

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061021\
 Data File : BP005831.D
 Acq On : 10 Jun 2021 11:16
 Operator : CG/JU
 Sample : PB136969TB
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB136969TB

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-BP060821.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005831.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.128	3	7	12	rVB	42250	49337	1.32%	0.276%
2	5.052	329	334	342	rBV	35407	50091	1.34%	0.280%
3	5.516	406	413	426	rBV	1374464	2031210	54.37%	11.370%
4	7.116	677	685	700	rBV	1224184	1933505	51.75%	10.823%
5	7.952	820	827	836	rBV	263578	421545	11.28%	2.360%
6	9.116	1017	1025	1041	rBV	721051	1205890	32.28%	6.750%
7	10.757	1295	1304	1319	rBV	289230	516928	13.84%	2.894%
8	13.204	1713	1720	1734	rBV	1743412	2676589	71.64%	14.983%
9	14.581	1946	1954	1965	rBV	397242	626635	16.77%	3.508%
10	16.063	2195	2206	2226	rBV	1413364	2139297	57.26%	11.975%
11	17.322	2412	2420	2432	rBV	469267	701013	18.76%	3.924%
12	19.869	2847	2853	2863	rBV	2998714	3736091	100.00%	20.914%
13	21.392	3107	3112	3124	rBV	619489	828053	22.16%	4.635%
14	23.815	3517	3524	3539	rVB2	451601	947961	25.37%	5.307%

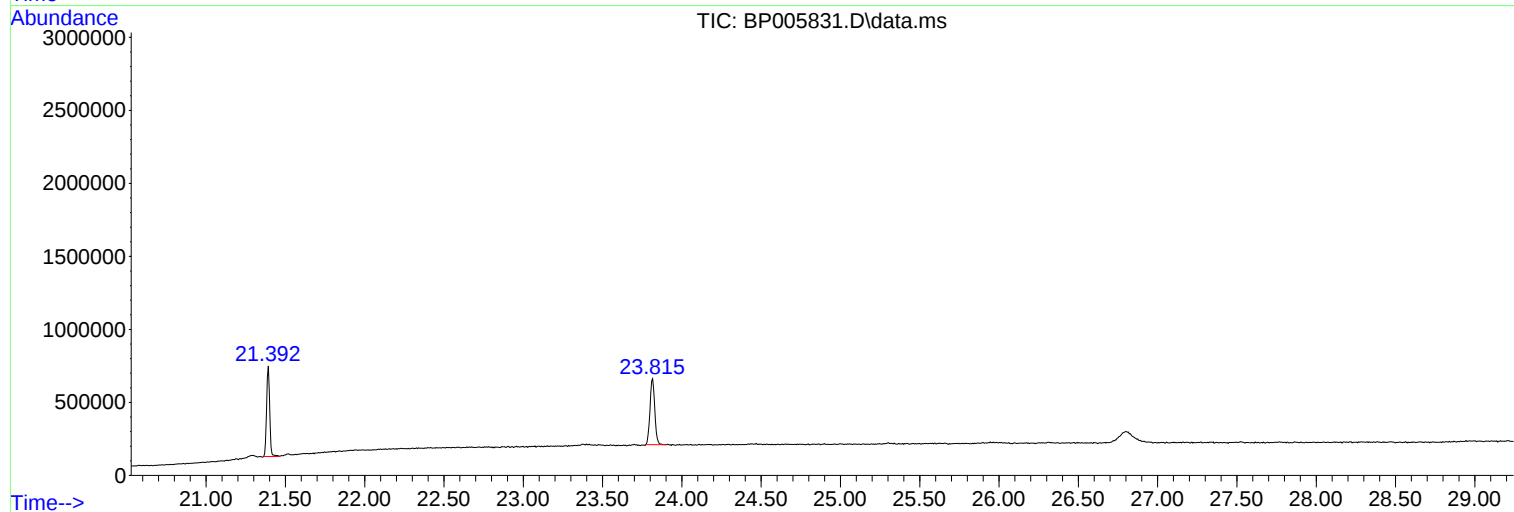
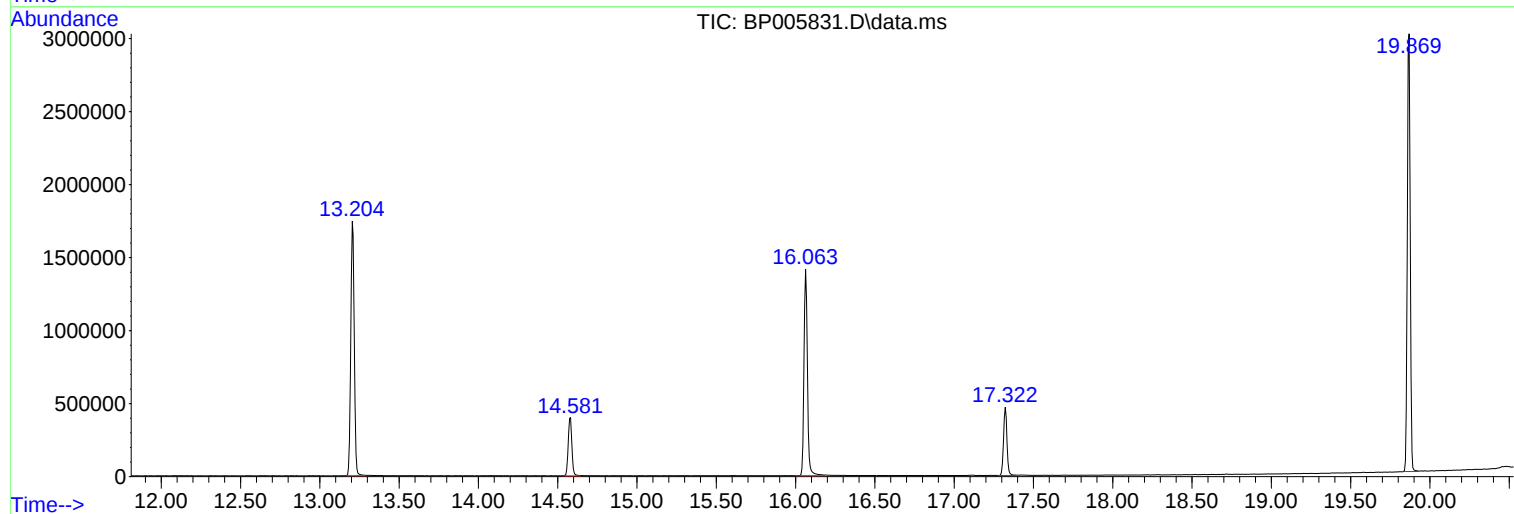
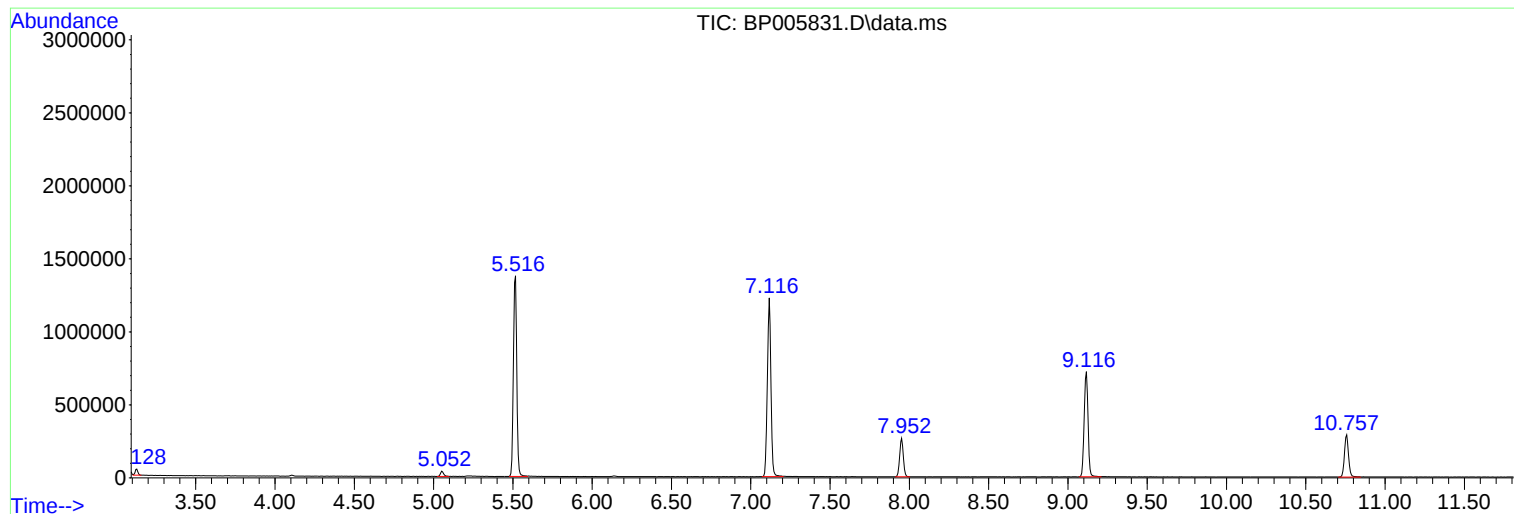
Sum of corrected areas: 17864145

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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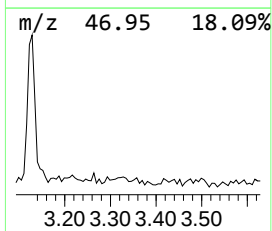
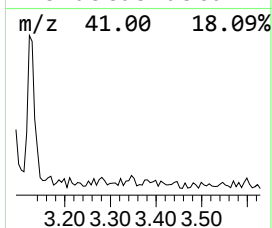
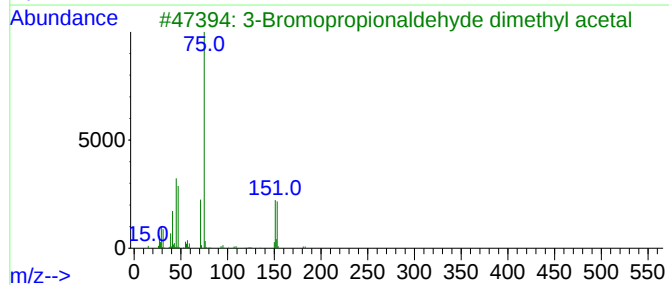
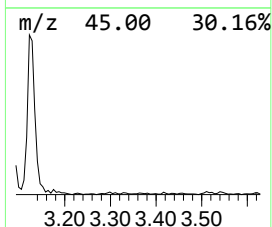
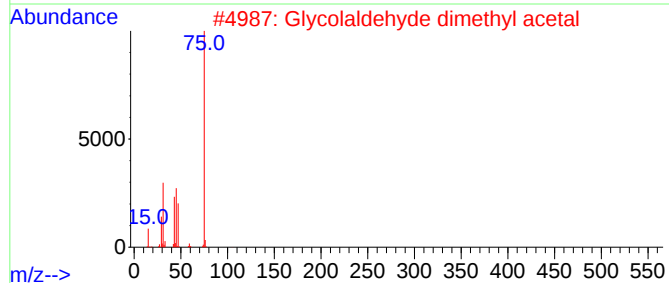
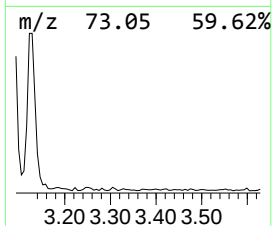
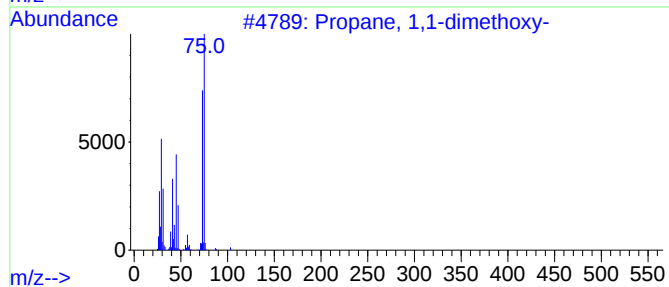
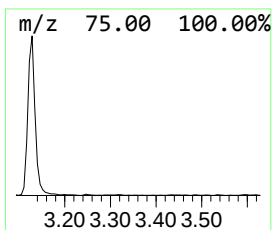
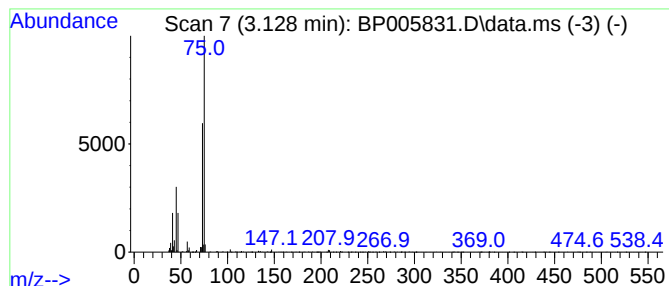
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.128	2.34 ng	49337	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	74
2			Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	50
3			3-Bromopropionaldehyde dimethyl ...	182	C5H11BrO2	036255-44-4	38
4			Propane, 1,1,3,3-tetramethoxy-	164	C7H16O4	000102-52-3	36
5			2-Propanol, 1,1-dimethoxy-, acetate	162	C7H14O4	074495-37-7	36



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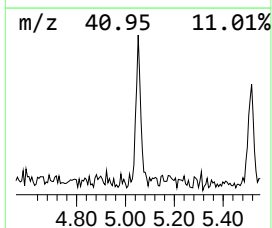
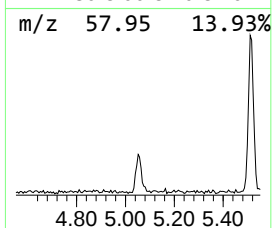
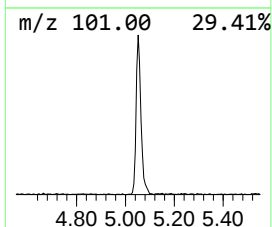
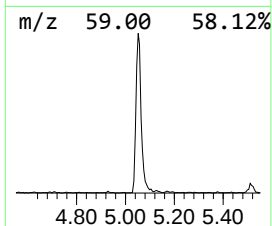
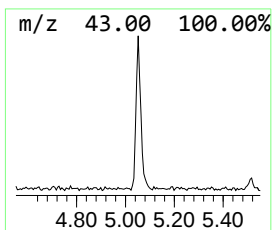
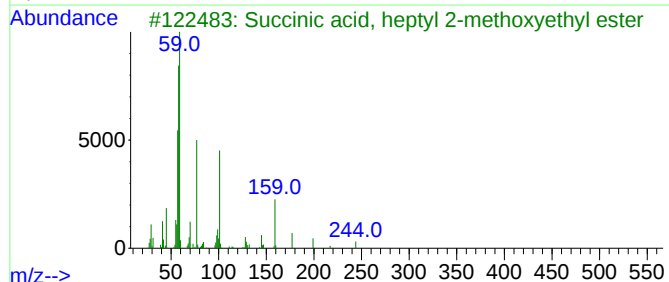
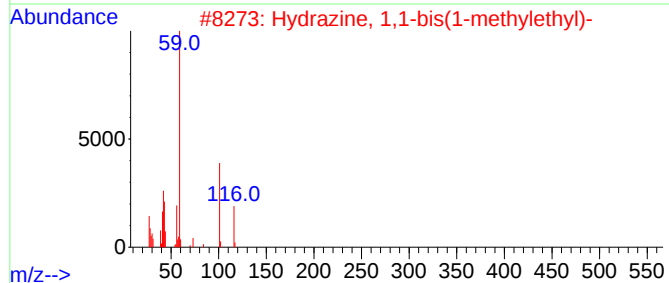
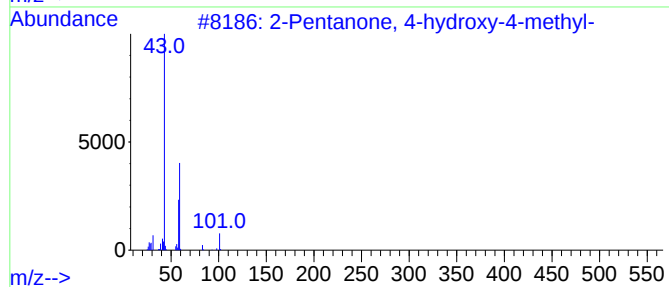
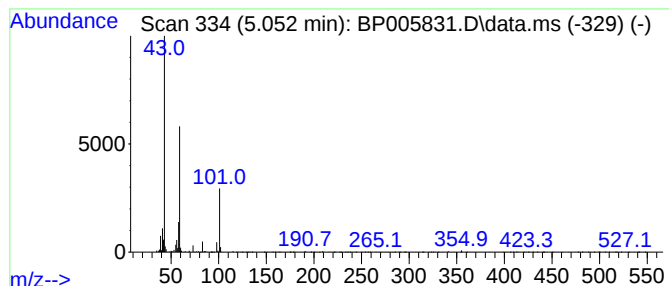
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.052	2.38 ng	50091	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38		
3	Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	28		
4	Hexanamide	115	C6H13NO	000628-02-4	27		
5	Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	27		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	3.128	2.3	ng	49337	1	7.952	421545	20.0
2-Pentanone, 4-...	5.052	2.4	ng	50091	1	7.952	421545	20.0