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SUPPLEMENTARY MATERIAL

N⁶-Cyclopentyl-8-(ethylamino)adenosine (3b): ¹H NMR (DMSO-d₆): δ 7.93 (s, 1H, H-2), 6.85 (t, J = 5.2 Hz, 1H, 8-NH, exchangeable with D₂O), 6.64 (d, J = 7.8 Hz, 1H, 6-NH, exchangeable with D₂O), 5.87 (m, 2H, H-1' and 5'-OH) which changed by addition of D₂O to (d, J_{1',2'} = 7.4 Hz, 1H, H-1'), 5.21 (d, J = 6.4 Hz, 1H, 2'-OH, exchangeable with D₂O), 5.14 (d, J = 3.6 Hz, 1H, 3'-OH, exchangeable with D₂O), 4.64 (AB, which changed by addition of D₂O in dd, J_{1',2'} = 7.5 Hz, J_{2',3'} = 5.3 Hz, 1H, H-2'), 4.58-4.46 (m, 1H, 6-N-CH, cyclopentyl), 4.10 (m, which changed by addition of D₂O in dd, J_{2',3'} = 5.3 Hz, J_{3',4'} = 1.8 Hz, 1H, H-3'), 3.95 (m, 1H, H-4'), 3.62 (m, which changed by addition of D₂O in bs, 2H, H-5'/H-5''), 3.41-3.31 (m, 2H, N-CH₂, ethyl), 1.95-1.88 (m, 2H, N-CH-CHH, cyclopentyl), 1.69-1.52 (m, 6H, N-CH-CHH-CH₂, cyclopentyl), 1.17 (t, J = 7.2 Hz, 3H, N-CH₂-CH₃, ethyl); ¹³C NMR (DMSO-d₆): δ 151.4 (C-6 or C-4), 151.2 (C-6 or C-4), 149.2 (C-8), 149.1 (C-2), 117.0 (C-5), 86.7 (C-1'), 85.9 (C-4'), 71.1 (C-2'), 71.0 (C-3'), 61.8 (C-5'), 52.0 (N-CH, cyclopentyl), 37.4 (N-CH₂, ethyl), 32.9 and 32.8 (N-CH-CH₂, cyclopentyl), 23.7 (N-CH-CH₂-CH₂, cyclopentyl), 14.8 (N-CH₂-CH₃, ethyl);

N⁶-Cyclopentyl-8-(n-propylamino)adenosine (3c): ¹H NMR (DMSO-d₆): δ 7.93 (s, 1H, H-2), 6.87 (t, J = 5.4 Hz, 1H, 8-NH, exchangeable with D₂O), 6.63 (d, J = 7.8 Hz, 1H, 6-NH, exchangeable with D₂O), 5.89 (d, J_{1',2'} = 7.5 Hz, 1H, 1H, H-1'), 5.87 (t, J = 4.6 Hz, 1H, 5'-OH, exchangeable with D₂O), 5.21 (d, J = 6.8 Hz, 1H, 2'-OH, exchangeable with D₂O), 5.14 (d, J = 4.0 Hz, 1H, 3'-OH, exchangeable with D₂O), 4.63 (AB, which changed by addition of D₂O in dd, J_{1',2'} = 7.5 Hz, J_{2',3'} = 5.4 Hz, 1H, H-2'), 4.53-4.49 (m, 1H, 6-N-CH, cyclopentyl), 4.10 (m, which changed by addition of D₂O in dd, J_{2',3'} = 5.3 Hz, J_{3',4'} = 1.5 Hz, 1H, H-3'), 3.95 (m, 1H, H-4'), 3.61 (m, which changed by addition of D₂O in bs, 2H, H-5'/H-5''), 3.39-3.25 (m, 2H, N-CH₂, propyl), 1.95-1.89 (m, 2H, N-CH-CHH, cyclopentyl), 1.69-1.52 (m, 8H, N-CH₂-CH₂, propyl and N-CH-CHH-CH₂, cyclopentyl), 0.89 (t, J = 7.4 Hz, 3H, N-CH₂-CH₂-CH₃, propyl); ¹³C NMR (DMSO-d₆): δ 151.4 (C-6 or C-4), 151.2 (C-6 or C-4), 149.1 (C-8), 149.0 (C-2), 117.0 (C-5), 86.6 (C-1'), 85.9 (C-4'), 71.1 (C-2'), 70.9 (C-3'), 61.8 (C-5'), 52.0 (N-CH, cyclopentyl), 44.3 (N-CH₂, propyl), 32.9 and 32.8 (N-CH-CH₂, cyclopentyl), 23.6 (N-CH-CH₂-CH₂, cyclopentyl), 22.2 (N-CH₂-CH₂, propyl), 11.6 (N-CH₂-CH₂-CH₃, propyl);

N⁶-Cyclopentyl-8-(n-butylamino)adenosine (3d): ¹H NMR (DMSO-d₆): δ 7.93 (s, 1H, H-2), 6.82 (t, J = 5.4 Hz, 1H, 8-NH, exchangeable with D₂O), 6.60 (d, J = 7.8 Hz, 1H, 6-NH, exchangeable with D₂O), 5.89 (d, J_{1',2'} = 7.5 Hz, 1H, 1H, H-1'), 5.85 (t, J = 4.6 Hz, 1H, 5'-OH, exchangeable with D₂O), 5.20 (d, J = 6.8 Hz, 1H, 2'-OH, exchangeable with D₂O), 5.13 (d, J = 4.1 Hz, 1H, 3'-OH, exchangeable with D₂O), 4.62 (AB, which changed

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by addition of D₂O in dd, $J_{1',2'} = 7.5$ Hz, $J_{2',3'} = 5.4$ Hz, 1H, H-2'), 4.54-4.50 (m, 1H, 6-N-CH, cyclopentyl), 4.11 (m, which changed by addition of D₂O in dd, $J_{2',3'} = 5.3$ Hz, $J_{3',4'} = 1.7$ Hz, 1H, H-3'), 3.96 (m, 1H, H-4'), 3.62 (m, which changed by addition of D₂O in bs, 2H, H-5'/H-5''), 3.40-3.26 (m, 2H, N-CH₂, butyl), 1.96-1.88 (m, 2H, N-CH-CHH, cyclopentyl), 1.70-1.48 (m, 8H, N-CH₂-CH₂, butyl and N-CH-CHH-CH₂, cyclopentyl), 1.33 (m, 2H, N-CH₂-CH₂-CH₂, butyl), 0.89 (t, $J = 7.3$ Hz, 3H, N-CH₂-CH₂-CH₂-CH₃, butyl); ¹³C NMR (DMSO-d₆): δ 151.3 (C-6 or C-4), 151.1 (C-6 or C-4), 149.1 (C-8), 148.8 (C-2), 117.0 (C-5), 86.5 (C-1'), 85.7 (C-4'), 70.9 (C-2'), 70.8 (C-3'), 61.6 (C-5'), 51.9 (N-CH, cyclopentyl), 42.0 (N-CH₂, butyl), 32.8 and 32.7 (N-CH-CH₂, cyclopentyl), 30.9 (N-CH₂-CH₂, butyl), 23.5 (N-CH-CH₂-CH₂, cyclopentyl), 19.7 (N-CH₂-CH₂-CH₂, butyl), 13.8 (N-CH₂-CH₂-CH₂-CH₃, butyl);

N⁶-Cyclopentyl-8-(cyclopentylamino)adenosine (3e): ¹H NMR (DMSO-d₆): δ 7.94 (s, 1H, H-2), 6.70 (d, $J = 7.0$ Hz, 1H, 8-NH, exchangeable with D₂O), 6.64 (d, $J = 7.8$ Hz, 1H, 6-NH, exchangeable with D₂O), 5.91 (d, $J_{1',2'} = 7.5$ Hz, 1H, 1H, H-1'), 5.87 (t, $J = 4.6$ Hz, 1H, 5'-OH, exchangeable with D₂O), 5.21 (d, $J = 6.1$ Hz, 1H, 2'-OH, exchangeable with D₂O), 5.18 (d, $J = 5.4$ Hz, 1H, 3'-OH, exchangeable with D₂O), 4.59 (AB, which changed by addition of D₂O in dd, $J_{1',2'} = 7.5$ Hz, $J_{2',3'} = 5.5$ Hz, 1H, H-2'), 4.58-4.52 (m, 1H, 6-N-CH, cyclopentyl), 4.22 (q, $J = 6.9$ Hz, 1H, 8-N-CH, cyclopentyl), 4.10 (m, which changed by addition of D₂O in dd, $J_{2',3'} = 5.3$ Hz, $J_{3',4'} = 1.8$ Hz, 1H, H-3'), 3.95 (m, 1H, H-4'), 3.62 (m, which changed by addition of D₂O in bs, 2H, H-5'/H-5''), 1.96-1.88 (m, 4H, N-CH-CHH, 2 x cyclopentyl), 1.71-1.45 (m, 12H, N-CH-CHH-CH₂, 2 x cyclopentyl); ¹³C NMR (DMSO-d₆): δ 151.2 (C-6 or C-4), 151.1 (C-6 or C-4), 149.2 (C-8), 149.1 (C-2), 117.0 (C-5), 86.6 (C-1'), 86.0 (C-4'), 71.1 (C-2'), 71.0 (C-3'), 61.8 (C-5'), 54.3 and 52.1 (N-CH, 2 x cyclopentyl), 33.1, 33.0, 32.6 and 32.5 (N-CH-CH₂, 2 x cyclopentyl), 23.8 and 23.7 (N-CH-CH₂-CH₂, 2 x cyclopentyl);

8-Cyclopentylamino-N⁶-(*n*-propyl)adenosine (9c): ¹H NMR (DMSO-d₆): δ 7.92 (s, 1H, H-2), 6.82 (t, $J = 6.1$ Hz, 1H, 6-NH, exchangeable with D₂O), 6.66 (d, $J = 6.9$ Hz, 1H, 8-NH, exchangeable with D₂O), 5.90 (d, $J_{1',2'} = 7.4$ Hz, 1H, H-1'), 5.82 (t, $J = 4.7$ Hz, 1H, 5'-OH, exchangeable with D₂O), 5.17 (d, $J = 6.7$ Hz, 1H, 2'-OH, exchangeable with D₂O), 5.11 (d, $J = 4.0$ Hz, 1H, 3'-OH, exchangeable with D₂O), 4.64 (AB, which changed by addition of D₂O in dd, $J_{1',2'} = 7.5$ Hz, $J_{2',3'} = 5.4$ Hz, 1H, H-2'), 4.11 (m, 1H, 8-N-CH, cyclopentyl), 4.07 (m, which changed by addition of D₂O in dd, $J_{2',3'} = 5.4$ Hz, $J_{3',4'} = 1.9$ Hz, 1H, H-3'), 3.95 (m, 1H, H-4'), 3.60 (m, which changed by addition of D₂O in d, $J = 2.2$ Hz, 2H, H-5'/H-5''), 3.46-3.38 (m, 2H, N-CH₂, propyl), 1.95-1.88 (m, 2H, N-CH-CHH, cyclopentyl), 1.67-1.46 (m, 8H, N-CH₂-CH₂, propyl and N-CH-CHH-CH₂, cyclopentyl), 0.86 (t, $J = 7.4$ Hz, 3H, N-CH₂-CH₂-CH₃, propyl); ¹³C NMR (DMSO-d₆): δ 151.5 (C-

6 or C-4), 150.7 (C-6 or C-4), 149.0 (C-8), 148.3 (C-2), 117.0 (C-5), 86.3 (C-1'), 86.0 (C-4'), 70.7 (C-2' and C-3'), 61.5 (C-5'), 53.9 (N-CH, cyclopentyl), 41.9 (N-CH₂, propyl), 32.2 and 32.1 (N-CH-CH₂, cyclopentyl), 23.3 and 23.2 (N-CH-CH₂-CH₂, cyclopentyl), 22.7 (N-CH₂-CH₂, propyl), 11.1 (N-CH₂-CH₂-CH₃, propyl);

8-Cyclopentylamino-N⁶-(*n*-butyl)adenosine (9d): ¹H NMR (DMSO-d₆): δ 7.92 (s, 1H, H-2), 6.77 (t, J = 5.9 Hz, 1H, 6-NH, exchangeable with D₂O), 6.65 (d, J = 6.9 Hz, 1H, 8-NH, exchangeable with D₂O), 5.90 (d, J_{1',2'} = 7.5 Hz, 1H, H-1'), 5.80 (t, J = 5.1 Hz, 1H, 5'-OH, exchangeable with D₂O), 5.15 (d, J = 6.9 Hz, 1H, 2'-OH, exchangeable with D₂O), 5.10 (d, J = 4.2 Hz, 1H, 3'-OH, exchangeable with D₂O), 4.60 (AB, which changed by addition of D₂O in dd, J_{1',2'} = 7.6 Hz, J_{2',3'} = 5.5 Hz, 1H, H-2'), 4.21 (m, 1H, 8-N-CH, cyclopentyl), 4.11 (m, which changed by addition of D₂O in dd, J_{2',3'} = 5.4 Hz, J_{3',4'} = 1.6 Hz, 1H, H-3'), 3.94 (m, 1H, H-4'), 3.62 (m, which changed by addition of D₂O in bs, 2H, H-5'/H-5''), 3.59-3.40 (m, 2H, N-CH₂, butyl), 1.97-1.90 (m, 2H, N-CH-CHH, cyclopentyl), 1.73-1.45 (m, 8H, N-CH₂-CH₂, butyl and N-CH-CHH-CH₂, cyclopentyl), 1.31 (m, 2H, N-CH₂-CH₂-CH₂, butyl), 0.88 (t, J = 7.3 Hz, 3H, N-CH₂-CH₂-CH₂-CH₃, butyl); ¹³C NMR (DMSO-d₆): δ 151.5 (C-6 or C-4), 150.7 (C-6 or C-4), 149.0 (C-8), 148.3 (C-2), 117.1 (C-5), 86.3 (C-1'), 85.5 (C-4'), 70.7 (C-2' and C-3'), 61.5 (C-5'), 53.9 (N-CH, cyclopentyl), 39.8 (N-CH₂, butyl), 32.2 and 32.1 (N-CH-CH₂, cyclopentyl), 31.7 (N-CH₂-CH₂, butyl), 23.2 (N-CH-CH₂-CH₂, cyclopentyl), 19.4 (N-CH₂-CH₂-CH₂, butyl), 13.6 (N-CH₂-CH₂-CH₂-CH₃, butyl);

8-Cyclopentylamino-N⁶-ethyladenosine (9b): ¹H NMR (DMSO-d₆): δ 7.93 (s, 1H, H-2), 6.78 (t, J = 5.6 Hz, 1H, 6-NH, exchangeable with D₂O), 6.66 (d, J = 6.7 Hz, 1H, 8-NH, exchangeable with D₂O), 5.90 (d, J_{1',2'} = 7.4 Hz, 1H, H-1'), 5.80 (t, J = 4.5 Hz, 1H, 5'-OH, exchangeable with D₂O), 5.15 (d, J = 6.9 Hz, 1H, 2'-OH, exchangeable with D₂O), 5.10 (d, J = 4.0 Hz, 1H, 3'-OH, exchangeable with D₂O), 4.59 (AB, which changed by addition of D₂O in dd, J_{1',2'} = 7.5 Hz, J_{2',3'} = 5.5 Hz, 1H, H-2'), 4.21 (m, 1H, 8-N-CH, cyclopentyl), 4.10 (m, which changed by addition of D₂O in dd, J_{2',3'} = 5.3 Hz, J_{3',4'} = 1.4 Hz, 1H, H-3'), 3.95 (m, 1H, H-4'), 3.62 (m, which changed by addition of D₂O in bs, 2H, H-5'/H-5''), 3.49 (m, 2H, N-CH₂, ethyl), 1.95-1.87 (m, 2H, N-CH-CHH, cyclopentyl), 1.74-1.52 (m, 6H, N-CH-CHH-CH₂, cyclopentyl), 1.14 (t, J = 7.1 Hz, 3H, N-CH₂-CH₃, ethyl); ¹³C NMR (DMSO-d₆): δ 151.4 (C-6 or C-4), 151.2 (C-6 or C-4), 149.0 (C-8), 148.2 (C-2), 117.1 (C-5), 86.7 (C-1'), 85.9 (C-4'), 71.0 (C-2' and C-3'), 61.7 (C-5'), 53.8 (N-CH, cyclopentyl), 37.3 (N-CH₂, ethyl), 32.4 and 32.2 (N-CH-CH₂, cyclopentyl), 23.4 and 23.3 (N-CH-CH₂-CH₂, cyclopentyl), 14.7 (N-CH₂-CH₃, ethyl);

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Crystal data of **8-cyclopentylamino-N⁶-ethyladenosine (9b)**.

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S U P P L E M E N T A R Y M A T E R I A L

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B E L O N G I N G T O T H E P A P E R

N⁶, C8-disubstituted adenosine derivatives as
partial agonists for adenosine A₁ receptors

b y

Harlof Roelen, Nora Veldman, Anthony L. Spek,
Jacobien von Frijtag Drabbe K^{unzel},
Ron A.A. Math^{not} and Ad P. IJzerman

C o n t e n t s

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Table S1 - Crystal Data and Details of the Structure Determination
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S2 - Final Coordinates and Equivalent Isotropic Thermal
Parameters of the non-Hydrogen atoms
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S3 - Hydrogen Atom Positions and Isotropic Thermal
Parameters
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S4 - (An)isotropic Thermal Parameters
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S5 - Bond Distances (Angstrom)
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S6 - Bond Angles (Degrees)
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S7 - Torsion Angles (Degrees)
for: s962a C17 H26 N6 O4 * C4 H10 O1

Table S1 - Crystal Data and Details of the Structure Determination
for: s962a C17 H26 N6 O4 * C4 H10 O1

Crystal Data

Empirical Formula	C17 H26 N6 O4 * C4 H10 O1		
Formula Weight	452.55		
Crystal System	Orthorhombic		
Space group	P212121	(No. 19)	
a, b, c [Angstrom]	11.039(3)	8.708(2)	24.815(12)
alpha, beta, gamma [deg]	90	90	90
V [Ang ³]	2385.4(14)		
Z	4		
D(calc) [g/cm ³]	1.260		
F(000) [Electrons]	976		
Mu(MoKa) [/cm]	0.9		
Crystal Size [mm]	0.03 x 0.50 x 0.63		

Data Collection

Temperature (K)	295		
Radiation [Angstrom]	MoKa (with monochromator)	0.71073	
Theta Min-Max [Deg]	1.6, 24.2		
Scan type	Omega/2Theta		
Scan, [Deg]	1.69 + 0.35 Tan(Theta)		
Hor. and vert. aperture [mm]	4.53	4.00	
Reference Reflection(s)	-2, 0, 5; -1, -2, -2; 3 0 -2		
Dataset	-12: 0 ; -10: 0 ; -27: 27		
Tot., Uniq. Data	3913, 2874		
R(int) = 0.0815 R(sigma) = 0.1992 Friedel opposites not merged			

Refinement

wR2 = 0.2455 before cycle 4 for 3424 data and 2 / 298 parameters	
R1 = 0.0974 for 1450 Fo > 4.sigma(Fo) and 0.2247 for all 3426 data	
GooF = S = 0.947; Restrained GooF = 0.947 for 1 restraints	
Weight	1/[sigma ² (Fo ²) + (0.1000 * [(Max(Fo ² , 0) + 2*Fc ²)/3]) ²]
Max. and Av. Shift/Error	0.004, 0.001
Min. and Max. resd. dens. [e/Ang ³]	-0.25, 0.23

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Table S2 - Final Coordinates and Equivalent Isotropic Thermal Parameters of the non-Hydrogen atoms
for: s962a C17 H26 N6 O4 * C4 H10 O1

Atom	x	y	z	U(eq) [Ang**2]
O(1)	0.1215(6)	0.5856(6)	0.4866(3)	0.044(3)
O(12)	0.0738(6)	0.7682(8)	0.3613(3)	0.056(3)
O(13)	0.1332(7)	0.9294(6)	0.4531(3)	0.055(3)
O(15)	-0.1367(7)	0.6161(8)	0.5017(3)	0.059(3)
N(1)	0.4140(9)	0.2168(11)	0.3335(4)	0.073(4)
N(3)	0.3473(9)	0.4616(9)	0.3705(4)	0.056(4)
N(6)	0.2978(9)	-0.0011(10)	0.3371(4)	0.071(4)
N(7)	0.1075(8)	0.1959(8)	0.3950(3)	0.042(3)
N(8)	-0.0430(8)	0.3432(8)	0.4368(3)	0.044(3)
N(9)	0.1441(7)	0.4466(7)	0.4082(3)	0.036(3)
C(2)	0.4262(11)	0.3683(13)	0.3464(5)	0.071(5)
C(4)	0.2479(9)	0.3861(10)	0.3846(4)	0.038(4)
C(5)	0.2205(10)	0.2320(10)	0.3744(4)	0.044(4)
C(6)	0.3126(12)	0.1537(12)	0.3481(5)	0.058(4)
C(8)	0.0675(10)	0.3261(9)	0.4137(4)	0.037(3)
C(11)	0.1404(10)	0.6004(9)	0.4306(4)	0.042(4)
C(12)	0.0418(8)	0.7059(9)	0.4107(4)	0.034(3)
C(13)	0.0347(9)	0.8217(8)	0.4584(4)	0.038(3)
C(14)	0.0588(10)	0.7211(9)	0.5063(5)	0.053(4)
C(15)	-0.0502(9)	0.6750(12)	0.5372(5)	0.055(4)
C(61)	0.3920(12)	-0.0924(15)	0.3086(7)	0.105(7)
C(62)	0.3441(17)	-0.207(2)	0.2749(8)	0.176(11)
C(81)	-0.1200(10)	0.2106(10)	0.4463(5)	0.055(4)
C(82)	-0.1727(11)	0.1494(15)	0.3945(5)	0.078(5)
C(83)	-0.2866(15)	0.209(3)	0.3891(7)	0.174(12)
C(84)	-0.3248(14)	0.291(2)	0.4382(7)	0.121(8)
C(85)	-0.2311(11)	0.2557(14)	0.4782(5)	0.071(5)

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Table S2 - Final Coordinates and Equivalent Isotropic Thermal Parameters of the non-Hydrogen atoms (continued)
for: s962a C17 H26 N6 O4 * C4 H10 O1

Atom	x	y	z	U(eq) [Ang**2]
O(33)	0.1314(9)	0.3346(10)	0.2015(3)	0.085(4)
C(31)	0.2793(14)	0.199(2)	0.1514(7)	0.121(8)
C(32)	0.2522(16)	0.295(2)	0.1984(8)	0.132(10)
C(34)	0.108(2)	0.436(2)	0.2462(6)	0.133(9)
C(35)	-0.011(2)	0.457(3)	0.2562(10)	0.168(11)

U(eq) = 1/3 of the trace of the orthogonalized U

Table S3 - Hydrogen Atom Positions and Isotropic Thermal Parameters
for: s962a C17 H26 N6 O4 * C4 H10 O1

Atom	x	y	z	U(iso) [Ang**2]
H(2)	0.4999(11)	0.4128(13)	0.3371(5)	0.085
H(6)	0.2365	-0.0655	0.3435	0.107
H(8)	-0.043(10)	0.412(8)	0.463(3)	0.066
H(11)	0.2190(10)	0.6498(9)	0.4245(4)	0.051
H(12)	0.014(2)	0.773(13)	0.3418(17)	0.084
H(12A)	-0.0347(8)	0.6494(9)	0.4074(4)	0.041
H(13)	0.122(5)	0.983(9)	0.427(3)	0.083
H(13A)	-0.0443(9)	0.8728(8)	0.4605(4)	0.046
H(14)	0.1134(10)	0.7762(9)	0.5307(5)	0.064
H(15)	-0.2034(19)	0.617(13)	0.5162(17)	0.088
H(15A)	-0.0288(9)	0.5976(12)	0.5637(5)	0.066
H(15B)	-0.0832(9)	0.7631(12)	0.5561(5)	0.066
H(61A)	0.4414(12)	-0.0235(15)	0.2871(7)	0.126
H(61B)	0.4443(12)	-0.1404(15)	0.3351(7)	0.126
H(62A)	0.395(10)	-0.296(8)	0.276(6)	0.2640
H(62B)	0.340(16)	-0.169(8)	0.2386(16)	0.2640
H(62C)	0.264(7)	-0.234(16)	0.287(5)	0.2640
H(81)	-0.0749(10)	0.1299(10)	0.4651(5)	0.065
H(82A)	-0.1226(11)	0.1795(15)	0.3642(5)	0.093
H(82B)	-0.1765(11)	0.0381(15)	0.3956(5)	0.093
H(83A)	-0.3434(15)	0.127(3)	0.3816(7)	0.2110
H(83B)	-0.2879(15)	0.280(3)	0.3589(7)	0.2110
H(84A)	-0.3294(14)	0.401(2)	0.4318(7)	0.1450
H(84B)	-0.4033(14)	0.255(2)	0.4502(7)	0.1450
H(85A)	-0.2568(11)	0.1720(14)	0.5013(5)	0.085
H(85B)	-0.2145(11)	0.3449(14)	0.5004(5)	0.085

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Table S3 - Hydrogen Atom Positions and Isotropic Thermal Parameters (continued)
for: s962a C17 H26 N6 O4 * C4 H10 O1

Atom	x	y	z	U(iso) [Ang**2]
H(31A)	0.362(4)	0.163(11)	0.154(3)	0.1820
H(31B)	0.269(10)	0.258(4)	0.1191(7)	0.1820
H(31C)	0.225(7)	0.112(7)	0.151(3)	0.1820
H(32A)	0.2751(16)	0.240(2)	0.2309(8)	0.158
H(32B)	0.3006(16)	0.388(2)	0.1965(8)	0.158
H(34A)	0.146(2)	0.395(2)	0.2782(6)	0.16
H(34B)	0.144(2)	0.535(2)	0.2386(6)	0.16
H(35A)	-0.021(2)	0.500(18)	0.291(3)	0.255
H(35B)	-0.052(3)	0.359(4)	0.255(7)	0.255
H(35C)	-0.045(4)	0.525(15)	0.230(4)	0.255

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The Temperature Factor has the Form of $\text{Exp}(-T)$ Where
 $T = 8 * (\text{Pi}^2) * U * (\text{Sin}(\text{Theta}) / \text{Lambda})^2$ for Isotropic Atoms

Table S4 - (An)isotropic Thermal Parameters
for: s962a C17 H26 N6 O4 * C4 H10 O1

Atom	U(1,1) or U	U(2,2)	U(3,3)	U(2,3)	U(1,3)	U(1,2)
O(1)	0.053(5)	0.034(3)	0.046(5)	0.000(3)	-0.012(4)	0.006(3)
O(12)	0.065(5)	0.050(4)	0.053(5)	0.017(4)	0.002(4)	0.004(4)
O(13)	0.049(4)	0.037(3)	0.080(6)	0.016(4)	-0.011(5)	-0.006(4)
O(15)	0.047(5)	0.066(4)	0.063(5)	-0.010(4)	0.008(5)	-0.009(4)
N(1)	0.056(7)	0.063(6)	0.101(9)	-0.015(6)	0.019(6)	0.008(5)
N(3)	0.049(6)	0.063(6)	0.055(7)	-0.018(5)	0.017(5)	0.002(5)
N(6)	0.050(6)	0.055(5)	0.108(9)	-0.029(6)	0.001(6)	0.003(5)
N(7)	0.046(6)	0.034(4)	0.046(6)	-0.003(4)	0.004(5)	0.004(4)
N(8)	0.051(6)	0.030(4)	0.051(6)	-0.012(4)	0.005(5)	-0.008(4)
N(9)	0.028(5)	0.031(4)	0.050(6)	0.000(4)	-0.004(4)	0.001(4)
C(2)	0.057(8)	0.064(7)	0.092(10)	-0.010(7)	0.025(8)	-0.027(7)
C(4)	0.029(6)	0.048(6)	0.038(7)	0.011(5)	0.010(5)	0.009(5)
C(5)	0.056(8)	0.033(5)	0.042(7)	-0.002(4)	-0.010(6)	0.005(5)
C(6)	0.074(9)	0.059(7)	0.041(7)	-0.015(6)	0.001(7)	0.001(7)
C(8)	0.051(7)	0.036(5)	0.023(5)	0.002(4)	0.000(5)	0.010(5)
C(11)	0.045(7)	0.039(5)	0.043(7)	-0.002(5)	0.001(6)	-0.004(5)
C(12)	0.027(6)	0.032(4)	0.043(7)	-0.002(4)	-0.020(5)	-0.004(4)
C(13)	0.044(6)	0.021(4)	0.049(6)	-0.007(4)	-0.012(6)	-0.003(4)
C(14)	0.065(8)	0.031(5)	0.063(8)	-0.005(5)	-0.004(7)	0.007(5)
C(15)	0.035(6)	0.060(6)	0.069(8)	-0.009(6)	0.014(7)	0.003(6)
C(61)	0.065(10)	0.080(9)	0.169(17)	-0.059(10)	0.034(11)	0.004(8)
C(62)	0.158(19)	0.19(2)	0.18(2)	-0.090(17)	0.088(17)	0.000(17)
C(81)	0.046(7)	0.047(6)	0.071(9)	-0.001(5)	-0.004(7)	-0.007(5)
C(82)	0.059(9)	0.088(8)	0.086(11)	-0.008(8)	0.021(8)	-0.016(7)
C(83)	0.084(13)	0.35(3)	0.089(15)	-0.071(18)	-0.017(12)	0.075(18)
C(84)	0.092(13)	0.153(15)	0.118(16)	0.006(13)	0.040(12)	0.021(12)
C(85)	0.062(9)	0.065(7)	0.087(11)	0.021(7)	0.004(8)	-0.029(7)

Table S4 - (An)isotropic Thermal Parameters (continued)
for: s962a C17 H26 N6 O4 * C4 H10 O1

Atom	U(1,1) or U	U(2,2)	U(3,3)	U(2,3)	U(1,3)	U(1,2)
O(33)	0.091(7)	0.099(6)	0.065(7)	-0.008(5)	-0.012(6)	-0.002(6)
C(31)	0.097(13)	0.156(15)	0.111(15)	-0.018(13)	0.027(12)	0.031(11)
C(32)	0.097(14)	0.142(16)	0.156(19)	-0.018(14)	-0.039(13)	0.001(13)
C(34)	0.19(2)	0.153(16)	0.055(11)	-0.010(11)	-0.003(14)	-0.024(17)
C(35)	0.151(19)	0.164(19)	0.19(2)	-0.094(17)	0.02(2)	-0.020(17)

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The Temperature Factor has the Form of $\text{Exp}(-T)$ Where

$T = 8 * (\pi^2) * U * (\sin(\theta) / \lambda)^2$ for Isotropic Atoms

$T = 2 * (\pi^2) * \sum_i h(i) * h(j) * U(i,j) * Astar(i) * Astar(j)$, for Anisotropic Atoms. Astar(i) are Reciprocal Axial Lengths and h(i) are the Reflection Indices.

Table S5 - Bond Distances (Angstrom)
for: s962a C17 H26 N6 O4 * C4 H10 O1

O(1)	-C(11)	1.411(12)	C(84)	-C(85)	1.47(2)
O(1)	-C(14)	1.453(11)	O(12)	-H(12)	0.82(3)
O(12)	-C(12)	1.386(12)	O(13)	-H(13)	0.81(8)
O(13)	-C(13)	1.442(11)	O(15)	-H(15)	0.82(3)
O(15)	-C(15)	1.397(13)	N(6)	-H(6)	0.893(9)
N(1)	-C(2)	1.364(15)	N(8)	-H(8)	0.88(7)
N(1)	-C(6)	1.299(16)	C(2)	-H(2)	0.930(17)
N(3)	-C(2)	1.333(15)	C(11)	-H(11)	0.980(15)
N(3)	-C(4)	1.326(14)	C(12)	-H(12A)	0.981(12)
N(6)	-C(6)	1.385(14)	C(13)	-H(13A)	0.980(13)
N(6)	-C(61)	1.488(17)	C(14)	-H(14)	0.980(15)
N(7)	-C(5)	1.384(14)	C(15)	-H(15A)	0.971(16)
N(7)	-C(8)	1.302(11)	C(15)	-H(15B)	0.970(15)
N(8)	-C(8)	1.356(14)	C(61)	-H(61A)	0.97(2)
N(8)	-C(81)	1.453(12)	C(61)	-H(61B)	0.97(2)
N(9)	-C(4)	1.390(12)	C(62)	-H(62A)	0.96(9)
N(9)	-C(8)	1.355(11)	C(62)	-H(62B)	0.96(5)
N(9)	-C(11)	1.451(10)	C(62)	-H(62C)	0.96(9)
C(4)	-C(5)	1.399(13)	C(81)	-H(81)	0.979(15)
C(5)	-C(6)	1.387(16)	C(82)	-H(82A)	0.969(18)
C(11)	-C(12)	1.508(13)	C(82)	-H(82B)	0.970(18)
C(12)	-C(13)	1.557(13)	C(83)	-H(83A)	0.97(3)
C(13)	-C(14)	1.500(14)	C(83)	-H(83B)	0.97(3)
C(14)	-C(15)	1.482(16)	C(84)	-H(84A)	0.97(2)
C(61)	-C(62)	1.41(2)	C(84)	-H(84B)	0.97(2)
C(81)	-C(82)	1.508(17)	C(85)	-H(85A)	0.970(17)
C(81)	-C(85)	1.512(17)	C(85)	-H(85B)	0.970(17)
C(82)	-C(83)	1.37(2)	O(33)	-C(32)	1.38(2)
C(83)	-C(84)	1.47(3)	O(33)	-C(34)	1.441(18)

Table S5 - Bond Distances (Angstrom) (continued)
for: s962a C17 H26 N6 O4 * C4 H10 O1

C(31)	-C(32)	1.47(3)	C(32)	-H(32B)	0.97(2)
C(34)	-C(35)	1.35(3)	C(34)	-H(34A)	0.97(2)
C(31)	-H(31A)	0.97(5)	C(34)	-H(34B)	0.97(3)
C(31)	-H(31B)	0.96(3)	C(35)	-H(35A)	0.95(10)
C(31)	-H(31C)	0.97(7)	C(35)	-H(35B)	0.97(4)
C(32)	-H(32A)	0.97(3)	C(35)	-H(35C)	0.96(11)

Table S6 - Bond Angles (Degrees)
for: s962a C17 H26 N6 O4 * C4 H10 O1

C(11)	-O(1)	-C(14)	109.1(7)	O(13)	-C(13)	-C(14)	108.6(8)
C(2)	-N(1)	-C(6)	115.4(10)	C(12)	-C(13)	-C(14)	102.4(6)
C(2)	-N(3)	-C(4)	110.9(9)	O(1)	-C(14)	-C(13)	107.0(9)
C(6)	-N(6)	-C(61)	122.1(10)	O(1)	-C(14)	-C(15)	109.9(7)
C(5)	-N(7)	-C(8)	103.9(8)	C(13)	-C(14)	-C(15)	115.1(9)
C(8)	-N(8)	-C(81)	120.5(7)	O(15)	-C(15)	-C(14)	109.2(10)
C(4)	-N(9)	-C(8)	105.3(7)	N(6)	-C(61)	-C(62)	113.5(12)
C(4)	-N(9)	-C(11)	122.3(8)	N(8)	-C(81)	-C(82)	111.6(9)
C(8)	-N(9)	-C(11)	131.2(8)	N(8)	-C(81)	-C(85)	110.7(8)
N(1)	-C(2)	-N(3)	129.1(11)	C(82)	-C(81)	-C(85)	103.0(9)
N(3)	-C(4)	-N(9)	127.2(8)	C(81)	-C(82)	-C(83)	107.7(12)
N(3)	-C(4)	-C(5)	127.4(9)	C(82)	-C(83)	-C(84)	111.5(14)
N(9)	-C(4)	-C(5)	105.2(8)	C(83)	-C(84)	-C(85)	104.8(13)
N(7)	-C(5)	-C(4)	110.2(8)	C(81)	-C(85)	-C(84)	105.8(11)
N(7)	-C(5)	-C(6)	136.3(9)	C(12)	-O(12)	-H(12)	110(4)
C(4)	-C(5)	-C(6)	113.5(10)	C(13)	-O(13)	-H(13)	109(4)
N(1)	-C(6)	-N(6)	117.3(11)	C(15)	-O(15)	-H(15)	109(4)
N(1)	-C(6)	-C(5)	123.7(10)	C(6)	-N(6)	-H(6)	131.7(11)
N(6)	-C(6)	-C(5)	119.0(11)	C(61)	-N(6)	-H(6)	106.2(9)
N(7)	-C(8)	-N(8)	123.5(8)	C(8)	-N(8)	-H(8)	113(7)
N(7)	-C(8)	-N(9)	115.3(9)	C(81)	-N(8)	-H(8)	115(6)
N(8)	-C(8)	-N(9)	121.3(7)	N(1)	-C(2)	-H(2)	115.5(13)
O(1)	-C(11)	-N(9)	107.3(7)	N(3)	-C(2)	-H(2)	115.4(13)
O(1)	-C(11)	-C(12)	105.8(8)	O(1)	-C(11)	-H(11)	108.8(10)
N(9)	-C(11)	-C(12)	117.2(8)	N(9)	-C(11)	-H(11)	108.7(10)
O(12)	-C(12)	-C(11)	110.1(8)	C(12)	-C(11)	-H(11)	108.7(9)
O(12)	-C(12)	-C(13)	115.6(7)	O(12)	-C(12)	-H(12A)	110.0(10)
C(11)	-C(12)	-C(13)	100.5(7)	C(11)	-C(12)	-H(12A)	110.1(9)
O(13)	-C(13)	-C(12)	108.3(8)	C(13)	-C(12)	-H(12A)	110.2(10)

Table S6 - Bond Angles for: s962a			(Degrees)	(continued)			
			C17 H26 N6	O4 * C4 H10 O1			
O(13)	-C(13)	-H(13A)	112.4(8)	H(82A) -C(82)	-H(82B)	108.4(16)	
C(12)	-C(13)	-H(13A)	112.3(10)	C(82) -C(83)	-H(83A)	110(3)	
C(14)	-C(13)	-H(13A)	112.4(11)	C(82) -C(83)	-H(83B)	109.3(18)	
O(1)	-C(14)	-H(14)	108.2(11)	C(84) -C(83)	-H(83A)	109.3(18)	
C(13)	-C(14)	-H(14)	108.2(10)	C(84) -C(83)	-H(83B)	109(3)	
C(15)	-C(14)	-H(14)	108.2(13)	H(83A) -C(83)	-H(83B)	108(2)	
O(15)	-C(15)	-H(15A)	109.8(11)	C(83) -C(84)	-H(84A)	111(2)	
O(15)	-C(15)	-H(15B)	109.8(11)	C(83) -C(84)	-H(84B)	110.7(19)	
C(14)	-C(15)	-H(15A)	109.9(11)	C(85) -C(84)	-H(84A)	110.7(17)	
C(14)	-C(15)	-H(15B)	109.9(11)	C(85) -C(84)	-H(84B)	110.8(18)	
H(15A)	-C(15)	-H(15B)	108.2(15)	H(84A) -C(84)	-H(84B)	109(2)	
N(6)	-C(61)	-H(61A)	108.9(14)	C(81) -C(85)	-H(85A)	110.6(13)	
N(6)	-C(61)	-H(61B)	108.9(17)	C(81) -C(85)	-H(85B)	110.6(13)	
C(62)	-C(61)	-H(61A)	108.9(19)	C(84) -C(85)	-H(85A)	110.6(14)	
C(62)	-C(61)	-H(61B)	108.8(15)	C(84) -C(85)	-H(85B)	110.5(14)	
H(61A)	-C(61)	-H(61B)	107.7(17)	H(85A) -C(85)	-H(85B)	108.8(16)	
C(61)	-C(62)	-H(62A)	110(7)	C(32) -O(33)	-C(34)	111.7(13)	
C(61)	-C(62)	-H(62B)	109(6)	O(33) -C(32)	-C(31)	112.6(14)	
C(61)	-C(62)	-H(62C)	109(8)	O(33) -C(34)	-C(35)	113.5(17)	
H(62A)	-C(62)	-H(62B)	109(12)	C(32) -C(31)	-H(31A)	109(5)	
H(62A)	-C(62)	-H(62C)	109(10)	C(32) -C(31)	-H(31B)	110(3)	
H(62B)	-C(62)	-H(62C)	109(13)	C(32) -C(31)	-H(31C)	109(5)	
N(8)	-C(81)	-H(81)	110.5(11)	H(31A) -C(31)	-H(31B)	110(8)	
C(82)	-C(81)	-H(81)	110.4(11)	H(31A) -C(31)	-H(31C)	109(7)	
C(85)	-C(81)	-H(81)	110.4(13)	H(31B) -C(31)	-H(31C)	110(7)	
C(81)	-C(82)	-H(82A)	110.2(13)	O(33) -C(32)	-H(32A)	109.2(19)	
C(81)	-C(82)	-H(82B)	110.2(13)	O(33) -C(32)	-H(32B)	109.0(18)	
C(83)	-C(82)	-H(82A)	110.2(15)	C(31) -C(32)	-H(32A)	109.0(19)	
C(83)	-C(82)	-H(82B)	110.0(16)	C(31) -C(32)	-H(32B)	109(2)	

Table S6 - Bond Angles for: s962a			(Degrees)	(continued)			
			C17 H26 N6	O4 * C4 H10 O1			
H(32A)	-C(32)	-H(32B)	108(2)	C(34) -C(35)	-H(35A)	110(3)	
O(33)	-C(34)	-H(34A)	109.2(19)	C(34) -C(35)	-H(35B)	109(4)	
O(33)	-C(34)	-H(34B)	108.8(17)	C(34) -C(35)	-H(35C)	110(4)	
C(35)	-C(34)	-H(34A)	109(2)	H(35A) -C(35)	-H(35B)	109(13)	
C(35)	-C(34)	-H(34B)	108(2)	H(35A) -C(35)	-H(35C)	109(11)	
H(34A)	-C(34)	-H(34B)	108(2)	H(35B) -C(35)	-H(35C)	110(10)	

Table S7 - Torsion Angles (Degrees)
for: s962a C17 H26 N6 O4 * C4 H10 O1

C(14)	-O(1)	-C(11)	-C(12)	-25.9(9)
C(11)	-O(1)	-C(14)	-C(15)	127.6(9)
C(11)	-O(1)	-C(14)	-C(13)	2.0(10)
C(14)	-O(1)	-C(11)	-N(9)	-151.8(8)
C(2)	-N(1)	-C(6)	-N(6)	-178.3(10)
C(2)	-N(1)	-C(6)	-C(5)	0.8(17)
C(6)	-N(1)	-C(2)	-N(3)	-2.1(19)
C(4)	-N(3)	-C(2)	-N(1)	3.2(18)
C(2)	-N(3)	-C(4)	-N(9)	-176.7(10)
C(2)	-N(3)	-C(4)	-C(5)	-3.2(16)
C(61)	-N(6)	-C(6)	-C(5)	179.4(11)
C(6)	-N(6)	-C(61)	-C(62)	-145.6(14)
C(61)	-N(6)	-C(6)	-N(1)	-1.4(17)
C(8)	-N(7)	-C(5)	-C(6)	-178.8(13)
C(5)	-N(7)	-C(8)	-N(8)	178.6(9)
C(5)	-N(7)	-C(8)	-N(9)	-0.8(11)
C(8)	-N(7)	-C(5)	-C(4)	3.5(11)
C(81)	-N(8)	-C(8)	-N(7)	5.5(15)
C(81)	-N(8)	-C(8)	-N(9)	-175.0(9)
C(8)	-N(8)	-C(81)	-C(82)	-72.6(12)
C(8)	-N(8)	-C(81)	-C(85)	173.3(9)
C(8)	-N(9)	-C(4)	-C(5)	4.0(10)
C(11)	-N(9)	-C(4)	-N(3)	-12.6(15)
C(8)	-N(9)	-C(4)	-N(3)	178.6(10)
C(11)	-N(9)	-C(8)	-N(8)	11.1(16)
C(4)	-N(9)	-C(11)	-O(1)	-115.2(9)
C(4)	-N(9)	-C(11)	-C(12)	126.2(10)
C(8)	-N(9)	-C(11)	-O(1)	50.4(13)
C(8)	-N(9)	-C(11)	-C(12)	-68.3(13)

Table S7 - Torsion Angles (Degrees) (continued)
for: s962a C17 H26 N6 O4 * C4 H10 O1

C(11)	-N(9)	-C(4)	-C(5)	172.7(8)
C(4)	-N(9)	-C(8)	-N(8)	178.5(9)
C(11)	-N(9)	-C(8)	-N(7)	-169.4(9)
C(4)	-N(9)	-C(8)	-N(7)	-2.1(11)
N(3)	-C(4)	-C(5)	-N(7)	-179.4(10)
N(3)	-C(4)	-C(5)	-C(6)	2.3(16)
N(9)	-C(4)	-C(5)	-C(6)	177.0(9)
N(9)	-C(4)	-C(5)	-N(7)	-4.7(11)
N(7)	-C(5)	-C(6)	-N(6)	0(2)
C(4)	-C(5)	-C(6)	-N(1)	-0.9(17)
N(7)	-C(5)	-C(6)	-N(1)	-178.6(11)
C(4)	-C(5)	-C(6)	-N(6)	178.2(10)
O(1)	-C(11)	-C(12)	-C(13)	37.9(8)
O(1)	-C(11)	-C(12)	-O(12)	160.3(7)
N(9)	-C(11)	-C(12)	-O(12)	-80.2(10)
N(9)	-C(11)	-C(12)	-C(13)	157.4(8)
O(12)	-C(12)	-C(13)	-O(13)	-39.3(10)
C(11)	-C(12)	-C(13)	-C(14)	-35.4(9)
O(12)	-C(12)	-C(13)	-C(14)	-153.9(8)
C(11)	-C(12)	-C(13)	-O(13)	79.2(8)
O(13)	-C(13)	-C(14)	-O(1)	-92.9(8)
C(12)	-C(13)	-C(14)	-C(15)	-100.9(9)
O(13)	-C(13)	-C(14)	-C(15)	144.7(8)
C(12)	-C(13)	-C(14)	-O(1)	21.5(10)
O(1)	-C(14)	-C(15)	-O(15)	-70.5(11)
C(13)	-C(14)	-C(15)	-O(15)	50.4(11)
N(8)	-C(81)	-C(82)	-C(83)	-97.0(14)
C(85)	-C(81)	-C(82)	-C(83)	21.7(15)
N(8)	-C(81)	-C(85)	-C(84)	92.6(12)

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Table S7 - Torsion Angles (Degrees) (continued)
for: s962a C17 H26 N6 O4 * C4 H10 O1

C(82)	-C(81)	-C(85)	-C(84)	-26.9(12)
C(81)	-C(82)	-C(83)	-C(84)	-8(2)
C(82)	-C(83)	-C(84)	-C(85)	-9(2)
C(83)	-C(84)	-C(85)	-C(81)	22.3(16)
C(34)	-O(33)	-C(32)	-C(31)	-176.7(13)
C(32)	-O(33)	-C(34)	-C(35)	-169.5(17)

L.I.S.T.I.N.G O.F

O.B.S.E.R.V.E.D A.N.D C.A.L.C.U.L.A.T.E.D

S.T.R.U.C.T.U.R.E. F.A.C.T.O.R.S

F.O.R

S962A MOKA 60KV150MA 22C MON 260993

B E L O N G I N G T O T H E P A P E R

N⁶\$, C8-disubstituted adenosine derivatives as
partial agonists for adenosine A₁\$ receptors

b y

Harlof Roelen, Nora Veldman, Anthony L. Spek,
Jacobien von Frijtag Drabbe K\"unzel,
Ron A.A. Math\^ot and Ad P. IJzerman

Fc,Fo,s(Fo) * 10 for - S962A MOKA 60KV150MA 22C MON 260993 - Page

== H,	0,	0	==	6	155	162	17	-6	102	92	31	8	295	304	14
2	328	313	8	7	64	123	21	-5	67	14	37	9	136	86	40
4	904	948	8	8	55	89	28	-4	293	315	10	10	92	0	43
6	253	290	14	9	45	64	37	-3	299	299	9	11	113	142	23
8	148	137	32	10	11	0	43	-2	1071	966	5	== H,	4,	1	==
10	253	280	19	== H,	6,	0	==	-1	594	654	5	-10	70	76	41
12	140	148	32	0	83	69	31	0	190	191	6	-9	78	103	24
== H,	1,	0	==	1	75	104	19	1	594	662	5	-8	90	117	23
1	204	189	10	2	145	168	14	2	1071	1014	5	-7	73	70	36
2	299	254	9	3	68	91	23	3	299	280	10	-6	159	171	16
3	96	73	30	4	230	241	11	4	292	271	11	-5	159	137	17
4	541	544	8	5	110	102	25	5	67	79	33	-4	250	275	10
5	311	282	12	6	68	49	35	6	102	125	23	-3	192	186	11
6	321	323	12	7	77	79	32	7	245	245	16	-2	235	220	10
7	131	127	27	8	222	203	16	8	51	0	39	-1	273	270	9
8	56	0	39	9	5	55	36	9	35	78	44	0	382	376	5
9	107	104	42	== H,	7,	0	==	10	126	97	45	1	273	261	9
10	21	0	42	1	7	67	32	11	91	146	27	2	235	227	10
11	146	136	32	2	124	150	15	12	99	91	43	3	192	188	13
12	10	73	44	3	19	50	32	== H,	2,	1	==	4	250	248	11
== H,	2,	0	==	4	76	105	21	-11	87	86	42	5	159	116	22
0	1624	1610	5	5	2	0	42	-10	121	120	33	6	159	150	19
1	236	247	9	6	88	0	45	-9	81	128	26	7	73	0	40
2	89	159	13	7	129	96	26	-8	109	110	32	8	90	94	31
3	429	389	8	8	63	63	34	-7	94	126	24	9	78	99	27
4	448	401	8	== H,	8,	0	==	-6	417	399	11	10	70	71	41
5	226	212	14	0	117	114	19	-5	67	48	34	== H,	5,	1	==
6	538	546	9	1	212	224	12	-4	142	118	21	-10	52	105	25
7	106	109	30	2	160	148	17	-3	109	129	17	-9	149	120	26
8	229	233	16	3	49	29	34	-2	308	306	8	-8	100	114	24
9	26	102	35	4	119	113	21	-1	440	442	6	-7	193	179	16
10	21	77	42	5	18	58	33	0	870	913	4	-6	148	166	14
11	10	0	43	6	50	75	24	1	440	467	6	-5	153	170	15
== H,	3,	0	==	7	29	0	45	2	308	315	8	-4	61	69	29
1	84	100	18	== H,	9,	0	==	3	109	109	20	-3	345	335	9
2	2	49	30	1	34	0	42	4	142	139	17	-2	192	208	11
3	224	245	10	2	20	0	43	5	67	60	35	-1	151	139	13
4	419	379	9	3	48	69	28	6	417	418	10	0	108	80	24
5	60	76	28	4	24	0	44	7	94	24	39	1	151	144	15
6	155	151	18	5	82	93	23	8	109	78	41	2	192	209	12
7	98	84	34	== H,	10,	0	==	9	81	109	33	3	345	346	9
8	46	0	40	0	37	74	29	10	121	112	34	4	61	108	20
9	69	81	38	1	100	103	22	11	87	39	43	5	154	142	19
10	23	115	27	== H,	0,	1	==	== H,	3,	1	==	6	148	143	18
11	16	0	44	1	652	593	18	-11	113	147	23	7	193	175	16
== H,	4,	0	==	2	426	407	5	-10	92	139	20	8	100	46	39
0	64	83	19	3	438	448	10	-9	136	126	25	9	149	136	23
1	40	71	25	4	99	109	29	-8	295	298	14	10	52	88	32
2	110	127	15	5	297	300	19	-7	98	149	19	== H,	6,	1	==
3	47	61	32	6	123	132	16	-6	115	59	37	-9	37	0	45
4	252	252	11	7	235	234	12	-5	8	0	34	-8	71	91	31
5	27	0	34	8	25	0	33	-4	485	439	8	-7	45	0	42
6	120	113	24	9	4	73	37	-3	227	210	11	-6	154	152	16
7	75	31	39	10	97	123	26	-2	484	497	7	-5	181	153	18
8	110	120	24	11	118	134	25	-1	243	245	9	-4	70	0	39
9	22	0	42	12	34	83	43	0	449	461	10	-3	109	116	19
10	89	122	23	== H,	1,	1	==	1	243	244	9	-2	293	285	10
== H,	5,	0	==	-12	99	140	27	2	484	475	8	-1	23	31	30
1	275	286	9	-11	91	98	42	3	227	223	11	0	8	63	20
2	252	215	11	-10	126	119	34	4	485	506	8	1	23	42	30
3	34	0	37	-9	35	91	43	5	8	59	33	2	293	284	10
4	147	169	16	-8	51	48	43	6	115	76	36	3	109	95	26
5	101	138	17	-7	245	255	16	7	98	117	24	4	70	102	20

5	181	188	15	7	101	113	39	-5	101	122	18	-8	131	136	19
6	154	120	23	8	100	76	38	-4	448	408	9	-7	150	145	18
7	45	0	44	9	25	103	30	-3	347	384	9	-6	94	127	20
8	71	55	36	10	228	249	14	-2	307	301	9	-5	75	74	27
9	37	58	37	11	75	93	35	-1	178	175	12	-4	244	230	11
= H, 7,	1	==	12	149	178	21	0	322	312	7	-3	164	158	16	
-8	53	60	35	== H, 1,	2	==	1	178	185	11	-2	250	225	12	
-7	28	0	45	-12	55	69	45	2	307	326	9	-1	141	145	14
-6	69	88	26	-11	35	0	44	3	347	380	8	0	156	162	10
-5	117	145	16	-10	250	220	23	4	448	452	8	1	141	159	14
-4	302	304	11	-9	147	131	30	5	101	73	33	2	250	227	12
-3	148	127	20	-8	118	43	44	6	251	278	12	3	164	171	15
-2	40	65	30	-7	40	79	39	7	134	120	25	4	243	234	13
-1	79	0	42	-6	104	78	36	8	102	85	37	5	75	65	38
0	172	189	21	-5	36	0	33	9	142	116	29	6	94	120	21
1	79	87	25	-4	87	85	28	10	52	85	40	7	150	133	21
2	40	0	42	-3	165	152	16	11	34	61	41	8	131	91	30
3	147	126	19	-2	538	536	6	== H, 4,	2	==	9	77	116	23	
4	302	296	11	-1	262	295	8	-10	117	125	23	== H, 7,	2	==	
5	117	159	16	0	7	0	23	-9	57	0	44	-8	49	48	36
6	69	64	38	1	262	236	9	-8	51	71	35	-7	16	64	35
7	28	49	35	2	538	542	6	-7	136	160	16	-6	173	156	15
8	53	73	29	3	165	185	13	-6	82	59	35	-5	115	82	28
= H, 8,	1	==	4	87	107	22	-5	88	133	15	-4	59	56	33	
-7	28	0	46	5	36	0	33	-4	78	84	24	-3	186	186	14
-6	25	0	44	6	104	119	23	-3	247	266	10	-2	104	118	17
-5	34	29	32	7	41	88	35	-2	49	0	34	-1	129	133	15
-4	32	0	43	8	118	139	26	-1	127	160	11	0	98	98	14
-3	81	92	25	9	147	136	32	0	328	329	6	1	129	115	19
-2	188	182	14	10	250	246	20	1	127	140	15	2	104	86	25
-1	123	149	15	11	35	24	46	2	49	64	30	3	186	189	13
0	100	98	17	12	55	55	46	3	247	266	11	4	59	0	43
1	123	177	12	== H, 2,	2	==	4	78	0	38	5	115	143	17	
2	188	196	12	-11	38	92	38	5	88	125	17	6	173	183	16
3	81	96	24	-10	75	82	43	6	82	90	29	7	16	0	46
4	32	0	43	-9	156	132	29	7	136	173	16	8	49	0	45
5	34	24	32	-8	154	142	25	8	51	64	37	== H, 8,	2	==	
6	25	54	34	-7	133	153	20	9	57	0	42	-7	36	79	24
7	28	0	47	-6	68	124	22	10	117	103	31	-6	11	0	43
= H, 9,	1	==	5	538	526	9	== H, 5,	2	==	-5	75	87	24		
-5	14	89	22	-4	486	454	8	-10	62	80	33	-4	79	0	44
-4	11	0	41	-3	726	678	7	-9	92	111	23	-3	138	137	17
-3	6	72	26	-2	466	479	7	-8	90	131	20	-2	91	108	17
-2	10	63	31	-1	755	799	6	-7	47	58	34	-1	168	154	15
-1	38	0	43	0	539	549	4	-6	149	120	22	0	121	129	12
0	159	157	11	1	755	778	6	-5	109	108	22	1	168	163	15
1	38	0	45	2	466	487	7	-4	69	87	22	2	91	106	22
2	10	102	21	3	726	726	7	-3	108	102	22	3	138	111	22
3	6	0	46	4	486	466	8	-2	196	204	11	4	79	71	31
4	11	0	43	5	538	534	9	-1	174	133	15	5	76	85	27
5	14	0	46	6	68	100	28	0	422	399	6	6	11	84	28
= H, 10,	1	==	7	133	139	23	1	174	172	13	7	36	0	48	
-1	12	36	32	8	154	164	22	2	196	223	10	== H, 9,	2	==	
0	35	70	20	9	156	167	25	3	108	105	21	-5	79	54	33
1	13	0	44	10	75	31	43	4	69	104	23	-4	40	0	43
= H, 0,	2	==	11	38	63	43	5	109	122	21	-3	97	95	23	
0	1042	1024	17	== H, 3,	2	==	6	149	156	17	-2	97	98	24	
1	1014	996	4	-11	34	98	33	7	47	0	40	-1	53	71	32
2	366	351	13	-10	52	19	41	8	90	64	37	0	81	81	26
3	1167	1136	28	-9	142	121	27	9	92	107	27	1	53	77	28
4	328	327	8	-8	102	74	39	10	62	73	37	2	97	105	21
5	144	115	19	-7	134	125	22	== H, 6,	2	==	3	98	98	25	
6	58	47	30	-6	252	256	13	-9	77	89	32	4	40	94	19

5	79	0	47	11	47	49	42	5	69	64	37	-3	8	78	25
== H,	10,	2	==	== H,	3,	3	==	6	91	92	27	-2	52	83	25
0	30	42	31	-11	77	103	31	7	143	175	15	-1	63	56	31
== H,	0,	3	==	-10	104	76	39	8	89	0	43	0	61	0	37
1	1034	1022	16	-9	42	39	39	9	81	99	31	1	63	0	44
2	926	917	22	-8	180	157	20	10	25	0	44	2	52	0	46
3	296	262	28	-7	68	29	36	== H,	6,	3	==	3	8	89	22
4	34	0	28	-6	105	36	36	-9	51	0	43	4	36	34	33
5	334	357	8	-5	214	183	14	-8	24	51	34	5	28	0	46
6	284	285	11	-4	298	275	10	-7	123	85	29	== H,	0,	4	==
7	79	0	31	-3	146	125	18	-6	217	205	14	0	549	379	61
8	68	86	32	-2	278	298	8	-5	89	133	16	1	6	0	25
9	0	80	36	-1	315	318	9	-4	202	189	13	2	754	723	13
10	81	41	37	0	792	816	4	-3	218	237	12	3	90	94	19
11	146	183	19	1	315	311	9	-2	54	31	32	4	285	311	8
12	82	80	42	2	278	284	10	-1	28	0	37	5	161	175	30
== H,	1,	3	==	3	146	131	17	0	30	0	31	6	154	151	15
-11	72	118	34	4	298	326	10	1	28	0	39	7	96	107	34
-10	41	0	42	5	214	201	13	2	54	86	24	8	249	284	12
-9	132	151	27	6	105	118	21	3	218	238	12	9	336	362	11
-8	86	149	23	7	68	108	26	4	202	216	14	10	16	78	40
-7	262	280	13	8	180	108	33	5	89	86	31	11	28	0	36
-6	98	57	37	9	42	0	43	6	217	245	12	== H,	1,	4	==
-5	249	281	12	10	104	62	41	7	123	135	19	-11	127	112	38
-4	299	292	10	11	77	106	31	8	24	51	35	-10	144	149	26
-3	353	334	9	== H,	4,	3	==	9	51	42	35	-9	180	162	27
-2	438	433	7	-10	27	35	38	== H,	7,	3	==	-8	142	86	43
-1	624	616	6	-9	62	92	33	-8	113	106	20	-7	198	168	20
0	475	505	5	-8	190	180	16	-7	85	76	29	-6	10	0	36
1	625	597	6	-7	107	81	33	-6	40	32	34	-5	374	371	10
2	437	433	7	-6	88	98	25	-5	175	160	16	-4	107	85	27
3	353	329	9	-5	80	73	27	-4	69	0	41	-3	140	132	18
4	299	275	11	-4	287	305	10	-3	95	120	18	-2	447	419	8
5	249	276	11	-3	174	164	14	-2	72	92	23	-1	625	636	6
6	98	95	29	-2	410	436	8	-1	50	0	39	0	666	679	4
7	262	284	14	-1	306	314	8	0	3	64	25	1	625	585	6
8	86	159	20	0	306	300	6	1	50	0	41	2	447	446	7
9	132	113	37	1	306	282	10	2	72	105	21	3	140	148	16
10	41	121	30	2	410	444	8	3	95	0	44	4	107	111	22
11	72	159	26	3	174	186	13	4	69	61	35	5	373	398	10
== H,	2,	3	==	4	287	314	10	5	175	162	17	6	10	75	38
-11	47	0	44	5	80	100	24	6	40	61	34	7	198	190	19
-10	183	152	26	6	88	113	24	7	85	75	35	8	142	147	20
-9	52	0	41	7	107	42	38	8	113	71	36	9	180	193	20
-8	157	138	27	8	190	204	15	== H,	8,	3	==	10	144	179	23
-7	160	147	22	9	62	0	44	-7	39	0	39	11	127	154	23
-6	124	86	32	10	27	88	33	-6	28	17	34	== H,	2,	4	==
-5	187	150	18	== H,	5,	3	==	-5	71	60	36	-11	65	44	44
-4	408	418	8	-10	25	0	43	-4	38	17	31	-10	110	114	30
-3	316	306	9	-9	81	63	35	-3	48	71	28	-9	45	0	43
-2	529	562	7	-8	89	103	26	-2	95	114	17	-8	28	0	36
-1	593	625	6	-7	143	108	23	-1	132	133	17	-7	109	98	30
0	194	200	11	-6	91	82	25	0	194	190	9	-6	333	333	11
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= H,	6,	9	==	4	263	232	33	-2	191	176	14	-1	121	84	29
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= H,	7,	9	==	-2	748	714	8	-6	22	0	39	1	66	0	46
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= H,	8,	9	==	-9	50	105	28	10	55	41	39	3	42	49	35
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9	50	125	31	9	41	89	37	6	190	218	14	6	161	166	24
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== H,	2,	14	==	-8	17	100	29	8	142	131	23	8	77	0	45
-10	55	126	28	-7	48	42	38	9	15	92	45	9	31	0	46
-9	42	0	43	-6	75	91	32	10	103	99	50	== H,	4,	15	==
-8	60	56	43	-5	87	103	28	== H,	1,	15	==	-8	93	112	30
-7	148	133	28	-4	129	122	21	-10	83	147	23	-7	135	124	28
-6	48	75	41	-3	77	39	34	-9	83	0	43	-6	130	104	34
-5	70	0	41	-2	26	108	17	-8	101	101	40	-5	88	88	35
-4	31	42	41	-1	69	111	19	-7	122	155	27	-4	196	211	16
-3	231	233	16	0	17	0	33	-6	130	145	28	-3	51	36	35
-2	44	0	36	1	69	52	37	-5	94	104	36	-2	68	0	38
-1	78	112	24	2	26	24	37	-4	287	307	15	-1	40	66	33
0	118	75	27	3	77	0	44	-3	119	53	44	0	161	159	13
1	78	0	41	4	129	116	26	-2	268	210	24	1	40	43	38
2	44	57	38	5	87	44	39	-1	73	0	41	2	68	73	38
3	231	205	18	6	75	46	38	0	303	336	92	3	51	124	22
4	31	0	40	7	48	70	38	1	73	0	44	4	196	196	18
5	70	61	41	8	17	35	38	2	268	304	15	5	88	0	46
6	48	122	26	== H,	6,	14	==	3	119	195	19	6	130	115	28
7	148	154	23	-7	11	52	38	4	287	238	17	7	135	158	22
8	60	95	36	-6	20	0	41	5	94	153	22	8	93	108	34
9	42	0	46	-5	89	115	21	6	130	106	37	== H,	5,	15	==
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== H,	3,	14	==	-3	24	35	32	8	101	0	43	-6	158	150	20
-9	25	0	41	-2	67	57	33	9	84	84	43	-5	103	108	25
-8	37	0	42	-1	138	157	15	10	84	78	47	-4	109	118	22
-7	116	47	43	0	69	62	26	== H,	2,	15	==	-3	165	174	16
-6	64	108	29	1	138	168	17	-9	51	0	43	-2	149	148	16
-5	80	0	39	2	67	95	28	-8	84	31	42	-1	179	145	20
-4	60	105	29	3	24	0	45	-7	112	145	26	0	58	72	23
-3	211	189	18	4	54	86	29	-6	41	82	39	1	179	187	15
-2	43	0	37	5	88	36	37	-5	62	0	41	2	148	140	21
-1	101	96	24	6	20	0	47	-4	45	0	39	3	165	153	19
0	227	226	10	7	12	75	37	-3	120	137	27	4	109	137	20
1	101	36	38	== H,	7,	14	==	-2	114	24	42	5	103	132	22
2	43	40	37	-5	43	0	42	-1	163	174	16	6	158	119	28
3	211	212	17	-4	33	0	41	0	243	263	9	7	56	56	40
4	60	100	30	-3	83	58	33	1	163	119	27	== H,	6,	15	==
5	80	83	38	-2	54	26	32	2	114	124	29	-6	43	0	40
6	64	99	34	-1	92	93	26	3	120	142	25	-5	21	0	42
7	116	125	27	0	191	178	12	4	45	0	42	-4	93	51	34
8	37	77	41	1	92	105	25	5	62	38	41	-3	94	85	29
9	25	0	44	2	54	62	35	6	41	61	41	-2	128	118	21
== H,	4,	14	==	3	83	83	30	7	111	104	34	-1	64	114	19
-9	41	0	44	4	33	61	35	8	84	42	43	0	102	106	18
-8	72	0	44	5	43	82	30	9	51	93	39	1	64	54	36
-7	23	140	22	== H,	8,	14	==	== H,	3,	15	==	2	128	113	24
-6	46	0	41	-4	59	48	30	-9	31	0	45	3	94	99	26
-5	90	43	41	-3	15	44	33	-8	77	64	40	4	93	26	39
-4	97	99	27	-2	129	149	17	-7	92	106	32	5	21	0	47
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1	231	231	15	3	15	61	35	-2	101	60	36	-2	101	95	24
2	37	107	22	4	59	81	30	-1	63	48	35	-1	119	116	21
3	175	155	19	== H,	0,	15	==	0	90	38	32	0	120	124	16

1	118	120	22	== H,	3,	16	==	2	29	0	45	1	83	76	41
2	101	106	25	-9	23	138	22	3	100	119	21	2	46	0	45
3	36	17	36	-8	59	70	43	4	50	50	37	3	149	184	22
4	15	0	46	-7	63	43	42	5	24	74	36	4	160	115	33
H, 8,	15	==	-6	65	95	33		6	33	0	48	5	124	0	44
-3	89	84	24	-5	70	0	41	== H,	7,	16	==	6	39	0	43
-2	87	116	19	-4	102	107	34	-4	79	19	34	7	55	85	39
-1	3	0	41	-3	78	42	40	-3	96	63	31	8	75	110	33
0	8	74	31	-2	139	103	29	-2	86	88	24	9	63	0	46
1	3	94	27	-1	34	0	38	-1	43	41	34	== H,	3,	17	==
2	87	0	48	0	98	81	24	0	57	64	28	-8	18	0	43
3	89	78	30	1	34	0	42	1	43	71	36	-7	93	140	24
H, 0,	16	==	2	140	144	23		2	86	108	24	-6	62	53	41
0	80	63	34	3	78	93	32	3	96	120	22	-5	35	36	41
1	9	44	36	4	102	131	25	4	79	81	30	-4	167	177	21
2	113	138	46	5	70	113	30	== H,	8,	16	==	-3	47	0	40
3	106	59	35	6	65	76	40	-3	49	19	31	-2	58	85	32
4	96	81	34	7	63	19	43	-2	41	57	31	-1	68	95	25
5	111	103	26	8	59	98	37	-1	35	38	34	0	251	242	10
6	16	43	36	9	23	43	41	0	101	113	37	1	68	0	43
7	76	109	26	== H,	4,	16	==	1	35	0	47	2	58	75	39
8	148	187	16	-8	74	106	34	2	41	74	31	3	47	0	44
9	50	77	37	-7	92	95	39	== H,	0,	17	==	4	167	167	22
H, 1,	16	==	-6	71	56	40		1	102	105	26	5	35	58	41
-9	8	0	43	-5	88	78	37	2	113	155	42	6	62	92	37
-8	37	46	44	-4	38	50	38	3	25	69	35	7	93	0	45
-7	97	110	34	-3	35	88	28	4	161	165	20	8	18	61	40
-6	100	131	29	-2	155	169	16	5	31	72	38	== H,	4,	17	==
-5	85	145	27	-1	98	142	19	6	40	39	37	-7	69	106	29
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-2	32	0	48	2	155	179	17	9	49	17	38	-4	153	146	22
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0	302	280	90	4	38	0	44	-9	58	0	44	-2	63	44	36
1	171	170	21	5	88	68	40	-8	50	116	33	-1	45	0	39
2	32	147	27	6	71	92	36	-7	109	99	44	0	222	203	11
3	148	220	16	7	92	139	23	-6	157	168	26	1	45	0	43
4	242	261	16	8	74	0	49	-5	36	64	44	2	63	105	26
5	85	118	32	== H,	5,	16	==	-4	137	165	25	3	47	0	43
6	100	92	41	-7	27	31	39	-3	109	60	45	4	153	168	19
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9	8	0	43	-4	144	164	16	0	237	220	80	7	69	36	40
H, 2,	16	==	-3	120	142	17		1	30	0	47	== H,	5,	17	==
-9	52	76	41	-2	82	75	32	2	120	125	30	-7	21	0	43
-8	63	122	31	-1	39	83	26	3	109	140	28	-6	40	0	42
-7	105	109	35	0	88	60	31	4	137	148	28	-5	52	76	36
-6	118	102	40	1	39	17	37	5	36	0	43	-4	40	34	36
-5	53	87	37	2	82	69	37	6	157	176	23	-3	36	0	42
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-3	238	249	15	4	144	132	23	8	50	0	44	-1	83	54	36
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-1	87	47	40	6	103	106	30	== H,	2,	17	==	1	84	120	24
0	78	64	31	7	28	69	39	-9	63	0	45	2	187	138	26
1	87	29	41	== H,	6,	16	==	-8	75	0	44	3	36	80	33
2	220	206	19	-6	33	0	42	-7	55	86	43	4	40	0	45
3	238	211	19	-5	24	0	42	-6	39	125	28	5	52	57	37
4	70	0	43	-4	50	32	34	-5	124	31	45	6	40	0	47
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6	118	166	22	-2	29	80	26	-3	149	87	46	== H,	6,	17	==
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9	52	0	44	1	45	0	45	0	46	114	17	-3	22	52	33

-2	26	0	40	5	119	127	28	-1	9	0	39	7	36	61	42
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0	96	90	24	7	76	104	34	1	9	0	47	-7	28	0	44
1	114	137	20	8	88	78	42	== H,	0,	19	==	-6	53	0	46
2	26	40	35	== H,	3,	18	==	1	202	170	21	-5	25	57	39
3	22	19	35	-8	34	96	32	2	107	92	48	-4	94	122	27
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== H,	7,	17	==	-5	39	38	43	5	85	101	52	-1	152	143	21
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0	33	73	27	-1	81	76	35	== H,	1,	19	==	3	112	81	41
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== H,	8,	17	==	3	159	121	30	-5	168	150	35	7	28	0	46
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1	68	71	35	6	44	0	46	-2	39	29	44	-4	46	0	43
== H,	0,	18	==	7	31	96	35	-1	97	160	24	-3	74	65	36
0	55	0	35	8	34	0	45	0	75	0	40	-2	112	96	30
1	49	84	44	== H,	4,	18	==	1	97	118	36	-1	39	0	41
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3	170	193	14	-6	149	163	23	3	45	132	28	1	39	73	38
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5	184	190	20	-4	42	86	34	5	168	133	34	3	74	122	21
6	16	38	37	-3	80	91	34	6	43	89	47	4	46	50	38
7	158	184	18	-2	204	153	21	7	33	81	44	5	63	71	39
8	55	0	37	-1	58	94	28	8	39	70	45	== H,	6,	19	==
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-1	152	124	31	-6	79	99	35	0	158	180	14	== H,	0,	20	==
0	256	249	34	-5	60	29	40	1	116	124	29	0	268	232	31
1	152	128	31	-4	47	81	34	2	113	96	43	1	42	0	38
2	55	99	36	-3	62	0	43	3	89	120	33	2	76	69	37
3	162	139	27	-2	36	38	36	4	184	204	22	3	148	153	21
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== H,	2,	18	==	4	47	0	47	-7	36	47	44	== H,	1,	20	==
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1	112	134	24	2	24	93	27	3	210	203	19	1	100	98	42
2	189	170	21	3	56	54	35	4	119	116	31	2	31	59	43
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4	145	138	30	== H,	7,	18	==	6	110	133	27	4	18	0	45

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6	149	156	26	-2	44	48	34	-3	14	0	42	-6	20	0	44	
7	7	0	44	-1	45	48	35	-2	111	119	26	-5	87	40	44	
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H,	2,	20	==	1	45	0	46	0	40	58	33	-3	36	76	43	
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-1	104	108	35	4	165	130	52	-4	65	79	36	4	67	66	42	
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3	126	184	23	==	H,	1,	21	==	0	40	44	35	-5	25	129	24
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H,	3,	20	==	-3	66	64	47	==	H,	6,	21	==	0	91	87	23
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-4	64	61	41	1	21	0	45	==	H,	0,	22	==	4	27	68	42
-3	39	66	39	2	105	79	45	0	8	0	38	5	25	51	41	
-2	117	97	32	3	67	0	47	1	94	95	35	==	H,	5,	22	==
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1	135	151	23	6	114	109	43	4	54	0	37	-1	114	96	30	
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H,	4,	20	==	-3	126	153	25	-5	36	0	43	2	69	0	39	
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0	85	91	20	4	165	181	23	2	44	0	44	-5	37	0	43	
1	118	56	40	5	98	129	31	3	43	0	44	-4	45	34	46	
2	88	125	22	6	60	22	44	4	75	150	25	-3	48	62	46	
3	46	0	45	7	53	0	48	5	36	0	44	-2	66	122	30	
4	60	0	45	==	H,	3,	21	==	6	30	101	43	-1	102	46	45
5	90	113	30	-6	43	24	43	7	93	105	42	0	128	124	24	
6	70	0	46	-5	65	32	45	==	H,	2,	22	==	1	102	138	28
H,	5,	20	==	-4	134	145	25	-6	65	88	49	2	66	91	44	
-5	57	0	43	-3	114	117	31	-5	71	0	44	3	48	0	45	
-4	71	71	38	-2	53	86	35	-4	94	73	45	4	45	64	46	
-3	14	40	37	-1	42	0	42	-3	18	50	48	5	37	0	46	
-2	28	34	36	0	82	88	27	-2	128	124	29	6	47	71	45	
-1	82	139	17	1	42	0	45	-1	61	100	39	==	H,	2,	23	==
0	164	164	14	2	53	48	41	0	33	81	36	-5	91	154	28	
1	82	64	39	3	114	111	31	1	61	0	46	-4	38	0	44	
2	28	56	37	4	134	134	26	2	128	108	35	-3	156	172	23	
3	13	0	46	5	65	92	37	3	18	91	44	-2	40	46	45	
4	71	0	48	6	43	103	33	4	94	105	39	-1	126	119	31	
5	57	40	40	==	H,	4,	21	==	5	71	103	42	0	79	85	28
H,	6,	20	==	-5	87	81	43	6	65	32	46	1	126	97	44	

2	40	22	45	-3	33	0	43	2	32	0	46	
3	156	167	24	-2	56	46	42	==	H,	0,	27	==
4	38	116	33	-1	29	0	43	1	18	0	37	
5	91	78	48	0	10	41	37	==	H,	1,	27	==
==	H,	3,	23	==	1	29	0	46	0	33	69	41
-5	63	86	44	2	56	0	48					
-4	71	46	43	3	33	29	43					
-3	22	44	41	4	33	82	43					
-2	55	91	34	==	H,	4,	24	==				
-1	56	97	33	-2	26	0	42					
0	55	16	36	-1	52	65	40					
1	56	0	46	0	38	73	32					
2	54	0	46	1	52	0	47					
3	22	0	48	2	26	0	47					
4	71	67	43	==	H,	0,	25	==				
5	63	44	43	1	93	25	39					
==	H,	4,	23	==	2	82	124	39				
-4	31	0	44	3	22	0	38					
-3	38	76	40	4	12	83	39					
-2	14	0	44	==	H,	1,	25	==				
-1	30	0	41	-4	65	0	47					
0	2	51	33	-3	2	86	47					
1	30	60	39	-2	49	100	41					
2	14	72	41	-1	23	0	42					
3	38	49	40	0	50	65	39					
4	31	0	48	1	23	0	47					
==	H,	5,	23	==	2	49	0	48				
-1	53	0	42	3	2	0	47					
0	96	96	22	4	65	103	43					
1	53	71	39	==	H,	2,	25	==				
==	H,	0,	24	==	-4	73	35	46				
0	92	96	44	-3	44	116	34					
1	56	86	44	-2	23	82	44					
2	58	83	37	-1	50	0	43					
3	103	76	39	0	82	59	38					
4	12	55	39	1	50	81	44					
5	41	59	41	2	23	98	39					
==	H,	1,	24	==	3	44	49	45				
-5	27	26	45	4	72	57	47					
-4	57	35	46	==	H,	3,	25	==				
-3	58	106	41	-2	23	0	43					
-2	69	41	46	-1	28	0	41					
-1	10	32	45	0	9	80	29					
0	57	58	39	1	28	84	41					
1	10	0	47	2	23	36	41					
2	69	0	46	==	H,	0,	26	==				
3	58	123	33	0	8	53	39					
4	57	103	39	1	62	61	40					
5	27	0	46	2	36	64	39					
==	H,	2,	24	==	3	4	70	39				
-5	99	129	34	==	H,	1,	26	==				
-4	71	0	43	-3	50	0	44					
-3	37	103	40	-2	28	0	45					
-2	80	87	44	-1	50	0	43					
-1	53	72	46	0	12	108	28					
0	16	85	37	1	50	125	33					
1	53	0	47	2	28	83	45					
2	80	61	44	3	50	78	44					
3	37	82	46	==	H,	2,	26	==				
4	71	0	48	-2	32	98	39					
5	99	73	46	-1	24	0	44					
==	H,	3,	24	==	0	17	106	25				
-4	34	24	43	1	24	32	46					