

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF040722\
 Data File : BF127677.D
 Acq On : 07 Apr 2022 11:53
 Operator : CG\JU
 Sample : PB143795BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB143795BL

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-BF040622.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF127677.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.222	496	499	507	rVB	59072	81363	1.64%	0.291%
2	5.593	556	562	565	rBV	3157801	3476278	69.92%	12.452%
3	6.581	723	730	733	rBV	3091724	3553969	71.48%	12.730%
4	6.963	791	795	798	rBV	982311	754075	15.17%	2.701%
5	7.528	885	891	894	rBV	2353837	2388001	48.03%	8.554%
6	8.245	1008	1013	1017	rBV	1116450	970505	19.52%	3.476%
7	9.328	1190	1197	1200	rBV	4537064	4323168	86.96%	15.485%
8	10.010	1307	1313	1316	rBV	1281844	1151866	23.17%	4.126%
9	10.798	1440	1447	1450	rBV	3047234	2920190	58.74%	10.460%
10	11.498	1562	1566	1570	rBV	1375260	1230004	24.74%	4.406%
11	13.098	1832	1838	1841	rBV	4953908	4971637	100.00%	17.808%
12	14.151	2012	2017	2020	rBV	1270787	1144827	23.03%	4.101%
13	15.674	2270	2276	2287	rBV	748115	952187	19.15%	3.411%

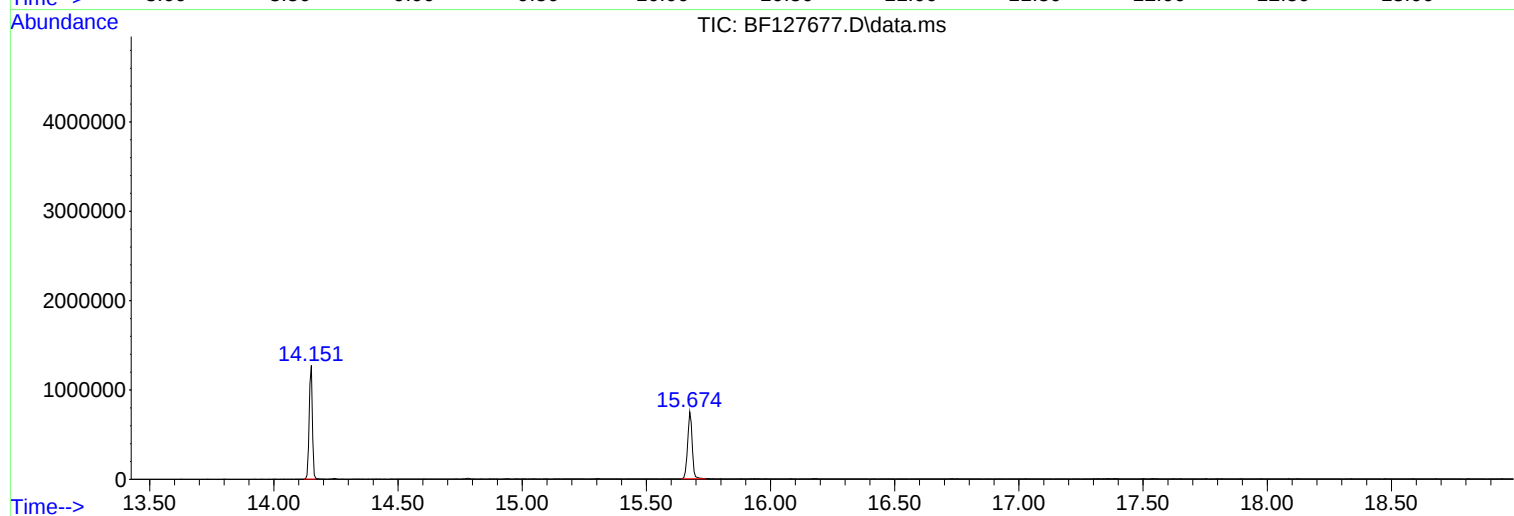
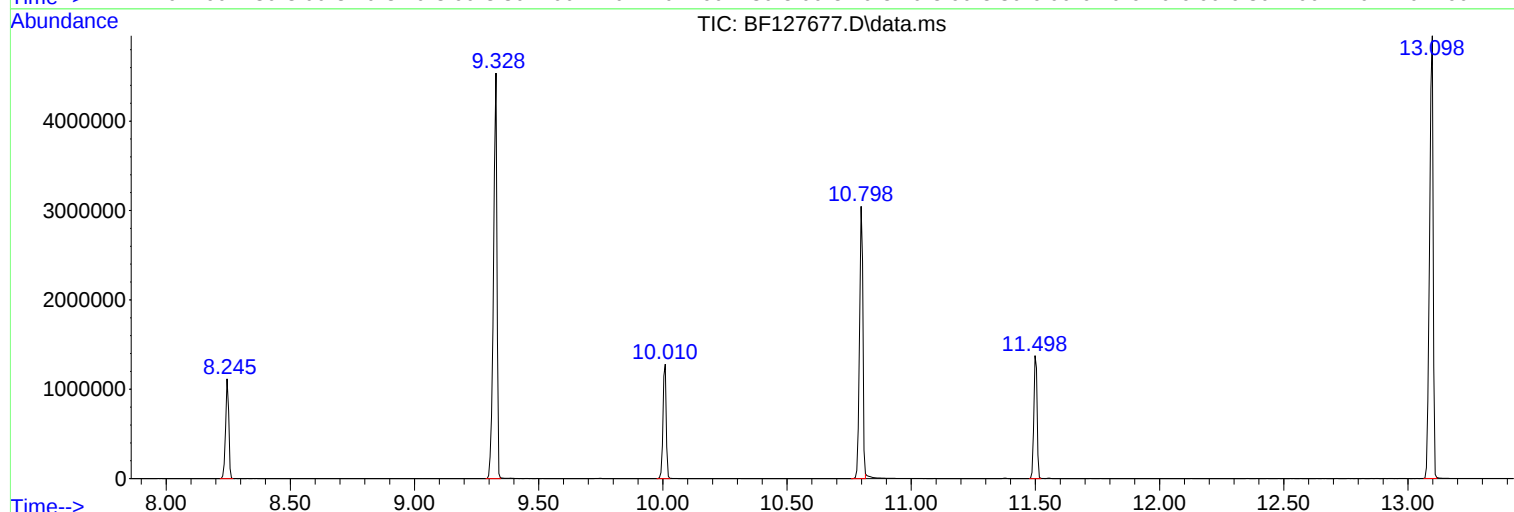
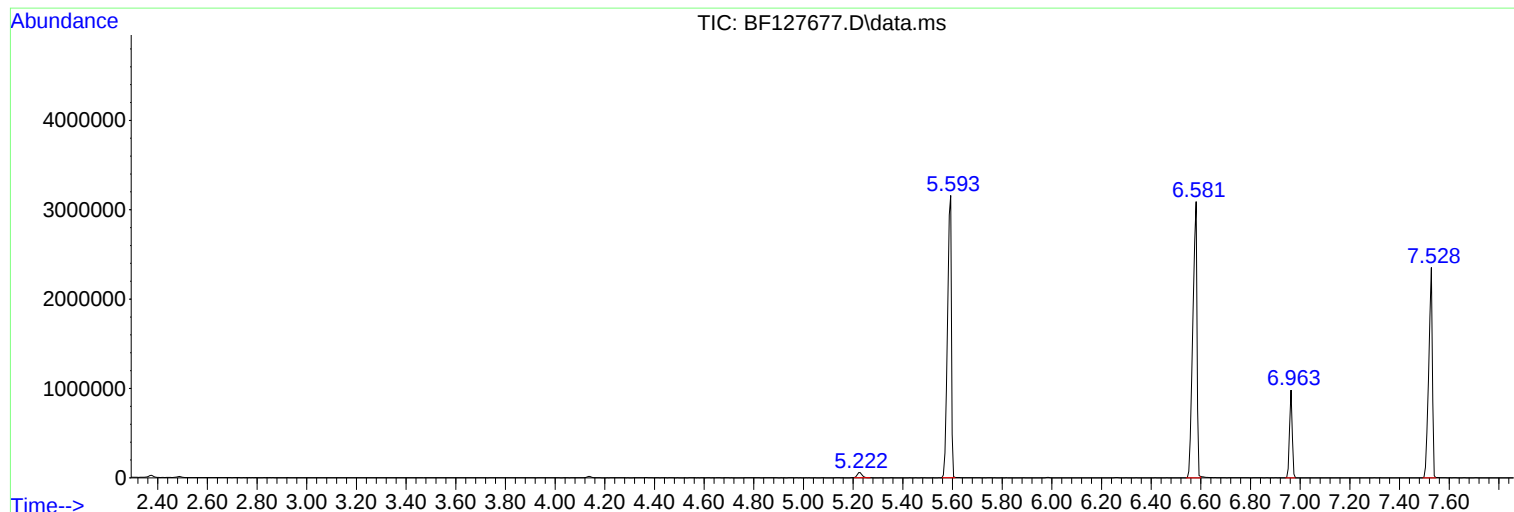
Sum of corrected areas: 27918070

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.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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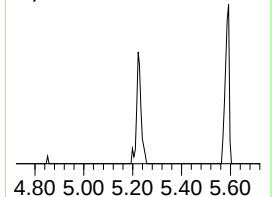
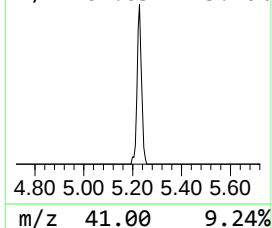
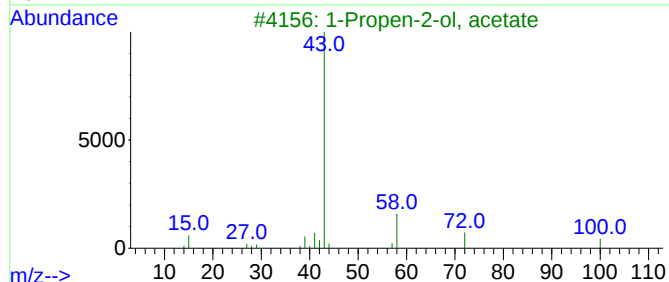
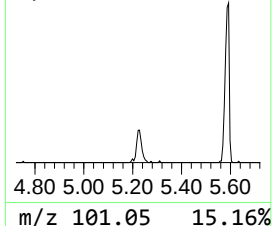
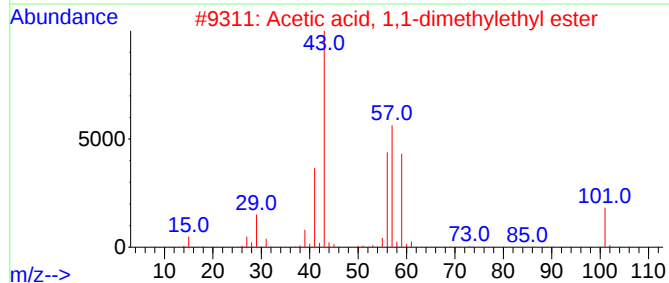
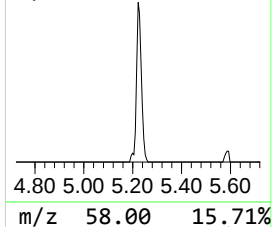
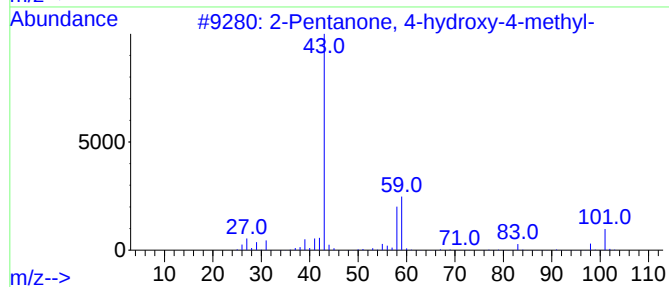
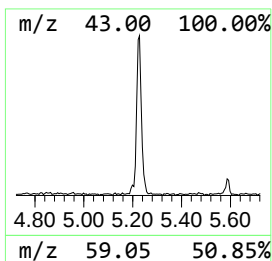
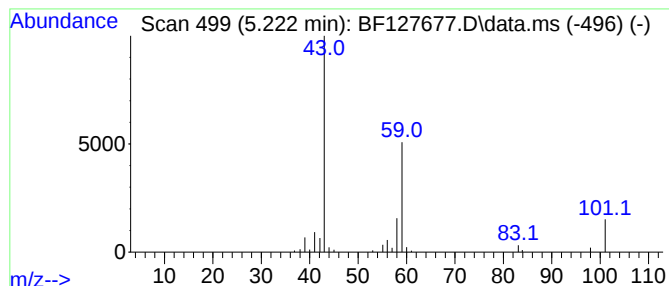
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.222	2.16 ng	81363	1,4-Dichlorobenzene-d4	6.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	45		
3	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	Acetamide, N-(cyanomethyl)-	98	C4H6N2O	004814-80-6	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.222	2.2	ng	81363	1	6.963	754075	20.0