

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF013125\
 Data File : BF141337.D
 Acq On : 31 Jan 2025 17:46
 Operator : RC/JU
 Sample : PB166354BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB166354BL

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-BF012925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF141337.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.187	10	16	28	rVB	87009	142392	2.98%	0.482%
2	5.028	491	499	508	rBV	51047	86402	1.81%	0.292%
3	5.416	558	565	570	rBV	2904645	4062975	85.13%	13.751%
4	6.422	730	736	742	rBV	2715302	4053215	84.93%	13.718%
5	6.799	794	800	805	rBV	656301	831802	17.43%	2.815%
6	7.357	888	895	900	rBV	1990075	2724663	57.09%	9.222%
7	8.075	1012	1017	1041	rVB	844042	1128329	23.64%	3.819%
8	9.157	1194	1201	1206	rBV	3443241	4772682	100.00%	16.153%
9	9.834	1310	1316	1336	rVB	996193	1294779	27.13%	4.382%
10	10.622	1444	1450	1476	rBV	2110132	2840595	59.52%	9.614%
11	11.322	1563	1569	1576	rBV	1030731	1299491	27.23%	4.398%
12	12.910	1833	1839	1845	rBV	3401036	4745735	99.44%	16.062%
13	13.963	2013	2018	2030	rBV	643551	847486	17.76%	2.868%
14	15.427	2261	2267	2281	rBV	451233	715764	15.00%	2.423%

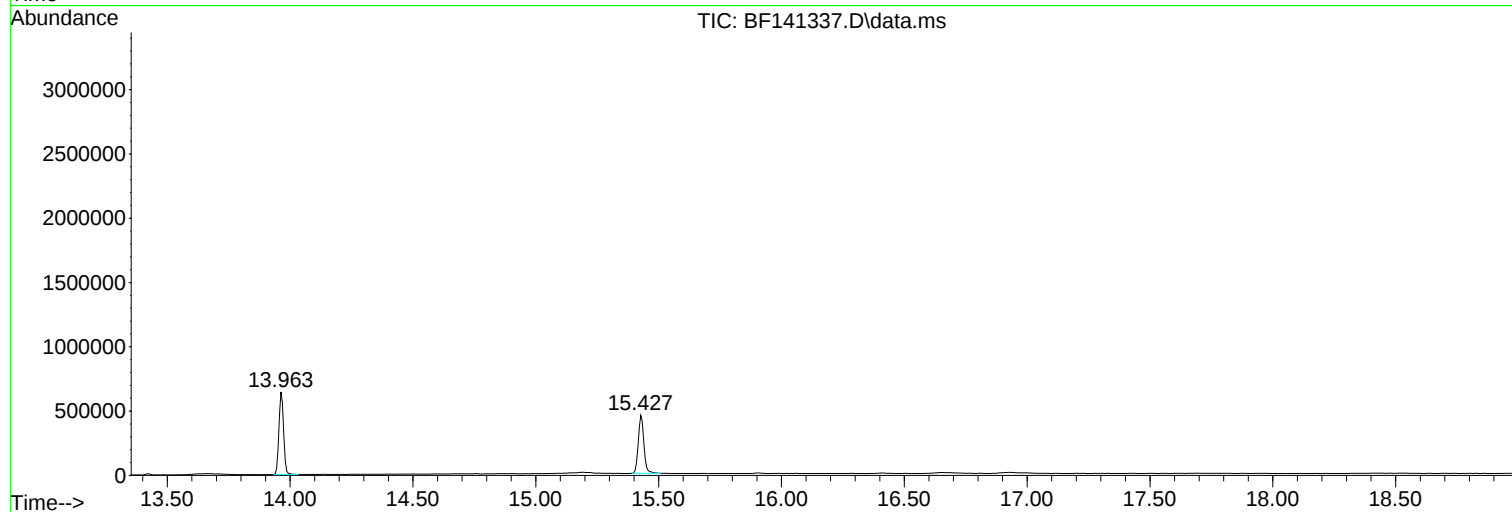
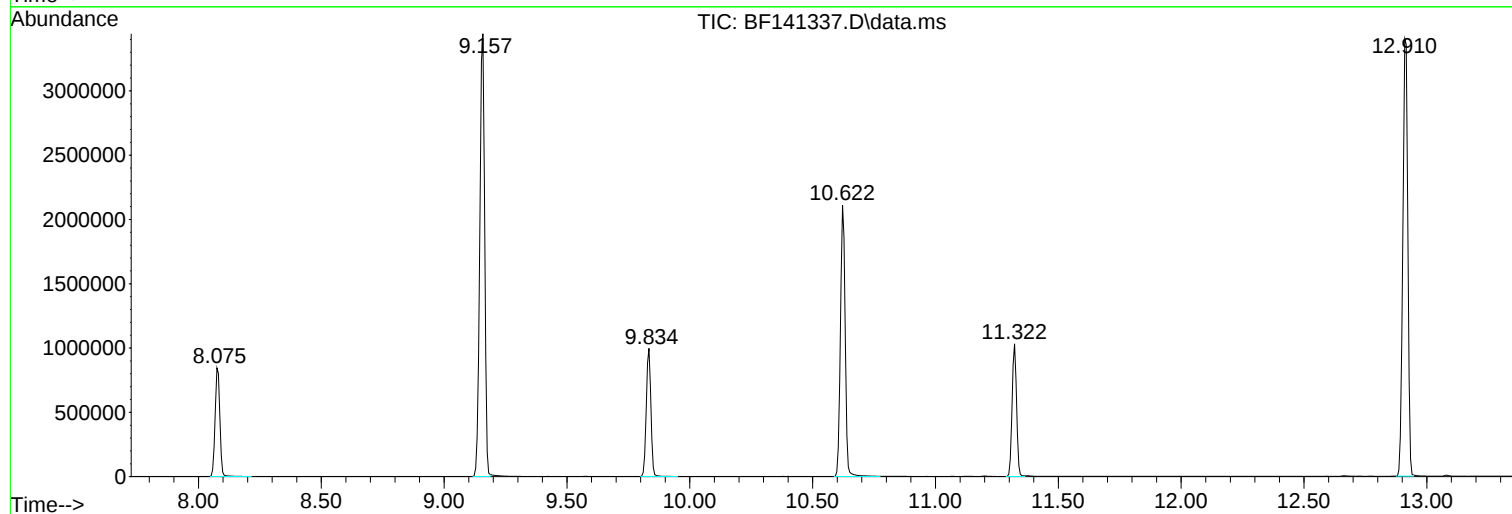
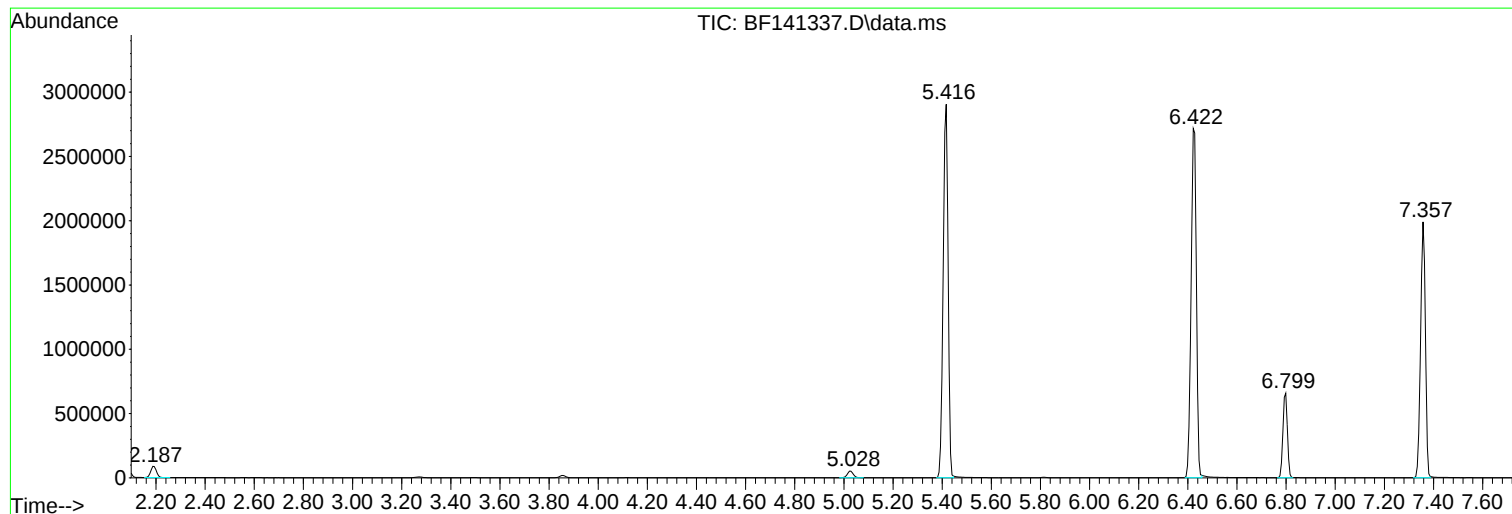
Sum of corrected areas: 29546310

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.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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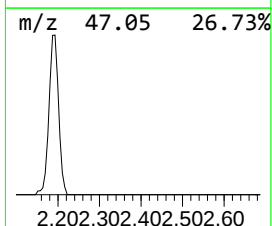
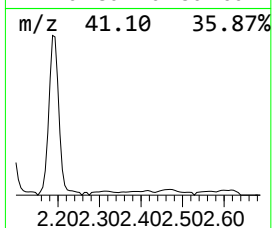
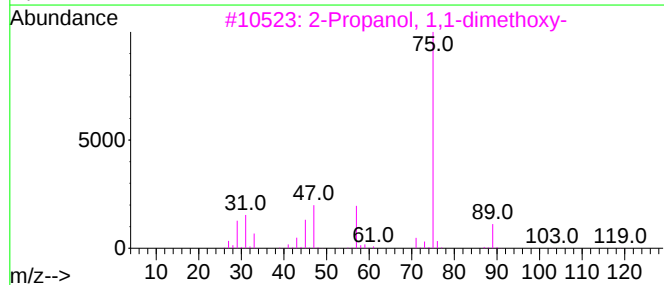
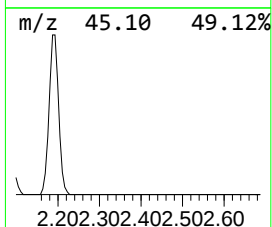
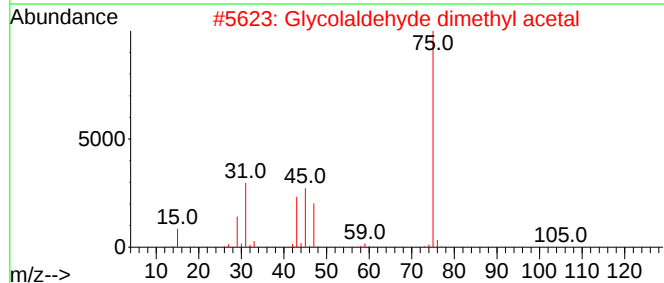
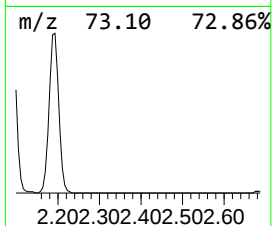
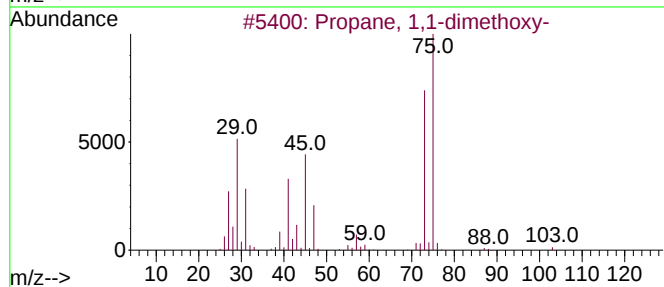
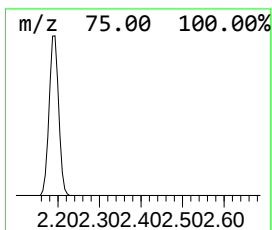
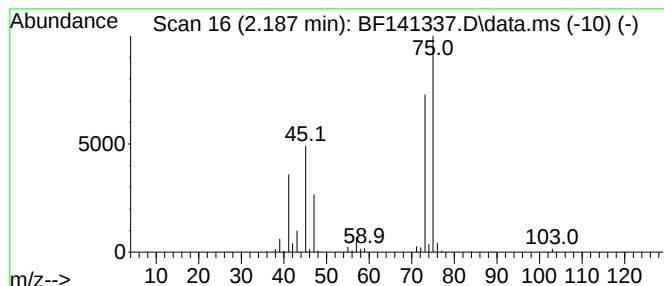
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.187	3.42 ng	142392	1,4-Dichlorobenzene-d4	6.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2			Glycolaldehyde dimethyl acetal	106	C4H10O3	030934-97-5	45
3			2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	42
4			Acetic acid, dimethoxy-, methyl ...	134	C5H10O4	000089-91-8	40
5			Aminoacetaldehyde dimethyl acetal	105	C4H11NO2	022483-09-6	36



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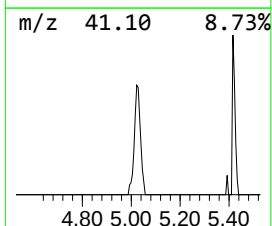
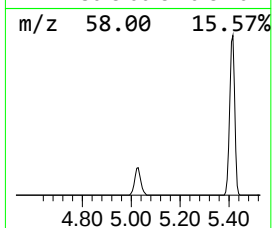
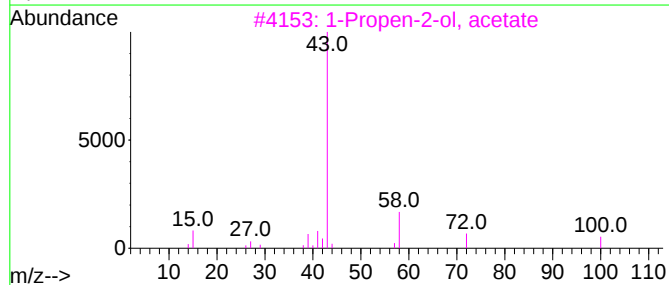
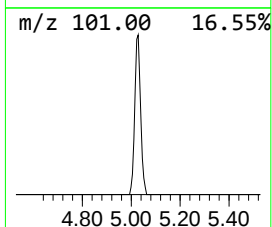
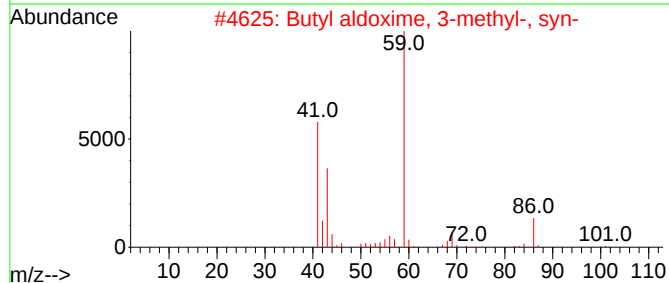
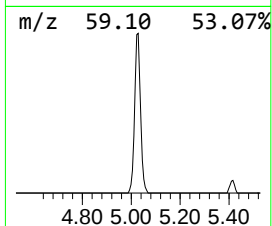
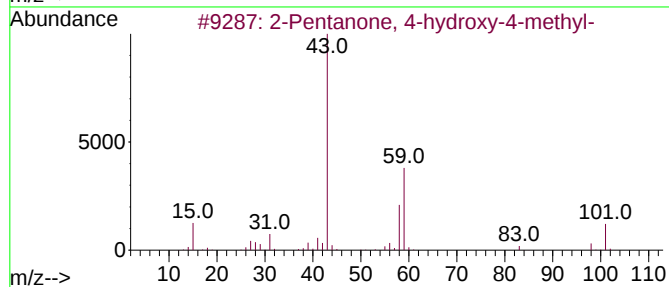
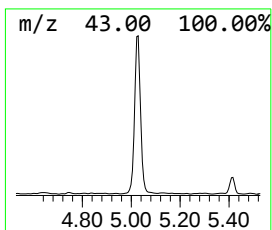
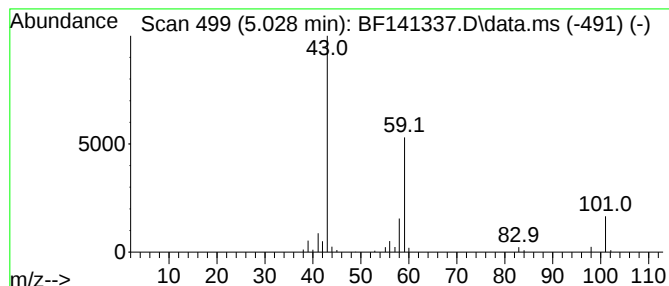
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.028	2.08 ng	86402	1,4-Dichlorobenzene-d4	6.799

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
4		Acetone	58	C3H6O	000067-64-1	12
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	2.187	3.4	ng	142392	1	6.799	831802	20.0
2-Pentanone, 4-...	5.028	2.1	ng	86402	1	6.799	831802	20.0