



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:36 AM UTC

PDB ID : 3AGR / pdb_00003agr
Title : Crystal structure of nucleoside triphosphate hydrolases from *Neospora caninum*
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Deposited on : 2010-04-06
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

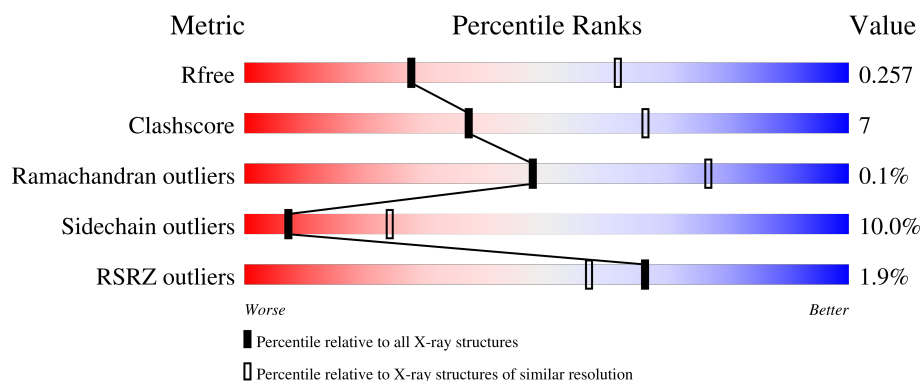
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	602	 2% 80% 15% . .
1	B	602	 % 77% 17% . .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9179 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nucleoside triphosphate hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	591	Total	C	N	O	S	0	0	0
			4586	2885	807	867	27			
1	B	591	Total	C	N	O	S	0	0	0
			4586	2885	807	867	27			

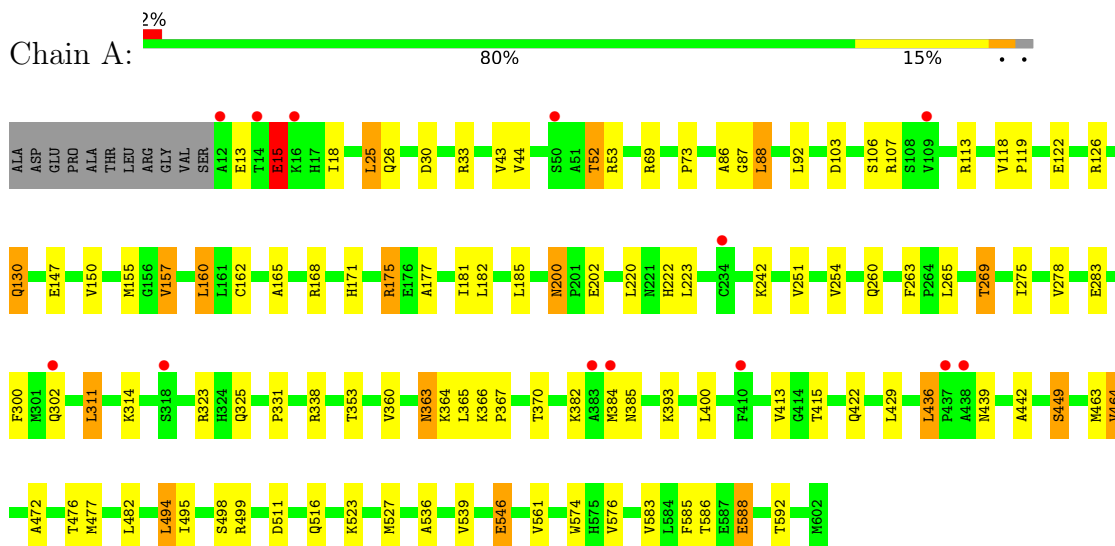
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	O	0	0
			4	4		
2	B	3	Total	O	0	0
			3	3		

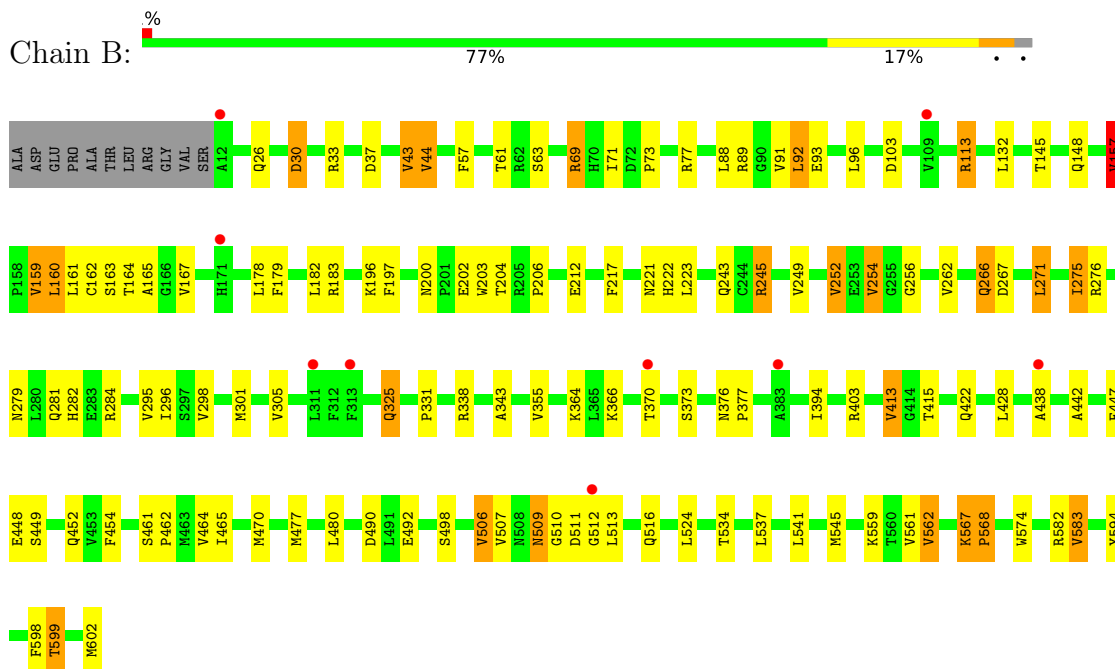
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Nucleoside triphosphate hydrolase



• Molecule 1: Nucleoside triphosphate hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	93.66Å 140.78Å 301.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.51 – 2.80 42.51 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.2 (42.51-2.80) 93.1 (42.51-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.78 (at 2.81Å)	Xtriage
Refinement program	REFMAC refmac_5.5.0109	Depositor
R, R_{free}	0.228 , 0.265 0.220 , 0.257	Depositor DCC
R_{free} test set	2354 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	72.2	Xtriage
Anisotropy	0.554	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9179	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.53	0/4684	0.88	6/6343 (0.1%)
1	B	0.53	0/4684	0.89	4/6343 (0.1%)
All	All	0.53	0/9368	0.89	10/12686 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	567	LYS	CA-C-N	8.48	130.44	119.84
1	B	567	LYS	C-N-CA	8.48	130.44	119.84
1	A	165	ALA	N-CA-C	6.44	118.30	111.28
1	A	157	VAL	N-CA-C	6.09	114.18	108.15
1	A	586	THR	N-CA-C	5.57	117.79	111.11
1	A	449	SER	N-CA-C	5.43	115.72	108.38
1	B	157	VAL	N-CA-C	5.08	113.18	108.15
1	B	157	VAL	CB-CA-C	5.07	115.90	111.08
1	A	157	VAL	CB-CA-C	5.06	115.89	111.08
1	A	15	GLU	N-CA-C	-5.03	108.89	114.62

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4586	0	4494	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4586	0	4494	81	0
2	A	4	0	0	0	0
2	B	3	0	0	0	0
All	All	9179	0	8988	129	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (129) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:THR:HG22	1:B:165:ALA:H	1.04	1.19
1:B:567:LYS:HB3	1:B:568:PRO:HD2	1.18	1.15
1:B:567:LYS:HB3	1:B:568:PRO:CD	1.77	1.11
1:B:245:ARG:HG2	1:B:245:ARG:HH21	1.22	1.05
1:B:567:LYS:CB	1:B:568:PRO:HD2	1.99	0.92
1:A:302:GLN:HB3	1:A:436:LEU:HD11	1.53	0.91
1:B:164:THR:HG22	1:B:165:ALA:N	1.86	0.88
1:A:363:ASN:HD22	1:A:365:LEU:H	1.21	0.87
1:B:249:VAL:HG12	1:B:462:PRO:HG2	1.58	0.84
1:B:164:THR:CG2	1:B:165:ALA:H	1.90	0.81
1:B:221:ASN:HD21	1:B:249:VAL:H	1.28	0.79
1:B:279:ASN:HD22	1:B:282:HIS:H	1.28	0.77
1:B:506:VAL:HG21	1:B:524:LEU:HD22	1.67	0.77
1:B:200:ASN:HD22	1:B:202:GLU:H	1.33	0.74
1:B:438:ALA:HB2	1:B:448:GLU:HG2	1.73	0.70
1:B:245:ARG:HG2	1:B:245:ARG:NH2	1.96	0.68
1:B:43:VAL:HG22	1:B:159:VAL:HB	1.77	0.66
1:B:221:ASN:HD21	1:B:249:VAL:N	1.93	0.66
1:A:585:PHE:HB3	1:A:588:GLU:HG3	1.78	0.65
1:A:200:ASN:HD22	1:A:202:GLU:H	1.45	0.63
1:A:168:ARG:HA	1:A:175:ARG:HH21	1.62	0.63
1:B:254:VAL:HG13	1:B:470:MET:HG3	1.81	0.62
1:A:52:THR:HG23	1:A:86:ALA:O	2.00	0.61
1:A:363:ASN:ND2	1:A:365:LEU:H	1.96	0.61
1:B:221:ASN:ND2	1:B:249:VAL:H	1.98	0.61
1:B:162:CYS:HB2	1:B:574:TRP:CH2	2.36	0.60
1:B:366:LYS:HE2	1:B:370:THR:OG1	2.02	0.60
1:B:449:SER:OG	1:B:452:GLN:HG3	2.03	0.59
1:B:266:GLN:NE2	1:B:267:ASP:H	2.01	0.58
1:B:196:LYS:HD3	1:B:598:PHE:CE1	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:CYS:HB2	1:A:574:TRP:CH2	2.38	0.58
1:B:203:TRP:CD1	1:B:582:ARG:HG2	2.39	0.57
1:B:145:THR:H	1:B:148:GLN:HE21	1.52	0.57
1:B:394:ILE:HG23	1:B:403:ARG:HG2	1.85	0.57
1:B:477:MET:HE1	1:B:498:SER:HA	1.85	0.57
1:B:161:LEU:HB3	1:B:204:THR:HG23	1.86	0.57
1:B:279:ASN:ND2	1:B:282:HIS:H	2.02	0.56
1:A:168:ARG:HA	1:A:175:ARG:NH2	2.20	0.56
1:A:26:GLN:O	1:A:30:ASP:HB2	2.06	0.56
1:A:200:ASN:HD22	1:A:200:ASN:C	2.14	0.55
1:B:245:ARG:NH2	1:B:245:ARG:CG	2.68	0.55
1:B:276:ARG:HH12	1:B:442:ALA:HA	1.72	0.55
1:B:509:ASN:HD22	1:B:510:GLY:H	1.55	0.55
1:B:92:LEU:HD13	1:B:178:LEU:HD22	1.89	0.54
1:A:13:GLU:C	1:A:15:GLU:H	2.15	0.54
1:B:298:VAL:HG11	1:B:447:PHE:HB2	1.89	0.54
1:A:472:ALA:O	1:A:476:THR:HG23	2.09	0.53
1:A:69:ARG:CD	1:A:155:MET:HE1	2.40	0.52
1:A:360:VAL:O	1:A:366:LYS:HG3	2.10	0.51
1:B:449:SER:HG	1:B:452:GLN:HG3	1.73	0.51
1:B:164:THR:CG2	1:B:165:ALA:N	2.59	0.51
1:A:73:PRO:HB2	1:A:222:HIS:NE2	2.25	0.51
1:B:145:THR:H	1:B:148:GLN:NE2	2.08	0.51
1:B:562:VAL:HG13	1:B:567:LYS:HA	1.92	0.51
1:B:266:GLN:HE21	1:B:267:ASP:H	1.58	0.51
1:A:494:LEU:HD13	1:A:539:VAL:HG22	1.93	0.50
1:A:52:THR:HG21	1:A:88:LEU:HD22	1.93	0.50
1:A:265:LEU:HD22	1:A:269:THR:HG21	1.93	0.50
1:B:256:GLY:HA2	1:B:305:VAL:HG22	1.92	0.50
1:B:599:THR:O	1:B:602:MET:HG2	2.11	0.50
1:B:179:PHE:HB3	1:B:183:ARG:NH1	2.26	0.50
1:B:394:ILE:CG2	1:B:403:ARG:HG2	2.41	0.50
1:A:200:ASN:ND2	1:A:202:GLU:H	2.08	0.50
1:A:382:LYS:HE3	1:A:384:MET:HB3	1.93	0.50
1:B:243:GLN:HA	1:B:243:GLN:OE1	2.12	0.50
1:A:200:ASN:HD22	1:A:202:GLU:N	2.10	0.49
1:B:254:VAL:HG13	1:B:470:MET:CG	2.41	0.49
1:A:331:PRO:HA	1:A:415:THR:HB	1.94	0.49
1:B:200:ASN:ND2	1:B:202:GLU:HB2	2.28	0.49
1:B:373:SER:HB3	1:B:403:ARG:NH2	2.27	0.49
1:B:223:LEU:HB3	1:B:561:VAL:HG21	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:PRO:HB2	1:B:222:HIS:CE1	2.48	0.48
1:B:163:SER:HB3	1:B:179:PHE:HZ	1.76	0.48
1:A:69:ARG:HD3	1:A:155:MET:HE1	1.95	0.48
1:B:69:ARG:HG2	1:B:594:TYR:CE2	2.49	0.48
1:B:89:ARG:O	1:B:93:GLU:HG3	2.13	0.47
1:A:177:ALA:O	1:A:181:ILE:HG12	2.14	0.47
1:A:588:GLU:O	1:A:592:THR:HG23	2.15	0.47
1:A:477:MET:HE1	1:A:498:SER:HA	1.97	0.47
1:B:37:ASP:HA	1:B:61:THR:O	2.14	0.47
1:B:477:MET:CE	1:B:498:SER:HA	2.44	0.47
1:B:44:VAL:HG13	1:B:57:PHE:CE2	2.50	0.47
1:A:422:GLN:HA	1:A:536:ALA:HB1	1.97	0.47
1:A:546:GLU:OE2	1:A:546:GLU:HA	2.15	0.46
1:B:276:ARG:NH1	1:B:442:ALA:HA	2.31	0.46
1:A:283:GLU:OE1	1:A:442:ALA:HB1	2.16	0.46
1:B:343:ALA:O	1:B:364:LYS:HB2	2.16	0.46
1:B:454:PHE:CE1	1:B:545:MET:HG2	2.50	0.46
1:B:490:ASP:OD1	1:B:492:GLU:HG2	2.16	0.45
1:A:300:PHE:CZ	1:A:449:SER:HA	2.52	0.45
1:A:338:ARG:HG2	1:A:413:VAL:HG22	1.98	0.45
1:A:220:LEU:HD21	1:A:464:VAL:HG13	1.99	0.44
1:B:325:GLN:HE21	1:B:325:GLN:HA	1.82	0.44
1:A:103:ASP:O	1:A:107:ARG:HB2	2.18	0.44
1:B:113:ARG:H	1:B:113:ARG:HG2	1.50	0.44
1:B:331:PRO:HA	1:B:415:THR:OG1	2.17	0.44
1:A:160:LEU:HD23	1:A:160:LEU:N	2.32	0.44
1:B:262:VAL:HA	1:B:296:ILE:O	2.17	0.44
1:B:252:VAL:HG13	1:B:465:ILE:HD13	1.99	0.43
1:B:266:GLN:HE21	1:B:266:GLN:HA	1.83	0.43
1:A:200:ASN:C	1:A:200:ASN:ND2	2.76	0.43
1:B:279:ASN:HD21	1:B:281:GLN:HB2	1.83	0.43
1:B:217:PHE:CD1	1:B:262:VAL:HG12	2.53	0.43
1:B:203:TRP:NE1	1:B:582:ARG:HG2	2.34	0.43
1:B:71:ILE:HD12	1:B:583:VAL:HG22	2.01	0.43
1:B:157:VAL:HG22	1:B:197:PHE:CD1	2.54	0.43
1:B:26:GLN:O	1:B:30:ASP:HB2	2.19	0.42
1:B:271:LEU:HB3	1:B:275:ILE:HG22	2.02	0.42
1:B:164:THR:HG23	1:B:212:GLU:OE1	2.20	0.42
1:B:338:ARG:HG2	1:B:413:VAL:HG12	2.01	0.42
1:A:126:ARG:O	1:A:130:GLN:HB2	2.20	0.42
1:B:157:VAL:HG22	1:B:197:PHE:HD1	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:GLU:HA	1:A:181:ILE:HG21	2.02	0.41
1:B:376:ASN:HA	1:B:377:PRO:HD3	1.95	0.41
1:A:314:LYS:HE2	1:A:393:LYS:O	2.20	0.41
1:A:220:LEU:HD23	1:A:251:VAL:HB	2.03	0.41
1:A:366:LYS:HB3	1:A:367:PRO:HD3	2.03	0.41
1:B:509:ASN:HB3	1:B:511:ASP:H	1.86	0.41
1:A:263:PHE:HD2	1:A:463:MET:HE1	1.85	0.41
1:A:527:MET:HE2	1:A:527:MET:HA	2.03	0.40
1:B:160:LEU:HD23	1:B:160:LEU:N	2.36	0.40
1:B:512:GLY:O	1:B:513:LEU:HB2	2.22	0.40
1:A:52:THR:HG23	1:A:87:GLY:HA2	2.03	0.40
1:A:118:VAL:N	1:A:119:PRO:CD	2.84	0.40
1:A:25:LEU:HD12	1:A:25:LEU:HA	1.96	0.40
1:A:495:ILE:O	1:A:499:ARG:HG3	2.21	0.40
1:A:311:LEU:HD11	1:A:436:LEU:HD21	2.04	0.40
1:B:161:LEU:O	1:B:204:THR:HA	2.21	0.40
1:B:179:PHE:CE2	1:B:206:PRO:HG3	2.57	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	589/602 (98%)	564 (96%)	25 (4%)	0	100	100
1	B	589/602 (98%)	556 (94%)	32 (5%)	1 (0%)	43	72
All	All	1178/1204 (98%)	1120 (95%)	57 (5%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	568	PRO

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	494/502 (98%)	442 (90%)	52 (10%)	6	22
1	B	494/502 (98%)	447 (90%)	47 (10%)	8	26
All	All	988/1004 (98%)	889 (90%)	99 (10%)	7	24

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLU
1	A	18	ILE
1	A	25	LEU
1	A	33	ARG
1	A	43	VAL
1	A	44	VAL
1	A	52	THR
1	A	53	ARG
1	A	88	LEU
1	A	92	LEU
1	A	106	SER
1	A	113	ARG
1	A	130	GLN
1	A	147	GLU
1	A	150	VAL
1	A	157	VAL
1	A	160	LEU
1	A	171	HIS
1	A	175	ARG
1	A	182	LEU
1	A	185	LEU
1	A	200	ASN
1	A	223	LEU
1	A	242	LYS
1	A	254	VAL
1	A	260	GLN
1	A	269	THR

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Mol	Chain	Res	Type
1	A	275	ILE
1	A	278	VAL
1	A	311	LEU
1	A	323	ARG
1	A	325	GLN
1	A	353	THR
1	A	363	ASN
1	A	364	LYS
1	A	370	THR
1	A	385	ASN
1	A	400	LEU
1	A	429	LEU
1	A	436	LEU
1	A	439	ASN
1	A	464	VAL
1	A	482	LEU
1	A	494	LEU
1	A	511	ASP
1	A	516	GLN
1	A	523	LYS
1	A	546	GLU
1	A	561	VAL
1	A	576	VAL
1	A	583	VAL
1	A	588	GLU
1	B	30	ASP
1	B	33	ARG
1	B	43	VAL
1	B	44	VAL
1	B	63	SER
1	B	69	ARG
1	B	77	ARG
1	B	88	LEU
1	B	91	VAL
1	B	92	LEU
1	B	96	LEU
1	B	103	ASP
1	B	113	ARG
1	B	132	LEU
1	B	157	VAL
1	B	159	VAL
1	B	160	LEU

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Mol	Chain	Res	Type
1	B	167	VAL
1	B	182	LEU
1	B	245	ARG
1	B	252	VAL
1	B	254	VAL
1	B	266	GLN
1	B	271	LEU
1	B	275	ILE
1	B	284	ARG
1	B	295	VAL
1	B	301	MET
1	B	325	GLN
1	B	355	VAL
1	B	413	VAL
1	B	422	GLN
1	B	428	LEU
1	B	461	SER
1	B	464	VAL
1	B	480	LEU
1	B	506	VAL
1	B	507	VAL
1	B	509	ASN
1	B	516	GLN
1	B	534	THR
1	B	537	LEU
1	B	541	LEU
1	B	559	LYS
1	B	562	VAL
1	B	583	VAL
1	B	599	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	200	ASN
1	A	243	GLN
1	A	319	ASN
1	A	325	GLN
1	A	339	GLN
1	A	363	ASN
1	A	375	ASN

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Mol	Chain	Res	Type
1	A	422	GLN
1	A	439	ASN
1	A	478	GLN
1	A	516	GLN
1	A	522	GLN
1	A	543	GLN
1	A	544	HIS
1	B	116	GLN
1	B	148	GLN
1	B	153	GLN
1	B	200	ASN
1	B	221	ASN
1	B	260	GLN
1	B	266	GLN
1	B	279	ASN
1	B	325	GLN
1	B	385	ASN
1	B	422	GLN
1	B	424	GLN
1	B	478	GLN
1	B	488	GLN
1	B	508	ASN
1	B	509	ASN
1	B	516	GLN
1	B	522	GLN
1	B	543	GLN
1	B	558	GLN
1	B	575	HIS
1	B	601	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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, 95th 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 ²)	Q<0.9
1	A	591/602 (98%)	0.19	13 (2%) 62 52	60, 86, 146, 184	0
1	B	591/602 (98%)	0.20	9 (1%) 72 63	59, 87, 134, 195	0
All	All	1182/1204 (98%)	0.20	22 (1%) 66 57	59, 86, 139, 195	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	12	ALA	5.6
1	B	12	ALA	4.8
1	A	109	VAL	4.3
1	A	438	ALA	3.4
1	A	234	CYS	3.1
1	A	437	PRO	2.9
1	B	313	PHE	2.7
1	B	109	VAL	2.7
1	A	410	PHE	2.6
1	B	171	HIS	2.5
1	B	383	ALA	2.4
1	A	383	ALA	2.3
1	B	438	ALA	2.3
1	B	512	GLY	2.2
1	A	384	MET	2.2
1	A	302	GLN	2.2
1	A	318	SER	2.2
1	B	311	LEU	2.2
1	A	14	THR	2.1
1	A	50	SER	2.1
1	A	16	LYS	2.1
1	B	370	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.