

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091225\
 Data File : BP025735.D
 Acq On : 11 Sep 2025 17:35
 Operator : CG/JU
 Sample : PB169633BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169633BL

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: BP025735.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.920	331	337	350	rBV	419082	600163	5.33%	1.302%
2	5.384	409	416	439	rBV	2857838	4551423	40.42%	9.873%
3	6.943	674	681	711	rBV	2695066	4700145	41.74%	10.195%
4	7.755	812	819	831	rBV	513330	818910	7.27%	1.776%
5	8.896	1005	1013	1038	rBV	1623188	3116719	27.68%	6.761%
6	10.531	1283	1291	1310	rBV	483128	1088266	9.66%	2.361%
7	12.996	1702	1710	1734	rBV	3909498	6816043	60.53%	14.785%
8	14.390	1939	1947	1967	rBV	794280	1495742	13.28%	3.244%
9	15.907	2196	2205	2227	rBV	2994883	5416612	48.10%	11.749%
10	17.195	2417	2424	2435	rBV	959535	1838093	16.32%	3.987%
11	19.907	2878	2885	2897	rBV	7158587	11260811	100.00%	24.426%
12	21.636	3173	3179	3192	rBV	1182419	2195110	19.49%	4.762%
13	25.001	3743	3751	3771	rVB2	714379	2203107	19.56%	4.779%

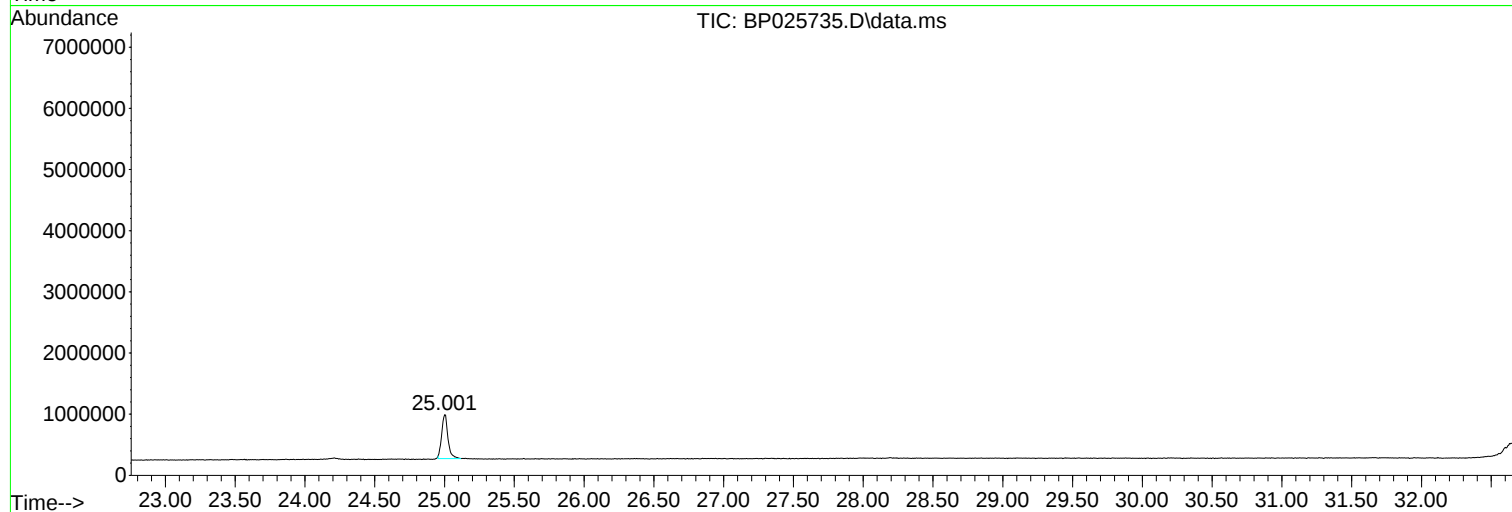
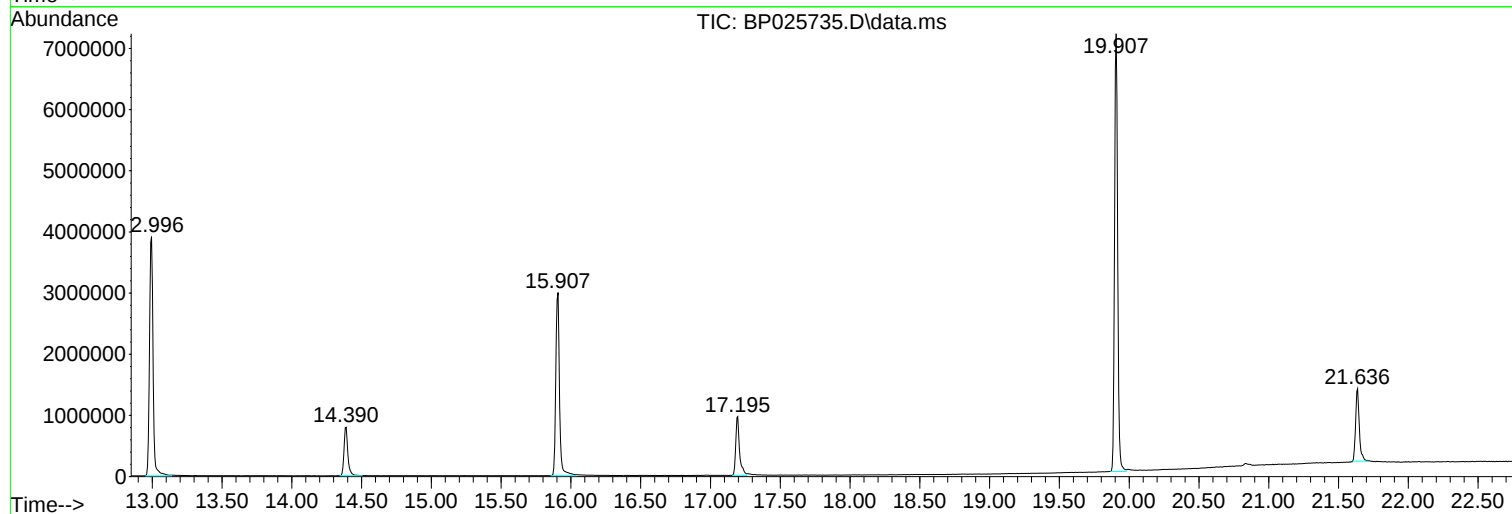
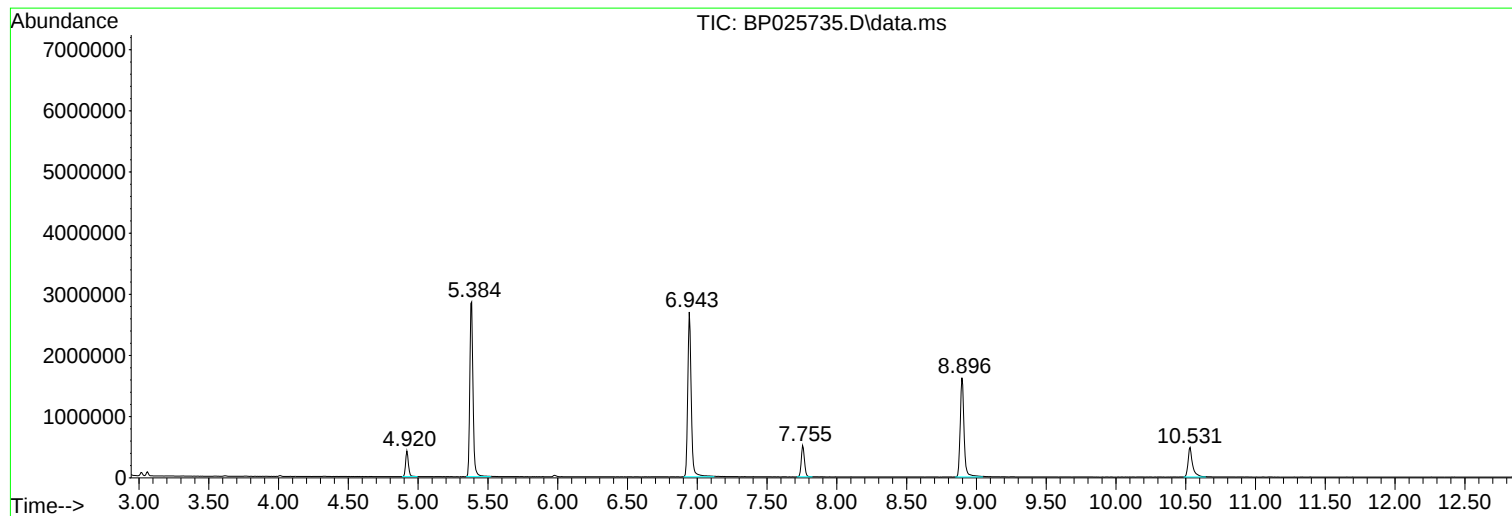
Sum of corrected areas: 46101144

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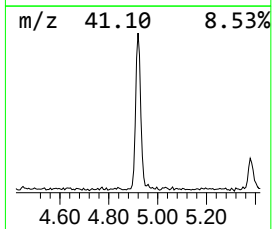
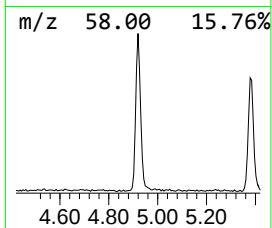
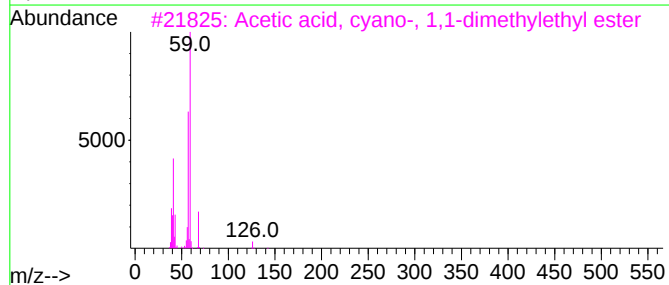
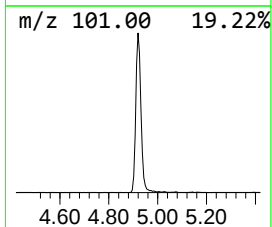
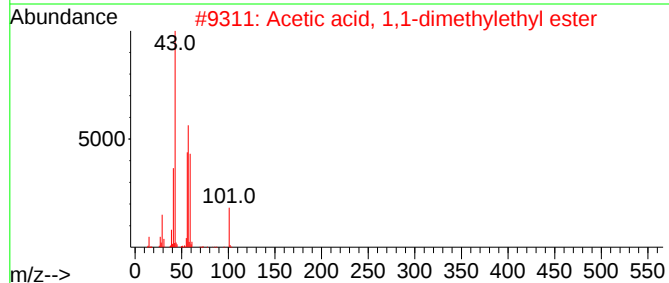
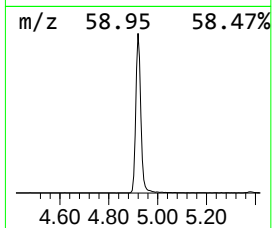
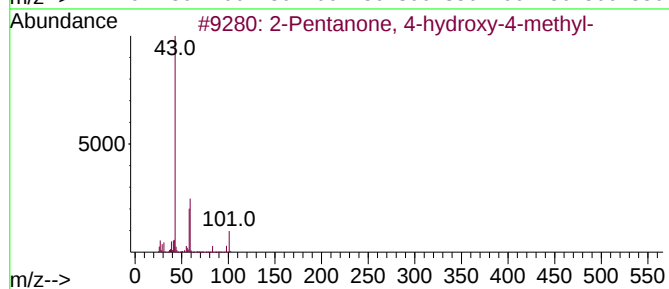
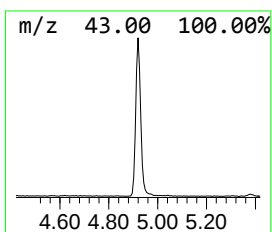
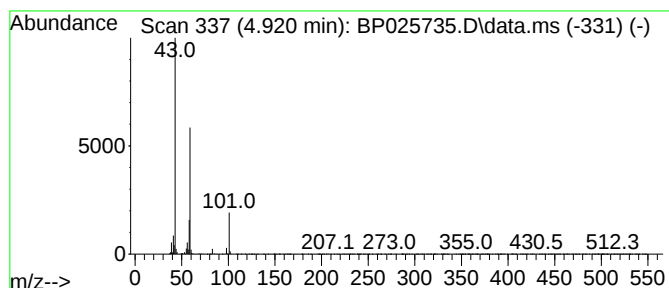
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.920	14.66 ng	600163	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	64		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39		
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-...	4.920	14.7	ng	600163	1	7.755	818910	20.0