

Are the molecular non-covalent interactions of alkali cation-benzene ($M^+XC_6H_5$) complexes confined to a specific carbon atom or propagated through space effects of the substituents? An intense reinvestigation using LFER

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Supplementary Figures:

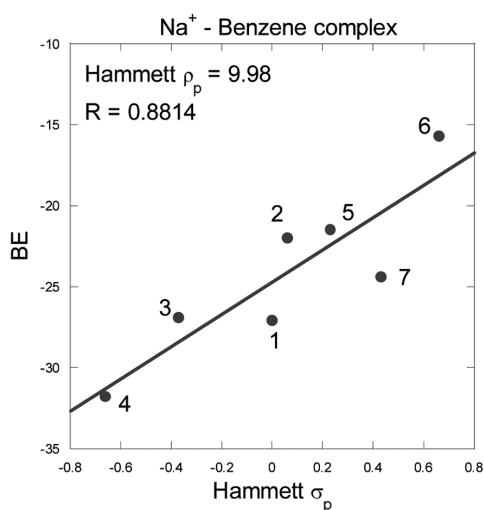


Fig. 5

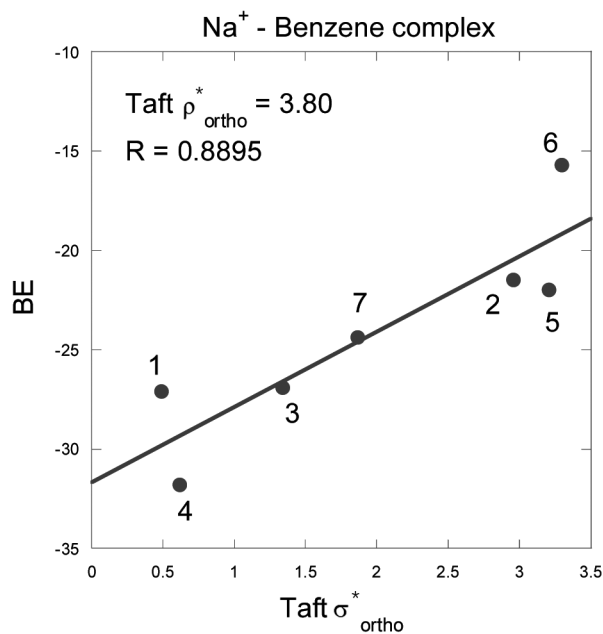


Fig. 7

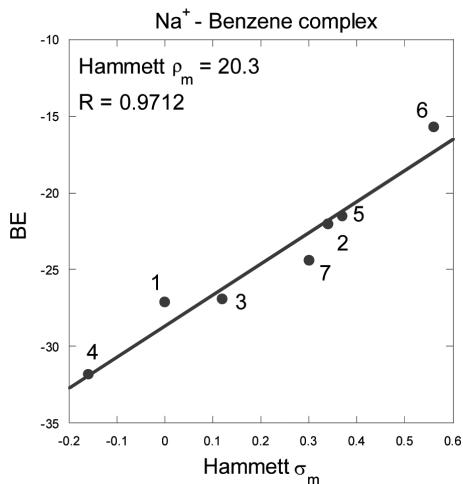


Fig. 6

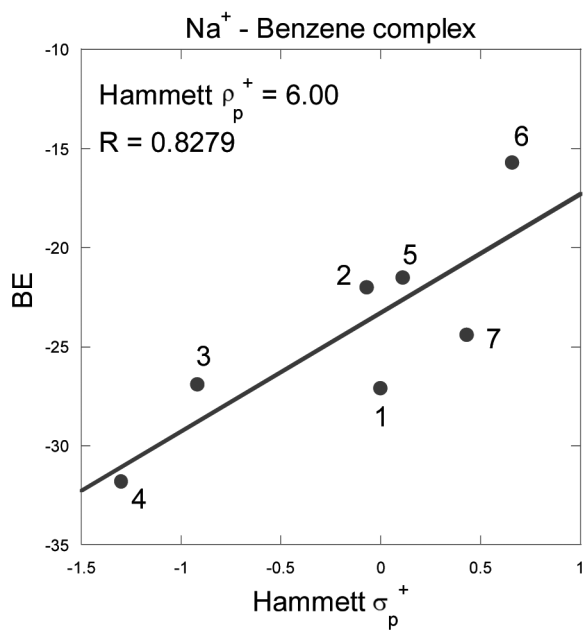


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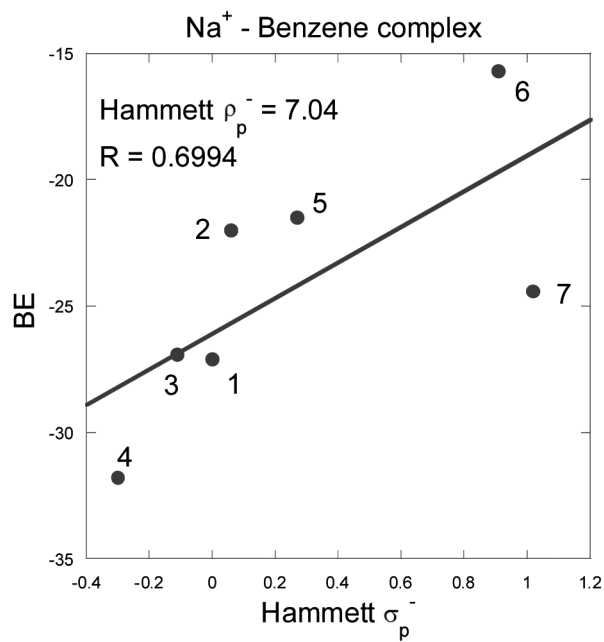


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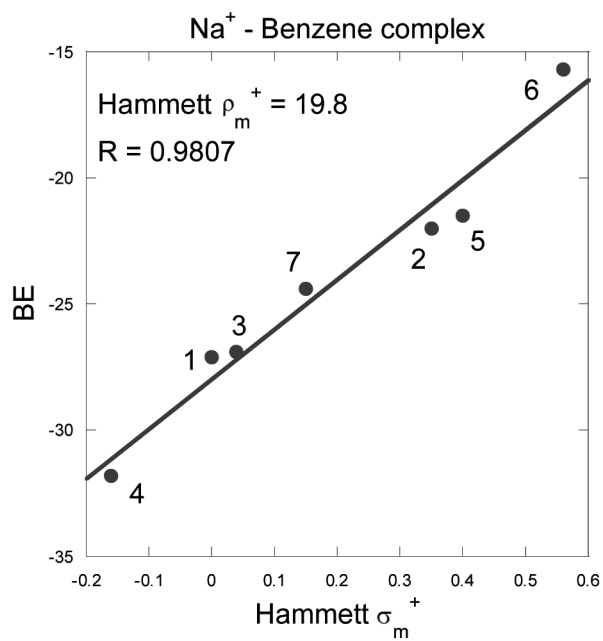


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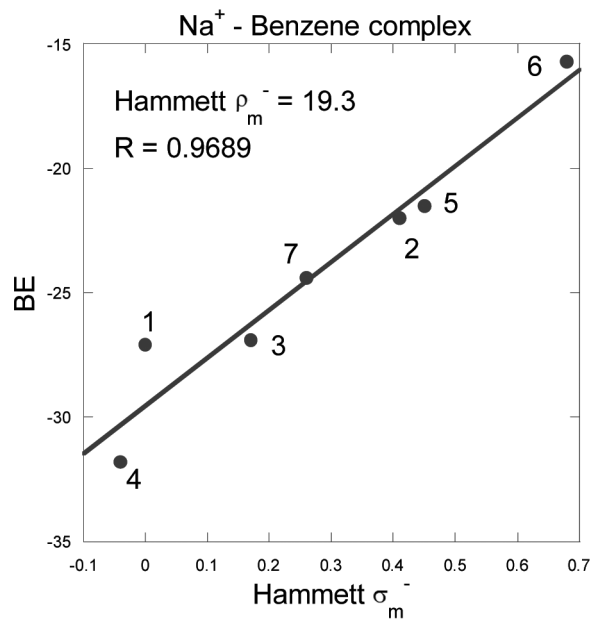


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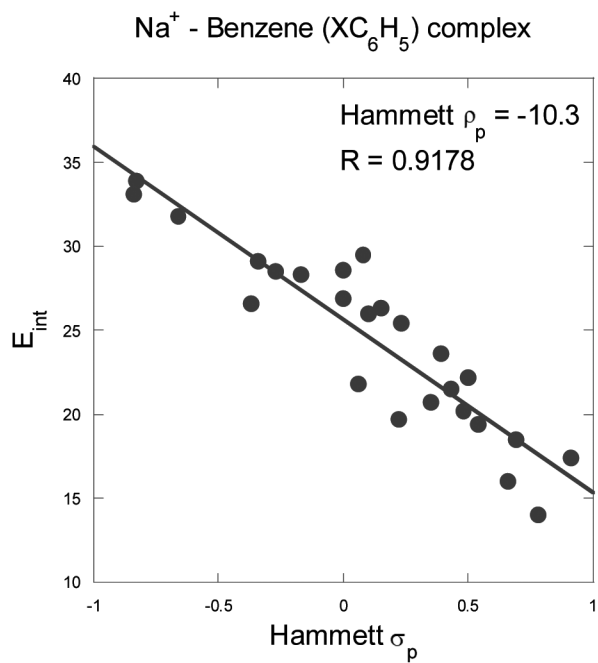


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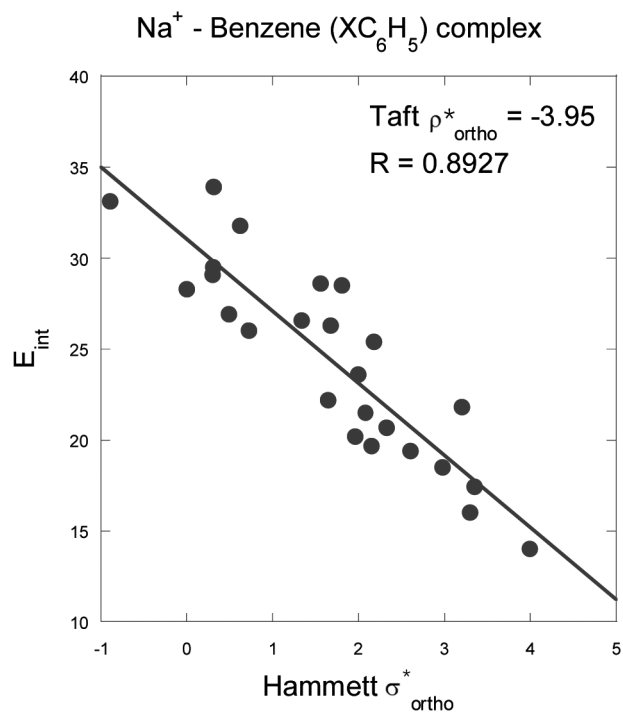


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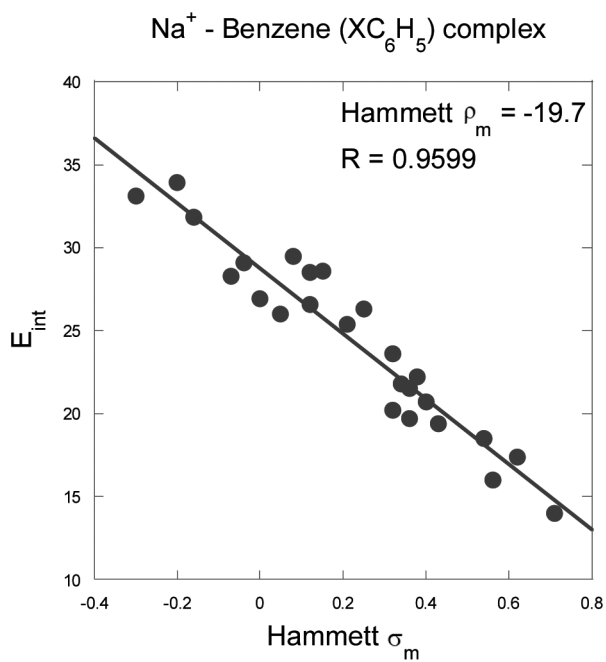


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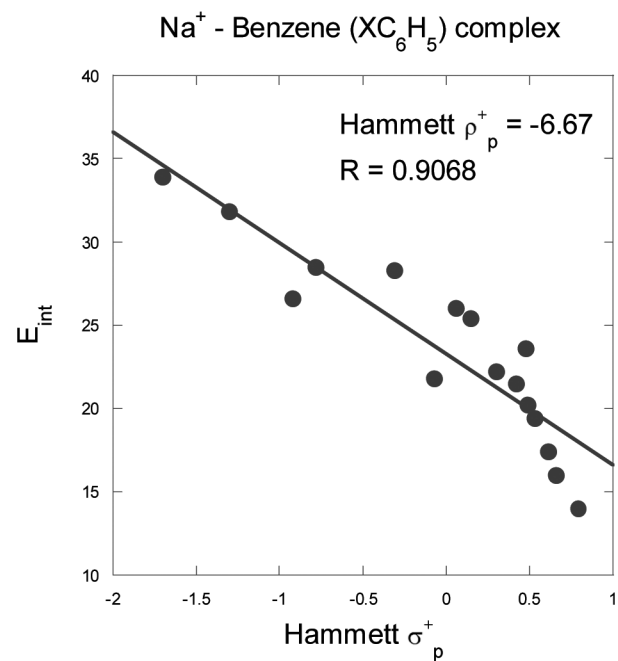


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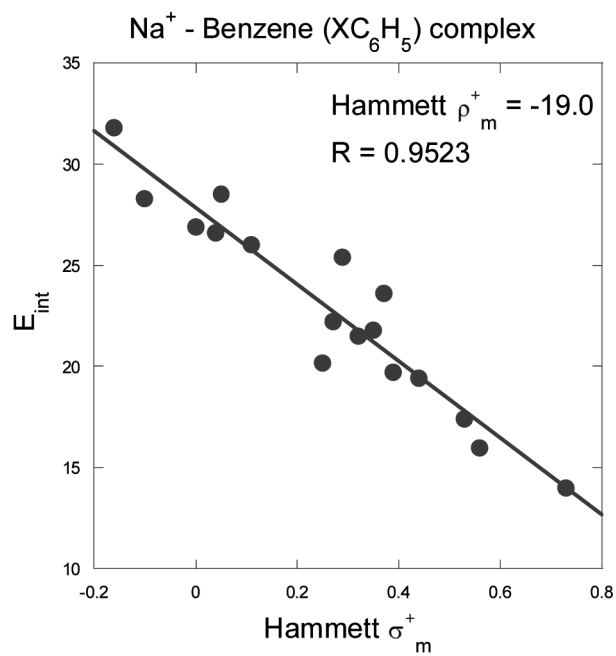


Fig. 16

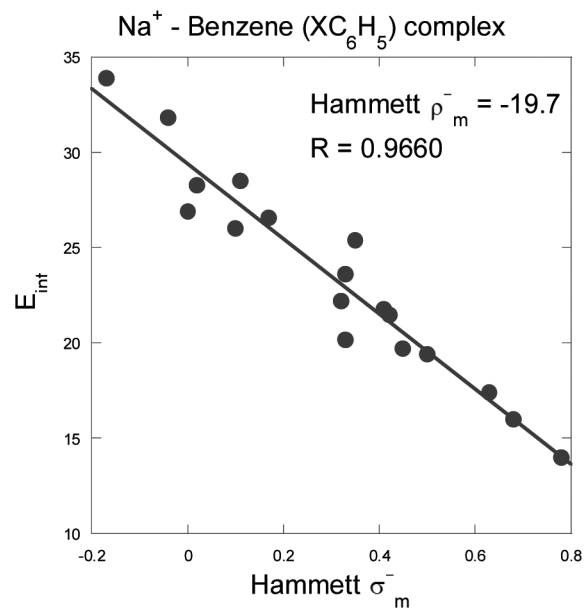


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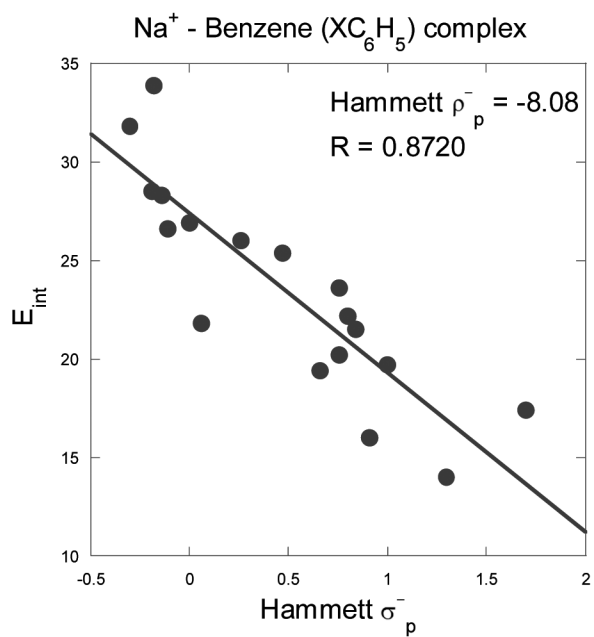


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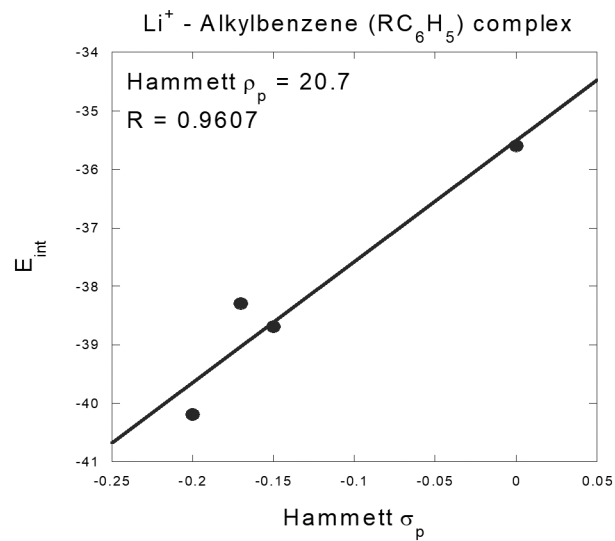


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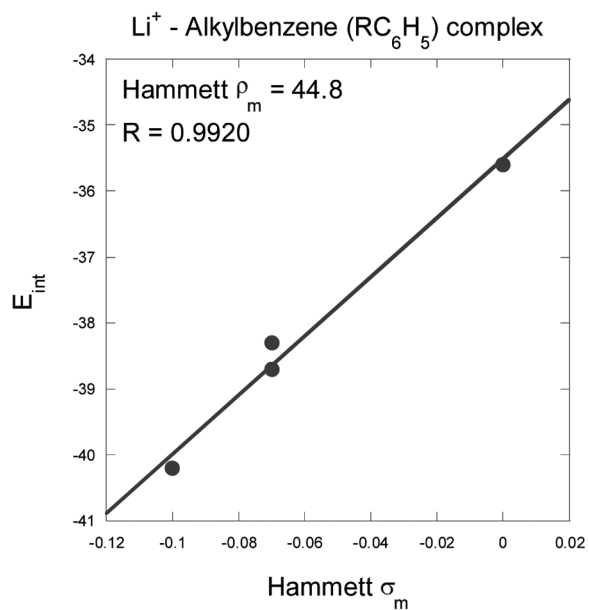


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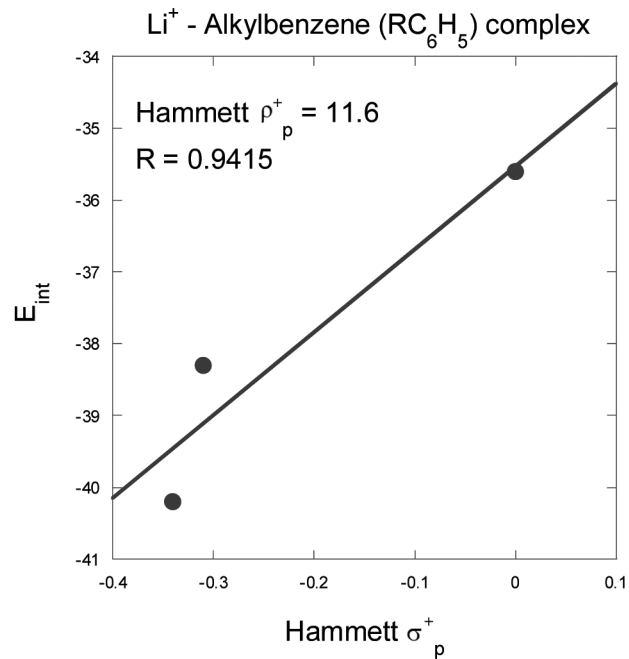


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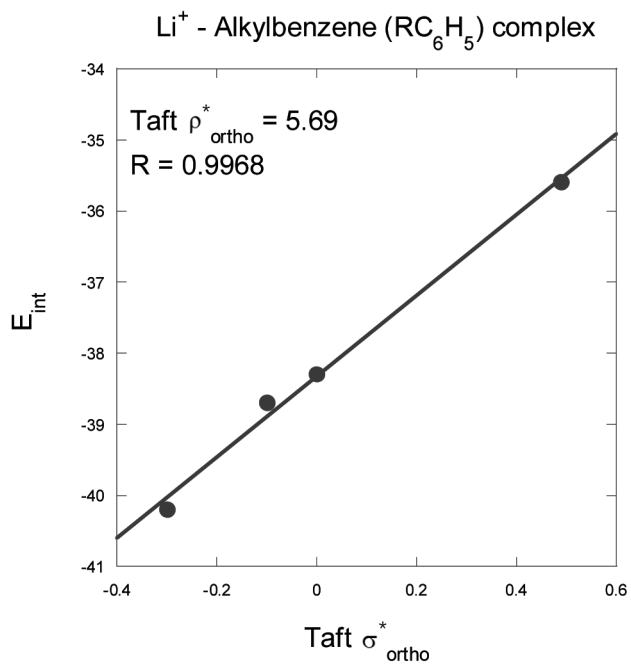


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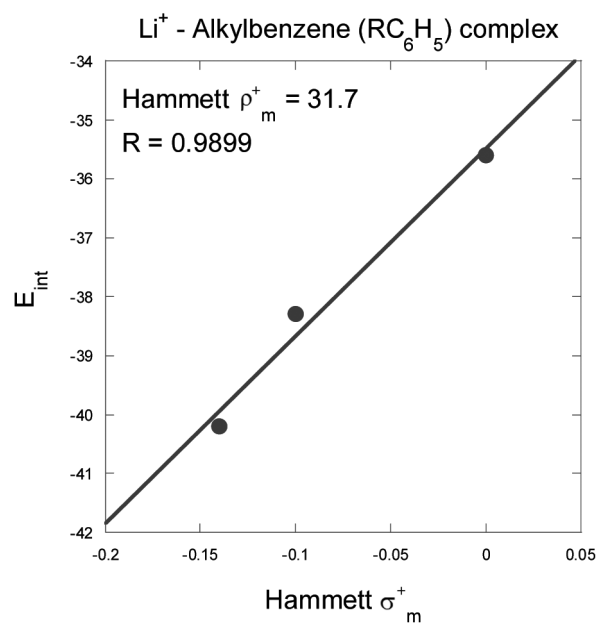


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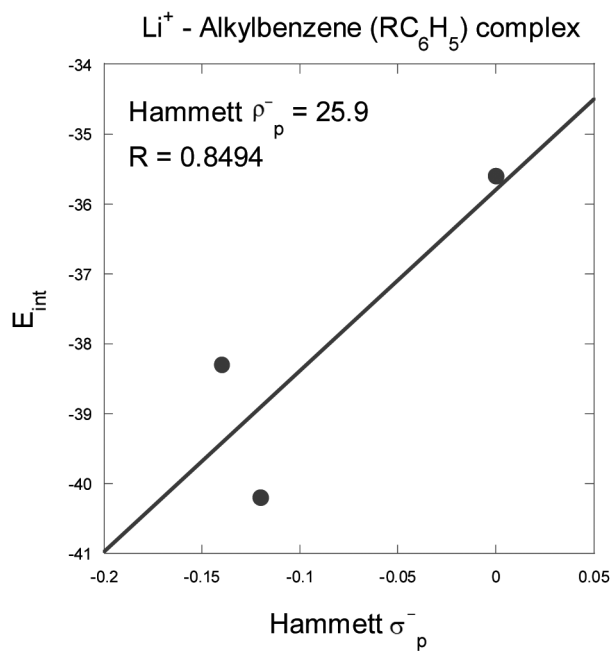


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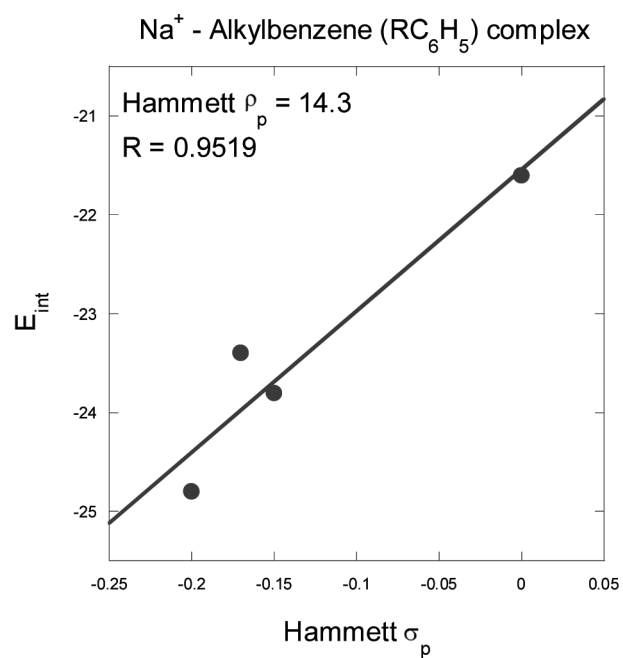


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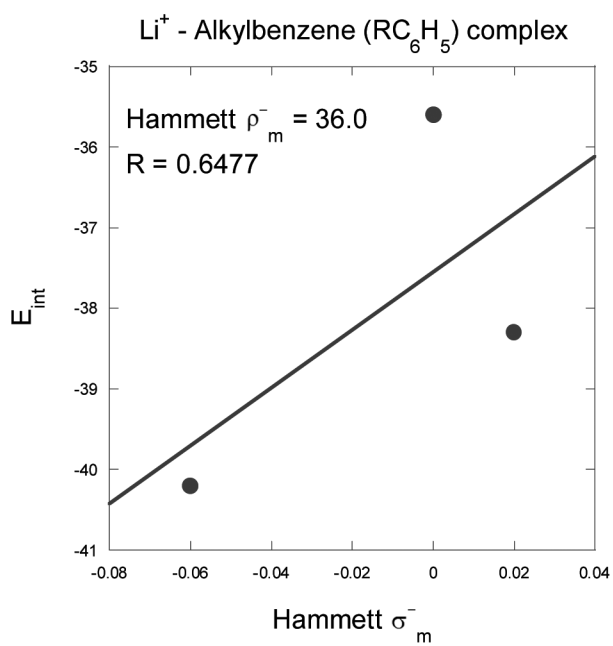


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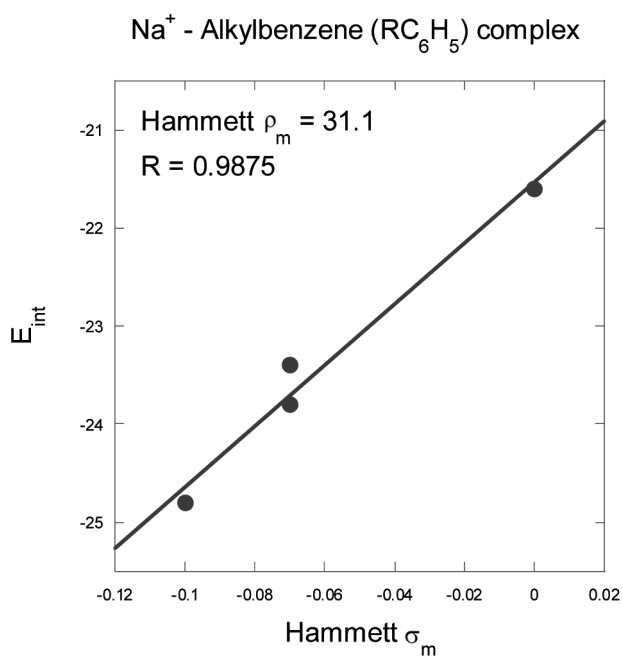


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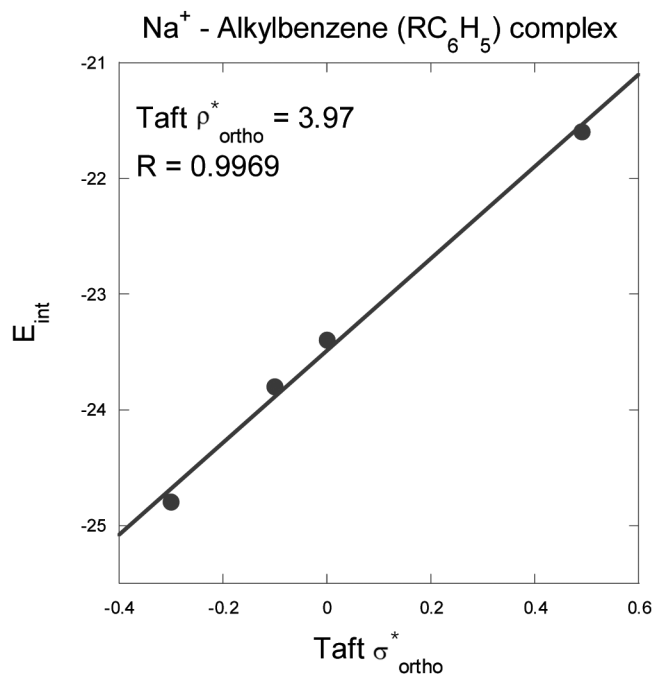


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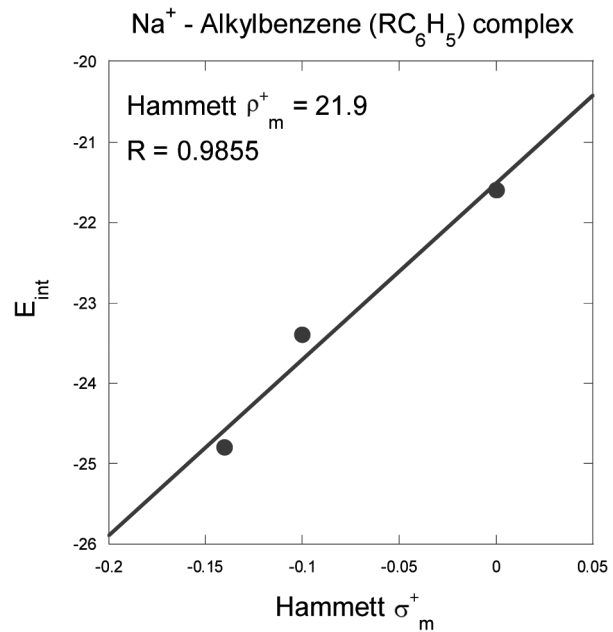


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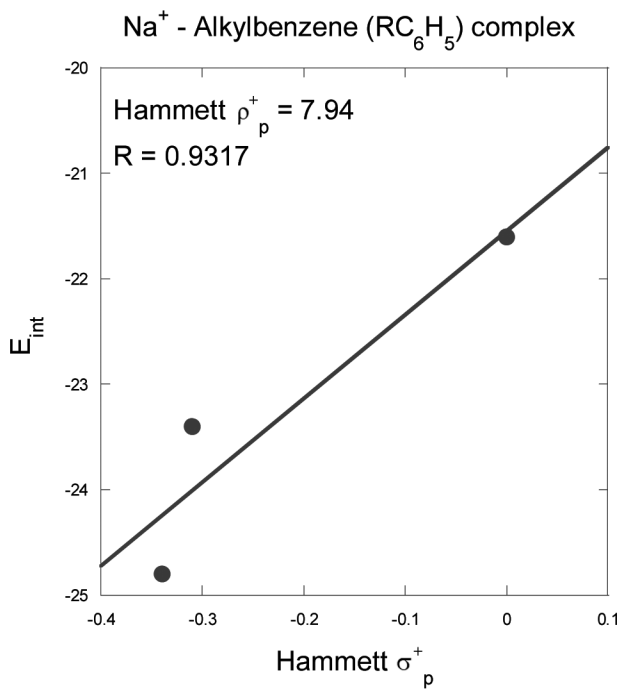


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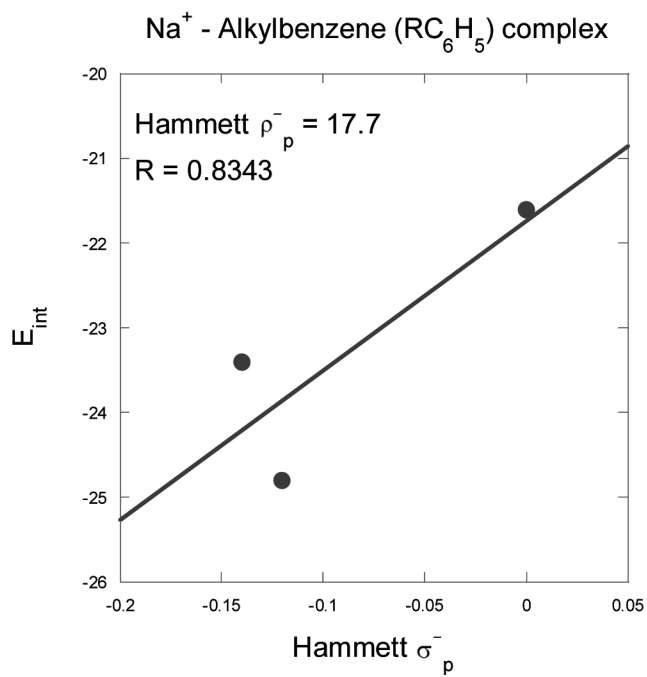


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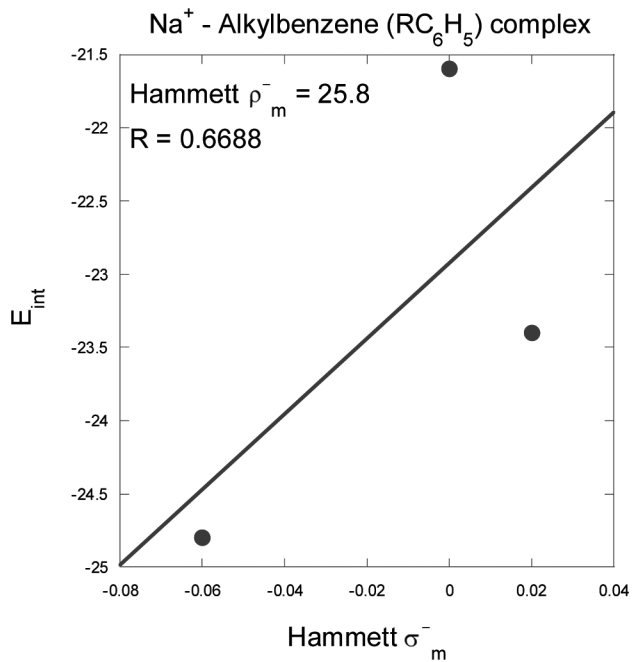


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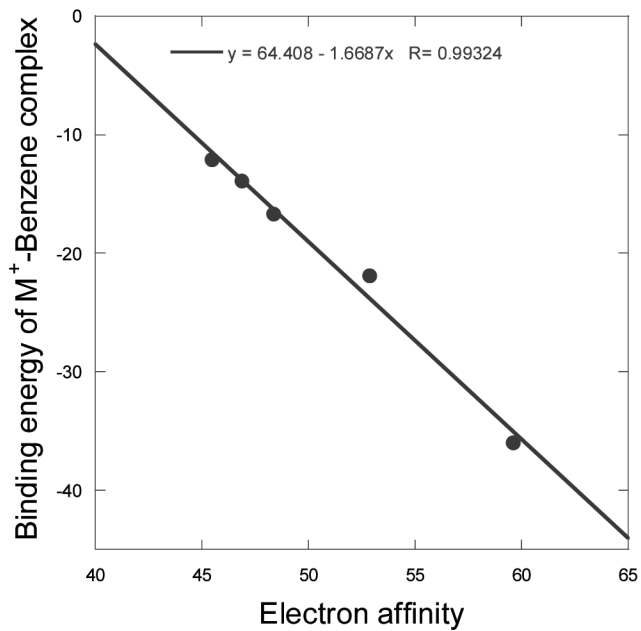


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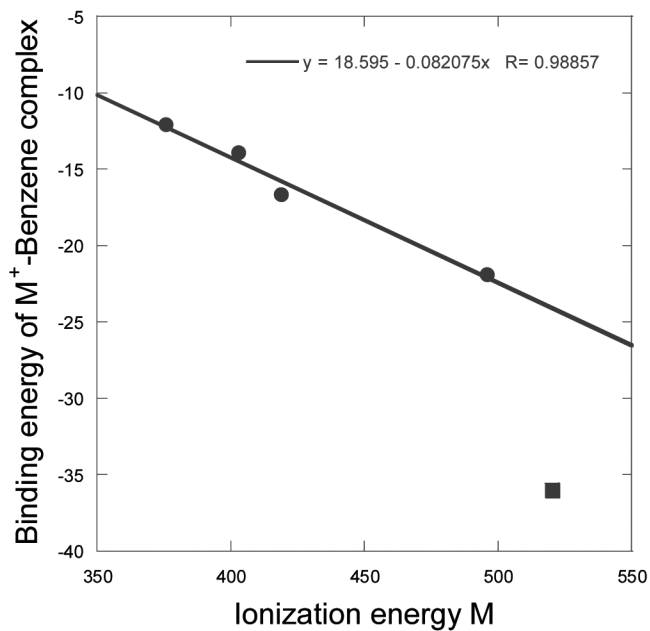


Fig. 33