

**Table 1S.** Selected crystal data for the structures not included in Table 2.

	0.0%P ( <b>1a</b> )	15.0%P ( <b>1a</b> )	28.6%P ( <b>1a</b> )	55.6%P ( <b>1a</b> )	79.9%P ( <b>1a</b> )
Chemical formula	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>				
Formula weight	513.65	513.65	513.65	513.65	513.65
Crystal system	Orthorhombic	Orthorhombic	Orthorhombic	Orthorhombic	Orthorhombic
Space group	P <sub>2</sub> 1 <sub>2</sub> 1 <sub>2</sub> 1				
<i>a</i> /Å	9.375 (2)	9.3750 (15)	9.382 (2)	9.3702 (13)	9.3006 (12)
<i>b</i> /Å	10.734 (3)	10.7698 (19)	10.864 (3)	11.016 (2)	11.3515 (16)
<i>c</i> /Å	29.950 (6)	29.857 (4)	29.571 (7)	29.134 (5)	28.334 (4)
<i>V</i> /Å <sup>3</sup>	3013.9 (12)	3014.6 (8)	3014.1 (13)	3007.3 (9)	2991.4 (7)
<i>Z</i>	4	4	4	4	4
<i>D</i> <sub>x</sub> /Mg m <sup>-3</sup>	1.132	1.132	1.132	1.135	1.141
μ/mm <sup>-1</sup>	0.07	0.07	0.07	0.07	0.07
T/K	299(2)	299(2)	299(2)	299(2)	299(2)
Relections collected	14678	14671	14682	14678	14608
Reflections independent	3010	3006	2994	2990	2979
<i>R</i> <sub>int</sub>	0.038	0.038	0.071	0.077	0.056
<i>R</i> , <i>wR</i> ( <i>F</i> <sup>2</sup> >2σ( <i>F</i> <sup>2</sup> ))	0.055, 0.156	0.073, 0.217	0.087, 0.266	0.101, 0.316	0.082, 0.243
Δρ <sub>max</sub> , Δρ <sub>min</sub> /e Å <sup>-3</sup>	0.19, -0.20	0.21, -0.26	0.20, -0.20	0.17, -0.18	0.18, -0.15

	81.2%P ( <b>1a</b> )	87.4%P ( <b>1a</b> )	100.0%P ( <b>1a</b> )
Chemical formula	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>
Formula weight	513.65	513.65	513.65
Crystal system	Orthorhombic	Orthorhombic	Orthorhombic
Space group	P <sub>2</sub> 1 <sub>2</sub> 1 <sub>2</sub> 1	P <sub>2</sub> 1 <sub>2</sub> 1 <sub>2</sub> 1	P <sub>2</sub> 1 <sub>2</sub> 1 <sub>2</sub> 1
<i>a</i> /Å	9.2818 (15)	9.2671 (14)	9.2509 (16)
<i>b</i> /Å	11.418 (2)	11.4608 (18)	11.481 (2)
<i>c</i> /Å	28.183 (5)	28.078 (4)	28.074 (5)
<i>V</i> /Å <sup>3</sup>	2986.8 (9)	2982.1 (8)	2981.7 (9)
<i>Z</i>	4	4	4
<i>D</i> <sub>x</sub> /Mg m <sup>-3</sup>	1.142	1.144	1.144
μ/mm <sup>-1</sup>	0.07	0.07	0.07
T/K	299(2)	299(2)	299(2)
Relections collected	14591	14559	14562
Reflections independent	2968	2960	2959
<i>R</i> <sub>int</sub>	0.053	0.052	0.049
<i>R</i> , <i>wR</i> ( <i>F</i> <sup>2</sup> >2σ( <i>F</i> <sup>2</sup> ))	0.080, 0.246	0.077, 0.226	0.067, 0.178
Δρ <sub>max</sub> , Δρ <sub>min</sub> /e Å <sup>-3</sup>	0.20, -0.21	0.17, -0.20	0.18, -0.14

	15.9%P for A 17.4%P for B ( <b>1b</b> )	20.4%P for A 30.6%P for B ( <b>1b</b> )	25.5%P for A 54.9%P for B ( <b>1b</b> )
Chemical formula	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>	C <sub>33</sub> H <sub>39</sub> NO <sub>4</sub>
Formula weight	513.65	513.65	513.65
Crystal system	Orthorhombic	Orthorhombic	Orthorhombic
Space group	<i>Pna</i> 2 <sub>1</sub>	<i>Pna</i> 2 <sub>1</sub>	<i>Pna</i> 2 <sub>1</sub>
<i>a</i> /Å	22.093 (4)	22.142 (4)	22.179 (4)
<i>b</i> /Å	9.3747 (19)	9.3891 (17)	9.3739 (17)
<i>c</i> /Å	28.939 (6)	28.954 (5)	28.978 (6)
<i>V</i> /Å <sup>3</sup>	5994 (2)	6019.3 (19)	6025 (2)
<i>Z</i>	8	8	8
<i>D</i> <sub>x</sub> /Mg m <sup>-3</sup>	1.138	1.134	1.133
μ/mm <sup>-1</sup>	0.07	0.07	0.07
T/K	299(2)	299(2)	299(2)
Relections collected	27901	28026	28066
Reflections independent	5369	5390	5393
<i>R</i> <sub>int</sub>	0.049	0.052	0.054
<i>R</i> , <i>wR</i> ( <i>F</i> <sup>2</sup> >2σ( <i>F</i> <sup>2</sup> ))	0.076, 0.223	0.077, 0.229	0.088, 0.272
Δρ <sub>max</sub> , Δρ <sub>min</sub> /e Å <sup>-3</sup>	0.18, -0.27	0.18, -0.22	0.19, -0.20