



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 05:54 AM GMT

PDB ID : 3C4Y
Title : Crystal Structure of Apo form of G protein coupled receptor kinase 1 at 7.51Å
Authors : Singh, P.; Tesmer, J.J.G.
Deposited on : 2008-01-30
Resolution : 7.51 Å(reported)

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

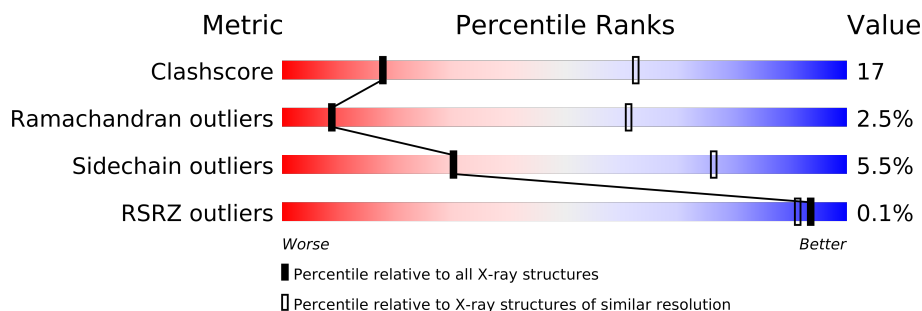
The following versions of software and data (see [references](#)) were used in the production of this report:

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 7.51 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))
Clashscore	79885	1039 (10.00-3.52)
Ramachandran outliers	78287	1301 (11.50-3.50)
Sidechain outliers	78261	1275 (11.50-3.50)
RSRZ outliers	66119	1105 (11.50-3.50)

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 ≥ 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	543	
1	B	543	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7629 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 Rhodopsin kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	478	Total	C	N	O	S	0	0	0
			3854	2474	666	695	19			
1	B	466	Total	C	N	O	S	0	0	0
			3775	2427	649	680	19			

There are 16 discrepancies between the modelled and reference sequences:

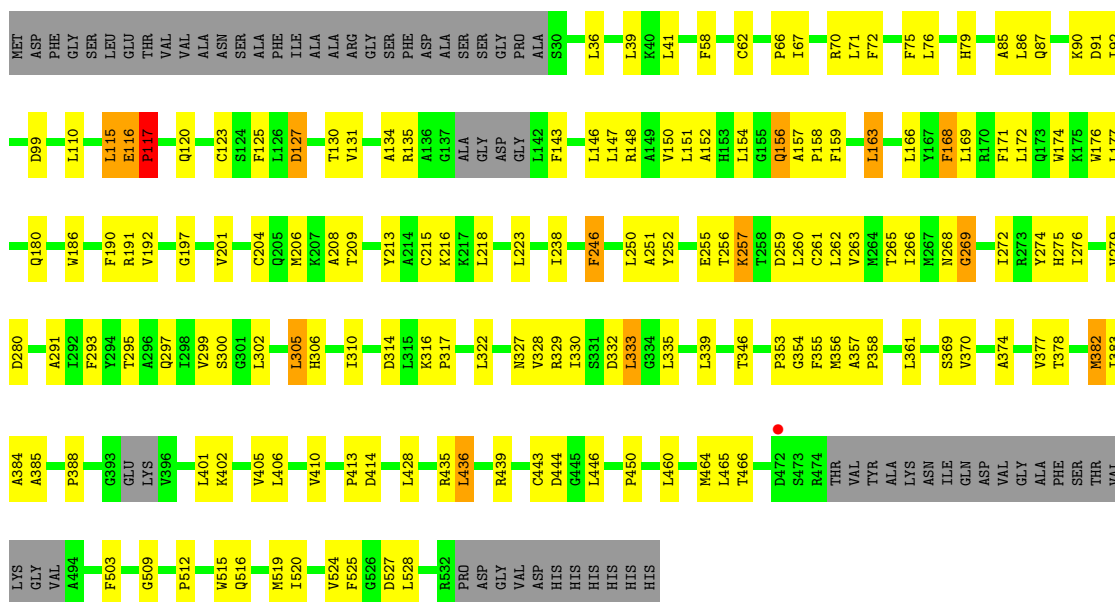
Chain	Residue	Modelled	Actual	Comment	Reference
A	536	VAL	-	EXPRESSION TAG	UNP P28327
A	537	ASP	-	EXPRESSION TAG	UNP P28327
A	538	HIS	-	EXPRESSION TAG	UNP P28327
A	539	HIS	-	EXPRESSION TAG	UNP P28327
A	540	HIS	-	EXPRESSION TAG	UNP P28327
A	541	HIS	-	EXPRESSION TAG	UNP P28327
A	542	HIS	-	EXPRESSION TAG	UNP P28327
A	543	HIS	-	EXPRESSION TAG	UNP P28327
B	536	VAL	-	EXPRESSION TAG	UNP P28327
B	537	ASP	-	EXPRESSION TAG	UNP P28327
B	538	HIS	-	EXPRESSION TAG	UNP P28327
B	539	HIS	-	EXPRESSION TAG	UNP P28327
B	540	HIS	-	EXPRESSION TAG	UNP P28327
B	541	HIS	-	EXPRESSION TAG	UNP P28327
B	542	HIS	-	EXPRESSION TAG	UNP P28327
B	543	HIS	-	EXPRESSION TAG	UNP P28327

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.

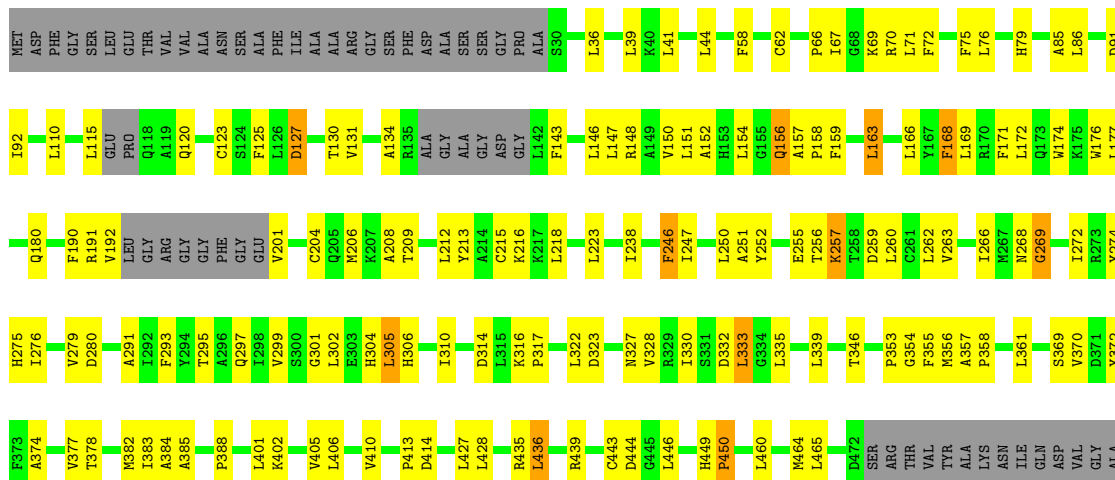
• Molecule 1: Rhodopsin kinase

Chain A:



• Molecule 1: Rhodopsin kinase

Chain B:



PHE	SER	THR	VAL	LYS	GLY	VAL	A494	F503	G509	F512	W515	Q516	M519	T520	V524	F525	Q526	Q527	L528	R532	PRO	ASP	GLY	VAL	ASP	HIS	HIS	HIS	HIS	HIS
-----	-----	-----	-----	-----	-----	-----	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	123.31Å 123.31Å 192.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 7.51 19.94 – 7.51	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-7.51) 100.0 (19.94-7.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.03 (at 7.78Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.182 , (Not available) 0.200 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	540.3	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.14 , 162.3	EDS
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 2230 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7629	wwPDB-VP
Average B, all atoms (Å ²)	224.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.85% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.39	0/3944	0.53	0/5316
1	B	0.38	0/3862	0.53	0/5205
All	All	0.39	0/7806	0.53	0/10521

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.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	LEU	Peptide
1	A	116	GLU	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	3854	0	3820	133	0
1	B	3775	0	3748	130	0
All	All	7629	0	7568	262	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 17.

All (262) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:209:THR:HG22	1:B:524:VAL:HG21	1.43	1.00
1:A:209:THR:HG22	1:A:524:VAL:HG21	1.46	0.95
1:B:275:HIS:HA	1:B:279:VAL:HG23	1.51	0.90
1:B:134:ALA:HB2	1:B:146:LEU:HD13	1.52	0.90
1:A:275:HIS:HA	1:A:279:VAL:HG23	1.53	0.88
1:A:134:ALA:HB2	1:A:146:LEU:HD13	1.54	0.87
1:A:250:LEU:HD12	1:A:263:VAL:O	1.76	0.85
1:B:250:LEU:HD12	1:B:263:VAL:O	1.77	0.84
1:B:314:ASP:HB2	1:B:335:LEU:HD13	1.61	0.83
1:B:361:LEU:HD12	1:B:405:VAL:HG11	1.61	0.82
1:A:314:ASP:HB2	1:A:335:LEU:HD13	1.61	0.81
1:B:356:MET:HE1	1:B:361:LEU:HD21	1.60	0.81
1:A:361:LEU:HD12	1:A:405:VAL:HG11	1.64	0.80
1:A:410:VAL:HG21	1:A:428:LEU:HD13	1.63	0.78
1:B:410:VAL:HG21	1:B:428:LEU:HD13	1.68	0.76
1:A:116:GLU:HB3	1:A:117:PRO:HD2	1.70	0.73
1:A:250:LEU:HD21	1:A:262:LEU:HD22	1.74	0.69
1:A:250:LEU:HD11	1:A:262:LEU:HD22	1.75	0.68
1:B:358:PRO:HB3	1:B:406:LEU:HD23	1.75	0.68
1:B:250:LEU:HD21	1:B:262:LEU:HD22	1.75	0.68
1:A:410:VAL:HG21	1:A:428:LEU:CD1	2.24	0.68
1:B:250:LEU:HD11	1:B:262:LEU:HD22	1.76	0.67
1:A:356:MET:HE1	1:A:361:LEU:HD21	1.76	0.67
1:B:238:ILE:HG21	1:B:310:ILE:CD1	2.24	0.67
1:A:353:PRO:HA	1:A:356:MET:HE2	1.77	0.67
1:A:358:PRO:HB3	1:A:406:LEU:HD23	1.75	0.67
1:A:150:VAL:HG22	1:A:154:LEU:CD1	2.25	0.67
1:B:356:MET:CE	1:B:361:LEU:HD21	2.25	0.66
1:A:168:PHE:CZ	1:A:172:LEU:HD11	2.31	0.65
1:A:192:VAL:HG13	1:A:201:VAL:O	1.97	0.65
1:A:356:MET:CE	1:A:361:LEU:HD21	2.27	0.64
1:B:353:PRO:HA	1:B:356:MET:HE2	1.80	0.64
1:B:401:LEU:O	1:B:405:VAL:HG23	1.98	0.64
1:B:410:VAL:HG21	1:B:428:LEU:CD1	2.28	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:305:LEU:HD13	1:A:306:HIS:NE2	2.13	0.63
1:B:168:PHE:CZ	1:B:172:LEU:HD11	2.33	0.63
1:A:377:VAL:HG13	1:A:388:PRO:CD	2.29	0.63
1:A:401:LEU:O	1:A:405:VAL:HG23	1.99	0.62
1:B:305:LEU:HD13	1:B:306:HIS:NE2	2.14	0.62
1:B:150:VAL:HG22	1:B:154:LEU:CD1	2.28	0.62
1:B:212:LEU:HD12	1:B:266:ILE:HD12	1.83	0.61
1:B:192:VAL:HG13	1:B:201:VAL:O	2.00	0.61
1:B:213:TYR:CG	1:B:263:VAL:HG13	2.35	0.61
1:B:302:LEU:HD23	1:B:305:LEU:HD12	1.82	0.61
1:A:302:LEU:HD23	1:A:305:LEU:HD12	1.83	0.60
1:A:238:ILE:HG21	1:A:310:ILE:CD1	2.31	0.60
1:B:251:ALA:HB3	1:B:263:VAL:HG12	1.83	0.60
1:B:305:LEU:HD13	1:B:306:HIS:CD2	2.36	0.60
1:B:305:LEU:HD22	1:B:310:ILE:O	2.02	0.60
1:A:357:ALA:CB	1:A:370:VAL:HG12	2.32	0.60
1:B:377:VAL:HG13	1:B:388:PRO:CD	2.32	0.59
1:A:251:ALA:HB3	1:A:263:VAL:HG12	1.83	0.59
1:B:166:LEU:HA	1:B:169:LEU:HD12	1.82	0.59
1:A:213:TYR:CG	1:A:263:VAL:HG13	2.38	0.59
1:A:76:LEU:HB2	1:A:86:LEU:HD22	1.83	0.58
1:A:305:LEU:HD13	1:A:306:HIS:CD2	2.37	0.58
1:B:275:HIS:HA	1:B:279:VAL:CG2	2.30	0.58
1:A:166:LEU:HA	1:A:169:LEU:HD12	1.85	0.58
1:B:92:ILE:HG21	1:B:147:LEU:HD13	1.86	0.58
1:B:357:ALA:HB2	1:B:370:VAL:HA	1.85	0.58
1:A:305:LEU:HD22	1:A:310:ILE:O	2.03	0.58
1:A:116:GLU:HB3	1:A:117:PRO:CD	2.34	0.58
1:B:279:VAL:HG12	1:B:280:ASP:N	2.19	0.57
1:A:295:THR:O	1:A:299:VAL:HG23	2.05	0.57
1:A:357:ALA:HB2	1:A:370:VAL:HA	1.86	0.57
1:B:76:LEU:HB2	1:B:86:LEU:HD22	1.87	0.57
1:B:115:LEU:CD2	1:B:131:VAL:HG13	2.35	0.57
1:A:130:THR:HG22	1:A:146:LEU:HD12	1.87	0.56
1:B:134:ALA:CB	1:B:146:LEU:HD13	2.30	0.56
1:B:295:THR:O	1:B:299:VAL:HG23	2.05	0.56
1:B:130:THR:HG22	1:B:146:LEU:HD12	1.86	0.56
1:A:250:LEU:HD11	1:A:252:TYR:O	2.05	0.56
1:B:212:LEU:HD12	1:B:266:ILE:CD1	2.36	0.56
1:A:157:ALA:HB3	1:A:158:PRO:HD3	1.88	0.56
1:A:436:LEU:HD21	1:A:446:LEU:HA	1.88	0.56
1:A:361:LEU:HD22	1:A:402:LYS:HD2	1.88	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:357:ALA:HB3	1:A:370:VAL:HG12	1.88	0.55
1:B:250:LEU:HD11	1:B:252:TYR:O	2.07	0.55
1:A:115:LEU:CD2	1:A:131:VAL:HG13	2.37	0.55
1:B:317:PRO:HD3	1:B:378:THR:HG23	1.88	0.54
1:B:322:LEU:HG	1:B:328:VAL:HG12	1.88	0.54
1:B:157:ALA:HB3	1:B:158:PRO:HD3	1.89	0.54
1:A:134:ALA:HB1	1:A:143:PHE:CD2	2.42	0.54
1:A:238:ILE:HD12	1:A:310:ILE:HD11	1.89	0.54
1:B:134:ALA:HB1	1:B:143:PHE:CD2	2.43	0.53
1:B:357:ALA:CB	1:B:370:VAL:HG12	2.38	0.53
1:B:361:LEU:HD22	1:B:402:LYS:HD2	1.90	0.53
1:B:215:CYS:SG	1:B:263:VAL:HG23	2.48	0.52
1:A:92:ILE:HG21	1:A:147:LEU:HD13	1.91	0.52
1:A:147:LEU:O	1:A:150:VAL:HG12	2.09	0.52
1:B:71:LEU:HD12	1:B:171:PHE:HD1	1.75	0.52
1:A:275:HIS:HA	1:A:279:VAL:CG2	2.33	0.52
1:A:377:VAL:HG13	1:A:388:PRO:CG	2.40	0.52
1:A:322:LEU:HG	1:A:328:VAL:HG12	1.91	0.52
1:A:174:TRP:CH2	1:A:525:PHE:HA	2.45	0.52
1:B:174:TRP:CH2	1:B:525:PHE:HA	2.45	0.51
1:A:297:GLN:OE1	1:A:328:VAL:HG22	2.11	0.51
1:B:519:MET:SD	1:B:519:MET:N	2.83	0.51
1:B:251:ALA:HB3	1:B:263:VAL:CG1	2.41	0.51
1:A:150:VAL:HG22	1:A:154:LEU:HD11	1.90	0.51
1:B:67:ILE:HD12	1:B:516:GLN:HG2	1.93	0.51
1:A:251:ALA:HB3	1:A:263:VAL:CG1	2.40	0.51
1:A:238:ILE:HG21	1:A:310:ILE:HD13	1.92	0.50
1:B:247:ILE:HD11	1:B:304:HIS:HB3	1.92	0.50
1:B:91:ASP:HB3	1:B:110:LEU:HD21	1.94	0.50
1:B:436:LEU:HD21	1:B:446:LEU:HA	1.92	0.50
1:A:279:VAL:HG12	1:A:280:ASP:N	2.25	0.50
1:A:250:LEU:CD2	1:A:262:LEU:HD22	2.41	0.50
1:A:71:LEU:HD12	1:A:171:PHE:HD1	1.76	0.50
1:B:150:VAL:HG22	1:B:154:LEU:HD11	1.94	0.50
1:A:297:GLN:CG	1:A:460:LEU:HD11	2.42	0.50
1:B:377:VAL:HG13	1:B:388:PRO:CG	2.42	0.50
1:A:215:CYS:SG	1:A:263:VAL:HG23	2.52	0.49
1:A:317:PRO:HD3	1:A:378:THR:HG23	1.94	0.49
1:A:333:LEU:N	1:A:333:LEU:HD22	2.28	0.49
1:A:134:ALA:CB	1:A:146:LEU:HD13	2.32	0.49
1:B:357:ALA:HB3	1:B:370:VAL:HG12	1.93	0.49
1:A:443:CYS:HB3	1:A:446:LEU:HB3	1.94	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:246:PHE:HB2	1:B:301:GLY:HA2	1.94	0.49
1:A:246:PHE:CE2	1:A:300:SER:HB2	2.48	0.49
1:B:355:PHE:O	1:B:374:ALA:HB2	2.13	0.49
1:B:317:PRO:CD	1:B:378:THR:HG23	2.42	0.49
1:B:213:TYR:CD2	1:B:263:VAL:HG13	2.47	0.49
1:B:67:ILE:HG13	1:B:520:ILE:HD11	1.95	0.49
1:B:148:ARG:HA	1:B:151:LEU:HD12	1.95	0.49
1:B:361:LEU:HD12	1:B:405:VAL:CG1	2.38	0.48
1:A:218:LEU:O	1:A:259:ASP:HA	2.13	0.48
1:A:250:LEU:HD13	1:A:262:LEU:HB3	1.94	0.48
1:B:218:LEU:O	1:B:259:ASP:HA	2.14	0.48
1:A:91:ASP:HB3	1:A:110:LEU:HD21	1.94	0.48
1:A:355:PHE:O	1:A:374:ALA:HB2	2.13	0.48
1:A:36:LEU:HD12	1:A:39:LEU:HD12	1.96	0.48
1:B:250:LEU:HD13	1:B:262:LEU:HB3	1.96	0.48
1:A:176:TRP:CZ3	1:A:177:LEU:HD13	2.49	0.48
1:B:85:ALA:HB1	1:B:154:LEU:HD21	1.96	0.48
1:B:166:LEU:O	1:B:169:LEU:N	2.47	0.47
1:A:168:PHE:CE1	1:A:172:LEU:HD21	2.50	0.47
1:B:168:PHE:CE1	1:B:172:LEU:HD21	2.49	0.47
1:A:67:ILE:HD12	1:A:516:GLN:HG2	1.96	0.47
1:A:150:VAL:HG22	1:A:154:LEU:HG	1.97	0.47
1:B:260:LEU:HD11	1:B:503:PHE:CD2	2.49	0.47
1:B:272:ILE:HG22	1:B:276:ILE:HG13	1.95	0.47
1:B:512:PRO:O	1:B:515:TRP:HB3	2.15	0.47
1:B:333:LEU:N	1:B:333:LEU:HD22	2.30	0.47
1:B:159:PHE:CE2	1:B:163:LEU:HD21	2.49	0.47
1:B:159:PHE:CZ	1:B:163:LEU:HD21	2.49	0.47
1:B:71:LEU:HD12	1:B:171:PHE:CD1	2.50	0.47
1:A:213:TYR:CD2	1:A:263:VAL:HG13	2.50	0.47
1:A:67:ILE:HG13	1:A:520:ILE:HD11	1.97	0.47
1:B:297:GLN:CG	1:B:460:LEU:HD11	2.45	0.47
1:A:159:PHE:CZ	1:A:163:LEU:HD21	2.50	0.47
1:A:297:GLN:HG3	1:A:460:LEU:HD11	1.97	0.46
1:A:519:MET:N	1:A:519:MET:SD	2.88	0.46
1:A:272:ILE:HG22	1:A:276:ILE:HG13	1.96	0.46
1:A:314:ASP:CB	1:A:335:LEU:HD13	2.40	0.46
1:A:250:LEU:CD1	1:A:262:LEU:HD22	2.43	0.46
1:B:297:GLN:OE1	1:B:328:VAL:HG22	2.15	0.46
1:A:91:ASP:CB	1:A:110:LEU:HD21	2.45	0.46
1:B:176:TRP:CZ3	1:B:177:LEU:HD13	2.51	0.46
1:A:354:GLY:O	1:A:377:VAL:HG21	2.16	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:443:CYS:HB3	1:B:446:LEU:HB3	1.97	0.46
1:A:159:PHE:CE2	1:A:163:LEU:HD21	2.50	0.46
1:B:62:CYS:SG	1:B:72:PHE:HB2	2.56	0.46
1:B:152:ALA:O	1:B:156:GLN:HG3	2.15	0.46
1:A:260:LEU:HD11	1:A:503:PHE:CD2	2.50	0.46
1:B:147:LEU:O	1:B:150:VAL:HG12	2.16	0.46
1:B:143:PHE:CD2	1:B:146:LEU:HD22	2.51	0.46
1:B:91:ASP:CB	1:B:110:LEU:HD21	2.45	0.46
1:B:204:CYS:SG	1:B:215:CYS:HB2	2.56	0.46
1:A:152:ALA:O	1:A:156:GLN:HG3	2.16	0.46
1:B:150:VAL:HG22	1:B:154:LEU:HG	1.97	0.45
1:B:115:LEU:HD23	1:B:131:VAL:HG13	1.98	0.45
1:A:377:VAL:HG13	1:A:388:PRO:HG3	1.98	0.45
1:A:317:PRO:CD	1:A:378:THR:HG23	2.46	0.45
1:A:238:ILE:CG2	1:A:310:ILE:CD1	2.95	0.45
1:B:250:LEU:CD2	1:B:262:LEU:HD22	2.43	0.45
1:B:314:ASP:CB	1:B:335:LEU:HD13	2.40	0.45
1:A:71:LEU:HD12	1:A:171:PHE:CD1	2.52	0.45
1:A:166:LEU:O	1:A:169:LEU:N	2.48	0.45
1:B:293:PHE:CE1	1:B:465:LEU:HD22	2.51	0.45
1:B:250:LEU:CD1	1:B:262:LEU:HD22	2.44	0.45
1:A:339:LEU:HD21	1:A:346:THR:HG22	1.99	0.45
1:A:213:TYR:HB3	1:A:263:VAL:HG13	1.98	0.45
1:B:354:GLY:O	1:B:377:VAL:HG21	2.17	0.45
1:B:297:GLN:HG3	1:B:460:LEU:HD11	1.99	0.45
1:A:148:ARG:HA	1:A:151:LEU:HD12	1.99	0.45
1:A:265:THR:HG21	1:A:329:ARG:NH2	2.32	0.45
1:A:384:ALA:O	1:A:385:ALA:HB3	2.17	0.45
1:A:361:LEU:HD13	1:A:402:LYS:HA	2.00	0.44
1:A:62:CYS:SG	1:A:72:PHE:HB2	2.56	0.44
1:A:177:LEU:HD12	1:A:180:GLN:HG3	1.98	0.44
1:B:291:ALA:CB	1:B:383:ILE:HD11	2.47	0.44
1:A:306:HIS:HD2	1:A:310:ILE:O	2.00	0.44
1:B:377:VAL:HG13	1:B:388:PRO:HG3	1.99	0.44
1:A:204:CYS:SG	1:A:215:CYS:HB2	2.57	0.44
1:B:302:LEU:HD22	1:B:306:HIS:CE1	2.51	0.44
1:A:512:PRO:O	1:A:515:TRP:HB3	2.18	0.44
1:B:41:LEU:HD11	1:B:528:LEU:O	2.18	0.44
1:A:36:LEU:CD1	1:A:208:ALA:HB1	2.48	0.44
1:B:256:THR:OG1	1:B:257:LYS:N	2.50	0.44
1:A:291:ALA:CB	1:A:383:ILE:HD11	2.47	0.44
1:B:361:LEU:HD13	1:B:402:LYS:HA	1.98	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:384:ALA:O	1:B:385:ALA:HB3	2.18	0.44
1:B:168:PHE:O	1:B:171:PHE:HB3	2.17	0.44
1:B:213:TYR:HB3	1:B:263:VAL:HG13	2.00	0.43
1:B:302:LEU:CD2	1:B:305:LEU:HD12	2.48	0.43
1:A:39:LEU:HD21	1:A:186:TRP:CZ3	2.54	0.43
1:A:75:PHE:CE1	1:A:79:HIS:CE1	3.06	0.43
1:A:168:PHE:O	1:A:171:PHE:HB3	2.18	0.43
1:A:143:PHE:CD2	1:A:146:LEU:HD22	2.54	0.43
1:A:85:ALA:HB1	1:A:154:LEU:HD21	2.00	0.43
1:B:127:ASP:N	1:B:127:ASP:OD1	2.49	0.43
1:B:36:LEU:HD12	1:B:39:LEU:HD12	2.00	0.43
1:B:369:SER:HB2	1:B:435:ARG:HD2	2.00	0.43
1:A:85:ALA:HB2	1:A:125:PHE:CZ	2.54	0.43
1:B:201:VAL:HG22	1:B:216:LYS:HG3	2.01	0.43
1:A:150:VAL:HG22	1:A:154:LEU:CG	2.49	0.43
1:B:339:LEU:HD21	1:B:346:THR:HG22	2.01	0.43
1:B:268:ASN:O	1:B:269:GLY:C	2.57	0.43
1:B:134:ALA:HB1	1:B:143:PHE:HD2	1.84	0.43
1:A:201:VAL:HG22	1:A:216:LYS:HG3	2.01	0.43
1:A:36:LEU:HD13	1:A:208:ALA:C	2.39	0.43
1:B:85:ALA:HB2	1:B:125:PHE:CZ	2.54	0.43
1:A:41:LEU:HD11	1:A:528:LEU:O	2.18	0.43
1:A:213:TYR:CG	1:A:263:VAL:CG1	3.02	0.42
1:B:361:LEU:HD13	1:B:402:LYS:HG3	2.01	0.42
1:A:361:LEU:HD13	1:A:402:LYS:HG3	2.01	0.42
1:A:302:LEU:HD22	1:A:306:HIS:CE1	2.55	0.42
1:B:177:LEU:HD12	1:B:180:GLN:HG3	2.01	0.42
1:B:213:TYR:CG	1:B:263:VAL:CG1	3.01	0.42
1:B:66:PRO:HB2	1:B:520:ILE:CD1	2.49	0.42
1:A:256:THR:OG1	1:A:257:LYS:N	2.50	0.42
1:B:36:LEU:CD1	1:B:208:ALA:HB1	2.50	0.42
1:A:134:ALA:HB1	1:A:143:PHE:HD2	1.83	0.42
1:A:115:LEU:HD23	1:A:131:VAL:HG13	2.02	0.42
1:B:323:ASP:OD1	1:B:327:ASN:N	2.52	0.42
1:A:131:VAL:HG12	1:A:135:ARG:NH2	2.35	0.42
1:A:369:SER:HB2	1:A:435:ARG:HD2	2.02	0.42
1:A:465:LEU:HD23	1:A:466:THR:N	2.35	0.41
1:A:291:ALA:HB1	1:A:382:MET:HE3	2.02	0.41
1:A:66:PRO:HB2	1:A:520:ILE:CD1	2.51	0.41
1:B:75:PHE:CE1	1:B:79:HIS:CE1	3.08	0.41
1:A:268:ASN:O	1:A:269:GLY:C	2.58	0.41
1:B:150:VAL:HG22	1:B:154:LEU:CG	2.50	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:36:LEU:HD13	1:B:208:ALA:C	2.40	0.41
1:B:238:ILE:HG21	1:B:310:ILE:HD13	2.01	0.41
1:A:316:LYS:HB2	1:A:317:PRO:CD	2.51	0.41
1:A:291:ALA:HB3	1:A:383:ILE:HD11	2.03	0.41
1:B:66:PRO:HB2	1:B:520:ILE:HD13	2.02	0.41
1:A:127:ASP:N	1:A:127:ASP:OD1	2.53	0.41
1:B:358:PRO:O	1:B:406:LEU:HD21	2.21	0.41
1:A:215:CYS:SG	1:A:261:CYS:HB3	2.61	0.41
1:A:169:LEU:HD11	1:B:44:LEU:CD1	2.51	0.41
1:A:297:GLN:HE22	1:A:327:ASN:HB3	1.85	0.41
1:B:67:ILE:CG1	1:B:520:ILE:HD11	2.51	0.41
1:B:316:LYS:HB2	1:B:317:PRO:CD	2.51	0.40
1:B:449:HIS:CG	1:B:450:PRO:HD2	2.57	0.40
1:B:66:PRO:HA	1:B:69:LYS:HB3	2.04	0.40
1:A:293:PHE:CE1	1:A:465:LEU:HD22	2.56	0.40
1:B:372:TYR:O	1:B:427:LEU:HD22	2.21	0.40
1:A:250:LEU:HD22	1:A:262:LEU:HD13	2.03	0.40
1:A:87:GLN:HA	1:A:90:LYS: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	470/543 (87%)	401 (85%)	56 (12%)	13 (3%)	8	59
1	B	456/543 (84%)	394 (86%)	52 (11%)	10 (2%)	10	64
All	All	926/1086 (85%)	795 (86%)	108 (12%)	23 (2%)	9	61

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	PRO
1	A	266	ILE

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Mol	Chain	Res	Type
1	A	127	ASP
1	A	257	LYS
1	A	269	GLY
1	B	127	ASP
1	B	257	LYS
1	B	269	GLY
1	A	190	PHE
1	A	332	ASP
1	B	190	PHE
1	B	332	ASP
1	A	436	LEU
1	B	436	LEU
1	A	58	PHE
1	A	413	PRO
1	B	58	PHE
1	B	413	PRO
1	A	450	PRO
1	A	509	GLY
1	B	450	PRO
1	B	509	GLY
1	A	197	GLY

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	401/449 (89%)	378 (94%)	23 (6%)	29	75
1	B	395/449 (88%)	374 (95%)	21 (5%)	32	77
All	All	796/898 (89%)	752 (94%)	44 (6%)	30	76

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ARG
1	A	99	ASP
1	A	117	PRO
1	A	120	GLN

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Mol	Chain	Res	Type
1	A	123	CYS
1	A	156	GLN
1	A	163	LEU
1	A	168	PHE
1	A	191	ARG
1	A	206	MET
1	A	223	LEU
1	A	246	PHE
1	A	255	GLU
1	A	274	TYR
1	A	305	LEU
1	A	330	ILE
1	A	333	LEU
1	A	382	MET
1	A	414	ASP
1	A	439	ARG
1	A	444	ASP
1	A	464	MET
1	A	527	ASP
1	B	70	ARG
1	B	120	GLN
1	B	123	CYS
1	B	156	GLN
1	B	163	LEU
1	B	168	PHE
1	B	191	ARG
1	B	206	MET
1	B	223	LEU
1	B	246	PHE
1	B	255	GLU
1	B	274	TYR
1	B	305	LEU
1	B	330	ILE
1	B	333	LEU
1	B	382	MET
1	B	414	ASP
1	B	439	ARG
1	B	444	ASP
1	B	464	MET
1	B	527	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	173	GLN
1	A	268	ASN
1	A	278	ASN
1	A	306	HIS
1	B	120	GLN
1	B	173	GLN
1	B	268	ASN
1	B	278	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 ⓘ

There are no ligands in this entry.

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, 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	478/543 (88%)	-0.32	1 (0%) 93 89	223, 224, 225, 238	1 (0%)
1	B	466/543 (85%)	-0.31	0 100 100	223, 224, 225, 238	0
All	All	944/1086 (86%)	-0.32	1 (0%) 93 91	223, 224, 225, 238	1 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	472	ASP	2.4

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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.