

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112320\
 Data File : BP004057.D
 Acq On : 23 Nov 2020 16:14
 Operator : CG/JU
 Sample : PB133168BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB133168BL

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-BP110320.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP004057.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.917	3	5	17	rVB	300879	350745	3.72%	1.318%
2	5.264	398	404	413	rBV	121794	176746	1.88%	0.664%
3	5.799	488	495	508	rBV	890133	1399145	14.86%	5.258%
4	7.452	767	776	794	rBV	868971	1452288	15.42%	5.457%
5	8.263	905	914	924	rBV	407342	645645	6.86%	2.426%
6	9.428	1103	1112	1122	rBV	567149	928862	9.86%	3.490%
7	11.093	1387	1395	1408	rBV	499638	859997	9.13%	3.232%
8	13.510	1796	1806	1820	rBV	1617071	2406850	25.56%	9.044%
9	14.887	2033	2040	2054	rBV	760852	1139943	12.10%	4.284%
10	16.375	2285	2293	2322	rBV	2393381	3619334	38.43%	13.600%
11	17.622	2498	2505	2518	rBV	949692	1355480	14.39%	5.094%
12	20.122	2924	2930	2942	rBV	7767136	9417199	100.00%	35.387%
13	21.657	3186	3191	3201	rVB	1105580	1440070	15.29%	5.411%
14	24.127	3604	3611	3625	rVB2	682170	1419527	15.07%	5.334%

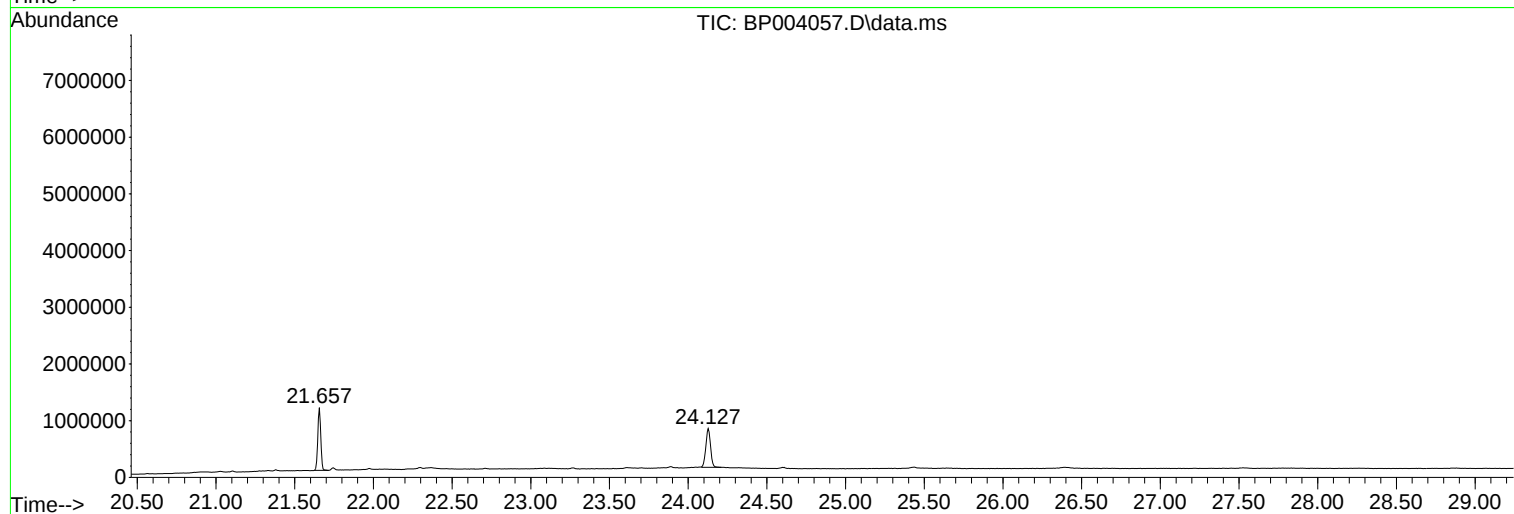
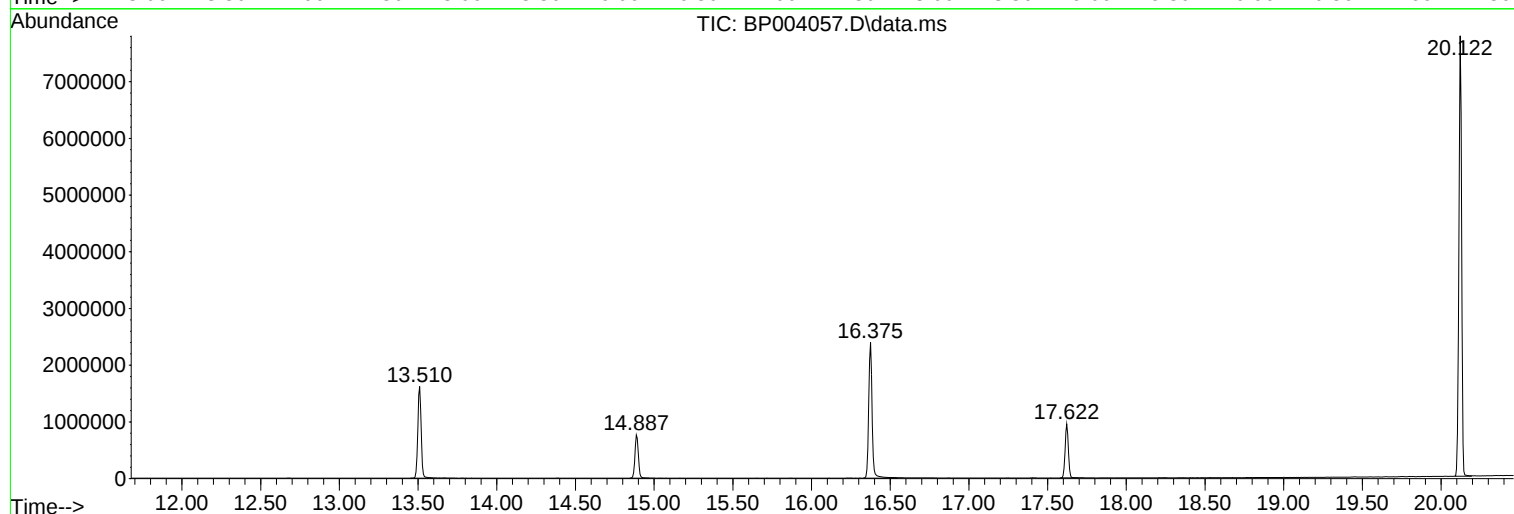
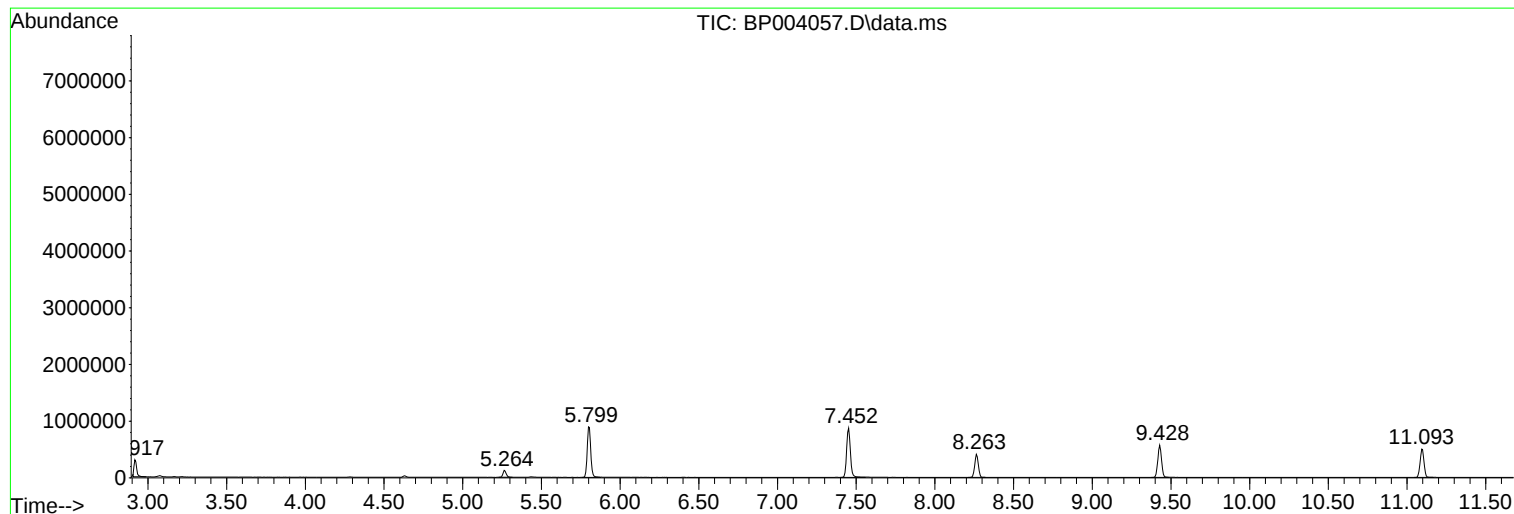
Sum of corrected areas: 26611831

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.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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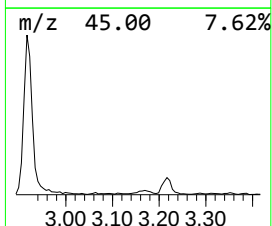
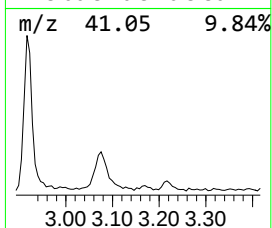
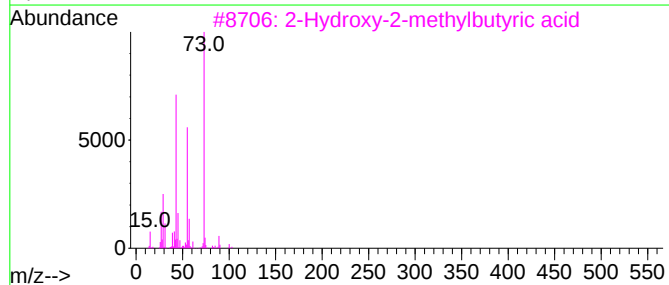
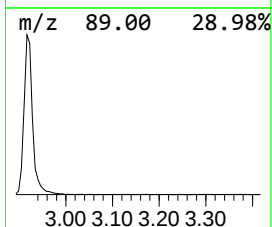
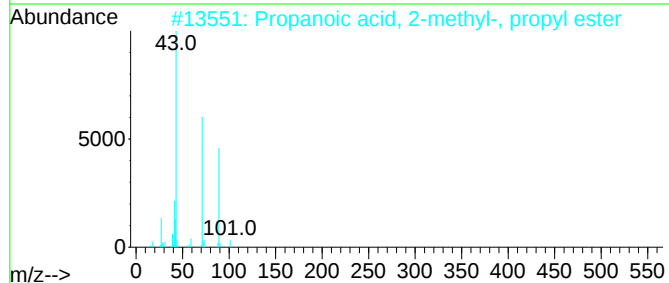
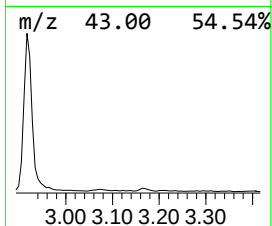
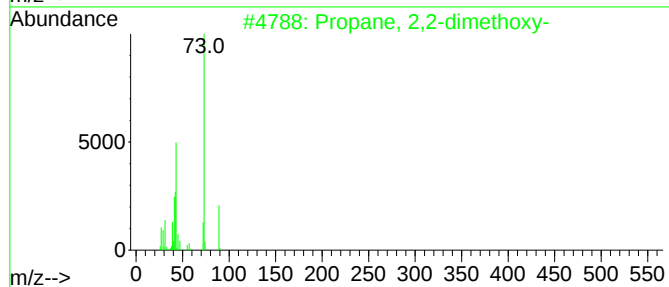
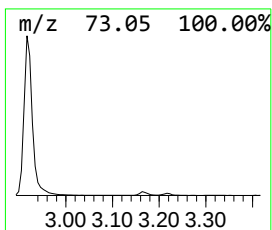
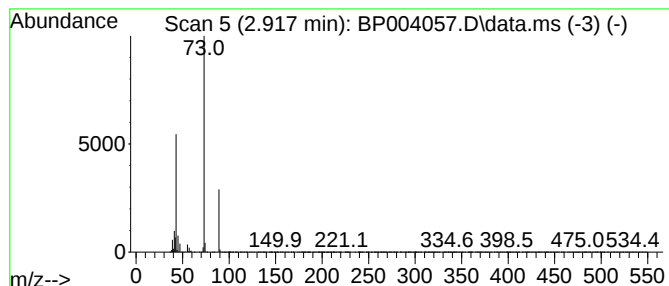
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP110320.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.917	10.86 ng	350745	1,4-Dichlorobenzene-d4	8.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	78
2			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
3			2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5			1-(2-Methoxyethoxy)-2-methyl-2-p...	162	C8H18O3	1000364-24-4	9



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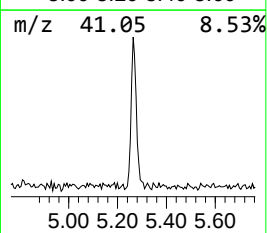
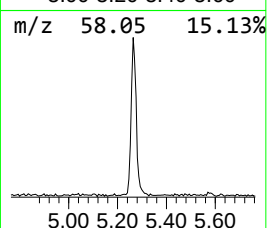
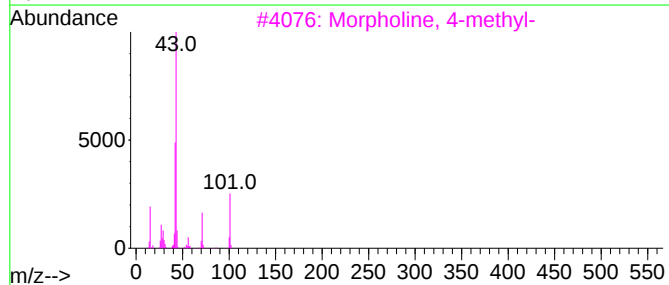
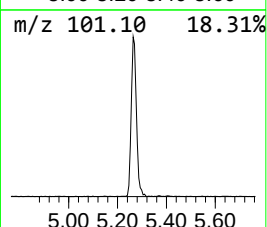
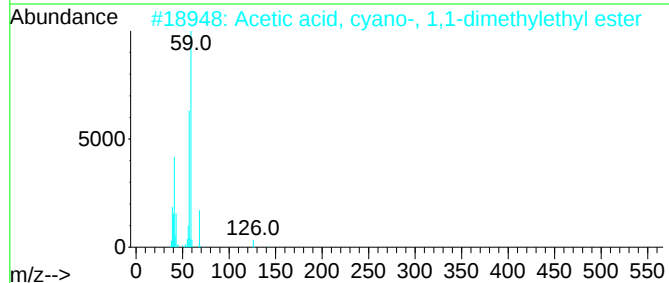
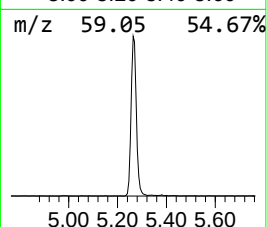
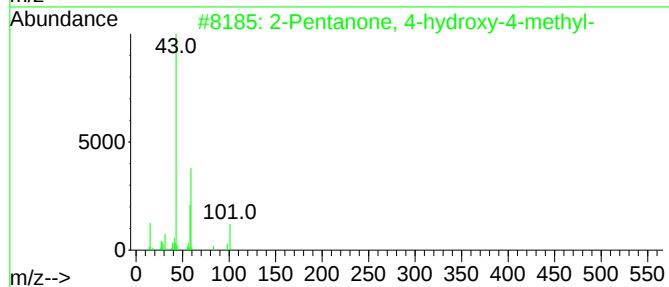
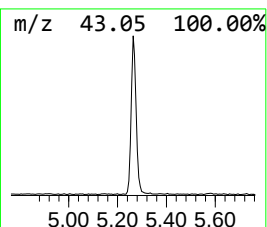
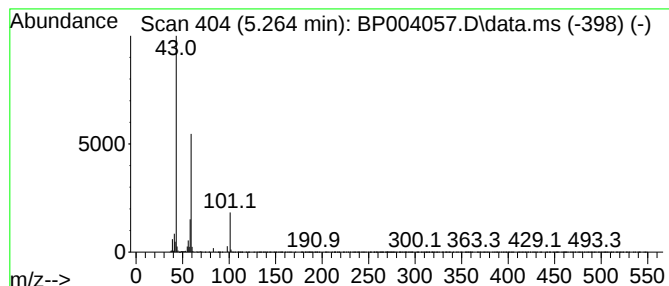
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.264	5.48 ng	176746	1,4-Dichlorobenzene-d4	8.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-di...	2.917	10.9	ng	350745	1	8.263	645645	20.0
2-Pentanone, 4-...	5.264	5.5	ng	176746	1	8.263	645645	20.0