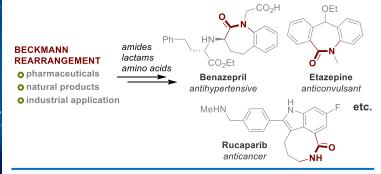
Regioselectivity study of Beckman reaction as a pathway for synthesis of bicyclic anilines and benzylamines



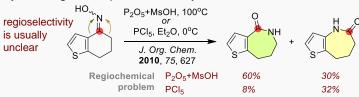
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Background of the project

The Beckmann rearrangement is a highly versatile and industrially important reaction that plays a crucial role in the production of polymers & pharmaceuticals due to its ability to efficiently produce lactams and amides



OBSTACLE: Low regioselectivity of the Beckmann rearrangement of bicyclic aryl ketones poses a great problem to its wide application in synthetic organic and process chemistry:



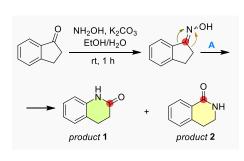
THIS PROJECT:

- * bicyclic aryl ketones with cyclopentane and cyclohexane moieties
- * A wide range of EWG, 'neutral', and EDG substituents placed in the aromatic ring in different positions
- thoroughly optimized rearrangement conditions

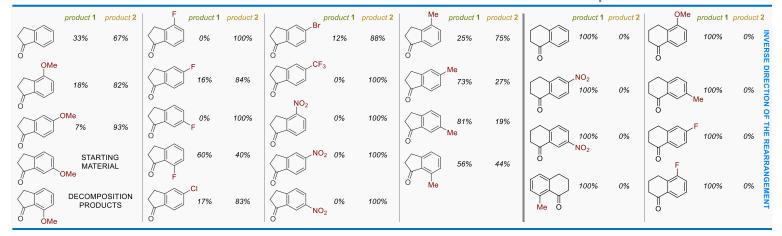


Contact

Optimization of the rearrangement conditions and 'structure/direction' relationships



- CONDITIONS A 1) MsCl, DCM, rt, 1 h (0%) 2) MsCl, DCM, rt, 3 h (3%) 3) MsCl, DCM, rt, 12 h (7%) MsCl. DCM, rt. 24 h (9%) 5) TsCl, Pyridine, rt, 3 h (4%) 6) TsCl. Pyridine, rt, 16 h (15%) 7) TsCl, Pyridine, 60 °C, 3 h (18%) 8) TsCl. Pvridine, 60 °C, 16 h (34%) 9) SOCl₂, THF, rt, 1 h (7%) 10) SOCI₂, THF, rt, 3 h (13%) 11) SOCI₂, THF, 50 °C, 1 h (68%) 12) SOCI₂ THF, 80 °C, 1 h (84%) 13) H₂SO₄ (conc.), rt, 3 h (5%)
 - 14) H₂SO₄ (conc.), rt, 16 h (8%) 15) H₂SO₄ (conc.), 60 °C, 3 h (15%) Multiparameter optimization 16) H₂SO₄ (conc.), 60 °C, 16 h (33%) of rearrangement conditions 17) H₂SO₄ (conc.), 100 °C, 1 h (68%) includes different reagents, 18) H₂SO₄ /H₂O (20%), rt, 16 h (5%) solvents, temperature modes, 19) H₂SO₄ /H₂O (20%), 60 °C, 3 h (7%) 20) H₂SO₄ /H₂O (20%), 60 °C, 16 h (17%) 21) PPA, 60 °C, 3 h (4%) 22) PPA, 60 °C, 16 h (15%) 23) PPA, 100 °C, 1 h (54%) 24) PPA, 100 °C, 3 h (75%) 25) PPA, 120 °C, 15 min (86%) 26) PPA, 150 °C, 5 min (87%)
 - and reaction times Identified the three reaction conditions providing the best results, which were selected for the further substrates screening



Future horizons

- DFT calculations and building a predicting model for the studied aryl ketones allowing for the efficient expansion of the lactams chemical space
- * To involve in the project a wide range of heteroaromatic ketones (208 substrates in total, n = 1, 2, Q = heteroatom, A = heteroatom or carbon)

CURENTLY ONGOING