

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP073125\
Data File : BP025279.D
Acq On : 31 Jul 2025 14:04
Operator : CG/JU
Sample : PB169066BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB169066BL

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-BP072125.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025279.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.637	284	289	302	rBV	461890	709869	2.83%	0.716%
2	5.096	361	367	390	rBV	5175020	9187595	36.64%	9.264%
3	6.643	623	630	653	rBV	3947022	8913891	35.55%	8.988%
4	7.396	751	758	771	rBV	1223795	1929800	7.70%	1.946%
5	8.531	943	951	965	rBV	3249447	5829522	23.25%	5.878%
6	10.137	1216	1224	1251	rBV	1168901	2579476	10.29%	2.601%
7	12.637	1641	1649	1674	rBV	9345747	14546872	58.02%	14.667%
8	14.042	1878	1888	1904	rBV	1811623	3428521	13.67%	3.457%
9	15.584	2142	2150	2171	rBV	6174895	11431599	45.59%	11.526%
10	16.866	2361	2368	2388	rBV2	2178886	4273180	17.04%	4.309%
11	19.630	2833	2838	2848	rBV	19064131	25072247	100.00%	25.280%
12	21.301	3116	3122	3133	rBV2	2825125	5407663	21.57%	5.452%
13	24.395	3642	3648	3668	rVB2	2067802	5867717	23.40%	5.916%

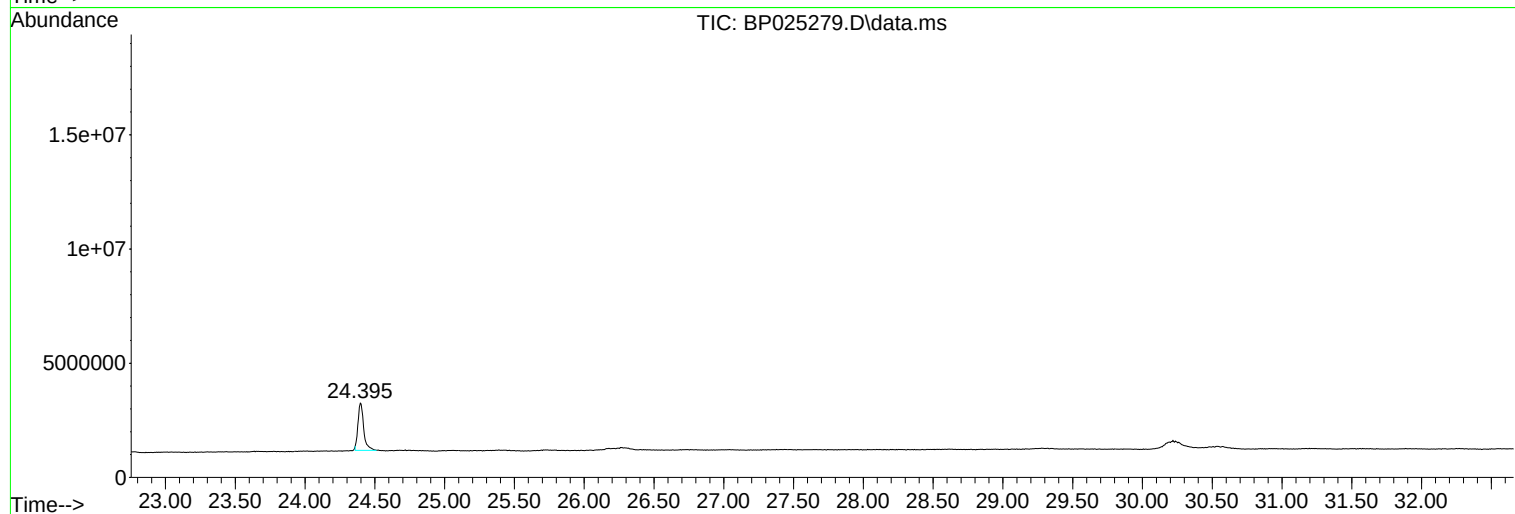
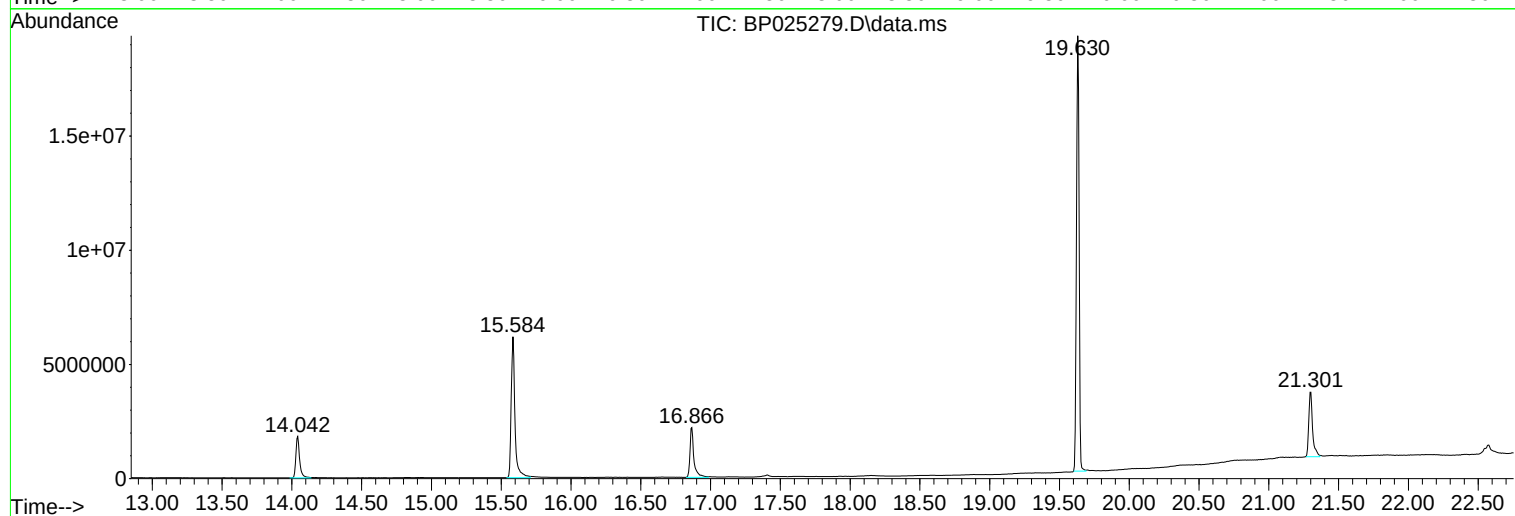
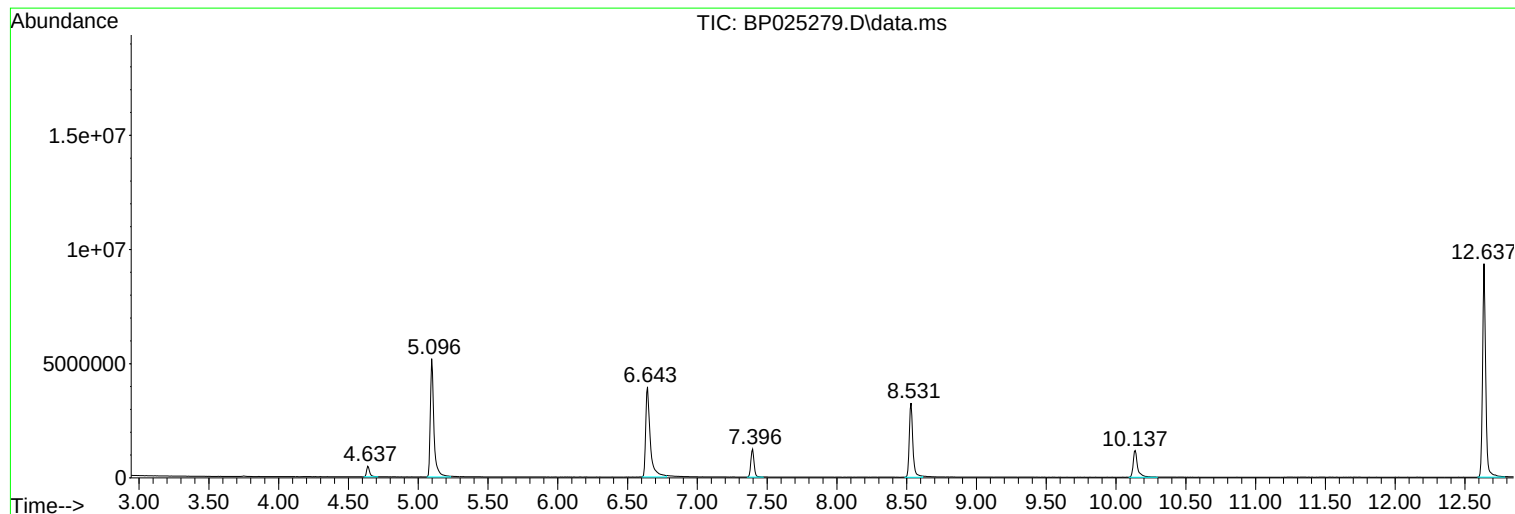
Sum of corrected areas: 99177952

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072125.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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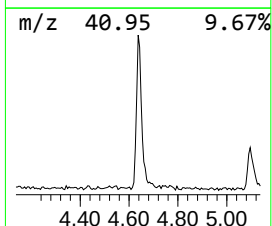
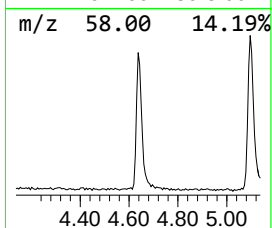
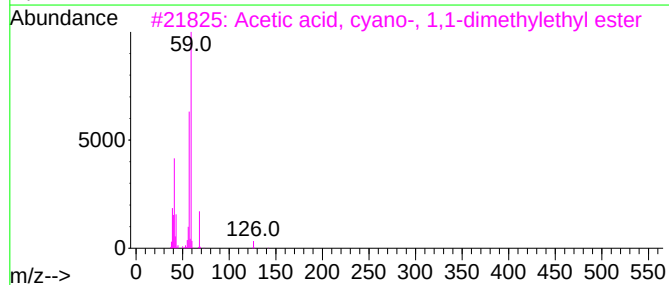
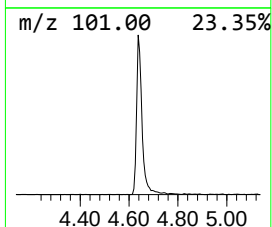
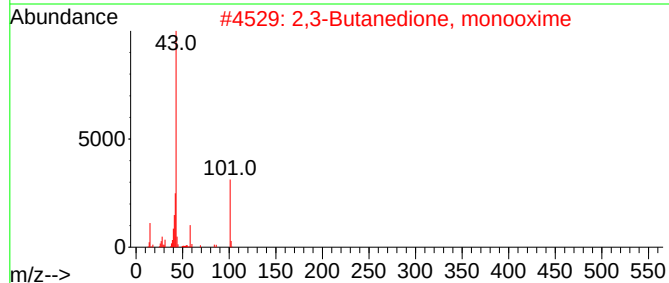
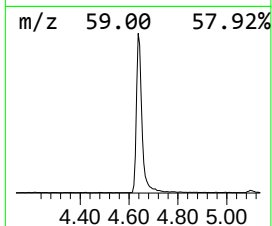
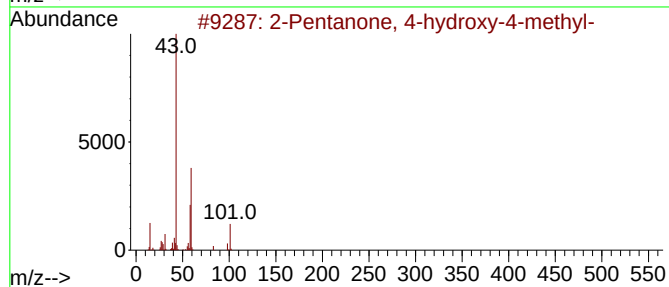
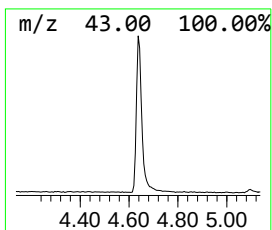
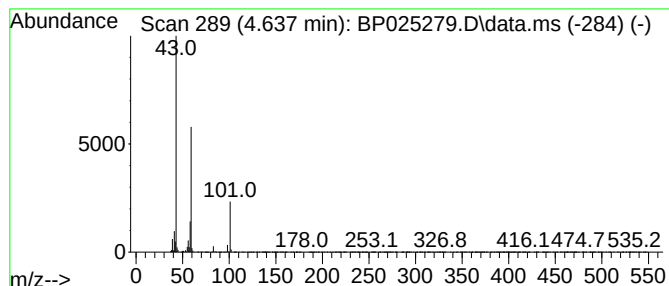
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.637	7.36 ng	709869	1,4-Dichlorobenzene-d4	7.396

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	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.637	7.4	ng	709869	1	7.396	1929800	20.0