

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101023\
Data File : BF135678.D
Acq On : 10 Oct 2023 11:16
Operator : CG\JU
Sample : PB156189BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB156189BL

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: BF135678.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.981	488	492	503	rBV	314427	474906	13.37%	2.116%
2	5.381	557	560	582	rBV	2835789	3091476	87.00%	13.775%
3	6.393	728	732	748	rBV	2905182	3147866	88.59%	14.027%
4	6.740	787	791	798	rBV	662439	600114	16.89%	2.674%
5	7.304	883	887	899	rBV	1910193	2042941	57.49%	9.103%
6	8.016	1004	1008	1015	rBV	820217	750404	21.12%	3.344%
7	9.098	1186	1192	1195	rBV	3335735	3395416	95.56%	15.130%
8	9.775	1301	1307	1315	rBV	949037	847496	23.85%	3.776%
9	10.575	1438	1443	1456	rBV	2188342	2247118	63.24%	10.013%
10	11.269	1557	1561	1570	rBV	1080903	960985	27.05%	4.282%
11	12.869	1828	1833	1837	rBV	3379287	3553265	100.00%	15.833%
12	13.933	2010	2014	2022	rBV	754439	730827	20.57%	3.256%
13	15.404	2259	2264	2274	rBV	421215	599301	16.87%	2.670%

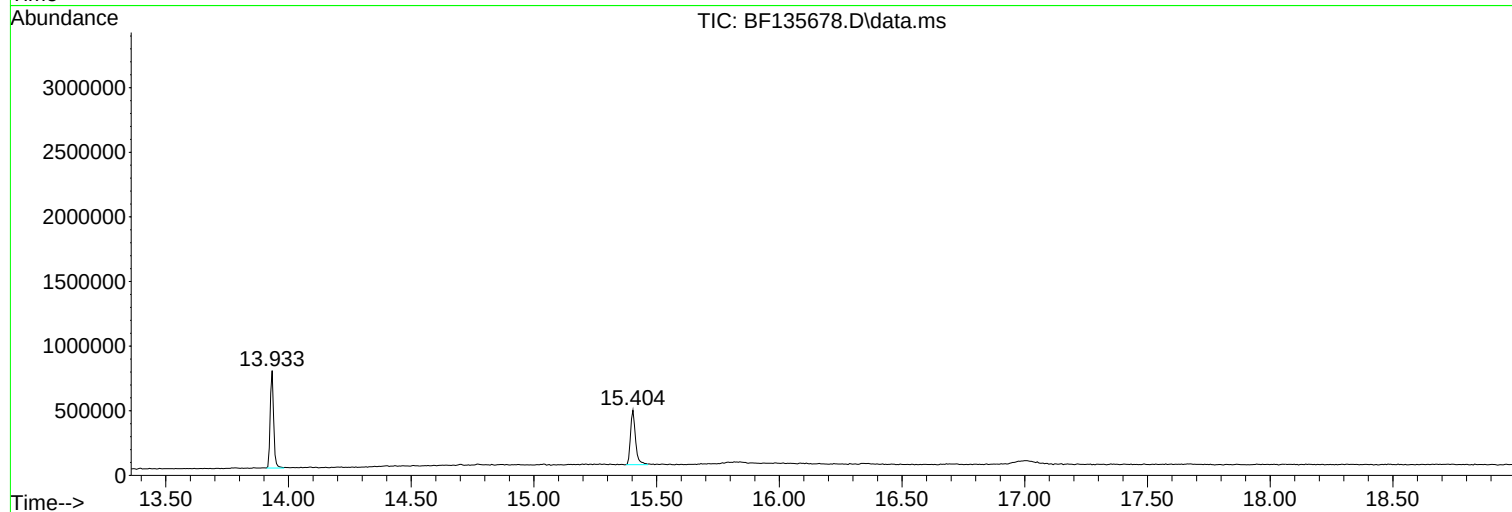
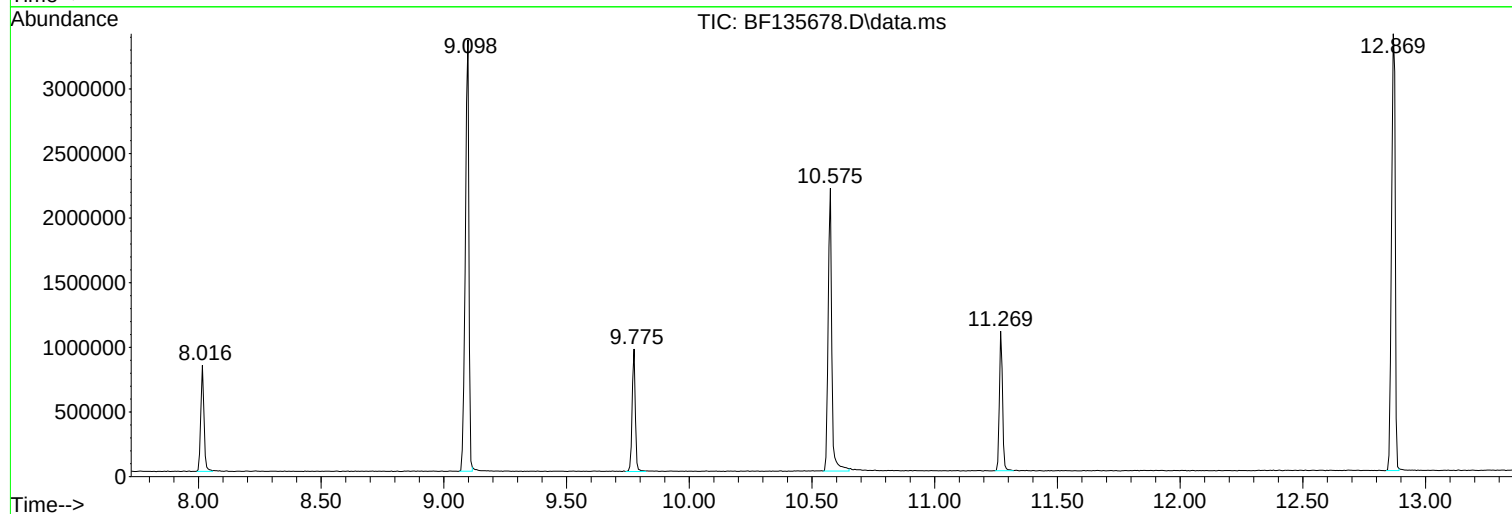
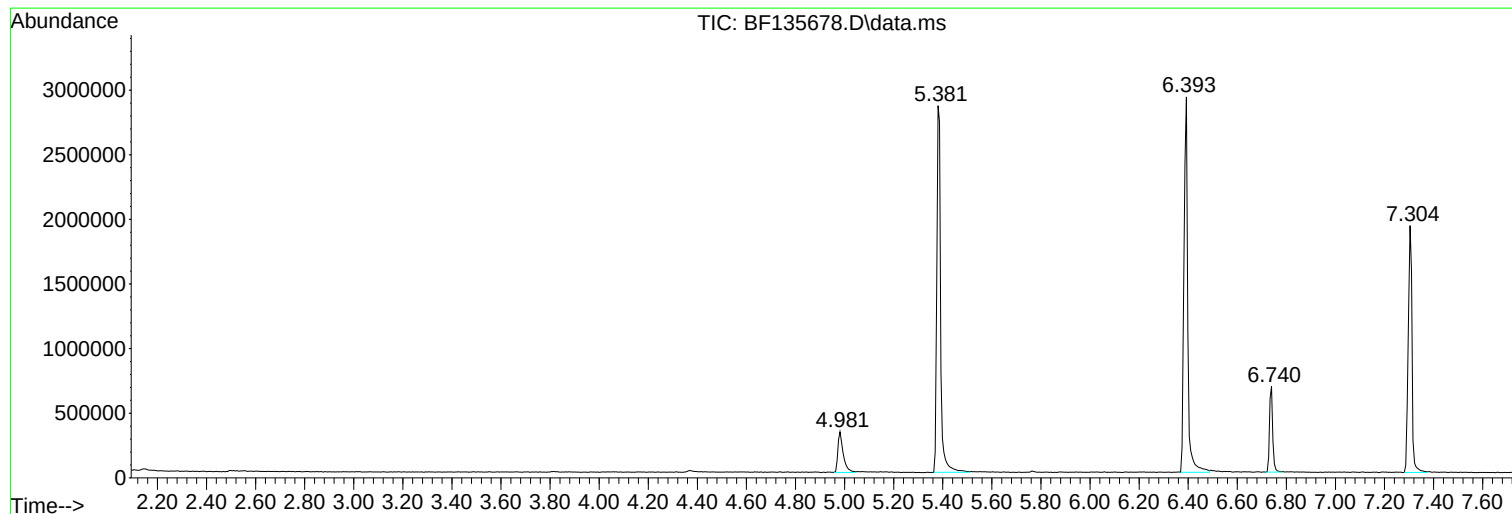
Sum of corrected areas: 22442115

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.981	15.83 ng	474906	1,4-Dichlorobenzene-d4	6.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			3-Hexanol	102	C6H14O	000623-37-0	33
5			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.981	15.8	ng	474906	1	6.740	600114	20.0