



# Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 12:23 AM GMT

PDB ID : 4A5Y  
Title : Intermediate state of human kinesin Eg5 in complex with Ispinesib  
Authors : Kaan, H.Y.K.; Kozielski, F.  
Deposited on : 2011-10-29  
Resolution : 2.45 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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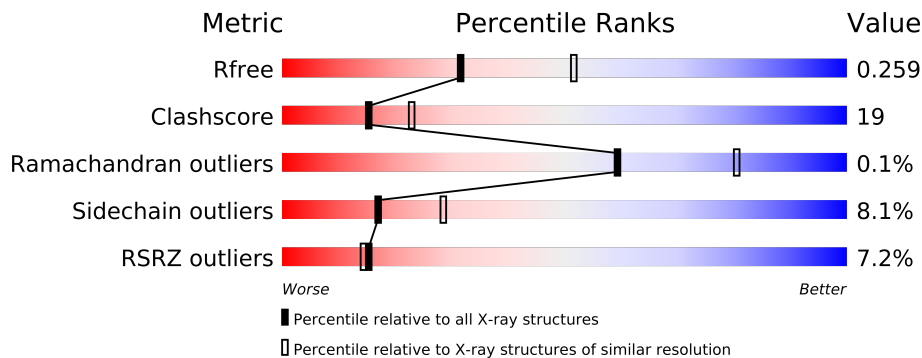
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.45 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	3566 (2.50-2.42)
Clashscore	79885	4471 (2.50-2.42)
Ramachandran outliers	78287	4383 (2.50-2.42)
Sidechain outliers	78261	4385 (2.50-2.42)
RSRZ outliers	66119	3568 (2.50-2.42)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	368	
1	B	368	
1	C	368	

## 2 Entry composition i

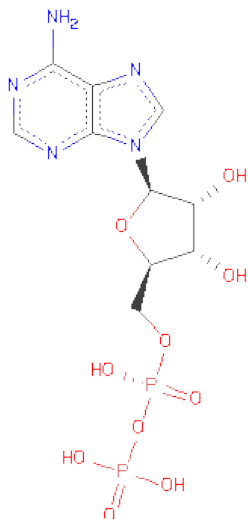
There are 7 unique types of molecules in this entry. The entry contains 8127 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called KINESIN-LIKE PROTEIN KIF11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	4	0
			2598	1631	448	509	10			
1	B	302	Total	C	N	O	S	0	0	0
			2364	1489	404	461	10			
1	C	339	Total	C	N	O	S	0	1	0
			2618	1643	458	507	10			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).

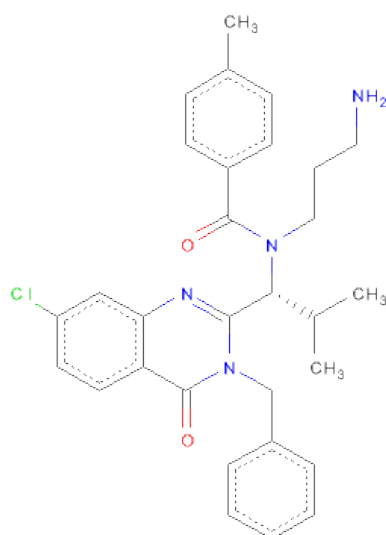


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	C	1	Total	Cl	0	0
			1	1		

- Molecule 4 is ISPINESIB MESILATE (three-letter code: G7X) (formula: C<sub>30</sub>H<sub>33</sub>ClN<sub>4</sub>O<sub>2</sub>).

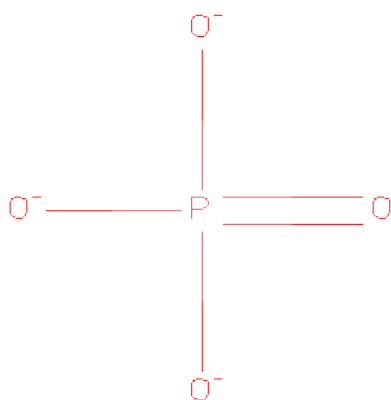


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	0	0
			37	30	1	4	2		
4	B	1	Total	C	Cl	N	O	0	0
			37	30	1	4	2		
4	C	1	Total	C	Cl	N	O	0	0
			37	30	1	4	2		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Mg	0	0
			1	1		

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	O	P	0	0
			5	4	1		
6	C	1	Total	O	P	0	0
			5	4	1		

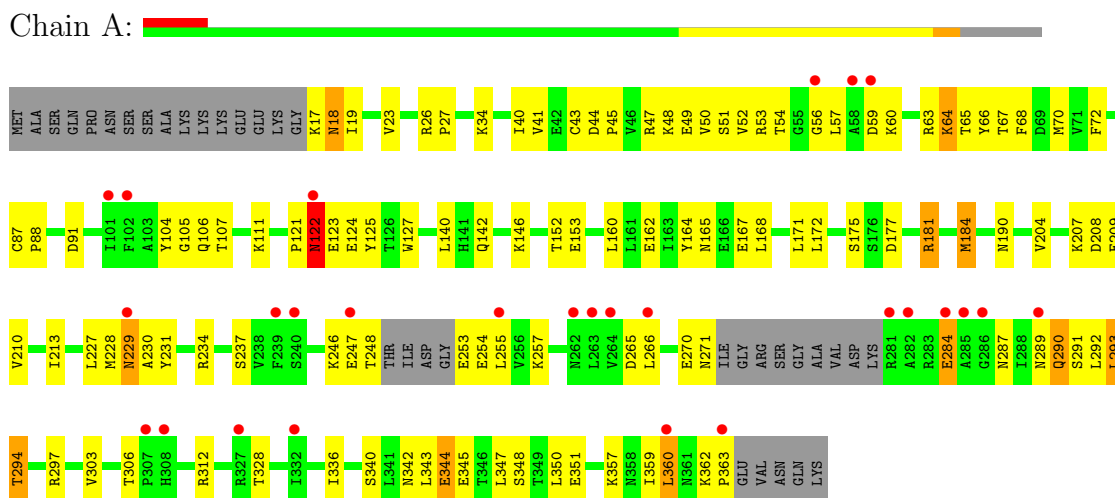
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	93	Total	O	0	0
			93	93		
7	B	81	Total	O	0	0
			81	81		
7	C	168	Total	O	0	0
			168	168		

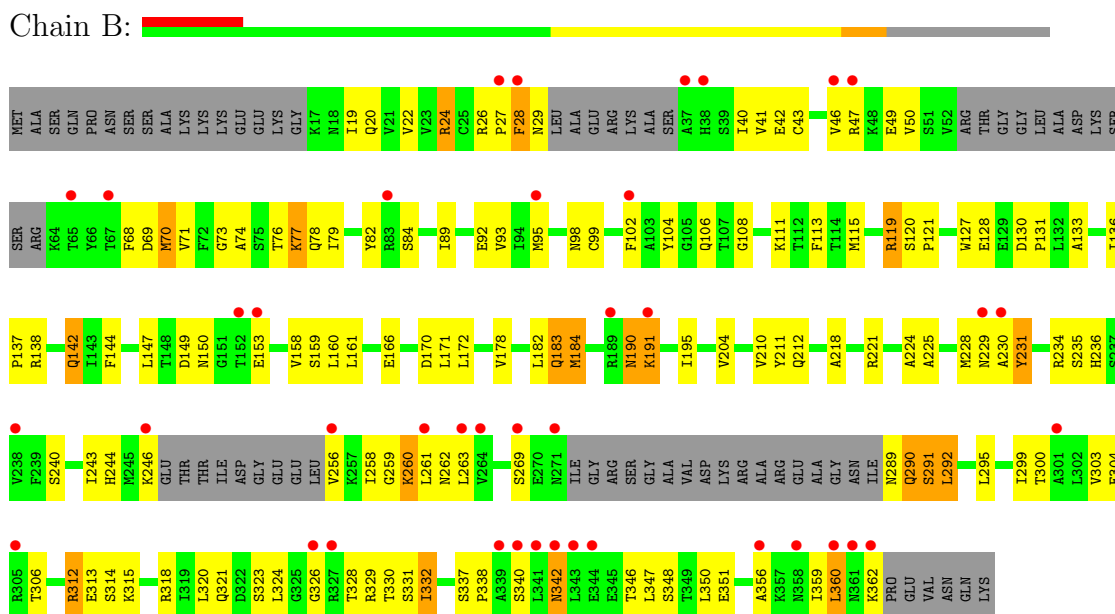
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: KINESIN-LIKE PROTEIN KIF11

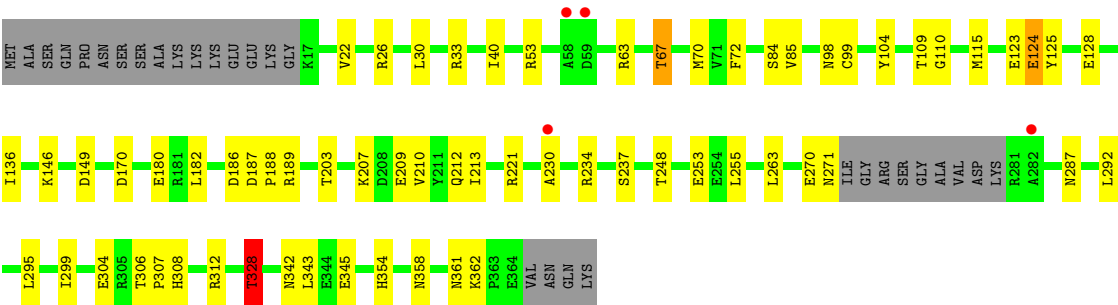


#### • Molecule 1: KINESIN-LIKE PROTEIN KIF11



#### • Molecule 1: KINESIN-LIKE PROTEIN KIF11





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.63Å 91.93Å 163.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.60 – 2.45 29.60 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.60-2.45) 99.6 (29.60-2.45)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.74 (at 2.45Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.201 , 0.268 0.195 , 0.259	Depositor DCC
$R_{free}$ test set	2428 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	45.3	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 45.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48117 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8127	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.27 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.7156e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: G7X, MG, PO4, ADP, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.67	0/2648	0.81	0/3586
1	B	0.58	0/2399	0.71	1/3245 (0.0%)
1	C	0.78	0/2661	0.81	2/3603 (0.1%)
All	All	0.68	0/7708	0.78	3/10434 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	328	THR	CB-CA-C	-5.59	96.50	111.60
1	C	110	GLY	N-CA-C	5.50	126.84	113.10
1	B	326	GLY	N-CA-C	-5.03	100.52	113.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121	PRO	Peptide
1	A	122	ASN	Peptide

## 5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2598	0	2588	102	0
1	B	2364	0	2363	125	0
1	C	2618	0	2612	59	0
2	A	27	0	12	2	0
2	B	27	0	12	2	0
2	C	27	0	12	0	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	37	0	33	2	0
4	B	37	0	33	8	0
4	C	37	0	33	3	0
5	C	1	0	0	0	0
6	C	10	0	0	0	0
7	A	93	0	0	5	0
7	B	81	0	0	9	0
7	C	168	0	0	8	0
All	All	8127	0	7698	291	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 19.

All (291) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:244:HIS:HD2	1:B:258:ILE:HG12	1.21	1.04
1:B:153:GLU:HG3	1:B:246:LYS:HB3	1.39	1.01
1:A:227:LEU:HG	1:A:228:MET:HG3	1.48	0.96
1:A:344:GLU:H	1:A:344:GLU:CD	1.68	0.93
1:B:304:GLU:HG3	1:B:306:THR:HG23	1.48	0.92
1:A:142:GLN:HE21	1:A:146:LYS:HD2	1.35	0.91
1:C:115:MET:HE1	1:C:263:LEU:HB3	1.54	0.90
1:B:106:GLN:HE21	1:B:342:ASN:ND2	1.69	0.89
1:B:244:HIS:CD2	1:B:258:ILE:HG12	2.07	0.89
1:A:57:LEU:HA	1:A:60:LYS:HB2	1.58	0.85
1:A:107:THR:OG1	1:A:270:GLU:OE2	1.96	0.83

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:82:TYR:HE2	1:B:142:GLN:HE21	1.23	0.82
1:A:294:THR:HA	1:A:297:ARG:HG3	1.62	0.81
1:A:190:ASN:HB2	7:A:2066:HOH:O	1.79	0.81
1:C:115:MET:CE	1:C:263:LEU:HB3	2.09	0.80
1:C:98:ASN:O	1:C:328:THR:CG2	2.29	0.80
1:B:289:ASN:OD1	1:B:290:GLN:N	2.13	0.80
1:C:287:ASN:ND2	7:C:2144:HOH:O	2.15	0.79
1:A:167:GLU:OE2	1:A:181:ARG:HD3	1.83	0.79
1:A:122:ASN:HB3	1:A:124:GLU:H	1.47	0.78
1:B:342:ASN:OD1	1:B:346:THR:OG1	1.99	0.77
1:A:270:GLU:O	1:A:271:ASN:ND2	2.18	0.77
1:A:142:GLN:HE21	1:A:146:LYS:CD	1.97	0.77
1:B:166:GLU:O	1:B:315:LYS:NZ	2.17	0.77
1:B:106:GLN:NE2	1:B:342:ASN:ND2	2.32	0.77
4:C:1369:G7X:OAV	4:C:1369:G7X:HBD2	1.86	0.76
1:B:291:SER:OG	7:B:2039:HOH:O	2.05	0.74
1:A:248:THR:HA	1:A:253:GLU:O	1.85	0.74
1:A:340:SER:HA	1:A:343:LEU:CD1	2.17	0.74
1:B:289:ASN:CG	1:B:290:GLN:H	1.91	0.73
1:B:76:THR:O	7:B:2004:HOH:O	2.05	0.73
1:A:140:LEU:HD13	1:A:210:VAL:HG11	1.71	0.73
1:B:40:ILE:CD1	1:B:340:SER:HA	2.20	0.72
1:B:28:PHE:O	1:B:29:ASN:HB2	1.89	0.71
1:B:150:ASN:OD1	1:C:308[B]:HIS:CD2	2.44	0.70
1:C:307:PRO:HG2	1:C:308[A]:HIS:CD2	2.27	0.69
1:A:227:LEU:HG	1:A:228:MET:N	2.06	0.69
1:A:152:THR:HG21	1:A:246:LYS:O	1.92	0.69
1:B:359:ILE:O	7:B:2002:HOH:O	2.10	0.69
1:C:221:ARG:HD2	7:C:2119:HOH:O	1.94	0.68
1:B:92:GLU:HG2	1:B:329:ARG:HD2	1.74	0.68
1:A:19:ILE:HD12	1:A:303:VAL:HG23	1.76	0.68
1:A:359:ILE:HG12	1:A:360:LEU:H	1.58	0.68
1:B:360:LEU:HD12	1:B:360:LEU:H	1.57	0.68
1:A:122:ASN:HB3	1:A:124:GLU:N	2.08	0.68
1:A:56:GLY:O	1:A:57:LEU:HB2	1.94	0.67
1:C:67:THR:HG21	1:C:361:ASN:HA	1.77	0.67
1:A:152:THR:CG2	1:A:153:GLU:H	2.07	0.67
1:B:28:PHE:H	1:B:28:PHE:HD1	1.42	0.66
1:C:123:GLU:O	1:C:124:GLU:HG2	1.93	0.66
1:B:98:ASN:ND2	1:B:260:LYS:HG2	2.10	0.66
1:B:314:SER:O	1:B:318:ARG:HG3	1.96	0.66
1:A:43:CYS:O	1:A:45:PRO:HD3	1.95	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:340:SER:O	1:A:343:LEU:HD13	1.97	0.65
1:C:307:PRO:HG2	1:C:308[A]:HIS:HD2	1.62	0.65
1:A:53:ARG:O	7:A:2014:HOH:O	2.15	0.65
1:A:140:LEU:HD13	1:A:210:VAL:CG1	2.27	0.65
1:B:24:ARG:NH2	2:B:601:ADP:N6	2.45	0.64
1:B:41:VAL:HG12	1:B:42:GLU:H	1.63	0.64
1:A:142:GLN:NE2	1:A:146:LYS:HD2	2.11	0.63
1:C:98:ASN:O	1:C:328:THR:HG23	1.98	0.63
7:A:2043:HOH:O	1:C:180:GLU:OE2	2.16	0.63
1:B:170:ASP:HB2	1:B:182:LEU:HD11	1.80	0.63
1:C:98:ASN:O	1:C:328:THR:HG22	1.99	0.62
1:B:115:MET:CE	1:B:263:LEU:HB3	2.30	0.62
1:A:344:GLU:N	1:A:344:GLU:CD	2.49	0.62
1:C:123:GLU:O	1:C:124:GLU:CG	2.47	0.62
1:C:187:ASP:OD1	1:C:189:ARG:HD3	2.00	0.62
1:A:152:THR:CG2	1:A:153:GLU:N	2.62	0.62
1:B:41:VAL:HG12	1:B:42:GLU:N	2.14	0.62
1:A:51:SER:HA	1:A:64:LYS:O	1.99	0.61
1:C:128:GLU:OE2	1:C:207:LYS:HE2	2.01	0.61
1:A:40:ILE:CD1	1:A:343:LEU:HG	2.30	0.61
1:A:290:GLN:O	1:A:294:THR:HG23	2.01	0.61
1:B:178:VAL:CG2	1:B:224:ALA:HB2	2.30	0.61
1:A:17:LYS:HD2	1:A:17:LYS:N	2.14	0.61
1:A:231:TYR:O	1:A:234:ARG:N	2.32	0.61
1:B:218:ALA:HB2	4:B:1363:G7X:HAF	1.82	0.61
1:C:67:THR:CG2	1:C:361:ASN:HA	2.31	0.60
1:B:40:ILE:HD12	1:B:340:SER:HA	1.83	0.60
1:A:227:LEU:CG	1:A:228:MET:HG3	2.28	0.60
1:A:340:SER:HA	1:A:343:LEU:HD12	1.83	0.60
1:A:175:SER:HB2	1:C:203:THR:HB	1.82	0.60
1:C:124:GLU:HG2	1:C:125:TYR:H	1.66	0.60
1:B:28:PHE:N	1:B:28:PHE:CD1	2.70	0.60
1:B:136:ILE:N	1:B:137:PRO:HD2	2.15	0.60
1:C:312:ARG:NH2	7:C:2154:HOH:O	2.34	0.60
1:C:124:GLU:CG	1:C:125:TYR:H	2.15	0.59
1:C:186:ASP:O	1:C:188:PRO:HD3	2.02	0.59
1:A:344:GLU:HG2	1:A:345:GLU:OE1	2.02	0.59
1:C:328:THR:HG21	7:C:2045:HOH:O	2.01	0.59
1:B:150:ASN:OD1	1:C:308[B]:HIS:CG	2.56	0.59
1:B:289:ASN:CG	1:B:290:GLN:N	2.54	0.58
1:A:142:GLN:HG3	1:A:146:LYS:HD2	1.85	0.58
1:C:230:ALA:HB3	1:C:234:ARG:HD3	1.84	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:340:SER:O	1:A:343:LEU:CD1	2.50	0.58
1:B:130:ASP:OD1	1:B:131:PRO:HD2	2.04	0.58
1:A:152:THR:HG22	1:A:153:GLU:N	2.19	0.58
1:B:191:LYS:HD2	1:B:191:LYS:N	2.19	0.57
1:B:40:ILE:O	1:B:40:ILE:HG22	2.04	0.57
1:B:98:ASN:O	1:B:328:THR:OG1	2.20	0.57
1:A:53:ARG:HH11	1:A:63:ARG:HG3	1.70	0.57
1:C:271:ASN:HB2	7:C:2140:HOH:O	2.05	0.57
1:A:287:ASN:O	1:A:290:GLN:HB3	2.04	0.57
1:C:123:GLU:O	1:C:124:GLU:OE1	2.22	0.57
1:B:98:ASN:HD21	1:B:260:LYS:HG2	1.70	0.56
1:A:164:TYR:HE2	1:A:228:MET:HB3	1.70	0.56
1:B:40:ILE:HD11	1:B:340:SER:HA	1.87	0.56
1:B:133:ALA:O	1:B:138:ARG:NE	2.39	0.56
1:B:41:VAL:CG2	1:B:338:PRO:HA	2.36	0.56
1:A:127:TRP:HB2	4:A:1365:G7X:HAZ3	1.88	0.56
1:B:337:SER:HB3	1:B:342:ASN:ND2	2.21	0.55
1:B:115:MET:HE1	1:B:263:LEU:HB3	1.89	0.55
1:B:225:ALA:HA	1:B:231:TYR:CD1	2.42	0.55
1:A:254:GLU:O	1:A:255:LEU:HD22	2.06	0.55
1:A:152:THR:HG23	1:A:153:GLU:H	1.72	0.55
1:B:229:ASN:O	1:B:230:ALA:C	2.43	0.55
1:A:177:ASP:HA	7:A:2057:HOH:O	2.07	0.55
1:A:142:GLN:NE2	1:A:146:LYS:HG3	2.22	0.55
1:A:181:ARG:HH21	1:A:228:MET:HE1	1.72	0.54
1:B:70:MET:HE3	1:B:84:SER:HB2	1.90	0.54
1:C:295:LEU:O	1:C:299:ILE:HG12	2.07	0.54
1:B:147:LEU:HD11	1:B:243:ILE:HG21	1.89	0.54
1:B:231:TYR:C	1:B:231:TYR:CD1	2.80	0.54
1:B:43:CYS:HB3	1:B:71:VAL:CG1	2.38	0.54
1:B:20:GLN:OE1	1:B:69:ASP:HB3	2.07	0.54
1:A:290:GLN:O	1:A:294:THR:CG2	2.56	0.54
1:B:229:ASN:OD1	1:B:234:ARG:NE	2.40	0.54
1:B:70:MET:CE	1:B:84:SER:HB2	2.37	0.54
1:B:159:SER:HA	1:B:172:LEU:HD12	1.89	0.53
4:B:1363:G7X:CAL	4:B:1363:G7X:HBD1	2.38	0.53
1:A:165[B]:ASN:OD1	1:A:284:GLU:HB3	2.09	0.53
1:B:43:CYS:HB3	1:B:71:VAL:HG12	1.91	0.53
1:A:26:ARG:HB2	1:A:27:PRO:HD2	1.90	0.53
1:A:67:THR:O	1:A:359:ILE:CG2	2.57	0.52
1:A:40:ILE:HD13	1:A:343:LEU:HG	1.89	0.52
1:A:53:ARG:NH1	1:A:63:ARG:HG3	2.25	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:87:CYS:HB2	1:A:88:PRO:HD3	1.90	0.52
1:A:23:VAL:HG21	1:A:68:PHE:CE2	2.44	0.52
1:B:98:ASN:OD1	1:B:323:SER:HB2	2.09	0.52
1:B:144:PHE:HZ	1:B:204:VAL:HG22	1.75	0.52
1:B:89:ILE:HD11	1:B:331:SER:OG	2.10	0.52
1:C:354:HIS:NE2	1:C:358:ASN:OD1	2.43	0.52
1:B:171:LEU:HD13	1:B:221:ARG:HB2	1.91	0.52
1:A:44:ASP:OD2	1:A:47:ARG:HD3	2.10	0.52
1:A:67:THR:O	1:A:359:ILE:HG21	2.10	0.52
1:C:67:THR:HG21	7:C:2019:HOH:O	2.09	0.52
1:B:204:VAL:HG21	1:B:210:VAL:HG23	1.92	0.52
1:A:26:ARG:HD3	2:A:601:ADP:C8	2.45	0.51
1:B:41:VAL:HG23	1:B:338:PRO:HA	1.92	0.51
1:A:342:ASN:ND2	7:A:2010:HOH:O	2.11	0.51
1:C:40:ILE:HD12	1:C:343:LEU:HD13	1.93	0.51
1:B:28:PHE:O	1:B:29:ASN:CB	2.57	0.51
1:B:178:VAL:HG21	1:B:224:ALA:HB2	1.93	0.51
1:A:359:ILE:HG12	1:A:360:LEU:N	2.26	0.51
1:B:26:ARG:NH1	1:B:27:PRO:O	2.43	0.51
1:B:144:PHE:HZ	1:B:204:VAL:CG2	2.24	0.51
1:B:77:LYS:HA	7:B:2005:HOH:O	2.10	0.51
1:B:73:GLY:HA3	7:B:2003:HOH:O	2.10	0.50
1:C:345:GLU:HG2	7:C:2163:HOH:O	2.10	0.50
1:A:340:SER:HA	1:A:343:LEU:HD11	1.92	0.50
1:B:131:PRO:HA	7:B:2026:HOH:O	2.10	0.50
1:B:102:PHE:HB2	1:B:332:ILE:HB	1.94	0.50
1:A:160:LEU:HB3	1:A:172:LEU:HG	1.94	0.50
1:B:93:VAL:O	1:B:259:GLY:HA3	2.12	0.50
1:C:115:MET:HE2	1:C:263:LEU:HB3	1.92	0.49
1:C:99:CYS:HA	1:C:328:THR:HG22	1.94	0.49
1:A:64:LYS:HD3	1:A:65:THR:H	1.77	0.49
1:B:22:VAL:CG1	1:B:70:MET:HB2	2.43	0.49
1:A:209:GLU:O	1:A:213:ILE:HD12	2.13	0.49
4:C:1369:G7X:CAL	4:C:1369:G7X:HBD2	2.42	0.49
1:C:67:THR:CG2	1:C:361:ASN:OD1	2.61	0.49
1:A:91:ASP:OD1	1:A:146:LYS:NZ	2.42	0.48
1:B:191:LYS:HZ3	1:B:191:LYS:HA	1.77	0.48
1:B:191:LYS:HA	1:B:191:LYS:NZ	2.28	0.48
1:C:253:GLU:OE2	1:C:255:LEU:HD11	2.13	0.48
1:B:47:ARG:O	1:B:49:GLU:HG3	2.13	0.48
1:B:347:LEU:O	1:B:351:GLU:HG2	2.13	0.48
1:C:170:ASP:HB2	1:C:182:LEU:HD11	1.96	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:256:VAL:O	1:B:256:VAL:HG12	2.13	0.48
1:A:237:SER:HB3	1:A:265:ASP:HB3	1.96	0.48
1:B:144:PHE:CZ	1:B:204:VAL:CG2	2.96	0.48
1:B:312:ARG:HG2	7:B:2071:HOH:O	2.13	0.47
1:B:160:LEU:HD21	4:B:1363:G7X:CLD	2.51	0.47
1:A:303:VAL:HG21	1:A:357:LYS:HB2	1.94	0.47
1:A:66:TYR:OH	1:A:351:GLU:OE2	2.28	0.47
1:B:342:ASN:OD1	1:B:342:ASN:C	2.53	0.47
4:B:1363:G7X:NAH	4:B:1363:G7X:HBC1	2.28	0.47
1:A:70:MET:HE3	1:A:72:PHE:HZ	1.78	0.47
1:B:128:GLU:OE2	7:B:2024:HOH:O	2.20	0.47
1:A:18:ASN:HB2	1:A:328:THR:O	2.15	0.47
1:C:306:THR:HG23	1:C:307:PRO:HD2	1.96	0.47
1:A:247:GLU:OE1	1:A:257:LYS:HG2	2.14	0.47
1:B:104:TYR:OH	1:B:269:SER:HB2	2.15	0.47
1:A:41:VAL:HG22	1:A:52:VAL:HG12	1.97	0.46
1:B:68:PHE:CZ	1:B:350:LEU:CD2	2.99	0.46
1:A:289:ASN:O	1:A:293:LEU:HB2	2.15	0.46
1:B:228:MET:O	1:B:231:TYR:HB3	2.16	0.46
1:A:44:ASP:N	1:A:49:GLU:O	2.36	0.46
1:B:183:GLN:HA	1:B:183:GLN:OE1	2.14	0.46
1:B:41:VAL:O	1:B:42:GLU:HB2	2.15	0.46
1:B:225:ALA:HB2	1:B:231:TYR:CE1	2.50	0.46
1:A:254:GLU:N	1:A:254:GLU:OE1	2.49	0.46
1:C:70:MET:CE	1:C:85:VAL:HG22	2.46	0.46
1:C:115:MET:O	1:C:136:ILE:HG13	2.16	0.46
1:A:168:LEU:HD11	1:A:184:MET:HE2	1.97	0.46
1:B:106:GLN:NE2	1:B:342:ASN:HD22	2.10	0.45
1:B:184:MET:HA	1:B:195:ILE:O	2.16	0.45
1:A:204:VAL:CG1	1:A:213:ILE:CD1	2.95	0.45
1:A:162:GLU:CG	1:A:171:LEU:HD11	2.46	0.45
1:A:164:TYR:CE2	1:A:228:MET:HB3	2.51	0.45
1:C:209:GLU:O	1:C:213:ILE:HD13	2.17	0.45
1:A:19:ILE:O	1:A:357:LYS:NZ	2.43	0.45
1:C:124:GLU:HG3	1:C:125:TYR:CD2	2.52	0.45
1:A:70:MET:HE3	1:A:72:PHE:CZ	2.51	0.45
4:C:1369:G7X:OAV	4:C:1369:G7X:CBD	2.59	0.45
1:C:312:ARG:NH1	7:C:2105:HOH:O	2.45	0.45
1:A:104:TYR:HB2	1:A:266:LEU:HD12	1.99	0.45
1:C:70:MET:HB2	1:C:70:MET:HE3	1.65	0.45
1:C:362:LYS:HE2	1:C:362:LYS:HB2	1.75	0.45
1:B:119:ARG:NH2	1:B:211:TYR:OH	2.50	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:601:ADP:O2A	2:A:601:ADP:H3'	2.16	0.45
1:B:360:LEU:HD12	1:B:360:LEU:N	2.29	0.45
1:B:300:THR:O	1:B:304:GLU:HG2	2.17	0.44
1:C:270:GLU:O	1:C:271:ASN:HB2	2.16	0.44
1:A:229:ASN:O	1:A:230:ALA:HB3	2.17	0.44
1:B:299:ILE:HG21	1:B:356:ALA:HB1	1.99	0.44
1:B:149:ASP:HB3	1:C:308[B]:HIS:CD2	2.53	0.44
1:B:26:ARG:O	1:B:338:PRO:CG	2.65	0.44
1:B:190:ASN:C	1:B:191:LYS:HD2	2.38	0.44
1:A:344:GLU:N	1:A:344:GLU:OE1	2.50	0.44
1:B:24:ARG:NH2	2:B:601:ADP:HN61	2.13	0.44
1:C:67:THR:HG23	1:C:361:ASN:OD1	2.18	0.44
1:A:50:VAL:HG13	1:A:66:TYR:HB2	2.00	0.44
1:A:142:GLN:NE2	1:A:146:LYS:CG	2.82	0.43
1:B:224:ALA:HB3	1:B:231:TYR:OH	2.18	0.43
1:A:270:GLU:O	1:A:271:ASN:CB	2.66	0.43
1:B:295:LEU:O	1:B:299:ILE:HG12	2.18	0.43
1:B:19:ILE:HA	1:B:330:THR:O	2.17	0.43
1:C:70:MET:HE2	1:C:85:VAL:HG22	2.01	0.43
1:A:105:GLY:O	1:A:111:LYS:NZ	2.34	0.43
1:B:27:PRO:HG3	1:B:74:ALA:O	2.19	0.43
1:B:243:ILE:O	1:B:258:ILE:HA	2.19	0.43
1:B:160:LEU:CD2	4:B:1363:G7X:CLD	3.03	0.43
1:B:236:HIS:HE1	1:B:292:LEU:HD13	1.83	0.43
4:B:1363:G7X:OAV	4:B:1363:G7X:HBD1	2.19	0.43
1:B:24:ARG:HD2	1:B:76:THR:HG21	2.01	0.42
1:B:78:GLN:HB2	1:B:138:ARG:HH21	1.84	0.42
1:C:124:GLU:HG2	1:C:125:TYR:N	2.33	0.42
1:B:324:LEU:HD23	1:B:324:LEU:HA	1.85	0.42
1:B:120:SER:HA	1:B:121:PRO:HD3	1.90	0.42
1:B:92:GLU:HG2	1:B:329:ARG:CD	2.44	0.42
1:B:41:VAL:HG21	1:B:338:PRO:HA	2.01	0.42
1:B:299:ILE:O	1:B:303:VAL:HG23	2.20	0.42
1:C:53:ARG:HH11	1:C:63:ARG:CG	2.32	0.42
1:B:144:PHE:CZ	1:B:204:VAL:HG22	2.55	0.42
1:B:127:TRP:HB2	4:B:1363:G7X:HAZ3	2.01	0.42
1:C:26:ARG:HE	1:C:109:THR:HA	1.84	0.42
1:B:158:VAL:HA	1:B:240:SER:O	2.20	0.42
1:C:40:ILE:CD1	1:C:343:LEU:HD13	2.49	0.42
1:B:24:ARG:HH22	1:B:113:PHE:HE2	1.66	0.41
1:C:306:THR:CG2	1:C:307:PRO:HD2	2.49	0.41
1:A:106:GLN:NE2	1:A:345:GLU:HG2	2.35	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:136:ILE:N	1:B:137:PRO:CD	2.82	0.41
1:A:44:ASP:O	1:A:48:LYS:N	2.53	0.41
1:B:244:HIS:HD2	1:B:258:ILE:CG1	2.11	0.41
1:B:99:CYS:O	1:B:261:LEU:HA	2.19	0.41
1:A:142:GLN:HE21	1:A:146:LYS:CG	2.33	0.41
4:A:1365:G7X:HBB2	4:A:1365:G7X:CAW	2.51	0.41
1:A:293:LEU:HA	1:A:293:LEU:HD12	1.81	0.41
1:C:70:MET:HE1	1:C:72:PHE:HZ	1.85	0.41
1:B:320:LEU:O	1:B:321:GLN:C	2.59	0.41
1:A:362:LYS:CB	1:A:363:PRO:HD2	2.49	0.41
4:B:1363:G7X:HBB2	4:B:1363:G7X:CAW	2.50	0.41
1:A:59:ASP:O	1:A:60:LYS:HB3	2.21	0.41
1:A:40:ILE:HD13	1:A:343:LEU:CG	2.49	0.41
1:C:124:GLU:CG	1:C:125:TYR:N	2.82	0.41
1:A:64:LYS:HA	1:A:64:LYS:HD3	1.64	0.41
1:A:50:VAL:CG1	1:A:68:PHE:HE2	2.34	0.41
1:B:342:ASN:O	1:B:342:ASN:OD1	2.40	0.41
1:B:184:MET:HB2	1:B:184:MET:HE2	1.90	0.41
1:A:336:ILE:HD13	1:A:350:LEU:HD21	2.03	0.40
1:B:191:LYS:HD3	7:B:2052:HOH:O	2.21	0.40
1:B:50:VAL:HG22	1:B:71:VAL:HG11	2.03	0.40
1:A:162:GLU:OE2	1:A:171:LEU:HD11	2.21	0.40
1:A:362:LYS:O	1:A:363:PRO:C	2.59	0.40
1:C:342:ASN:O	1:C:343:LEU:C	2.60	0.40
1:C:104:TYR:CD1	1:C:104:TYR:C	2.94	0.40
1:B:26:ARG:NE	1:B:108:GLY:O	2.54	0.40
1:B:22:VAL:HG12	1:B:70:MET:HB2	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	332/368 (90%)	319 (96%)	12 (4%)	1 (0%)	50 70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	292/368 (79%)	280 (96%)	12 (4%)	0	100	100
1	C	336/368 (91%)	332 (99%)	4 (1%)	0	100	100
All	All	960/1104 (87%)	931 (97%)	28 (3%)	1 (0%)	59	80

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	GLU

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	287/322 (89%)	264 (92%)	23 (8%)	17	29
1	B	266/322 (83%)	236 (89%)	30 (11%)	9	14
1	C	287/322 (89%)	272 (95%)	15 (5%)	32	53
All	All	840/966 (87%)	772 (92%)	68 (8%)	17	28

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	34	LYS
1	A	54	THR
1	A	64	LYS
1	A	122	ASN
1	A	125	TYR
1	A	181	ARG
1	A	184	MET
1	A	207	LYS
1	A	208	ASP
1	A	229	ASN
1	A	284	GLU
1	A	290	GLN
1	A	291	SER

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Mol	Chain	Res	Type
1	A	292	LEU
1	A	293	LEU
1	A	294	THR
1	A	306	THR
1	A	312	ARG
1	A	344	GLU
1	A	347	LEU
1	A	348	SER
1	A	360	LEU
1	B	24	ARG
1	B	28	PHE
1	B	46	VAL
1	B	70	MET
1	B	77	LYS
1	B	79	ILE
1	B	95	MET
1	B	111	LYS
1	B	119	ARG
1	B	142	GLN
1	B	161	LEU
1	B	183	GLN
1	B	184	MET
1	B	190	ASN
1	B	191	LYS
1	B	212	GLN
1	B	231	TYR
1	B	235	SER
1	B	260	LYS
1	B	262	ASN
1	B	290	GLN
1	B	291	SER
1	B	292	LEU
1	B	312	ARG
1	B	313	GLU
1	B	332	ILE
1	B	342	ASN
1	B	348	SER
1	B	360	LEU
1	B	362	LYS
1	C	22	VAL
1	C	30	LEU
1	C	33	ARG

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Mol	Chain	Res	Type
1	C	67	THR
1	C	84	SER
1	C	124	GLU
1	C	146	LYS
1	C	149	ASP
1	C	210	VAL
1	C	212	GLN
1	C	237	SER
1	C	248	THR
1	C	292	LEU
1	C	304	GLU
1	C	328	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	142	GLN
1	B	98	ASN
1	B	106	GLN
1	B	244	HIS
1	B	262	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	G7X	A	1365	-	40,40,40	1.23	5 (12%)	53,56,56	1.67	6 (11%)
2	ADP	A	601	-	29,29,29	1.32	4 (13%)	45,45,45	2.24	12 (26%)
4	G7X	B	1363	-	40,40,40	1.34	6 (15%)	53,56,56	1.36	6 (11%)
2	ADP	B	601	-	29,29,29	1.24	4 (13%)	45,45,45	1.86	9 (20%)
6	PO4	C	1366	-	4,4,4	0.20	0	6,6,6	0.35	0
6	PO4	C	1367	-	4,4,4	0.19	0	6,6,6	0.31	0
4	G7X	C	1369	-	40,40,40	1.35	6 (15%)	53,56,56	1.69	9 (16%)
2	ADP	C	601	5	29,29,29	1.14	3 (10%)	45,45,45	1.95	10 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	G7X	A	1365	-	-	0/25/28/28	0/2/4/4
2	ADP	A	601	-	-	0/16/32/32	0/1/3/3
4	G7X	B	1363	-	-	0/25/28/28	0/2/4/4
2	ADP	B	601	-	-	0/16/32/32	0/1/3/3
6	PO4	C	1366	-	-	0/0/0/0	0/0/0/0
6	PO4	C	1367	-	-	0/0/0/0	0/0/0/0
4	G7X	C	1369	-	-	0/25/28/28	0/2/4/4
2	ADP	C	601	5	-	0/16/32/32	0/1/3/3

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1365	G7X	CAJ-NAK	3.50	1.51	1.46
4	B	1363	G7X	CAM-CAL	3.26	1.55	1.50
2	B	601	ADP	C5-C4	3.25	1.47	1.40
4	B	1363	G7X	CAI-CAJ	3.20	1.55	1.50
4	A	1365	G7X	CAR-CAG	-3.19	1.38	1.45
2	A	601	ADP	C4-N9	-3.15	1.33	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1369	G7X	CAI-CAJ	3.12	1.55	1.50
2	A	601	ADP	C5-C4	3.02	1.47	1.40
2	A	601	ADP	C2'-C1'	-2.96	1.49	1.53
2	C	601	ADP	C4-N9	-2.94	1.33	1.37
4	A	1365	G7X	CAF-CAG	-2.93	1.38	1.41
4	B	1363	G7X	CAJ-NAK	2.89	1.51	1.46
2	B	601	ADP	C4-N9	-2.84	1.33	1.37
4	C	1369	G7X	CAF-CAE	2.67	1.42	1.36
4	C	1369	G7X	CAM-CAL	2.64	1.54	1.50
4	C	1369	G7X	CAR-CAG	-2.56	1.40	1.45
4	C	1369	G7X	CAJ-NAK	2.52	1.50	1.46
4	C	1369	G7X	CAU-NAK	2.40	1.50	1.46
2	C	601	ADP	C5-C4	2.36	1.45	1.40
4	B	1363	G7X	CAF-CAG	-2.33	1.39	1.41
2	B	601	ADP	PB-O3A	2.28	1.64	1.60
4	B	1363	G7X	CAR-CAG	-2.22	1.40	1.45
4	A	1365	G7X	CAM-CAL	2.16	1.53	1.50
4	A	1365	G7X	CAS-NAT	2.14	1.41	1.35
2	B	601	ADP	PA-O3A	2.11	1.63	1.59
2	A	601	ADP	C2-N3	2.11	1.36	1.32
4	B	1363	G7X	CAI-NAT	2.10	1.41	1.36
2	C	601	ADP	PB-O3A	2.02	1.63	1.60

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1365	G7X	CAI-NAH-CAG	7.41	122.59	118.09
2	A	601	ADP	N3-C2-N1	-6.54	123.24	128.71
2	B	601	ADP	N3-C2-N1	-6.48	123.29	128.71
4	C	1369	G7X	CAI-NAH-CAG	6.13	121.81	118.09
4	C	1369	G7X	CAI-CAJ-CAB	-6.08	105.01	112.69
2	C	601	ADP	O4'-C1'-N9	-6.04	102.82	108.44
2	C	601	ADP	N3-C2-N1	-5.67	123.97	128.71
2	A	601	ADP	N3-C4-N9	5.61	135.56	125.43
4	B	1363	G7X	CAI-NAH-CAG	5.58	121.48	118.09
2	C	601	ADP	N3-C4-N9	5.15	134.73	125.43
2	B	601	ADP	N3-C4-N9	4.97	134.41	125.43
2	A	601	ADP	PA-O3A-PB	-4.69	117.92	131.68
2	B	601	ADP	C8-N9-C4	4.31	110.19	106.90
2	A	601	ADP	O4'-C1'-N9	4.08	112.24	108.44
2	A	601	ADP	C8-N9-C4	3.95	109.92	106.90
4	A	1365	G7X	CAQ-CAR-CAG	3.71	122.11	118.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	ADP	C4-C5-N7	-3.60	106.44	109.52
2	A	601	ADP	O2'-C2'-C1'	-3.52	100.58	111.23
2	A	601	ADP	O3B-PB-O2B	3.51	121.26	107.61
2	C	601	ADP	C5-C4-N3	-3.41	118.27	125.70
2	A	601	ADP	C5-C4-N3	-3.13	118.89	125.70
2	C	601	ADP	C4-C5-N7	-3.01	106.94	109.52
2	A	601	ADP	C3'-C2'-C1'	2.98	105.57	100.91
2	C	601	ADP	O3'-C3'-C2'	-2.96	102.20	111.83
2	A	601	ADP	C4-C5-N7	-2.81	107.12	109.52
2	C	601	ADP	O3'-C3'-C4'	-2.74	103.00	111.08
4	A	1365	G7X	CAI-CAJ-CAB	-2.71	109.27	112.69
2	C	601	ADP	C4'-O4'-C1'	2.66	112.64	109.75
4	B	1363	G7X	CBF-CBG-CBH	2.58	122.50	118.16
2	B	601	ADP	C5-C4-N3	-2.56	120.11	125.70
4	C	1369	G7X	OAV-CAL-CAM	-2.56	115.35	120.17
2	B	601	ADP	PA-O3A-PB	-2.48	124.40	131.68
4	A	1365	G7X	CAJ-NAK-CAL	2.45	127.54	118.96
4	C	1369	G7X	CAC-CAB-CAJ	-2.44	107.45	110.57
2	C	601	ADP	O2B-PB-O1B	2.42	118.35	110.44
2	A	601	ADP	C2-N3-C4	2.41	120.88	114.01
4	B	1363	G7X	CAI-CAJ-CAB	-2.34	109.74	112.69
4	C	1369	G7X	CAM-CAL-NAK	2.34	121.97	118.55
2	C	601	ADP	C2-N3-C4	2.28	120.50	114.01
4	B	1363	G7X	CBG-CBB-NAT	-2.26	109.03	112.62
2	A	601	ADP	C2'-C1'-N9	-2.21	107.60	113.27
4	B	1363	G7X	CAB-CAJ-NAK	-2.19	109.79	113.33
4	C	1369	G7X	CBG-CBB-NAT	-2.19	109.14	112.62
4	C	1369	G7X	CAJ-NAK-CAL	2.18	126.61	118.96
4	A	1365	G7X	CAF-CAG-CAR	-2.18	117.90	119.45
2	B	601	ADP	C2-N3-C4	2.18	120.21	114.01
2	B	601	ADP	O2B-PB-O1B	2.17	117.53	110.44
4	C	1369	G7X	CAQ-CAR-CAG	2.11	120.47	118.32
2	B	601	ADP	C2-N1-C6	2.05	122.47	118.77
4	A	1365	G7X	CAJ-CAI-NAH	2.04	119.80	116.48
4	C	1369	G7X	CAA-CAB-CAJ	2.01	113.14	110.57
4	B	1363	G7X	CAR-CAS-NAT	-2.00	116.97	119.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	334/368 (90%)	0.30	27 (8%) 12 11	26, 49, 91, 110	0
1	B	302/368 (82%)	0.49	40 (13%) 4 3	37, 62, 98, 118	0
1	C	339/368 (92%)	-0.09	4 (1%) 75 76	18, 37, 65, 99	0
All	All	975/1104 (88%)	0.22	71 (7%) 15 14	18, 49, 91, 118	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	58	ALA	6.9
1	B	46	VAL	6.3
1	B	343	LEU	6.1
1	C	230	ALA	5.6
1	A	360	LEU	5.2
1	B	341	LEU	5.2
1	A	285	ALA	4.9
1	B	38	HIS	4.9
1	B	340	SER	4.9
1	B	47	ARG	4.3
1	A	282	ALA	4.3
1	B	360	LEU	4.2
1	B	342	ASN	4.0
1	A	307	PRO	3.9
1	B	95	MET	3.7
1	A	229	ASN	3.7
1	B	37	ALA	3.7
1	B	189	ARG	3.3
1	B	339	ALA	3.3
1	B	27	PRO	3.3
1	A	363	PRO	3.3
1	B	28	PHE	3.2
1	B	229	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	271	ASN	3.2
1	B	362	LYS	3.2
1	B	327	ARG	3.1
1	A	264	VAL	3.1
1	B	326	GLY	3.0
1	B	263	LEU	3.0
1	A	56	GLY	3.0
1	B	361	ASN	2.9
1	A	102	PHE	2.9
1	C	58	ALA	2.9
1	A	327	ARG	2.9
1	B	246	LYS	2.9
1	B	191	LYS	2.8
1	B	264	VAL	2.8
1	A	284	GLU	2.8
1	B	65	THR	2.8
1	A	101	ILE	2.8
1	A	289	ASN	2.7
1	B	152	THR	2.7
1	C	59	ASP	2.6
1	A	308	HIS	2.6
1	B	67	THR	2.6
1	C	282	ALA	2.5
1	A	281	ARG	2.4
1	B	305	ARG	2.4
1	B	301	ALA	2.4
1	B	230	ALA	2.3
1	B	256	VAL	2.3
1	A	263	LEU	2.3
1	A	239	PHE	2.3
1	B	83	ARG	2.3
1	A	266	LEU	2.2
1	A	247	GLU	2.2
1	B	261	LEU	2.1
1	A	122	ASN	2.1
1	B	344	GLU	2.1
1	A	286	GLY	2.1
1	B	238	VAL	2.1
1	B	102	PHE	2.1
1	B	356	ALA	2.1
1	A	59	ASP	2.1
1	A	255	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	332	ILE	2.1
1	B	358	ASN	2.0
1	A	240	SER	2.0
1	B	153	GLU	2.0
1	A	262	ASN	2.0
1	B	269	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	G7X	A	1365	37/37	0.15	0.12	30,38,47,51	0
4	G7X	B	1363	37/37	0.12	-0.49	36,45,53,56	0
6	PO4	C	1367	5/5	0.12	-0.64	63,71,86,86	0
6	PO4	C	1366	5/5	0.11	-0.70	53,54,57,60	0
4	G7X	C	1369	37/37	0.10	-0.88	22,27,37,39	0
2	ADP	B	601	27/27	0.10	-0.99	69,76,84,88	0
2	ADP	C	601	27/27	0.13	-1.23	26,36,45,48	0
2	ADP	A	601	27/27	0.12	-1.30	42,52,61,62	0
3	CL	A	1364	1/1	0.09	-1.55	74,74,74,74	0
3	CL	C	1368	1/1	0.10	-1.57	38,38,38,38	0
5	MG	C	1365	1/1	0.12	-1.95	48,48,48,48	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.