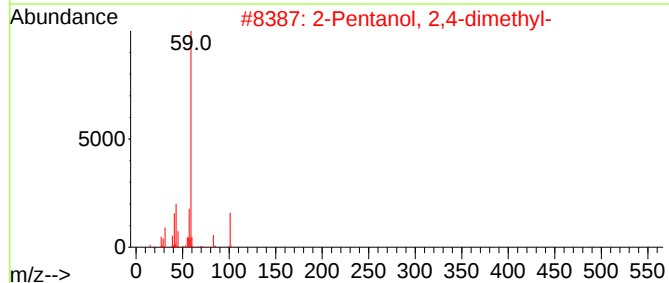
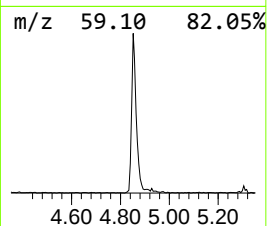
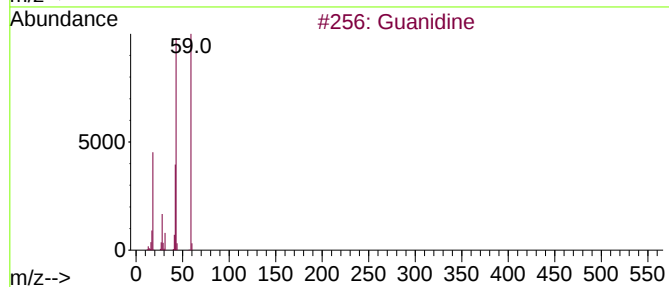
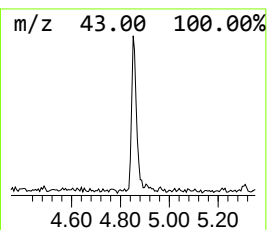
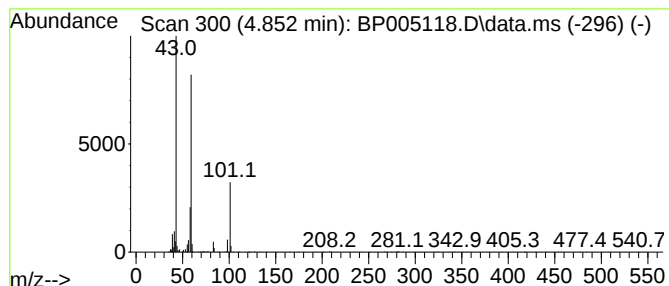
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Peak Number 1 Guanidine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.852	4.44 ng	34201	1,4-Dichlorobenzene-d4	7.710

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Guanidine	59	CH5N3	000113-00-8	50
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	45
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
4			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	40
5			Dimethadione	129	C5H7NO3	000695-53-4	38



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Guanidine	4.852	4.4	ng	34201	1	7.710	153992	20.0