

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
Data File : BP021983.D
Acq On : 23 Sep 2024 15:28
Operator : RC/JU
Sample : PB163571BL
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PB163571BL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BP091924.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP021983.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.834	309	314	324	rVB	80967	111093	1.42%	0.402%
2	5.293	385	392	405	rBV	1689805	2393484	30.66%	8.665%
3	6.851	647	657	669	rBV	1555041	2400398	30.75%	8.690%
4	7.669	788	796	804	rBV	323272	513158	6.57%	1.858%
5	8.810	982	990	1001	rBV	920324	1567085	20.07%	5.673%
6	10.439	1258	1267	1279	rBV	373357	629180	8.06%	2.278%
7	12.922	1681	1689	1701	rBV	2124387	3591899	46.01%	13.004%
8	14.310	1917	1925	1937	rBV	503873	791770	10.14%	2.866%
9	15.833	2177	2184	2207	rBV	3108821	3720907	47.66%	13.471%
10	17.127	2397	2404	2418	rBV	702415	1080208	13.84%	3.911%
11	19.851	2860	2867	2875	rBV	5611148	7806660	100.00%	28.262%
12	21.557	3151	3157	3165	rBV2	861564	1476048	18.91%	5.344%
13	24.856	3710	3718	3730	rVB	522343	1540269	19.73%	5.576%

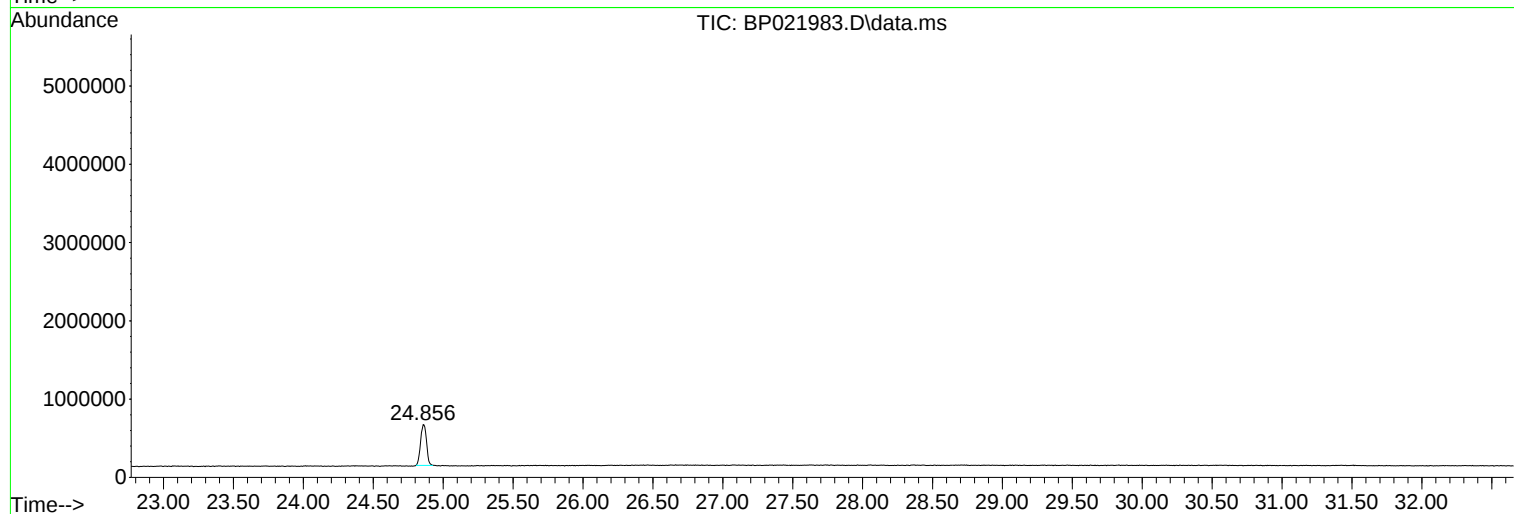
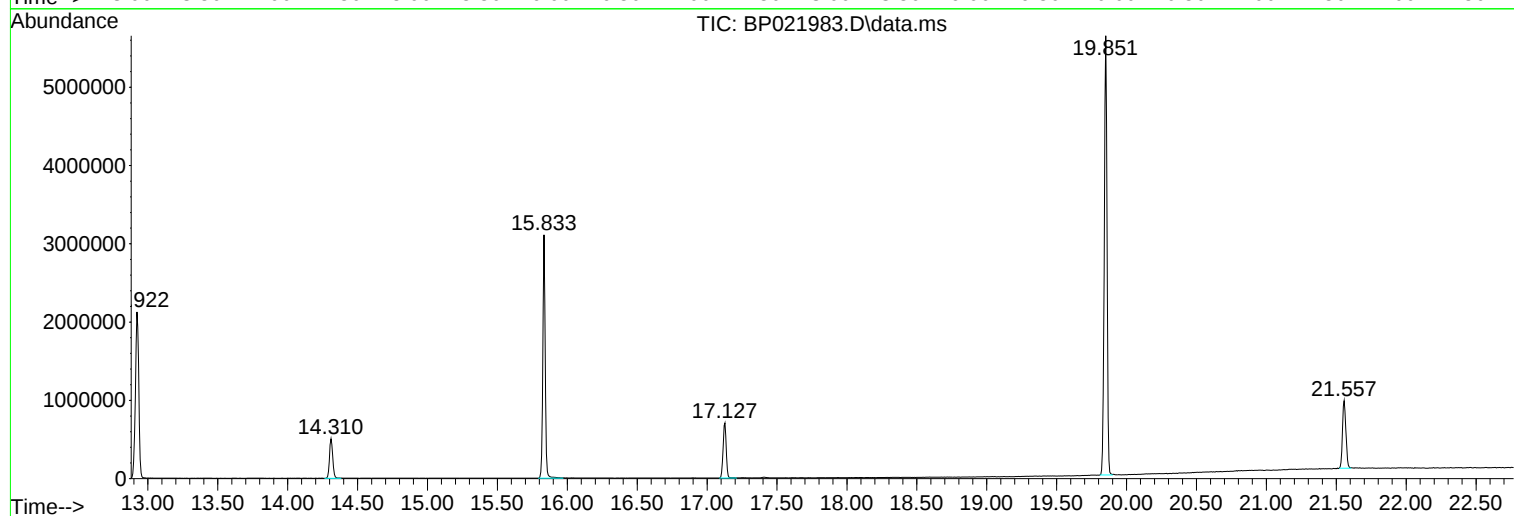
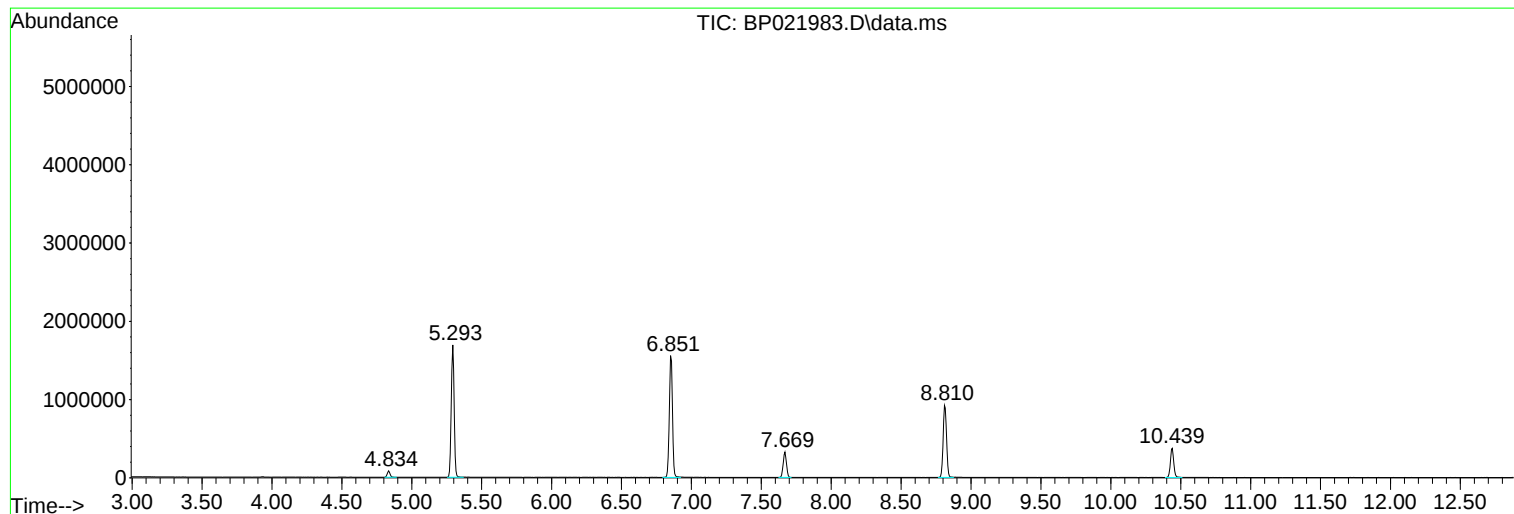
Sum of corrected areas: 27622159

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.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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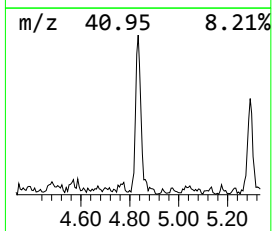
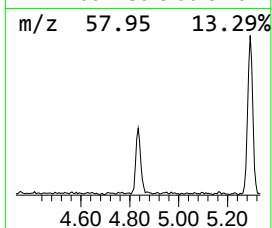
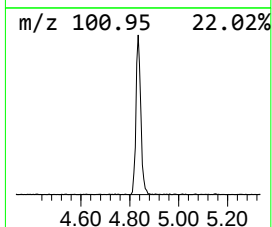
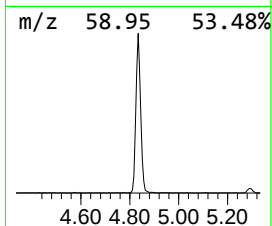
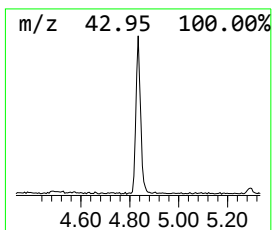
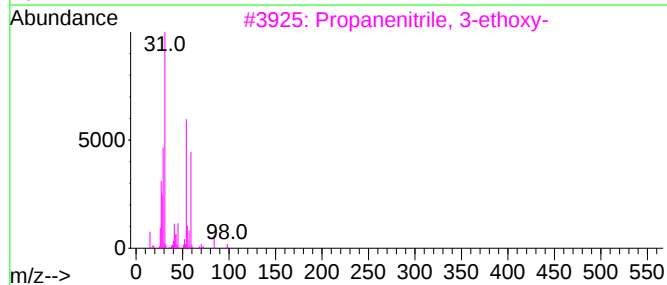
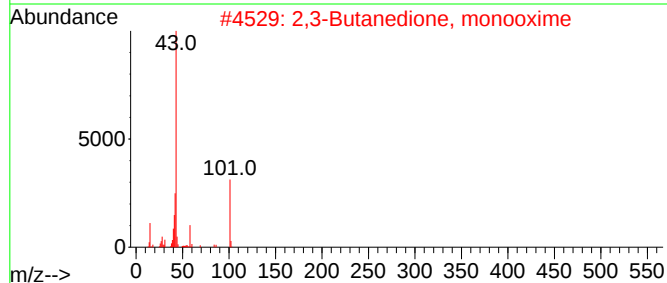
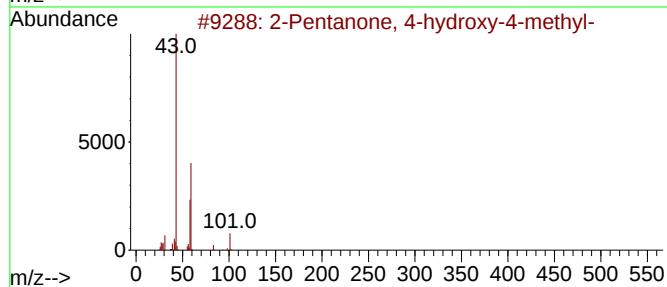
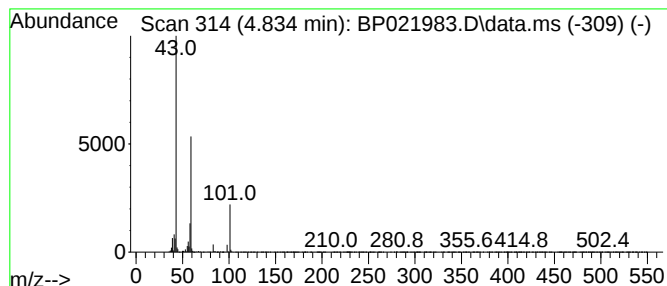
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.834	4.33 ng	111093	1,4-Dichlorobenzene-d4	7.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	Propanenitrile, 3-ethoxy-	99	C5H9NO	002141-62-0	9		
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.834	4.3	ng	111093	1	7.669	513158	20.0