

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
 Data File : BP007929.D
 Acq On : 10 Nov 2021 20:32
 Operator : CG/JU
 Sample : M4578-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB

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-BP110221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007929.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.975	280	287	297	rVB	120816	173464	1.98%	0.545%
2	5.440	360	366	375	rBV	1541448	2207398	25.16%	6.934%
3	7.028	627	636	652	rBV	904666	1447675	16.50%	4.548%
4	7.851	769	776	784	rBV	406052	644984	7.35%	2.026%
5	9.010	966	973	983	rBV	1222840	2095132	23.88%	6.582%
6	10.651	1244	1252	1262	rBV	470202	788855	8.99%	2.478%
7	13.116	1664	1671	1686	rBV	3258674	4914144	56.02%	15.437%
8	14.492	1897	1905	1917	rBV	672494	1052077	11.99%	3.305%
9	15.986	2150	2159	2179	rBV	2824858	4383721	49.97%	13.771%
10	17.245	2367	2373	2384	rBV	918535	1357320	15.47%	4.264%
11	19.809	2804	2809	2820	rBV	6704069	8772165	100.00%	27.557%
12	21.339	3064	3069	3074	rBV	1440545	1888016	21.52%	5.931%
13	23.750	3472	3479	3490	rVB2	985966	2107672	24.03%	6.621%

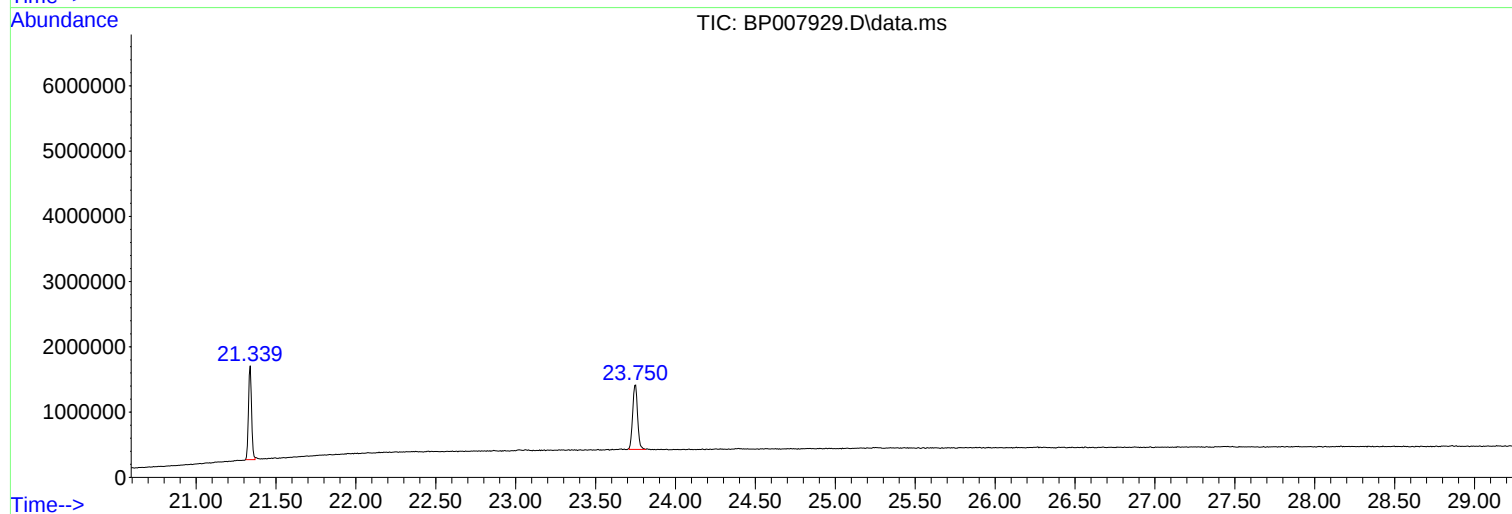
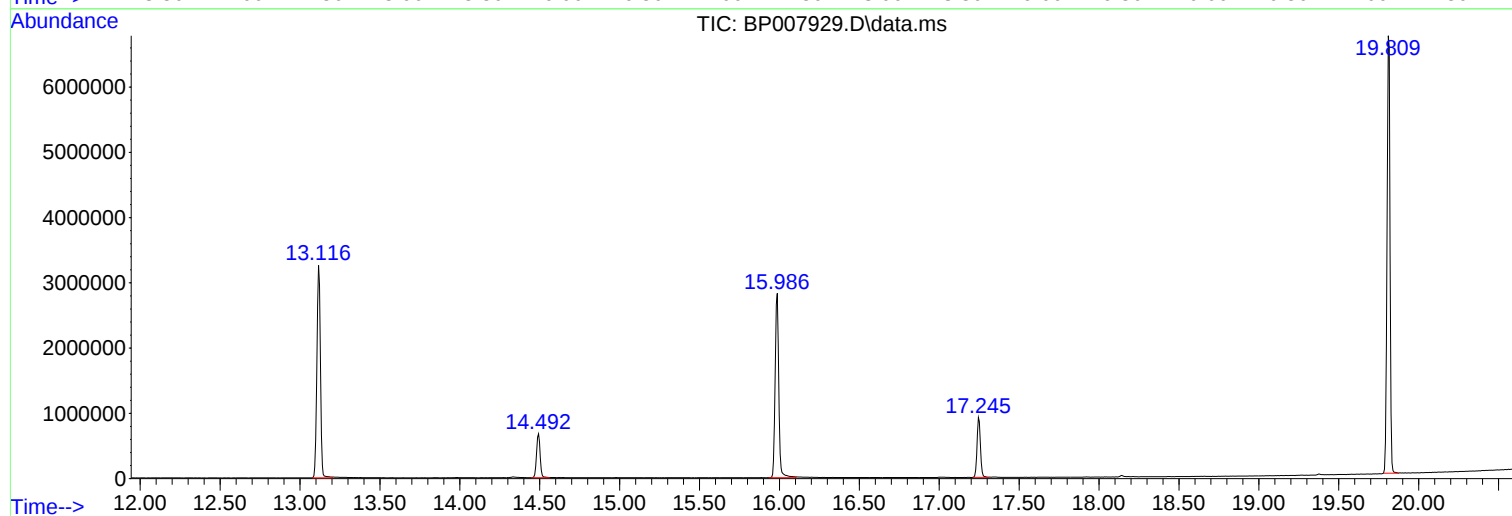
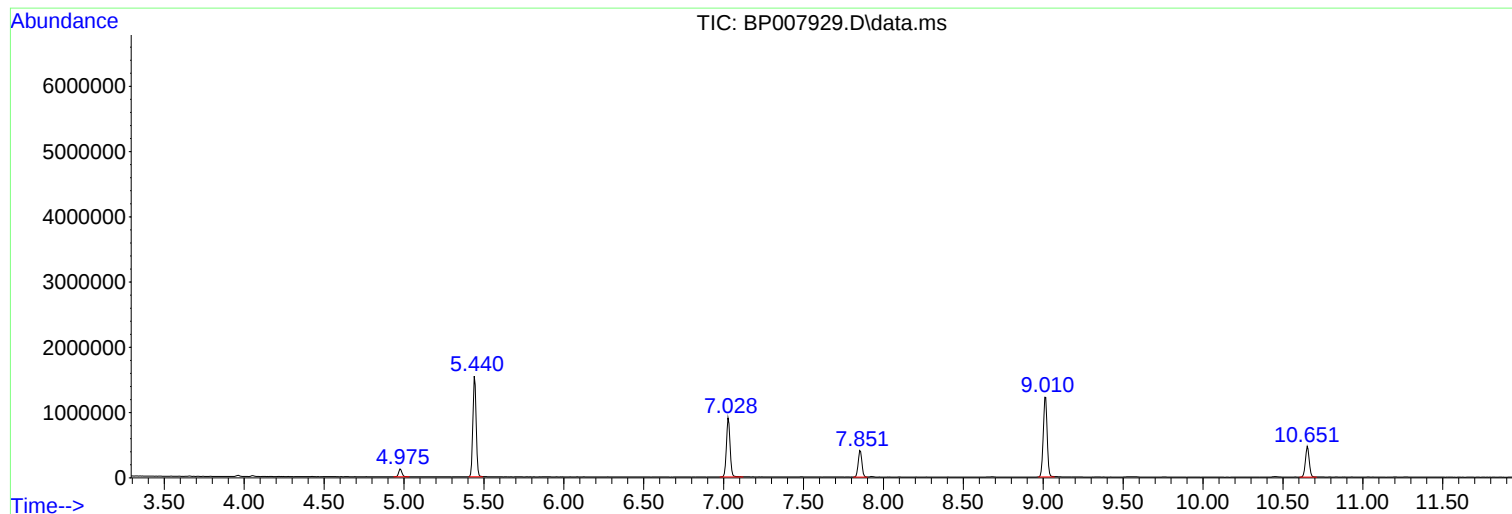
Sum of corrected areas: 31832623

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.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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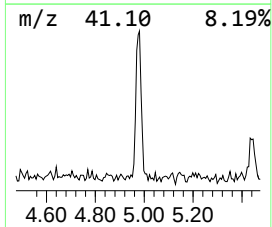
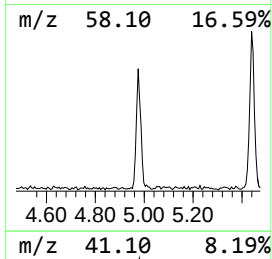
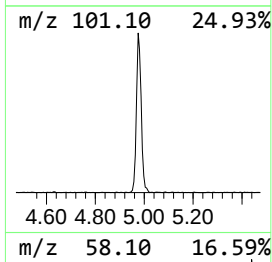
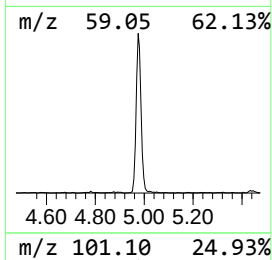
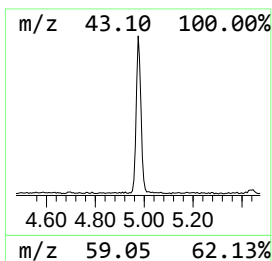
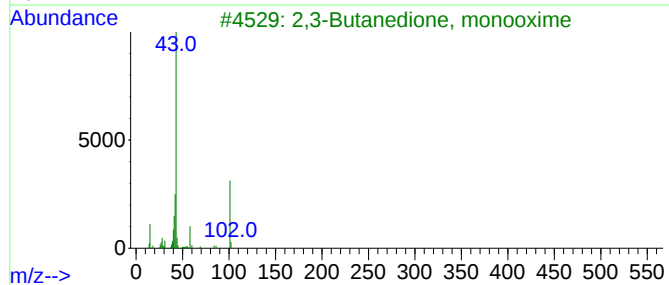
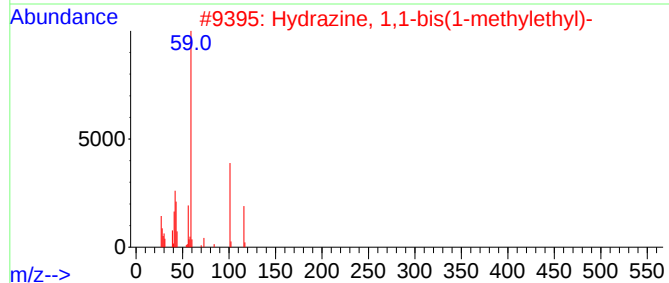
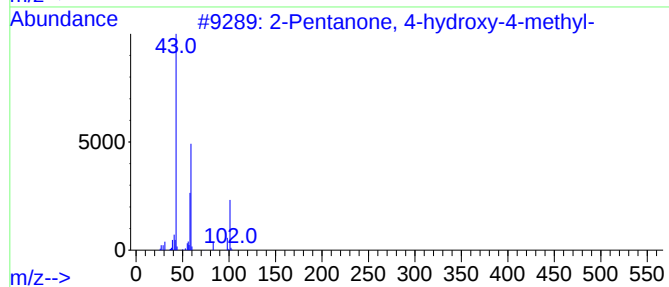
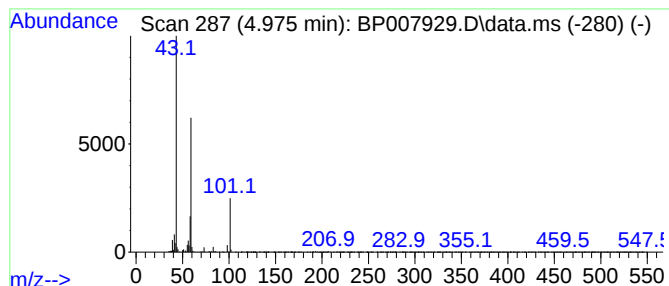
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.975	5.38 ng	173464	1,4-Dichlorobenzene-d4	7.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
4	Hexanamide	115	C6H13NO	000628-02-4	25		
5	Tetraglyme	222	C10H22O5	000143-24-8	25		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.975	5.4	ng	173464	1	7.851	644984	20.0