

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090722\
 Data File : BF130162.D
 Acq On : 07 Sep 2022 15:08
 Operator : CG\JU
 Sample : PB147513BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB147513BL

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: BF130162.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.916	442	447	462	rBV	366357	617123	6.94%	1.248%
2	5.310	508	514	517	rBV	5234971	6364075	71.58%	12.865%
3	6.334	681	688	691	rBV	4759527	6125120	68.89%	12.382%
4	6.692	746	749	753	rVB	1560969	1301339	14.64%	2.631%
5	7.263	840	846	849	rBV	3732474	4134018	46.50%	8.357%
6	7.975	962	967	971	rBV	1985217	1776807	19.99%	3.592%
7	9.057	1144	1151	1154	rBV	6977822	7532106	84.72%	15.226%
8	9.727	1259	1265	1268	rBV	2093126	2057957	23.15%	4.160%
9	10.516	1392	1399	1402	rBV	5183978	5197406	58.46%	10.507%
10	11.204	1512	1516	1520	rBV	2489560	2103735	23.66%	4.253%
11	12.798	1781	1787	1790	rBV	7625662	8890596	100.00%	17.972%
12	13.833	1959	1963	1967	rBV	2167355	1846138	20.77%	3.732%
13	15.233	2196	2201	2206	rBV	1342190	1521710	17.12%	3.076%

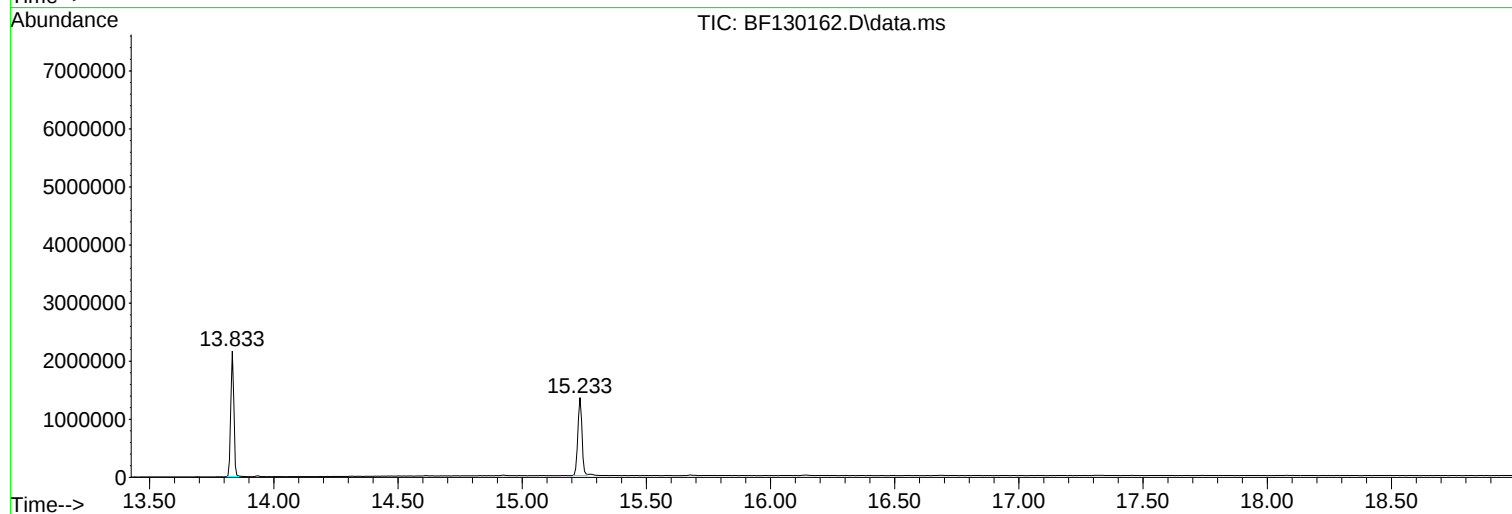
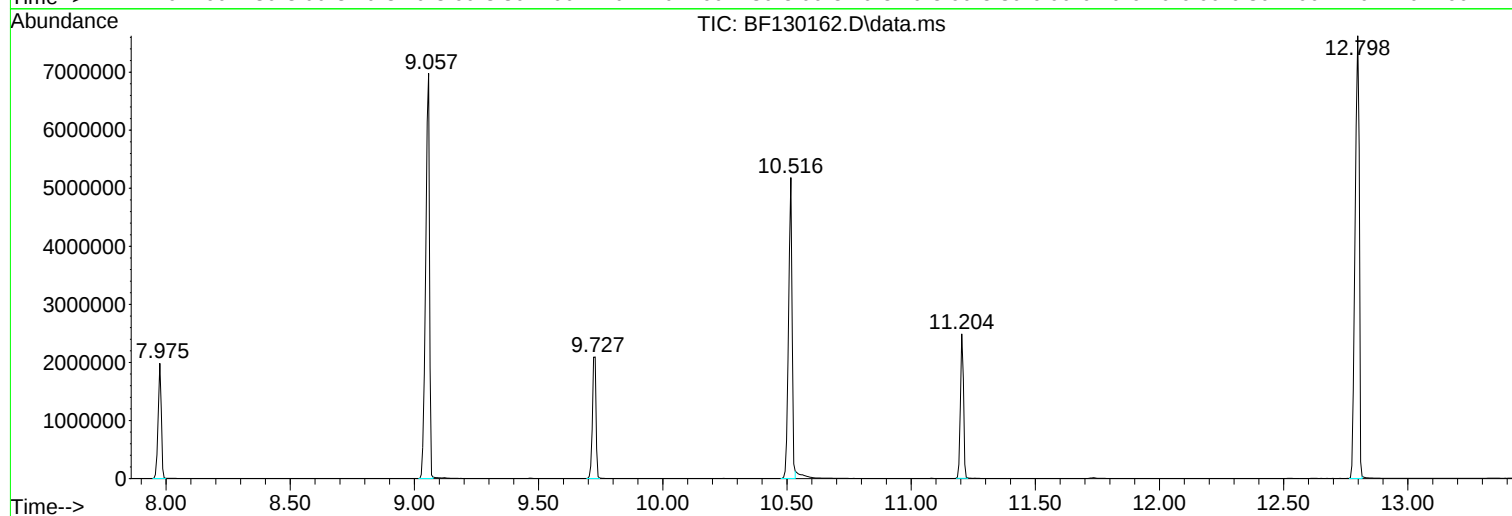
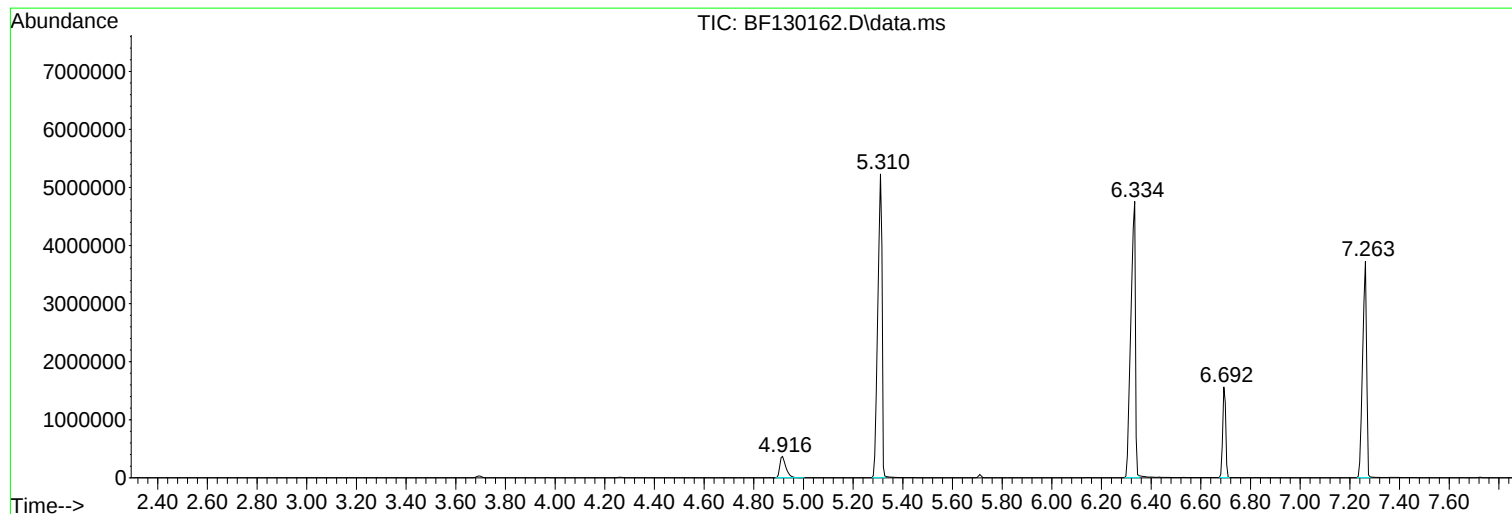
Sum of corrected areas: 49468130

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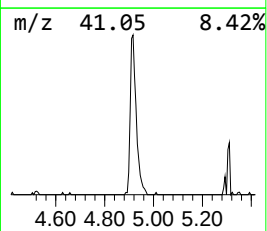
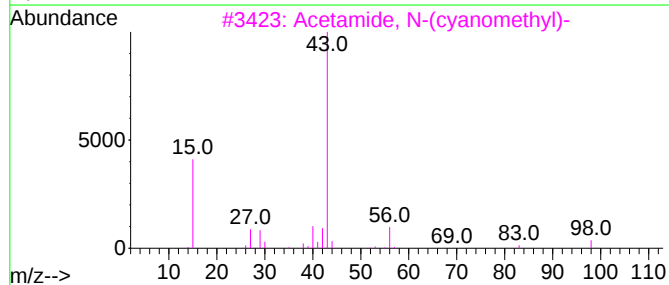
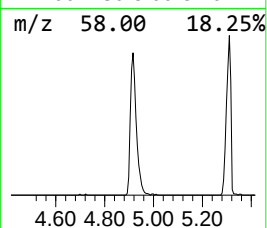
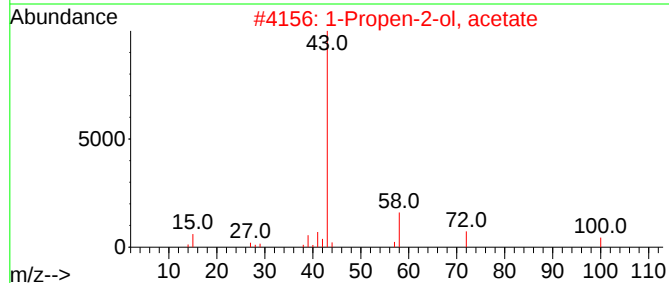
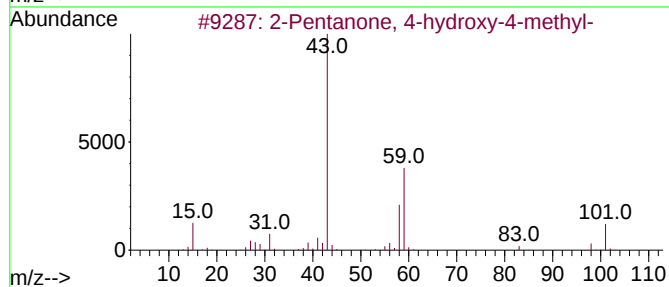
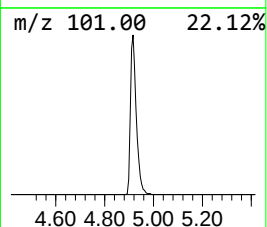
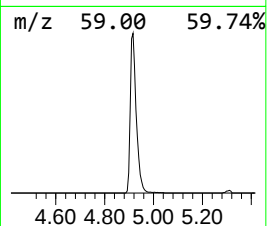
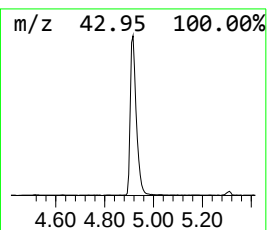
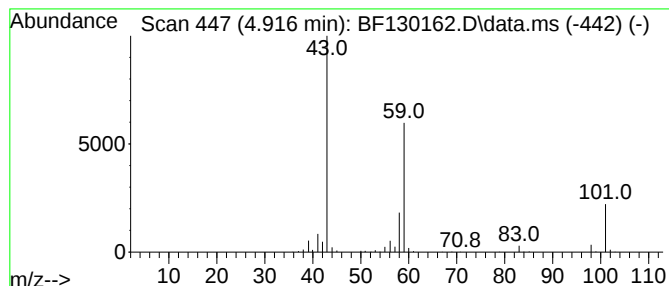
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.916	9.48 ng	617123	1,4-Dichlorobenzene-d4	6.692

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10	
3	Acetamide, N-(cyanomethyl)-	98	C4H6N2O	004814-80-6	9	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	
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
2-Pentanone, 4-...	4.916	9.5	ng	617123	1	6.692	1301340	20.0