Electronic Supplementary Material (ESI) for CrystEngComm.

Table S1. Atomic coordinates and equivalent isotropic displacement parameters (Å)

Atom	x	y	z	U(eq)
Co(1)	0.7458(1)	0.7458(1)	0.3115(1)	0.005(1)
Co(2)	0.0	0.7310(1)	0.5	0.006(1)
P(1)	0.0	0.5	0.25	0.007(1)
P(2)	0.5	0.5	0.2119(1)	0.004(1)
Na(1)	0.5	0.5	0.5	0.013(1)
Cl(1)	0.7685(1)	0.7685(1)	0.0727(1)	0.011(1)
O(1)	0.5963(2)	0.4037(2)	0.1196(3)	0.007(1)
O(2)	0.9227(2)	0.6110(2)	0.3396(2)	0.011(1)
O(3)	0.5949(2)	0.5949(2)	0.2994(3)	0.008(1)