

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011623\
Data File : BP013484.D
Acq On : 16 Jan 2023 16:37
Operator : CG/JU
Sample : PB150193BL
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB150193BL

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-BP011223.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP013484.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.993	302	307	320	rBV	3939639	5521726	16.74%	3.807%
2	5.463	380	387	402	rBV	9340316	13228759	40.10%	9.120%
3	7.075	653	661	677	rBV	8984816	14141564	42.87%	9.749%
4	7.905	795	802	814	rVB	1574814	2418221	7.33%	1.667%
5	9.075	993	1001	1020	rBV	5263124	8618818	26.13%	5.942%
6	10.716	1271	1280	1291	rBV	2008508	3479084	10.55%	2.398%
7	13.175	1691	1698	1714	rBV	14037039	20600136	62.44%	14.201%
8	14.557	1924	1933	1941	rBV	3374355	4927644	14.94%	3.397%
9	16.051	2180	2187	2206	rBV	12602355	17492451	53.02%	12.059%
10	17.316	2394	2402	2414	rBV	4342824	6257527	18.97%	4.314%
11	19.874	2831	2837	2842	rBV	25993630	32989928	100.00%	22.743%
12	21.404	3092	3097	3107	rBV	5797894	7393533	22.41%	5.097%
13	23.868	3510	3516	3527	rVB2	4042997	7986852	24.21%	5.506%

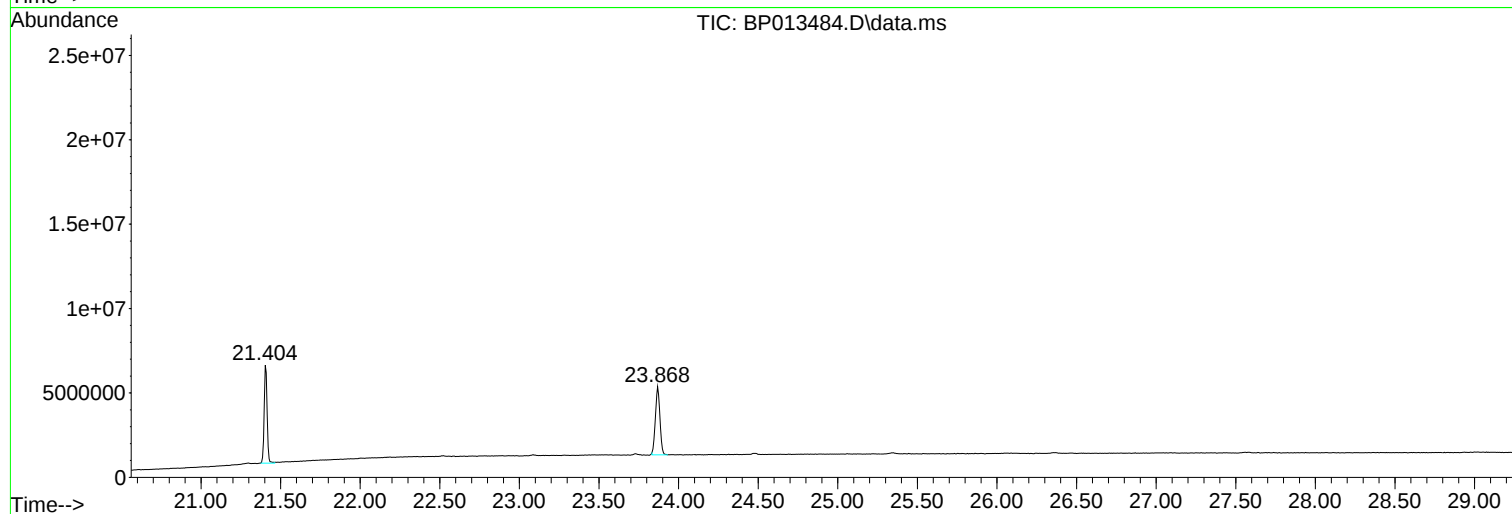
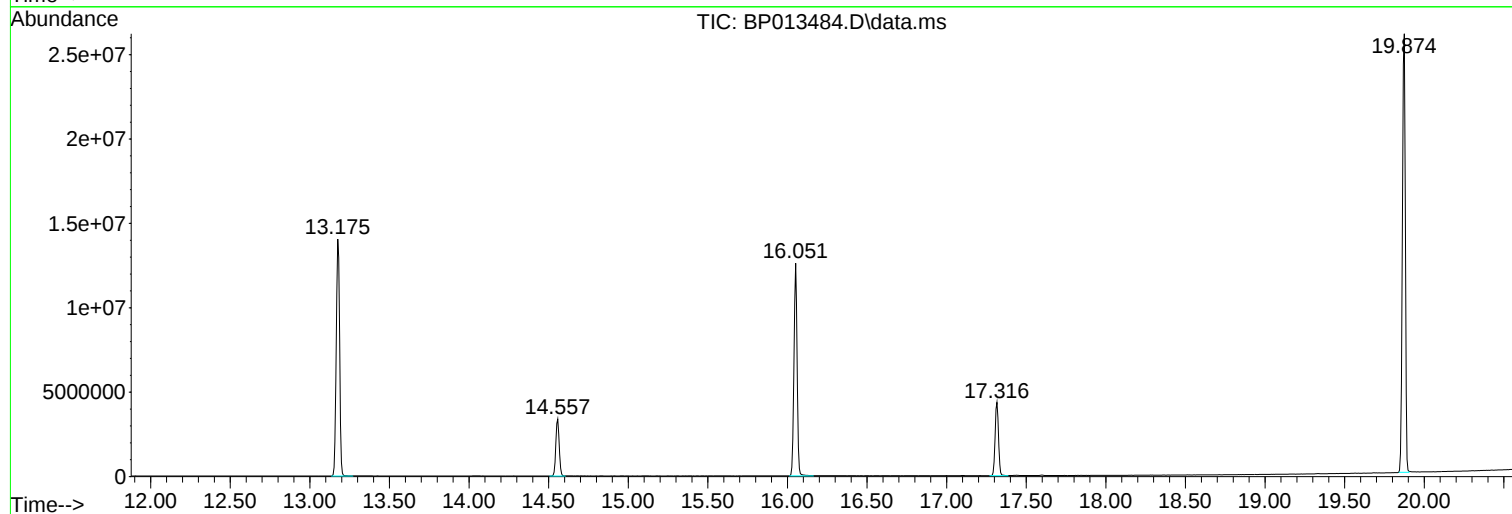
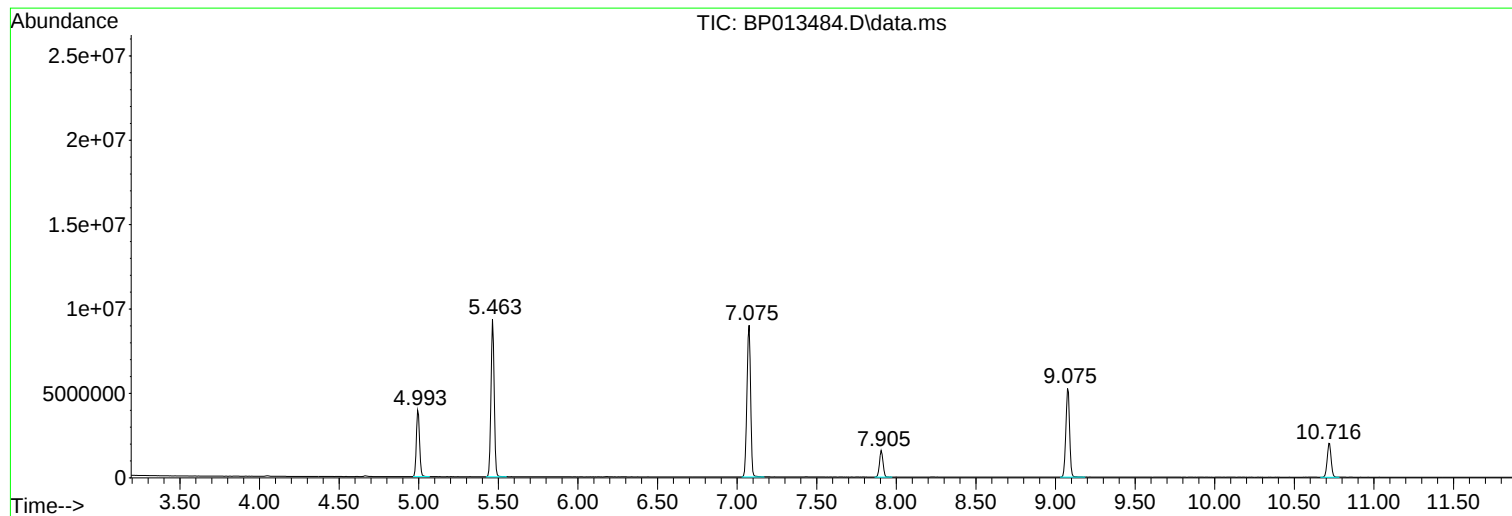
Sum of corrected areas: 145056243

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011223.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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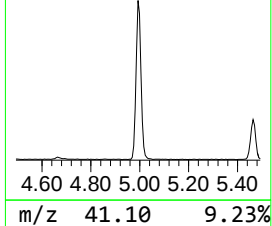
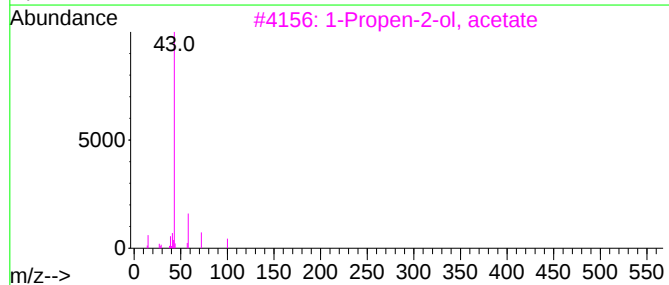
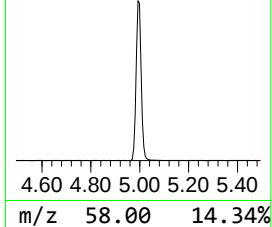
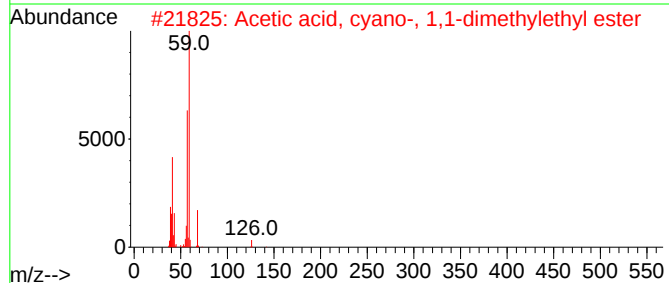
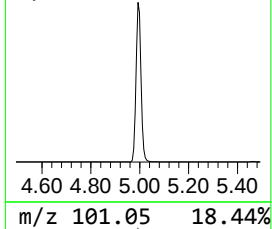
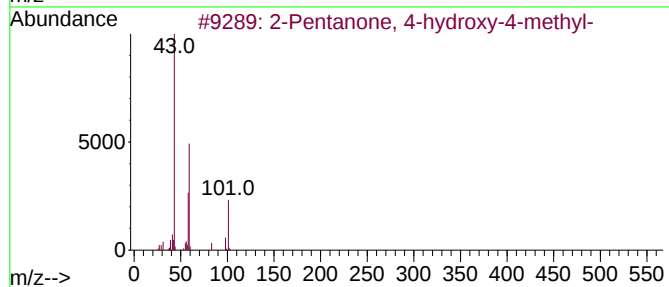
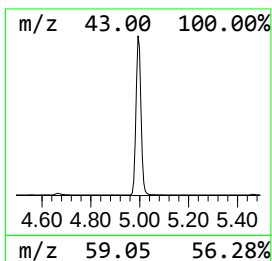
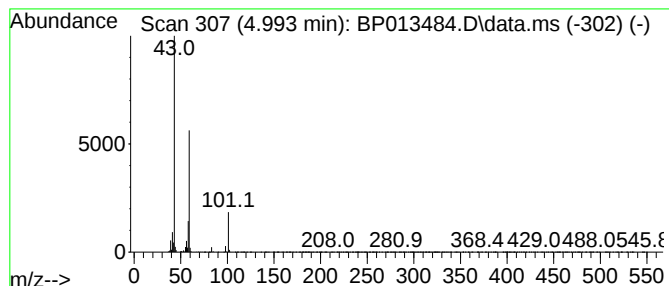
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.993	45.67 ng	5521730	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.993	45.7	ng	5521730	1	7.905	2418220	20.0