

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101220\
 Data File : BF121922.D
 Acq On : 12 Oct 2020 17:08
 Operator : JU/CG
 Sample : PB132280BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB132280BL

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_F\METHODS\8270-BF101220.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF121922.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.193	14	18	29	rVB	77446	104005	1.99%	0.348%
2	4.992	491	494	509	rBV	71803	121756	2.33%	0.408%
3	5.381	555	560	563	rBV	3714253	3788772	72.64%	12.692%
4	6.392	727	732	735	rBV	3280581	3925698	75.27%	13.151%
5	6.745	788	792	806	rVB	865583	749517	14.37%	2.511%
6	7.316	883	889	892	rBV	2506503	2772738	53.16%	9.288%
7	8.028	1005	1010	1031	rBV	969475	989683	18.97%	3.315%
8	9.110	1187	1194	1197	rBV	4333471	4682303	89.77%	15.685%
9	9.780	1303	1308	1318	rBV	1117504	1134635	21.75%	3.801%
10	10.574	1437	1443	1466	rBV	2677771	3016653	57.84%	10.106%
11	11.269	1555	1561	1569	rBV	1337163	1192880	22.87%	3.996%
12	12.863	1824	1832	1835	rBV	4841757	5215834	100.00%	17.473%
13	13.910	2006	2010	2020	rVB	1193522	1126314	21.59%	3.773%
14	15.351	2249	2255	2269	rBV	816108	1030559	19.76%	3.452%

Sum of corrected areas: 29851347

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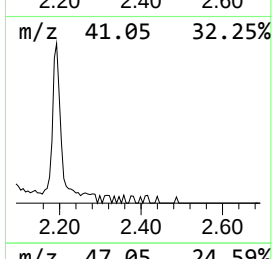
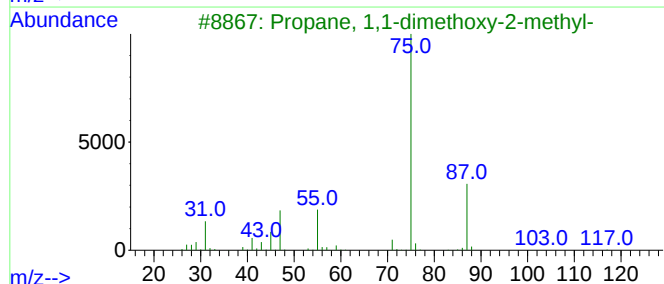
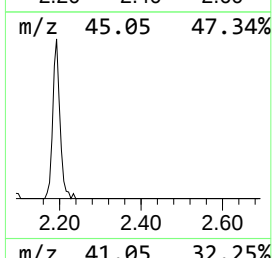
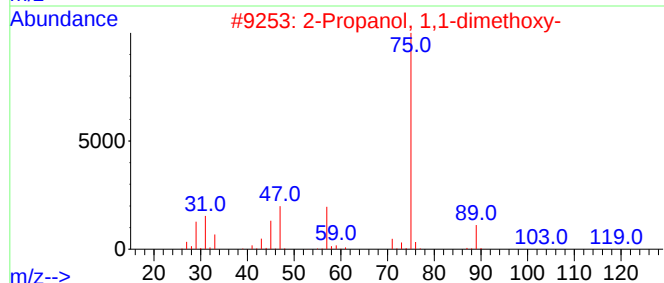
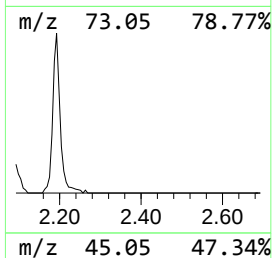
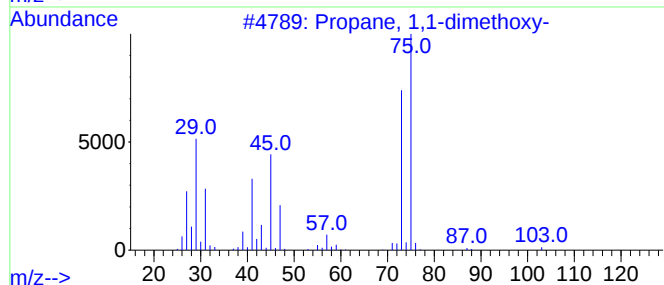
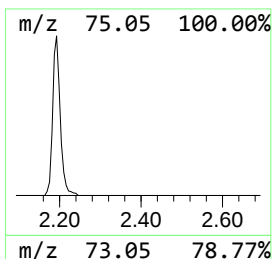
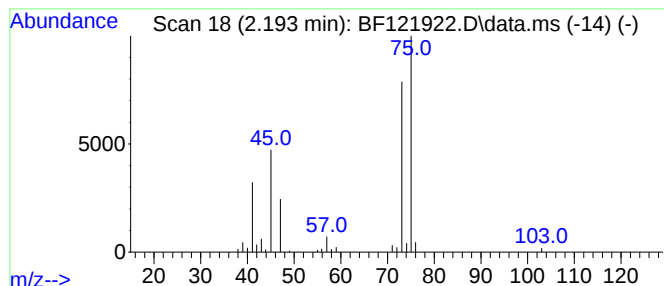
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.193	2.78 ng	104005	1,4-Dichlorobenzene-d4	6.745

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2			2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	33
3			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	28
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	16
5			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9



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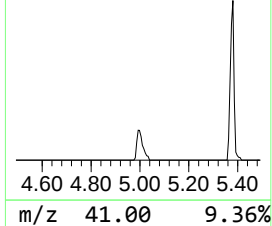
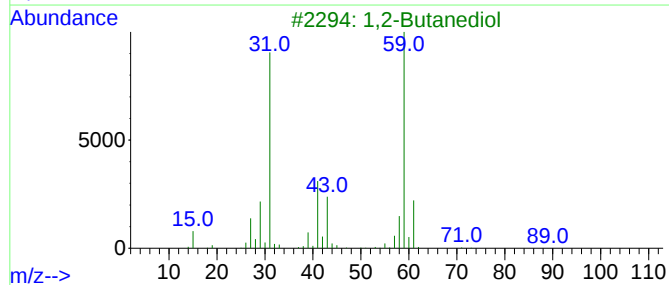
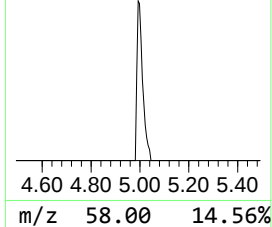
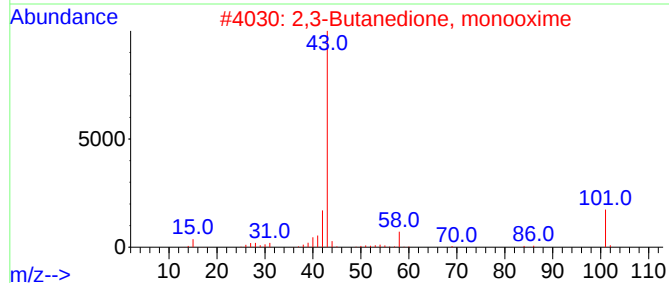
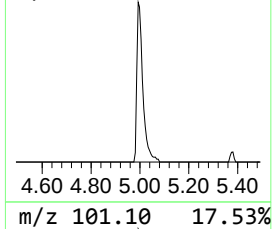
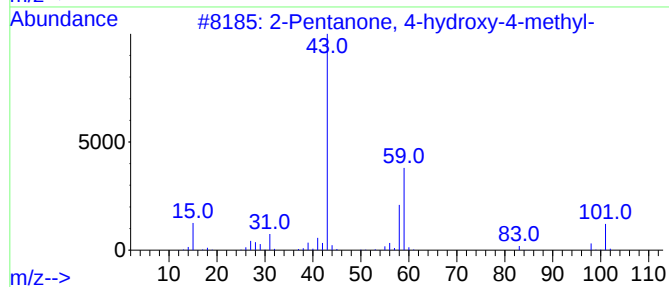
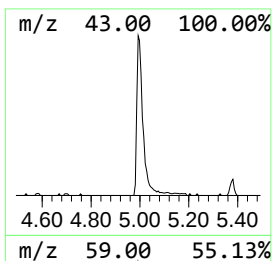
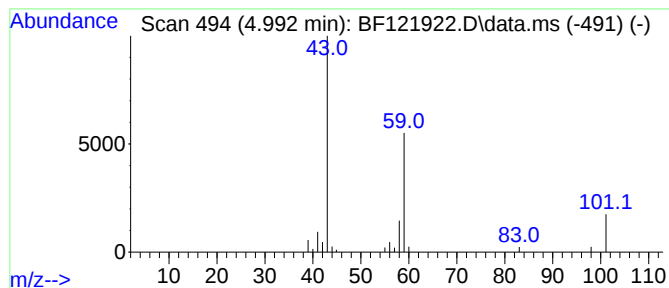
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
4.992	3.25 ng	121756	1,4-Dichlorobenzene-d4			6.745
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	64
2	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	9
3	1,2-Butanediol		90	C4H10O2	000584-03-2	9
4	2-Propanone, 1-(1-methylethoxy)-		116	C6H12O2	042781-12-4	9
5	Butyl aldoxime, 3-methyl-, syn-		101	C5H11NO	005780-40-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	2.193	2.8	ng	104005	1	6.745	749517	20.0
2-Pentanone, 4-...	4.992	3.3	ng	121756	1	6.745	749517	20.0