

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060925\
 Data File : BP024872.D
 Acq On : 09 Jun 2025 11:24
 Operator : RC/JU
 Sample : PB168323BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168323BL

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-BP060625.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP024872.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.778	308	313	326	rBV	510577	718345	4.10%	0.915%
2	5.243	385	392	419	rBV	5252634	8227174	46.92%	10.484%
3	6.814	652	659	680	rBV	5181034	8289401	47.27%	10.563%
4	7.608	787	794	814	rBV	1050014	1608526	9.17%	2.050%
5	8.760	980	990	1017	rBV	3138266	5402410	30.81%	6.884%
6	10.378	1258	1265	1287	rBV	1117329	2188080	12.48%	2.788%
7	12.860	1680	1687	1700	rBV	8116404	11956272	68.18%	15.236%
8	14.248	1915	1923	1940	rBV	1870333	2816175	16.06%	3.589%
9	15.778	2175	2183	2203	rBV	6677976	9467932	53.99%	12.065%
10	17.066	2395	2402	2418	rBV2	1929882	3204288	18.27%	4.083%
11	19.795	2859	2866	2877	rBV	13858273	17536338	100.00%	22.347%
12	21.489	3148	3154	3165	rBV2	1767894	3429892	19.56%	4.371%
13	24.730	3697	3705	3722	rVB	1140049	3628218	20.69%	4.624%

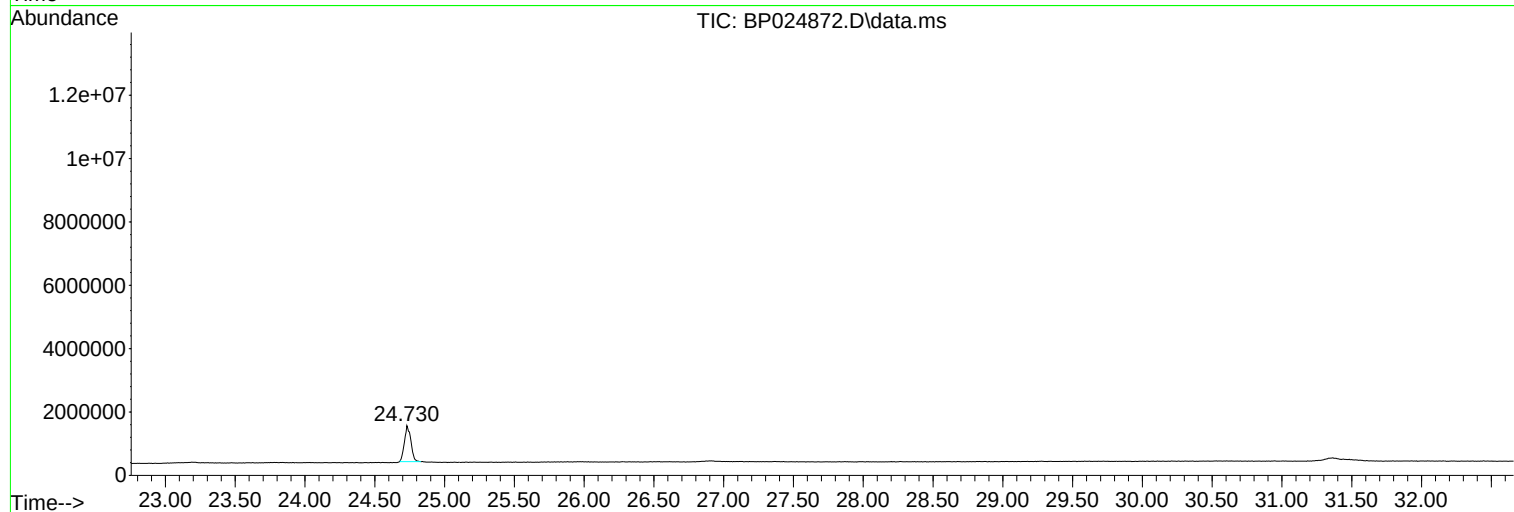
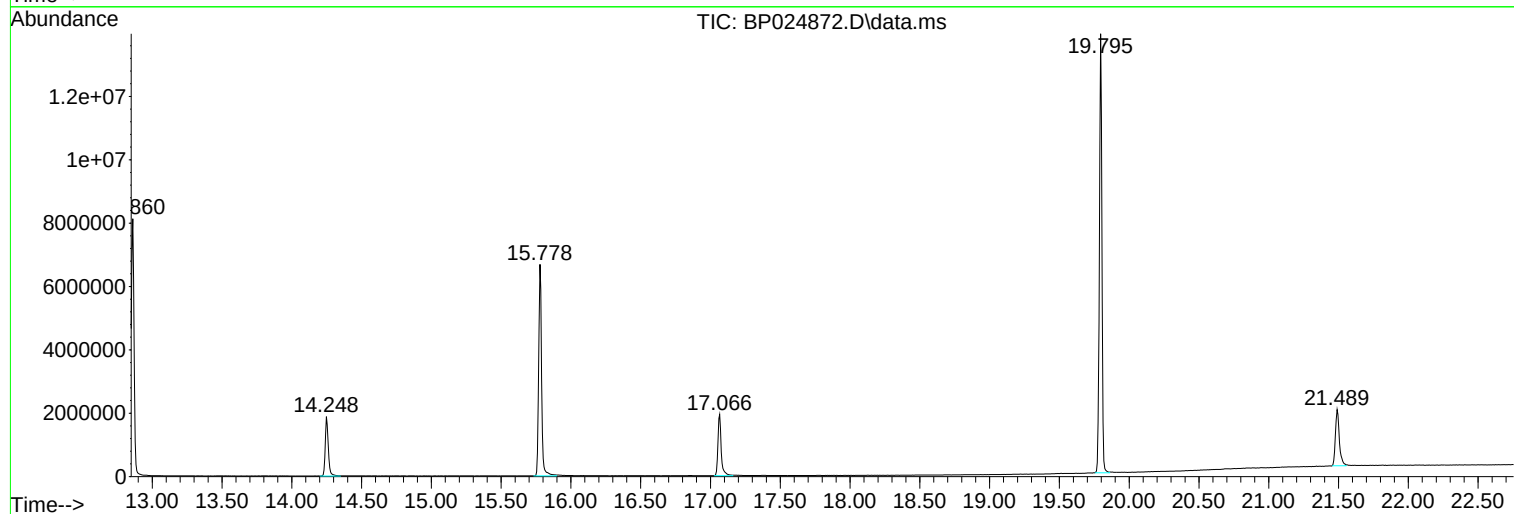
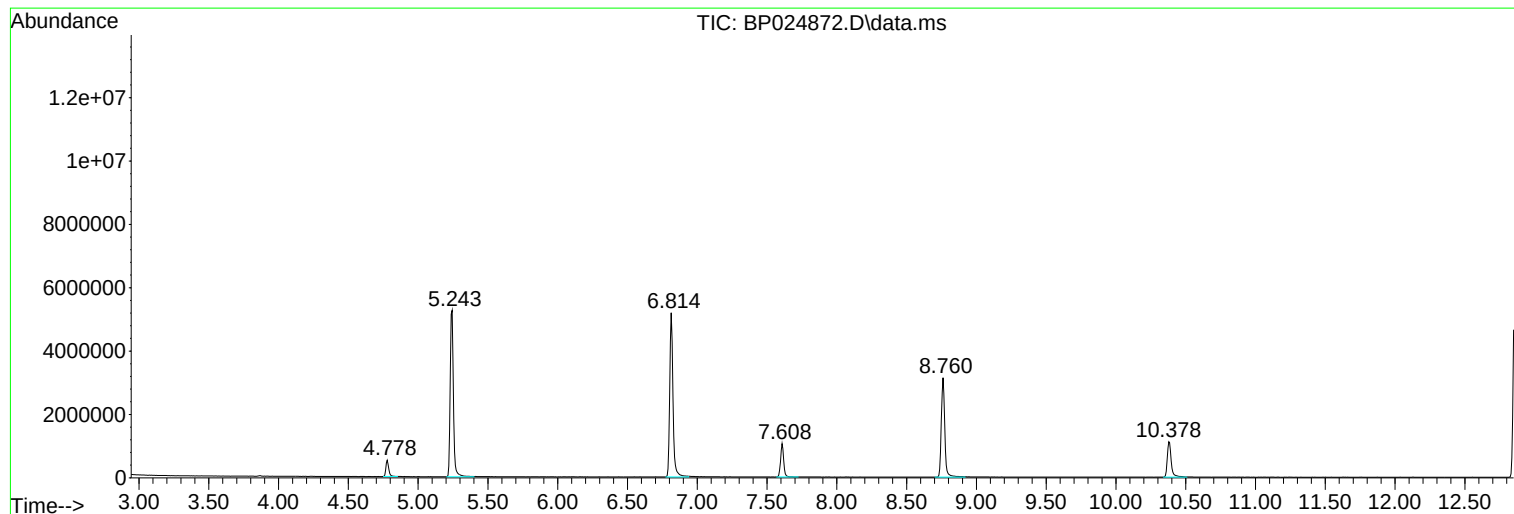
Sum of corrected areas: 78473051

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.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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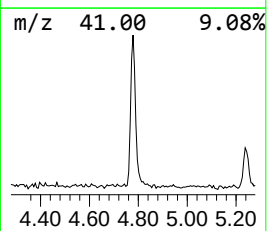
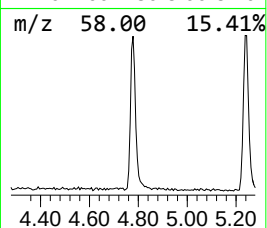
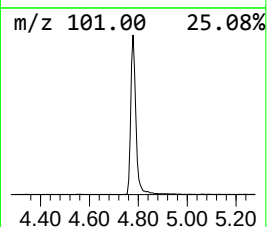
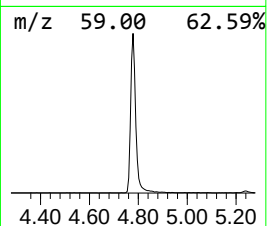
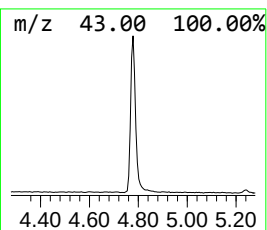
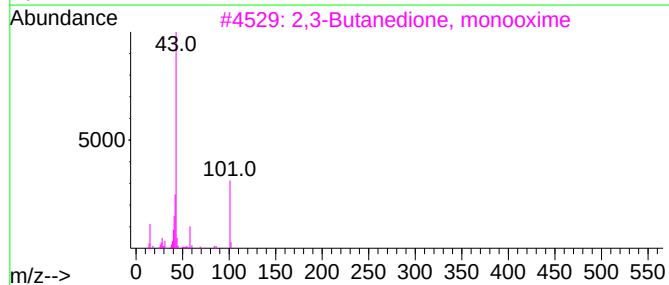
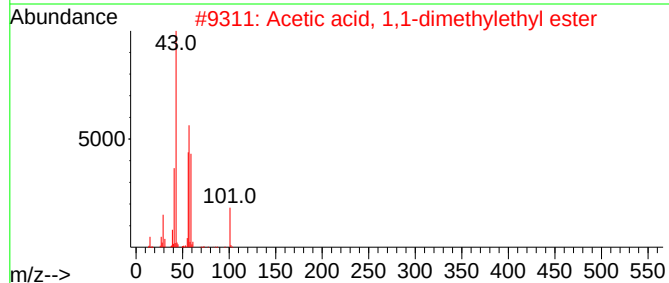
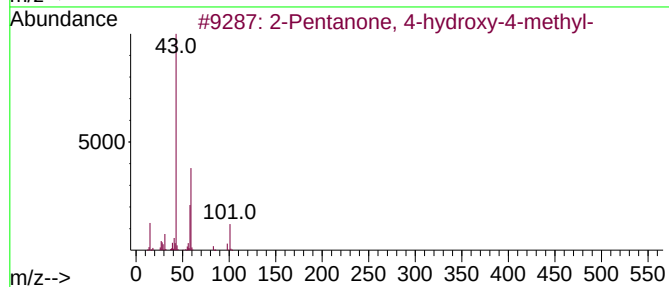
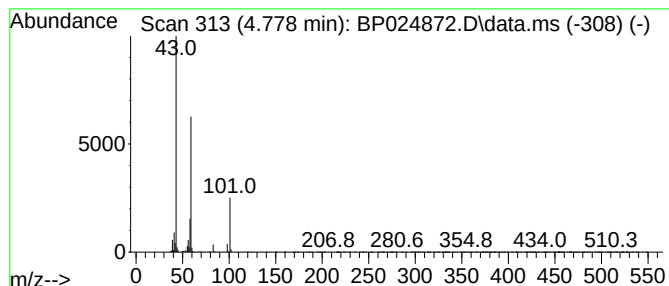
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.778	8.93 ng	718345	1,4-Dichlorobenzene-d4	7.608

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17		
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
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-...	4.778	8.9	ng	718345	1	7.608	1608530	20.0