

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091923\
 Data File : BF135340.D
 Acq On : 19 Sep 2023 17:53
 Operator : CG\JU
 Sample : PB155405BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB155405BL

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-BF091123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135340.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.175	9	15	22	rVB	58340	82513	2.01%	0.312%
2	5.016	492	498	507	rBV	90222	140285	3.42%	0.530%
3	5.398	558	563	566	rBV	3627491	3748682	91.42%	14.161%
4	6.404	728	734	753	rBV	3246028	3890203	94.87%	14.696%
5	6.757	790	794	797	rBV	927479	823308	20.08%	3.110%
6	7.322	884	890	893	rBV	2409858	2573672	62.77%	9.723%
7	8.034	1006	1011	1025	rBV	1220915	1109095	27.05%	4.190%
8	9.116	1189	1195	1198	rBV	4037334	4100394	100.00%	15.490%
9	9.792	1304	1310	1313	rBV	1296472	1182546	28.84%	4.467%
10	10.586	1440	1445	1462	rBV	2415314	2500250	60.98%	9.445%
11	11.280	1560	1563	1567	rBV	1284019	1240407	30.25%	4.686%
12	12.886	1830	1836	1839	rBV	3410268	3638772	88.74%	13.746%
13	13.939	2011	2015	2025	rBV	732193	704849	17.19%	2.663%
14	15.404	2259	2264	2283	rVB	501324	736180	17.95%	2.781%

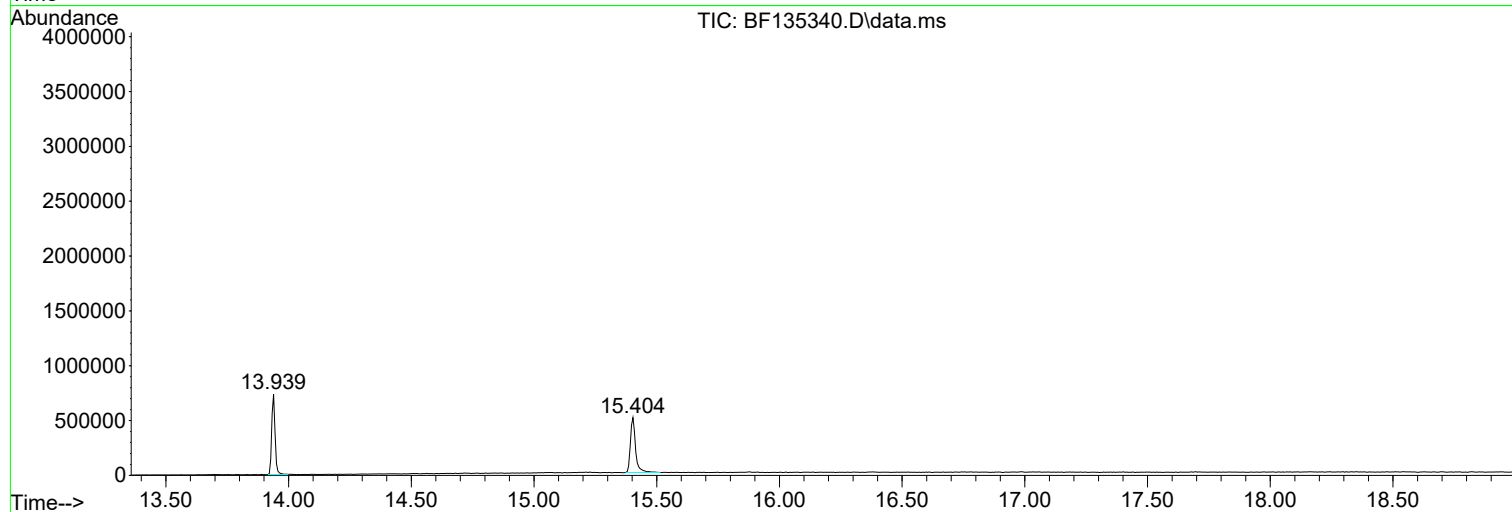
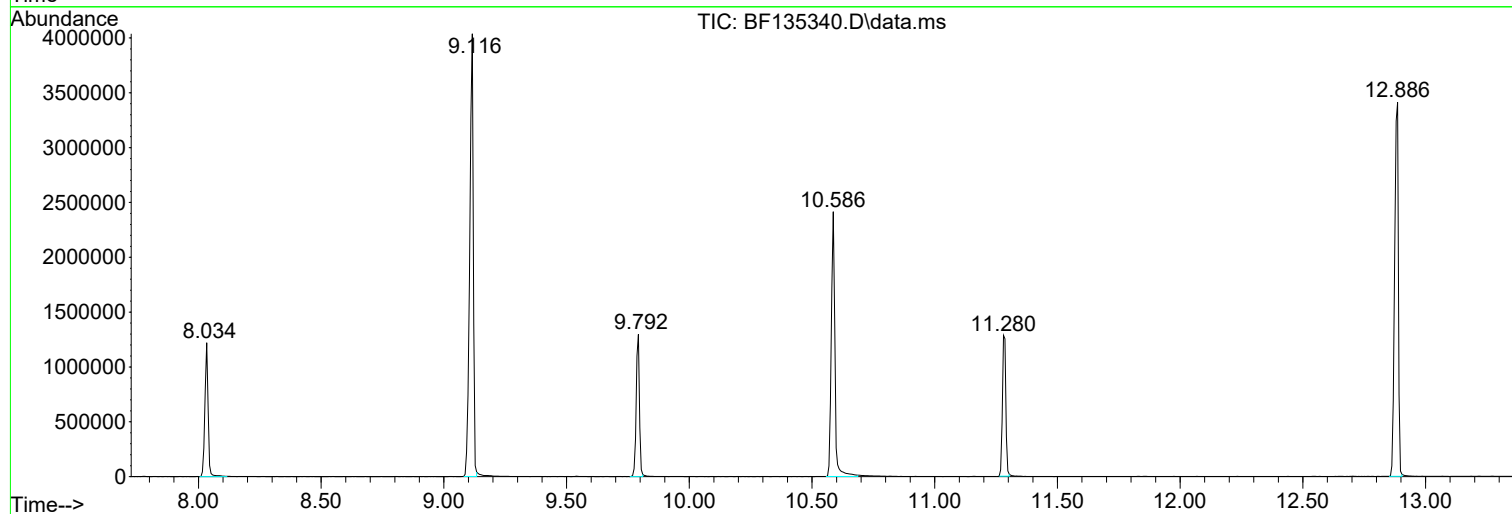
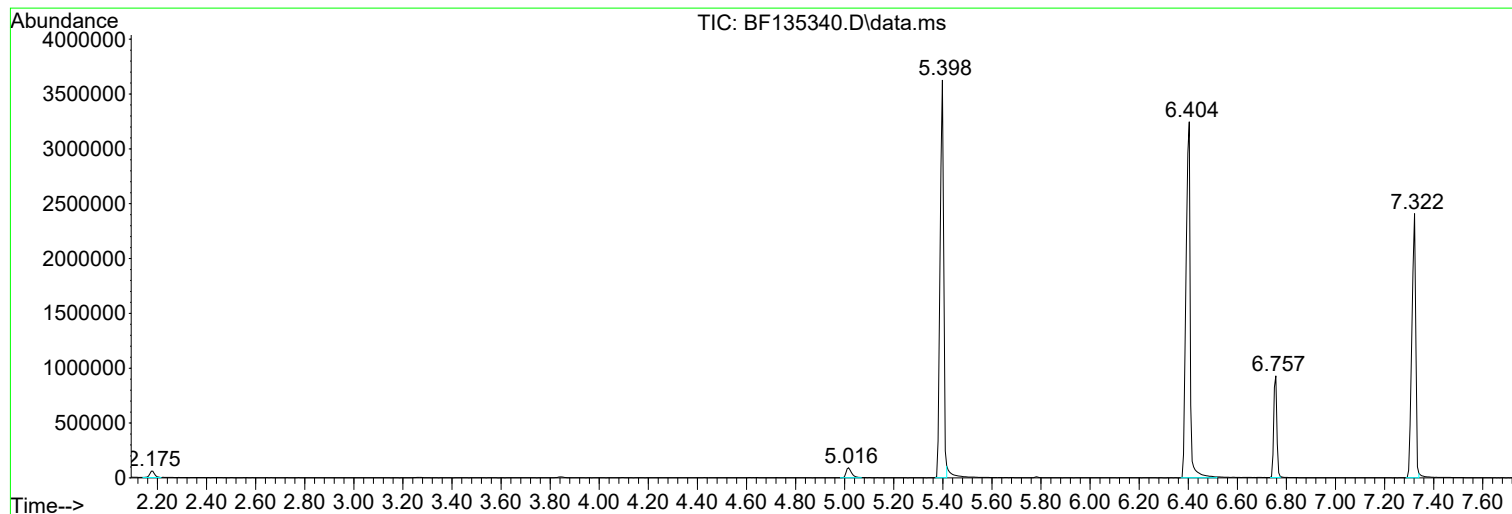
Sum of corrected areas: 26471156

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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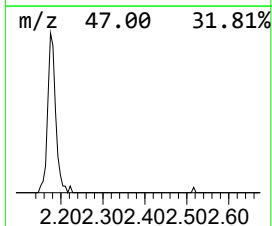
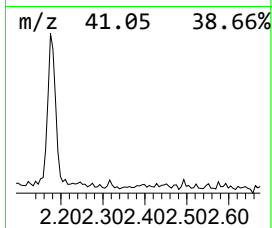
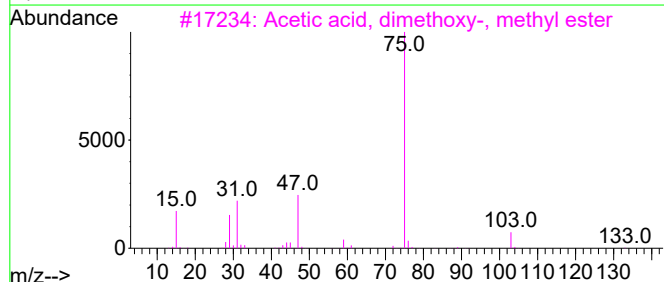
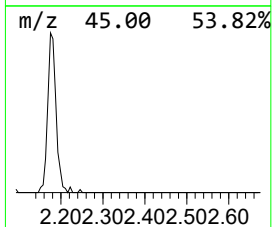
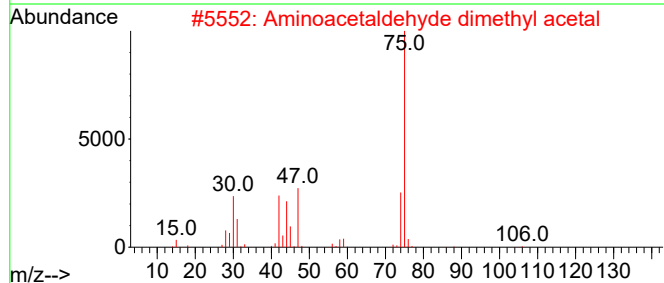
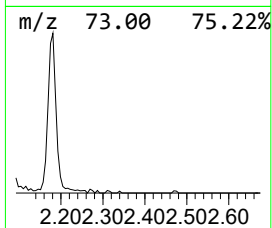
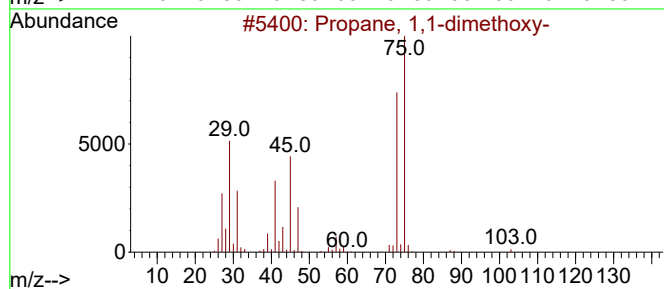
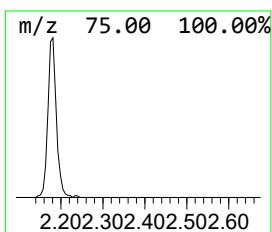
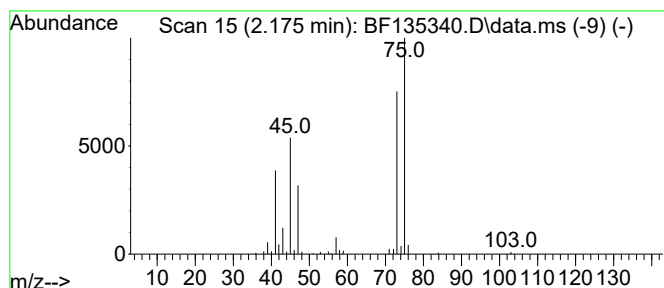
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.175	2.00 ng	82513	1,4-Dichlorobenzene-d4	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	74
2			Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	28
3			Acetic acid, dimethoxy-, methyl ...	134	C5H10O4	000089-91-8	28
4			Methane, trimethoxy-	106	C4H10O3	000149-73-5	28
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	22



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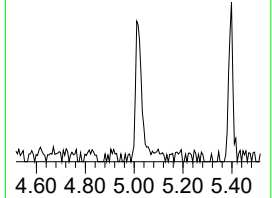
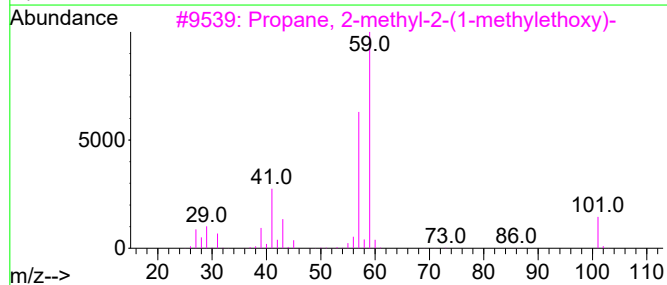
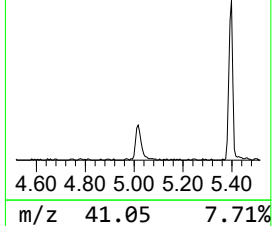
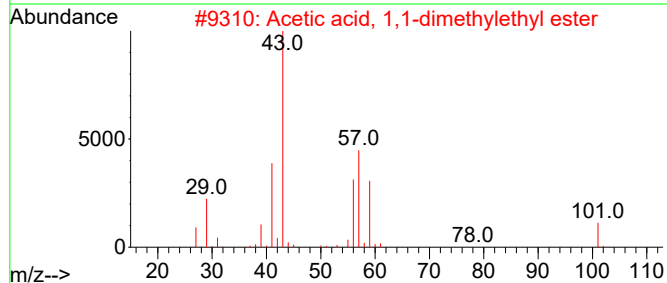
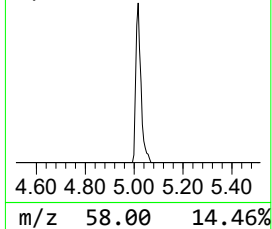
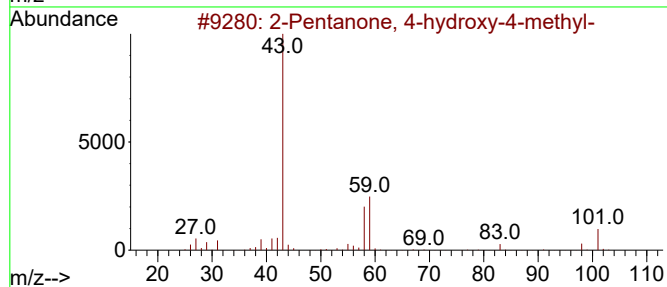
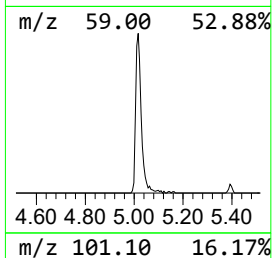
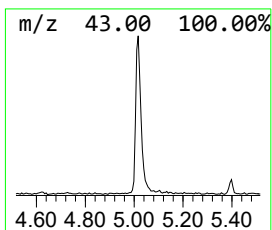
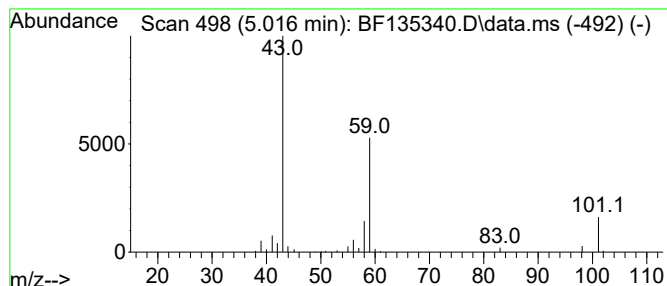
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.016	3.41 ng	140285	1,4-Dichlorobenzene-d4	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38		
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36		
4	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	2.175	2.0	ng	82513	1	6.757	823308	20.0
2-Pentanone, 4-...	5.016	3.4	ng	140285	1	6.757	823308	20.0