Supporting Information

Base Catalysed Domino and Self-domino Michael-Aldol Reactions: One-pot Synthesis of Dispirocyclopentaneoxindoles Containing Multiple Chiral Stereocenters

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1. NMR Spectra of products



Figure 1. ¹H NMR Spectrum of 3a



Figure 2. ¹³C NMR Spectrum of 3a







Figure 4. ¹H NMR Spectrum of 4a



Figure 5. ¹³C NMR Spectrum of 4a



Figure 6. DEPT-135 Spectrum of 4a



Figure 7. ¹H NMR Spectrum of **3b**



Figure 8. ¹³C NMR Spectrum of **3b**



Figure 9. DEPT-135 Spectrum of 3b



Figure 10. ¹H NMR Spectrum of 4b



Figure 11. ¹³C NMR Spectrum of 4b



Figure 12. DEPT-135 Spectrum of 4b



Figure 13. ¹H NMR Spectrum of 3c



Figure 14. ¹³C NMR Spectrum of 3c



Figure 15. DEPT-135 Spectrum of 3c



Figure 16. ¹H NMR Spectrum of 4c



Figure 17. ¹³C NMR Spectrum of 4c



Figure 18. DEPT-135 Spectrum of 4c







Figure 20. ¹³C NMR Spectrum of 3d



Figure 21. DEPT-135 Spectrum of 3d



Figure 22. ¹H NMR Spectrum of 4d



Figure 23. ¹³C NMR Spectrum of 4d



Figure 24. DEPT-135 Spectrum of 4d



Figure 25. ¹H NMR Spectrum of 3a'



Figure 26. ¹³C NMR Spectrum of 3a'





Figure 28. ¹H NMR Spectrum of 3b'



Figure 29. ¹³C NMR Spectrum of 3b'



Figure 30. DEPT-135 Spectrum of 3b'



Figure 32. ¹³C NMR Spectrum of 3c'





Figure 34. ¹H NMR Spectrum of 3d'



Figure 35. ¹³C NMR Spectrum of 3d'



Figure 36. DEPT-135 Spectrum of 3d'









Figure 41. ¹³C NMR Spectrum of 3f'



Figure 42. DEPT-135 Spectrum of 3f'



Figure 43. ¹H NMR Spectrum of 3g'



Figure 44. ¹³C NMR Spectrum of 3g'













Figure 49. ¹H NMR Spectrum of 3i'



Figure 50. ¹³C NMR Spectrum of 3i'



Figure 51. DEPT-135 Spectrum of 3i'









Figure 55. ¹H NMR Spectrum of 3k'



Figure 56. ¹³C NMR Spectrum of 3k'



Figure 57. DEPT-135 Spectrum of 3k'



Figure 58. ¹H NMR Spectrum of 3I'



Figure 59. ¹³C NMR Spectrum of 3I'



Figure 60. DEPT-135 Spectrum of 3l'



Figure 61. ¹H NMR Spectrum of 3m'



Figure 62. ¹³C NMR Spectrum of 3m'



Figure 63. DEPT-135 Spectrum of 3m'



Figure 64. ¹H NMR Spectrum of 3n'



Figure 65. ¹³C NMR Spectrum of 3n'



Figure 66. DEPT-135 Spectrum of 3n'



Figure 67. ¹H NMR Spectrum of 5a



Figure 68. ¹³C NMR Spectrum of 5a



Figure 69. DEPT-135 Spectrum of 5a



Figure 70. ¹H NMR Spectrum of 5b


Figure 71. ¹³C NMR Spectrum of 5b



Figure 72. DEPT-135 Spectrum of 5b





Figure 74. ¹³C NMR Spectrum of 5c







Figure 76. ¹H NMR Spectrum of 5d



Figure 77. ¹³C NMR Spectrum of 5d



Figure 78. DEPT-135 Spectrum of 5d



Figure 79. ¹H NMR Spectrum of 5e



Figure 80. ¹³C NMR Spectrum of 5e







Figure 82. ¹H NMR Spectrum of 5f



Figure 83. ¹³C NMR Spectrum of 5f



Figure 84. DEPT-135 Spectrum of 5f



Figure 85. ¹H NMR Spectrum of 5g



Figure 86. ¹³C NMR Spectrum of 5g



Figure 87. DEPT-135 Spectrum of 5g



Figure 88. ¹H NMR Spectrum of 5h





Figure 90. DEPT-135 Spectrum of 5h



Figure 91. ¹H NMR Spectrum of 5i



Figure 92. ¹³C NMR Spectrum of 5i



Figure 94. ¹H NMR Spectrum of 5j



Figure 95. ¹³C NMR Spectrum of 5j



Figure 96. DEPT-135 Spectrum of 5j



Figure 97. ¹H NMR Spectrum of 5k



Figure 98. ¹³C NMR Spectrum of 5k







Figure 100. ¹H NMR Spectrum of 51



Figure 101. ¹³C NMR Spectrum of 5l



Figure 102. DEPT-135 Spectrum of 51



Figure 103. ¹H NMR Spectrum of 5m



Figure 104. ¹³C NMR Spectrum of 5m



Figure 105. DEPT-135 Spectrum of 5m



Figure 106. ¹H NMR Spectrum of 5n



Figure 107. ¹³C NMR Spectrum of 5n



Figure 108. DEPT-135 Spectrum of 5n



Figure 109. ¹H NMR Spectrum of 50



Figure 110. ¹³C NMR Spectrum of 50







Figure 112. ¹H NMR Spectrum of 5p



Figure 113. ¹³C NMR Spectrum of 5p



Figure 114. DEPT-135 Spectrum of 5p



Figure 115. ¹H NMR Spectrum of 5q



Figure 116. ¹³C NMR Spectrum of 5q



Figure 117. DEPT-135 Spectrum of 5q



Figure 118. ¹H NMR Spectrum of 5r



Figure 120. DEPT-135 Spectrum of 5r



Figure 121. ¹H NMR Spectrum of 5s



Figure 122. ¹³C NMR Spectrum of 5s







Figure 124. ¹H NMR Spectrum of 5t



Figure 125. ¹³C NMR Spectrum of 5t



Figure 126. DEPT-135 Spectrum of 5t

2. HRMS (EI) Spectra of Compounds



Figure 127. Mass spectrum of 3a



Figure 128. Mass spectrum of 4a



Figure 129. Mass spectrum of 3b



Figure 130. Mass spectrum of 4b



Figure 131. Mass spectrum of 3c



Figure 132. Mass spectrum of 4c



Figure 133. Mass spectrum of 3d







Figure 135. Mass spectrum of 3a'



Figure 136. Mass spectrum of 3b'



Figure 137. Mass spectrum of 3c'



Figure 138. Mass spectrum of 3d'



Figure 139. Mass spectrum of 3e'



Figure 140. Mass spectrum of 3f'



Figure 141. Mass spectrum of 3g'



Figure 142. Mass spectrum of 3h'


Figure 143. Mass spectrum of 3i'



Figure 144. Mass spectrum of 3j'



Figure 145. Mass spectrum of 3k'



Figure 146. Mass spectrum of 3l'



Figure 147. Mass spectrum of 3m'



Figure 148. Mass spectrum of 3n'



Figure 149. Mass spectrum of 5a







Figure 151. Mass spectrum of 5c



Figure 152. Mass spectrum of 5d





735.2175

332.2307

255.3498

20-

Figure 153. Mass spectrum of 5f



Figure 154. Mass spectrum of 5g



Figure 155. Mass spectrum of 5h



Figure 156. Mass spectrum of 5i



Figure 157. Mass spectrum of 5j



Figure 158. Mass spectrum of 5k



Figure 159. Mass spectrum of 51



Figure 160. Mass spectrum of 5m



Figure 161. Mass spectrum of 5n



Figure 162. Mass spectrum of 50



Figure 163. Mass spectrum of 5p



Figure 164. Mass spectrum of 5q



Figure 165. Mass spectrum of 5r



Figure 166. Mass spectrum of 5s



Figure 167. Mass spectrum of 5t

3. Crystal data and structure refinement

Compound 3a':

Table 1. Crystal data and structure refinement for ksd3.

Identification code	ksd3
Empirical formula	C51 H38 F2 N4 O4
Formula weight	808.85
Temperature	296(2) K
Wavelength	0.71073 A
Crystal system, space group Monoclinic, P 21/c	
Unit cell dimensions	
a = 11.9700(5) A alpha = 90 deg.	
b = 15.5516(8) A beta = 103.1855(19) deg.	
c = 23.1725(11) A gamma = 90 deg.	
Volume	4199.9(3) A^3
Z, Calculated density	4, 1.279 Mg/m^3
Absorption coefficient	0.088 mm^-1
F(000)	1688
Crystal size	0.350 x 0.220 x 0.150 mm
Theta range for data collection 1.590 to 21.936 deg.	
Limiting indices	-12<=h<=11, -16<=k<=14, -22<=l<=24
Reflections collected / uniq	ue $18978 / 5092 [R(int) = 0.0289]$
Completeness to theta $= 25.242 66.9 \%$	
Absorption correction	None
Refinement method	Full-matrix least-squares on F^2
Data / restraints / parameter	s 5092 / 0 / 558
Goodness-of-fit on F^2	1.076
Final R indices [I>2sigma(I)] $R1 = 0.0470, wR2 = 0.1222$
R indices (all data)	R1 = 0.0629, wR2 = 0.1361
Extinction coefficient	n/a
Largest diff. peak and hole	0.704 and -0.186 e.A^-3

Compound 5f:

Table 2. Crystal data and structure refinement for fin .

Identification code	fin	
Empirical formula	C49 H34 Cl N O4	
Formula weight	736.22	
Temperature	293(2) K	
Wavelength	0.71073 A	
Crystal system, space group Triclinic, P -1		
Unit cell dimensions		
a = 8.890(5) A alpha = 91.701(5) deg.		
b = 10.257(5) A beta = 90.533(5) deg.		
c = 20.323(5) A gamma = 98.252(5) deg.		
Volume 18	33.0(14) A^3	
Z, Calculated density	2, 1.334 Mg/m^3	
Absorption coefficient	0.154 mm^-1	
F(000)	768	
Crystal size	0.20 x 0.15 x 0.10 mm	
Theta range for data collection 2.01 to 26.61 deg.		
Limiting indices	-10<=h<=11, -12<=k<=12, -25<=l<=25	
Reflections collected / uniqu	e 27769 / 7610 [R(int) = 0.0233]	
Completeness to theta = 26.61 98.9 %		
Absorption correction	Semi-empirical from equivalents	
Max. and min. transmission	0.9848 and 0.9698	
Refinement method	Full-matrix least-squares on F^2	
Data / restraints / parameter	s 7610 / 0 / 497	
Goodness-of-fit on F^2	1.013	
Final R indices [I>2sigma(I])] $R1 = 0.0395$, $wR2 = 0.1103$	
R indices (all data)	R1 = 0.0506, wR2 = 0.1189	
Largest diff. peak and hole	0.261 and -0.251 e.A^-3	