

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032621\
 Data File : BM029232.D
 Acq On : 26 Mar 2021 13:53
 Operator : CG/JU
 Sample : PB135047BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB135047BL

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_M\Methods\8270-BM032621.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM029232.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.016	19	22	26	rVV	45966	51856	1.23%	0.228%
2	3.057	26	29	38	rVB	106799	132427	3.14%	0.581%
3	4.910	340	344	353	rBV	60945	88128	2.09%	0.387%
4	5.357	413	420	433	rBV	1988238	2903681	68.91%	12.740%
5	6.922	679	686	698	rBV	1905934	3017142	71.60%	13.237%
6	7.757	821	828	835	rBV	375462	576394	13.68%	2.529%
7	8.904	1015	1023	1040	rBV	1100649	1846506	43.82%	8.101%
8	10.539	1292	1301	1310	rBV	436981	747101	17.73%	3.278%
9	13.004	1713	1720	1728	rBV	2595892	3724356	88.38%	16.340%
10	14.380	1948	1954	1962	rBV	607196	859784	20.40%	3.772%
11	15.863	2200	2206	2226	rVB	1583710	2097359	49.77%	9.202%
12	17.115	2413	2419	2431	rBV	639332	902084	21.41%	3.958%
13	19.739	2860	2865	2874	rBV	3479856	4213943	100.00%	18.488%
14	21.286	3123	3128	3135	rBV	660069	809357	19.21%	3.551%
15	23.521	3502	3508	3515	rBV	438316	822313	19.51%	3.608%

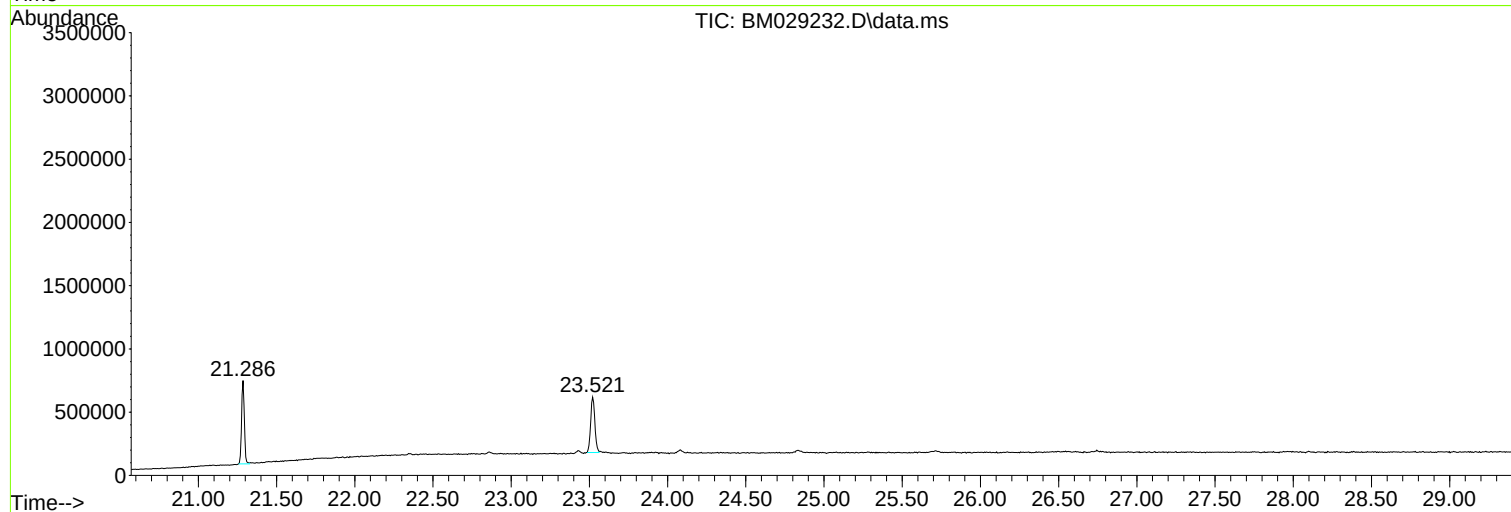
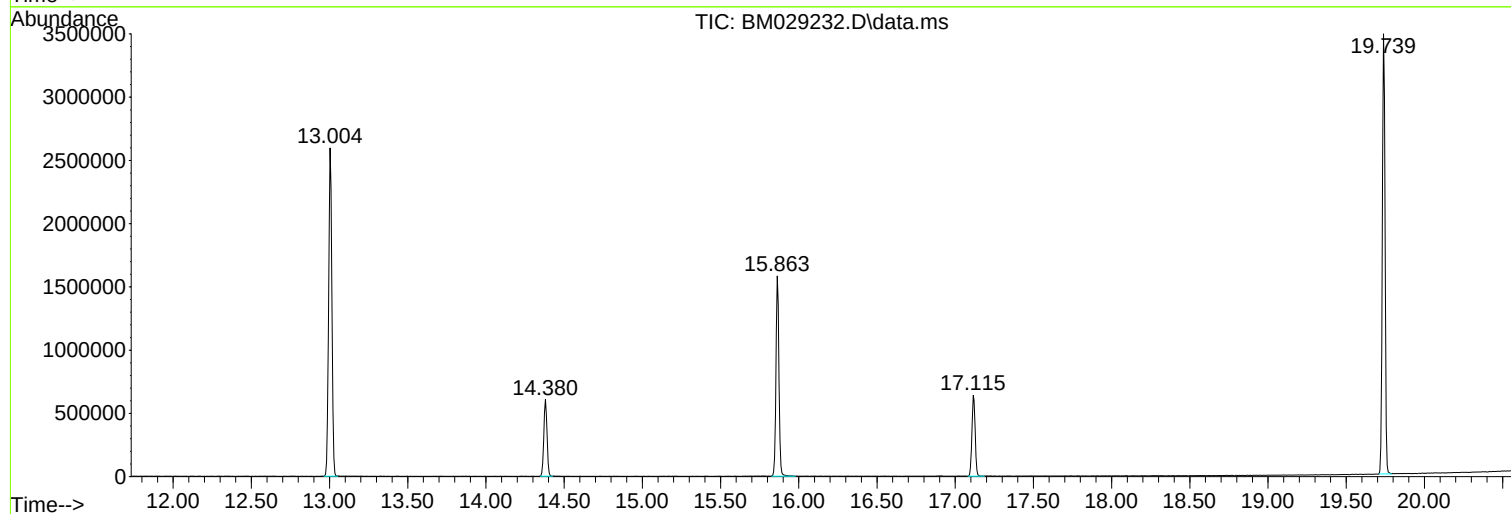
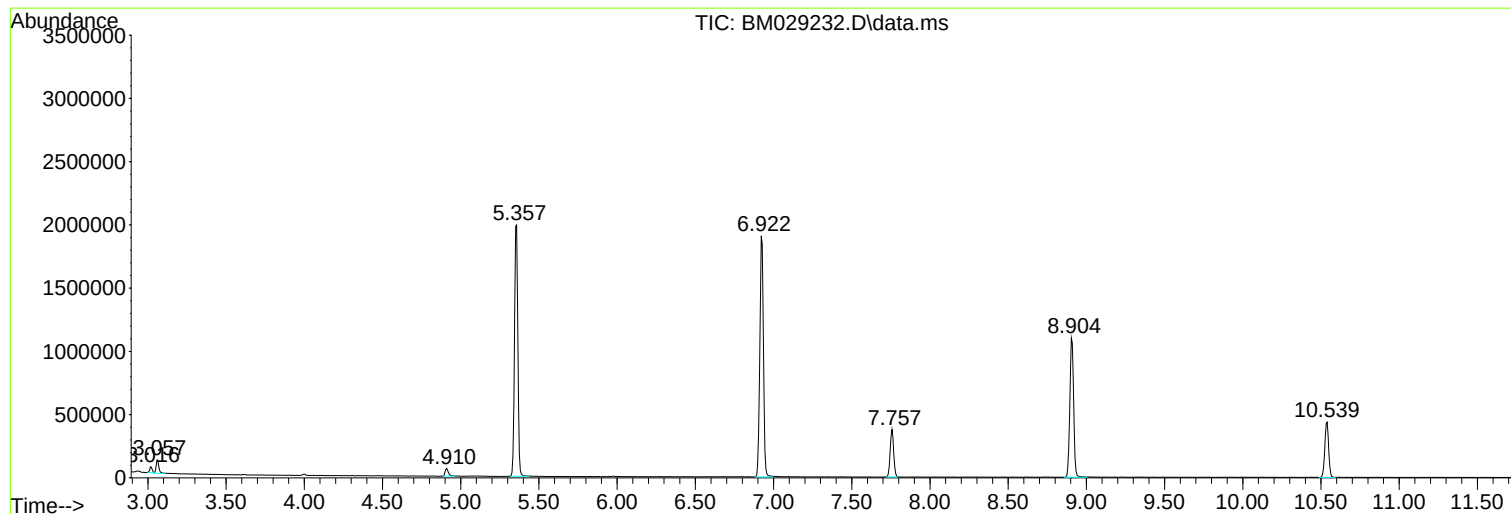
Sum of corrected areas: 22792431

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032621.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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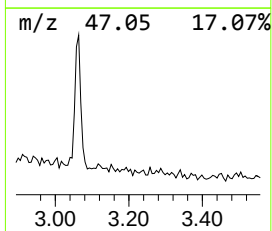
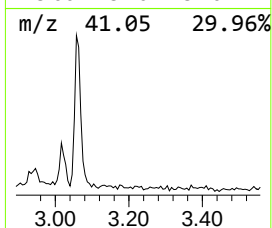
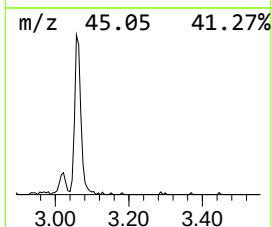
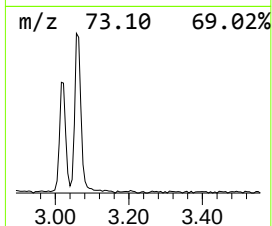
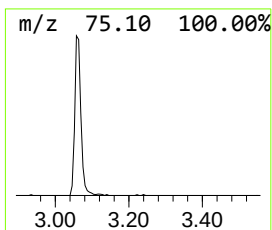
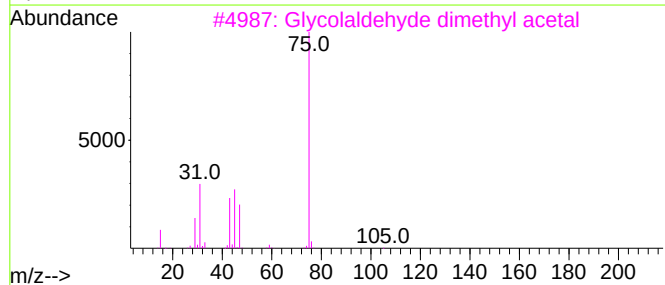
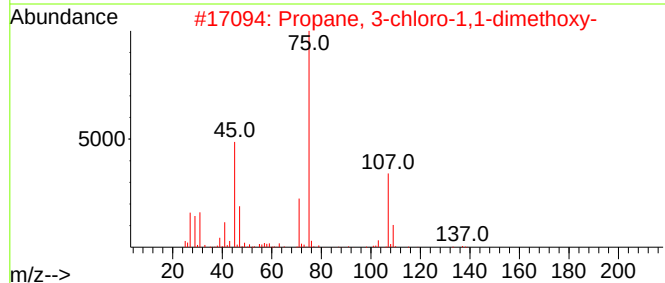
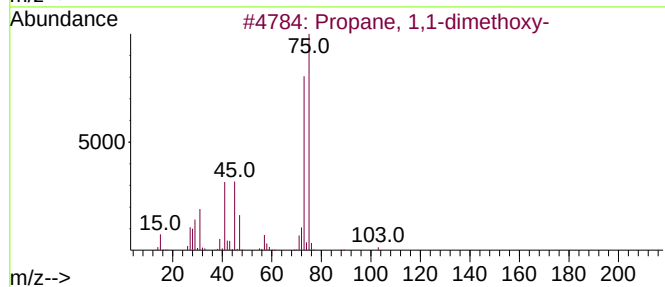
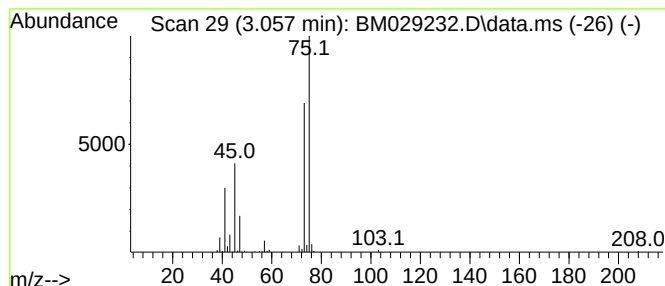
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Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.057	4.60 ng	132427	1,4-Dichlorobenzene-d4	7.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	64
2			Propane, 3-chloro-1,1-dimethoxy-	138	C5H11ClO2	035502-06-8	56
3			Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	45
4			3-Bromopropionaldehyde dimethyl ...	182	C5H11BrO2	036255-44-4	39
5			2-Propanol, 1,1-dimethoxy-, acetate	162	C7H14O4	074495-37-7	38



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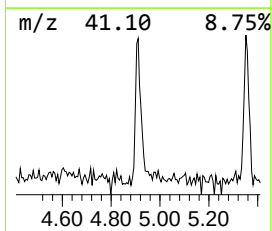
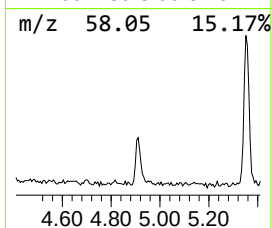
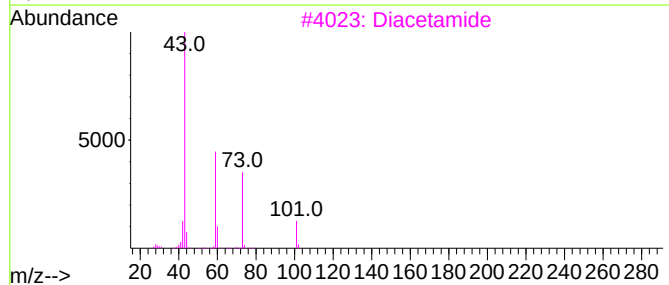
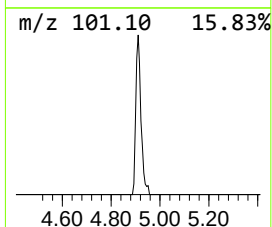
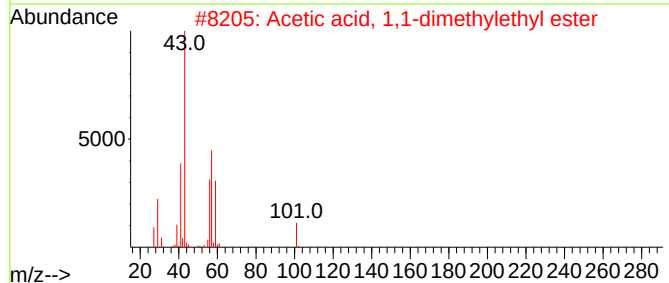
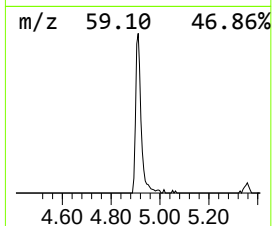
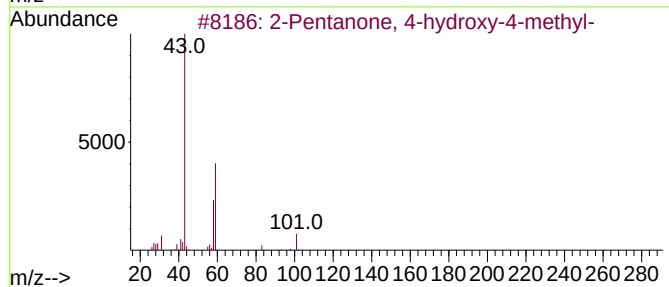
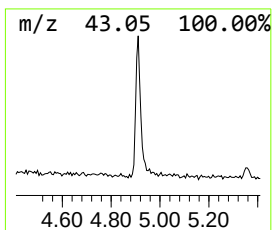
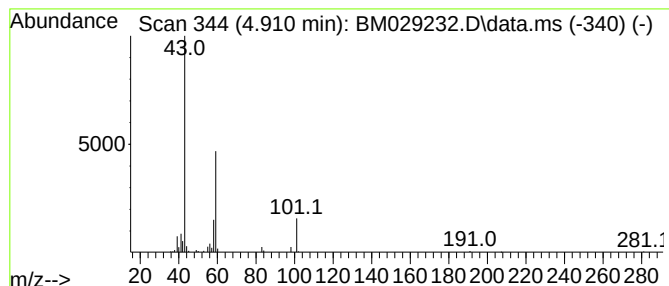
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.910	3.06 ng	88128	1,4-Dichlorobenzene-d4	7.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25
3			Diacetamide	101	C4H7NO2	000625-77-4	9
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Butane	58	C4H10	000106-97-8	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	3.057	4.6	ng	132427	1	7.757	576394	20.0
2-Pentanone, 4-...	4.910	3.1	ng	88128	1	7.757	576394	20.0