

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060821\
 Data File : BP005788.D
 Acq On : 08 Jun 2021 23:39
 Operator : CG/JU
 Sample : PB136919BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB136919BL

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: BP005788.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.129	4	7	13	rVB	50626	57494	1.37%	0.298%
2	5.058	330	335	341	rBV	41514	60931	1.45%	0.316%
3	5.523	407	414	429	rBV	1346177	2020836	48.11%	10.466%
4	7.122	676	686	707	rBV	1217408	1953193	46.50%	10.115%
5	7.958	820	828	836	rBV	289267	456529	10.87%	2.364%
6	9.122	1018	1026	1045	rBV	728542	1244331	29.63%	6.444%
7	10.763	1298	1305	1323	rBV	325177	570767	13.59%	2.956%
8	13.210	1714	1721	1732	rBV	1926983	2838947	67.59%	14.703%
9	14.581	1948	1954	1969	rBV	464378	723656	17.23%	3.748%
10	16.069	2200	2207	2227	rBV	1515899	2301218	54.79%	11.918%
11	17.328	2414	2421	2435	rBV	555593	837194	19.93%	4.336%
12	19.869	2847	2853	2864	rBV	3397068	4200121	100.00%	21.752%
13	21.392	3107	3112	3118	rBV	764738	964983	22.98%	4.998%
14	23.816	3517	3524	3533	rBV2	511302	1078761	25.68%	5.587%

Sum of corrected areas: 19308961

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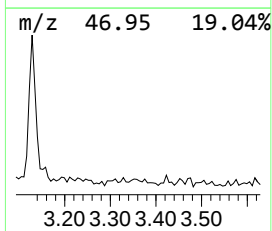
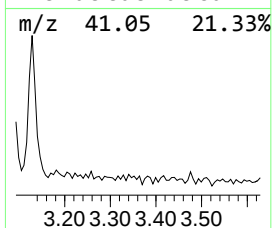
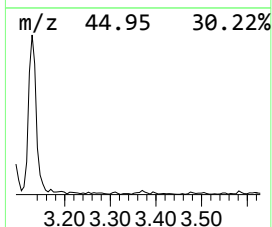
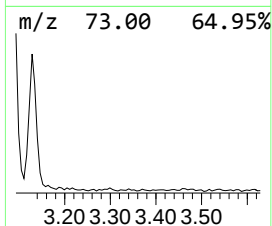
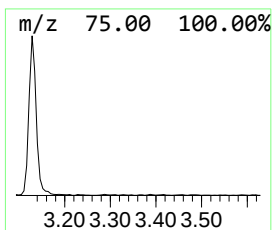
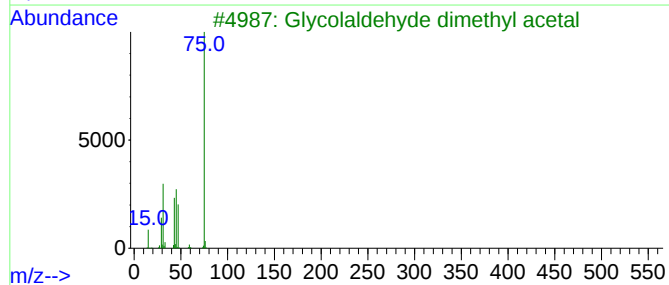
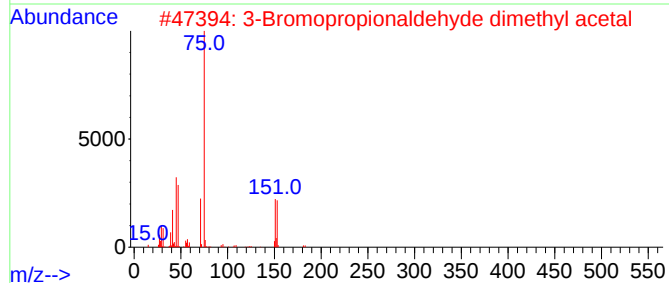
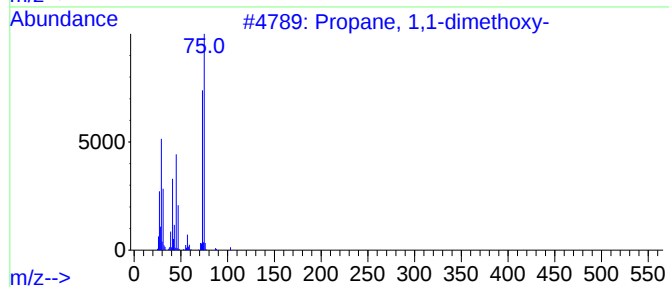
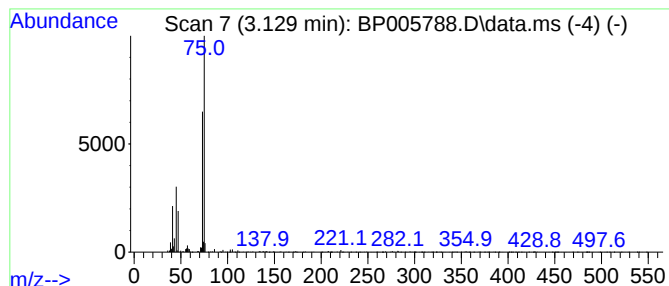
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060821.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.129	2.52 ng	57494	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2			3-Bromopropionaldehyde dimethyl ...	182	C5H11BrO2	036255-44-4	56
3			Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	56
4			4-Methylvaleric acid, trimethyls...	188	C9H20O2Si	959048-10-3	40
5			2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	39



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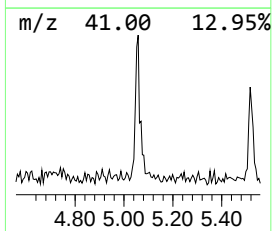
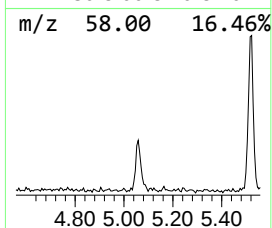
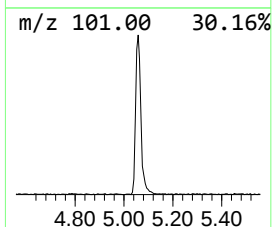
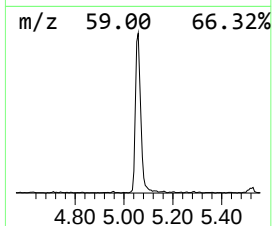
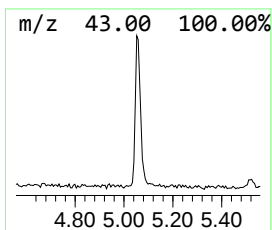
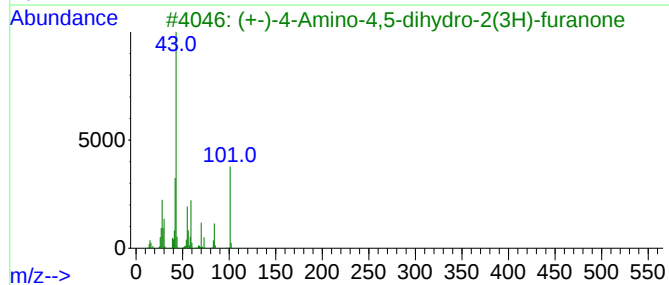
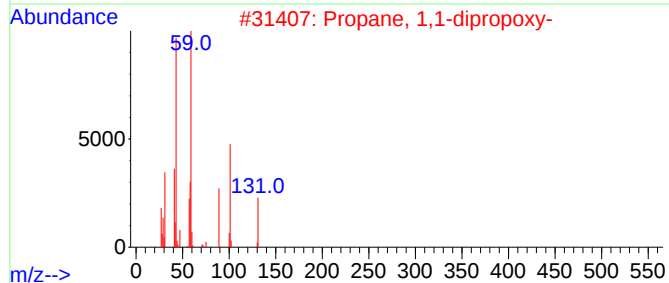
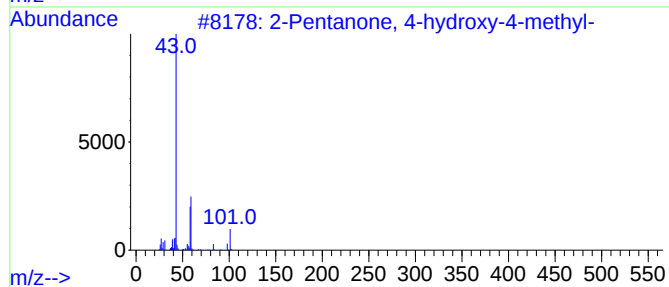
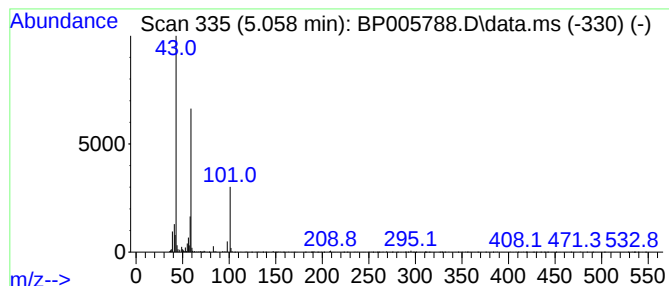
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.058	2.67 ng	60931	1,4-Dichlorobenzene-d4	7.958

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	38
3			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	32
4			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	28
5			Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	28



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	3.129	2.5	ng	57494	1	7.958	456529	20.0
2-Pentanone, 4-...	5.058	2.7	ng	60931	1	7.958	456529	20.0