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Table 1. Atomic coordinates [x 10^4] and equivalent isotropic displacement parameters [$Å^2$ x 10^3] for 2. U(eq) is defined as one third of the trace of the orthogonalized U_{ij} tensor.

	x	y	z	U(eq)
Ge	-4702(1)	7283(1)	514(1)	13(1)
C(1)	-5420(2)	6093(2)	558(2)	17(1)
N	-6011(2)	6318(2)	1045(1)	14(1)
C(2)	-6377(2)	5251(2)	1265(2)	19(1)
C(3)	-6969(2)	5538(3)	1743(2)	26(1)
O(15)	-7471(1)	6284(2)	1448(1)	29(1)
C(4)	-7132(2)	7318(3)	1270(2)	28(1)
C(5)	-6546(2)	7127(3)	766(2)	21(1)
O(1)	-5073(1)	8085(2)	1305(1)	15(1)
C(11)	-4649(2)	8033(2)	1819(2)	16(1)
O(5)	-4790(1)	8424(2)	2375(1)	26(1)
C(12)	-3926(2)	7505(2)	1680(1)	13(1)
O(2)	-3922(1)	7034(2)	1021(1)	14(1)
O(3)	-4236(1)	6686(2)	-246(1)	15(1)
C(21)	-4180(2)	7429(2)	-739(1)	15(1)
O(6)	-3881(1)	7247(2)	-1267(1)	22(1)
C(22)	-4531(2)	8572(2)	-599(2)	17(1)
O(4)	-4820(1)	8573(2)	68(1)	18(1)
O(7)	-2217(2)	8011(2)	2106(1)	39(1)
O(8)	-2415(2)	7748(2)	1039(1)	31(1)
O(9)	-4911(2)	5823(2)	1975(1)	25(1)
O(10)	-4111(2)	4768(2)	2519(1)	27(1)
O(11)	-5706(2)	7027(2)	-938(1)	34(1)
O(12)	-6072(2)	8192(3)	-1748(1)	50(1)
O(14)	-3018(2)	10332(2)	-70(1)	42(1)
C(13)	-3380(2)	8478(3)	1724(2)	19(1)
O(13)	-3179(1)	8480(2)	29(1)	26(1)
C(14)	-2633(2)	8075(3)	1640(2)	22(1)
C(15)	-3753(2)	6584(2)	2196(1)	16(1)
C(16)	-4322(2)	5693(2)	2209(2)	18(1)
C(23)	-5117(2)	8816(3)	-1113(2)	24(1)
C(24)	-5651(2)	7908(3)	-1236(2)	27(1)
C(25)	-3981(2)	9526(2)	-656(2)	20(1)
C(26)	-3362(2)	9385(3)	-200(2)	22(1)
O(1W)	-1839(2)	10023(2)	668(2)	49(1)

Table 2. Bond lengths [Å] and angles [°] for 2.

Ge-O(4)	1.781(2)	Ge-O(2)	1.803(2)
Ge-O(3)	1.884(2)	Ge-C(1)	1.955(3)
Ge-O(1)	1.964(2)	C(1)-N	1.500(4)
N-H	0.89(4)	N-C(5)	1.497(4)
N-C(2)	1.505(4)	C(2)-C(3)	1.503(5)
C(3)-O(15)	1.420(4)	O(15)-C(4)	1.426(4)
C(4)-C(5)	1.508(5)	O(1)-C(11)	1.299(4)
C(11)-O(5)	1.227(4)	C(11)-C(12)	1.521(5)
C(12)-O(2)	1.425(3)	C(12)-C(15)	1.532(4)
C(12)-C(13)	1.547(5)	O(3)-C(21)	1.321(3)
C(21)-O(6)	1.211(4)	C(21)-C(22)	1.534(4)
C(22)-O(4)	1.432(4)	C(22)-C(23)	1.530(5)
C(22)-C(25)	1.538(5)	O(7)-C(14)	1.215(4)
O(8)-H(8)	0.86(4)	O(8)-C(14)	1.321(4)
O(9)-C(16)	1.210(4)	O(10)-H(10)	0.84(4)
O(10)-C(16)	1.320(4)	O(11)-C(24)	1.205(4)
O(12)-H(12)	0.90(5)	O(12)-C(24)	1.334(4)
O(14)-H(14)	0.94(5)	O(14)-C(26)	1.321(4)
C(13)-C(14)	1.493(5)	O(13)-C(26)	1.215(4)
C(15)-C(16)	1.505(5)	C(23)-C(24)	1.493(5)
C(25)-C(26)	1.485(5)	O(1W)-H(2W)	0.78(3)
O(1W)-H(1W)	0.79(3)		

O(4)-Ge-O(2)	121.31(10)	O(4)-Ge-O(3)	88.92(9)
O(2)-Ge-O(3)	90.47(9)	O(4)-Ge-C(1)	123.76(12)
O(2)-Ge-C(1)	114.75(12)	O(3)-Ge-C(1)	94.87(11)
O(4)-Ge-O(1)	86.47(8)	O(2)-Ge-O(1)	85.54(9)
O(3)-Ge-O(1)	171.19(9)	C(1)-Ge-O(1)	93.93(11)
N-C(1)-Ge	114.4(2)	H-N-C(5)	113(2)
H-N-C(1)	103(2)	C(5)-N-C(1)	111.9(2)
H-N-C(2)	107(2)	C(5)-N-C(2)	109.8(3)
C(1)-N-C(2)	112.2(2)	C(3)-C(2)-N	109.4(3)
O(15)-C(3)-C(2)	111.9(3)	C(3)-O(15)-C(4)	109.9(3)
O(15)-C(4)-C(5)	111.3(3)	N-C(5)-C(4)	109.9(3)
C(11)-O(1)-Ge	112.8(2)	O(5)-C(11)-O(1)	123.9(3)
O(5)-C(11)-C(12)	120.8(3)	O(1)-C(11)-C(12)	115.2(3)
O(2)-C(12)-C(11)	109.4(3)	O(2)-C(12)-C(15)	109.5(2)
C(11)-C(12)-C(15)	111.2(3)	O(2)-C(12)-C(13)	109.9(2)
C(11)-C(12)-C(13)	106.1(2)	C(15)-C(12)-C(13)	110.7(3)
C(12)-O(2)-Ge	116.4(2)	C(21)-O(3)-Ge	112.4(2)
O(6)-C(21)-O(3)	124.2(3)	O(6)-C(21)-C(22)	121.0(3)
O(3)-C(21)-C(22)	114.8(3)	O(4)-C(22)-C(23)	110.2(3)
O(4)-C(22)-C(21)	109.4(2)	C(23)-C(22)-C(21)	110.9(3)
O(4)-C(22)-C(25)	108.9(2)	C(23)-C(22)-C(25)	107.2(2)
C(21)-C(22)-C(25)	110.2(3)	C(22)-O(4)-Ge	114.4(2)
H(8)-O(8)-C(14)	108(3)	H(10)-O(10)-C(16)	106(3)
H(12)-O(12)-C(24)	114(3)	H(14)-O(14)-C(26)	108(3)
C(14)-C(13)-C(12)	112.3(3)	O(7)-C(14)-O(8)	118.1(4)
O(7)-C(14)-C(13)	122.8(3)	O(8)-C(14)-C(13)	119.1(3)
C(16)-C(15)-C(12)	111.1(3)	O(9)-C(16)-O(10)	124.1(3)
O(9)-C(16)-C(15)	123.7(3)	O(10)-C(16)-C(15)	112.1(4)
C(24)-C(23)-C(22)	117.2(3)	O(11)-C(24)-O(12)	122.9(4)
O(11)-C(24)-C(23)	127.1(3)	O(12)-C(24)-C(23)	110.0(3)
C(26)-C(25)-C(22)	113.5(3)	O(13)-C(26)-O(14)	122.6(3)
O(13)-C(26)-C(25)	123.4(3)	O(14)-C(26)-C(25)	114.0(3)
H(2W)-O(1W)-H(1W)	103(6)		

Table 3. Anisotropic displacement parameters [$Å^2 \times 10^3$] for 2.

The anisotropic displacement factor exponent takes the form:

$-2\pi^2 [(ha^*)^2 U_{11} + ... + 2hka^* b^* U_{12}]$

	U11	U22	U33	U23	U13	U12
Ge	14(1)	12(1)	12(1)	0(1)	1(1)	1(1)
C(1)	16(3)	17(2)	19(2)	-2(1)	3(2)	0(1)
N	13(2)	15(1)	16(1)	2(1)	-1(1)	0(1)
C(2)	18(3)	13(1)	27(2)	5(1)	-1(2)	-4(1)
C(3)	24(3)	31(2)	24(2)	7(1)	-1(2)	-4(2)
O(15)	16(2)	33(1)	36(1)	3(1)	4(1)	-2(1)
C(4)	24(3)	25(2)	35(2)	-3(1)	0(2)	4(2)
C(5)	19(3)	18(2)	25(2)	6(1)	-2(2)	2(2)
O(1)	14(2)	15(1)	16(1)	-1(1)	0(1)	4(1)
C(11)	18(3)	12(1)	19(2)	-1(1)	-1(2)	-3(1)
O(5)	27(2)	32(1)	18(1)	-10(1)	0(1)	9(1)
C(12)	11(3)	15(1)	12(1)	-2(1)	1(2)	-1(1)
O(2)	13(2)	16(1)	12(1)	-1(1)	1(1)	3(1)
O(3)	19(2)	12(1)	14(1)	0(1)	3(1)	1(1)
C(21)	11(3)	18(2)	15(2)	-1(1)	-2(2)	-3(2)
O(6)	28(2)	21(1)	16(1)	-1(1)	7(1)	1(1)
C(22)	17(3)	19(2)	14(1)	3(1)	2(2)	2(2)
O(4)	22(2)	15(1)	15(1)	2(1)	5(1)	5(1)
O(7)	19(2)	71(2)	27(1)	4(1)	-6(1)	1(2)
O(8)	15(2)	54(2)	24(1)	-4(1)	-2(1)	3(1)
O(9)	16(2)	28(1)	30(1)	9(1)	-6(1)	-4(1)
O(10)	19(2)	23(1)	39(1)	14(1)	-2(1)	-2(1)
O(11)	31(2)	39(1)	33(1)	10(1)	-10(1)	-6(1)
O(12)	27(2)	80(2)	42(2)	36(2)	-19(2)	-19(2)
O(14)	44(2)	24(1)	57(2)	-2(1)	-16(2)	-11(1)
C(13)	13(3)	20(2)	24(2)	-3(1)	-2(2)	-3(2)
O(13)	28(2)	26(1)	23(1)	9(1)	-7(1)	-5(1)
C(14)	18(3)	25(2)	22(2)	5(1)	-3(2)	-7(2)
C(15)	15(3)	20(2)	13(1)	0(1)	1(2)	2(1)
C(16)	23(3)	18(2)	12(1)	0(1)	4(2)	2(2)
C(23)	21(3)	30(2)	20(2)	9(1)	1(2)	6(2)
C(24)	16(3)	45(2)	20(2)	9(2)	-1(2)	2(2)
C(25)	24(3)	17(2)	19(2)	2(1)	4(2)	1(2)
C(26)	26(3)	21(2)	18(2)	-3(1)	7(2)	-6(2)
O(1W)	47(3)	38(2)	63(2)	-4(2)	-16(2)	-21(2)

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

	x	y	z	U(eq)
H(1A)	-5626	5989	104	21
H(1B)	-5184	5378	685	21
H	-5780(2)	6590(3)	1403(17)	24(10)
H(2A)	-6571	4852	868	23
H(2B)	-6031	4747	1490	23
H(3A)	-7214	4835	1881	32
H(3B)	-6767	5892	2152	32
H(4A)	-6932	7673	1679	34
H(4B)	-7489	7841	1078	34
H(5A)	-6749	6819	344	25
H(5B)	-6312	7853	659	25
H(8)	-2740(2)	7910(3)	750(2)	37
H(10)	-4470(2)	4350(3)	2545(19)	32
H(12)	-6450(3)	7730(4)	-1800(2)	60
H(14)	-2600(3)	10150(4)	170(2)	50
H(13A)	-3487	9041	1370	23
H(13B)	-3427	8857	2165	23
H(15A)	-3291	6232	2081	19
H(15B)	-3709	6927	2648	19
H(23A)	-5375	9500	-965	28
H(23B)	-4886	8998	-1548	28
H(25A)	-3807	9561	-1126	24
H(25B)	-4216	10253	-555	24
H(1W)	-1660(3)	10590(3)	770(3)	74
H(2W)	-1590(3)	9570(4)	830(3)	74

Table 1. Atomic coordinates [x 10^4] and equivalent isotropic displacement parameters [$Å^2$ x 10^3] for 3. U(eq) is defined as one third of the trace of the orthogonalized U_{ij} tensor.

	x	y	z	U(eq)
Si	340(1)	7639(1)	4467(1)	20(1)
O(3)	760(1)	8248(1)	5180(1)	22(1)
N	-969(1)	8637(2)	3967(1)	21(1)
C(1)	-355(2)	8787(2)	4450(1)	24(1)
O(4)	207(1)	6457(1)	4901(1)	23(1)
C(2)	-1319(2)	9725(2)	3778(1)	28(1)
O(1)	-38(1)	6914(1)	3732(1)	21(1)
C(3)	-1929(2)	9501(2)	3300(1)	33(1)
O(2)	1067(1)	7901(1)	3991(1)	20(1)
C(4)	-2122(2)	7715(2)	3752(1)	33(1)
O(6)	1125(1)	7798(1)	6235(1)	28(1)
C(5)	-1519(2)	7851(2)	4247(1)	26(1)
O(5)	194(1)	6530(2)	2639(1)	33(1)
O(14)	1979(1)	4649(2)	5051(1)	49(1)
O(13)	1847(1)	6502(1)	4972(1)	32(1)
O(11)	-678(1)	7999(2)	5974(1)	41(1)
O(12)	-1070(1)	6773(2)	6745(1)	56(1)
O(9)	96(1)	9122(2)	2993(1)	31(1)
C(21)	826(2)	7566(2)	5708(1)	21(1)
O(10)	913(1)	10183(2)	2476(1)	34(1)
C(22)	477(2)	6429(2)	5581(1)	23(1)
O(8)	2610(1)	7224(2)	3960(1)	38(1)
C(25)	1024(2)	5474(2)	5646(1)	27(1)
O(7)	2795(1)	6996(2)	2878(1)	46(1)
C(26)	1649(2)	5599(2)	5192(1)	27(1)
O(15)	-2444(1)	8770(2)	3593(1)	36(1)
C(23)	-121(2)	6179(2)	6090(1)	30(1)
C(24)	-640(2)	7095(2)	6247(1)	34(1)
C(11)	364(2)	6932(2)	3189(1)	22(1)
C(12)	1087(2)	7449(2)	3319(1)	21(1)
C(15)	1265(2)	8372(2)	2805(1)	23(1)
C(16)	693(2)	9253(2)	2780(1)	23(1)
C(13)	1642(2)	6491(2)	3279(1)	25(1)
C(14)	2389(2)	6905(2)	3354(1)	27(1)
O(1W)	3166(1)	4935(2)	4312(1)	56(1)

Table 2. Bond lengths [Å] and angles [°] for 3.

Si-O(4)	1.666(2)	Si-O(2)	1.679(2)
Si-O(3)	1.763(2)	Si-O(1)	1.824(2)
Si-C(1)	1.888(3)	O(3)-C(21)	1.324(3)
N-H	0.89(3)	N-C(5)	1.497(3)
N-C(1)	1.500(3)	N-C(2)	1.500(3)
O(4)-C(22)	1.429(3)	C(2)-C(3)	1.501(4)
O(1)-C(11)	1.304(3)	C(3)-O(15)	1.422(3)
O(2)-C(12)	1.426(2)	C(4)-O(15)	1.429(3)
C(4)-C(5)	1.499(4)	O(6)-C(21)	1.209(3)
O(5)-C(11)	1.224(3)	O(14)-H(14)	0.89(4)
O(14)-C(26)	1.320(3)	O(13)-C(26)	1.218(3)
O(11)-C(24)	1.207(3)	O(12)-H(12)	0.93(4)
O(12)-C(24)	1.323(3)	O(9)-C(16)	1.203(3)
C(21)-C(22)	1.526(3)	O(10)-H(10)	0.86(3)
O(10)-C(16)	1.325(3)	C(22)-C(23)	1.529(4)
C(22)-C(25)	1.537(4)	O(8)-H(8)	0.88(3)
O(8)-C(14)	1.317(3)	C(25)-C(26)	1.478(4)
O(7)-C(14)	1.209(3)	C(23)-C(24)	1.495(4)
C(11)-C(12)	1.508(4)	C(12)-C(15)	1.529(3)
C(12)-C(13)	1.546(4)	C(15)-C(16)	1.502(4)
C(13)-C(14)	1.490(4)	O(1W)-H(1W)	0.83(3)
O(1W)-H(2W)	0.83(3)		

O(4)-Si-O(2)	124.29(9)	O(4)-Si-O(3)	90.52(8)
O(2)-Si-O(3)	90.15(8)	O(4)-Si-O(1)	86.91(8)
O(2)-Si-O(1)	87.87(8)	O(3)-Si-O(1)	175.11(9)
O(4)-Si-C(1)	121.31(11)	O(2)-Si-C(1)	114.36(10)
O(3)-Si-C(1)	91.28(9)	O(1)-Si-C(1)	93.61(9)
C(21)-O(3)-Si	114.2(2)	H-N-C(5)	109(2)
H-N-C(1)	105(2)	C(5)-N-C(1)	111.7(2)
H-N-C(2)	109(2)	C(5)-N-C(2)	109.4(2)
C(1)-N-C(2)	112.8(2)	N-C(1)-Si	116.9(2)
C(22)-O(4)-Si	116.49(14)	N-C(2)-C(3)	109.5(2)
C(11)-O(1)-Si	114.5(2)	O(15)-C(3)-C(2)	111.8(2)
C(12)-O(2)-Si	117.7(2)	O(15)-C(4)-C(5)	111.4(2)
N-C(5)-C(4)	110.3(2)	H(14)-O(14)-C(26)	110(2)
H(12)-O(12)-C(24)	111(2)	O(6)-C(21)-O(3)	125.0(2)
O(6)-C(21)-C(22)	122.7(2)	O(3)-C(21)-C(22)	112.3(2)
H(10)-O(10)-C(16)	107(2)	O(4)-C(22)-C(21)	106.4(2)
O(4)-C(22)-C(23)	110.9(2)	C(21)-C(22)-C(23)	112.4(2)
O(4)-C(22)-C(25)	109.3(2)	C(21)-C(22)-C(25)	111.1(2)
C(23)-C(22)-C(25)	106.8(2)	H(8)-O(8)-C(14)	111(2)
C(26)-C(25)-C(22)	113.7(2)	O(13)-C(26)-O(14)	122.8(3)
O(13)-C(26)-C(25)	123.0(2)	O(14)-C(26)-C(25)	114.2(2)
C(3)-O(15)-C(4)	110.0(2)	C(24)-C(23)-C(22)	117.9(2)
O(11)-C(24)-O(12)	123.5(3)	O(11)-C(24)-C(23)	126.9(2)
O(12)-C(24)-C(23)	109.6(2)	O(5)-C(11)-O(1)	124.4(3)
O(5)-C(11)-C(12)	122.8(2)	O(1)-C(11)-C(12)	112.7(2)
O(2)-C(12)-C(11)	106.7(2)	O(2)-C(12)-C(15)	110.2(2)
C(11)-C(12)-C(15)	112.3(2)	O(2)-C(12)-C(13)	110.1(2)
C(11)-C(12)-C(13)	106.9(2)	C(15)-C(12)-C(13)	110.6(2)
C(16)-C(15)-C(12)	111.6(2)	O(9)-C(16)-O(10)	123.6(3)
O(9)-C(16)-C(15)	124.1(2)	O(10)-C(16)-C(15)	112.3(3)
C(14)-C(13)-C(12)	112.4(2)	O(7)-C(14)-O(8)	118.3(3)
O(7)-C(14)-C(13)	122.9(2)	O(8)-C(14)-C(13)	118.7(2)
H(1W)-O(1W)-H(2W)	108(4)		

Table 3. Anisotropic displacement parameters [$Å^2 \times 10^3$] for 3.

The anisotropic displacement factor exponent takes the form:

$-2\pi^2 [(ha^*)^2 U_{11} + ... + 2hka^* b^* U_{12}]$

	U11	U22	U33	U23	U13	U12
Si	24(1)	20(1)	17(1)	0(1)	-1(1)	-2(1)
O(3)	28(1)	20(1)	18(1)	1(1)	-2(1)	-4(1)
N	21(2)	21(1)	21(1)	-1(1)	2(1)	1(1)
C(1)	24(2)	26(1)	21(1)	-3(1)	-4(1)	-5(1)
O(4)	27(1)	23(1)	18(1)	2(1)	-4(1)	-5(1)
C(2)	29(2)	22(1)	32(1)	4(1)	3(1)	5(1)
O(1)	21(1)	22(1)	20(1)	-1(1)	1(1)	-3(1)
C(3)	30(2)	38(2)	31(1)	8(1)	1(1)	9(1)
O(2)	23(1)	23(1)	14(1)	-2(1)	-1(1)	-2(1)
C(4)	30(2)	34(1)	36(1)	-1(1)	-2(1)	0(1)
O(6)	37(2)	29(1)	19(1)	-1(1)	-7(1)	-1(1)
C(5)	26(2)	25(1)	28(1)	4(1)	3(1)	-3(1)
O(5)	36(2)	40(1)	22(1)	-9(1)	0(1)	-9(1)
O(14)	54(2)	31(1)	62(1)	-2(1)	16(1)	11(1)
O(13)	36(2)	34(1)	26(1)	6(1)	5(1)	4(1)
O(11)	41(2)	44(1)	36(1)	8(1)	13(1)	3(1)
O(12)	34(2)	85(2)	48(1)	36(1)	17(1)	17(1)
O(9)	25(2)	35(1)	34(1)	10(1)	7(1)	3(1)
C(21)	23(2)	23(1)	18(1)	0(1)	3(1)	2(1)
O(10)	33(2)	27(1)	43(1)	13(1)	3(1)	2(1)
C(22)	25(2)	27(1)	17(1)	3(1)	-2(1)	-4(1)
O(8)	26(2)	59(1)	28(1)	0(1)	3(1)	-4(1)
C(25)	34(2)	22(1)	24(1)	2(1)	-2(1)	-1(1)
O(7)	29(2)	79(2)	31(1)	6(1)	6(1)	0(1)
C(26)	31(2)	28(1)	23(1)	1(1)	-6(1)	5(1)
O(15)	25(2)	44(1)	38(1)	5(1)	-2(1)	3(1)
C(23)	30(2)	37(1)	23(1)	7(1)	1(1)	-5(1)
C(24)	25(2)	53(2)	22(1)	9(1)	2(1)	-5(2)
C(11)	22(2)	21(1)	22(1)	-1(1)	-1(1)	3(1)
C(12)	22(2)	25(1)	15(1)	-5(1)	0(1)	0(1)
C(15)	24(2)	27(1)	19(1)	1(1)	0(1)	-2(1)
C(16)	28(3)	26(1)	16(1)	0(1)	-2(1)	-1(1)
C(13)	26(2)	26(1)	24(1)	-2(1)	1(1)	3(1)
C(14)	25(2)	30(1)	27(1)	4(1)	1(2)	7(1)
O(1W)	54(2)	48(1)	65(2)	-1(1)	15(1)	24(1)

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

	x	y	z	U(eq)
H	-782(15)	8330(2)	3596(13)	29(7)
H(1A)	-551	8869	4915	28
H(1B)	-112	9500	4334	28
H(2A)	-967	10225	3556	33
H(2B)	-1499	10103	4193	33
H(3A)	-2161	10219	3179	40
H(3B)	-1741	9161	2876	40
H(4A)	-1940	7366	3329	40
H(4B)	-2487	7209	3949	40
H(5A)	-1704	8147	4684	32
H(5B)	-1297	7112	4337	32
H(14)	2380(2)	4790(3)	4812(17)	59
H(12)	-1430(2)	7300(3)	6816(18)	67
H(10)	553(17)	10640(3)	2460(15)	41
H(8)	2299(18)	7050(3)	4277(15)	45
H(25A)	1193	5438	6123	32
H(25B)	784	4754	5542	32
H(23A)	-395	5529	5916	36
H(23B)	104	5944	6524	36
H(15A)	1725	8725	2932	28
H(15B)	1321	8035	2347	28
H(13A)	1541	5939	3643	30
H(13B)	1593	6103	2836	30
H(1W)	3400(2)	4340(3)	4260(2)	84
H(2W)	3420(2)	5460(3)	4180(2)	84