

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091225\
 Data File : BP025738.D
 Acq On : 12 Sep 2025 11:05
 Operator : CG/JU
 Sample : PB169490BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 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_P\Methods\8270E-BP091225.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025738.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.925	332	338	351	rBV	358805	526009	4.44%	1.138%
2	5.384	409	416	439	rBV	2914677	4420038	37.32%	9.563%
3	6.943	674	681	705	rBV	2566310	4527433	38.23%	9.795%
4	7.755	813	819	833	rBV	476424	810263	6.84%	1.753%
5	8.902	1006	1014	1035	rBV	1597220	3001560	25.35%	6.494%
6	10.531	1283	1291	1310	rBV	444235	1102059	9.31%	2.384%
7	12.990	1700	1709	1721	rBV	3891751	6507269	54.95%	14.079%
8	14.384	1938	1946	1963	rBV	757069	1531168	12.93%	3.313%
9	15.895	2195	2203	2238	rBV	3207381	5477144	46.25%	11.850%
10	17.184	2415	2422	2433	rBV2	948528	1936140	16.35%	4.189%
11	19.901	2878	2884	2896	rBV	8706833	11842248	100.00%	25.622%
12	21.630	3171	3178	3192	rBV	1300747	2383277	20.13%	5.156%
13	24.989	3740	3749	3768	rVB	658876	2155075	18.20%	4.663%

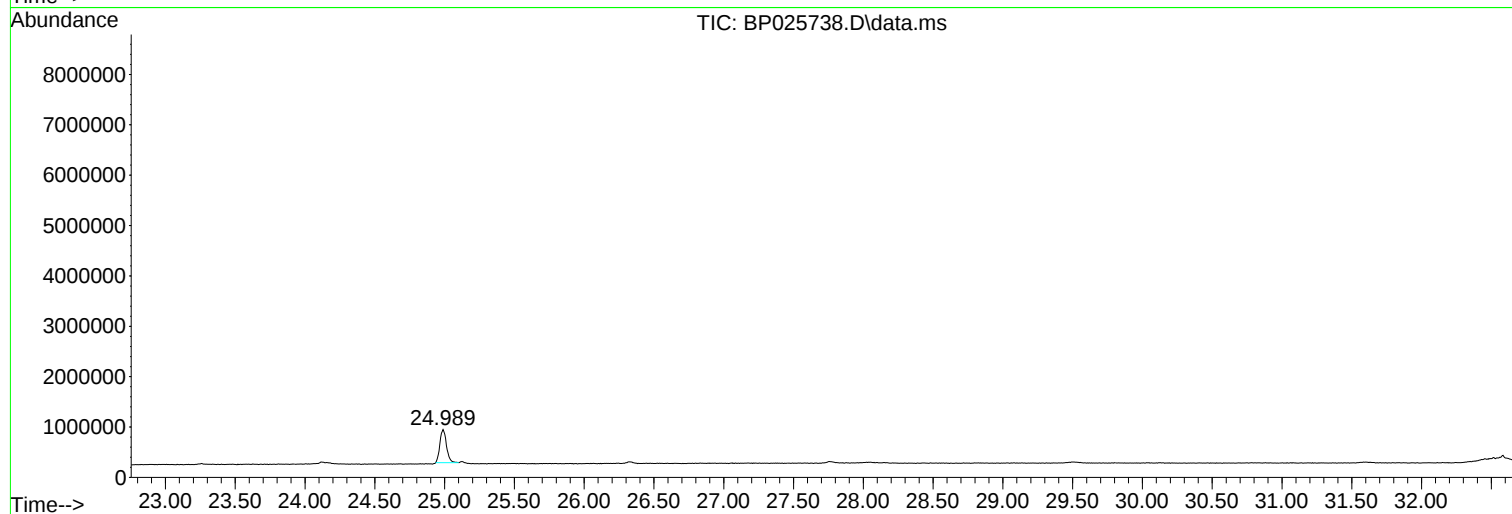
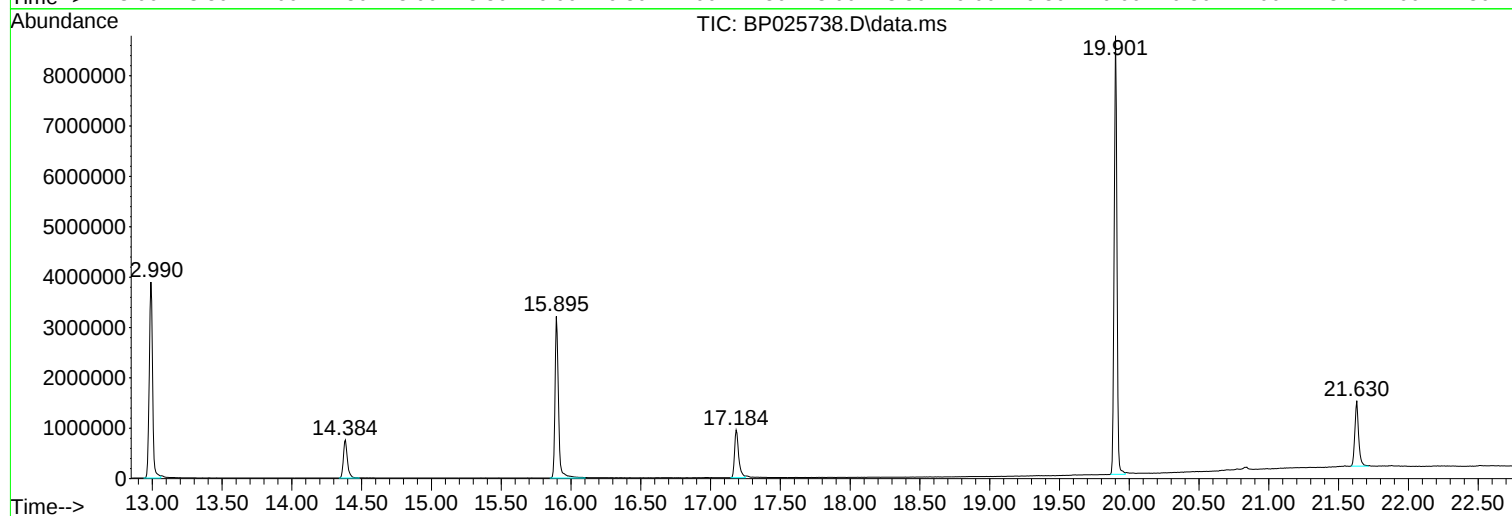
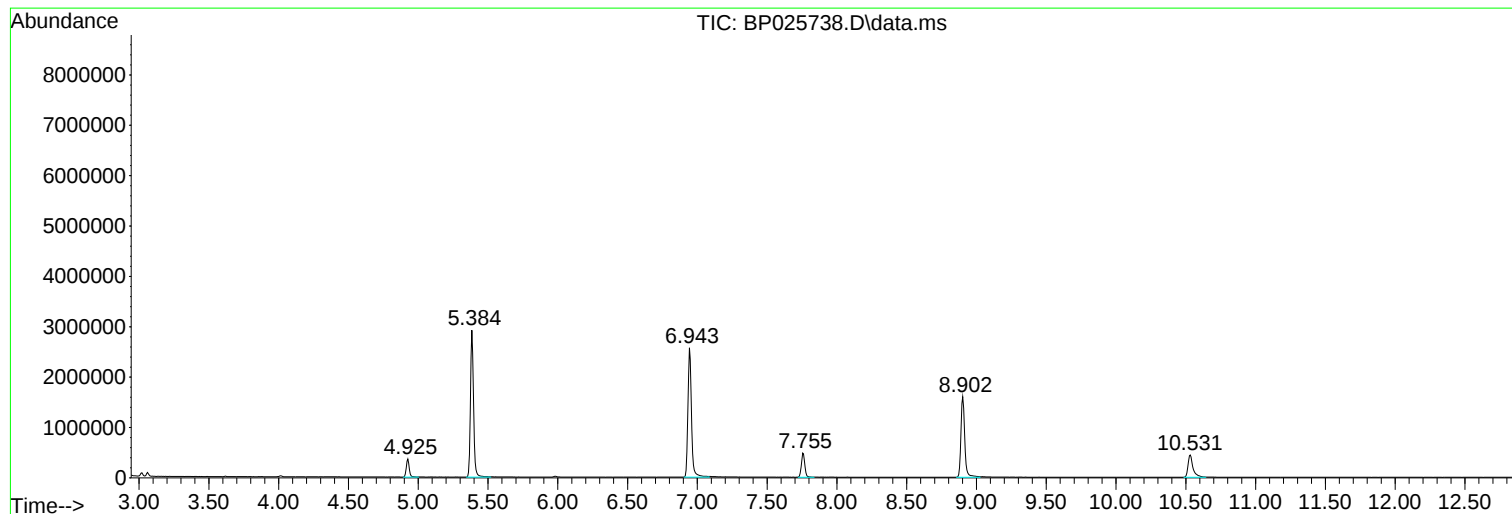
Sum of corrected areas: 46219683

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.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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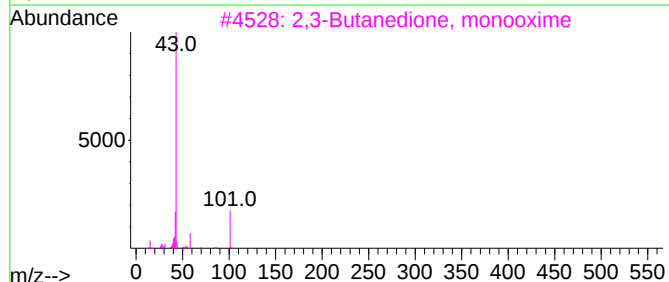
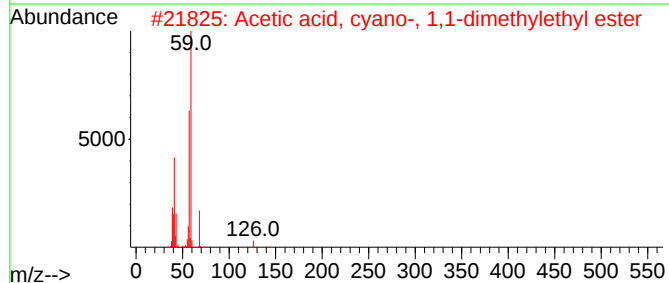
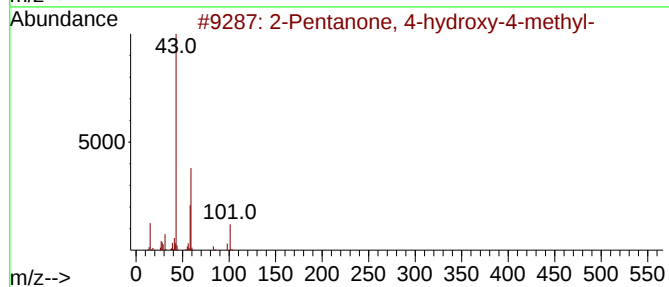
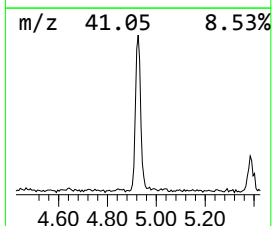
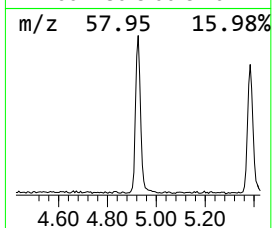
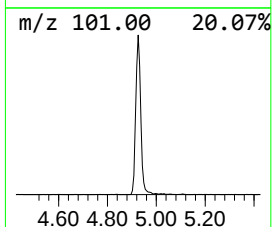
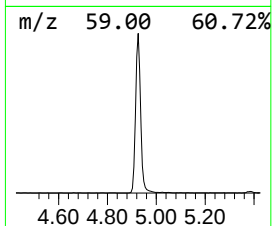
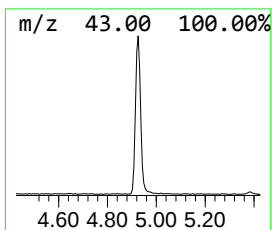
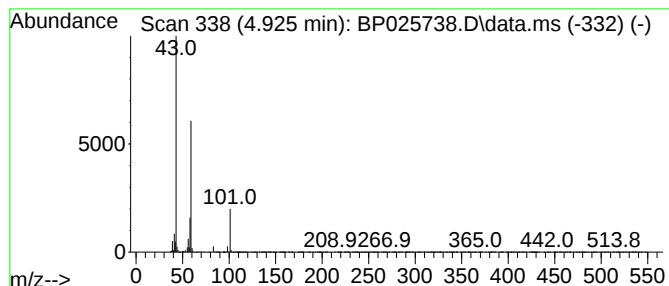
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.925	12.98 ng	526009	1,4-Dichlorobenzene-d4			7.755
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	72
2	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	23
3	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	9
4	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	9
5	3-Hexanol, 4-ethyl-		130	C8H18O	019780-44-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.925	13.0	ng	526009	1	7.755	810263	20.0