

Table S1. Atomic coordinates ($\times 10^4$) and equivalent isotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 19. U(eq) is defined as one third of the trace of the orthogonalized U_{ij} tensor.

	X	Y	Z	U (eq)
Si (1)	4431 (1)	746 (1)	1941 (1)	54 (1)
N (21)	3947 (3)	1425 (2)	5710 (2)	85 (1)
N (22)	-317 (4)	2062 (2)	4122 (2)	88 (1)
N (31)	3369 (3)	-950 (2)	5127 (2)	82 (1)
N (32)	-933 (3)	-422 (2)	3713 (2)	79 (1)
C (1)	2536 (3)	1490 (2)	3126 (2)	46 (1)
C (2)	1979 (3)	1147 (2)	4054 (2)	46 (1)
C (3)	1705 (3)	99 (2)	3821 (2)	43 (1)
C (4)	2086 (3)	-11 (2)	2770 (2)	42 (1)
C (5)	922 (3)	524 (2)	2063 (2)	44 (1)
C (6)	1209 (3)	1426 (2)	2266 (2)	46 (1)
C (7)	3385 (3)	651 (2)	2930 (2)	41 (1)
C (8)	324 (4)	2090 (3)	1740 (3)	64 (1)
C (9)	-865 (4)	1820 (3)	990 (3)	77 (1)
C (10)	-1149 (4)	922 (3)	799 (3)	76 (1)
C (11)	-274 (4)	248 (3)	1330 (2)	62 (1)
C (12)	3259 (3)	2411 (2)	3309 (2)	68 (1)
C (13)	2242 (3)	-989 (2)	2516 (2)	62 (1)
C (14)	3256 (3)	1116 (2)	730 (2)	76 (1)
C (15)	5242 (4)	-365 (2)	1797 (3)	90 (1)
C (16)	5918 (3)	1579 (3)	2348 (3)	88 (1)
C (21)	3093 (4)	1296 (2)	4994 (3)	58 (1)
C (22)	675 (4)	1653 (3)	4117 (2)	58 (1)
C (31)	2637 (3)	-480 (2)	4575 (3)	53 (1)
C (32)	205 (4)	-177 (2)	3764 (2)	52 (1)

Table S2. Complete Listing of Bond lengths [Å] and angles [deg] for 19.

Si (1) -C (15)	1.850 (3)	C (32) -C (3) -C (2)	112.8 (3)
Si (1) -C (14)	1.857 (3)	C (31) -C (3) -C (4)	110.8 (3)
Si (1) -C (16)	1.859 (3)	C (32) -C (3) -C (4)	111.8 (2)
Si (1) -C (7)	1.913 (3)	C (2) -C (3) -C (4)	102.7 (2)
N (21) -C (21)	1.133 (4)	C (13) -C (4) -C (5)	116.8 (3)
N (22) -C (22)	1.130 (4)	C (13) -C (4) -C (7)	120.9 (3)
N (31) -C (31)	1.136 (4)	C (5) -C (4) -C (7)	101.2 (2)
N (32) -C (32)	1.135 (4)	C (13) -C (4) -C (3)	111.8 (3)
C (1) -C (6)	1.511 (4)	C (5) -C (4) -C (3)	103.8 (2)
C (1) -C (12)	1.519 (4)	C (7) -C (4) -C (3)	99.7 (2)
C (1) -C (7)	1.547 (4)	C (6) -C (5) -C (11)	121.4 (3)
C (1) -C (2)	1.608 (4)	C (6) -C (5) -C (4)	107.3 (3)
C (2) -C (21)	1.479 (4)	C (11) -C (5) -C (4)	131.3 (3)
C (2) -C (22)	1.481 (5)	C (5) -C (6) -C (8)	121.1 (3)
C (2) -C (3)	1.591 (4)	C (5) -C (6) -C (1)	107.9 (3)
C (3) -C (31)	1.468 (4)	C (8) -C (6) -C (1)	131.0 (3)
C (3) -C (32)	1.479 (4)	C (1) -C (7) -C (4)	95.2 (2)
C (3) -C (4)	1.612 (4)	C (1) -C (7) -Si (1)	117.9 (2)
C (4) -C (13)	1.505 (4)	C (4) -C (7) -Si (1)	119.6 (2)
C (4) -C (5)	1.509 (4)	C (6) -C (8) -C (9)	118.0 (3)
C (4) -C (7)	1.554 (4)	C (10) -C (9) -C (8)	120.5 (4)
C (5) -C (6)	1.377 (4)	C (9) -C (10) -C (11)	122.1 (4)
C (5) -C (11)	1.387 (4)	C (5) -C (11) -C (10)	117.0 (3)
C (6) -C (8)	1.379 (4)	N (21) -C (21) -C (2)	178.9 (4)
C (8) -C (9)	1.393 (4)	N (22) -C (22) -C (2)	176.4 (4)
C (9) -C (10)	1.368 (5)	N (31) -C (31) -C (3)	177.1 (4)
C (10) -C (11)	1.388 (5)	N (32) -C (32) -C (3)	177.4 (4)
C (15) -Si (1) -C (14)	109.63 (17)		
C (15) -Si (1) -C (16)	108.03 (17)		
C (14) -Si (1) -C (16)	108.70 (17)		
C (15) -Si (1) -C (7)	109.15 (15)		
C (14) -Si (1) -C (7)	111.80 (14)		
C (16) -Si (1) -C (7)	109.46 (15)		
C (6) -C (1) -C (12)	116.3 (3)		
C (6) -C (1) -C (7)	100.5 (2)		
C (12) -C (1) -C (7)	120.7 (3)		
C (6) -C (1) -C (2)	103.5 (2)		
C (12) -C (1) -C (2)	112.3 (3)		
C (7) -C (1) -C (2)	101.0 (2)		
C (21) -C (2) -C (22)	107.5 (3)		
C (21) -C (2) -C (3)	111.7 (3)		
C (22) -C (2) -C (3)	114.0 (3)		
C (21) -C (2) -C (1)	110.7 (3)		
C (22) -C (2) -C (1)	110.6 (3)		
C (3) -C (2) -C (1)	102.3 (2)		
C (31) -C (3) -C (32)	106.3 (3)		
C (31) -C (3) -C (2)	112.5 (3)		

Table S3. Anisotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 19. The anisotropic displacement factor exponent takes the form: $-2\pi^2 [h^2 a^{*2} U11 + \dots + 2 h k a^* b^* U12]$

	U11	U22	U33	U23	U13	U12
Si (1)	62 (1)	58 (1)	51 (1)	5 (1)	28 (1)	2 (1)
N (21)	82 (2)	105 (3)	65 (2)	-18 (2)	17 (2)	-21 (2)
N (22)	92 (2)	94 (3)	86 (2)	-1• (2)	39 (2)	27 (2)
N (31)	78 (2)	96 (3)	75 (2)	28 (2)	24 (2)	15 (2)
N (32)	60 (2)	86 (3)	97 (3)	4 (2)	28 (2)	-15 (2)
C (1)	56 (2)	38 (2)	49 (2)	-5 (2)	23 (2)	-7 (2)
C (2)	52 (2)	46 (2)	43 (2)	-10 (2)	19 (2)	-4 (2)
C (3)	45 (2)	45 (2)	42 (2)	5 (2)	18 (2)	-3 (2)
C (4)	54 (2)	32 (2)	44 (2)	-4 (2)	20 (2)	-1 (2)
C (5)	50 (2)	46 (3)	36 (2)	1 (2)	14 (2)	0 (2)
C (6)	56 (2)	45 (3)	43 (2)	5 (2)	24 (2)	8 (2)
C (7)	44 (2)	40 (2)	41 (2)	2 (2)	16 (1)	-5 (2)
C (8)	84 (3)	52 (3)	66 (3)	9 (2)	36 (2)	16 (2)
C (9)	78 (3)	96 (4)	61 (3)	31 (3)	27 (2)	38 (3)
C (10)	67 (3)	104 (4)	54 (3)	11 (3)	8 (2)	1• (3)
C (11)	64 (2)	64 (3)	57 (2)	-1 (2)	15 (2)	-8 (2)
C (12)	88 (2)	45 (3)	80 (3)	-1• (2)	37 (2)	-17 (2)
C (13)	78 (2)	49 (3)	67 (2)	-7 (2)	31 (2)	-7 (2)
C (14)	94 (3)	89 (3)	51 (2)	14 (2)	31 (2)	7 (2)
C (15)	102 (3)	86 (3)	104 (3)	16 (3)	67 (2)	35 (3)
C (16)	76 (2)	113 (4)	87 (3)	4 (3)	45 (2)	-24 (3)
C (21)	62 (2)	64 (3)	54 (2)	-13 (2)	26 (2)	-8 (2)
C (22)	72 (3)	61 (3)	47 (2)	-11 (2)	24 (2)	2 (2)
C (31)	52 (2)	60 (3)	52 (2)	7 (2)	22 (2)	-2 (2)
C (32)	58 (2)	52 (3)	50 (2)	6 (2)	20 (2)	-2 (2)

Table S4. Hydrogen coordinates ($\times 10^4$) and isotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 19.

	x	y	z	U (eq)
H (7A)	4073	487	3556	49
H (8A)	514	2700	1882	77
H (9A)	-1471	2254	615	92
H (10A)	-1953	758	298	92
H (11A)	-480	-362	1200	74
H (12A)	3565	2591	2735	102
H (12B)	4081	2377	3871	102
H (12C)	2587	2847	3435	102
H (13A)	2467	-1028	1887	93
H (13B)	1354	-1302	2482	93
H (13C)	3004	-1260	3018	93
H (14A)	3816	1158	254	114
H (14B)	2850	1697	804	114
H (14C)	2493	685	506	114
H (15A)	5760	-322	1296	135
H (15B)	4493	-810	1606	135
H (15C)	5893	-538	2415	135
H (16A)	6436	1632	1848	131
H (16B)	6564	1377	2956	131
H (16C)	5524	2157	2450	131

Table S5. Atomic coordinates ($\times 10^4$) and equivalent isotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 22.
U(eq) is defined as one third of the trace of the orthogonalized U_{ij} tensor.

	x	y	z	U(eq)
Cr(1)	2689(1)	4432(1)	8836(1)	58(1)
Si(1)	2108(2)	888(2)	6507(2)	60(1)
O(55)	2048(6)	5998(6)	10846(6)	104(2)
O(44)	927(7)	6837(6)	6920(6)	117(2)
O(66)	5169(7)	6198(6)	7600(6)	96(2)
C(1)	2675(8)	2744(6)	6404(7)	51(2)
C(2)	4213(9)	2683(7)	6089(8)	69(2)
C(3)	4909(8)	2528(7)	7168(10)	66(2)
C(3A)	3869(8)	2525(6)	8322(7)	49(2)
C(4)	4004(10)	2364(7)	9715(10)	76(3)
C(5)	2770(16)	2365(8)	10595(9)	100(4)
C(6)	1443(12)	2584(9)	10143(11)	92(3)
C(7A)	2528(7)	2733(6)	7860(7)	43(2)
C(7)	1327(9)	2794(7)	8768(9)	69(2)
C(8)	191(8)	852(9)	6893(10)	114(3)
C(9)	2639(10)	686(9)	4802(9)	116(3)
C(10)	2967(9)	-590(7)	7890(9)	101(3)
C(11)	1844(9)	4014(7)	5364(8)	99(3)
C(33)	6464(8)	2313(9)	7266(10)	113(3)
C(44)	1615(9)	5893(8)	7633(8)	76(2)
C(55)	2292(7)	5389(8)	10060(8)	66(2)
C(66)	4201(9)	5504(8)	8077(8)	64(2)

Table S6. Complete Listing of Bond lengths [Å] and angles [deg] for 22.

Cr(1)-C(66)	1.822(9)	C(6)-Cr(1)-C(7A)	65.0(3)
Cr(1)-C(55)	1.826(9)	C(4)-Cr(1)-C(7A)	65.8(3)
Cr(1)-C(44)	1.836(8)	C(7)-Cr(1)-C(7A)	36.0(2)
Cr(1)-C(5)	2.172(8)	C(3A)-Cr(1)-C(7A)	36.0(2)
Cr(1)-C(6)	2.205(9)	C(10)-Si(1)-C(8)	109.9(4)
Cr(1)-C(4)	2.217(6)	C(10)-Si(1)-C(9)	110.4(4)
Cr(1)-C(7)	2.237(8)	C(8)-Si(1)-C(9)	109.3(4)
Cr(1)-C(3A)	2.280(7)	C(10)-Si(1)-C(1)	108.7(3)
Cr(1)-C(7A)	2.293(6)	C(8)-Si(1)-C(1)	109.8(3)
Si(1)-C(10)	1.831(7)	C(9)-Si(1)-C(1)	108.6(4)
Si(1)-C(8)	1.847(8)	C(2)-C(1)-C(7A)	100.9(5)
Si(1)-C(9)	1.853(9)	C(2)-C(1)-C(11)	116.1(7)
Si(1)-C(1)	1.938(7)	C(7A)-C(1)-C(11)	115.4(6)
O(55)-C(55)	1.168(8)	C(2)-C(1)-Si(1)	106.3(4)
O(44)-C(44)	1.154(7)	C(7A)-C(1)-Si(1)	107.4(4)
O(66)-C(66)	1.170(8)	C(11)-C(1)-Si(1)	109.8(5)
C(1)-C(2)	1.479(9)	C(3)-C(2)-C(1)	114.2(7)
C(1)-C(7A)	1.496(8)	C(2)-C(3)-C(3A)	107.0(7)
C(1)-C(11)	1.540(8)	C(2)-C(3)-C(33)	128.9(8)
C(2)-C(3)	1.339(10)	C(3A)-C(3)-C(33)	124.0(8)
C(3)-C(3A)	1.455(10)	C(7A)-C(3A)-C(4)	119.9(7)
C(3)-C(33)	1.508(10)	C(7A)-C(3A)-C(3)	108.5(6)
C(3A)-C(7A)	1.413(8)	C(4)-C(3A)-C(3)	131.6(8)
C(3A)-C(4)	1.419(10)	C(7A)-C(3A)-Cr(1)	72.5(4)
C(4)-C(5)	1.400(11)	C(4)-C(3A)-Cr(1)	69.2(4)
C(5)-C(6)	1.395(12)	C(3)-C(3A)-Cr(1)	130.1(4)
C(6)-C(7)	1.384(11)	C(5)-C(4)-C(3A)	117.3(8)
C(7A)-C(7)	1.403(9)	C(5)-C(4)-Cr(1)	69.7(4)
C(66)-Cr(1)-C(55)	88.6(3)	C(3A)-C(4)-Cr(1)	74.0(4)
C(66)-Cr(1)-C(44)	89.2(3)	C(4)-C(5)-C(6)	123.1(9)
C(55)-Cr(1)-C(44)	88.1(3)	C(4)-C(5)-Cr(1)	73.2(4)
C(66)-Cr(1)-C(5)	122.9(5)	C(6)-C(5)-Cr(1)	72.7(5)
C(55)-Cr(1)-C(5)	89.1(3)	C(7)-C(6)-C(5)	119.0(9)
C(44)-Cr(1)-C(5)	147.7(5)	C(7)-C(6)-Cr(1)	73.1(5)
C(66)-Cr(1)-C(6)	159.7(4)	C(5)-C(6)-Cr(1)	70.1(5)
C(55)-Cr(1)-C(6)	93.7(3)	C(7)-C(7A)-C(3A)	120.4(6)
C(44)-Cr(1)-C(6)	111.0(4)	C(7)-C(7A)-C(1)	130.4(7)
C(5)-Cr(1)-C(6)	37.2(3)	C(3A)-C(7A)-C(1)	109.1(6)
C(66)-Cr(1)-C(4)	92.9(3)	C(7)-C(7A)-Cr(1)	69.8(4)
C(55)-Cr(1)-C(4)	111.6(3)	C(3A)-C(7A)-Cr(1)	71.5(4)
C(44)-Cr(1)-C(4)	160.3(3)	C(1)-C(7A)-Cr(1)	135.0(4)
C(5)-Cr(1)-C(4)	37.2(3)	C(6)-C(7)-C(7A)	120.2(8)
C(6)-Cr(1)-C(4)	67.5(3)	C(6)-C(7)-Cr(1)	70.6(5)
C(66)-Cr(1)-C(7)	148.4(3)	C(7A)-C(7)-Cr(1)	74.1(4)
C(55)-Cr(1)-C(7)	122.8(3)	O(44)-C(44)-Cr(1)	177.1(7)
C(44)-Cr(1)-C(7)	89.1(3)	O(55)-C(55)-Cr(1)	179.5(6)
C(5)-Cr(1)-C(7)	65.8(4)	O(66)-C(66)-Cr(1)	179.3(7)
C(6)-Cr(1)-C(7)	36.3(3)	C(66)-Cr(1)-C(7A)	112.9(3)
C(4)-Cr(1)-C(7)	78.9(3)	C(55)-Cr(1)-C(7A)	158.2(3)
C(66)-Cr(1)-C(3A)	89.7(3)	C(44)-Cr(1)-C(7A)	95.3(3)
C(55)-Cr(1)-C(3A)	148.1(3)	C(5)-Cr(1)-C(7A)	76.5(3)
C(44)-Cr(1)-C(3A)	123.7(3)		
C(5)-Cr(1)-C(3A)	65.4(3)		
C(6)-Cr(1)-C(3A)	77.7(3)		
C(4)-Cr(1)-C(3A)	36.8(3)		
C(7)-Cr(1)-C(3A)	65.5(3)		

Table S7. Anisotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 22. The anisotropic displacement factor exponent takes the form: $-2\pi^2 [h^2 a^{*2} U_{11} + \dots + 2 h k a^* b^* U_{12}]$

	U11	U22	U33	U23	U13	U12
Cr (1)	72 (1)	53 (1)	51 (1)	-21 (1)	-13 (1)	6 (1)
Si (1)	63 (2)	56 (1)	67 (2)	-25 (1)	-24 (1)	3 (1)
O (55)	107 (5)	130 (5)	113 (5)	-92 (4)	-14 (4)	7 (4)
O (44)	160 (6)	74 (4)	125 (5)	-37 (4)	-80 (5)	41 (4)
O (66)	98 (5)	99 (4)	84 (4)	-24 (4)	2 (4)	-28 (4)
C (1)	60 (6)	51 (4)	41 (5)	-18 (3)	-22 (4)	14 (4)
C (2)	81 (7)	76 (5)	58 (5)	-31 (4)	10 (5)	-29 (5)
C (3)	52 (6)	69 (5)	95 (7)	-51 (5)	-4 (5)	-2 (4)
C (3A)	60 (6)	38 (4)	53 (5)	-17 (3)	-27 (4)	4 (3)
C (4)	103 (8)	49 (4)	85 (7)	-26 (5)	-59 (6)	21 (5)
C (5)	190 (13)	58 (5)	39 (5)	-9 (4)	11 (8)	-11 (7)
C (6)	96 (8)	78 (6)	95 (9)	-35 (6)	31 (7)	-17 (6)
C (7A)	46 (5)	41 (3)	43 (5)	-22 (3)	-2 (4)	5 (3)
C (7)	87 (7)	64 (5)	65 (6)	-33 (4)	-10 (5)	0 (4)
C (8)	72 (7)	124 (7)	162 (9)	-67 (7)	-9 (6)	-22 (6)
C (9)	144 (10)	131 (8)	112 (8)	-84 (7)	-11 (7)	-30 (7)
C (10)	141 (9)	46 (4)	125 (7)	-23 (5)	-76 (6)	9 (5)
C (11)	157 (9)	60 (5)	84 (6)	-21 (4)	-60 (6)	18 (5)
C (33)	57 (7)	137 (8)	182 (10)	-101 (8)	-10 (6)	-7 (6)
C (44)	100 (7)	59 (5)	76 (6)	-31 (4)	-32 (5)	10 (5)
C (55)	55 (6)	72 (5)	80 (6)	-36 (5)	-18 (5)	2 (4)
C (66)	72 (7)	55 (5)	63 (5)	-21 (4)	-7 (5)	4 (4)

Table S8. Hydrogen coordinates ($\times 10^4$) and isotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 22.

	x	y	z	U(eq)
H(2A)	4657(9)	2748(7)	5218(8)	83
H(4A)	4918(10)	2361(7)	10018(10)	92
H(5A)	2844(16)	2391(8)	11516(9)	120
H(6A)	616(12)	2741(9)	10747(11)	110
H(7A)	419(9)	3129(7)	8394(9)	83
H(8A)	-253(8)	1629(9)	6164(10)	171
H(8B)	-89(8)	979(9)	7766(10)	171
H(8C)	-83(8)	-72(9)	6949(10)	171
H(9A)	2180(10)	1467(9)	4088(9)	174
H(9B)	2378(10)	-237(9)	4843(9)	174
H(9C)	3633(10)	725(9)	4591(9)	174
H(10A)	3962(9)	-564(7)	7683(9)	152
H(10B)	2697(9)	-1516(7)	7948(9)	152
H(10C)	2691(9)	-464(7)	8765(9)	152
H(11A)	862(9)	3952(7)	5661(8)	149
H(11B)	2031(9)	3954(7)	4455(8)	149
H(11C)	2119(9)	4930(7)	5327(8)	149
H(33A)	6645(8)	2235(9)	8177(10)	170
H(33B)	6906(8)	3134(9)	6574(10)	170
H(33C)	6830(8)	1433(9)	7111(10)	170