



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 2WB7 / pdb_00002wb7
Title : pT26-6p
Authors : Keller, J.; Leulliot, N.; Soler, N.; Collinet, B.; Vincentelli, R.; Forterre, P.; van Tilbeurgh, H.
Deposited on : 2009-02-22
Resolution : 2.60 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

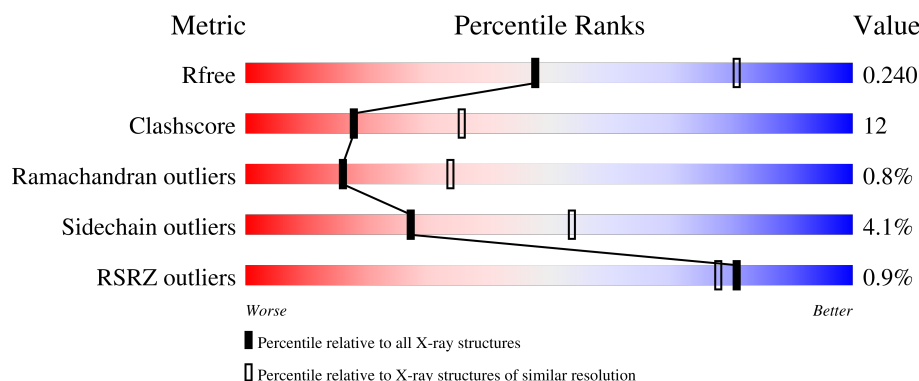
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	526	76% 22% .
1	B	526	69% 28% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8499 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 PT26-6P.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	525	Total	C	N	O	S	Se	0	0	0
			4137	2630	687	807	2	11			
1	B	525	Total	C	N	O	S	Se	0	0	0
			4137	2630	687	807	2	11			

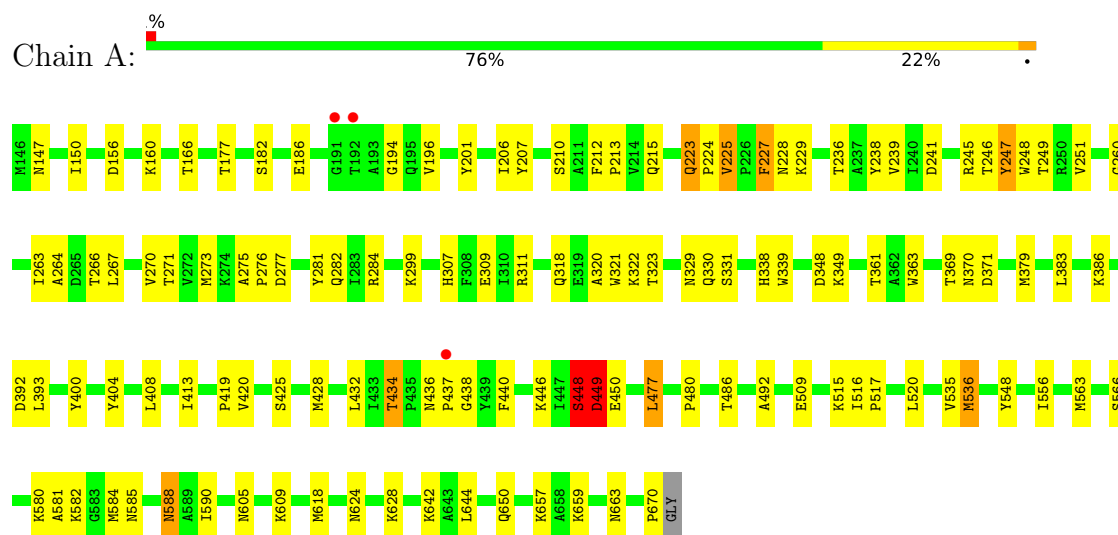
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	115	Total	O	0	0
			115	115		
2	B	110	Total	O	0	0
			110	110		

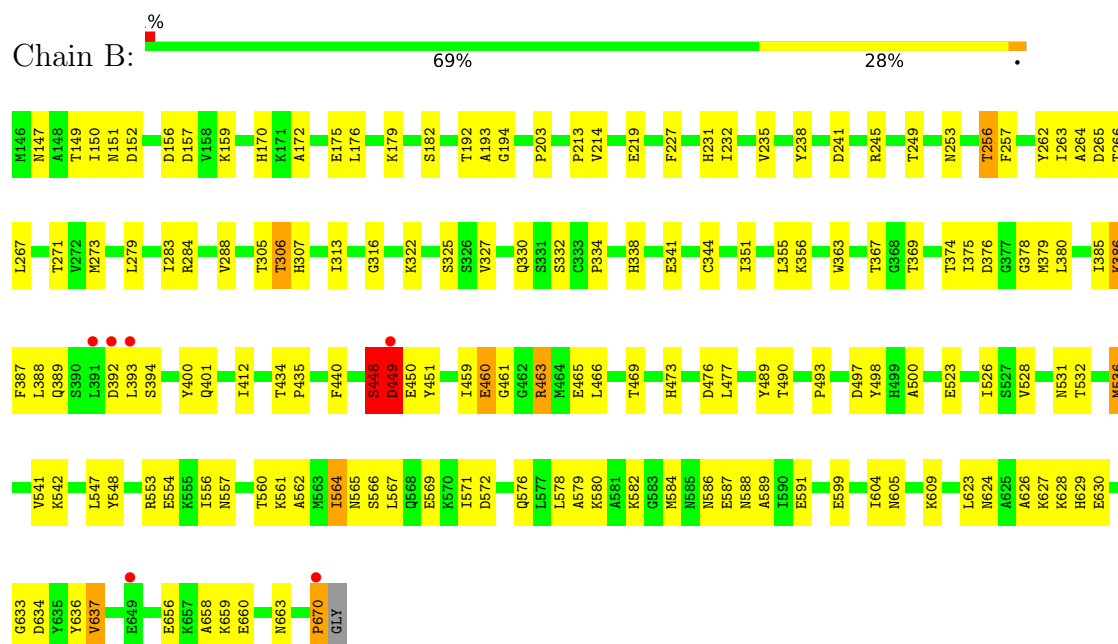
3 Residue-property plots

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: PT26-6P



• Molecule 1: PT26-6P



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	133.15Å 133.15Å 164.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.19 – 2.60 47.19 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.19-2.60) 99.3 (47.19-2.60)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, R_{free}	0.193 , 0.240 0.193 , 0.240	Depositor DCC
R_{free} test set	2576 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	48.7	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.048 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8499	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% 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

5.1 Standard geometry

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	0.59	0/4220	0.88	10/5723 (0.2%)
1	B	0.56	0/4220	0.86	2/5723 (0.0%)
All	All	0.58	0/8440	0.87	12/11446 (0.1%)

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	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	449	ASP	N-CA-C	10.07	132.24	110.80
1	A	449	ASP	N-CA-C	9.74	131.55	110.80
1	A	225	VAL	CA-C-N	7.01	128.61	119.84
1	A	225	VAL	C-N-CA	7.01	128.61	119.84
1	A	450	GLU	N-CA-C	6.01	119.19	109.40
1	A	448	SER	N-CA-C	5.80	118.92	108.48
1	A	277	ASP	CA-C-N	5.62	124.97	118.85
1	A	277	ASP	C-N-CA	5.62	124.97	118.85
1	A	492	ALA	CA-C-N	-5.57	113.82	119.83
1	A	492	ALA	C-N-CA	-5.57	113.82	119.83
1	A	225	VAL	N-CA-C	5.38	113.35	107.60
1	B	450	GLU	N-CA-C	5.38	118.50	109.95

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	448	SER	Peptide
1	B	448	SER	Peptide

5.2 Too-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 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	4137	0	4029	94	0
1	B	4137	0	4029	106	0
2	A	115	0	0	6	0
2	B	110	0	0	7	0
All	All	8499	0	8058	200	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 12.

All (200) 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:A:207:TYR:OH	1:A:370:ASN:HB3	1.59	1.00
1:A:446:LYS:HE2	1:A:509:GLU:HB2	1.52	0.91
1:A:282:GLN:HE22	1:A:299:LYS:NZ	1.67	0.90
1:B:599:GLU:OE1	1:B:629:HIS:HD2	1.57	0.87
1:B:307:HIS:HD2	1:B:363:TRP:CD2	1.95	0.84
1:B:147:ASN:O	1:B:151:ASN:HB2	1.86	0.76
1:A:282:GLN:HE22	1:A:299:LYS:HZ3	1.33	0.75
1:A:160:LYS:HD2	2:A:2003:HOH:O	1.85	0.74
1:A:642:LYS:NZ	1:A:650:GLN:HE21	1.85	0.74
1:A:392:ASP:HB2	2:A:2037:HOH:O	1.88	0.73
1:A:605:ASN:HD21	1:A:609:LYS:HE2	1.53	0.72
1:B:367:THR:OG1	1:B:369:THR:HG22	1.89	0.72
1:A:213:PRO:HG3	1:A:440:PHE:HA	1.73	0.70
1:B:567:LEU:O	1:B:571:ILE:HG13	1.92	0.70
1:A:177:THR:HG21	1:A:437:PRO:HD2	1.74	0.69
1:B:493:PRO:HA	2:B:2093:HOH:O	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:HIS:CD2	1:B:363:TRP:CE2	2.81	0.69
1:B:307:HIS:CD2	1:B:363:TRP:CD2	2.80	0.68
1:A:282:GLN:HE22	1:A:299:LYS:HZ1	1.42	0.66
1:A:236:THR:OG1	1:A:251:VAL:HG22	1.96	0.66
1:A:434:THR:HG23	1:A:480:PRO:HG3	1.78	0.66
1:B:463:ARG:HD2	2:B:2082:HOH:O	1.96	0.65
1:B:498:TYR:HB2	1:B:528:VAL:HG21	1.78	0.65
1:A:370:ASN:HD22	1:A:404:TYR:H	1.45	0.65
1:A:605:ASN:ND2	1:A:609:LYS:HE2	2.12	0.64
1:B:564:ILE:HG23	1:B:604:ILE:HG23	1.80	0.64
1:B:605:ASN:O	1:B:609:LYS:HG2	1.98	0.64
1:A:207:TYR:HE1	1:A:370:ASN:HA	1.63	0.64
1:A:428:MSE:HA	1:A:428:MSE:HE2	1.80	0.63
1:A:238:TYR:HB3	1:A:249:THR:HA	1.80	0.63
1:B:379:MSE:HG3	1:B:380:LEU:N	2.13	0.62
1:B:586:ASN:ND2	1:B:589:ALA:HB2	2.15	0.61
1:B:149:THR:HG22	1:B:541:VAL:HG11	1.82	0.61
1:B:656:GLU:O	1:B:660:GLU:HG3	1.99	0.61
1:A:330:GLN:HG2	1:A:339:TRP:CG	2.35	0.60
1:B:219:GLU:HG2	1:B:266:THR:HG22	1.83	0.60
1:A:207:TYR:CD2	1:A:408:LEU:HB3	2.36	0.60
1:A:428:MSE:HE1	1:A:486:THR:HG23	1.82	0.60
1:A:228:ASN:HD21	1:A:321:TRP:H	1.49	0.59
1:B:560:THR:O	1:B:564:ILE:HG13	2.02	0.59
1:A:477:LEU:H	1:A:477:LEU:HD23	1.69	0.57
1:A:330:GLN:HG2	1:A:339:TRP:CD2	2.38	0.57
1:A:659:LYS:HE2	1:A:663:ASN:HD21	1.69	0.57
1:B:341:GLU:OE2	1:B:341:GLU:HA	2.03	0.57
1:B:567:LEU:HD21	1:B:630:GLU:HG3	1.86	0.57
1:A:329:ASN:ND2	1:A:331:SER:HB2	2.20	0.57
1:A:446:LYS:HE2	1:A:509:GLU:CB	2.31	0.56
1:B:241:ASP:C	1:B:241:ASP:OD2	2.48	0.56
1:B:378:GLY:HA3	1:B:401:GLN:NE2	2.21	0.55
1:B:659:LYS:HE3	1:B:663:ASN:HD21	1.71	0.55
1:A:436:ASN:OD1	1:A:436:ASN:C	2.50	0.55
1:B:192:THR:HG23	1:B:355:LEU:HD22	1.89	0.55
1:B:271:THR:HG21	1:B:273:MSE:HE3	1.89	0.55
1:B:500:ALA:O	1:B:523:GLU:HG3	2.06	0.55
1:B:554:GLU:HA	1:B:554:GLU:OE1	2.06	0.55
1:A:281:TYR:CD1	1:A:284:ARG:NH1	2.74	0.54
1:A:206:ILE:HD13	1:A:212:PHE:CD2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:TYR:OH	1:A:370:ASN:CB	2.47	0.54
1:A:379:MSE:HG3	2:A:2003:HOH:O	2.07	0.54
1:A:307:HIS:CD2	1:A:363:TRP:CD2	2.96	0.53
1:A:548:TYR:HA	1:A:556:ILE:HD11	1.89	0.53
1:B:587:GLU:O	1:B:591:GLU:HG3	2.08	0.53
1:A:238:TYR:OH	1:A:311:ARG:HD2	2.08	0.53
1:A:177:THR:CG2	1:A:437:PRO:HD2	2.38	0.53
1:A:624:ASN:O	1:A:628:LYS:HG2	2.08	0.53
1:A:227:PHE:CD2	1:A:338:HIS:CD2	2.97	0.53
1:B:213:PRO:HG3	1:B:440:PHE:HA	1.89	0.53
1:B:214:VAL:HG21	1:B:273:MSE:HG3	1.91	0.53
1:B:548:TYR:CA	1:B:556:ILE:HD11	2.39	0.53
1:B:284:ARG:O	1:B:288:VAL:HG23	2.08	0.53
1:B:633:GLY:O	1:B:637:VAL:HG13	2.09	0.53
1:B:238:TYR:HB3	1:B:249:THR:HA	1.91	0.53
1:B:459:ILE:HD13	1:B:463:ARG:NH2	2.24	0.53
1:B:176:LEU:HD23	1:B:412:ILE:HD12	1.90	0.52
1:B:459:ILE:HD11	1:B:465:GLU:CD	2.34	0.52
1:B:262:TYR:CE2	1:B:267:LEU:HD21	2.45	0.52
1:B:152:ASP:OD1	1:B:628:LYS:HE2	2.10	0.51
1:A:247:TYR:N	1:A:247:TYR:CD1	2.77	0.51
1:A:624:ASN:HD21	1:A:628:LYS:NZ	2.08	0.51
1:A:223:GLN:HB2	1:A:263:ILE:HD11	1.93	0.51
1:A:223:GLN:HG3	1:A:224:PRO:HD2	1.93	0.51
1:B:562:ALA:O	1:B:565:ASN:HB2	2.11	0.51
1:A:322:LYS:HE3	1:A:348:ASP:OD2	2.11	0.51
1:B:256:THR:HG22	2:B:2031:HOH:O	2.10	0.50
1:A:536:MSE:HE3	1:A:563:MSE:SE	2.61	0.50
1:A:271:THR:HG21	1:A:273:MSE:HE3	1.92	0.50
1:A:215:GLN:HG2	1:A:270:VAL:HG22	1.94	0.50
1:A:448:SER:O	1:A:449:ASP:HB2	2.12	0.50
1:B:305:THR:HG22	1:B:306:THR:N	2.27	0.50
1:B:459:ILE:HD13	1:B:463:ARG:CZ	2.42	0.49
1:A:227:PHE:N	1:A:227:PHE:CD1	2.81	0.49
1:B:327:VAL:HG13	1:B:332:SER:HB2	1.94	0.49
1:A:241:ASP:OD1	1:A:245:ARG:HB3	2.13	0.49
1:B:587:GLU:OE2	1:B:587:GLU:HA	2.12	0.48
1:A:322:LYS:HG2	1:A:348:ASP:HB2	1.93	0.48
1:B:400:TYR:CE2	1:B:526:ILE:HG22	2.48	0.48
1:A:330:GLN:HG2	1:A:339:TRP:CD1	2.48	0.48
1:B:553:ARG:O	1:B:557:ASN:ND2	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:413:ILE:HG13	1:A:520:LEU:HD23	1.96	0.47
1:A:194:GLY:C	1:A:196:VAL:H	2.22	0.47
1:B:367:THR:HG1	1:B:369:THR:HG22	1.79	0.47
1:B:624:ASN:ND2	1:B:628:LYS:HE3	2.29	0.47
1:B:448:SER:O	1:B:449:ASP:HB2	2.15	0.47
1:A:281:TYR:HD1	1:A:284:ARG:NH1	2.12	0.47
1:A:369:THR:O	1:A:369:THR:HG23	2.14	0.47
1:B:150:ILE:HG12	1:B:542:LYS:HG3	1.97	0.47
1:B:170:HIS:HE1	2:B:2008:HOH:O	1.97	0.47
1:B:497:ASP:OD1	1:B:497:ASP:C	2.58	0.47
1:B:227:PHE:CE1	1:B:338:HIS:CD2	3.03	0.46
1:A:379:MSE:HE2	1:A:400:TYR:CZ	2.50	0.46
1:A:309:GLU:OE2	1:A:361:THR:N	2.47	0.46
1:B:235:VAL:HG23	1:B:257:PHE:CZ	2.51	0.46
1:B:451:TYR:CD1	1:B:451:TYR:C	2.92	0.46
1:A:147:ASN:OD1	1:A:150:ILE:N	2.46	0.46
1:A:438:GLY:HA2	2:A:2052:HOH:O	2.15	0.46
1:B:547:LEU:HB3	1:B:556:ILE:HG13	1.96	0.46
1:A:670:PRO:C	2:A:2115:HOH:O	2.58	0.46
1:B:170:HIS:HD2	2:B:2071:HOH:O	1.98	0.46
1:B:385:ILE:O	1:B:386:LYS:C	2.58	0.46
1:B:245:ARG:HB2	1:B:245:ARG:NH1	2.30	0.45
1:B:578:LEU:CD1	1:B:582:LYS:HE3	2.46	0.45
1:A:383:LEU:HD12	1:A:419:PRO:HB2	1.98	0.45
1:A:275:ALA:HB1	1:A:276:PRO:HD2	1.97	0.45
1:A:618:MSE:HE2	1:A:670:PRO:HG3	1.98	0.45
1:B:466:LEU:HD11	1:B:469:THR:HB	1.98	0.45
1:A:642:LYS:HZ3	1:A:650:GLN:HE21	1.64	0.45
1:A:509:GLU:HG3	1:A:515:LYS:HG3	1.99	0.45
1:A:581:ALA:HA	1:A:644:LEU:HD21	1.99	0.45
1:B:580:LYS:O	1:B:584:MSE:HG2	2.17	0.45
1:B:599:GLU:OE1	1:B:629:HIS:CD2	2.50	0.45
1:A:642:LYS:HZ1	1:A:650:GLN:HE21	1.62	0.45
1:B:548:TYR:HA	1:B:556:ILE:HD11	1.97	0.44
1:A:228:ASN:HD21	1:A:321:TRP:N	2.14	0.44
1:B:588:ASN:O	1:B:589:ALA:C	2.60	0.44
1:B:313:ILE:HG13	1:B:356:LYS:HG2	1.99	0.44
1:A:584:MSE:O	1:A:585:ASN:C	2.60	0.44
1:B:172:ALA:HB1	1:B:375:ILE:HG21	2.00	0.44
1:B:576:GLN:O	1:B:579:ALA:HB3	2.17	0.44
1:A:229:LYS:HD3	1:A:260:GLY:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:VAL:HG12	1:A:227:PHE:HD1	1.83	0.44
1:A:428:MSE:CE	1:A:486:THR:HG23	2.46	0.44
1:B:192:THR:O	1:B:194:GLY:N	2.51	0.43
1:B:578:LEU:HG	1:B:582:LYS:HE3	2.00	0.43
1:A:413:ILE:HA	1:A:432:LEU:O	2.17	0.43
1:B:192:THR:HG23	1:B:355:LEU:CD2	2.48	0.43
1:B:557:ASN:HB3	1:B:561:LYS:NZ	2.33	0.43
1:B:231:HIS:O	1:B:316:GLY:HA3	2.18	0.43
1:B:536:MSE:HE2	1:B:536:MSE:HB3	1.80	0.43
1:B:548:TYR:N	1:B:556:ILE:HD11	2.33	0.43
1:A:186:GLU:HG3	1:A:201:TYR:CE1	2.53	0.43
1:A:659:LYS:HE2	1:A:663:ASN:ND2	2.34	0.43
1:B:279:LEU:HB3	1:B:283:ILE:CD1	2.48	0.43
1:B:434:THR:OG1	1:B:435:PRO:HD2	2.19	0.43
1:B:489:TYR:O	1:B:490:THR:C	2.61	0.43
1:A:248:TRP:O	1:A:249:THR:C	2.61	0.43
1:A:266:THR:O	1:A:267:LEU:HD23	2.19	0.43
1:A:580:LYS:NZ	2:A:2092:HOH:O	2.45	0.43
1:A:223:GLN:CD	1:A:318:GLN:HE22	2.26	0.43
1:A:330:GLN:HG2	1:A:339:TRP:CE2	2.54	0.43
1:B:156:ASP:O	1:B:159:LYS:N	2.45	0.43
1:B:219:GLU:CG	1:B:266:THR:HG22	2.48	0.43
1:B:263:ILE:O	1:B:264:ALA:C	2.61	0.43
1:B:327:VAL:HB	1:B:344:CYS:HB2	2.01	0.43
1:B:182:SER:HB3	1:B:203:PRO:HB3	2.00	0.42
1:A:582:LYS:HA	1:A:590:ILE:HD11	2.02	0.42
1:B:392:ASP:C	1:B:394:SER:N	2.77	0.42
1:B:636:TYR:CE2	1:B:658:ALA:HB1	2.54	0.42
1:B:634:ASP:O	1:B:637:VAL:HG22	2.19	0.42
1:A:371:ASP:OD1	1:A:371:ASP:N	2.51	0.42
1:B:253:ASN:O	1:B:256:THR:HG23	2.20	0.42
1:A:329:ASN:HD22	1:A:331:SER:HB2	1.84	0.42
1:B:175:GLU:O	1:B:179:LYS:HG3	2.20	0.42
1:A:516:ILE:HA	1:A:517:PRO:HD3	1.87	0.42
1:B:460:GLU:O	1:B:461:GLY:C	2.62	0.42
1:B:566:SER:O	1:B:567:LEU:C	2.61	0.42
1:B:392:ASP:C	1:B:394:SER:H	2.28	0.42
1:A:657:LYS:HA	1:A:657:LYS:HD2	1.75	0.41
1:B:387:PHE:C	1:B:389:GLN:N	2.76	0.41
1:B:531:ASN:ND2	1:B:532:THR:O	2.53	0.41
1:B:553:ARG:HD2	1:B:553:ARG:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:253:ASN:OD1	1:B:253:ASN:C	2.62	0.41
1:B:219:GLU:HG2	1:B:266:THR:CG2	2.50	0.41
1:A:207:TYR:CE1	1:A:370:ASN:HA	2.50	0.41
1:A:535:VAL:HG22	1:A:566:SER:HB3	2.01	0.41
1:B:400:TYR:CZ	1:B:526:ILE:HG22	2.55	0.41
1:B:374:THR:HB	2:B:2058:HOH:O	2.21	0.41
1:B:623:LEU:O	1:B:626:ALA:HB3	2.20	0.41
1:A:239:VAL:O	1:A:246:THR:HA	2.21	0.41
1:A:275:ALA:HB1	1:A:276:PRO:CD	2.51	0.41
1:B:449:ASP:OD2	1:B:473:HIS:ND1	2.53	0.41
1:A:584:MSE:HE3	1:A:584:MSE:HB3	1.92	0.41
1:A:588:ASN:ND2	1:A:588:ASN:H	2.18	0.41
1:B:670:PRO:C	2:B:2110:HOH:O	2.63	0.41
1:A:320:ALA:HB3	1:A:349:LYS:HG2	2.03	0.40
1:A:215:GLN:CG	1:A:270:VAL:HG22	2.51	0.40
1:B:459:ILE:HG22	1:B:460:GLU:N	2.37	0.40
1:B:156:ASP:O	1:B:157:ASP:C	2.62	0.40
1:B:476:ASP:OD1	1:B:476:ASP:C	2.62	0.40
1:A:194:GLY:C	1:A:196:VAL:N	2.79	0.40
1:B:569:GLU:O	1:B:572:ASP:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	523/526 (99%)	503 (96%)	18 (3%)	2 (0%)	30	51
1	B	523/526 (99%)	483 (92%)	34 (6%)	6 (1%)	11	25
All	All	1046/1052 (99%)	986 (94%)	52 (5%)	8 (1%)	16	34

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	449	ASP
1	B	449	ASP
1	B	193	ALA
1	B	393	LEU
1	A	264	ALA
1	B	330	GLN
1	B	376	ASP
1	B	334	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	438/427 (103%)	421 (96%)	17 (4%)	28	55
1	B	438/427 (103%)	419 (96%)	19 (4%)	26	51
All	All	876/854 (103%)	840 (96%)	36 (4%)	27	54

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	156	ASP
1	A	166	THR
1	A	182	SER
1	A	210	SER
1	A	223	GLN
1	A	227	PHE
1	A	247	TYR
1	A	323	THR
1	A	386	LYS
1	A	393	LEU
1	A	420	VAL
1	A	425	SER
1	A	434	THR
1	A	448	SER
1	A	477	LEU
1	A	536	MSE

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Mol	Chain	Res	Type
1	A	588	ASN
1	B	232	ILE
1	B	256	THR
1	B	265	ASP
1	B	306	THR
1	B	322	LYS
1	B	325	SER
1	B	351	ILE
1	B	386	LYS
1	B	388	LEU
1	B	448	SER
1	B	449	ASP
1	B	460	GLU
1	B	463	ARG
1	B	477	LEU
1	B	536	MSE
1	B	564	ILE
1	B	627	LYS
1	B	637	VAL
1	B	670	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	ASN
1	A	215	GLN
1	A	223	GLN
1	A	228	ASN
1	A	244	ASN
1	A	282	GLN
1	A	318	GLN
1	A	338	HIS
1	A	576	GLN
1	A	588	ASN
1	A	605	ASN
1	A	613	GLN
1	A	624	ASN
1	A	650	GLN
1	A	663	ASN
1	B	170	HIS
1	B	195	GLN
1	B	243	ASN

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Mol	Chain	Res	Type
1	B	307	HIS
1	B	330	GLN
1	B	338	HIS
1	B	531	ASN
1	B	557	ASN
1	B	568	GLN
1	B	594	GLN
1	B	624	ASN
1	B	629	HIS
1	B	663	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	514/526 (97%)	-0.23	3 (0%) 85 83	28, 44, 69, 82	0
1	B	514/526 (97%)	-0.06	6 (1%) 76 73	31, 50, 72, 82	0
All	All	1028/1052 (97%)	-0.14	9 (0%) 81 78	28, 46, 71, 82	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	392	ASP	3.7
1	A	191	GLY	3.0
1	B	393	LEU	2.5
1	B	391	LEU	2.4
1	B	449	ASP	2.2
1	B	649	GLU	2.0
1	B	670	PRO	2.0
1	A	192	THR	2.0
1	A	437	PRO	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.