Dean J Tantillo

List of Publications by Year in descending order

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347 papers 11,970 citations

53 h-index 89 g-index

403 all docs 403 docs citations

403 times ranked

10934 citing authors

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| 3 | Evaluating the Accuracy of the QCEIMS Approach for Computational Prediction of Electron Ionization Mass Spectra of Purines and Pyrimidines. Metabolites, 2022, 12, 68. | 1.3 | 4 |
| 4 | Solvation Effects in Organic Chemistry. Journal of Organic Chemistry, 2022, 87, 1599-1601. | 1.7 | 11 |
| 5 | The Role of Through-Bond Stereoelectronic Effects in the Reactivity of 3-Azabicyclo[3.3.1]nonanes. Journal of Organic Chemistry, 2022, 87, 3378-3388. | 1.7 | 3 |
| 6 | Roads Not Taken: Mechanism and Origins of Regio- and Chemoselectivity of Directed Co ^{III} -Catalyzed Alkenylation of <i>N</i> -Pyridyl 2-Pyridone. Organometallics, 2022, 41, 937-947. | 1.1 | 2 |
| 7 | Divergent Asymmetric Synthesis of Panowamycins, TMâ€135, and Veramycin F using C–H Insertion with Donor/Donor Carbenes. Angewandte Chemie - International Edition, 2022, , . | 7.2 | 3 |
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| 16 | 1-BENZYLSPIRO[PIPERIDINE-4,1′-PYRIDO[3,4-b]indole] â€~co-potentiators' for minimal function CFTR mut European Journal of Medicinal Chemistry, 2021, 209, 112888. | ants. 2.6 | 7 |
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| 25 | Trapping a cross-linked lysine–tryptophan radical in the catalytic cycle of the radical SAM enzyme SuiB. Proceedings of the National Academy of Sciences of the United States of America, 2021, 118, . | 3.3 | 29 |
| 26 | Quantum Chemistry Calculations for Metabolomics. Chemical Reviews, 2021, 121, 5633-5670. | 23.0 | 47 |
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