

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081721\
 Data File : BP006714.D
 Acq On : 17 Aug 2021 12:51
 Operator : CG/JU
 Sample : PB138349BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB138349BL

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-BP080321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP006714.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	313	rVB	270878	384052	3.80%	0.832%
2	5.305	371	377	387	rBV	3499727	4862643	48.11%	10.535%
3	6.887	639	646	663	rVB	3263914	5048811	49.95%	10.939%
4	7.704	779	785	793	rVB	649420	999035	9.88%	2.165%
5	8.863	975	982	995	rVB	2060657	3304768	32.70%	7.160%
6	10.487	1251	1258	1266	rBV	887084	1400975	13.86%	3.035%
7	12.969	1671	1680	1691	rBV	4746213	6651442	65.81%	14.411%
8	14.345	1907	1914	1924	rBV	1199747	1709006	16.91%	3.703%
9	15.839	2161	2168	2192	rBV	3659866	5102889	50.49%	11.056%
10	17.104	2375	2383	2395	rBV	1393451	1992157	19.71%	4.316%
11	19.686	2816	2822	2827	rBV	8660051	10107212	100.00%	21.898%
12	21.210	3076	3081	3088	rBV	1785568	2258648	22.35%	4.894%
13	23.527	3468	3475	3486	rVB2	1167630	2333373	23.09%	5.056%

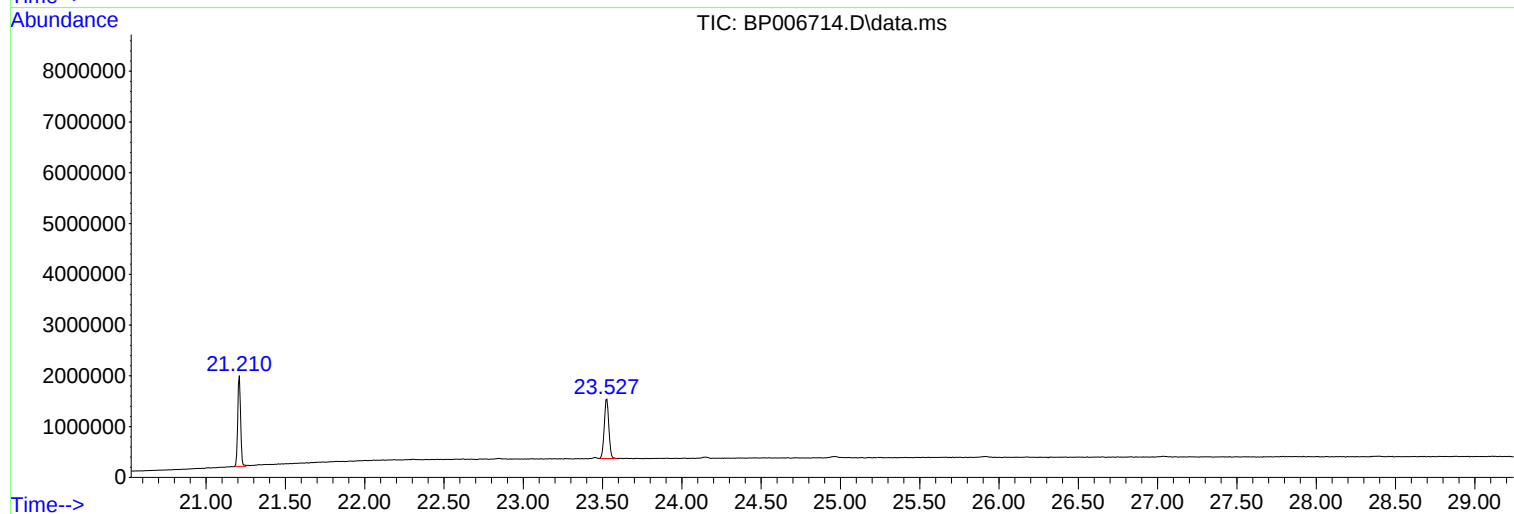
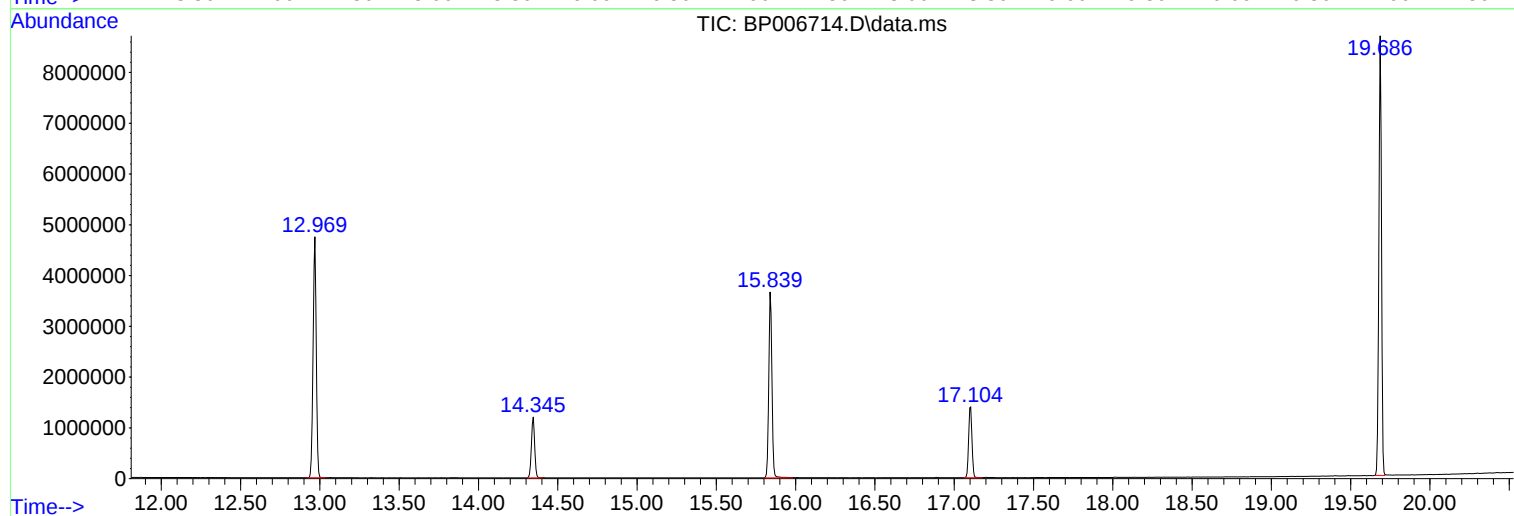
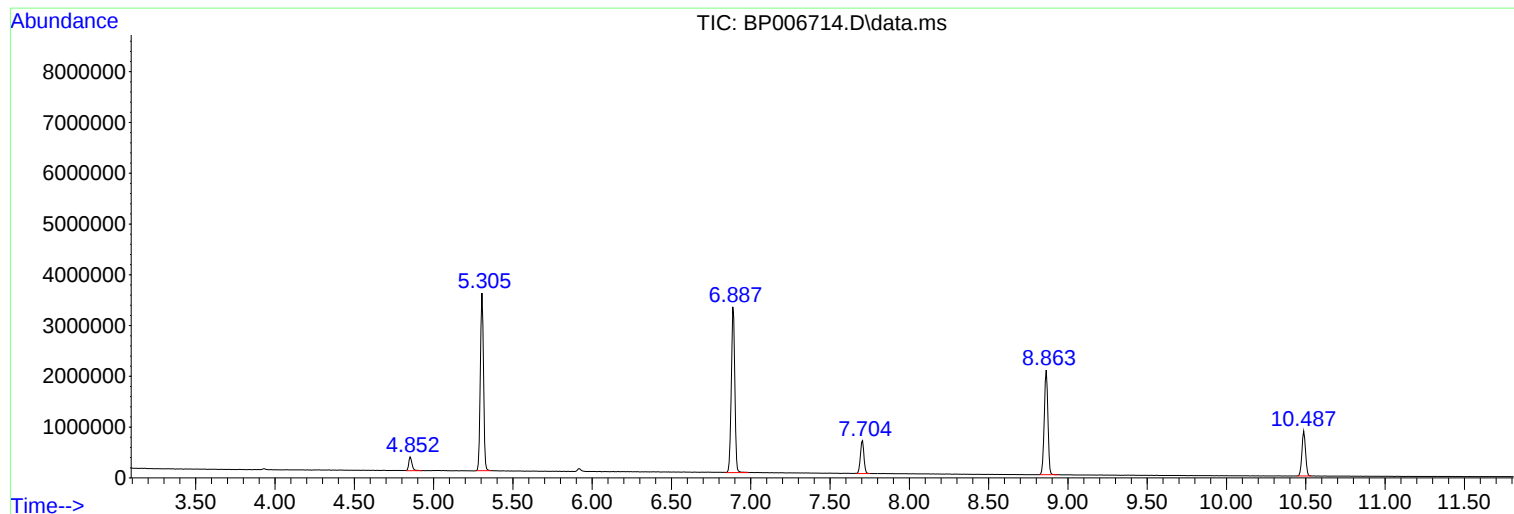
Sum of corrected areas: 46155011

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.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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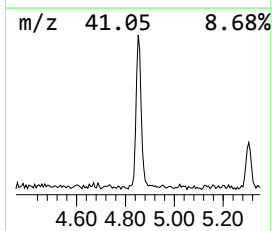
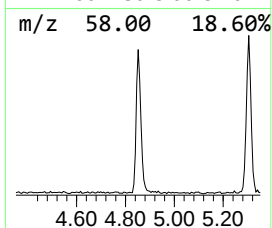
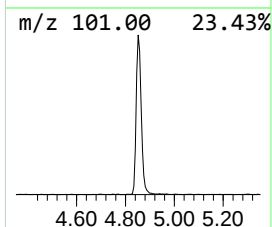
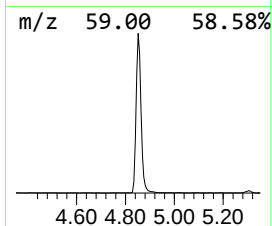
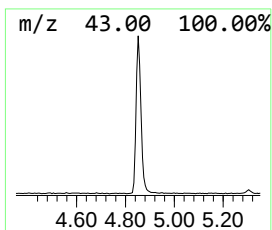
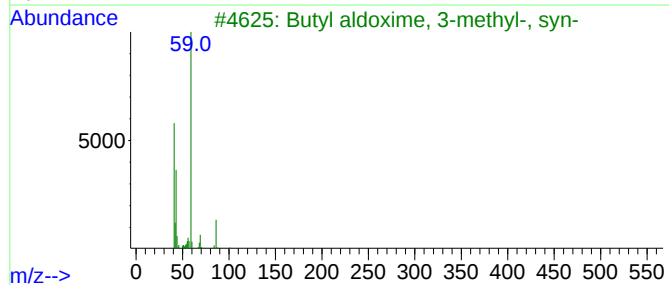
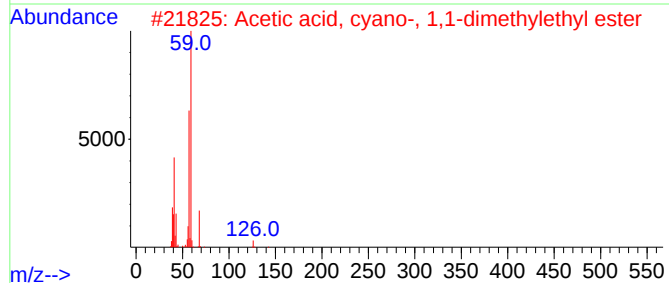
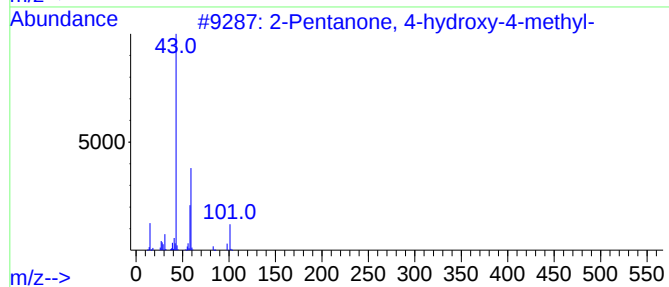
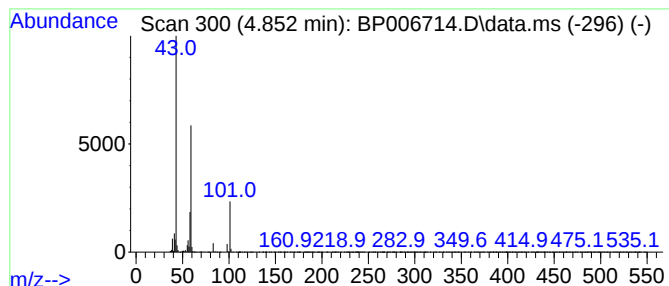
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.852	7.69 ng	384052	1,4-Dichlorobenzene-d4	7.704

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	12		
4	Acetone	58	C3H6O	000067-64-1	9		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.852	7.7	ng	384052	1	7.704	999035	20.0