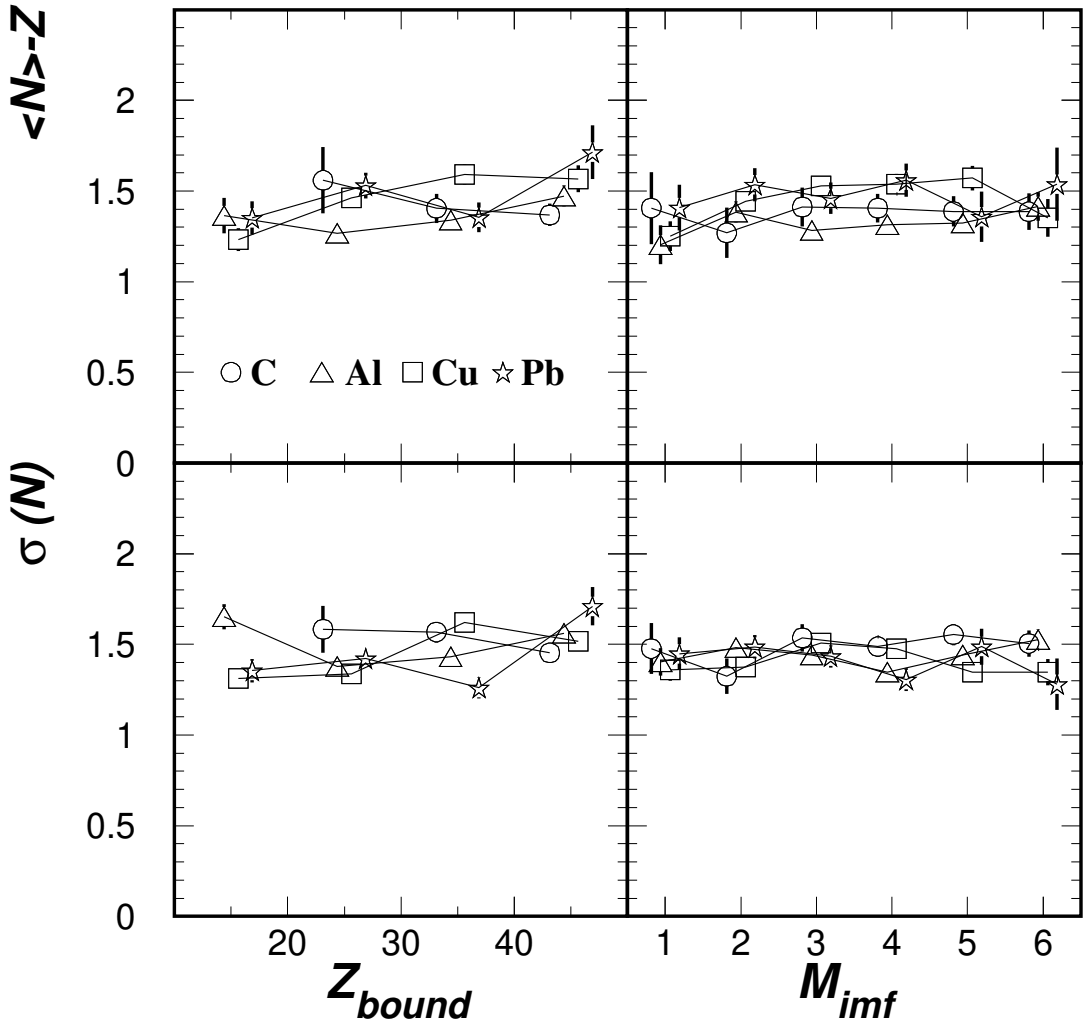
Z-<N>
σ (N)
C
Al
Cu
Pb
0
0.5
1
1.5
2
20
30
40
1
2
3
4
5
6
Z_bound
M_imf