

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052925\
 Data File : BP024829.D
 Acq On : 29 May 2025 14:41
 Operator : RC/JU
 Sample : PB168198BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168198BL

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-BP051325.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP024829.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.817	289	294	307	rBV	162104	242988	2.75%	0.722%
2	5.281	366	373	392	rBV	1919514	3004193	34.03%	8.931%
3	6.846	632	639	673	rBV	1593367	2962744	33.56%	8.808%
4	7.646	768	775	788	rBV	399741	660393	7.48%	1.963%
5	8.799	963	971	991	rBV	1022107	1920676	21.76%	5.710%
6	10.416	1238	1246	1267	rBV	450437	871249	9.87%	2.590%
7	12.887	1658	1666	1689	rBV	2765311	4621756	52.35%	13.739%
8	14.281	1893	1903	1923	rBV	723385	1121962	12.71%	3.335%
9	15.804	2154	2162	2183	rBV	2120834	4300631	48.72%	12.785%
10	17.087	2371	2380	2401	rBV	766245	1410024	15.97%	4.192%
11	19.804	2835	2842	2863	rBV	6469403	8828015	100.00%	26.244%
12	21.510	3126	3132	3148	rBV	1008877	1832262	20.76%	5.447%
13	24.798	3681	3691	3702	rBV	543173	1861761	21.09%	5.535%

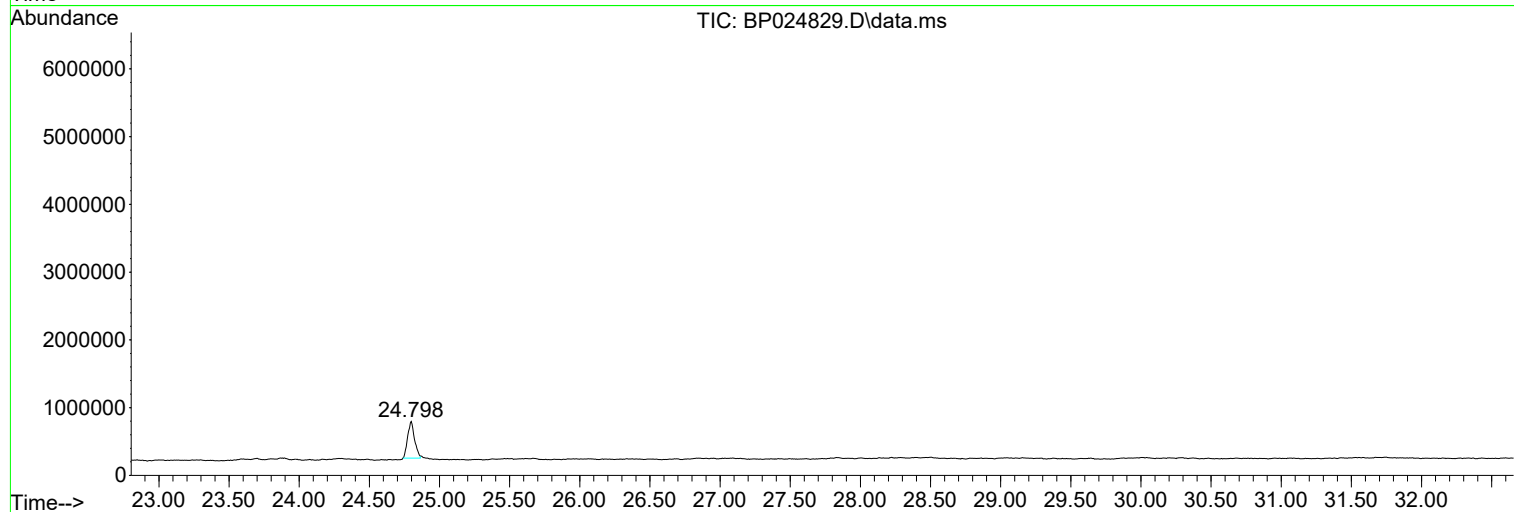
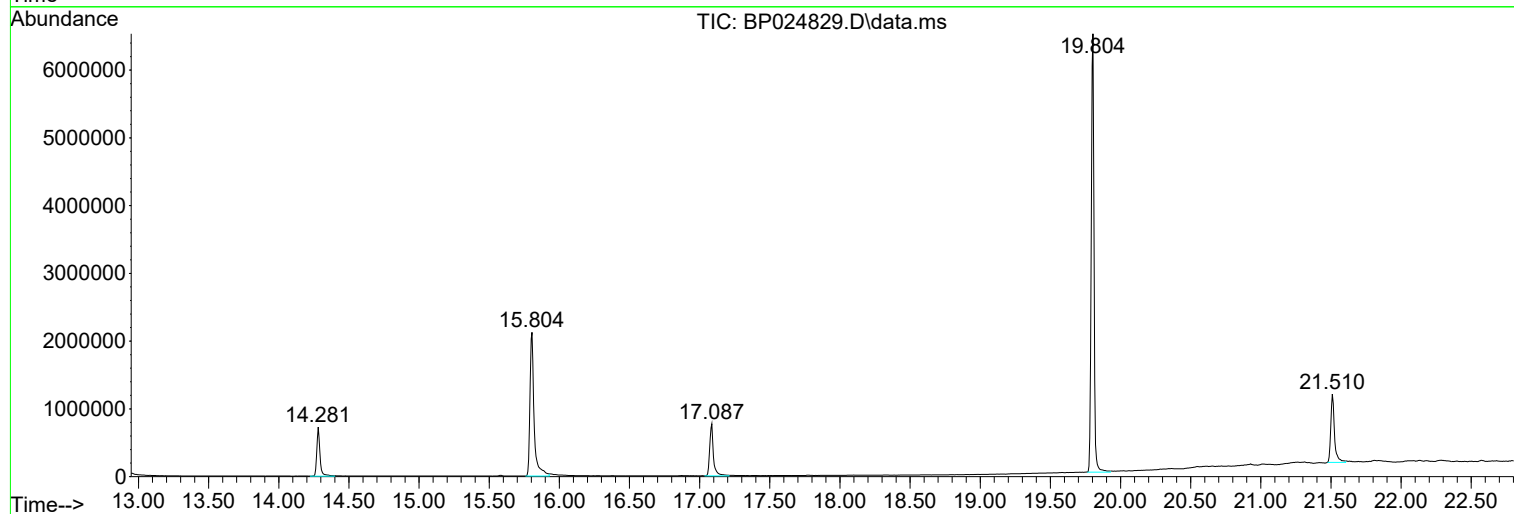
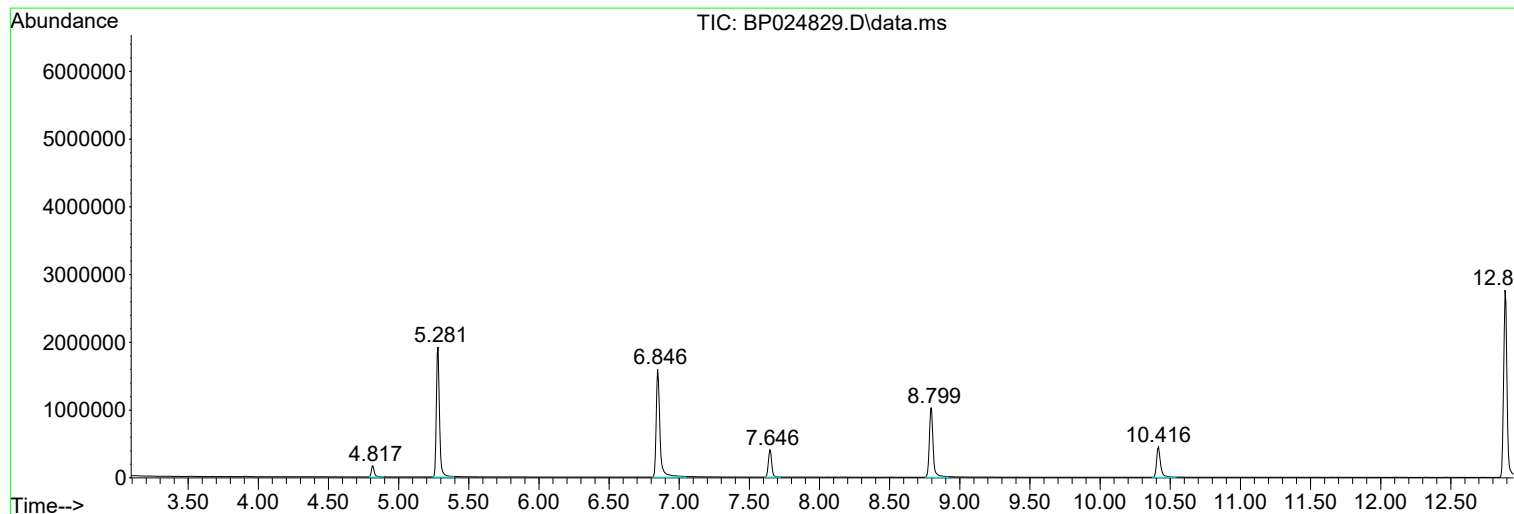
Sum of corrected areas: 33638654

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.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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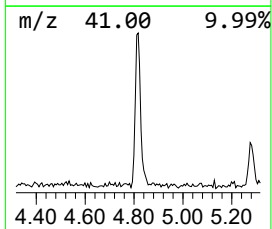
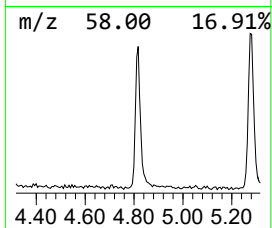
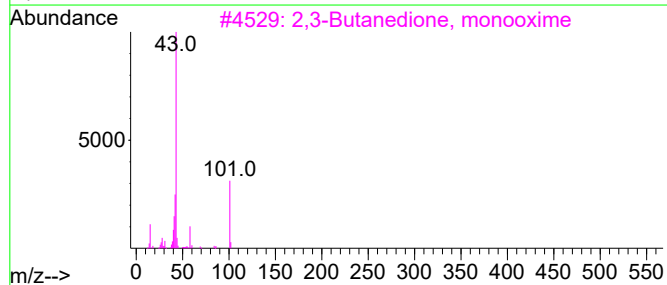
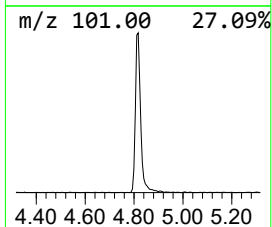
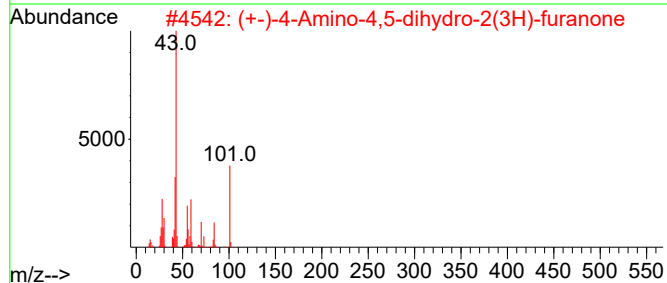
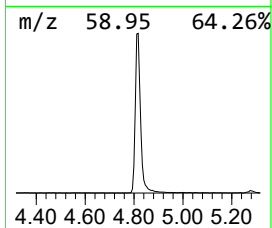
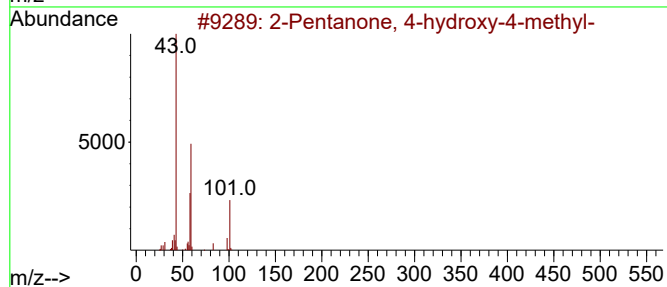
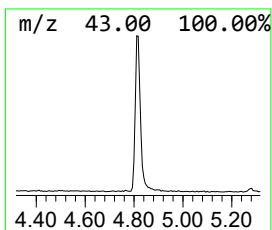
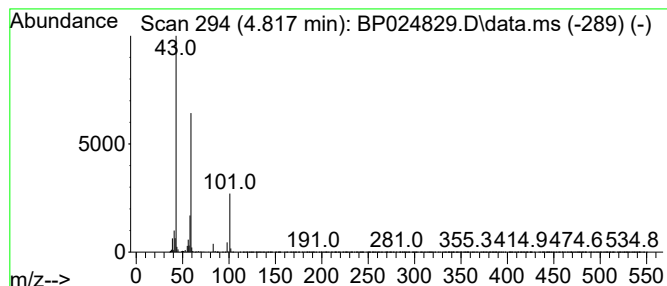
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
4.817	7.36 ng	242988	1,4-Dichlorobenzene-d4		7.646	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	56
2	(+-)-4-Amino-4,5-dihydro-2(3H)-f...		101	C4H7NO2	016504-58-8	47
3	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	25
4	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
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-...	4.817	7.4	ng	242988	1	7.646	660393	20.0