

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG101518\
 Data File : BG037333.D
 Acq On : 15 Oct 2018 14:48
 Operator : JU/SJ
 Sample : PB113787BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

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_G\METHODS\8270-BG101118.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

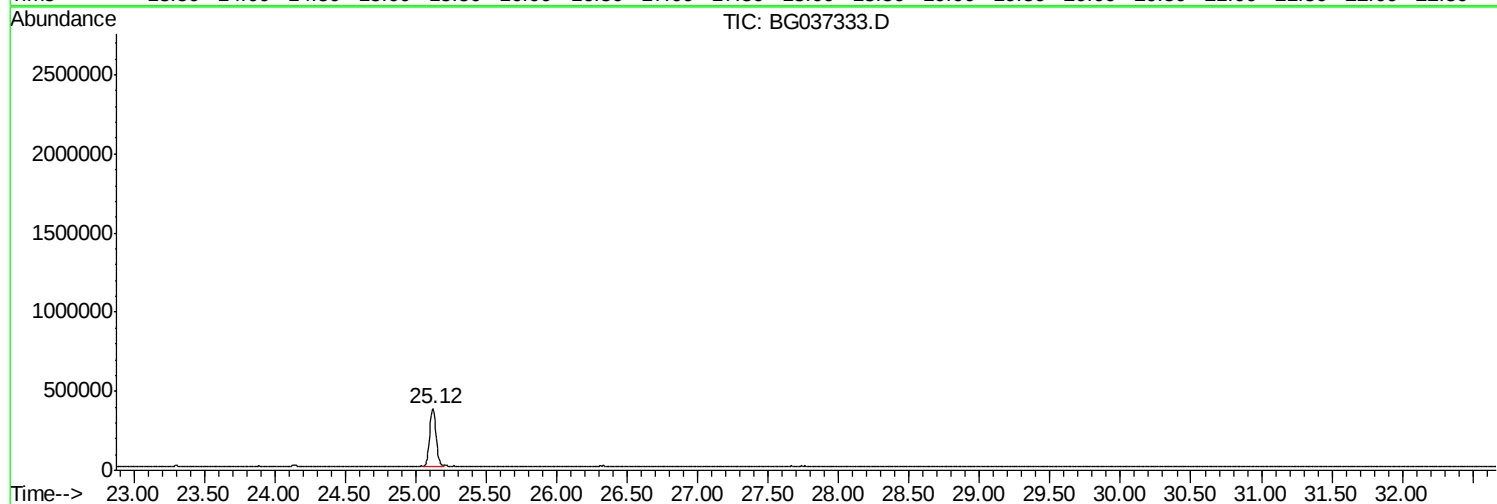
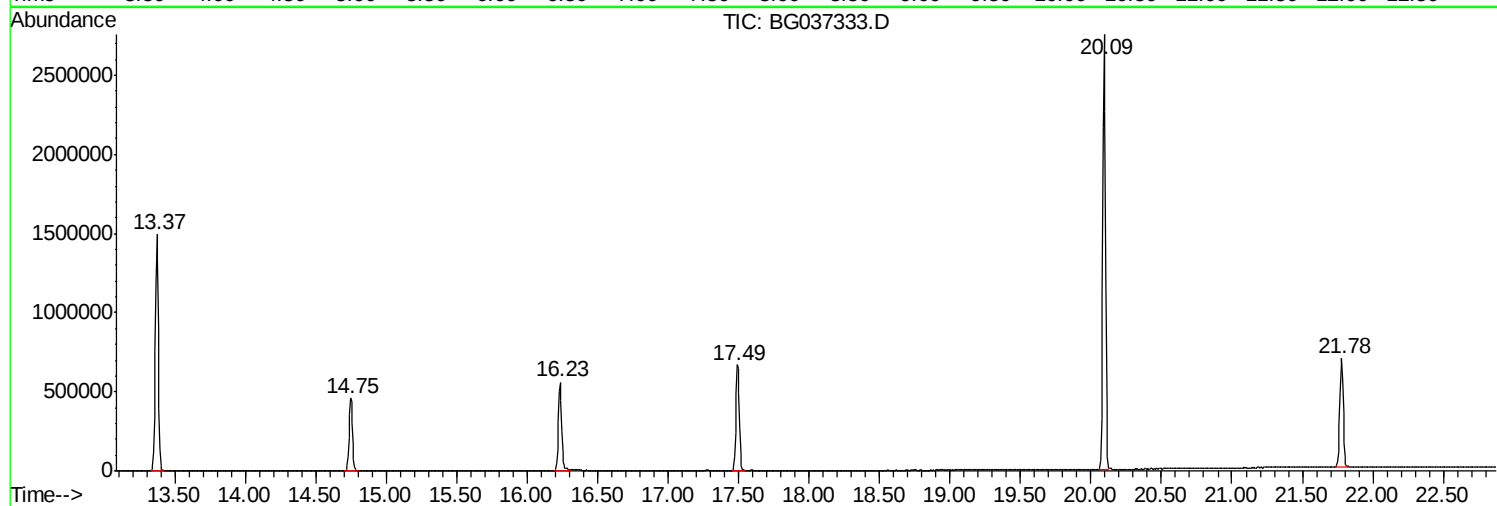
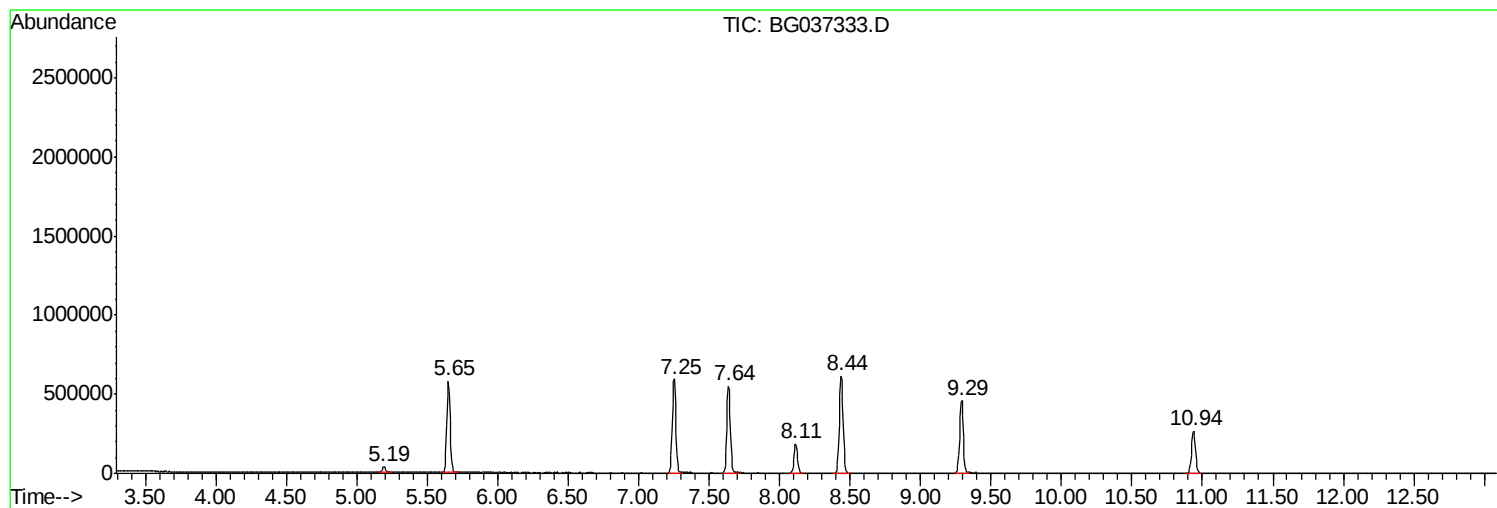
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.192	318	324	332	rVB	38621	61426	1.69%	0.365%
2	5.651	395	402	414	rBV	578962	948316	26.09%	5.634%
3	7.255	667	675	684	rBV	594346	1025996	28.22%	6.096%
4	7.636	733	740	753	rBV	551172	973394	26.78%	5.783%
5	8.112	814	821	829	rBV	183893	334065	9.19%	1.985%
6	8.436	867	876	886	rVB	615939	1113972	30.64%	6.619%
7	9.293	1013	1022	1030	rBV	454880	851133	23.41%	5.057%
8	10.938	1293	1302	1310	rBV	269034	505741	13.91%	3.005%
9	13.371	1709	1716	1725	rBV	1490903	2370615	65.21%	14.085%
10	14.746	1943	1950	1960	rBV2	458533	760950	20.93%	4.521%
11	16.232	2196	2203	2216	rBV	557532	892280	24.55%	5.301%
12	17.490	2411	2417	2427	rBV2	667337	1074115	29.55%	6.382%
13	20.092	2854	2860	2866	rBV	2749600	3635239	100.00%	21.599%
14	21.779	3139	3147	3154	rBV	686069	1181539	32.50%	7.020%
15	25.122	3705	3716	3730	rBV2	362916	1101932	30.31%	6.547%

Sum of corrected areas: 16830713

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG101118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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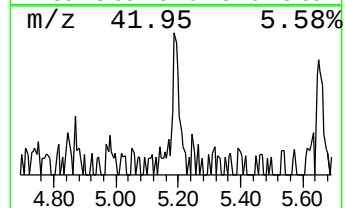
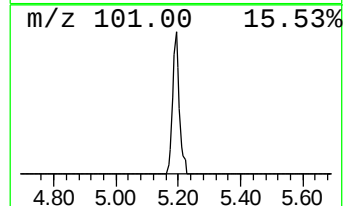
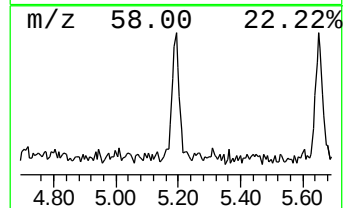
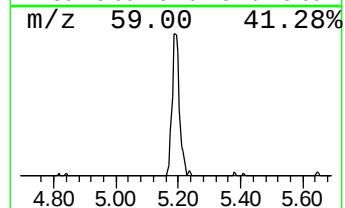
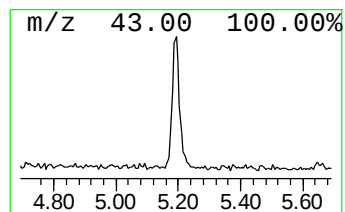
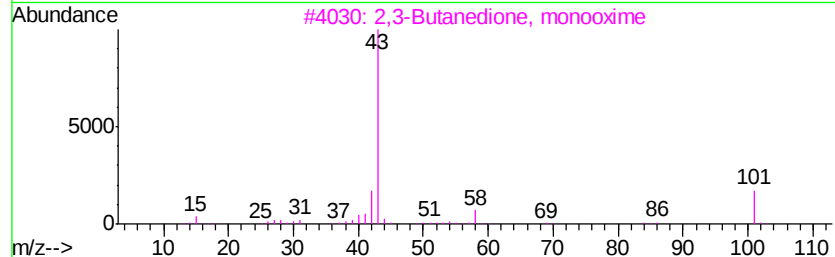
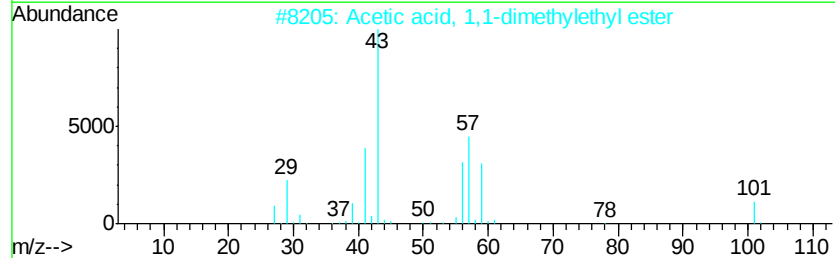
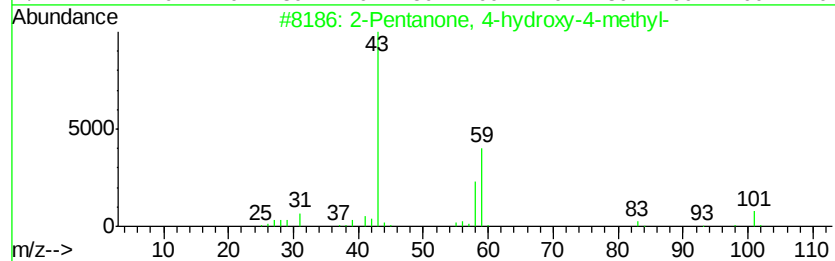
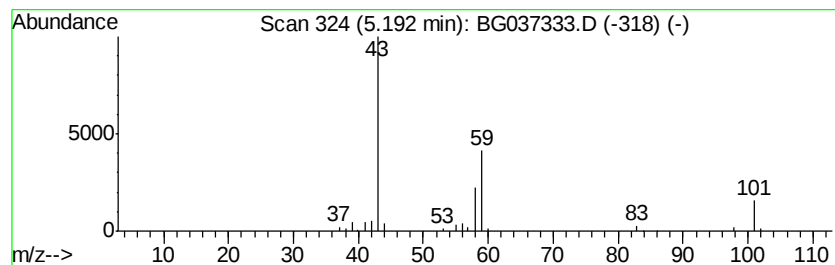
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	3.68 ng	61426	1,4-Dichlorobenzene-d4	8.11

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	78
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9
5			Acetone	58	C3H6O	000067-64-1	9



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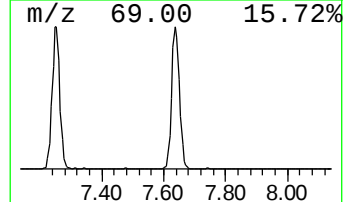
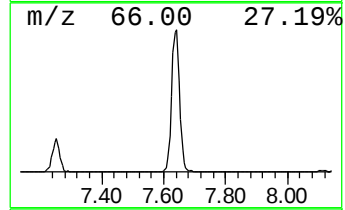
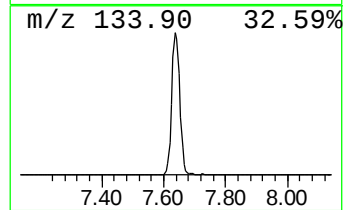
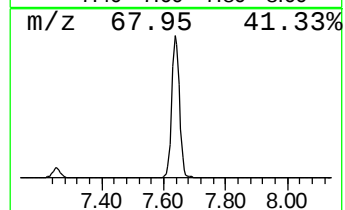
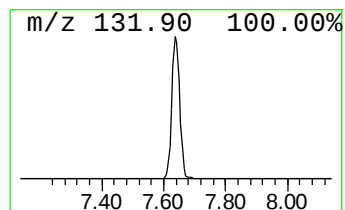
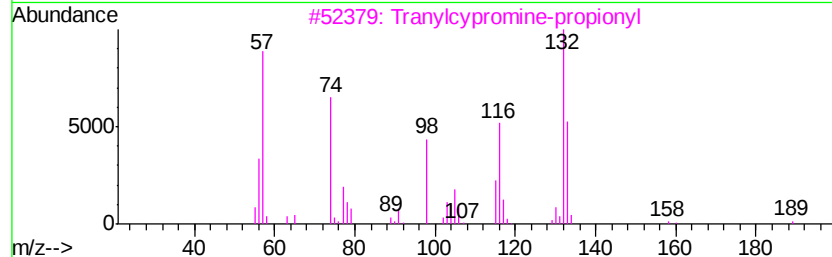
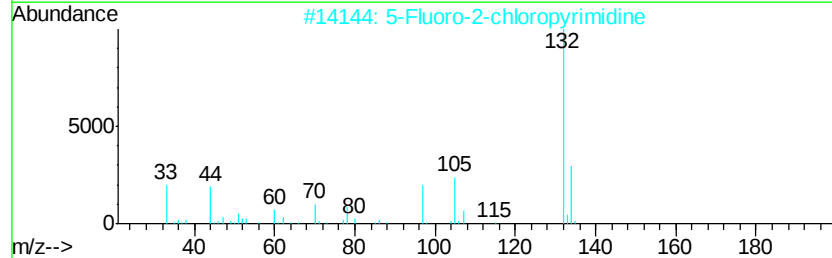
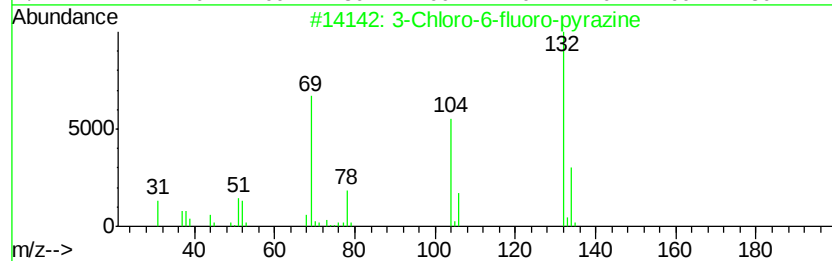
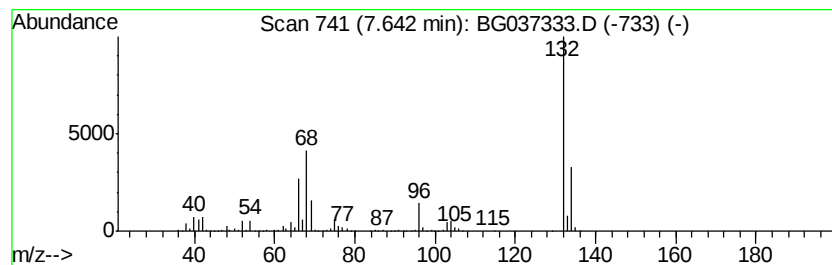
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Peak Number 2 unknown7.64 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	58.28 ng	973394	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	3.7	ng	61426	1	8.11	334065	20.0
unknown7.64	7.64	58.3	ng	973394	1	8.11	334065	20.0