

Supporting Information

**Five Types of Odd-Even Effect and Crystal Structure Changes Brought about by  $\omega$ -Phenylalkyl Groups in  $[\text{Ni}(\text{dmit})_2]$  Complex Salts**

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**Table S1.** Distances/ Å of S...S Interactions in (C2Ph)[Ni(dmit)<sub>2</sub>]

S4...S13 <sup>1)</sup>	3.406(1)
S5...S13 <sup>1)</sup>	3.604(1)
S18...S7 <sup>2)</sup>	3.534(1)

Symmetry codes; 1) x, y, 1+z; 2) -1/2+x, 1.5-y, 1-z;

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**Table S2.** Distances/ Å of S...S Interactions in (C3Ph)[Ni(dmit)<sub>2</sub>]

S1...S7 <sup>1)</sup>	3.631(2)
S2...S2 <sup>2)</sup>	3.614(2)
S2...S6 <sup>1)</sup>	3.500(3)
S3...S9 <sup>3)</sup>	3.298(2)
S4...S7 <sup>2)</sup>	3.617(2)
S5...S9 <sup>3)</sup>	3.577(2)
S11...S17 <sup>4)</sup>	3.644(2)
S12...S16 <sup>4)</sup>	3.474(2)
S11...S18 <sup>5)</sup>	3.649(2)
S13...S19 <sup>6)</sup>	3.237(2)
S13...S20 <sup>6)</sup>	3.537(2)
S19...S19 <sup>7)</sup>	3.310(2)
S19...S20 <sup>7)</sup>	3.556(2)

Symmetry codes; 1) 1-x, 1-y, 1-z; 2) 2-x, 1-y, 1-z; 3) x, 1+y, z; 4) 2-x, 1-y, -z; 5) 1-x, 1-y, -z; 6) x, -1+y, z; 7) 2-x, 2-y, -z;

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**Table S3.** Distances/ Å of S...S Interactions in (C4Ph)[Ni(dmit)<sub>2</sub>]

S3...S9 <sup>1)</sup>	3.540(1)
S6...S4 <sup>2)</sup>	3.624(1)

Symmetry codes; 1) -1+x, y, z; 2) 1-x, -1/2+y, 1/2-z;

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**Table S4.** Distances/ Å of S...S Interactions in (C5Ph)[Ni(dmit)<sub>2</sub>]

S5...S15 <sup>1)</sup>	3.449(3)
S6...S19 <sup>2)</sup>	3.634(2)
S8...S19 <sup>2)</sup>	3.451(2)
S11...S16 <sup>3)</sup>	3.642(2)
S11...S18 <sup>3)</sup>	3.545(2)
S16...S16 <sup>3)</sup>	3.571(3)

Symmetry codes; 1) 1.5-x, -1/2+y, 1.5-z; 2) 1-x, -y, 1-z; 3) 2-x, -y, 2-z;

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