# SPOTLIGHTS ...

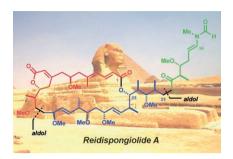
### Natural Products Synthesis

I. Paterson,\* K. Ashton, R. Britton,G. Cecere, G. Chouraqui,G. J. Florence, H. Knust, J. Stafford

Total Synthesis of (-)-Reidispongiolide A, an Actin-Targeting Macrolide Isolated from the Marine Sponge *Reidispongia* coerulea

Chem. Asian J.

DOI: 10.1002/asia.200700357



An answer fit for the Sphinx: The stereocontrolled total synthesis of (-)-reidispongiolide A uses an aldol-based strategy to construct the macrolactone and coupling of a derived aldehyde with an *N*-vinylformamide-containing ketone to install the full side chain. There are two possible routes to introduce the 2*E*,4*E* dienoate, and a fragment coupling by using a Mukaiyama aldol reaction installs the C13 stereocenter.

# Antitumor Agents

F. Wehner,\* A. Nören-Müller, O. Müller, I. Reis-Corrêa, Jr., A. Giannis, H. Waldmann

Indoloquinolizidine Derivatives as Novel and Potent Apoptosis Inducers and Cell-Cycle Blockers

ChemBioChem

DOI: 10.1002/cbic.200700558

A view to a kill: 11000 natural-product-derived and -inspired compounds were screened for potential apoptosis inducers in three human tumour cell lines. Seven indoloquinolizidine derivatives, one of which is shown in the figure, were identified. These compounds had  $IC_{50}$  values as low as 2  $\mu mol \, L^{-1}$  for inhibiting cell proliferation. Further analysis indicated that these effects were related to an arrest of cells in the  $G_2M$  phase of the cell cycle.

# Fluorescent Proteins

C. Blum,\* A. J. Meixner, V. Subramaniam\*

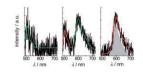
Spectral Versatility of Single Reef Coral Fluorescent Proteins Detected by Spectrally-Resolved Single Molecule Spectroscopy

ChemPhysChem

DOI: 10.1002/cphc.200700784



Single molecule spectroscopy



**Exciting chromophores:** Spectrally resolved single-molecule spectroscopy of members of the DsRed family of fluorescent proteins enables identification and characterization of different spectral forms for all variants (see picture).

The transitions between different spectral forms are also followed, drawing conclusions on the underlying molecular origins of the various spectral species.

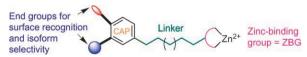
# Drug Design

A. P. Kozikowski,\* Y. Chen, A. M. Gaysin, D. N. Savoy, D. D. Billadeau, K. H. Kim

Chemistry, Biology, and QSAR Studies of Substituted Biaryl Hydroxamates and Mercaptoacetamides as HDAC Inhibitors—Nanomolar-Potency Inhibitors of Pancreatic Cancer Cell Growth

ChemMedChem

DOI: 10.1002/cmdc.200700314

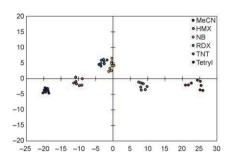


**Isoform selectivity**: structurally unique HDAC inhibitors are equipped with an amino acid residue that serves as a potential isoform-differentiating, surface-recognition element. The surface-rec-

ognition group is connected through the usual carbon linker to either a hydroxamate or a mercaptoacetamide group that chelates the catalytic site zinc ion.

# ... ON OUR SISTER JOURNALS





Sensing with suds: Using pattern recognition, commercially available fluorophores and surfactants are able to sense nitrated explosives with a detection limit of 19 µм (see figure).

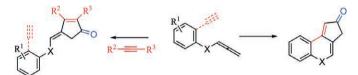
# **Explosive Detection**

A. D. Hughes, I. C. Glenn, A. D. Patrick, A. Ellington, E. V. Anslyn\*

**A Pattern Recognition Based** Fluorescence Quenching Assay for the **Detection and Identification of Nitrated Explosive Analytes** 

Chem. Eur. J.

DOI: 10.1002/chem.200701546



Pauson-Khand reactions of functionalized allenes with different alkynes give monocyclic cyclopentenones with generally high regio- and stereoselectivities. The allenes react with the external double bonds, giving cyclopentenones

with exocyclic double bonds at their β positions, mainly with E stereochemistry. Some intramolecular reactions with allenynes connected through aromatic rings are described.

### Allenic Pauson-Khand Reactions

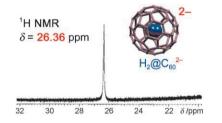
Á. González-Gómez, L. Añorbe, A. Poblador, G. Domínguez, J. Pérez-Castells\*

**Intermolecular and Intramolecular** Pauson-Khand Reactions of **Functionalized Allenes** 

Eur. J. Org. Chem.

DOI: 10.1002/ejoc.200701038

**Inside information**: The NMR signal of the molecular hydrogen in H<sub>2</sub>@C<sub>60</sub><sup>2-</sup> appears at extraordinarily low field  $(\delta = 26.36 \text{ ppm})$  relative to that of neutral  $H_2@C_{60}$ . This can be explained by the decrease in aromaticity of the C<sub>60</sub> cage upon 2e reduction. Spherical delocalization of the added two electrons over the  $\pi$  system of the fullerene cage is thought to be responsible for the drastic change in fullerene aromaticity.



### Aromaticity

Sustainable Fuels

M. Murata, Y. Ochi, F. Tanabe, K. Komatsu,\* Y. Murata\*

**Internal Magnetic Fields of Dianions** of Fullerene C<sub>60</sub> and Its Cage-Opened **Derivatives Studied with Encapsulated** H<sub>2</sub> as an NMR Probe

Angew. Chem. Int. Ed. DOI: 10.1002/anie.200705285

M. E. Domine, A. C. van Veen, Y. Schuurman,\* C. Mirodatos

DOI: 10.1002/cssc.200700049

**Coprocessing of Oxygenated Biomass** Compounds and Hydrocarbons for the **Production of Sustainable Fuel** 



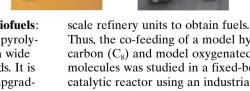
From biomass to tomorrow's biofuels: Pyrolysis oils, obtained by fast pyrolysis of solid biomass, comprise a wide range of oxygenated compounds. It is proposed that after adequate upgrading, these oils may be fed into large-

Thus, the co-feeding of a model hydrocarbon (C<sub>8</sub>) and model oxygenated molecules was studied in a fixed-bed catalytic reactor using an industrial

fluid catalytic cracking catalyst.

On these pages, we feature a selection of the excellent work that has recently been published in our sister journals. If you are reading these pages on a computer, click on any of

the items to read the full article. Otherwise please see the DOIs for easy online access through Wiley



InterScience.

ChemSusChem