

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011623\
 Data File : BF132082.D
 Acq On : 16 Jan 2023 22:07
 Operator : CG\JU
 Sample : PB150088BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB150088BL

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-BF011623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF132082.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.222	13	23	28	rBV	814440	971645	11.52%	1.857%
2	2.593	75	86	92	rVB	123005	182601	2.17%	0.349%
3	5.363	550	557	567	rVB	1630602	2173710	25.78%	4.154%
4	5.734	614	620	623	rBV	5073512	5617960	66.62%	10.736%
5	6.710	779	786	789	rBV	4656650	5782553	68.57%	11.051%
6	7.098	848	852	855	rBV	1573297	1438316	17.06%	2.749%
7	7.657	941	947	950	rBV	3558837	3753336	44.51%	7.173%
8	8.381	1065	1070	1074	rBV	1992672	1939430	23.00%	3.706%
9	9.463	1248	1254	1257	rBV	6921940	7403473	87.80%	14.149%
10	10.151	1365	1371	1374	rBV	2559225	2296026	27.23%	4.388%
11	10.945	1499	1506	1509	rBV	5187313	5017344	59.50%	9.589%
12	11.651	1620	1626	1629	rBV	2807853	2484889	29.47%	4.749%
13	13.251	1891	1898	1901	rBV	7981365	8432633	100.00%	16.116%
14	14.310	2074	2078	2082	rBV	2779247	2459096	29.16%	4.700%
15	15.915	2344	2351	2362	rBV	1688146	2372924	28.14%	4.535%

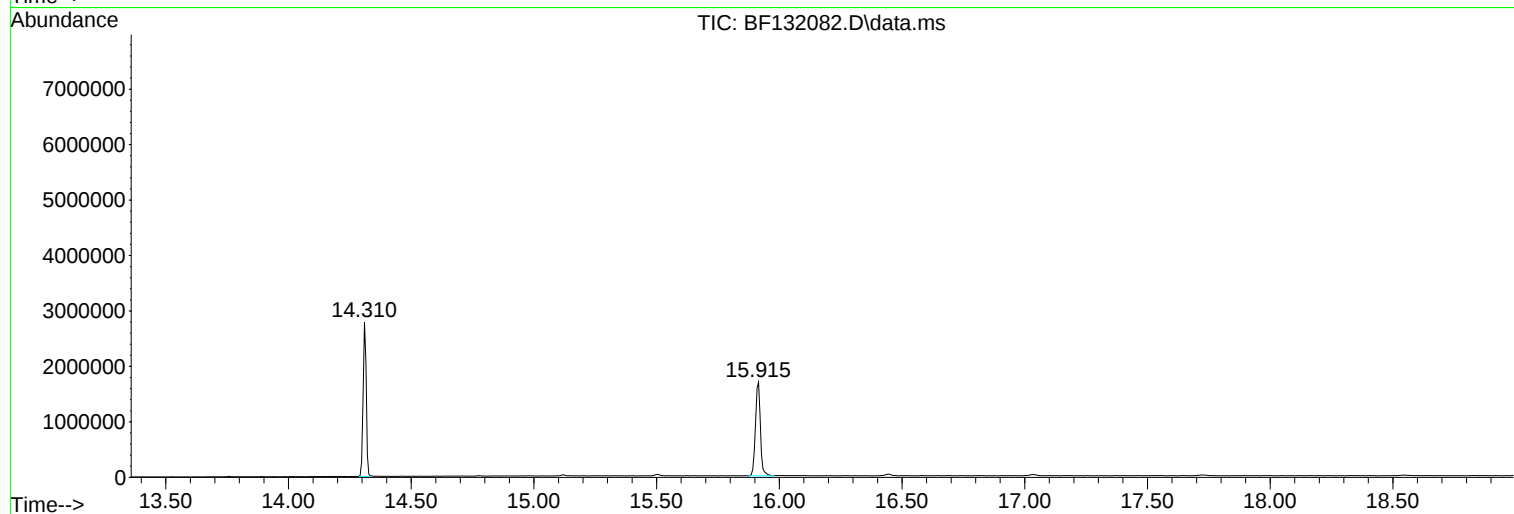
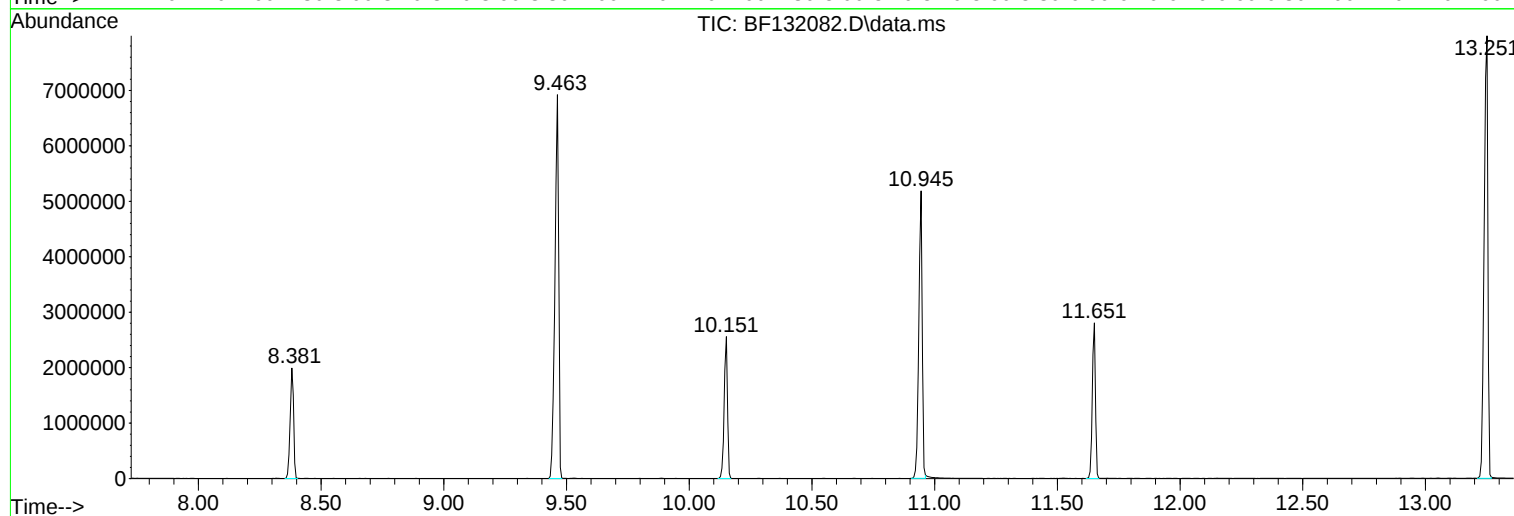
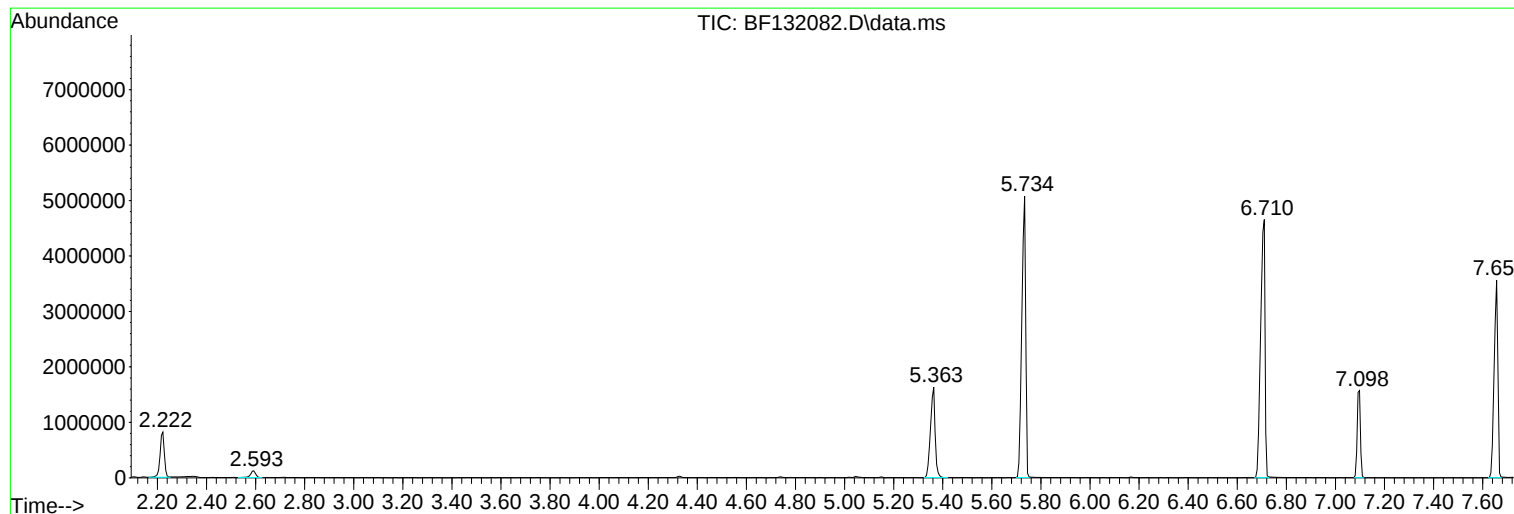
Sum of corrected areas: 52325936

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011623.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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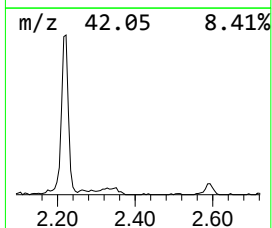
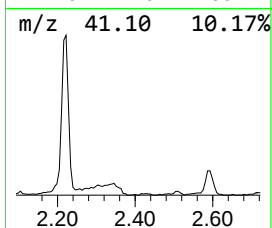
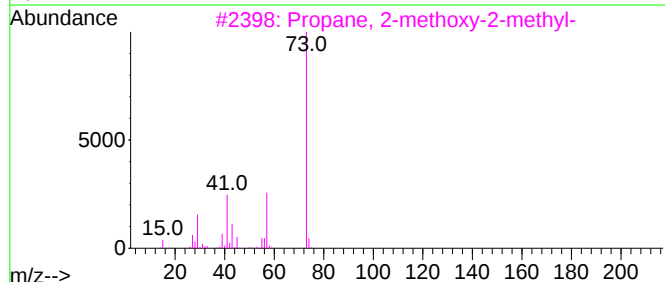
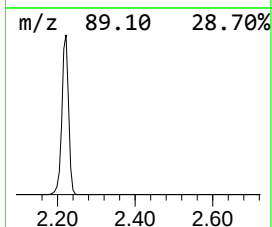
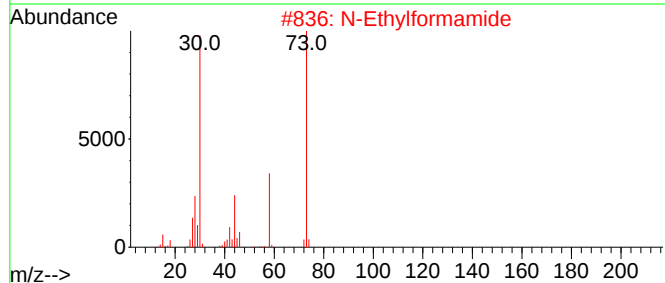
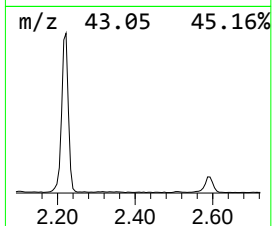
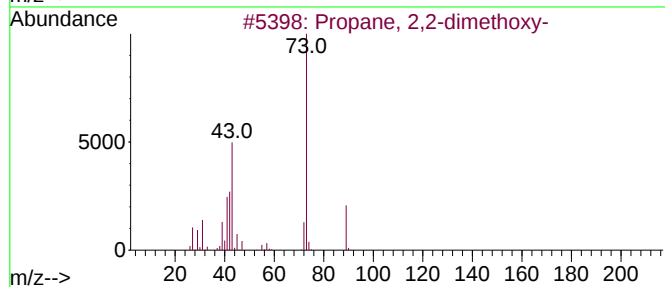
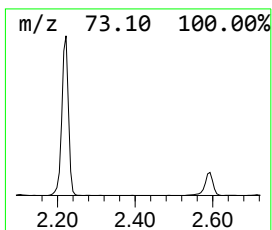
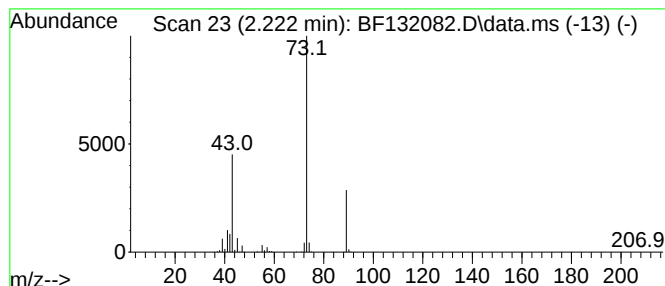
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.222	13.51 ng	971645	1,4-Dichlorobenzene-d4	7.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	50
2			N-Ethylformamide	73	C3H7NO	000627-45-2	37
3			Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4			2-Acetamido-N-methylacetamide	130	C5H10N2O2	007606-79-3	9
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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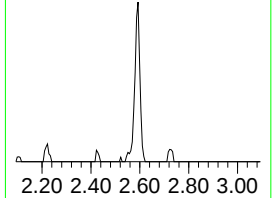
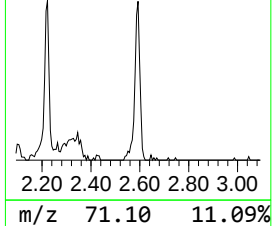
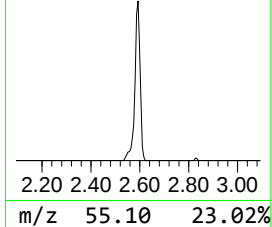
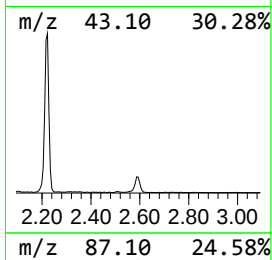
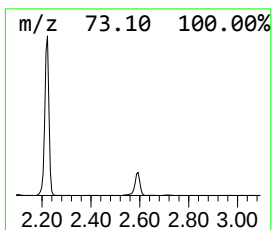
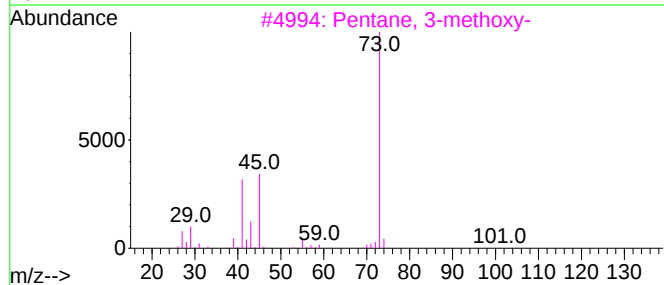
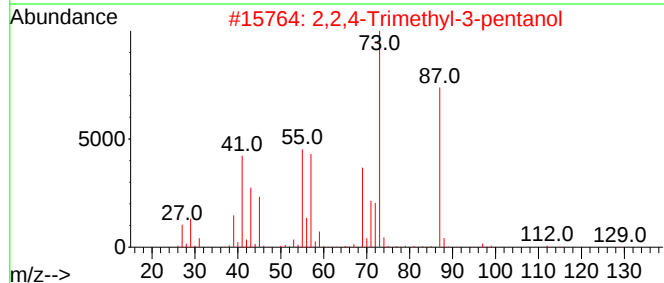
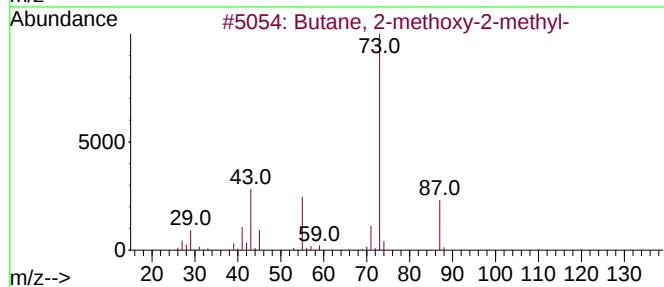
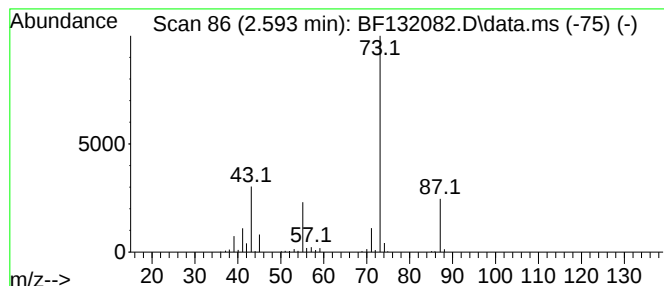
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TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.593	2.54 ng	182601	1,4-Dichlorobenzene-d4	7.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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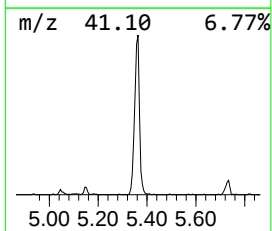
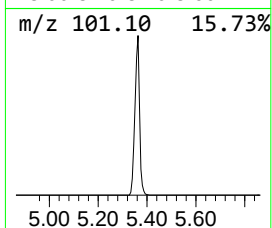
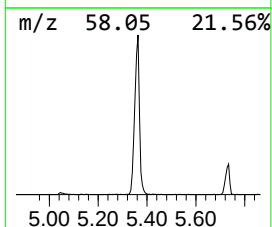
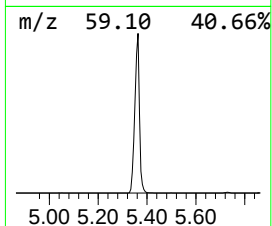
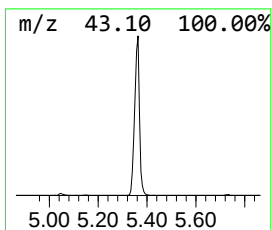
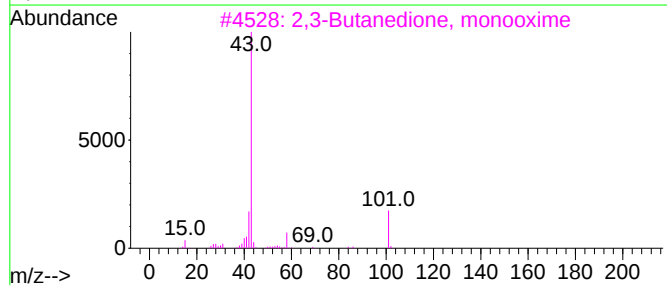
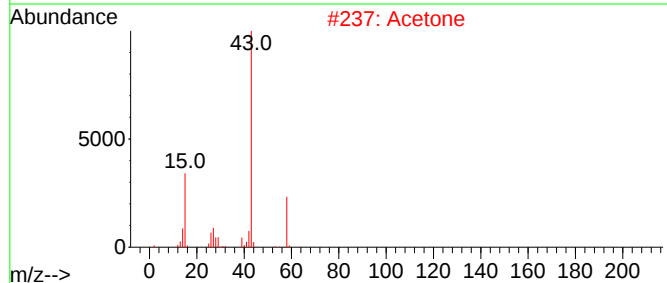
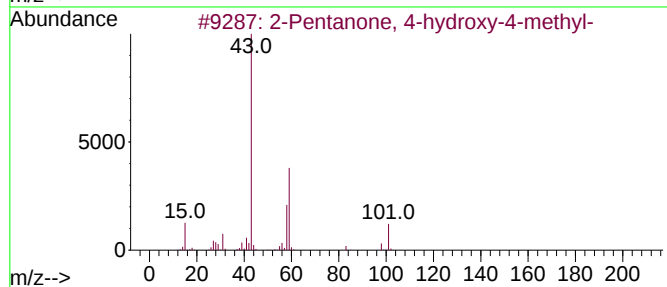
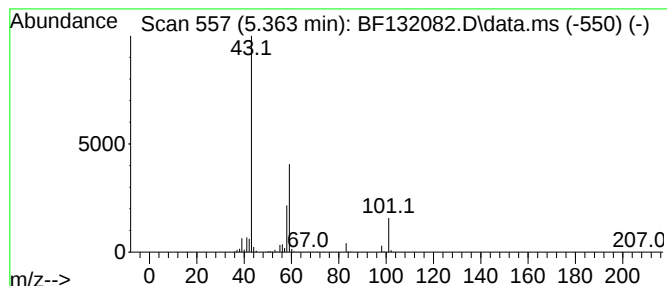
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.363	30.23 ng	2173710	1,4-Dichlorobenzene-d4	7.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2			Acetone	58	C3H6O	000067-64-1	16
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2,2-di...	2.222	13.5	ng	971645	1	7.098	1438320	20.0
Butane, 2-metho...	2.593	2.5	ng	182601	1	7.098	1438320	20.0
2-Pentanone, 4-...	5.363	30.2	ng	2173710	1	7.098	1438320	20.0