

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090622\
 Data File : BF130139.D
 Acq On : 06 Sep 2022 15:59
 Operator : CG\JU
 Sample : PB147437BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB147437BL

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-BF083122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130139.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.910	441	446	458	rBV	400965	593856	8.30%	1.366%
2	5.310	508	514	517	rBV	4957830	5435517	75.94%	12.506%
3	6.328	681	687	690	rBV	4370218	5437846	75.97%	12.511%
4	6.693	746	749	753	rVB	1548110	1312151	18.33%	3.019%
5	7.263	840	846	849	rBV	3467155	3468798	48.46%	7.981%
6	7.975	962	967	971	rBV	2042949	1797006	25.10%	4.134%
7	9.057	1145	1151	1154	rBV	6600987	6661355	93.06%	15.326%
8	9.728	1259	1265	1268	rBV	2508233	2116771	29.57%	4.870%
9	10.516	1392	1399	1402	rBV	4597872	4364062	60.97%	10.040%
10	11.204	1512	1516	1520	rBV	2551931	2198792	30.72%	5.059%
11	12.798	1782	1787	1790	rBV	7223258	7158113	100.00%	16.469%
12	13.833	1959	1963	1967	rBV	1968242	1692935	23.65%	3.895%
13	15.233	2196	2201	2207	rBV	1103984	1227718	17.15%	2.825%

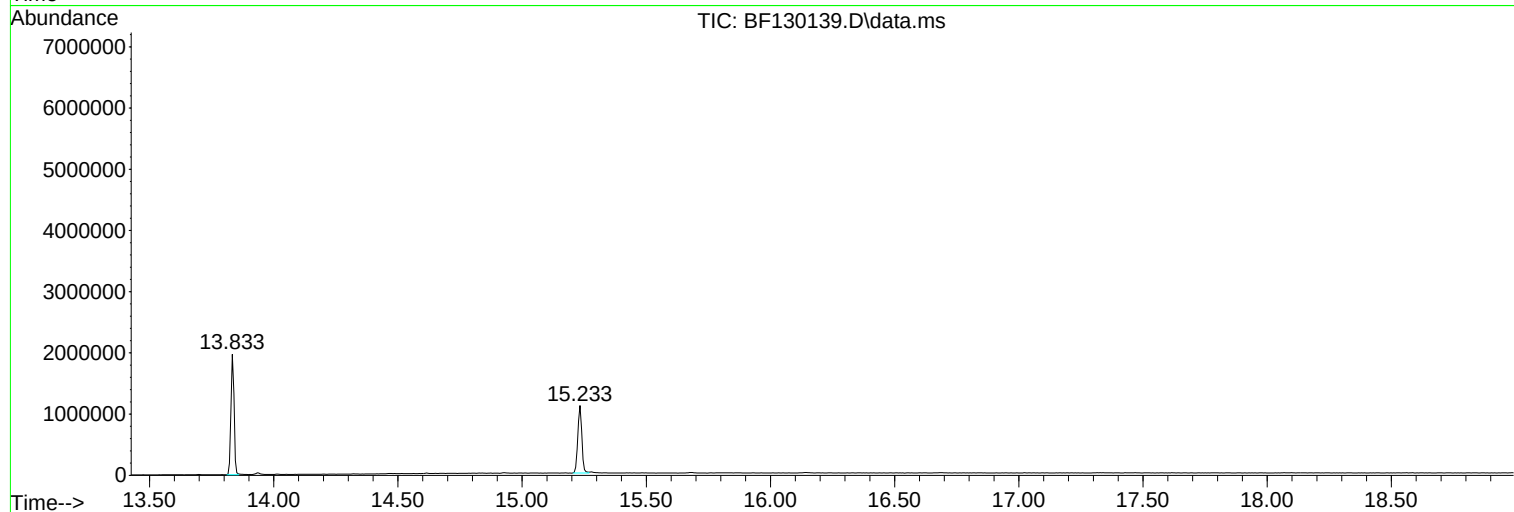
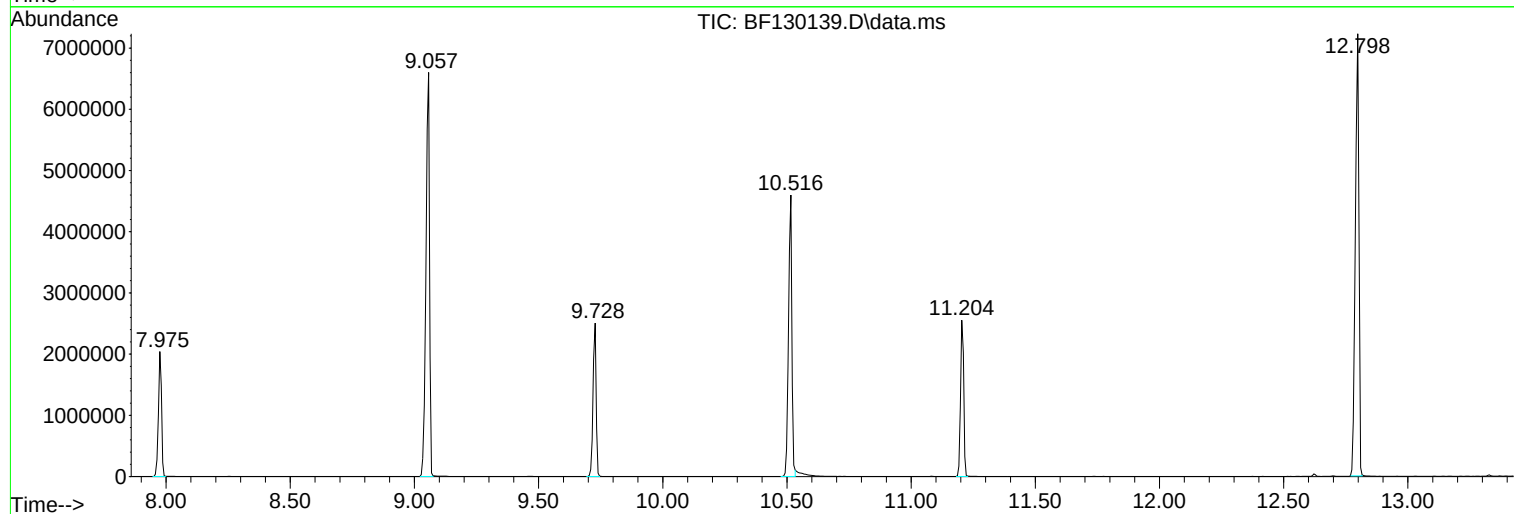
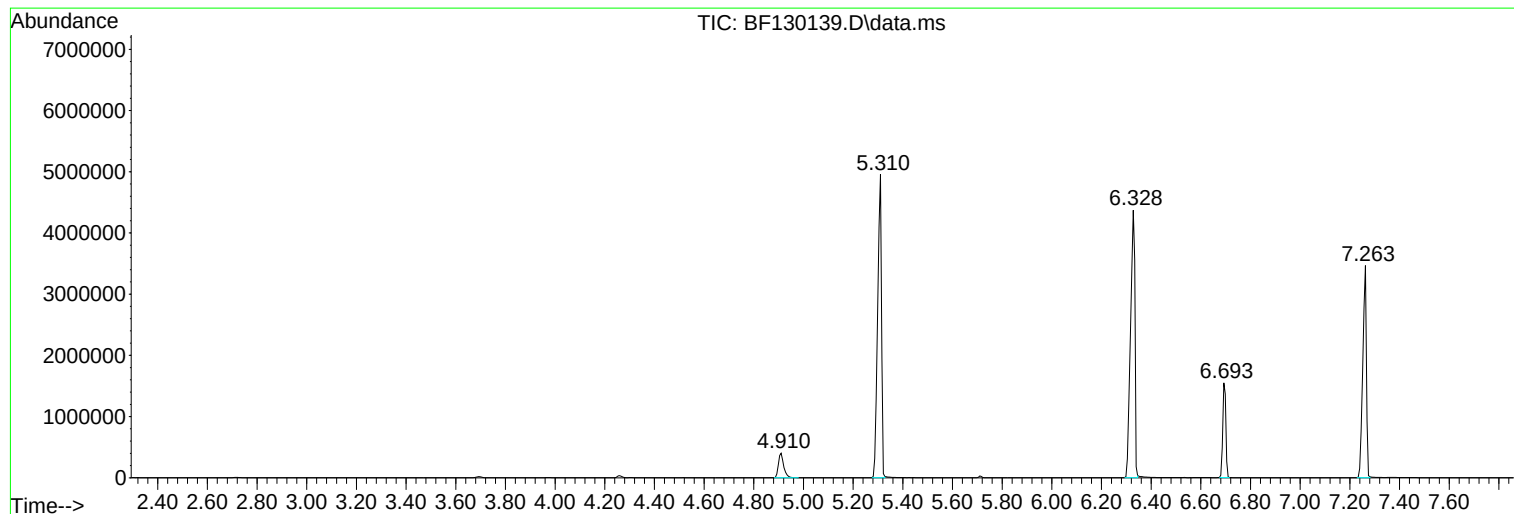
Sum of corrected areas: 43464920

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.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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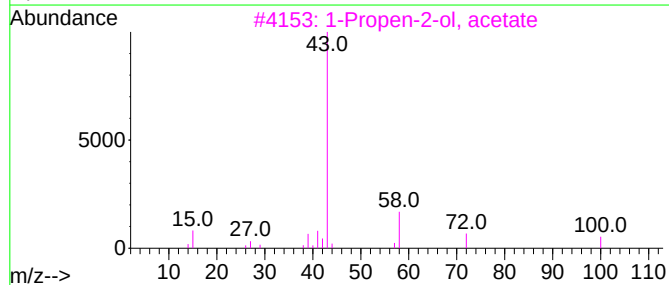
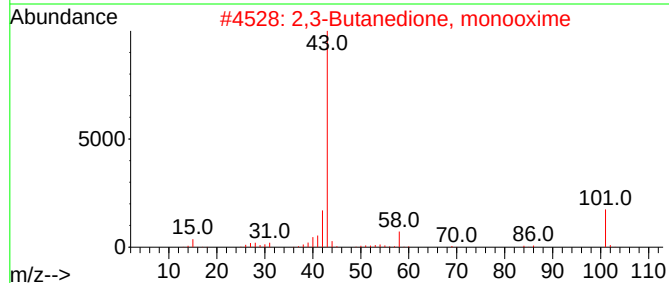
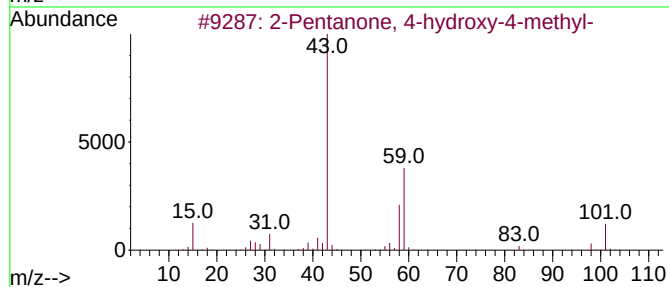
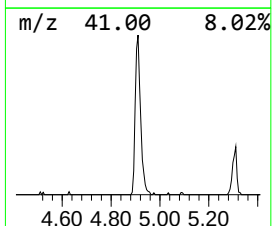
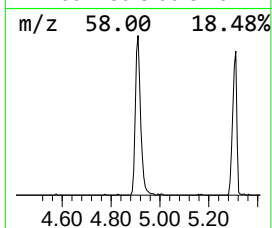
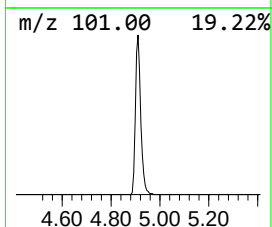
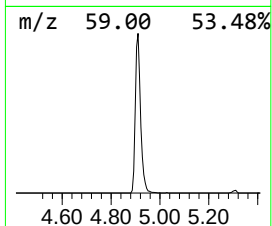
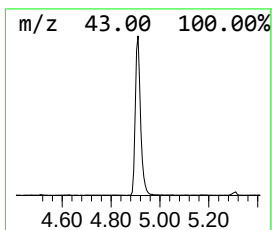
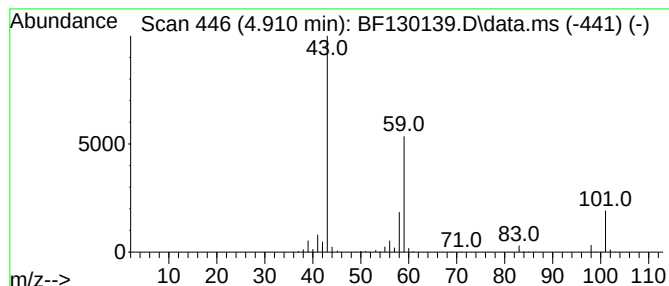
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.910	9.05 ng	593856	1,4-Dichlorobenzene-d4	6.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
4	Acetone	58	C3H6O	000067-64-1	9		
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.910	9.1	ng	593856	1	6.693	1312150	20.0