

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091423\
 Data File : BF135255.D
 Acq On : 14 Sep 2023 15:40
 Operator : CG\JU
 Sample : PB154922BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB154922BL

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-BF091123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135255.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.016	494	498	511	rBV	183882	295180	6.97%	1.093%
2	5.404	559	564	567	rBV	3339988	3774871	89.13%	13.979%
3	6.410	729	735	738	rBV	3152719	3556135	83.97%	13.169%
4	6.763	791	795	799	rBV	993679	801281	18.92%	2.967%
5	7.328	885	891	894	rBV	2043662	2298324	54.27%	8.511%
6	8.039	1007	1012	1016	rBV	1197582	1062006	25.08%	3.933%
7	9.122	1190	1196	1199	rBV	3838791	3947718	93.21%	14.619%
8	9.798	1305	1311	1315	rBV	1499722	1230334	29.05%	4.556%
9	10.598	1442	1447	1450	rBV	2569726	2665312	62.93%	9.870%
10	11.292	1561	1565	1568	rBV	1747743	1436301	33.91%	5.319%
11	12.892	1832	1837	1840	rBV	3783482	4235122	100.00%	15.683%
12	13.945	2012	2016	2026	rBV	997218	890683	21.03%	3.298%
13	15.415	2260	2266	2276	rBV	606987	811326	19.16%	3.004%

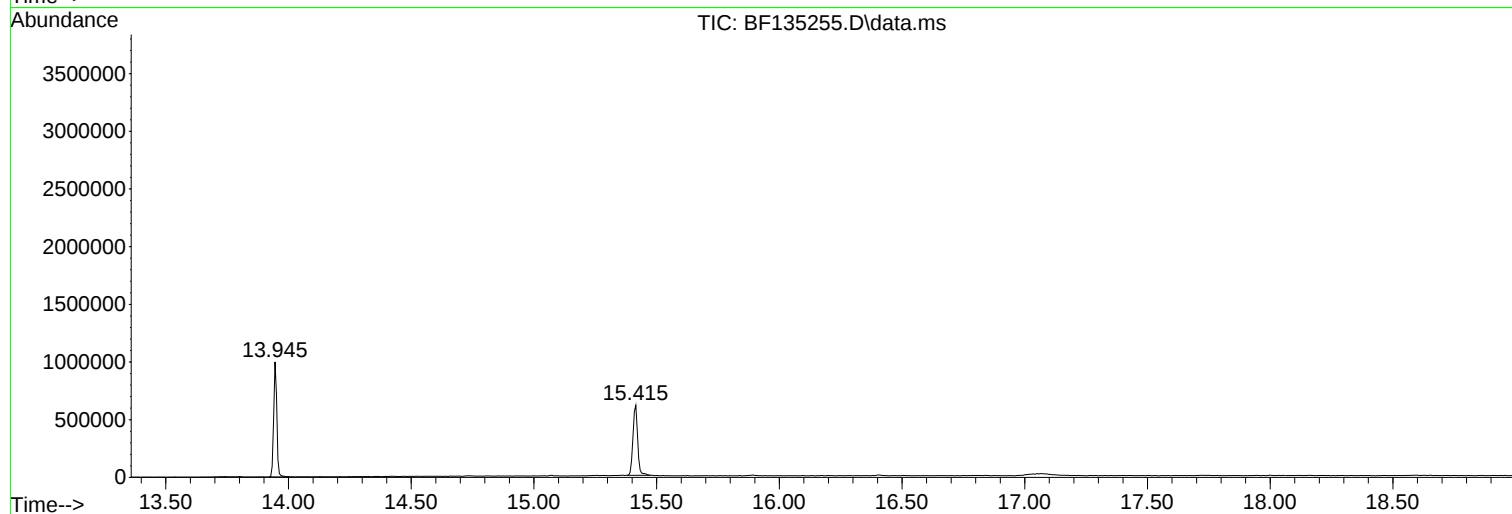
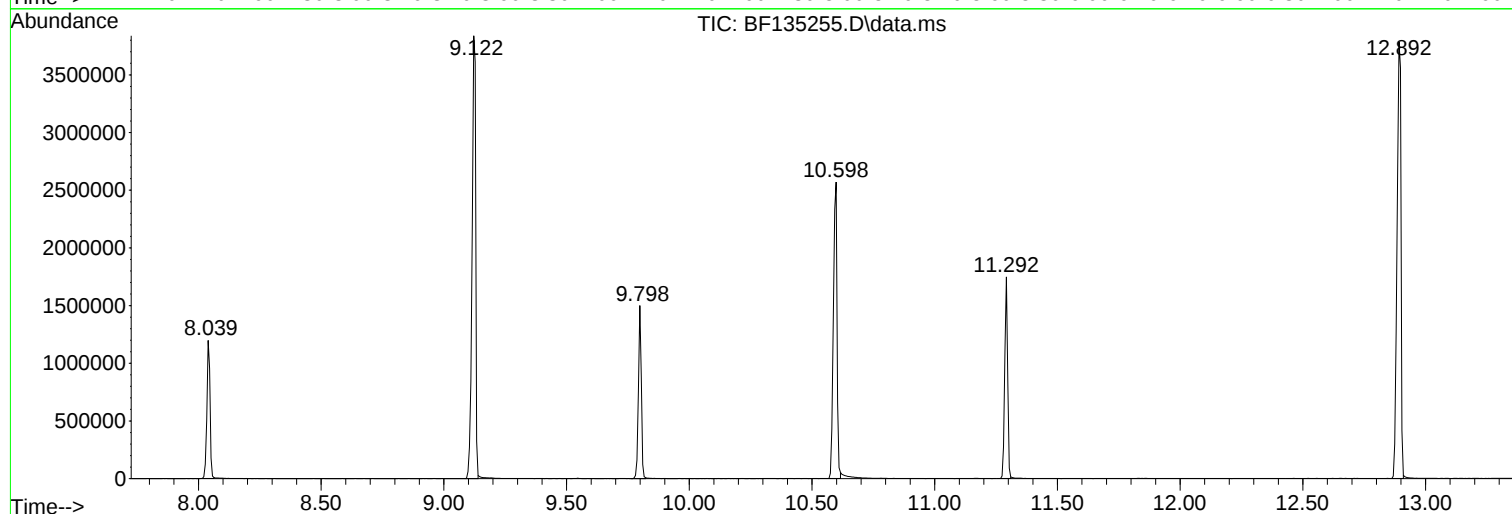
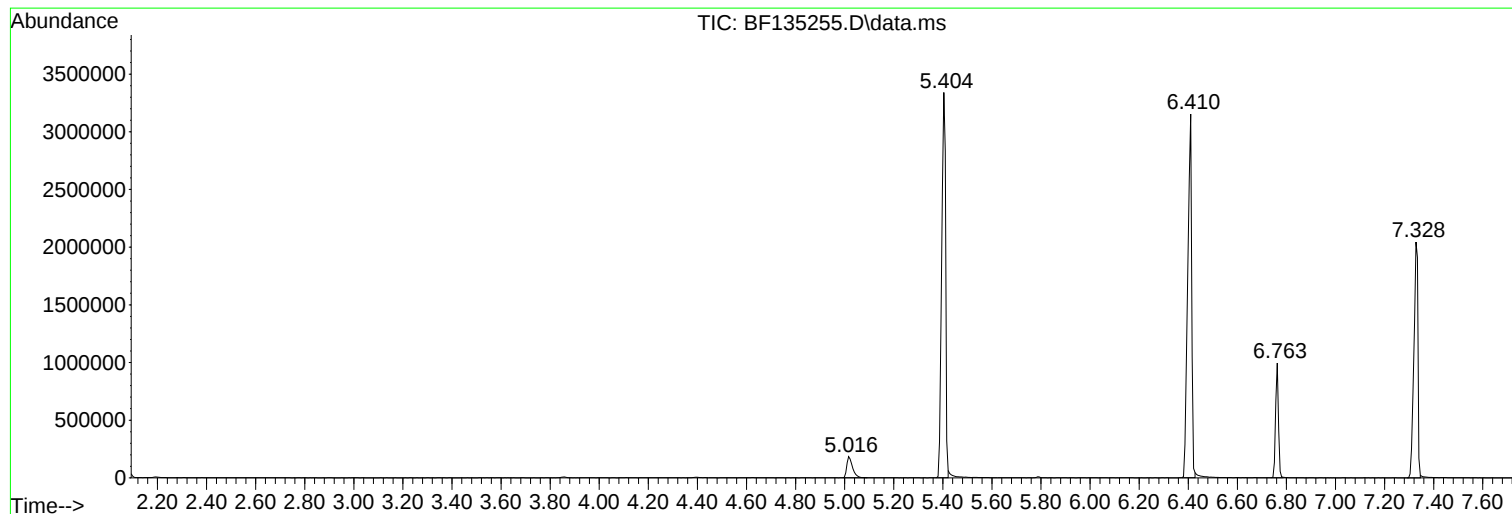
Sum of corrected areas: 27004593

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.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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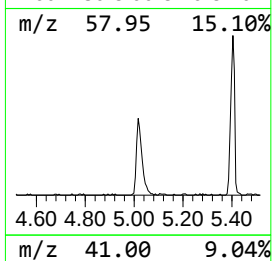
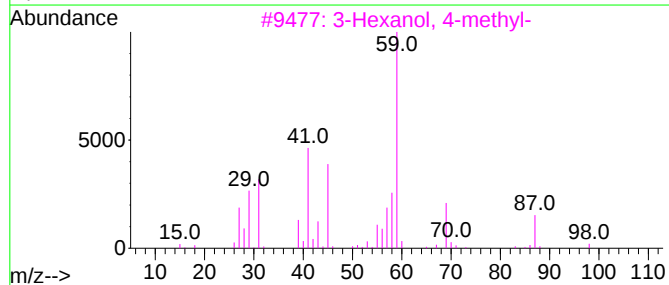
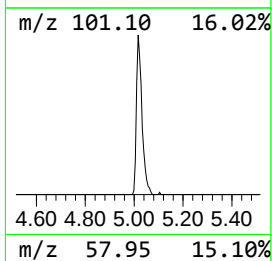
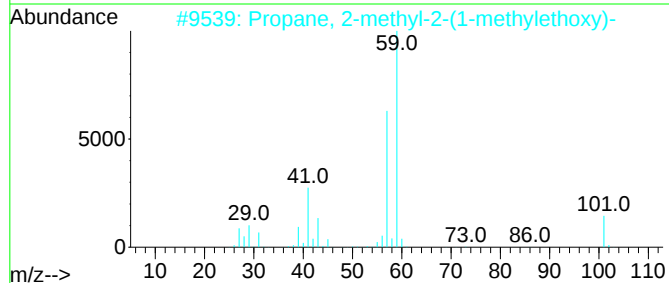
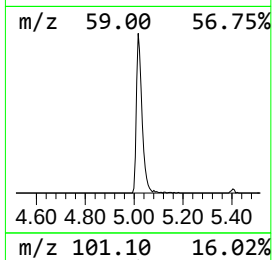
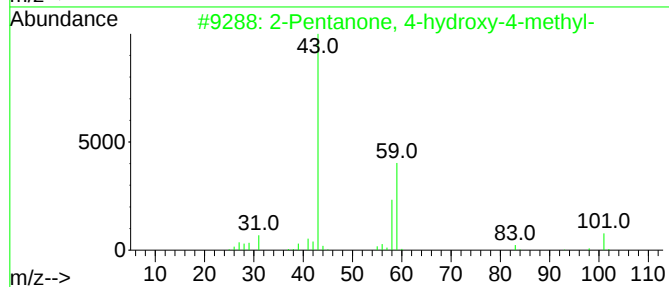
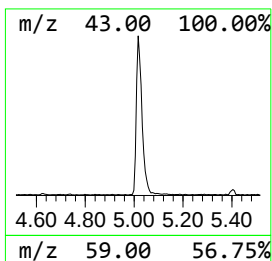
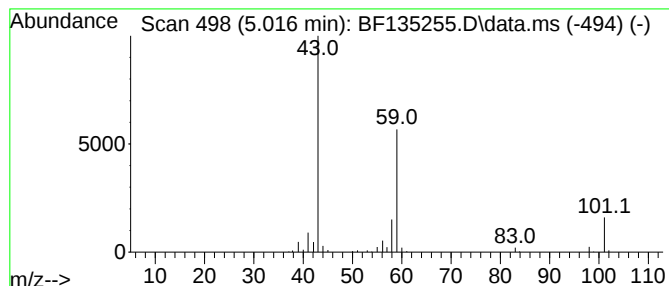
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.016	7.37 ng	295180	1,4-Dichlorobenzene-d4			6.763
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	50
2	Propane, 2-methyl-2-(1-methyleth...		116	C7H16O	017348-59-3	36
3	3-Hexanol, 4-methyl-		116	C7H16O	000615-29-2	33
4	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	28
5	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	23



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	5.016	7.4	ng	295180	1	6.763	801281	20.0