SAPT: License

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Dear Colleagues,

With regards to the ab initio symmetry-adapted perturbation theory (SAPT) codes, together with its associated auxilary and utility programs, with which you have supplied me a copy, I agree to the following conditions:

1. I will not supply a copy of the code to anyone outside my institution or corporation for any reason whatsoever. Instead, I will refer any requests for copies of the program to you at University of Delaware. This in no way limits my making of copies of the code for backup purposes, or for running on more than one computer system at my institution.

2. I understand that the copyright or ownership rights to SAPT are retained by the authors of SAPT. I will not incorporate any part of SAPT into any other program system, either for sale or for non-profit distribution, without written permission from the authors of SAPT.

3. I understand that no large program such as SAPT can be considered to be bug free, and accordingly the authors of SAPT supply the SAPT software on an "as is" basis, with no additional responsibility or liability.

4. If results obtained with SAPT are published in the scientific literature, I will reference the program as: "SAPT2020: An Ab Initio Program for Many-Body Symmetry-Adapted Perturbation Theory Calculations of Intermolecular Interaction Energies" by R. Bukowski, W. Cencek, J. Garcia, P. Jankowski, B. Jeziorski, M. Jeziorska, T. Korona, S. A. Kucharski, V. F. Lotrich, A. J. Misquitta, R. Moszynski, K. Patkowski, R. Podeszwa, F. Rob, S. Rybak, K. Szalewicz, H. L. Williams, R. J. Wheatley, P. E. S. Wormer, and P. S. Zuchowski.
See also:
B. Jeziorski, R. Moszynski, and K. Szalewicz "Perturbation Theory Approach to Intermolecular Potential Energy Surfaces of van der Waals Complexes", Chem. Rev. 94, 1887-1930 (1994).

Sincerely yours.

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