

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091225\
Data File : BP025743.D
Acq On : 12 Sep 2025 14:30
Operator : CG/JU
Sample : PB169499BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB169499BL

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-BP091225.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025743.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	332	337	345	rBV	362062	520563	4.69%	1.150%
2	5.384	409	416	437	rBV	2718157	4439313	39.98%	9.811%
3	6.943	674	681	702	rBV	2519358	4551720	40.99%	10.060%
4	7.754	812	819	829	rBV	490514	815027	7.34%	1.801%
5	8.901	1005	1014	1036	rBV	1548695	3005597	27.07%	6.643%
6	10.531	1279	1291	1310	rBV	461253	1106396	9.96%	2.445%
7	12.984	1701	1708	1734	rBV	4088197	6571167	59.18%	14.523%
8	14.378	1935	1945	1962	rBV	777101	1514496	13.64%	3.347%
9	15.889	2196	2202	2233	rBV	3146755	5228341	47.09%	11.555%
10	17.183	2414	2422	2440	rBV2	980261	1935444	17.43%	4.277%
11	19.907	2878	2885	2895	rBV	6710708	11103815	100.00%	24.540%
12	21.618	3170	3176	3189	rBV2	1137451	2313061	20.83%	5.112%
13	24.971	3739	3746	3765	rVB	689998	2142746	19.30%	4.736%

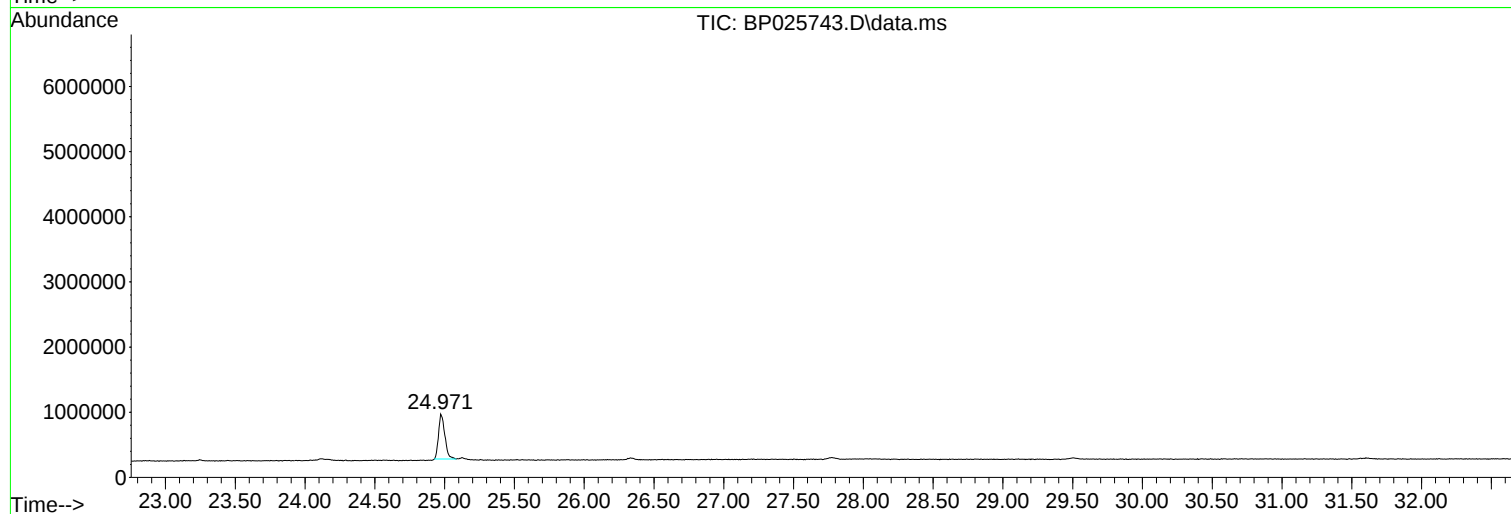
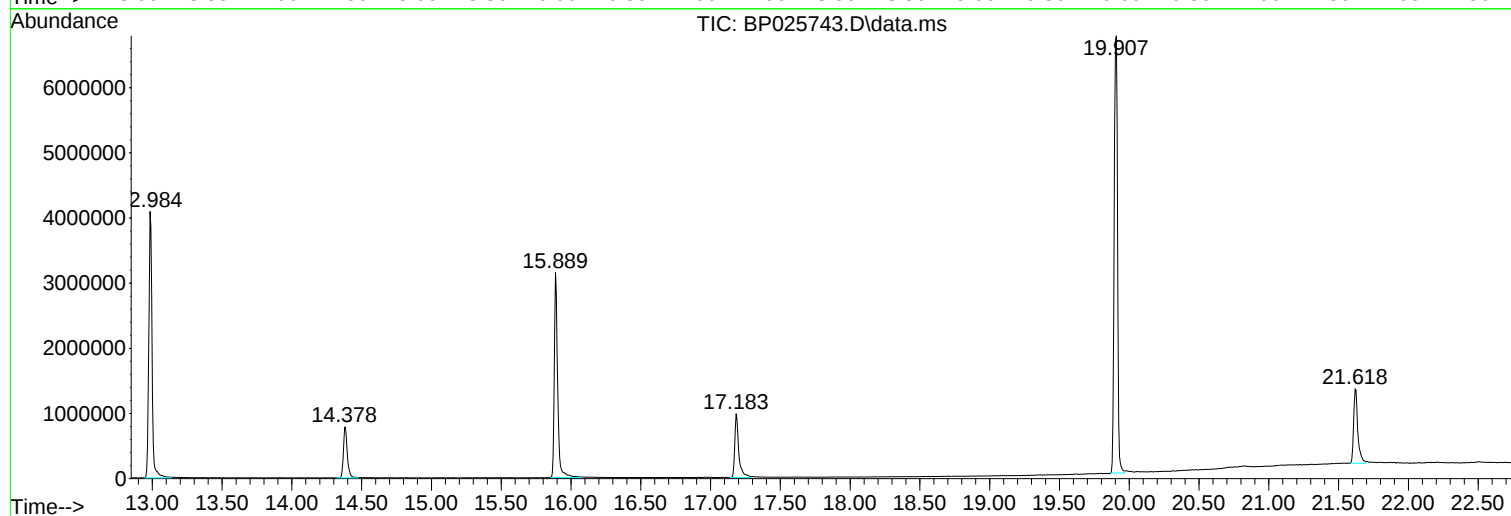
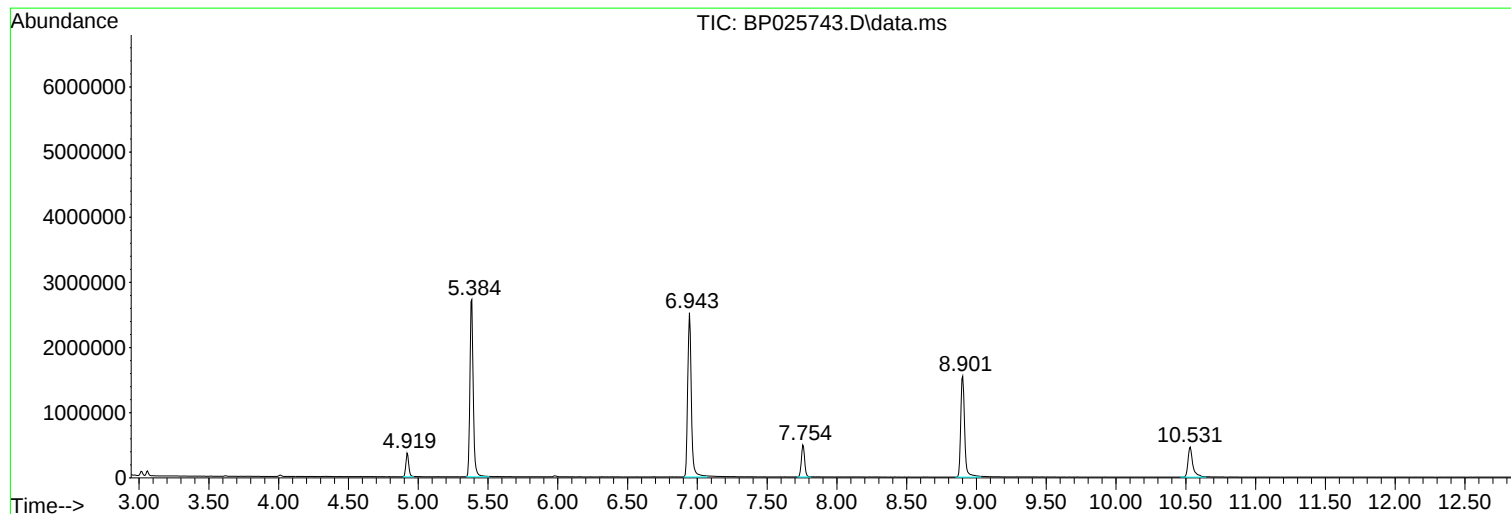
Sum of corrected areas: 45247686

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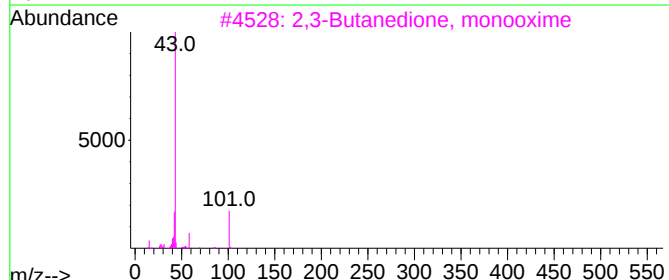
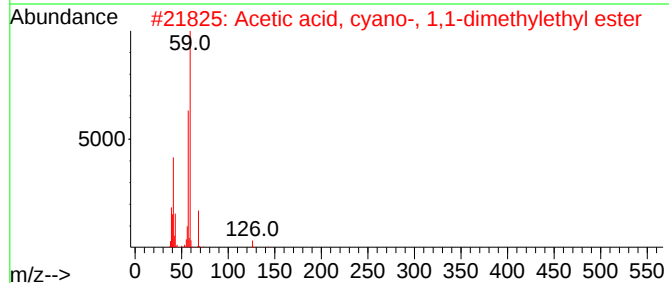
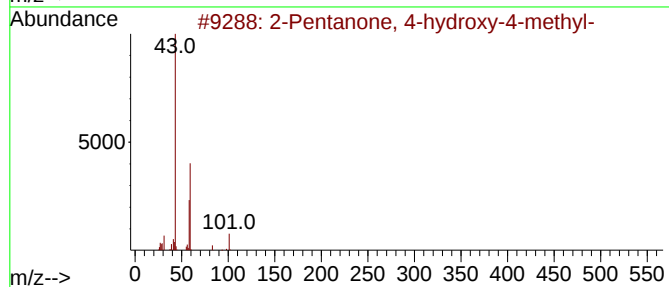
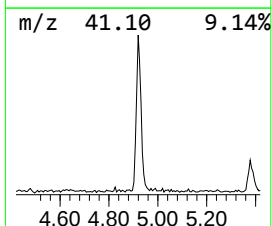
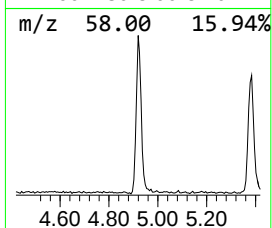
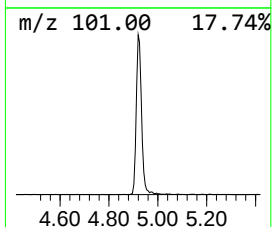
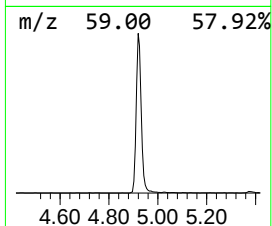
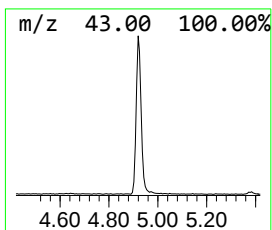
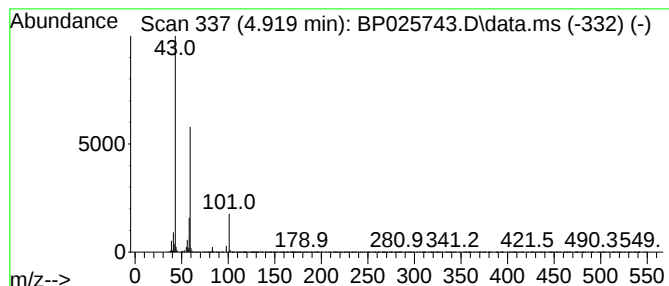
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.919	12.77 ng	520563	1,4-Dichlorobenzene-d4	7.754

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	23		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9		
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.919	12.8	ng	520563	1	7.754	815027	20.0