

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
 Data File : BG019127.D
 Acq On : 10 Oct 2015 21:15
 Operator : UM/NP
 Sample : G3929-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-03

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:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.144	47	52	60	rBV	47911	82328	4.29%	1.030%
2	3.215	60	64	71	rBV	190160	252135	13.15%	3.155%
3	5.136	385	391	404	rBV	51228	78124	4.07%	0.978%
4	5.606	464	471	481	rVB	141988	225745	11.77%	2.825%
5	7.192	734	741	748	rBV	91073	151127	7.88%	1.891%
6	7.562	797	804	816	rBB	284940	490728	25.59%	6.140%
7	8.032	878	884	891	rVB	61836	103988	5.42%	1.301%
8	8.356	932	939	947	rVB	280191	481241	25.10%	6.022%
9	9.190	1074	1081	1091	rBV	263690	458540	23.91%	5.738%
10	10.835	1354	1361	1370	rBV	88801	157240	8.20%	1.968%
11	13.285	1770	1778	1785	rBV	728475	1123716	58.60%	14.061%
12	14.660	2006	2012	2019	rVB2	157622	245353	12.80%	3.070%
13	16.146	2259	2265	2280	rBV	659027	982022	51.21%	12.288%
14	16.869	2383	2388	2393	rBV	14591	19617	1.02%	0.245%
15	17.404	2473	2479	2487	rBV2	225688	332395	17.33%	4.159%
16	19.713	2868	2872	2878	rBV2	29812	32684	1.70%	0.409%
17	20.018	2918	2924	2930	rBV	1503573	1917556	100.00%	23.994%
18	20.753	3043	3049	3053	rBV2	23717	25534	1.33%	0.320%
19	21.675	3199	3206	3214	rBV	274643	428713	22.36%	5.364%
20	24.901	3745	3755	3770	rBV2	145676	403055	21.02%	5.043%

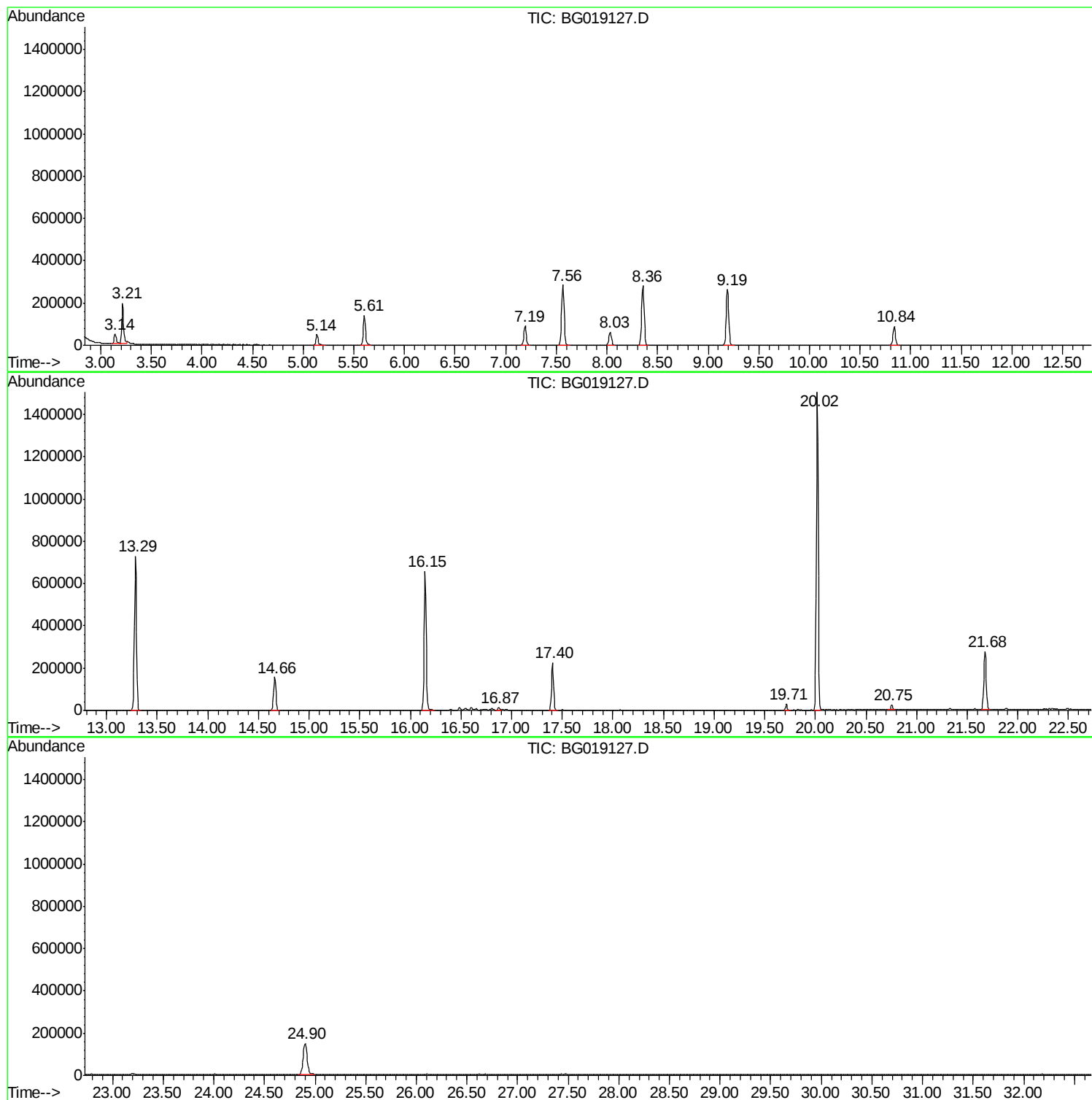
Sum of corrected areas: 7991841

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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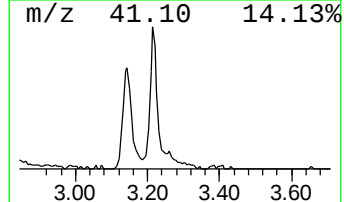
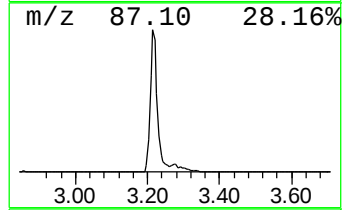
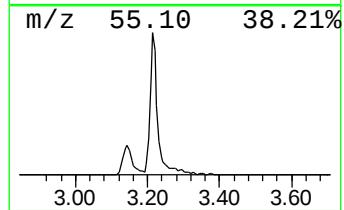
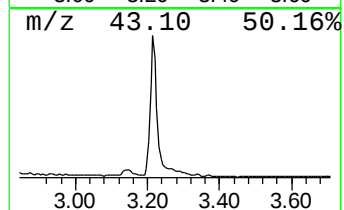
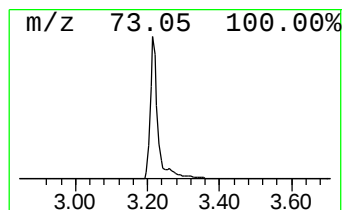
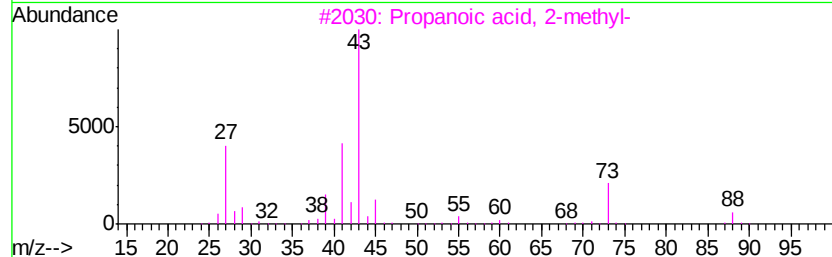
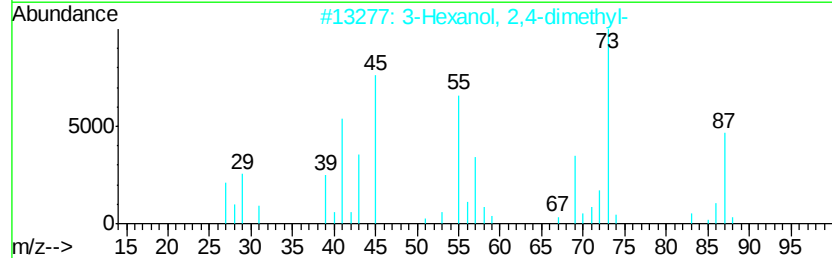
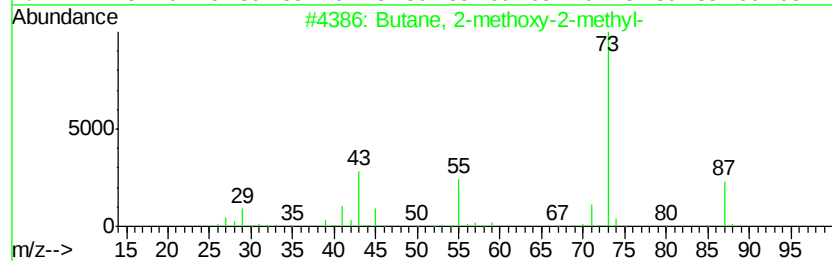
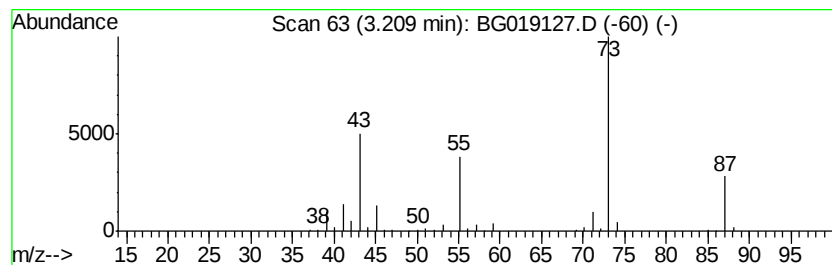
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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.
3.21	48.49 ng	252135	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	42
3		Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	10
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		3,4-Hexanediol, 2,5-dimethyl-	146	C8H18O2	022607-11-0	9



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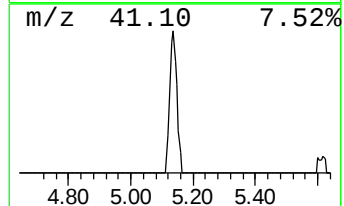
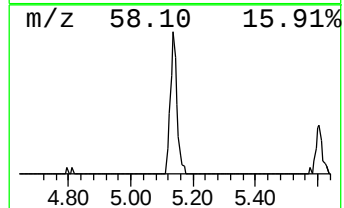
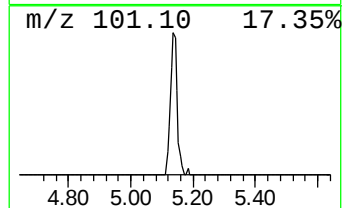
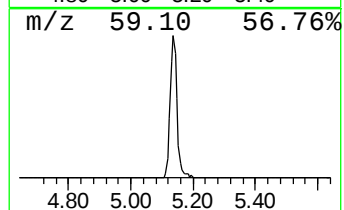
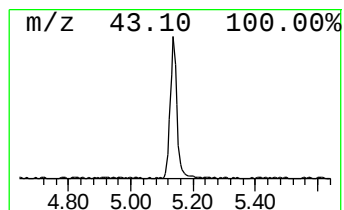
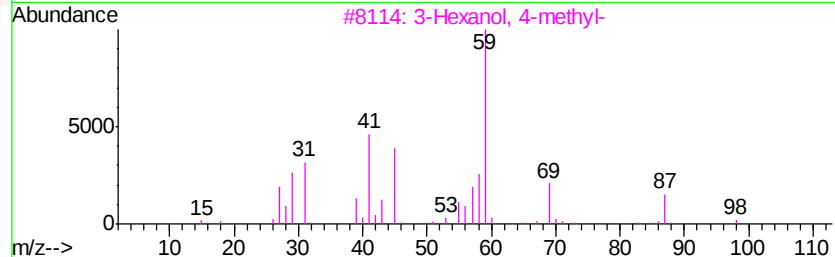
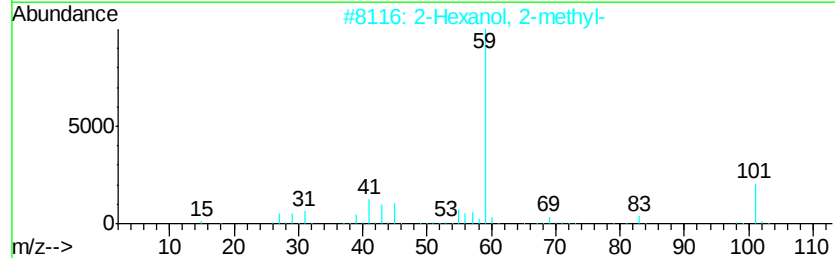
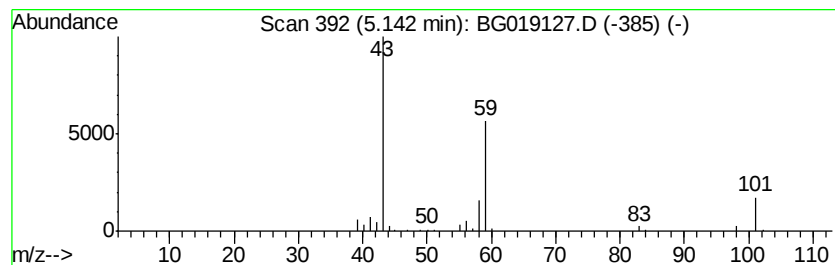
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	15.03 ng	78124	1,4-Dichlorobenzene-d4	8.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		3-Octanol	130	C8H18O	000589-98-0	33
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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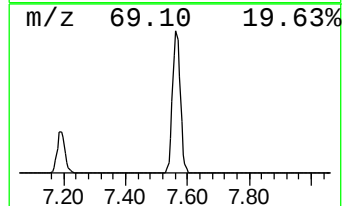
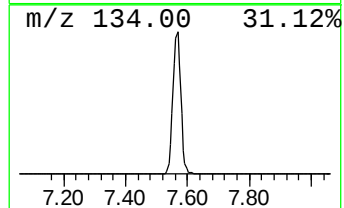
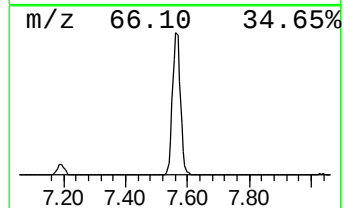
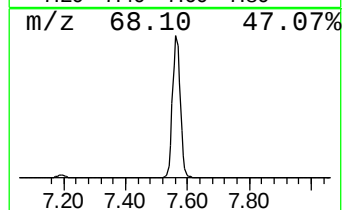
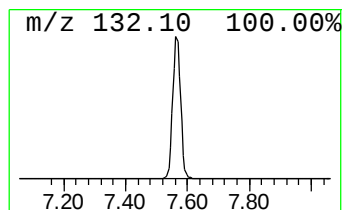
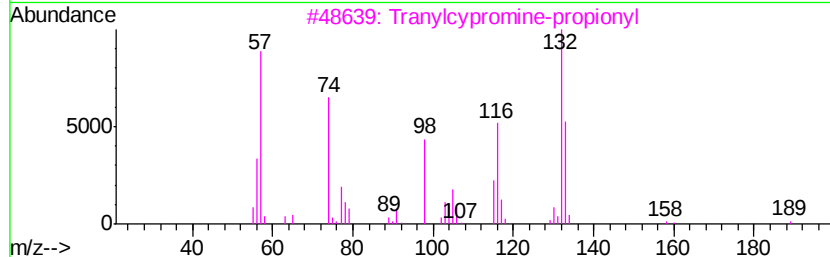
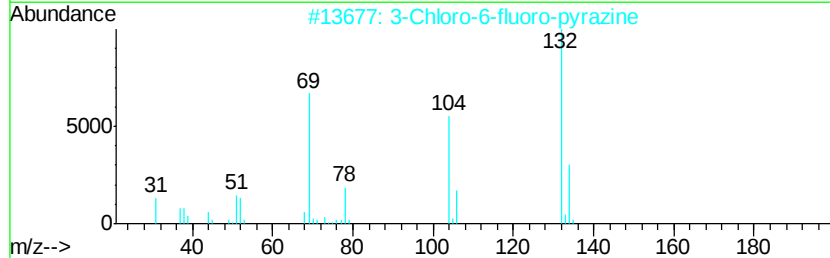
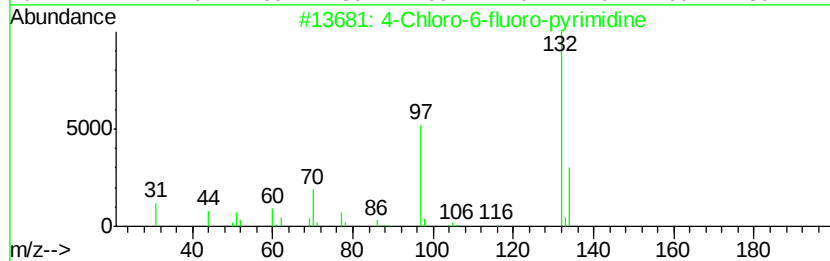
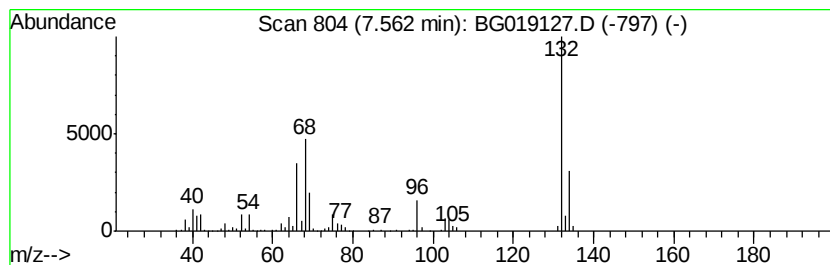
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Peak Number 4 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	94.38 ng	490728	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.21	48.5	ng	252135	1	8.03	103988	20.0
2-Pentanone, 4-hy...	5.14	15.0	ng	78124	1	8.03	103988	20.0
unknown7.56	7.56	94.4	ng	490728	1	8.03	103988	20.0