

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091525\
Data File : BP025752.D
Acq On : 15 Sep 2025 14:33
Operator : CG/JU
Sample : PB169526BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB169526BL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BP091225.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025752.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.919	330	337	348	rBV	386610	565092	5.10%	1.225%
2	5.384	409	416	440	rBV	2909537	4707225	42.46%	10.204%
3	6.943	674	681	711	rBV	2643422	4790095	43.21%	10.384%
4	7.755	812	819	831	rBV	555101	876893	7.91%	1.901%
5	8.902	1005	1014	1036	rBV	1709127	3150355	28.42%	6.829%
6	10.531	1283	1291	1310	rBV	521399	1178442	10.63%	2.555%
7	12.990	1701	1709	1741	rBV	3909484	6806940	61.41%	14.756%
8	14.378	1937	1945	1963	rBV	734885	1583894	14.29%	3.433%
9	15.895	2195	2203	2216	rBV	2947963	5123969	46.22%	11.107%
10	17.183	2415	2422	2432	rBV	892429	1890682	17.06%	4.098%
11	19.901	2878	2884	2896	rBV	7329965	11085277	100.00%	24.030%
12	21.630	3173	3178	3189	rBV2	1297746	2281900	20.58%	4.947%
13	24.995	3742	3750	3768	rVB	735137	2090721	18.86%	4.532%

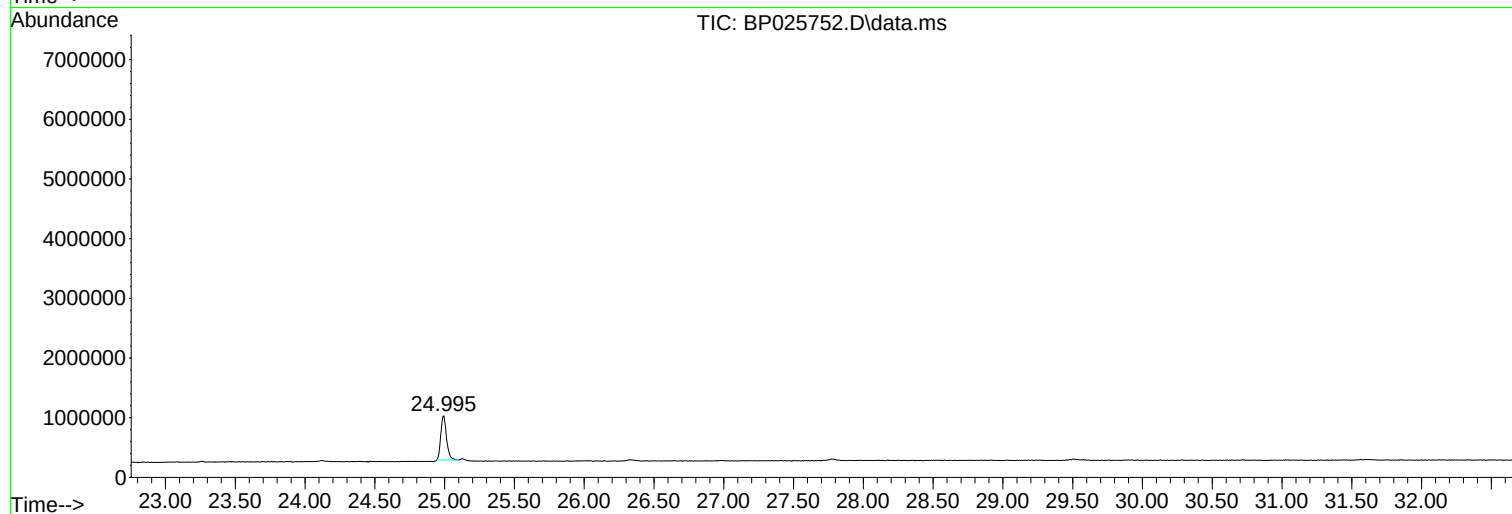
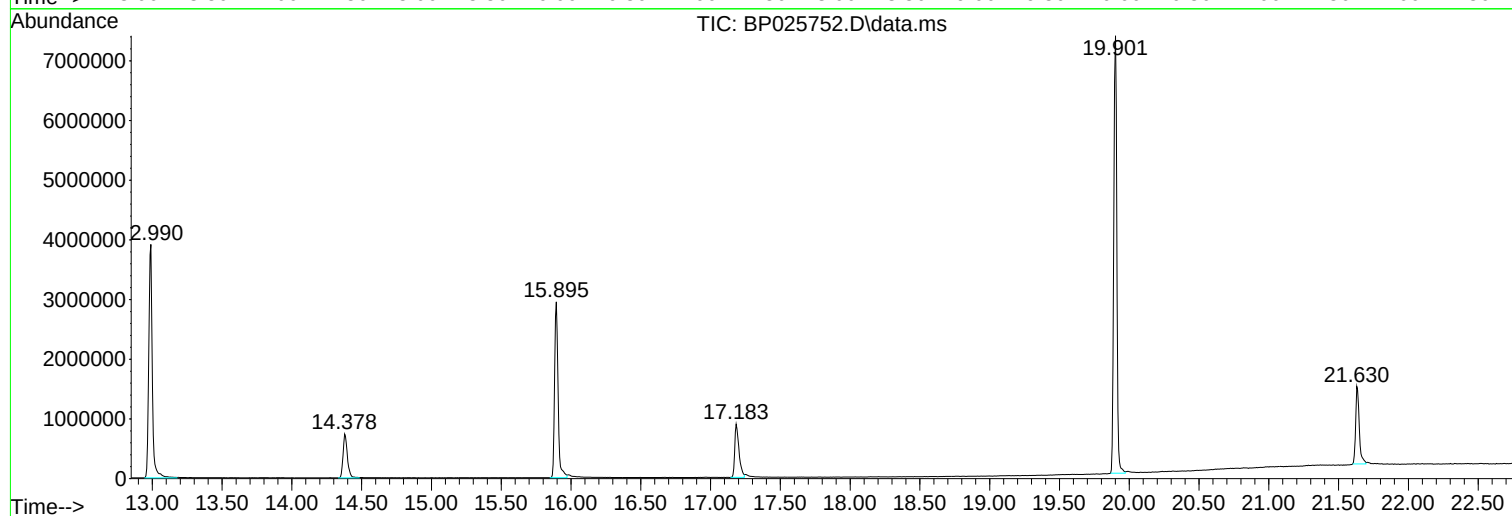
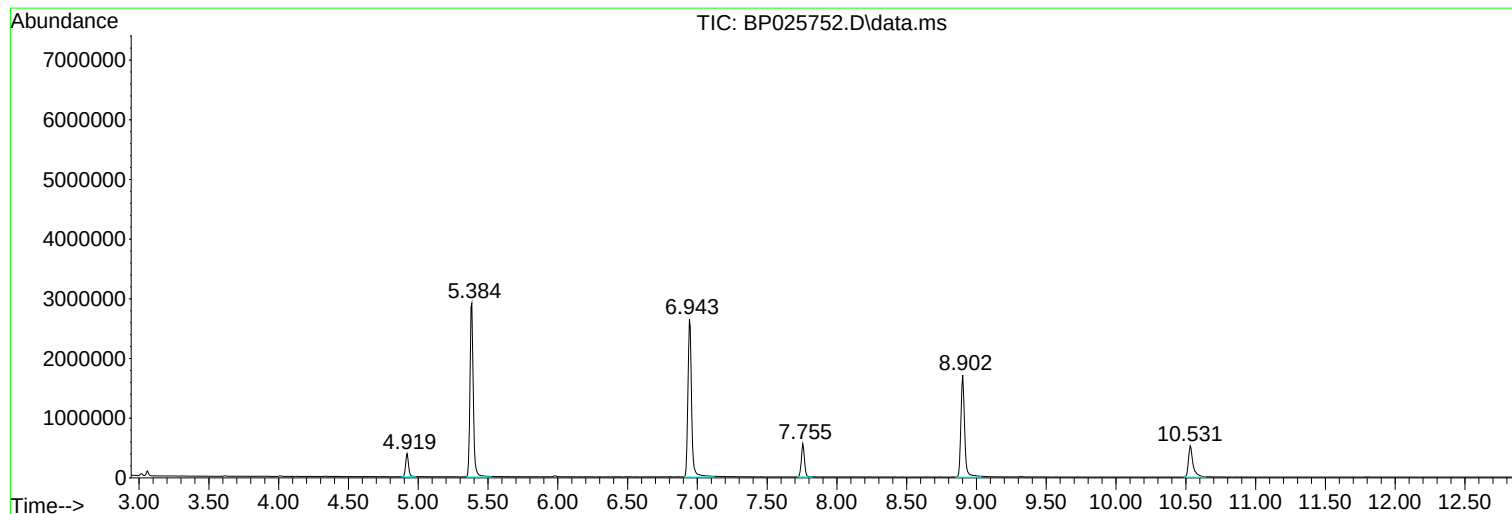
Sum of corrected areas: 46131485

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.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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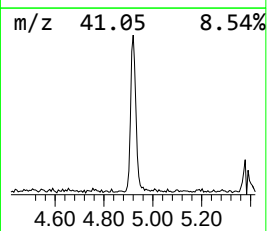
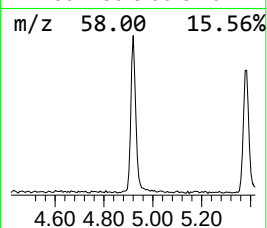
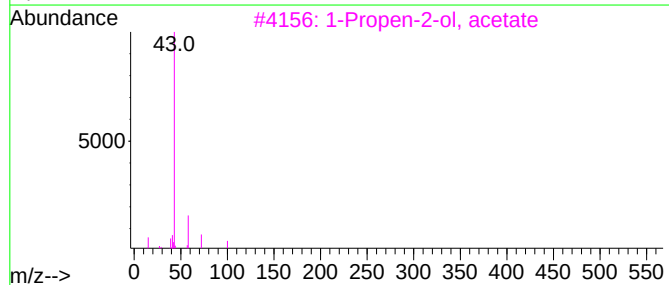
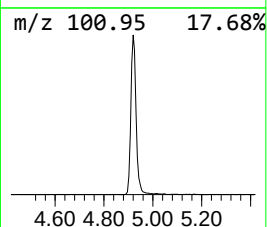
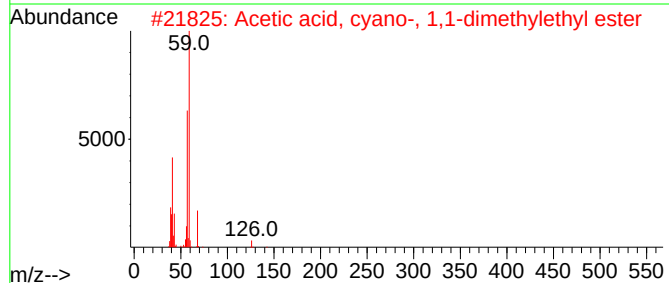
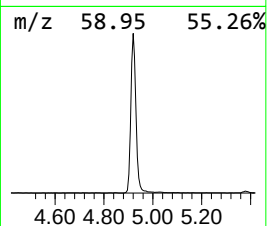
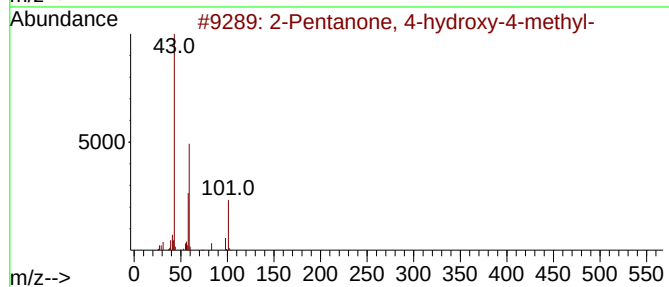
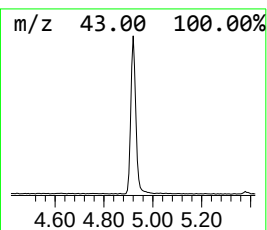
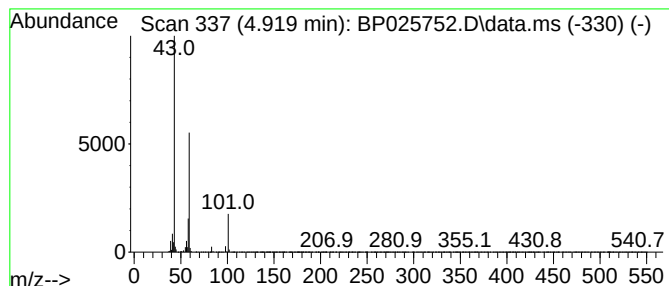
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.919	12.89 ng	565092	1,4-Dichlorobenzene-d4	7.755

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
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-...	4.919	12.9	ng	565092	1	7.755	876893	20.0