

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090525\
 Data File : BP025676.D
 Acq On : 05 Sep 2025 14:09
 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

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-BP081425.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025676.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.925	331	338	348	rBV	319866	457298	4.35%	1.106%
2	5.390	410	417	433	rBV	2669567	3982006	37.89%	9.627%
3	6.955	675	683	707	rBV	2333337	3998686	38.05%	9.667%
4	7.766	814	821	829	rBV	463425	768572	7.31%	1.858%
5	8.907	1007	1015	1037	rBV	1555329	2688600	25.58%	6.500%
6	10.537	1284	1292	1312	rBV	486872	991167	9.43%	2.396%
7	12.996	1701	1710	1735	rBV	3365407	5878828	55.94%	14.212%
8	14.395	1941	1948	1968	rBV	661281	1305871	12.43%	3.157%
9	15.913	2198	2206	2221	rBV	2823868	4994187	47.52%	12.074%
10	17.201	2418	2425	2446	rBV	940384	1680404	15.99%	4.062%
11	19.919	2880	2887	2899	rBV	6905492	10508715	100.00%	25.405%
12	21.648	3175	3181	3193	rVB2	1045018	2015691	19.18%	4.873%
13	25.018	3746	3754	3776	rVB	640691	2094440	19.93%	5.063%

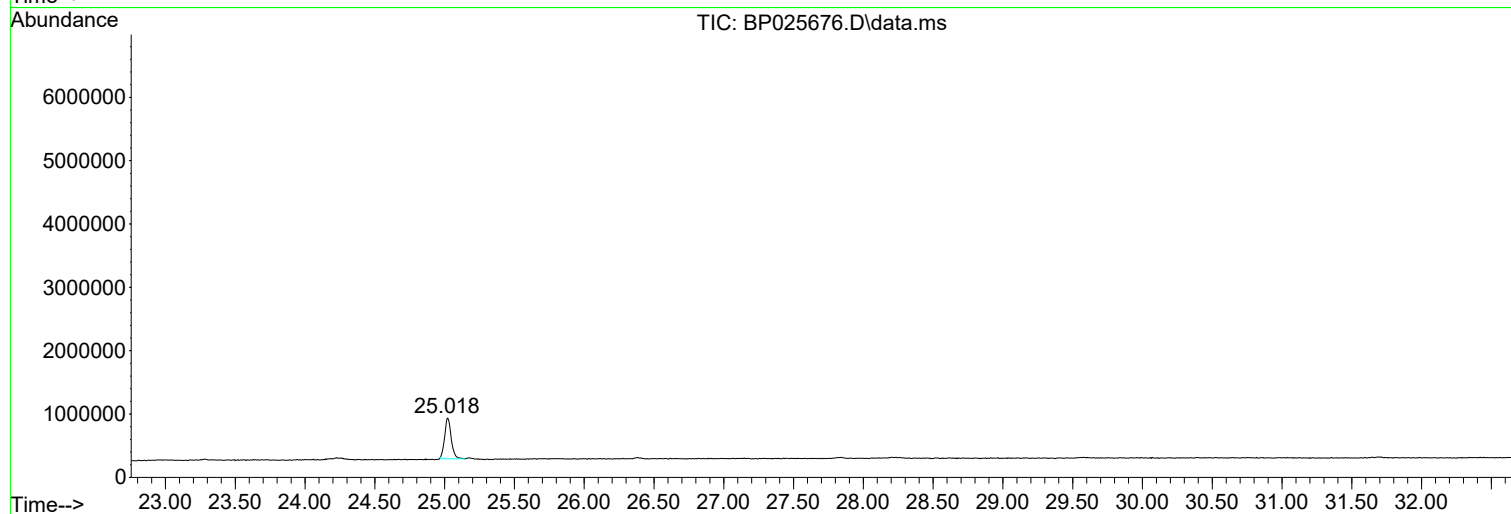
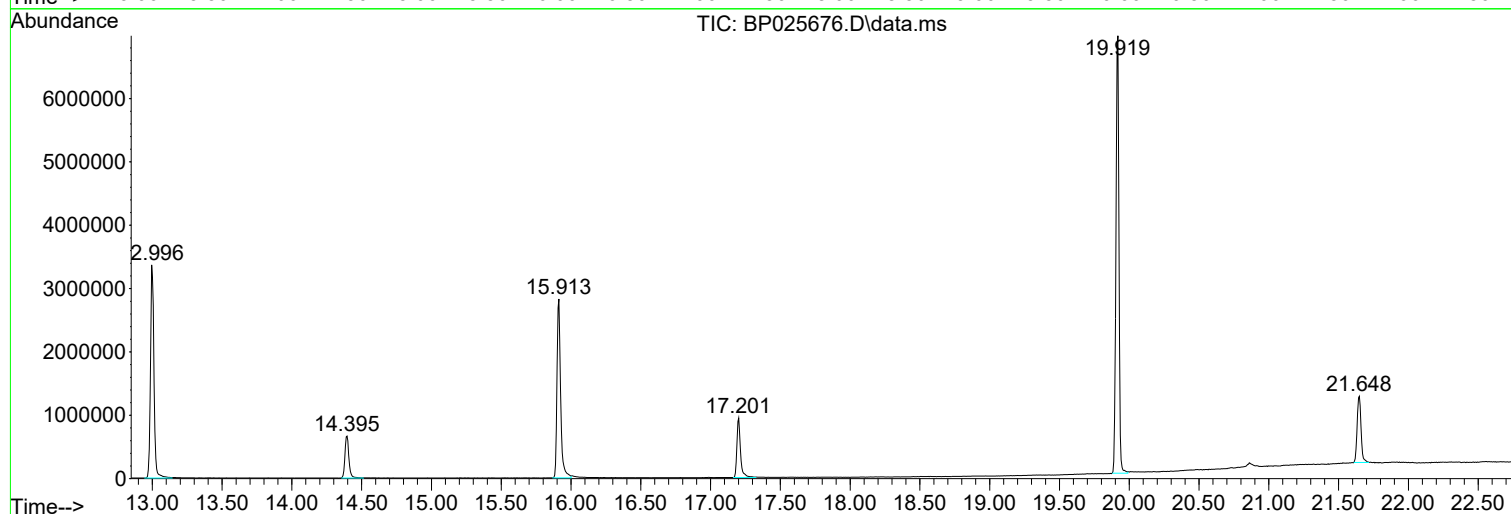
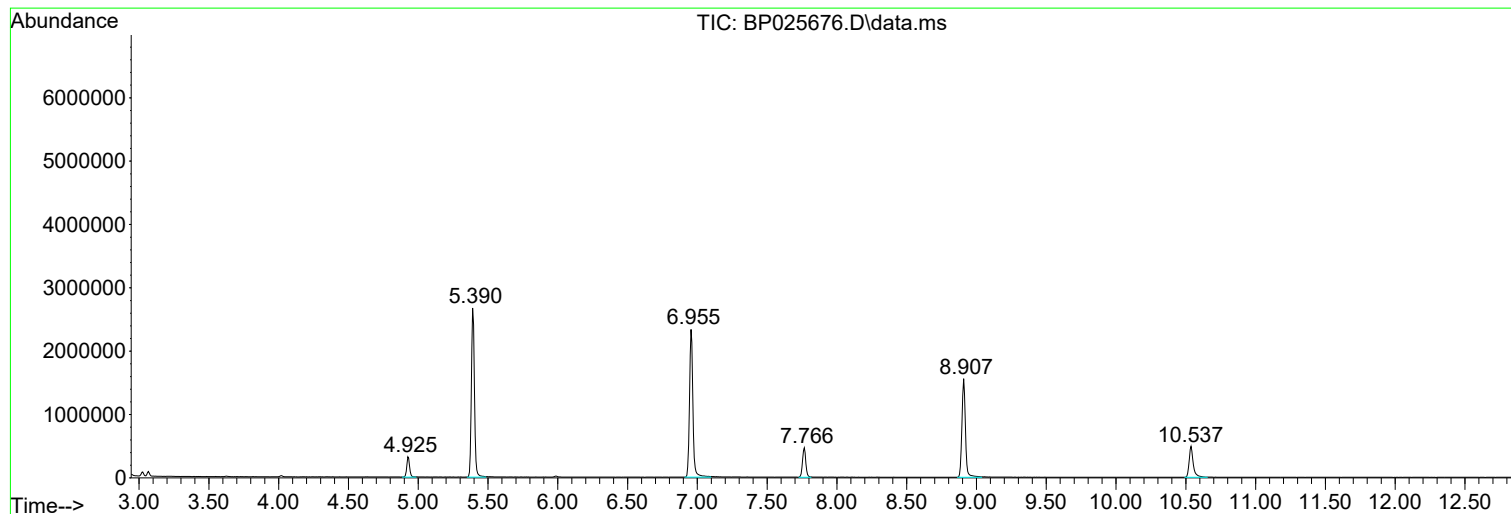
Sum of corrected areas: 41364465

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.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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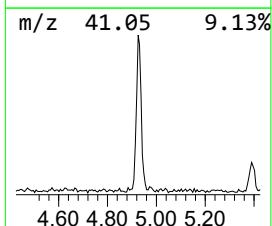
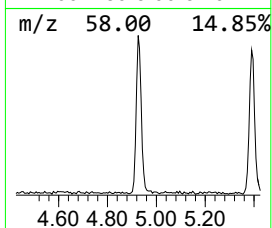
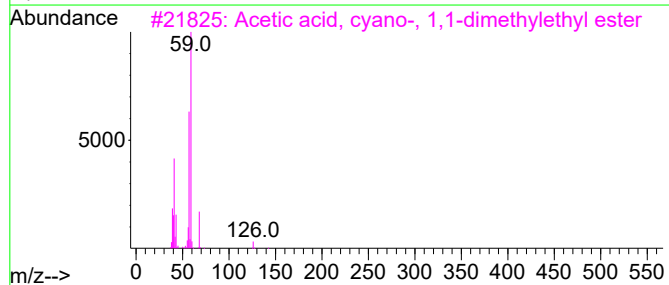
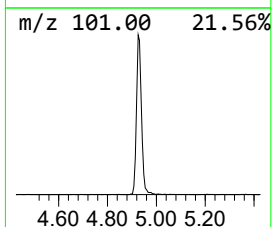
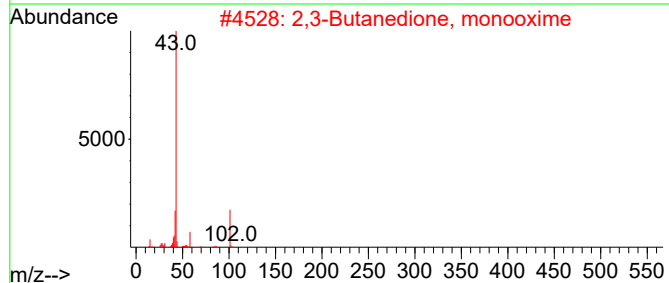
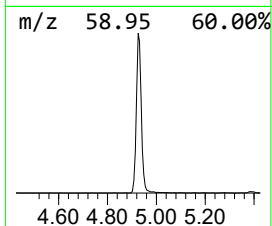
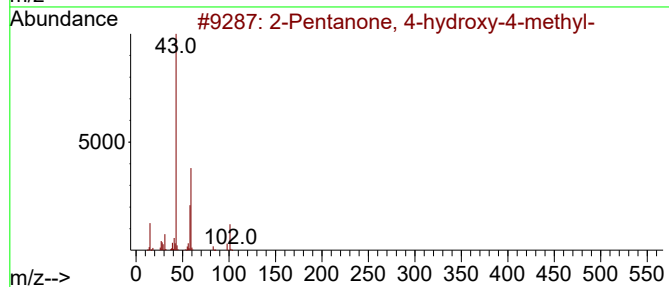
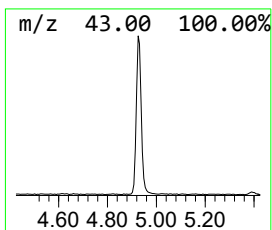
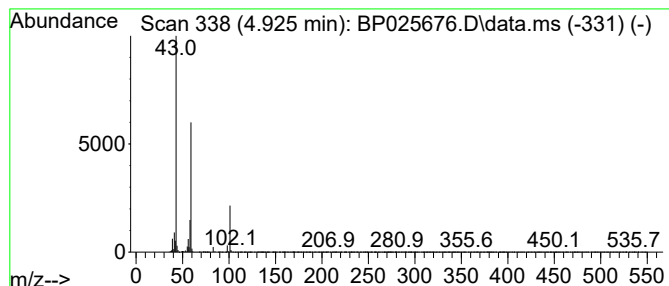
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.925	11.90 ng	457298	1,4-Dichlorobenzene-d4	7.766	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
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	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-...	4.925	11.9	ng	457298	1	7.766	768572	20.0