

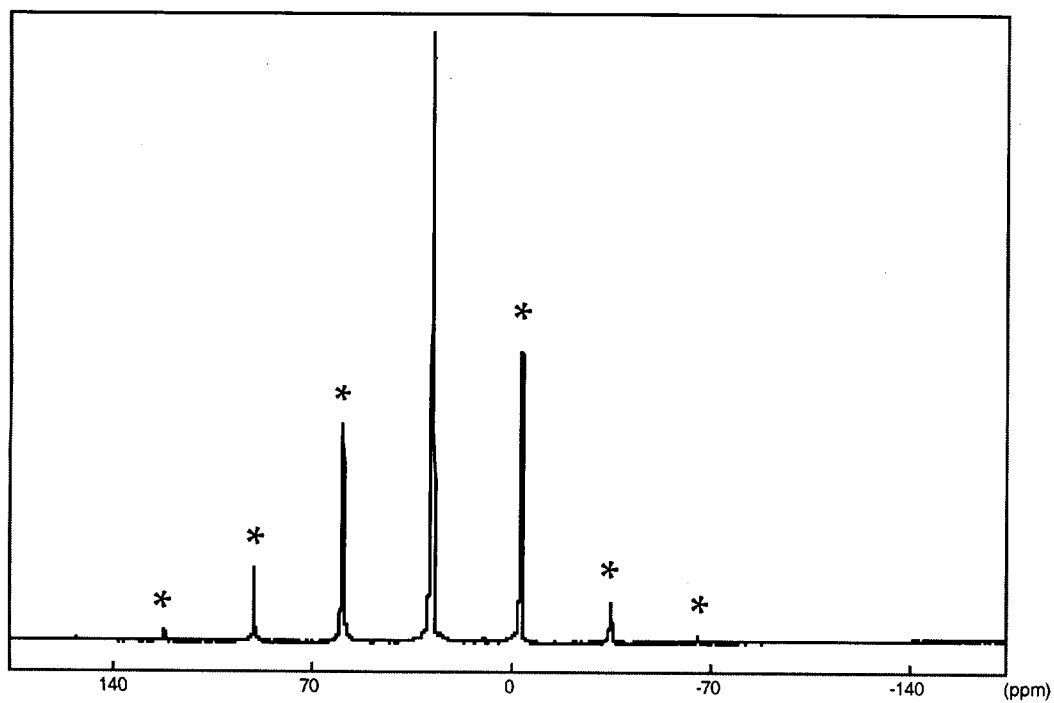
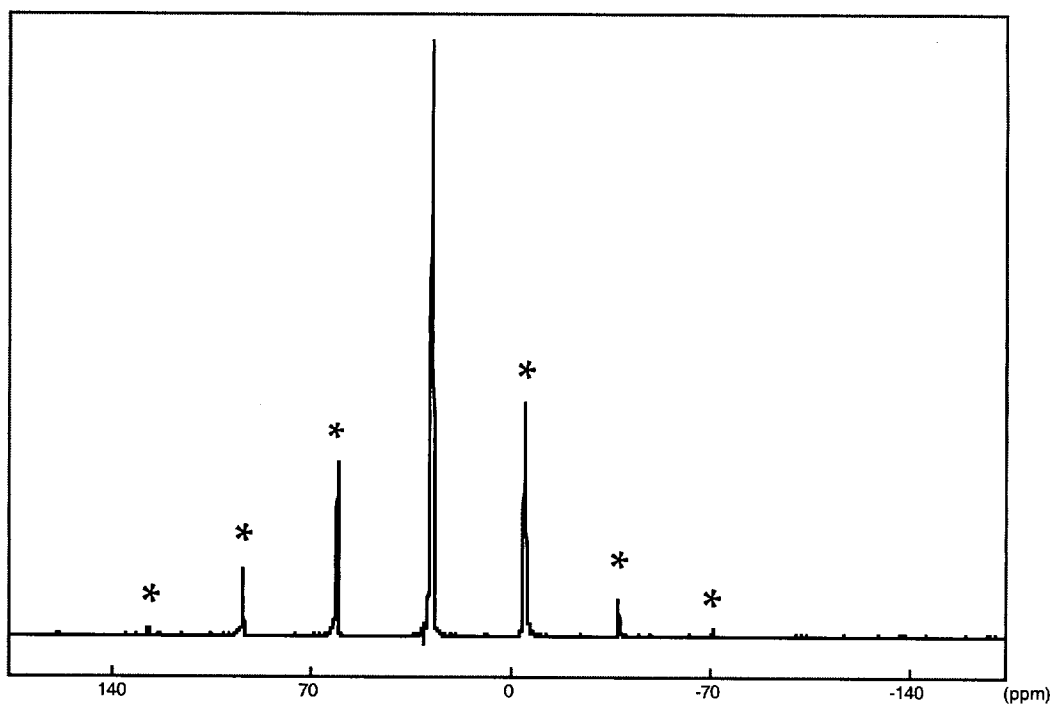


Figure S1. ^{31}P MAS NMR spectra of (a) $\text{Cd}(\text{O}_3\text{PC}_2\text{H}_4\text{NH}_2)$ and (b) $\text{Cd}(\text{O}_3\text{PCH}_3).n\text{-NH}_2\text{C}_4\text{H}_9$. Asterisks denote spinning side bands.

Table S1. Data collection and refinement parameters for Cd(O₃PC₂H₄NH₂)

Data collection	
<i>Diffractometer :</i>	Rigaku computer-automated diffractometer
Wavelengths (Å)	1.540598 ($\lambda_{K\alpha 1}$) - 1.544380 ($\lambda_{K\alpha 2}$)
2 θ angular range (°)	2-80
Step (°) ; time per step (s)	0.02 ; 20
Refinement	
<i>Diffractometer / sample :</i>	
No. of background parameters (shifted Chebyshev function)	12
K ratio	0.39(1)
Scale factor	52.0(2)
preferred orientation	1.016(4)
Sample shift	4.73(4)
<i>Crystal system :</i>	
Space group	<i>Pna2</i> ₁
Cell parameters (Å)	15.4643(2) 5.16512(7) 6.27650(8)
Volume (Å ³) ; Z	501.33(2) ; 4
M (g mol ⁻¹) ; ρ_{calcd} (g cm ⁻³)	253.46 ; 3.12
No. of reflections	368
No. of refined parameters	55
<i>Profiles :</i>	
Function	Pseudovoigt with Finger-Cox-Jephcoat asymmetry
Halfwidth parameters (centideg ²) (U ; V ; W)	1(1) ; 0 ; 5.1(2)
Strain broadening (Y, centideg)	17.7(2)
Particle size broadening (X)	0.43 (fixed)
Asymmetry parameters (S/L ; H/L)	0.0289 ; 0.0272 (fixed)
<i>Agreement factors :</i>	
R _p (%)	5.7
R _{wp} (%)	4.0
R(F ²) (%)	3.0

Table S2. Bond Lengths (Å) and Angles (deg) for the Non-Hydrogen Atoms of Cd(O₃PC₂H₄NH₂)

Cd-O(1) ^a	2.373(9)	P-O(1)	1.481(10)
Cd-O(1) ^b	2.339(8)	P-O(2)	1.559(5)
Cd-O(2) ^c	2.266(6)	P-O(3)	1.575(10)
Cd-O(3) ^a	2.516(9)	P-C(1)	1.826(5)
Cd-O(3) ^d	2.189(8)	C(1)-C(2)	1.524(6)
Cd-N	2.320(7)	C(2)-N	1.492(6)
O(1) ^a -Cd-O(1) ^b	146.8(5)	O(3) ^a -Cd-O(3) ^d	159.4(3)
O(1) ^a -Cd-O(2) ^c	88.4(5)	O(3) ^a -Cd-N	83.1(3)
O(1) ^a -Cd-O(3) ^a	60.9(2)	O(3) ^d -Cd-N ^c	98.5(3)
O(1) ^a -Cd-O(3) ^d	98.9(4)	O(1)-P-O(2)	107.2(9)
O(1) ^a -Cd-N	82.5(2)	O(1)-P-O(3)	108.4(4)
O(1) ^b -Cd-O(2) ^c	91.3(5)	O(1)-P-C(1)	105.9(5)
O(1) ^b -Cd-O(3) ^a	86.0(4)	O(2)-P-O(3)	113.9(9)
O(1) ^b -Cd-O(3) ^d	114.3(2)	O(2)-P-C(1)	106.6(4)
O(1) ^b -Cd-N	92.3(3)	O(3)-P-C(1)	105.7(5)
O(2) ^c -Cd-O(3) ^a	86.2(3)	P-C(1)-C(2)	123.1(3)
O(2) ^c -Cd-O(3) ^d	90.0(5)	C(1)-C(2)-N	114.2(3)
O(2) ^c -Cd-N	168.4(2)		

^aAtom related by $1/2 + x, 3/2 - y, z$; ^b $1 - x, 1 - y, -1/2 + z$; ^c $1/2 + x, 1/2 - y, z$; ^d $1 - x, 1 - y, 1/2 + z$

Table S3. Positional Parameters and Thermal Parameters for the Atoms of Cd(O₃PC₂H₄NH₂)

atom	<i>x</i>	<i>y</i>	<i>z</i>	<i>U</i> _{iso} , Å ²
Cd	0.73057(6)	0.3378(2)	0.3881(2)	0.0456(6)
P	0.3570(1)	0.7230(3)	0.3757(17)	0.037(1)
O(1)	0.3136(5)	0.839(2)	0.5629(12)	0.034(1)
O(2)	0.3482(4)	0.4235(6)	0.397(2)	0.034(1)
O(3)	0.3217(6)	0.859(2)	0.1687(12)	0.034(1)
N	0.6302(6)	0.6707(5)	0.3653(4)	0.039(1)
C(1)	0.4730(6)	0.8035(7)	0.3978(9)	0.058(2)
C(2)	0.5424(9)	0.6028(6)	0.4465(9)	0.058(2)