Molecular dynamics and Monte Carlo simulations for the structure of the aqueous trimethylammonium chloride solution in the 0.2-1 molar range

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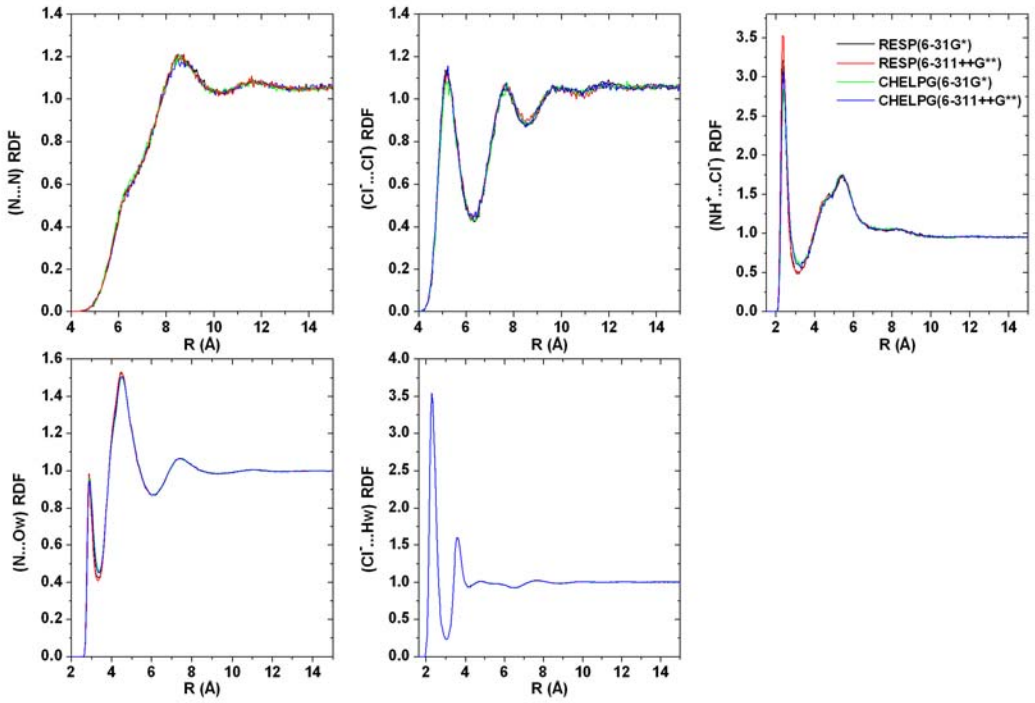

Figure S1. RDFs for case (1), about molar solution. Comparison of the different charge models.

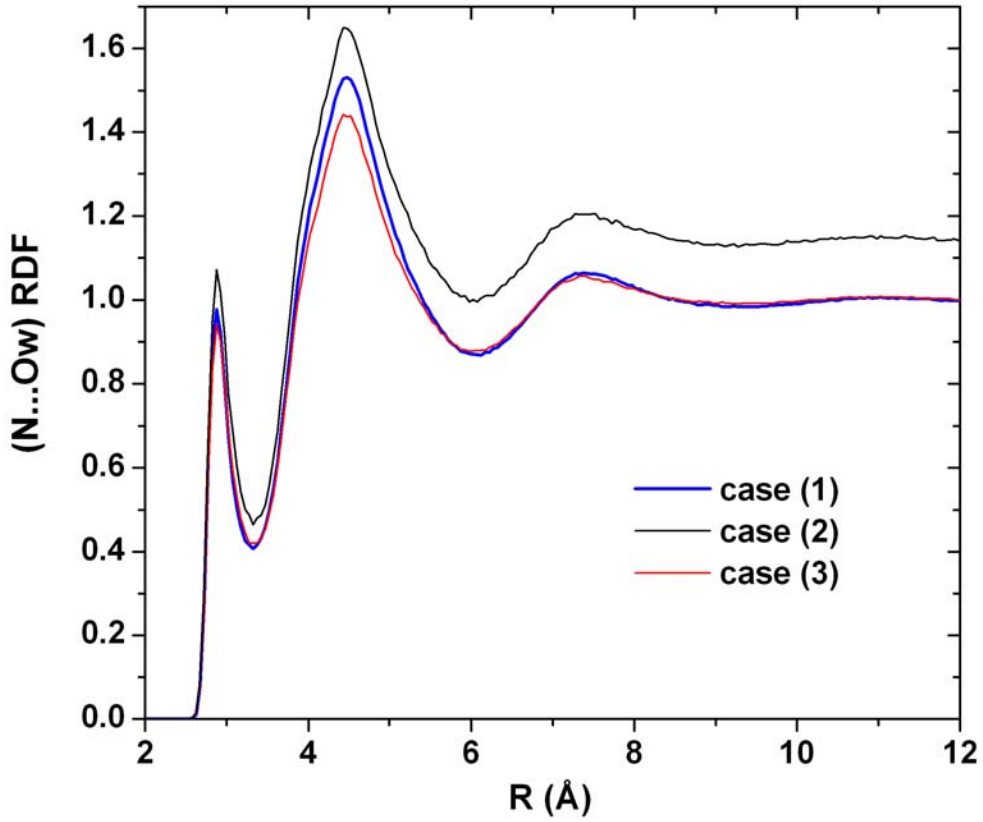

Figure S2: N—O_W Radial Distribution Functions for case (1) (blue), case (2) (black) and case (3) (red).