

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050421.D
 Acq On : 14 Jul 2025 11:30
 Operator : RC/JU
 Sample : PB168813BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168813BL

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_M\Methods\8270-BM070925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM050421.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.975	333	338	348	rVB	639578	887664	4.21%	0.900%
2	5.440	410	417	433	rBV	7779929	11073179	52.50%	11.225%
3	7.022	678	686	699	rBV	6883203	11073300	52.50%	11.225%
4	7.863	821	829	836	rBV	1551498	2404715	11.40%	2.438%
5	9.016	1016	1025	1037	rBV	4314705	7059162	33.47%	7.156%
6	10.657	1296	1304	1312	rBV	1748670	2976543	14.11%	3.017%
7	13.116	1714	1722	1737	rBV	11131546	15962949	75.68%	16.182%
8	14.486	1948	1955	1971	rBV	2495561	3613608	17.13%	3.663%
9	15.968	2200	2207	2224	rBV	7135189	10198477	48.35%	10.339%
10	17.221	2413	2420	2429	rBV2	2683723	3836931	18.19%	3.890%
11	19.833	2858	2864	2874	rBV	18070674	21092470	100.00%	21.382%
12	21.439	3132	3137	3148	rBV	2697882	4058513	19.24%	4.114%
13	24.474	3645	3653	3671	rVB	1643745	4408002	20.90%	4.469%

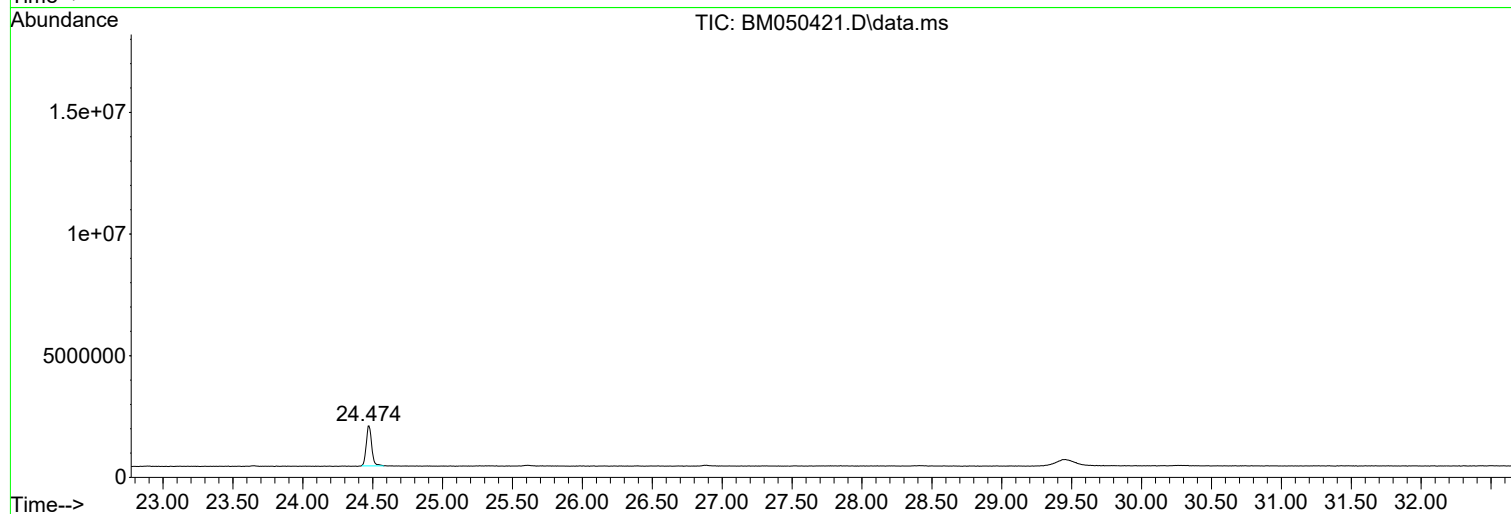
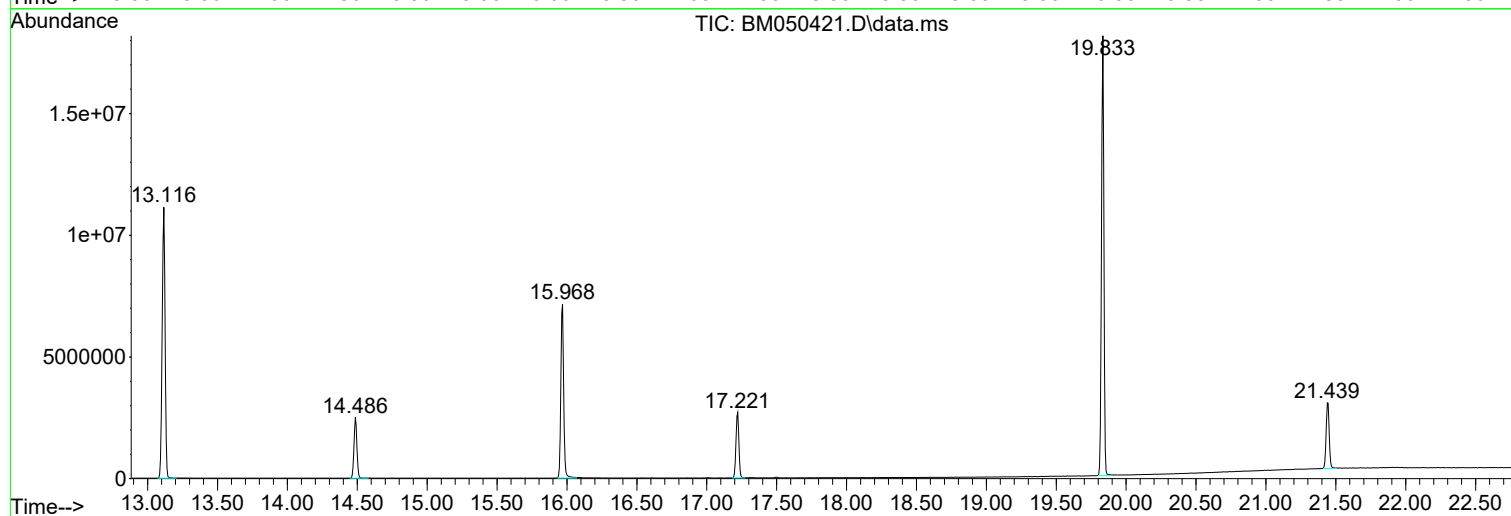
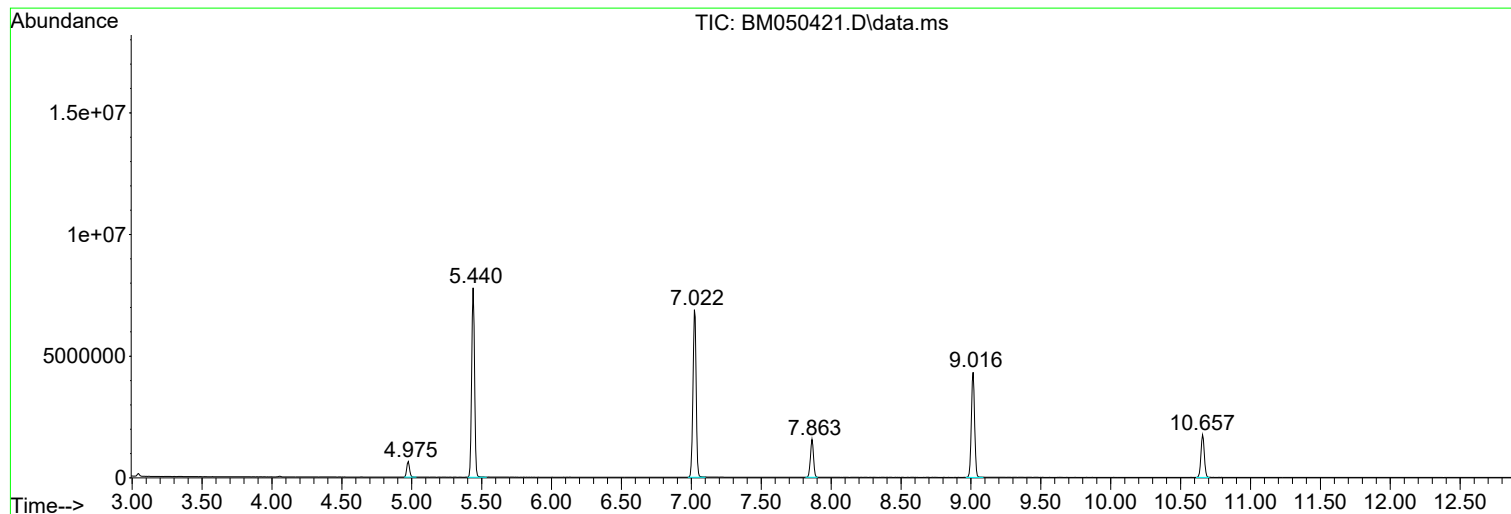
Sum of corrected areas: 98645513

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.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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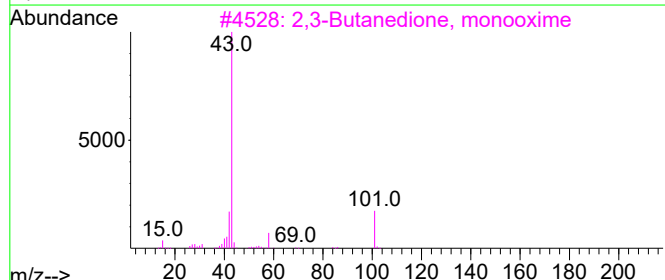
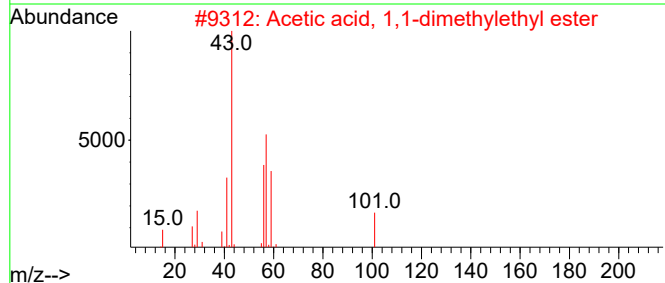
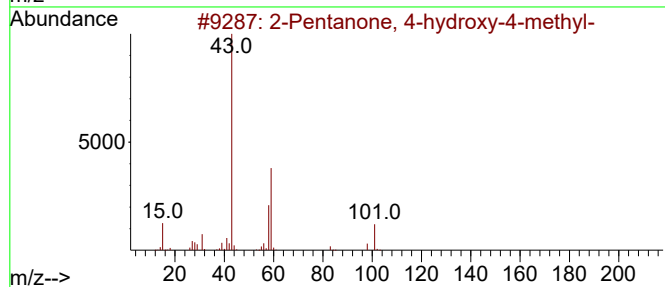
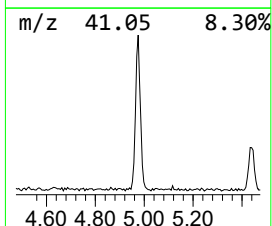
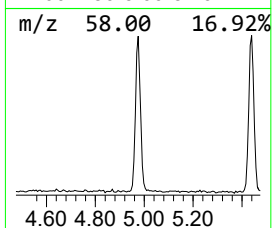
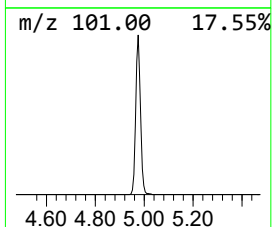
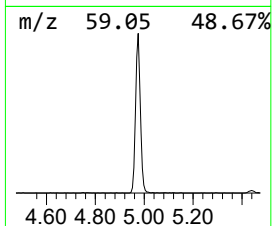
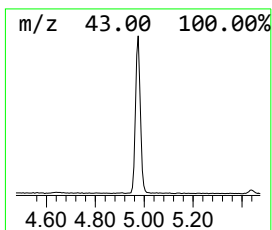
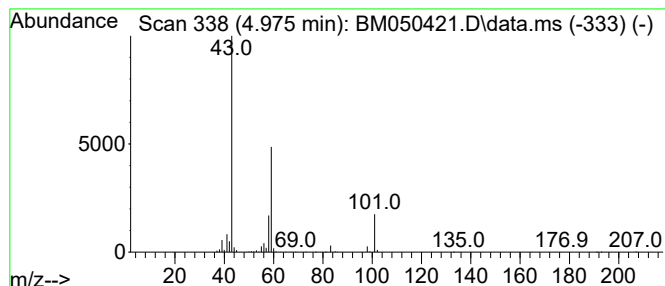
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.975	7.38 ng	887664	1,4-Dichlorobenzene-d4			7.863
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	59
2	Acetic acid, 1,1-dimethylethyl e...		116	C6H12O2	000540-88-5	38
3	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	16
4	Acetone		58	C3H6O	000067-64-1	12
5	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.975	7.4	ng	887664	1	7.863	2404720	20.0