



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:46 PM UTC

PDB ID : 4U0Q / pdb_00004u0q
Title : Plasmodium falciparum reticulocyte-binding protein homologue 5 (PfRH5) bound to basigin
Authors : Wright, K.E.; Hjerrild, K.A.; Bartlett, J.; Douglas, A.D.; Jin, J.; Brown, R.E.; Ashfield, R.; Clemmensen, S.B.; de Jongh, W.A.; Draper, S.J.; Higgins, M.K.
Deposited on : 2014-07-14
Resolution : 3.10 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

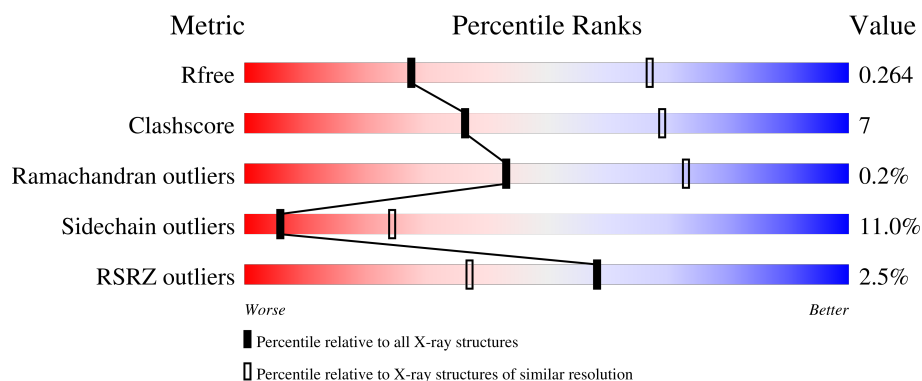
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	2% 41% 12% • 46%
1	C	526	39% 13% • 46%
2	B	269	2% 48% 17% • 33%
2	D	269	3% 50% 15% • 33%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7629 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 Reticulocyte binding protein 5.

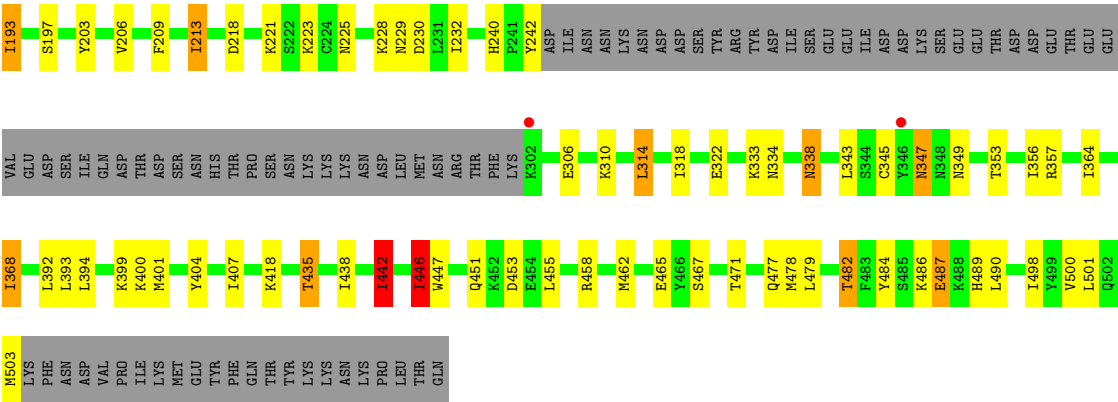
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	S	0	0	0
			2430	1569	407	439	15			
1	C	285	Total	C	N	O	S	0	0	0
			2421	1563	405	438	15			

There are 2 discrepancies between the modelled and reference sequences:

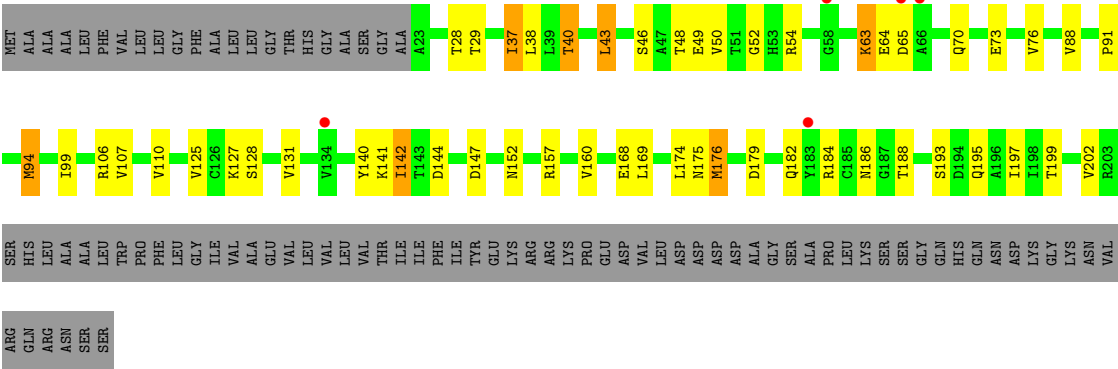
Chain	Residue	Modelled	Actual	Comment	Reference
A	216	ALA	THR	engineered mutation	UNP B2L3N7
C	216	ALA	THR	engineered mutation	UNP B2L3N7

- Molecule 2 is a protein called Basigin.

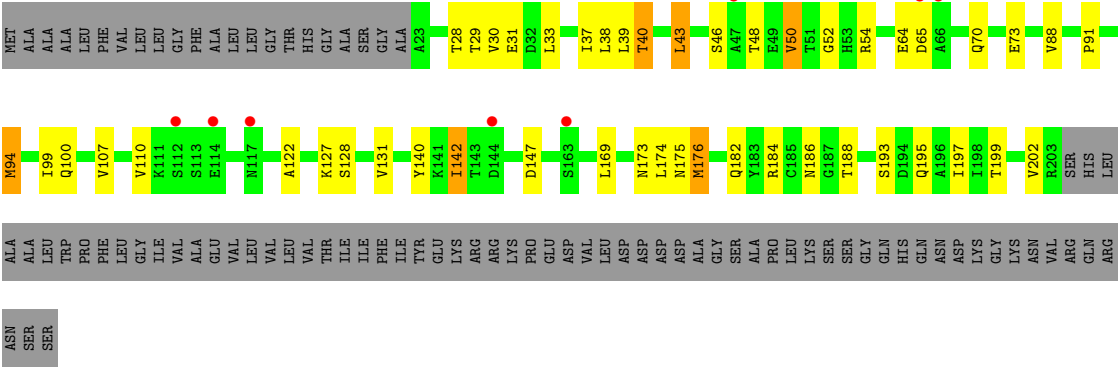
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	181	Total	C	N	O	S	0	0	0
			1389	862	238	281	8			
2	D	181	Total	C	N	O	S	0	0	0
			1389	862	238	281	8			



• Molecule 2: Basigin



• Molecule 2: Basigin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.28Å 109.36Å 151.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	62.40 – 3.10 62.40 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.7 (62.40-3.10) 99.7 (62.40-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 3.13Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.217 , 0.244 0.232 , 0.264	Depositor DCC
R_{free} test set	1186 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	48.0	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7629	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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	0.72	1/2481 (0.0%)	1.45	8/3327 (0.2%)
1	C	0.74	0/2472	1.46	6/3316 (0.2%)
2	B	0.73	0/1416	1.20	1/1918 (0.1%)
2	D	0.72	0/1416	1.18	1/1918 (0.1%)
All	All	0.73	1/7785 (0.0%)	1.36	16/10479 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	176	ILE	CA-CB	5.33	1.56	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	442	ILE	N-CA-CB	7.28	119.06	110.55
1	C	442	ILE	N-CA-CB	6.90	118.62	110.55
1	A	446	ILE	N-CA-C	6.24	122.32	109.34
1	C	446	ILE	N-CA-C	5.96	121.73	109.34
1	C	175	ILE	CA-C-N	5.87	126.07	120.43
1	C	175	ILE	C-N-CA	5.87	126.07	120.43
1	C	225	ASN	N-CA-C	5.73	117.99	111.11
1	A	175	ILE	CA-C-N	5.69	125.89	120.43
1	A	175	ILE	C-N-CA	5.69	125.89	120.43
2	D	50	VAL	N-CA-CB	5.62	117.23	110.31
1	A	225	ASN	N-CA-C	5.54	118.04	111.33
1	C	364	ILE	N-CA-C	5.29	116.11	111.56
1	A	165	GLN	CA-C-N	5.16	127.19	120.28
1	A	165	GLN	C-N-CA	5.16	127.19	120.28
2	B	142	ILE	N-CA-C	5.09	114.94	108.12
1	A	364	ILE	N-CA-C	5.08	115.93	111.56

There are no chirality outliers.

There are no planarity outliers.

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	2430	0	2455	31	0
1	C	2421	0	2442	37	0
2	B	1389	0	1343	28	0
2	D	1389	0	1343	23	0
All	All	7629	0	7583	110	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 (110) 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:179:TYR:HD2	1:A:471:THR:HG22	1.14	1.11
1:A:447:TRP:HE3	2:B:28:THR:HG22	1.30	0.96
1:A:179:TYR:CD2	1:A:471:THR:HG22	2.05	0.89
1:C:479:LEU:HA	1:C:482:THR:HG23	1.66	0.77
1:C:180:THR:HG22	1:C:471:THR:HG21	1.70	0.73
2:B:54:ARG:HD3	2:B:63:LYS:HD3	1.74	0.70
1:A:447:TRP:CE3	2:B:28:THR:HG22	2.22	0.68
1:C:179:TYR:HD2	1:C:471:THR:HG22	1.57	0.68
2:B:152:ASN:HD22	2:B:160:VAL:H	1.41	0.67
2:B:54:ARG:CD	2:B:63:LYS:HD3	2.28	0.64
1:A:180:THR:HG22	1:A:471:THR:HG21	1.80	0.64
1:A:442:ILE:HG23	1:A:462:MET:HE2	1.80	0.62
1:C:442:ILE:HG23	1:C:462:MET:HE2	1.81	0.61
1:C:447:TRP:HE3	2:D:28:THR:HG22	1.64	0.61
1:C:467:SER:O	1:C:471:THR:HG23	2.02	0.60
1:A:446:ILE:HG23	1:A:447:TRP:CD1	2.37	0.59
1:A:192:SER:HA	1:A:343:LEU:HD12	1.85	0.59
2:D:52:GLY:HA3	2:D:65:ASP:HA	1.84	0.58
2:D:107:VAL:HG12	2:D:128:SER:HB2	1.85	0.58
1:C:446:ILE:HG23	1:C:447:TRP:CD1	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:107:VAL:HG12	2:B:128:SER:HB2	1.85	0.58
1:A:218:ASP:HA	1:A:221:LYS:HG2	1.87	0.56
1:A:168:GLU:OE1	1:A:240:HIS:NE2	2.39	0.56
2:B:37:ILE:HG22	2:B:76:VAL:HB	1.87	0.56
1:C:218:ASP:HA	1:C:221:LYS:HG2	1.88	0.55
2:D:28:THR:HG23	2:D:99:ILE:HG12	1.88	0.55
1:C:165:GLN:HG2	1:C:171:LEU:HD23	1.87	0.55
2:B:140:TYR:HB3	2:B:147:ASP:HB3	1.88	0.55
2:D:140:TYR:HB3	2:D:147:ASP:HB3	1.89	0.55
2:D:40:THR:HG23	2:D:73:GLU:HG2	1.89	0.54
1:A:484:TYR:HA	1:A:487:GLU:HB3	1.89	0.54
2:B:40:THR:HG23	2:B:73:GLU:HG2	1.90	0.54
2:B:91:PRO:HD2	2:B:94:MET:HE3	1.90	0.53
1:C:368:ILE:HG13	1:C:435:THR:HG22	1.90	0.53
2:D:91:PRO:HD2	2:D:94:MET:HE3	1.90	0.53
1:C:401:MET:HE2	1:C:404:TYR:HB2	1.90	0.53
1:C:484:TYR:HA	1:C:487:GLU:HB3	1.90	0.53
1:A:357:ARG:HA	1:A:446:ILE:HD11	1.92	0.51
1:A:394:LEU:HA	1:A:397:LEU:HD12	1.92	0.51
1:C:357:ARG:HA	1:C:446:ILE:HD11	1.92	0.51
2:B:38:LEU:HD11	2:B:73:GLU:HB3	1.93	0.50
2:D:38:LEU:HD11	2:D:73:GLU:HB3	1.93	0.50
2:B:152:ASN:HD22	2:B:160:VAL:N	2.09	0.50
1:A:209:PHE:CE2	1:A:213:ILE:HD12	2.47	0.49
1:C:192:SER:HA	1:C:343:LEU:HD12	1.95	0.48
2:D:110:VAL:HG22	2:D:127:LYS:HE3	1.95	0.48
1:C:500:VAL:HG23	1:C:501:LEU:HD12	1.95	0.48
1:C:168:GLU:OE1	1:C:240:HIS:NE2	2.36	0.48
2:B:28:THR:HG23	2:B:99:ILE:HG12	1.94	0.48
2:D:182:GLN:HG2	2:D:199:THR:HG22	1.96	0.48
1:A:496:HIS:HE1	2:D:33:LEU:HD13	1.79	0.48
1:A:447:TRP:HE3	2:B:28:THR:CG2	2.15	0.48
1:C:209:PHE:CE2	1:C:213:ILE:HD12	2.49	0.47
2:B:54:ARG:HB3	2:B:88:VAL:HB	1.96	0.47
1:C:203:TYR:CE1	2:D:131:VAL:HG22	2.50	0.47
2:B:52:GLY:HA3	2:B:65:ASP:HA	1.97	0.47
2:D:54:ARG:HB3	2:D:88:VAL:HB	1.97	0.47
1:A:447:TRP:HB3	2:B:28:THR:O	2.15	0.47
2:B:110:VAL:HG22	2:B:127:LYS:HE3	1.95	0.47
2:B:182:GLN:HG2	2:B:199:THR:HG22	1.96	0.47
2:B:186:ASN:HD22	2:B:195:GLN:HG2	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:122:ALA:HB2	2:D:174:LEU:HD11	1.97	0.47
2:D:186:ASN:HD22	2:D:195:GLN:HG2	1.81	0.46
2:B:188:THR:HG23	2:B:193:SER:HB3	1.98	0.46
1:C:498:ILE:HG23	1:C:503:MET:HB2	1.98	0.46
1:C:353:THR:HG22	1:C:356:ILE:HD12	1.98	0.46
2:B:43:LEU:HD22	2:B:46:SER:HB2	1.98	0.46
1:A:467:SER:O	1:A:471:THR:HG23	2.16	0.45
1:C:404:TYR:HD2	1:C:407:ILE:HG13	1.81	0.45
2:D:176:MET:HG2	2:D:202:VAL:HG13	1.99	0.45
1:A:173:PHE:CZ	1:A:221:LYS:HB2	2.52	0.45
1:A:438:ILE:O	1:A:442:ILE:HG13	2.17	0.45
2:D:188:THR:HG23	2:D:193:SER:HB3	1.97	0.45
1:C:173:PHE:CZ	1:C:221:LYS:HB2	2.51	0.45
1:A:188:LEU:HD21	1:A:460:LEU:HD23	1.99	0.45
1:A:168:GLU:OE1	1:A:489:HIS:NE2	2.33	0.45
2:D:43:LEU:HD22	2:D:46:SER:HB2	1.98	0.44
1:C:179:TYR:CD2	1:C:471:THR:HG22	2.45	0.44
2:D:142:ILE:HD11	2:D:184:ARG:HB2	1.98	0.44
1:A:200:TYR:HB3	2:B:131:VAL:CG2	2.48	0.44
1:C:229:ASN:HA	1:C:232:ILE:HD12	2.00	0.44
1:C:438:ILE:O	1:C:442:ILE:HG13	2.17	0.44
1:C:334:ASN:O	1:C:338:ASN:HB2	2.18	0.43
1:C:197:SER:OG	2:D:100:GLN:NE2	2.37	0.43
1:A:217:TYR:HA	1:A:328:ILE:HD13	1.99	0.43
2:B:157:ARG:HH22	2:B:179:ASP:CG	2.27	0.43
1:A:163:ILE:HG23	1:A:171:LEU:HD21	2.00	0.43
1:A:306:GLU:HG2	1:A:310:LYS:HE3	2.00	0.43
1:C:163:ILE:HG23	1:C:171:LEU:HD21	2.00	0.43
1:C:306:GLU:HG2	1:C:310:LYS:HE3	2.00	0.43
1:A:353:THR:HG22	1:A:356:ILE:HD12	2.01	0.42
2:B:176:MET:HG2	2:B:202:VAL:HG13	2.00	0.42
1:C:478:MET:O	1:C:482:THR:CG2	2.67	0.42
1:C:347:ASN:HD22	1:C:349:ASN:H	1.67	0.42
2:B:125:VAL:HG22	2:B:168:GLU:HG2	2.02	0.42
1:A:185:TYR:CZ	1:A:333:LYS:HG3	2.54	0.42
1:C:314:LEU:HD22	1:C:318:ILE:HD11	2.02	0.42
2:D:39:LEU:HB3	2:D:99:ILE:HD13	2.00	0.42
2:B:142:ILE:HD11	2:B:184:ARG:HB2	2.01	0.42
2:D:30:VAL:HG23	2:D:39:LEU:HD23	2.03	0.41
1:C:185:TYR:CZ	1:C:333:LYS:HG3	2.56	0.41
1:A:229:ASN:HA	1:A:232:ILE:HD12	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:168:GLU:OE1	1:C:489:HIS:NE2	2.34	0.41
1:C:193:ILE:HD13	1:C:206:VAL:HG21	2.03	0.41
1:A:383:MET:HG2	1:A:483:PHE:CE2	2.56	0.41
1:C:478:MET:O	1:C:482:THR:HG22	2.21	0.41
1:C:479:LEU:CA	1:C:482:THR:HG23	2.43	0.40
2:B:184:ARG:HG3	2:B:197:ILE:HG13	2.03	0.40
1:A:180:THR:CG2	1:A:471:THR:HG21	2.50	0.40
2:D:184:ARG:HG3	2:D:197:ILE:HG13	2.02	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	282/526 (54%)	274 (97%)	6 (2%)	2 (1%)	18	49
1	C	281/526 (53%)	276 (98%)	5 (2%)	0	100	100
2	B	179/269 (66%)	165 (92%)	14 (8%)	0	100	100
2	D	179/269 (66%)	167 (93%)	12 (7%)	0	100	100
All	All	921/1590 (58%)	882 (96%)	37 (4%)	2 (0%)	43	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	241	PRO
1	A	399	LYS

5.3.2 Protein sidechains [i](#)

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	277/511 (54%)	250 (90%)	27 (10%)	8	30
1	C	276/511 (54%)	241 (87%)	35 (13%)	4	18
2	B	156/225 (69%)	138 (88%)	18 (12%)	5	23
2	D	156/225 (69%)	141 (90%)	15 (10%)	8	30
All	All	865/1472 (59%)	770 (89%)	95 (11%)	6	24

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	ASP
1	A	163	ILE
1	A	164	LEU
1	A	165	GLN
1	A	168	GLU
1	A	193	ILE
1	A	213	ILE
1	A	223	LYS
1	A	230	ASP
1	A	305	ASP
1	A	314	LEU
1	A	345	CYS
1	A	347	ASN
1	A	368	ILE
1	A	418	LYS
1	A	419	HIS
1	A	442	ILE
1	A	446	ILE
1	A	451	GLN
1	A	453	ASP
1	A	455	LEU
1	A	458	ARG
1	A	465	GLU
1	A	477	GLN
1	A	487	GLU
1	A	490	LEU
1	A	497	LEU
2	B	29	THR

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Mol	Chain	Res	Type
2	B	37	ILE
2	B	40	THR
2	B	43	LEU
2	B	48	THR
2	B	49	GLU
2	B	50	VAL
2	B	63	LYS
2	B	64	GLU
2	B	70	GLN
2	B	94	MET
2	B	106	ARG
2	B	141	LYS
2	B	144	ASP
2	B	169	LEU
2	B	174	LEU
2	B	175	ASN
2	B	176	MET
1	C	162	ASP
1	C	163	ILE
1	C	164	LEU
1	C	168	GLU
1	C	193	ILE
1	C	213	ILE
1	C	223	LYS
1	C	228	LYS
1	C	230	ASP
1	C	242	TYR
1	C	314	LEU
1	C	322	GLU
1	C	338	ASN
1	C	345	CYS
1	C	347	ASN
1	C	368	ILE
1	C	392	LEU
1	C	393	LEU
1	C	394	LEU
1	C	399	LYS
1	C	400	LYS
1	C	418	LYS
1	C	435	THR
1	C	442	ILE
1	C	446	ILE

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Mol	Chain	Res	Type
1	C	451	GLN
1	C	453	ASP
1	C	455	LEU
1	C	458	ARG
1	C	465	GLU
1	C	477	GLN
1	C	482	THR
1	C	486	LYS
1	C	487	GLU
1	C	490	LEU
2	D	29	THR
2	D	31	GLU
2	D	37	ILE
2	D	40	THR
2	D	43	LEU
2	D	48	THR
2	D	50	VAL
2	D	64	GLU
2	D	70	GLN
2	D	94	MET
2	D	142	ILE
2	D	169	LEU
2	D	173	ASN
2	D	175	ASN
2	D	176	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	191	ASN
1	A	225	ASN
1	A	321	HIS
1	A	347	ASN
1	A	375	ASN
1	A	439	GLN
1	A	496	HIS
2	B	115	HIS
2	B	152	ASN
2	B	186	ASN
1	C	191	ASN
1	C	321	HIS
1	C	347	ASN

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Mol	Chain	Res	Type
1	C	349	ASN
1	C	375	ASN
1	C	439	GLN
1	C	451	GLN
1	C	495	HIS
1	C	496	HIS
2	D	100	GLN
2	D	186	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 ⓘ

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	286/526 (54%)	0.16	8 (2%) 55 34	26, 43, 77, 116	0
1	C	285/526 (54%)	-0.03	2 (0%) 84 68	25, 36, 55, 96	0
2	B	181/269 (67%)	0.55	5 (2%) 55 34	23, 57, 97, 111	0
2	D	181/269 (67%)	0.69	8 (4%) 39 21	39, 66, 108, 128	0
All	All	933/1590 (58%)	0.28	23 (2%) 58 37	23, 45, 93, 128	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	66	ALA	5.8
2	B	66	ALA	5.6
2	B	65	ASP	5.1
2	D	65	ASP	4.0
1	A	242	TYR	3.3
2	D	144	ASP	2.9
2	B	134	VAL	2.7
1	C	346	TYR	2.4
2	B	183	TYR	2.3
2	D	47	ALA	2.3
1	A	240	HIS	2.2
1	A	302	LYS	2.2
2	D	163	SER	2.2
1	A	305	ASP	2.2
2	D	112	SER	2.2
1	A	419	HIS	2.2
2	D	117	ASN	2.1
2	B	58	GLY	2.1
1	A	303	MET	2.0
1	A	160	SER	2.0
1	A	395	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	302	LYS	2.0
2	D	114	GLU	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.