

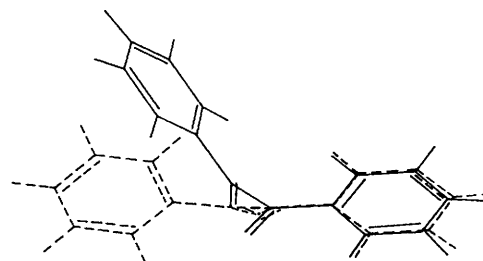
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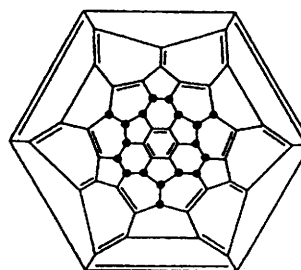
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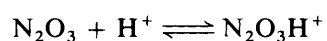
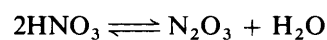



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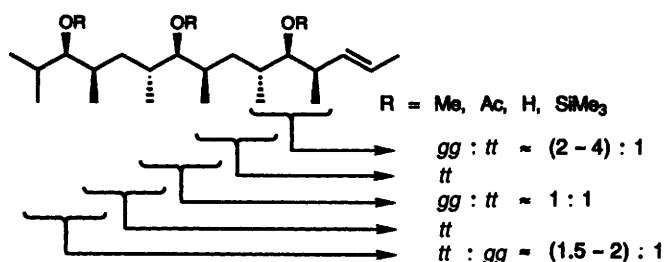
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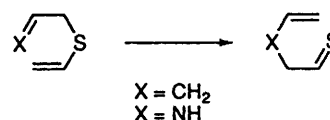
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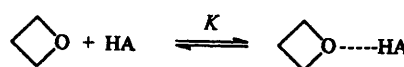
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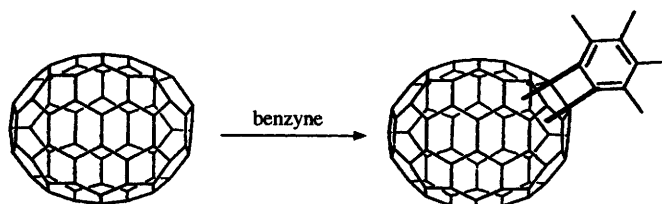
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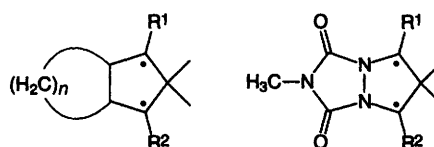
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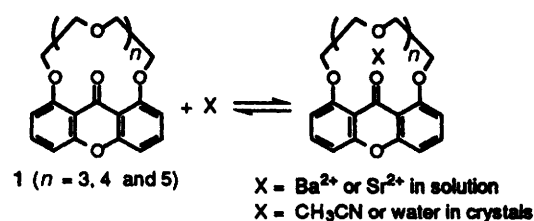
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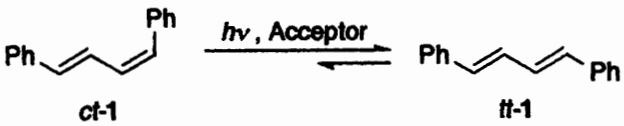
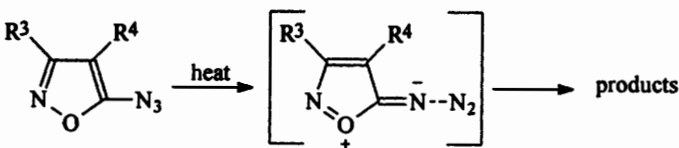
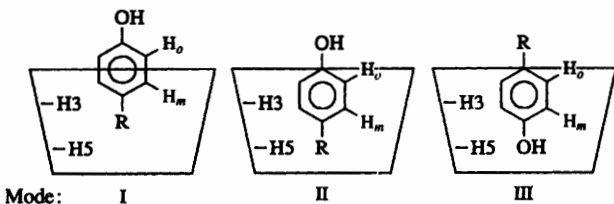
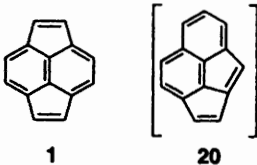
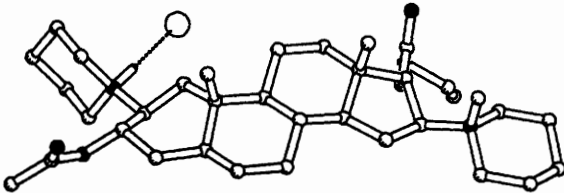
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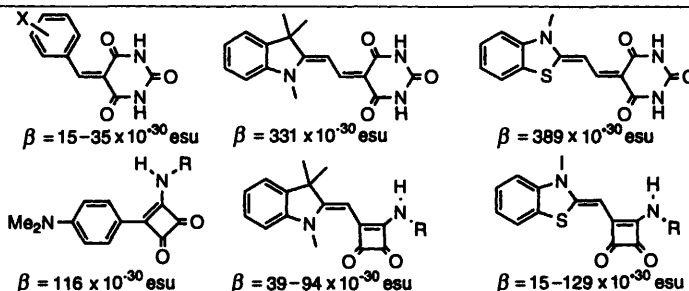
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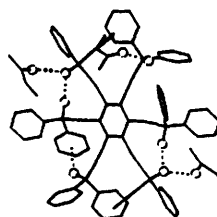
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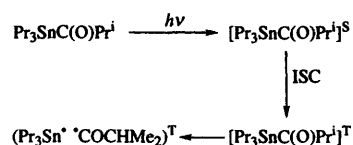
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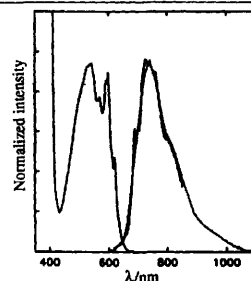
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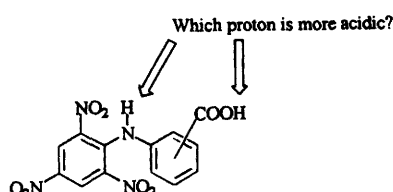
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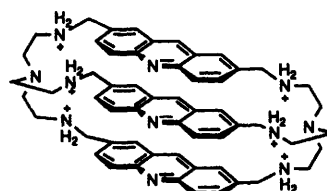
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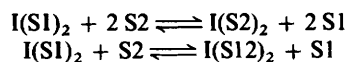
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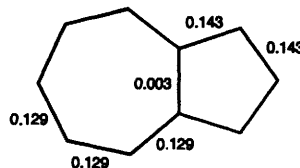
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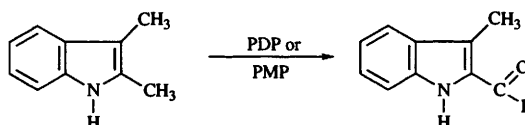
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The bond resonance energy defined for each  $\pi$  bond is useful for examining the geometric origin of aromaticity and reactivity in a polycyclic conjugated system

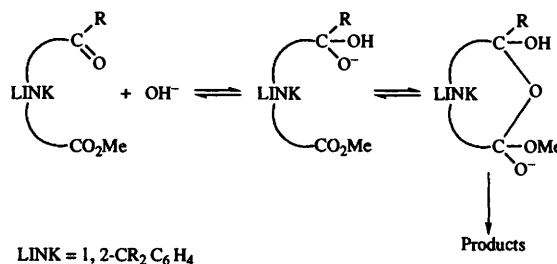
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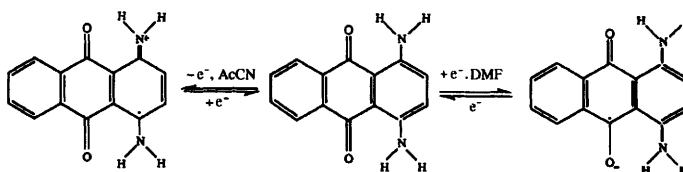
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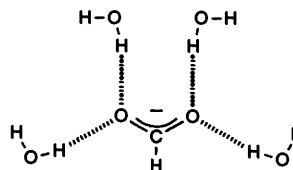
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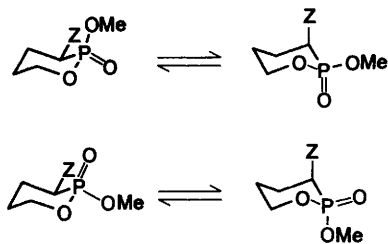
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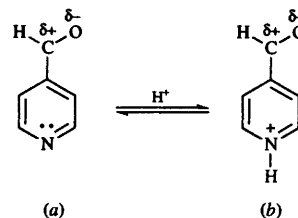
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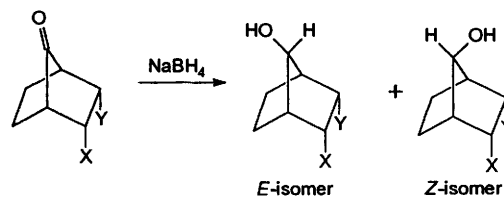
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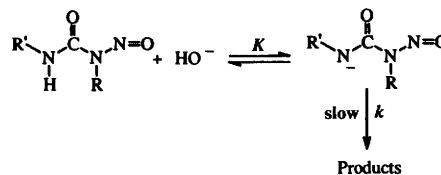
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