

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100423\
 Data File : BF135583.D
 Acq On : 04 Oct 2023 22:19
 Operator : CG\JU
 Sample : 04687-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-1-092923

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: BF135583.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.657	93	97	106	rBV	193966	282965	6.28%	0.968%
2	4.546	414	418	427	rBV	69238	84590	1.88%	0.289%
3	4.981	488	492	509	rBV2	524472	852164	18.91%	2.916%
4	5.381	556	560	592	rBV	3629756	3901972	86.56%	13.353%
5	6.387	727	731	759	rBV	3269792	3694128	81.95%	12.642%
6	6.740	787	791	800	rBV	940051	865552	19.20%	2.962%
7	7.304	883	887	890	rBV	2756801	2624641	58.23%	8.982%
8	8.016	1004	1008	1026	rVB	1248050	1141416	25.32%	3.906%
9	9.092	1186	1191	1195	rBV	4062354	4319499	95.83%	14.782%
10	9.769	1302	1306	1317	rBV	1438608	1271640	28.21%	4.352%
11	10.569	1438	1442	1445	rBV	2867498	2811485	62.37%	9.621%
12	11.263	1557	1560	1569	rBV	1563626	1348083	29.91%	4.613%
13	12.863	1827	1832	1835	rBV	4225796	4507593	100.00%	15.426%
14	13.921	2008	2012	2023	rVB	859820	782314	17.36%	2.677%
15	15.386	2256	2261	2274	rBV	539927	732911	16.26%	2.508%

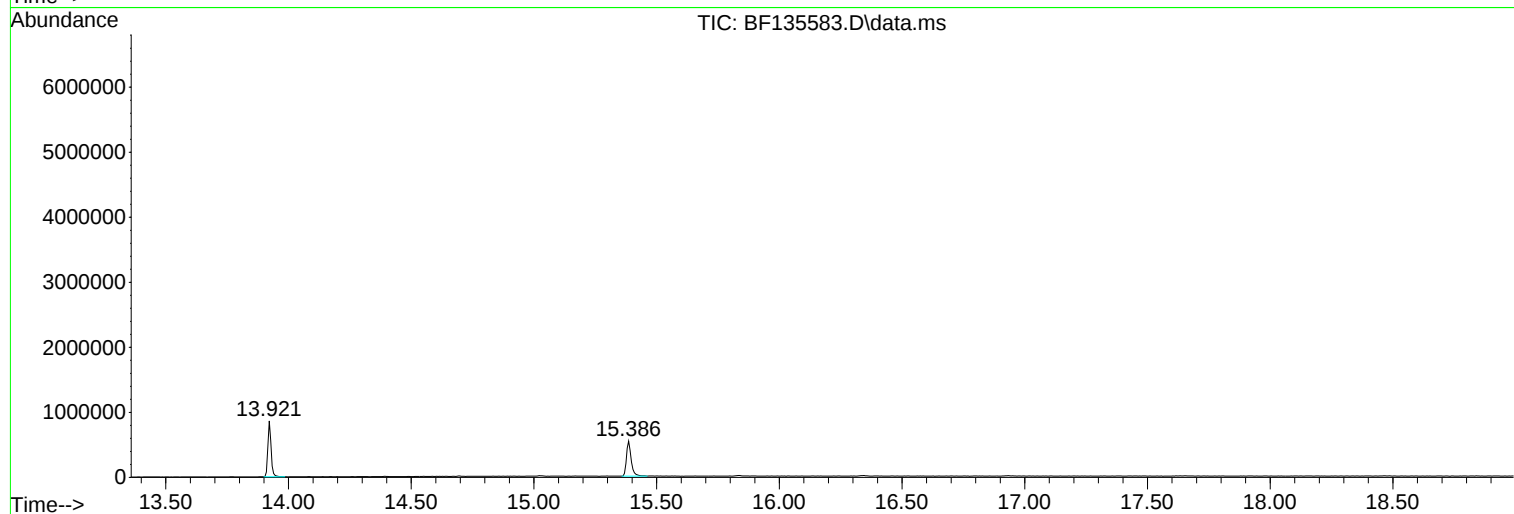
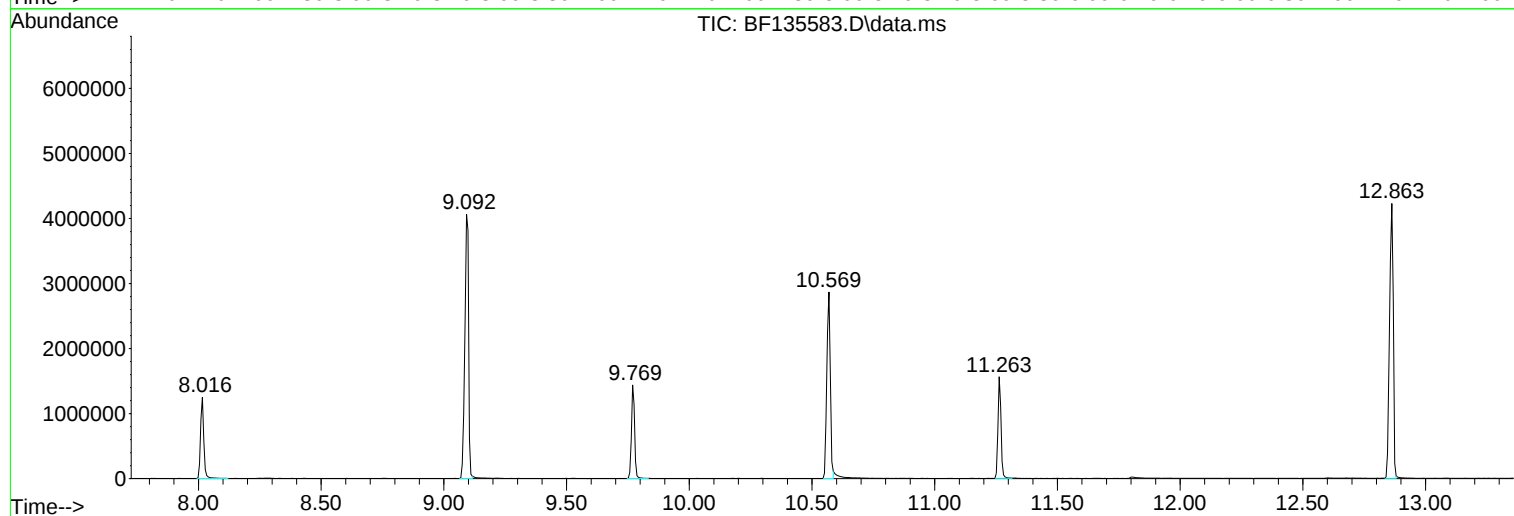
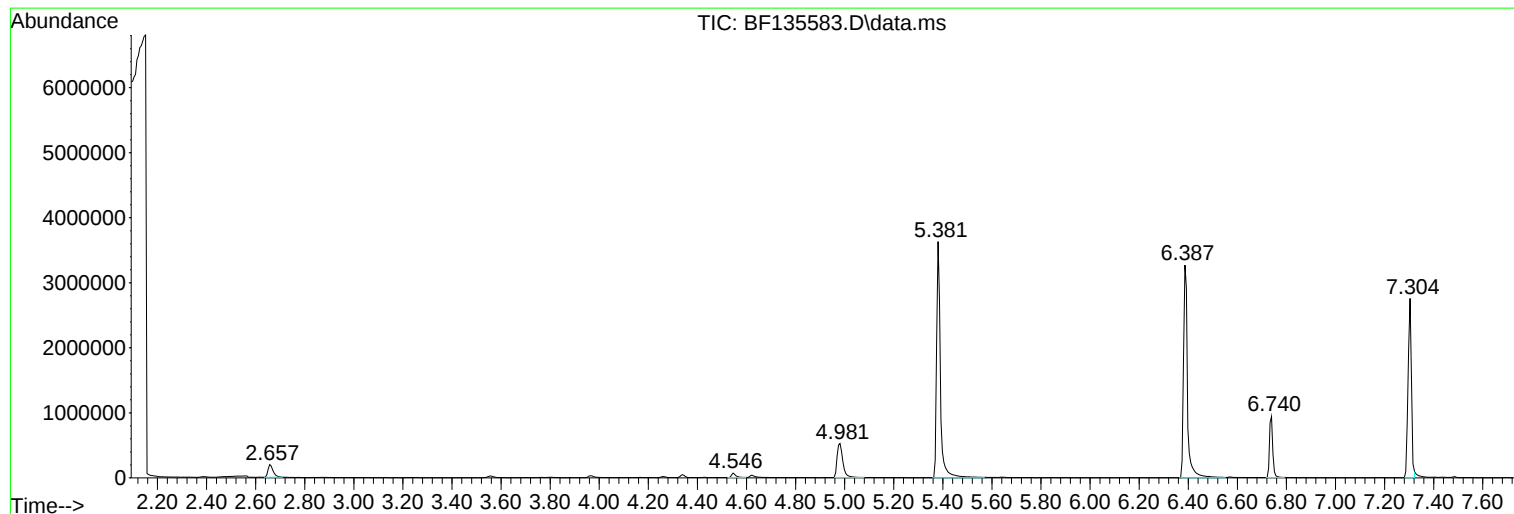
Sum of corrected areas: 29220953

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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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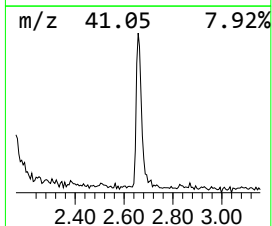
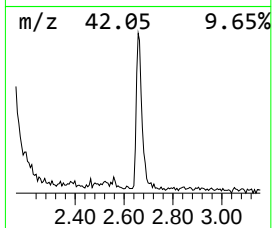
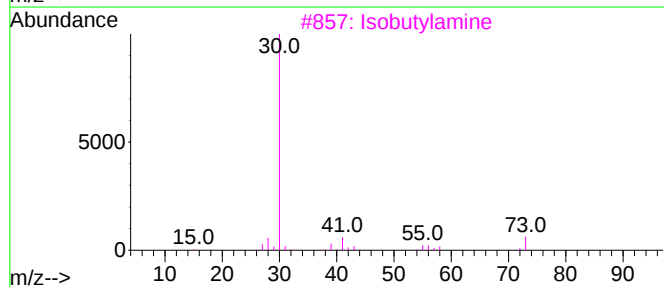
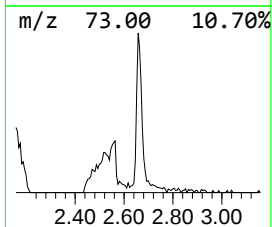
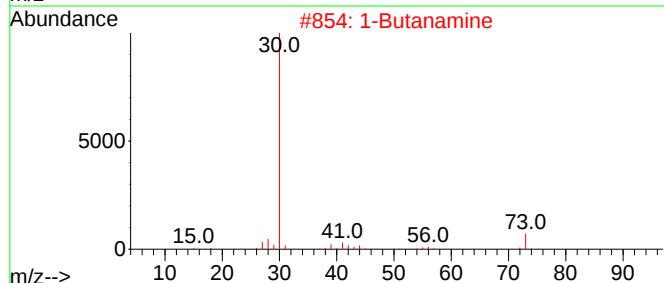
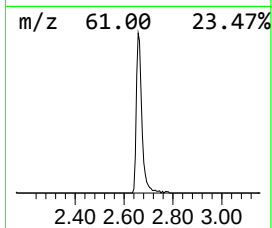
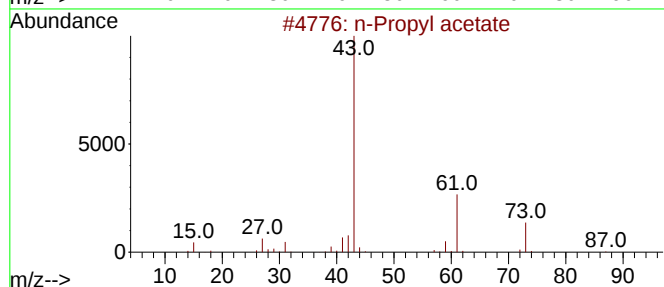
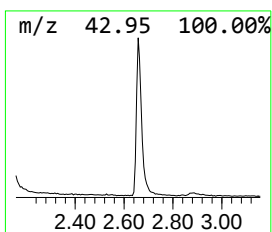
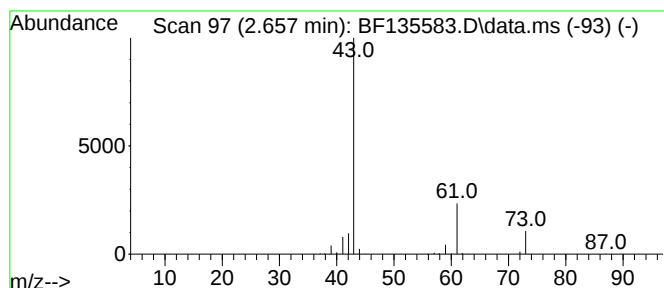
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 1 n-Propyl acetate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.657	6.54 ng	282965	1,4-Dichlorobenzene-d4	6.739

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Propyl acetate	102	C5H10O2	000109-60-4	83
2		1-Butanamine	73	C4H11N	000109-73-9	7
3		Isobutylamine	73	C4H11N	000078-81-9	5
4		Hydrogen azide	43	HN3	007782-79-8	4
5		1,4-Butanediol, diacetate	174	C8H14O4	000628-67-1	4



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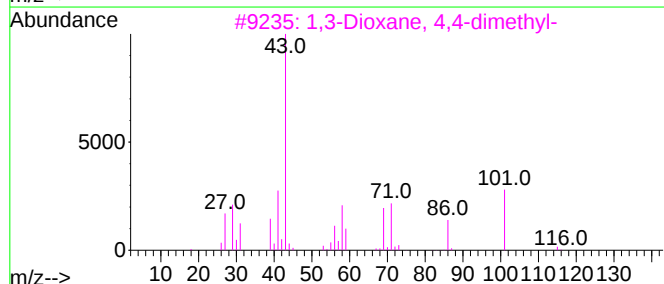
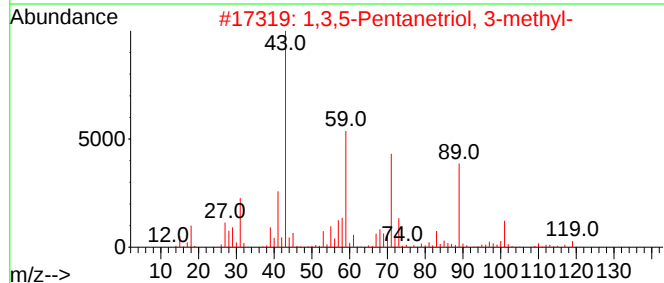
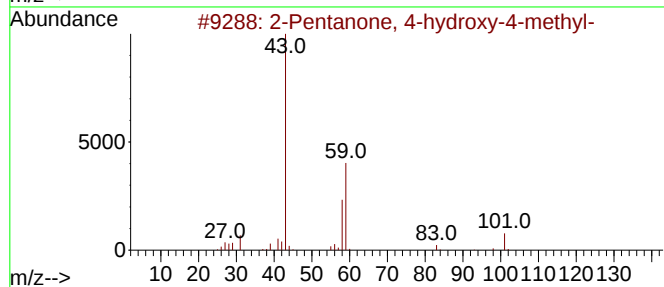
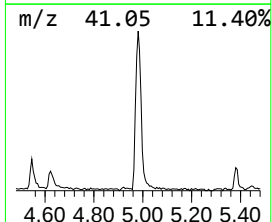
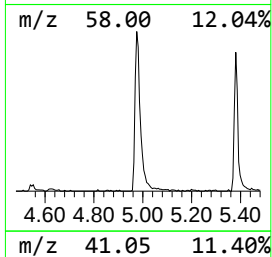
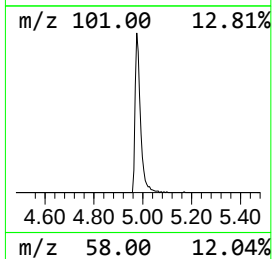
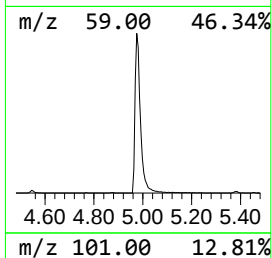
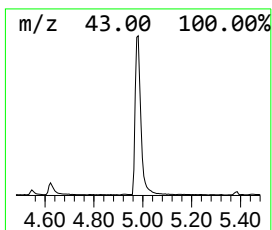
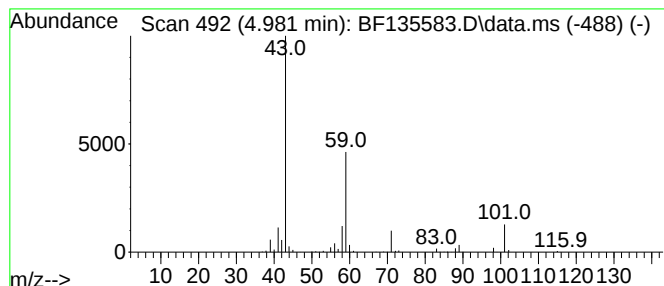
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.981	19.69 ng	852164	1,4-Dichlorobenzene-d4	6.739

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			1,3,5-Pentanetriol, 3-methyl-	134	C6H14O3	007564-64-9	38
3			1,3-Dioxane, 4,4-dimethyl-	116	C6H12O2	000766-15-4	27
4			Diacetamide	101	C4H7NO2	000625-77-4	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
n-Propyl acetate	2.657	6.5	ng	282965	1	6.739	865552	20.0
2-Pentanone, 4-...	4.981	19.7	ng	852164	1	6.739	865552	20.0