

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090425\
 Data File : BF143647.D
 Acq On : 04 Sep 2025 16:41
 Operator : RC/JU
 Sample : PB169499BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB169499BL

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-BF090325.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF143647.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.369	23	30	37	rVB	30887	49104	1.88%	0.328%
2	5.134	491	500	509	rBV	146053	236755	9.06%	1.581%
3	5.516	558	565	570	rBV	1380260	1967522	75.27%	13.141%
4	6.504	726	733	739	rBV	1257163	1877588	71.83%	12.541%
5	6.887	793	798	803	rBV	309309	386502	14.79%	2.582%
6	7.445	887	893	898	rBV	897002	1259049	48.17%	8.409%
7	8.169	1010	1016	1021	rBV	401027	521124	19.94%	3.481%
8	9.245	1192	1199	1204	rBV	1934275	2614030	100.00%	17.459%
9	9.922	1308	1314	1319	rBV	483072	614129	23.49%	4.102%
10	10.710	1442	1448	1453	rBV	1171393	1520957	58.18%	10.159%
11	11.410	1561	1567	1572	rBV	495872	611924	23.41%	4.087%
12	12.998	1831	1837	1842	rBV	1944232	2540739	97.20%	16.970%
13	14.045	2010	2015	2025	rBV	328611	409619	15.67%	2.736%
14	15.521	2259	2266	2278	rBV	237877	362947	13.88%	2.424%

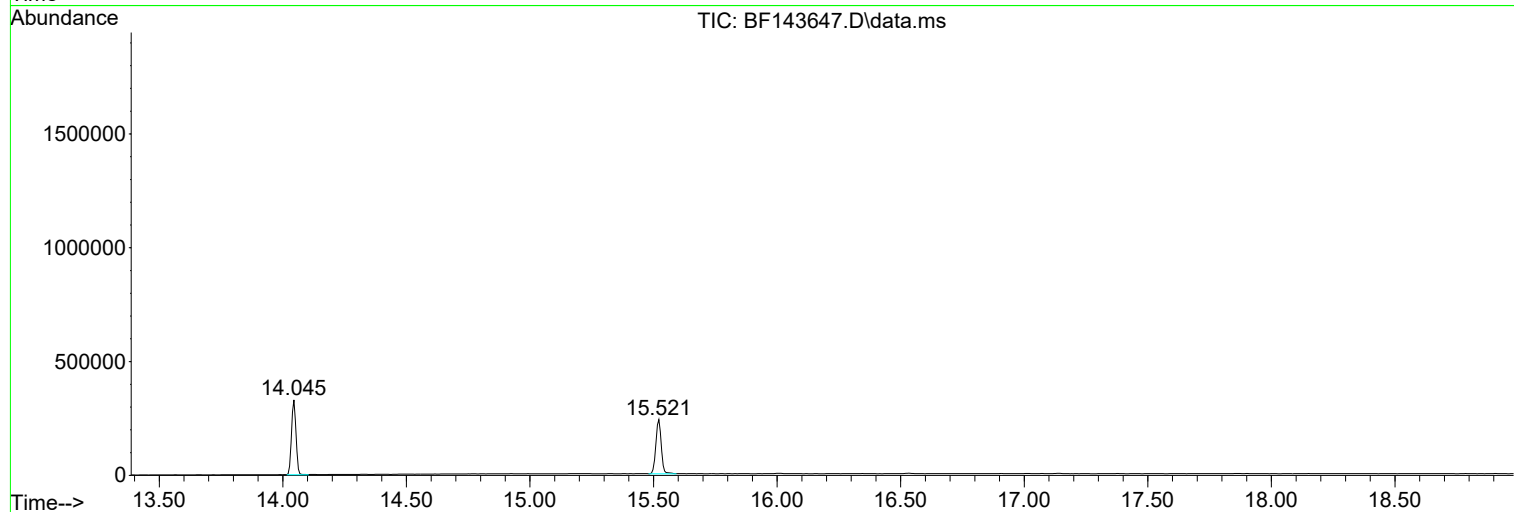
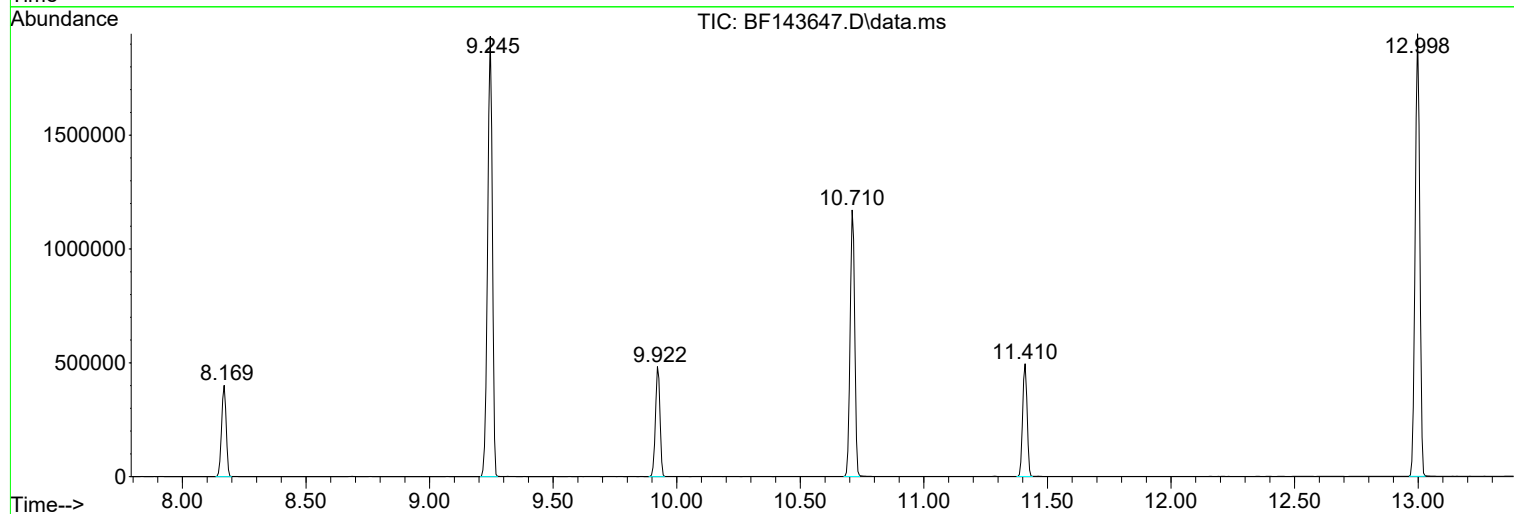
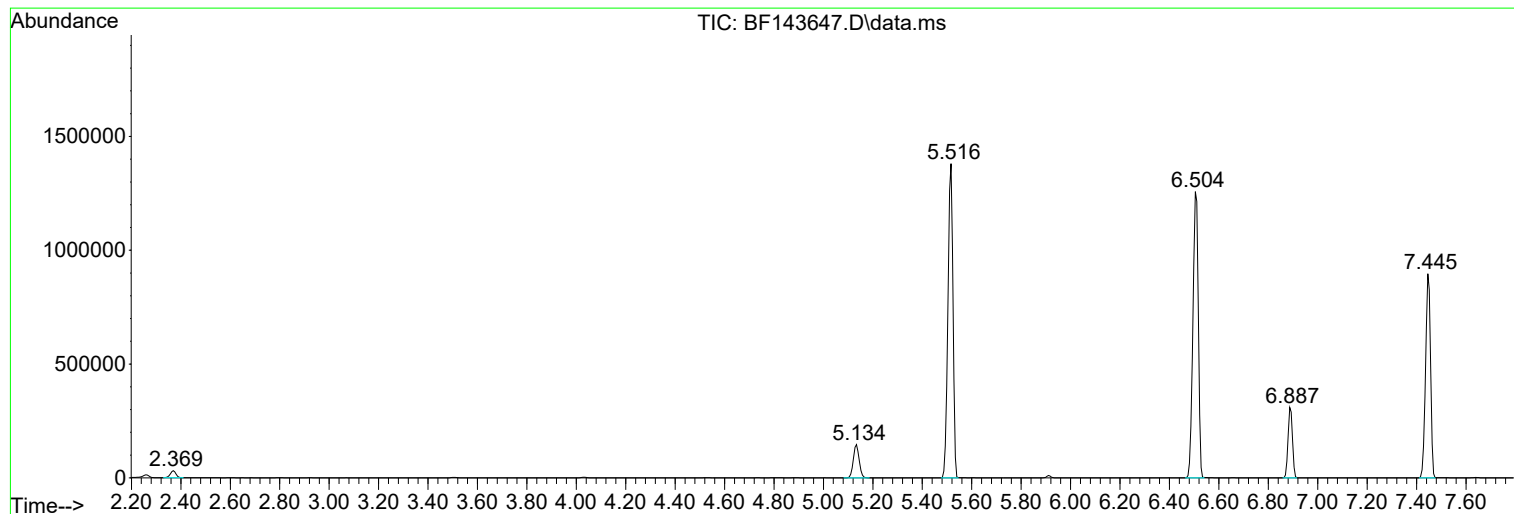
Sum of corrected areas: 14971989

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090325.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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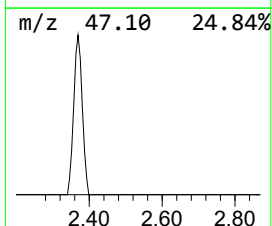
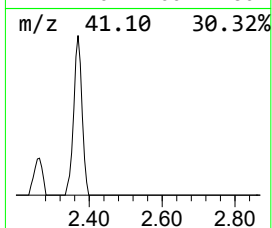
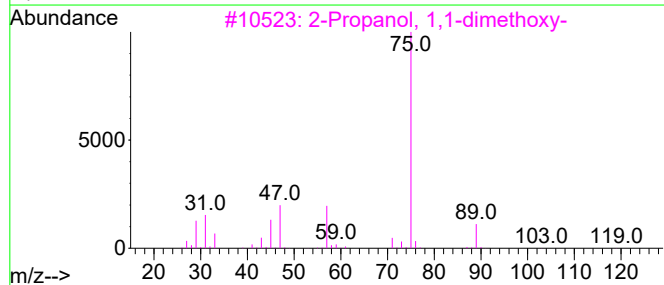
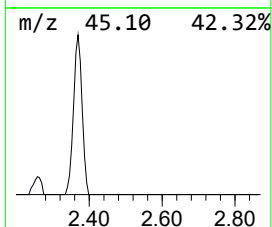
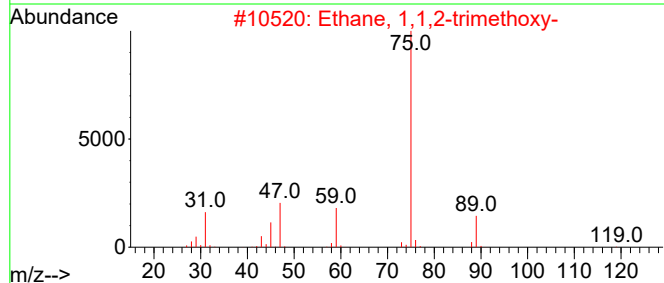
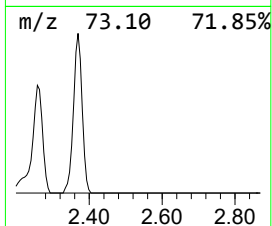
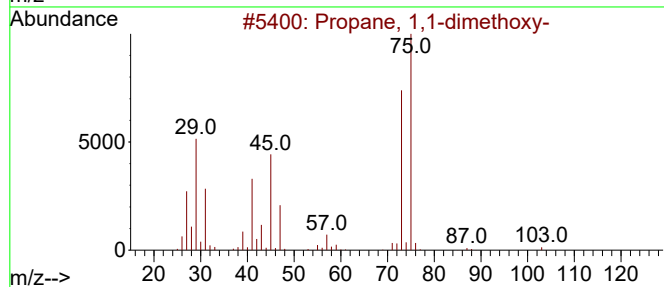
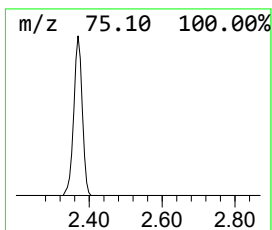
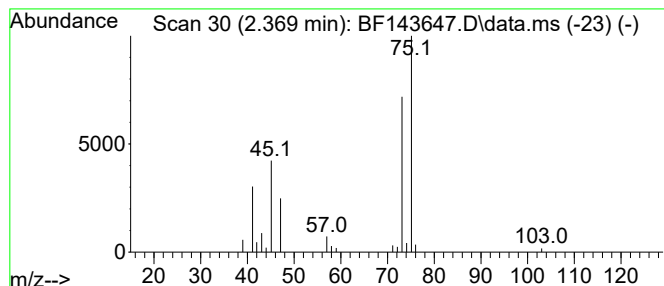
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090325.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.369	2.54 ng	49104	1,4-Dichlorobenzene-d4	6.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2			Ethane, 1,1,2-trimethoxy-	120	C5H12O3	024332-20-5	36
3			2-Propanol, 1,1-dimethoxy-	120	C5H12O3	042919-42-6	33
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9



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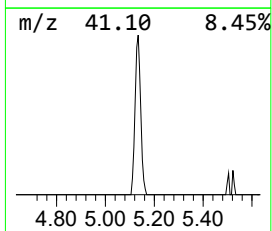
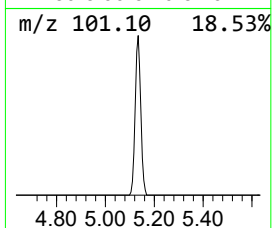
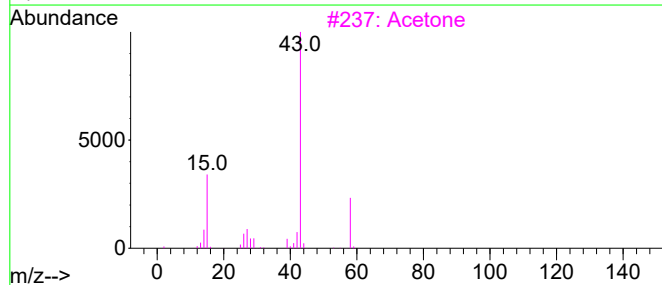
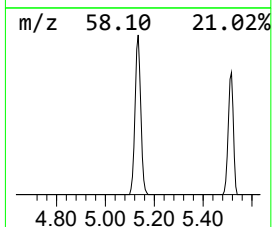
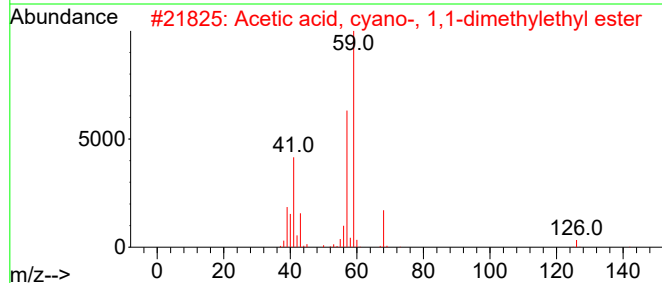
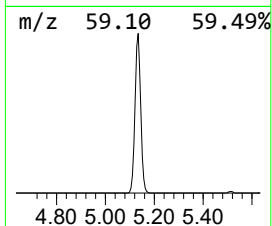
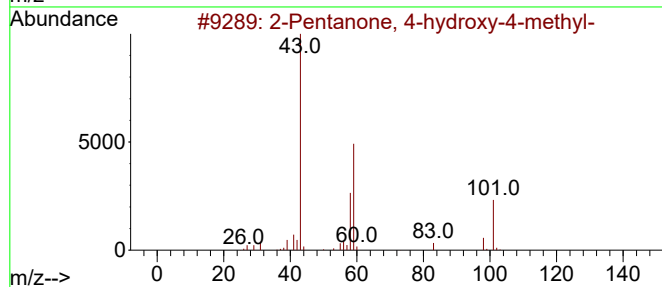
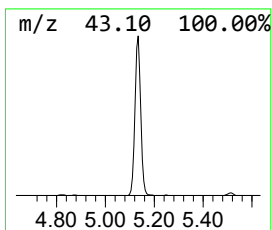
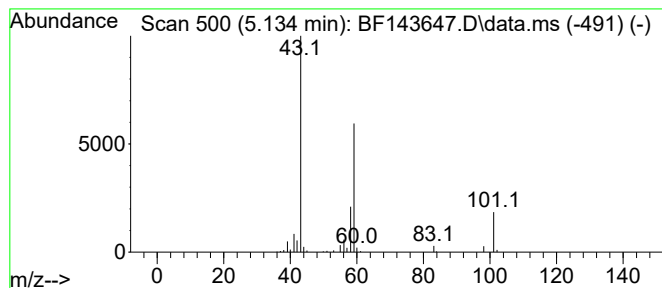
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.134	12.25 ng	236755	1,4-Dichlorobenzene-d4	6.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
3			Acetone	58	C3H6O	000067-64-1	16
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-di...	2.369	2.5	ng	49104	1	6.887	386502	20.0
2-Pentanone, 4-...	5.134	12.3	ng	236755	1	6.887	386502	20.0