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CORRECTION

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Correction: Enantioselective total synthesis of atisane diterpenoids: (+)-sapinsigin H, (+)-agallochaol C, and (+)-16α, 17-dihydroxy-atisan-3-one

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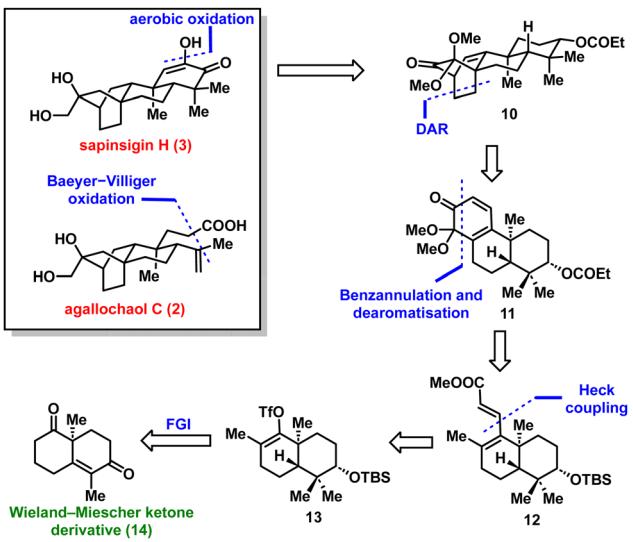
Correction for 'Enantioselective total synthesis of atisane diterpenoids: (+)-sapinsigin H, (+)-agallochaol C, and (+)-16α, 17-dihydroxy-atisan-3-one' by Dattatraya H. Dethe *et al.*, *Chem. Commun.*, 2024, **60**, 7866–7869, https://doi.org/10.1039/D4CC01982B.

The authors regret that Schemes 2 and 3 in the original article incorrectly showed mirror images of the molecular structures. The corrected Schemes 2 and 3 are provided below. The supplementary information file for the original article also contained incorrect mirror images, which have now been corrected in an updated version.



Scheme 2 Retrosynthetic analysis.

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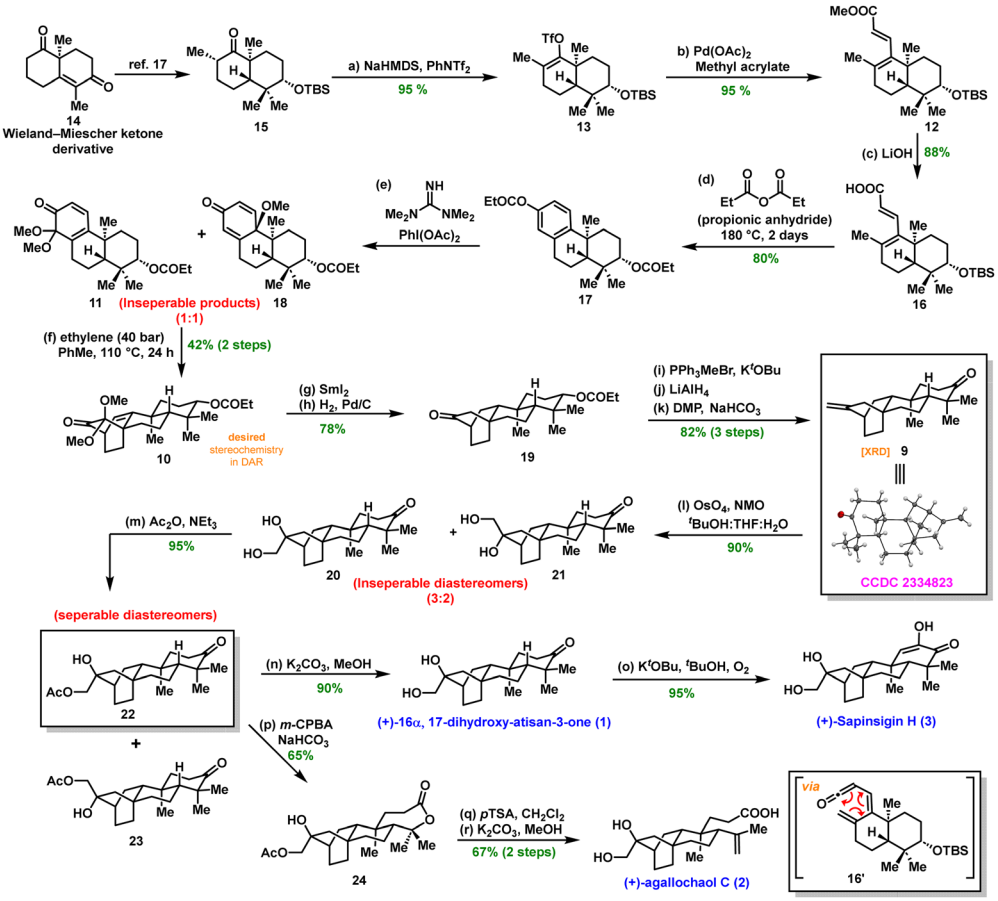

Scheme 3 Total synthesis of (+)-16α, 17-dihydroxy-atisan-3-one, (+)-sapinsigin H and (+)-agallochaol C. Reagent and conditions: (a) NaHMDS (1.9 M, 1.1 equiv.), $PhNTf_2$ (1.1 equiv.), THF, −78 °C, 4 h, 95%; (b) $Pd(OAc)_2$ (10 mol%), methyl acrylate (10.0 equiv.), Et_3N (10.0 equiv.), PPh_3 (0.1 equiv.), 95 °C, 14 h, 95%; (c) LiOH (6.0 equiv.), THF/MeOH/H_2O (3 : 1 : 2), rt, 12 h, 88%; (d) propionic anhydride, 180 °C, 2 days, 80%; (e) $HNC(N(CH_3)_2)_2$ (1.05 equiv.), PIDA (2.4 equiv.), MeOH, −7 °C, 4 h; (f) CH_2CH_2 (40 bar), toluene, 110 °C, 24 h, 40% (2 steps); (g) SmI_2, THF/MeOH (5: 1), rt, 30 min, 85%; (h) H_2, 10% Pd/C, EtOAc, rt, overnight, 92%; (i) PPh_3MeBr (3.0 equiv.), KO*t*-Bu (2.0 equiv.), THF, 0 °C to rt, 2 h, 91%; (j) $LiAlH_4$ (1.0 equiv.), THF, 0 °C to rt, 1 h, 95%; (k) DMP (1.2 equiv.), $NaHCO_3$ (8.0 equiv.), CH_2Cl_2, rt, 1 h, 95%; (l) OsO_4 (2.5 wt% in *t*-BuOH, 0.20 equiv.), NMO (2.0 equiv.), THF/*t*-BuOH/H_2O (1 : 1 : 0.2), rt, 30 h, 90%; (m) Ac_2O (3.0 equiv.), Et_3N (5.0 equiv.), CH_2Cl_2, rt, 24 h, separable diastereomers **22** (57%), **23** (42%); (*n*) K_2CO_3 (2.0 equiv.), MeOH, rt, 1 h, 90%; (o) KO*t*-Bu (7.0 equiv.), *t*-BuOH, O_2, rt, 4 h, 95%; (p) *m*CPBA (1.5 equiv.), $NaHCO_3$ (5.0 equiv.), 0 °C, 1 h, 65%; (q) *p*TSA, CH_2Cl_2, rt, 8 h, 75%; (r) K_2CO_3 (2.0 equiv.), MeOH, rt, 1 h, 90%.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.