## ADDITIONS AND CORRECTIONS

## 2009, Volume 113A

Pankaj Kulshrestha, N. Sukumar, Jane S. Murray,\*
Rossman F. Giese, and Troy D. Wood: Computational
Prediction of Antibody Binding Sites on Tetracycline Antibiotics: Electrostatic Potentials and Average Local Ionization
Energies on Molecular Surfaces

Page 756. Reference 47 is in error. The correct reference should be: Aga, D. S.; Goldfish, R.; Kulshrestha, P. *Analyst* **2003**, *128*, 658.

In addition, the binding isotherm data presented in Figure 1 were measured in Dr. Diana Aga's laboratory under NSF sponsorship (NSF Career Grant CHE-0233700) and should have been cited in the acknowledgments.

10.1021/jp903095g Published on Web 04/22/2009

## 2006, Volume 110A

**Edward F. C. Byrd\* and Betsy M. Rice:** Improved Prediction of Heats of Formation of Energetic Materials Using Quantum Mechanical Calculations

Pages 1005–1013. The names of the group equivalents in Table 3 were mislabeled. The results as published in the paper are consistent with the new, properly labeled, group equivalents shown in the corrected Table 3.

TABLE 3: Atom or Group Equivalent Energies and Parameters for Equations 3 and 4 using B3LYP/6-31G\* Geometries, B3LYP/6-311++G(2df, 2p) Energies, and Electrostatic Surface Potential Mappings

atom or group equivalent	$\varepsilon$ (hartree)	eq 3		eq 4	
C	-38.123748	a (kcal/mol-Å <sup>-1</sup> )	2.130167	a (kcal/mol-Å <sup>-4</sup> )	0.000267
Н	-0.597580	b (kcal/mol)	0.930065	b (kcal/mol)	1.650087
N	-54.785466	c (kcal/mol)	-17.843973	c (kcal/mol)	2.966078
0	-75.187087				
C'	-38.129456				
N'	-54.788487				
O'	-75.186033				
$C-NO_2$	-205.160396				
$N-NO_2$	-205.163484				
$O-NO_2$	-205.166631				
$C-N_3$	-164.364907				
X-NO	-129.976691				

We apologize for any inconvenience this might have caused to the readers. We would like to thank Dr. Didier Mathieu for bringing these typographical errors to our attention.

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