

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
 Data File : BF130284.D
 Acq On : 16 Sep 2022 10:43
 Operator : CG\JU
 Sample : PB147633BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB147633BL

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-BF091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130284.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.881	470	475	485	rBV	145383	214692	5.52%	0.961%
2	5.287	538	544	547	rBV	2742409	2854269	73.44%	12.776%
3	6.310	712	718	721	rBV	2389945	2882730	74.17%	12.903%
4	6.675	776	780	783	rBV	797215	605907	15.59%	2.712%
5	7.245	871	877	879	rBV	1698661	1868544	48.07%	8.364%
6	7.957	993	998	1001	rBV	961475	804755	20.70%	3.602%
7	9.039	1176	1182	1185	rBV	3725920	3634424	93.51%	16.268%
8	9.704	1290	1295	1299	rBV	1004767	916994	23.59%	4.105%
9	10.498	1424	1430	1433	rBV	2585420	2538129	65.30%	11.361%
10	11.186	1543	1547	1550	rBV	1158200	939077	24.16%	4.203%
11	12.774	1812	1817	1820	rBV	3928953	3886768	100.00%	17.398%
12	13.815	1990	1994	1997	rBV	820928	699621	18.00%	3.132%
13	15.209	2226	2231	2236	rBV	422828	494816	12.73%	2.215%

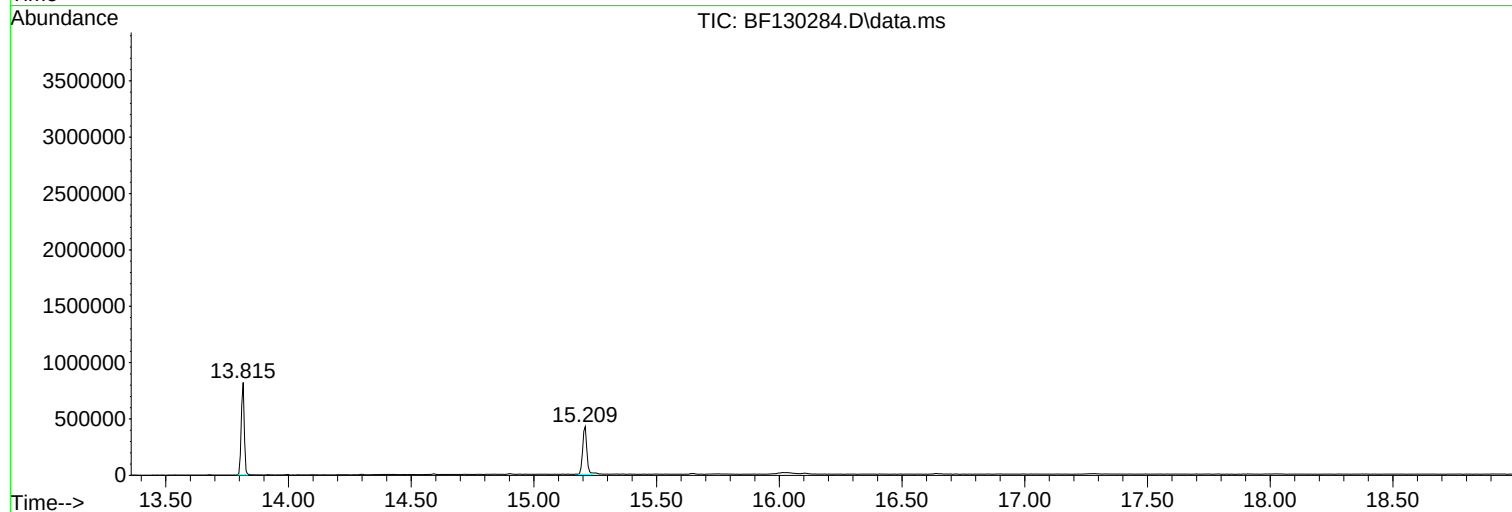
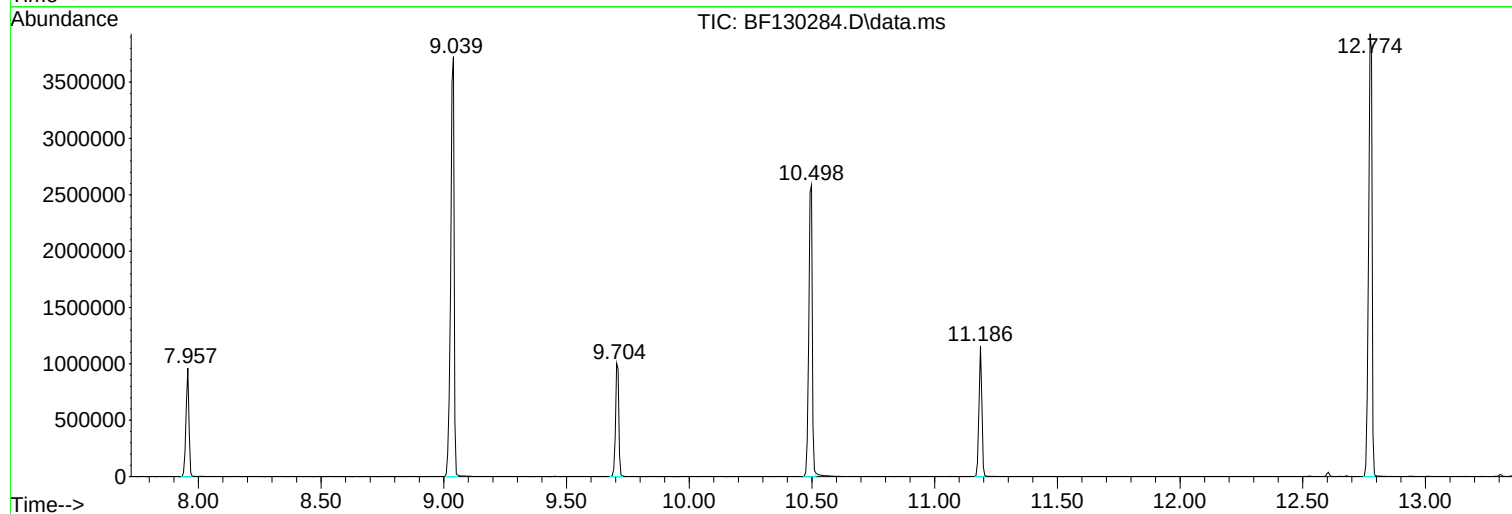
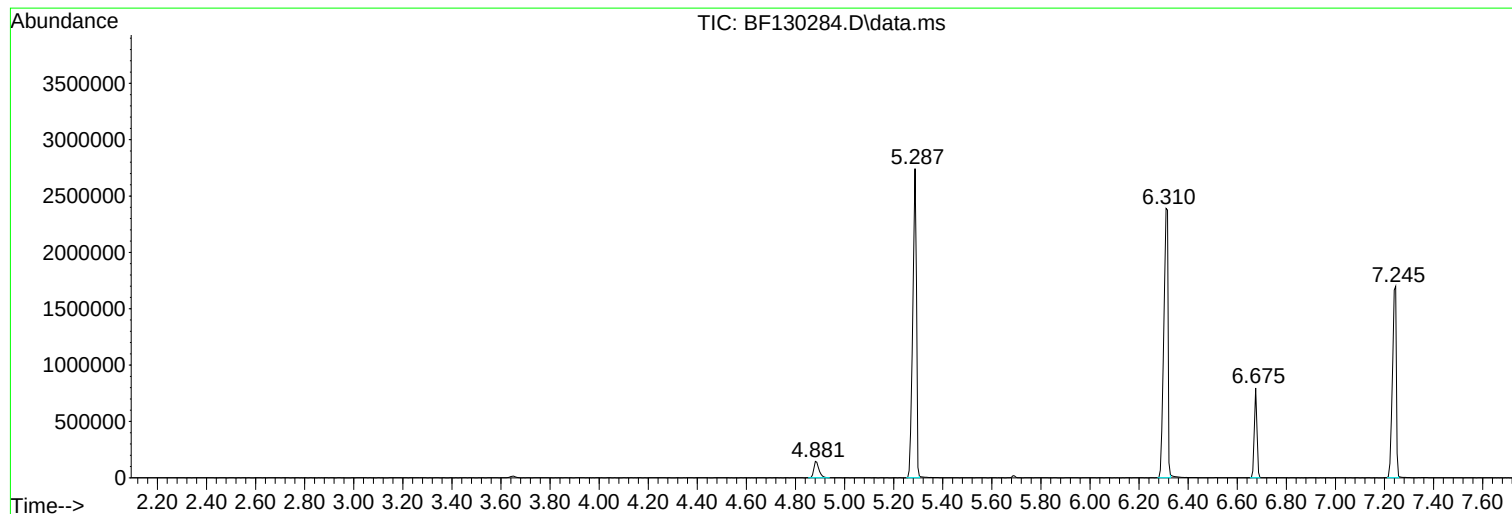
Sum of corrected areas: 22340726

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.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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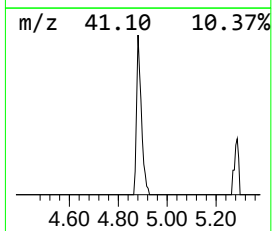
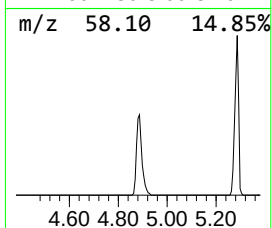
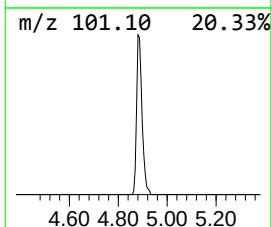
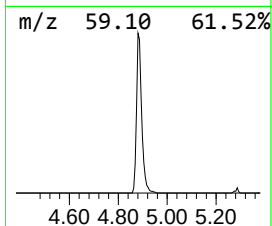
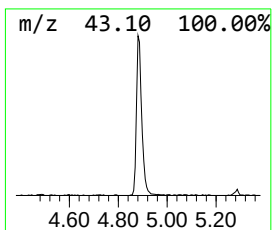
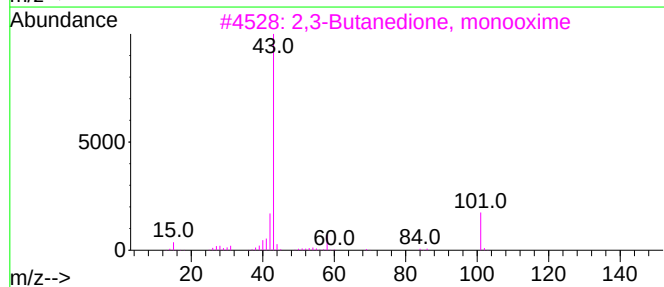
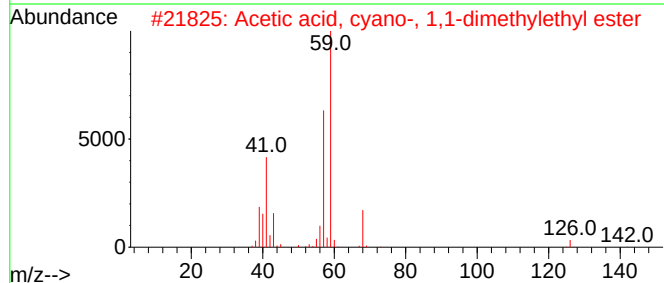
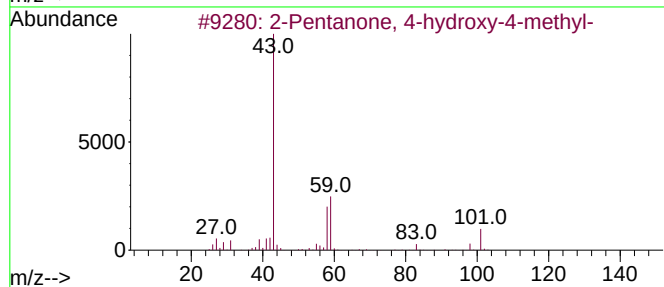
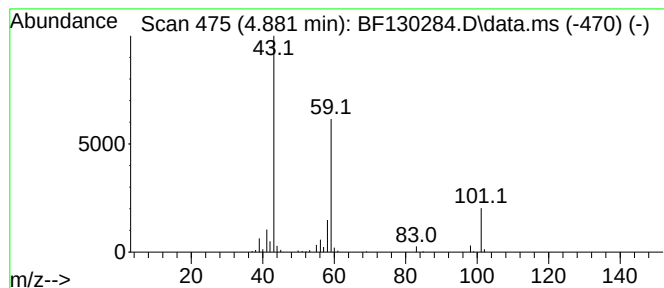
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.881	7.09 ng	214692	1,4-Dichlorobenzene-d4	6.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72		
2	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25		
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.881	7.1	ng	214692	1	6.675	605907	20.0