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Table S1: Crystal data and structure refinement for **14**

Empirical formula	$C_{13}H_{12}SiFO_6CoFe$
Molecular Weight	426.10
Description	parallelepiped
Dimensions, mm	0.02 x 0.04 x 0.04
Temperature, K	300(2)
Wavelength, Å	0.71073
Crystal system	Monoclinic
Space group	P21/n
Z	4
a, Å	8.2492(2)
b, Å	15.4888(2)
c, Å	13.6408(3)
β, deg	94.7430(10)
Volume, $Å^3$	1736.92(6)
Calcd Density, g/cm^3	1.629
Abs coeff, mm^{-1}	1.892
F(000)	856
θ-range for collection, deg.	1.99 to 26.58
Index ranges	$-10 \le h \le 10$, $-19 \le k \le 19$, $-16 \le l \le 16$
No. Reflections collected	14037
No. Independent reflection	3375
R(int)	0.0470
Refinement method	Full-matrix least-squares on F^2
Data / restraints / parameter	3375 / 0 / 209
Goodness-of-fit on F^2	1.059
Final R indices ($I>2\sigma(I)$)*	R1 = 0.0396, wR2 = 0.0890
R indices (all data)*	R1 = 0.0622, wR2 = 0.1016
Largest diff. Peak, $e/Å^3$	0.479
Largest diff. Hole, $e/Å^3$	-0.360

*$R_1 = \Sigma||Fo| - |Fc|| / \Sigma|Fo|$; $wR_2 = (\Sigma[w(F_o^2 - F_c^2)^2] / \Sigma[wF_o^4])^{1/2}$

Table S2: Atomic coordinates (x 10^4) and equivalent isotropic displacement parameters (A^2 x 10^3) for **14**. U(eq) is defined as one third of the trace of the orthogonalized Uij tensor.

Co(2)	800(1)	7041(1)	4073(1)	37(1)
Fe(1)	1263(1)	8169(1)	2764(1)	35(1)
Si(1)	4268(1)	6491(1)	2941(1)	39(1)
C(2)	2798(4)	7264(2)	3355(2)	34(1)
C(3)	2538(4)	7878(2)	4029(2)	34(1)
C(4)	2745(4)	8762(2)	4134(2)	38(1)
C(5)	1855(5)	9300(3)	4837(3)	55(1)
C(6)	4329(5)	9172(3)	3869(3)	51(1)
C(7)	5218(6)	5862(3)	3983(3)	67(1)
C(8)	5758(6)	7046(3)	2236(3)	64(1)
F(1)	3277(3)	5833(2)	2210(2)	74(1)
C(11)	353(5)	7499(3)	1810(3)	50(1)
O(11)	-195(4)	7067(2)	1188(2)	79(1)
C(12)	-384(5)	8925(3)	2884(3)	50(1)
O(12)	-1397(4)	9409(2)	2998(3)	78(1)
C(13)	2384(5)	8770(3)	1920(3)	48(1)
O(13)	3073(4)	9127(2)	1363(2)	79(1)
C(21)	1594(5)	6448(3)	5148(3)	55(1)
O(21)	2118(5)	6086(3)	5823(3)	92(1)
C(22)	-918(5)	7590(3)	4517(3)	54(1)
O(22)	-2025(4)	7917(3)	4797(3)	87(1)
C(23)	-148(5)	6149(3)	3346(3)	50(1)
O(23)	-718(4)	5600(2)	2902(3)	83(1)

Table S3: Selected bond lengths [Å] and angles [deg] for **14**.

Co(2)-C(22)	1.800(4)	C(22)-Co(2)-C(21)	102.1(2)
Co(2)-C(21)	1.807(5)	C(22)-Co(2)-C(23)	103.0(2)
Co(2)-C(23)	1.838(4)	C(21)-Co(2)-C(23)	99.5(2)
Co(2)-C(3)	1.937(4)	C(22)-Co(2)-C(3)	107.5(2)
Co(2)-C(2)	2.015(3)	C(21)-Co(2)-C(3)	98.4(2)
Co(2)-Fe(1)	2.5494(7)	C(23)-Co(2)-C(3)	140.3(2)
Fe(1)-C(11)	1.781(4)	C(22)-Co(2)-C(2)	141.4(2)
Fe(1)-C(13)	1.795(4)	C(21)-Co(2)-C(2)	103.0(2)
Fe(1)-C(12)	1.812(4)	C(23)-Co(2)-C(2)	101.3(2)
Fe(1)-C(3)	1.997(3)	C(3)-Co(2)-C(2)	39.95(14)
Fe(1)-C(2)	2.012(3)	C(22)-Co(2)-Fe(1)	94.59(14)
Fe(1)-C(4)	2.335(3)	C(21)-Co(2)-Fe(1)	148.40(14)
Si(1)-F(1)	1.601(3)	C(23)-Co(2)-Fe(1)	102.68(13)
Si(1)-C(2)	1.826(3)	C(3)-Co(2)-Fe(1)	50.63(10)
Si(1)-C(8)	1.837(4)	C(2)-Co(2)-Fe(1)	50.67(10)
Si(1)-C(7)	1.843(4)	C(11)-Fe(1)-C(13)	92.4(2)
C(2)-C(3)	1.352(5)	C(11)-Fe(1)-C(12)	99.7(2)
C(3)-C(4)	1.386(5)	C(13)-Fe(1)-C(12)	98.8(2)
C(4)-C(5)	1.506(5)	C(11)-Fe(1)-C(3)	131.3(2)
C(4)-C(6)	1.522(5)	C(13)-Fe(1)-C(3)	114.2(2)
C(11)-O(11)	1.144(5)	C(12)-Fe(1)-C(3)	114.4(2)
C(12)-O(12)	1.142(5)	C(11)-Fe(1)-C(2)	95.7(2)
C(13)-O(13)	1.130(5)	C(13)-Fe(1)-C(2)	106.1(2)
C(21)-O(21)	1.133(5)	C(12)-Fe(1)-C(2)	150.0(2)
C(22)-O(22)	1.138(5)	C(3)-Fe(1)-C(2)	39.42(14)
C(23)-O(23)	1.124(5)	C(11)-Fe(1)-C(4)	167.2(2)
		C(13)-Fe(1)-C(4)	92.6(2)
		C(12)-Fe(1)-C(4)	91.1(2)

C(3)-Fe(1)-C(4)	36.28(13)	C(3)-C(4)-Fe(1)	58.5(2)
C(2)-Fe(1)-C(4)	71.53(13)	C(5)-C(4)-Fe(1)	118.5(3)
C(11)-Fe(1)-Co(2)	91.93(13)	C(6)-C(4)-Fe(1)	112.2(2)
C(13)-Fe(1)-Co(2)	156.85(13)	O(11)-C(11)-Fe(1)	178.3(4)
C(12)-Fe(1)-Co(2)	102.85(13)	O(12)-C(12)-Fe(1)	177.2(4)
C(3)-Fe(1)-Co(2)	48.59(10)	O(13)-C(13)-Fe(1)	177.6(4)
C(2)-Fe(1)-Co(2)	50.78(9)	O(21)-C(21)-Co(2)	178.7(4)
C(4)-Fe(1)-Co(2)	78.93(9)	O(22)-C(22)-Co(2)	178.3(4)
F(1)-Si(1)-C(2)	106.9(2)	O(23)-C(23)-Co(2)	179.5(4)
F(1)-Si(1)-C(8)	107.5(2)		
C(2)-Si(1)-C(8)	110.2(2)		
F(1)-Si(1)-C(7)	107.9(2)		
C(2)-Si(1)-C(7)	111.0(2)		
C(8)-Si(1)-C(7)	113.0(2)		
C(3)-C(2)-Si(1)	144.9(3)		
C(3)-C(2)-Fe(1)	69.7(2)		
Si(1)-C(2)-Fe(1)	138.1(2)		
C(3)-C(2)-Co(2)	66.9(2)		
Si(1)-C(2)-Co(2)	128.9(2)		
Fe(1)-C(2)-Co(2)	78.54(12)		
C(2)-C(3)-C(4)	137.7(3)		
C(2)-C(3)-Co(2)	73.1(2)		
C(4)-C(3)-Co(2)	138.0(3)		
C(2)-C(3)-Fe(1)	70.9(2)		
C(4)-C(3)-Fe(1)	85.3(2)		
Co(2)-C(3)-Fe(1)	80.78(13)		
C(3)-C(4)-C(5)	123.3(3)		
C(3)-C(4)-C(6)	119.2(3)		
C(5)-C(4)-C(6)	113.2(3)		

Table S4: Anisotropic displacement parameters ($Å^2 \times 10^3$) for **18**. The anisotropic displacement factor exponent takes the form:

$-2p^2 [h^2 a^{*2} U11 + ... + 2 h k a^* b^* U12]$

	U11	U22	U33	U23	U13	U12
Co(2)	35(1)	37(1)	40(1)	1(1)	5(1)	0(1)
Fe(1)	34(1)	37(1)	34(1)	1(1)	-1(1)	4(1)
Si(1)	36(1)	38(1)	44(1)	-6(1)	3(1)	6(1)
C(2)	29(2)	36(2)	37(2)	2(2)	3(1)	0(1)
C(3)	33(2)	36(2)	32(2)	0(2)	-2(1)	4(1)
C(4)	40(2)	35(2)	36(2)	0(2)	-5(2)	2(2)
C(5)	63(3)	47(2)	54(2)	-13(2)	9(2)	4(2)
C(6)	47(2)	45(2)	60(3)	-6(2)	0(2)	-10(2)
C(7)	63(3)	66(3)	72(3)	12(2)	9(2)	24(2)
C(8)	54(3)	86(3)	55(3)	8(2)	19(2)	12(2)
F(1)	67(2)	63(2)	89(2)	-39(2)	-6(2)	1(1)
C(11)	45(2)	58(3)	46(2)	0(2)	-1(2)	-2(2)
O(11)	78(2)	97(3)	59(2)	-28(2)	-12(2)	-18(2)
C(12)	49(2)	53(2)	47(2)	5(2)	-2(2)	11(2)
O(12)	67(2)	85(3)	80(2)	2(2)	3(2)	41(2)
C(13)	48(2)	53(2)	43(2)	3(2)	-3(2)	0(2)
O(13)	82(2)	91(3)	66(2)	20(2)	18(2)	-16(2)
C(21)	55(3)	55(3)	56(3)	10(2)	12(2)	5(2)
O(21)	105(3)	97(3)	73(2)	38(2)	8(2)	24(2)
C(22)	45(2)	62(3)	57(3)	5(2)	10(2)	3(2)
O(22)	58(2)	107(3)	100(3)	-6(2)	32(2)	24(2)
C(23)	38(2)	45(2)	68(3)	-2(2)	6(2)	-3(2)
O(23)	66(2)	66(2)	118(3)	-30(2)	4(2)	-20(2)

Table S5: Hydrogen coordinates (x 10^4) and isotropic displacement parameters ($Å^2$ x 10^3) for **14**.

	x	y	z	U(eq)
H(5A)	2180(5)	9892(3)	4790(3)	82
H(5B)	2114(5)	9097(3)	5496(3)	82
H(5C)	705(5)	9252(3)	4672(3)	82
H(6A)	4291(5)	9783(3)	3980(3)	76
H(6B)	4474(5)	9063(3)	3189(3)	76
H(6C)	5223(5)	8927(3)	4272(3)	76
H(7A)	5984(6)	5462(3)	3747(3)	100
H(7B)	4393(6)	5552(3)	4293(3)	100
H(7C)	5771(6)	6247(3)	4450(3)	100
H(8A)	6524(6)	6634(3)	2022(3)	96
H(8B)	6324(6)	7473(3)	2643(3)	96
H(8C)	5206(6)	7321(3)	1672(3)	96