

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
 Data File : BM031785.D
 Acq On : 17 Aug 2021 17:19
 Operator : CG/JU
 Sample : PB138482BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB138482BL

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_M\Methods\8270-BM081621.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM031785.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.746	277	282	292	rBV	90422	126644	3.51%	0.844%
2	5.187	350	357	370	rBV	1076859	1550800	42.95%	10.337%
3	6.746	614	622	632	rBV	1083449	1646697	45.61%	10.976%
4	7.545	752	758	765	rBV	187381	286410	7.93%	1.909%
5	8.698	946	954	971	rBV	640228	1033564	28.63%	6.889%
6	10.310	1219	1228	1238	rBV	236814	396448	10.98%	2.642%
7	12.798	1644	1651	1661	rBV	1522199	2149587	59.54%	14.328%
8	14.180	1878	1886	1893	rBV	332922	507137	14.05%	3.380%
9	15.686	2135	2142	2157	rBV	1296867	1780056	49.30%	11.865%
10	16.939	2349	2355	2364	rVB	434172	609806	16.89%	4.065%
11	19.586	2799	2805	2810	rBV	3025823	3610569	100.00%	24.065%
12	21.133	3063	3068	3078	rBV	534684	686032	19.00%	4.573%
13	23.286	3428	3434	3444	rVB	351668	619344	17.15%	4.128%

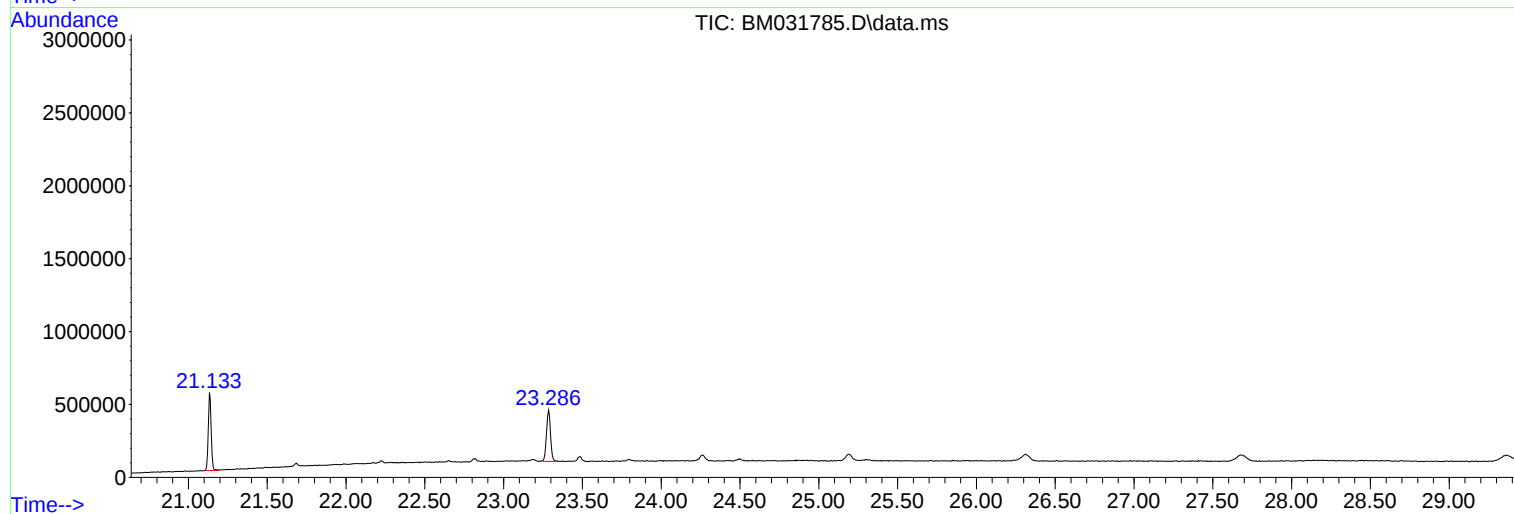
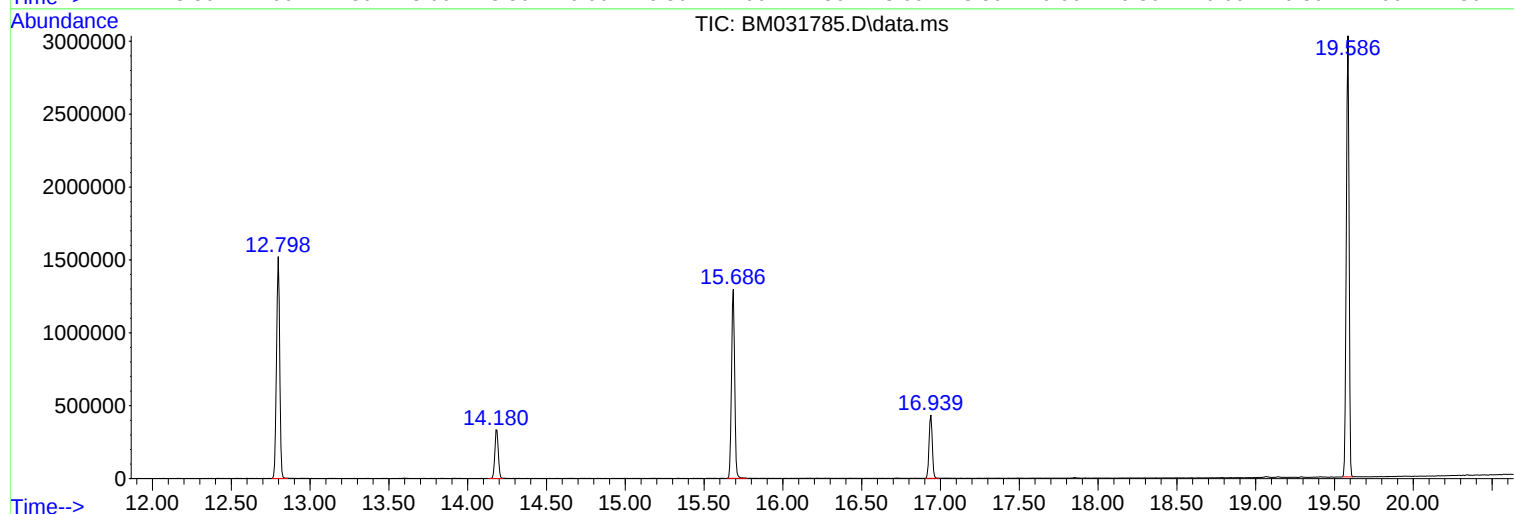
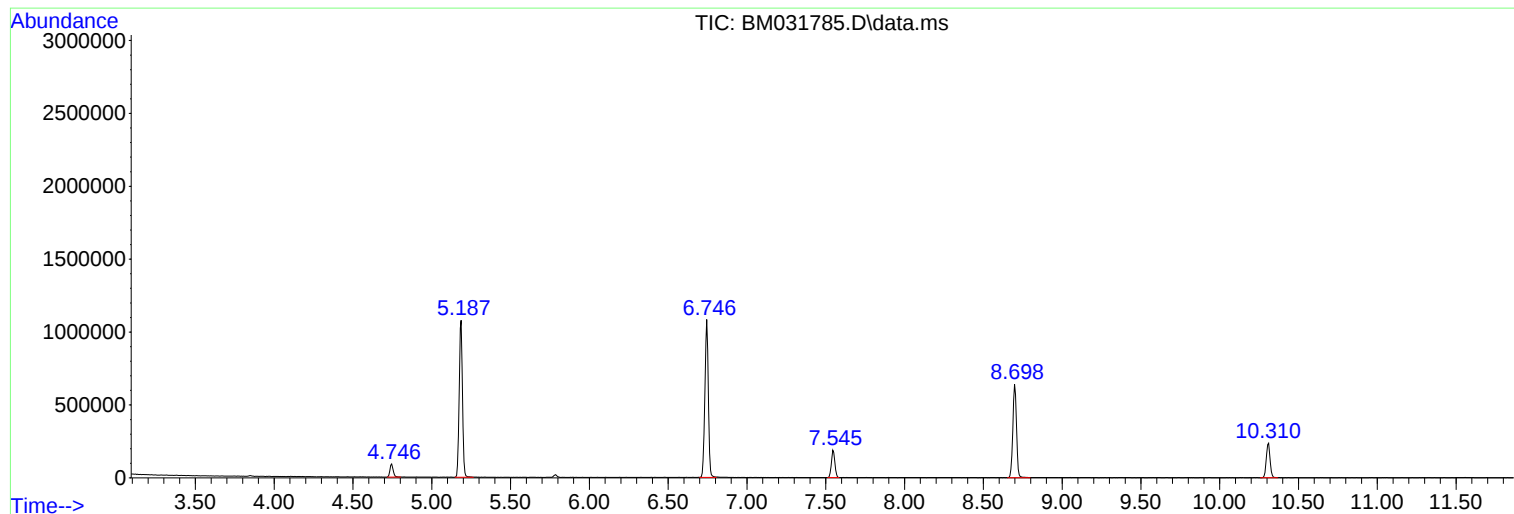
Sum of corrected areas: 15003094

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.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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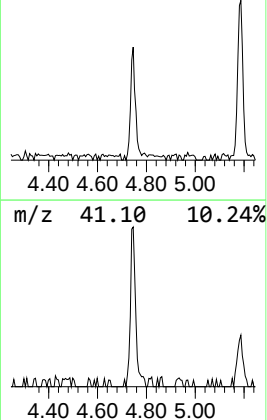
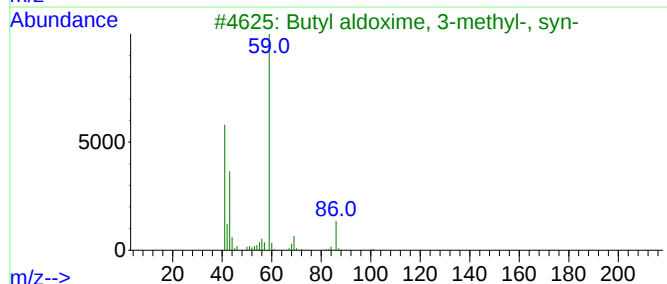
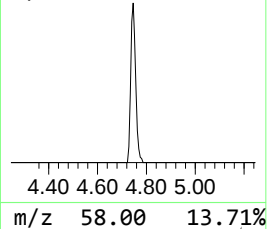
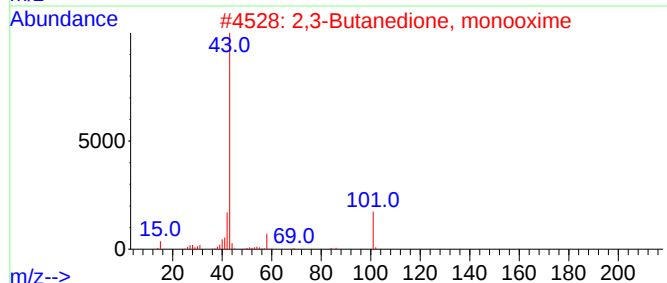
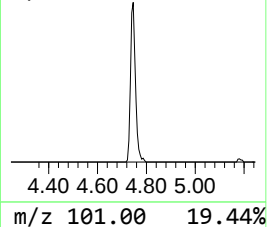
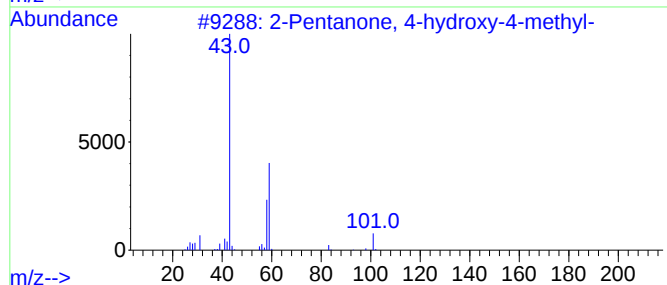
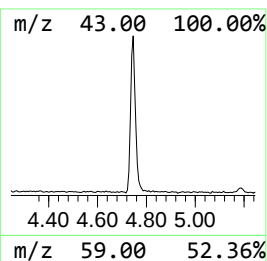
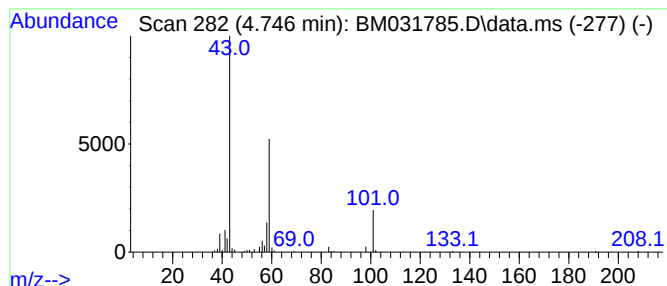
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.746	8.84 ng	126644	1,4-Dichlorobenzene-d4	7.545

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	25		
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.746	8.8	ng	126644	1	7.545	286410	20.0