

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090325\
 Data File : BG064251.D
 Acq On : 3 Sep 2025 13:02
 Operator : RC/JU
 Sample : PB169490BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB169490BL

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_G\Methods\8270-BG082825.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG064251.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.974	315	321	329	rVB	101669	154380	4.72%	1.118%
2	5.444	393	401	414	rBV	802933	1258337	38.48%	9.109%
3	7.025	662	670	682	rBV	876093	1447787	44.27%	10.480%
4	7.853	803	811	819	rBV	143595	236062	7.22%	1.709%
5	9.005	997	1007	1017	rBV	531554	938963	28.71%	6.797%
6	10.644	1278	1286	1293	rBV	190134	330465	10.10%	2.392%
7	13.106	1698	1705	1716	rVB	1377465	2124859	64.97%	15.382%
8	14.481	1930	1939	1946	rBV	300766	471404	14.41%	3.412%
9	15.967	2185	2192	2206	rBV	1087948	1675447	51.23%	12.128%
10	17.225	2398	2406	2416	rBV	369255	563812	17.24%	4.081%
11	19.845	2845	2852	2858	rBV	2388665	3270515	100.00%	23.675%
12	21.449	3118	3125	3135	rBV	420917	668322	20.43%	4.838%
13	24.457	3628	3637	3651	rBV2	240597	674023	20.61%	4.879%

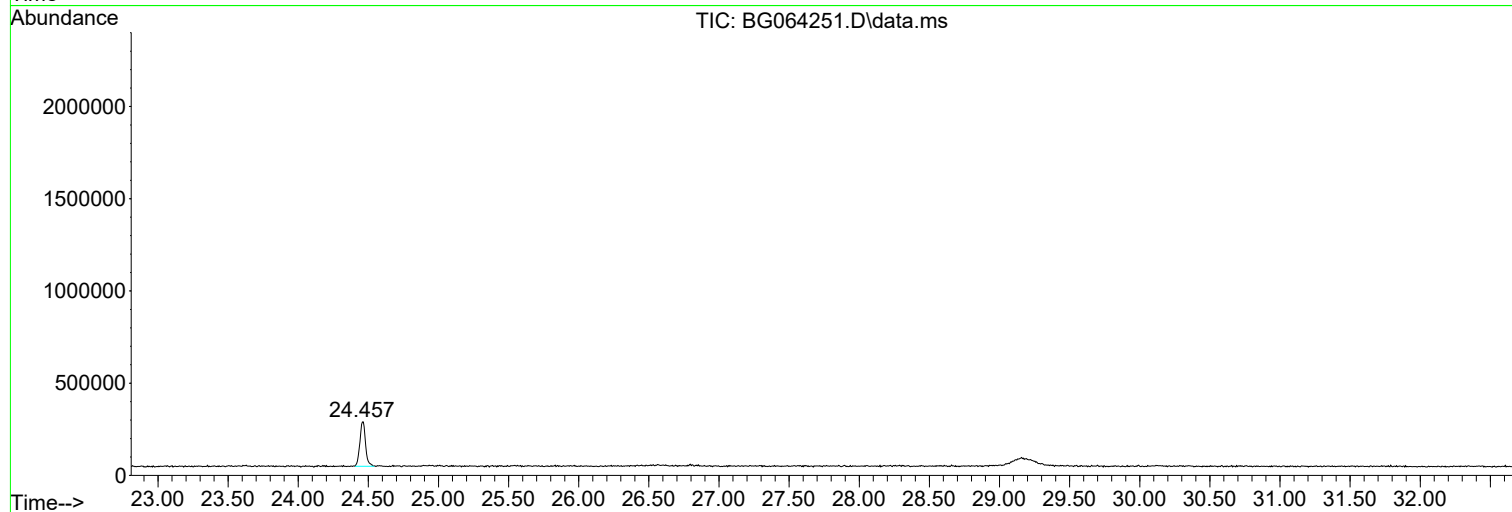
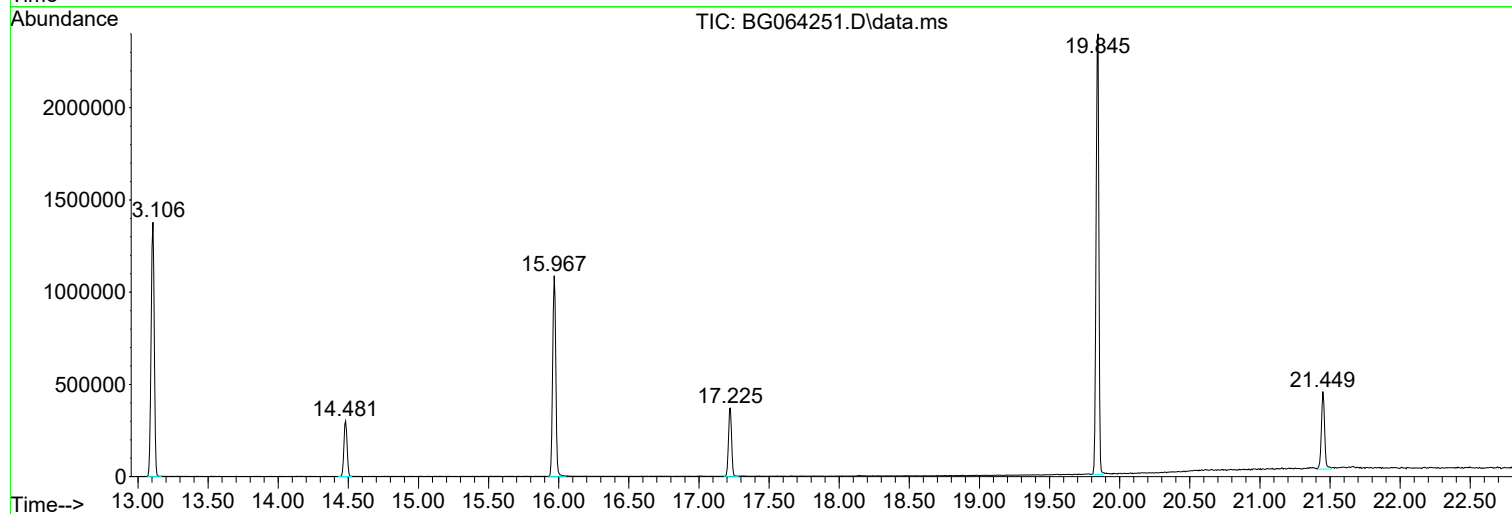
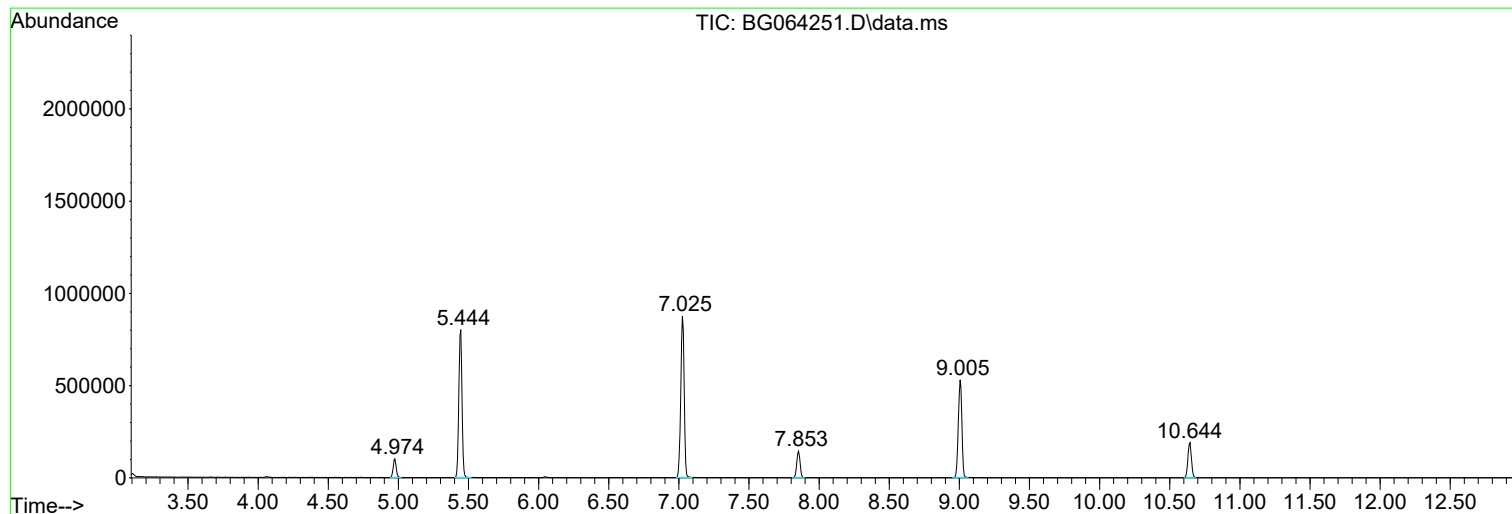
Sum of corrected areas: 13814376

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.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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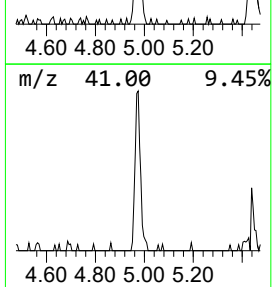
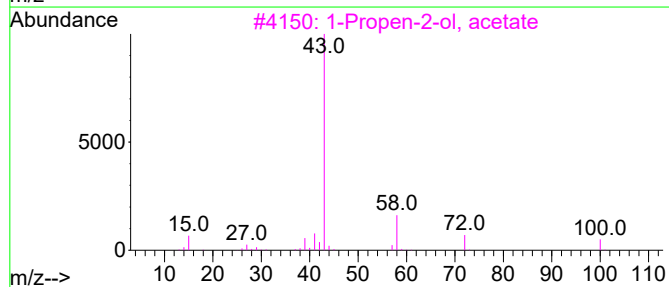
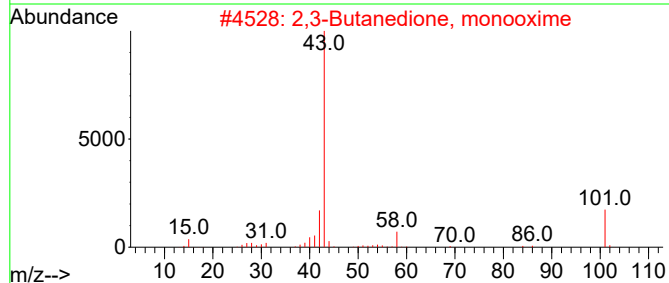
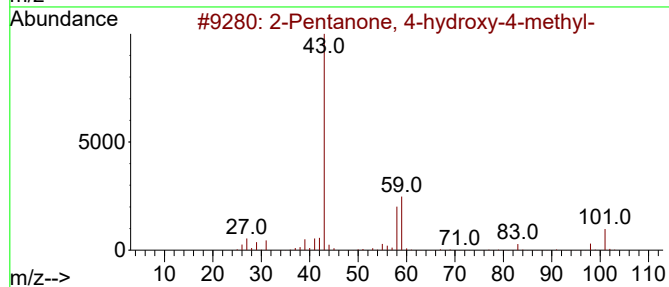
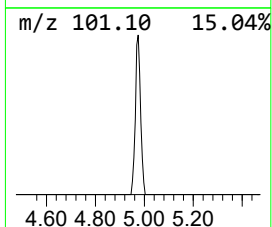
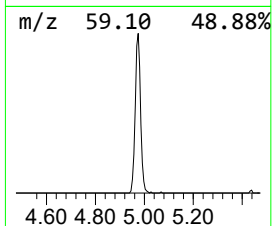
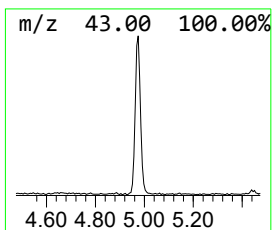
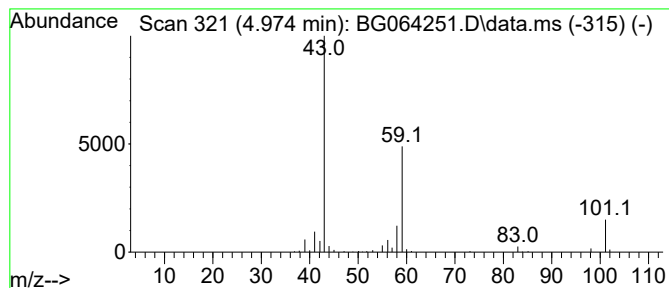
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.974	13.08 ng	154380	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
4	Acetamide, N-(aminocarbonyl)-	102	C3H6N2O2	000591-07-1	9		
5	5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.974	13.1	ng	154380	1	7.853	236062	20.0