



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:10 AM GMT

PDB ID : 3GDI
Title : Mammalian Clock Protein mPER2 - Crystal Structure of a PAS Domain Fragment
Authors : Hennig, S.; Wolf, E.
Deposited on : 2009-02-24
Resolution : 2.40 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>

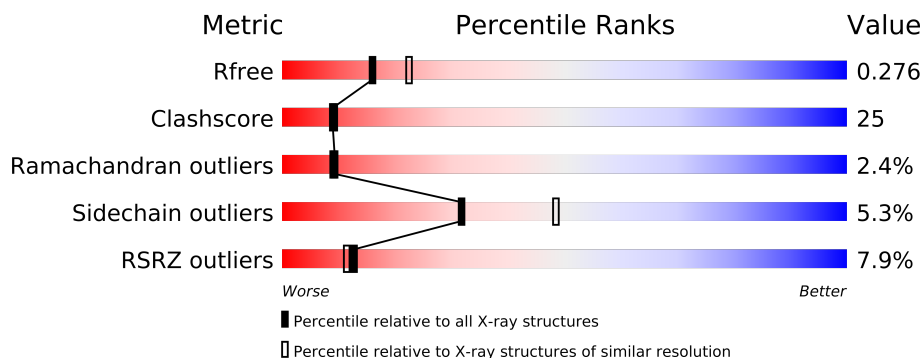
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.40 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	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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	309	
1	B	309	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4300 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 Period circadian protein homolog 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2148	1389	364	385	10			
1	B	250	Total	C	N	O	S	0	0	0
			1954	1264	335	346	9			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	165	GLY	-	EXPRESSION TAG	UNP O54943
A	166	PRO	-	EXPRESSION TAG	UNP O54943
A	167	LEU	-	EXPRESSION TAG	UNP O54943
A	168	GLY	-	EXPRESSION TAG	UNP O54943
A	169	SER	-	EXPRESSION TAG	UNP O54943
B	165	GLY	-	EXPRESSION TAG	UNP O54943
B	166	PRO	-	EXPRESSION TAG	UNP O54943
B	167	LEU	-	EXPRESSION TAG	UNP O54943
B	168	GLY	-	EXPRESSION TAG	UNP O54943
B	169	SER	-	EXPRESSION TAG	UNP O54943

- Molecule 2 is water.

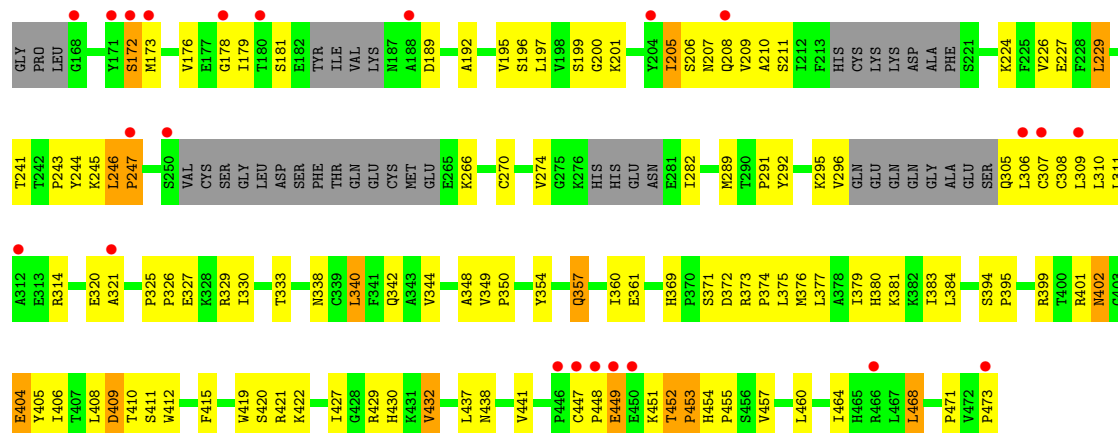
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	89	Total	O	0	0
			89	89		
2	B	109	Total	O	0	0
			109	109		

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.

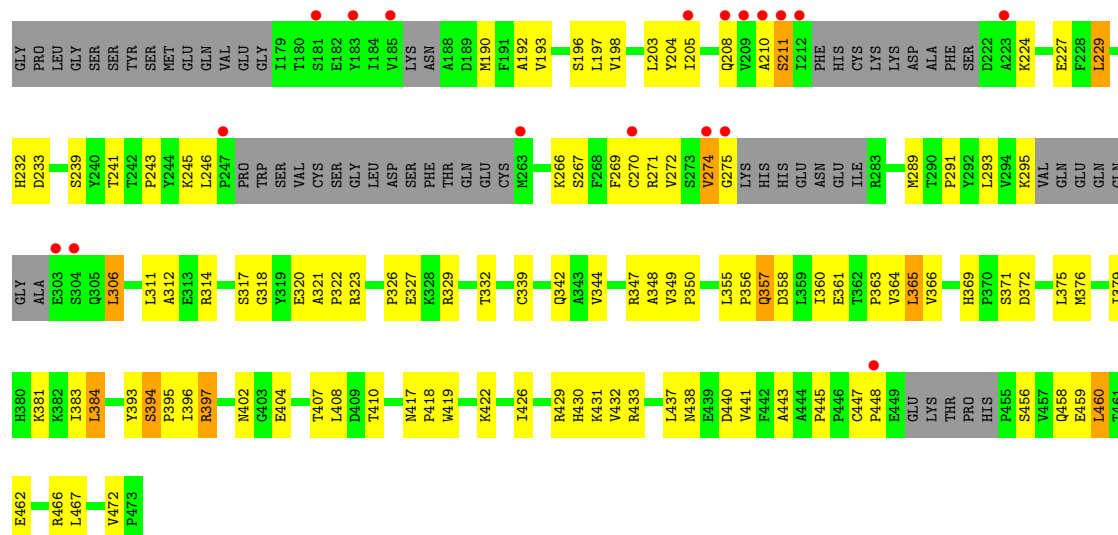
• Molecule 1: Period circadian protein homolog 2

Chain A: 



• Molecule 1: Period circadian protein homolog 2

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.67Å 63.38Å 72.51Å 90.00° 102.25° 90.00°	Depositor
Resolution (Å)	19.92 – 2.40 19.92 – 2.40	Depositor EDS
% Data completeness (in resolution range)	57.2 (19.92-2.40) 98.6 (19.92-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.41Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.272 0.232 , 0.276	Depositor DCC
R_{free} test set	1199 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.7	EDS
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 23980 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4300	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.61% 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	0.37	0/2208	0.62	1/3002 (0.0%)
1	B	0.36	0/2004	0.65	1/2725 (0.0%)
All	All	0.36	0/4212	0.64	2/5727 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	445	PRO	N-CA-CB	5.47	109.86	103.30
1	A	409	ASP	N-CA-C	-5.06	97.33	111.00

There are no chirality outliers.

There are no planarity outliers.

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	2148	0	2135	106	0
1	B	1954	0	1923	103	0
2	A	89	0	0	4	0
2	B	109	0	0	11	0
All	All	4300	0	4058	206	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 25.

All (206) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:394:SER:HB2	1:B:395:PRO:HD2	1.42	1.00
1:A:197:LEU:HD21	1:A:243:PRO:HG3	1.46	0.96
1:B:321:ALA:HB3	1:B:322:PRO:HD3	1.49	0.93
1:A:369:HIS:HD2	1:A:371:SER:H	1.10	0.90
1:A:241:THR:HA	1:A:246:LEU:HD21	1.59	0.83
1:B:323:ARG:HD3	2:B:62:HOH:O	1.79	0.82
1:B:197:LEU:HD12	1:B:306:LEU:HD12	1.61	0.82
1:B:320:GLU:HG2	1:B:321:ALA:H	1.43	0.81
1:A:226:VAL:HA	1:A:229:LEU:HD22	1.63	0.81
1:A:326:PRO:HA	1:A:329:ARG:HG3	1.65	0.78
1:B:396:ILE:HD11	1:B:408:LEU:HB2	1.65	0.78
1:A:452:THR:H	1:A:453:PRO:HD2	1.50	0.76
1:A:402:ASN:HD21	1:A:404:GLU:HB2	1.50	0.76
1:A:402:ASN:HD22	1:A:402:ASN:C	1.88	0.76
1:A:402:ASN:ND2	1:A:404:GLU:H	1.83	0.76
1:A:402:ASN:HD22	1:A:404:GLU:H	1.34	0.75
1:B:271:ARG:NH2	1:B:355:LEU:HD23	2.01	0.74
1:A:357:GLN:H	1:A:357:GLN:HE21	1.33	0.74
1:B:375:LEU:HD12	1:B:375:LEU:H	1.53	0.73
1:B:241:THR:HA	1:B:246:LEU:HD21	1.69	0.73
1:A:172:SER:HB3	1:A:296:VAL:HA	1.70	0.73
1:A:360:ILE:HG22	1:A:361:GLU:HG2	1.68	0.73
1:B:396:ILE:HD12	1:B:396:ILE:O	1.88	0.73
1:A:452:THR:H	1:A:453:PRO:CD	2.02	0.73
1:A:306:LEU:HD23	1:A:306:LEU:H	1.52	0.72
1:A:192:ALA:H	1:A:206:SER:HB3	1.55	0.72
1:B:372:ASP:HA	1:B:375:LEU:HD13	1.70	0.72
1:A:349:VAL:HB	1:A:350:PRO:HD3	1.70	0.72
1:B:349:VAL:HB	1:B:350:PRO:HD3	1.72	0.72
1:B:381:LYS:HG3	1:B:467:LEU:HD11	1.72	0.72
1:B:396:ILE:CD1	1:B:408:LEU:HB2	2.20	0.71
1:A:402:ASN:HD21	1:A:404:GLU:CB	2.02	0.71
1:B:295:LYS:HA	1:B:306:LEU:HD22	1.72	0.71
1:A:179:ILE:HD11	1:A:181:SER:HB2	1.72	0.71
1:B:394:SER:HB2	1:B:395:PRO:CD	2.18	0.71
1:A:309:LEU:HD23	1:A:310:LEU:N	2.06	0.70
1:B:274:VAL:HG12	1:B:275:GLY:H	1.56	0.70
1:A:402:ASN:ND2	1:A:404:GLU:HB2	2.07	0.69
1:B:404:GLU:OE2	1:B:438:ASN:HB3	1.92	0.69
1:B:190:MET:HB2	1:B:312:ALA:O	1.93	0.68
1:B:293:LEU:HD11	1:B:306:LEU:HD11	1.74	0.68
1:A:454:HIS:HB3	1:A:457:VAL:HB	1.75	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:357:GLN:H	1:A:357:GLN:NE2	1.92	0.67
1:A:199:SER:HB3	1:A:201:LYS:HG2	1.75	0.67
1:A:247:PRO:HB3	1:A:266:LYS:NZ	2.08	0.67
1:B:375:LEU:HD23	1:B:393:TYR:OH	1.95	0.66
1:B:369:HIS:HD2	1:B:371:SER:H	1.44	0.66
1:B:269:PHE:CB	1:B:355:LEU:HD12	2.26	0.66
1:B:472:VAL:HG23	2:B:144:HOH:O	1.95	0.65
1:B:375:LEU:HD12	1:B:375:LEU:N	2.11	0.65
1:B:267:SER:HB3	1:B:289:MET:O	1.97	0.64
1:A:208:GLN:O	1:A:211:SER:HB3	1.98	0.64
1:B:357:GLN:H	1:B:357:GLN:HE21	1.43	0.64
1:B:317:SER:HB3	1:B:320:GLU:HB2	1.78	0.64
1:A:369:HIS:CD2	1:A:371:SER:H	2.03	0.64
1:A:451:LYS:HB3	1:A:453:PRO:HD2	1.78	0.63
1:B:271:ARG:CZ	1:B:355:LEU:HD23	2.28	0.63
1:B:375:LEU:CD1	1:B:375:LEU:H	2.12	0.62
1:B:456:SER:HA	1:B:459:GLU:HG2	1.81	0.61
1:A:270:CYS:HB2	1:A:289:MET:CE	2.31	0.60
1:B:320:GLU:HG2	1:B:321:ALA:N	2.13	0.60
1:A:406:ILE:HG13	1:A:408:LEU:HD11	1.84	0.60
1:B:339:CYS:O	1:B:364:VAL:HG12	2.01	0.60
1:A:420:SER:O	1:A:421:ARG:HB2	2.02	0.59
1:B:229:LEU:HD12	1:B:272:VAL:HG22	1.85	0.59
1:B:269:PHE:HB2	1:B:355:LEU:HD12	1.86	0.58
1:A:205:ILE:HD12	1:A:209:VAL:HG11	1.85	0.58
1:A:270:CYS:HB2	1:A:289:MET:HE1	1.84	0.58
1:B:327:GLU:OE2	1:B:327:GLU:HA	2.04	0.58
1:B:233:ASP:HB3	1:B:270:CYS:HB2	1.85	0.58
1:A:394:SER:HB3	1:A:409:ASP:HA	1.85	0.58
1:A:192:ALA:H	1:A:206:SER:CB	2.16	0.57
1:A:247:PRO:HB3	1:A:266:LYS:HZ1	1.68	0.57
1:B:363:PRO:HB2	1:B:366:VAL:HG23	1.86	0.57
1:A:373:ARG:HB2	1:A:374:PRO:HD3	1.86	0.57
1:B:344:VAL:HG13	1:B:348:ALA:CB	2.35	0.57
1:B:447:CYS:SG	1:B:448:PRO:HD2	2.45	0.57
2:A:487:HOH:O	1:B:266:LYS:HG3	2.04	0.57
1:B:410:THR:HG22	1:B:430:HIS:ND1	2.20	0.56
1:B:196:SER:OG	1:B:198:VAL:HG22	2.06	0.56
1:A:452:THR:N	1:A:453:PRO:CD	2.69	0.56
1:B:271:ARG:NH2	1:B:358:ASP:OD2	2.39	0.55
1:A:282:ILE:HD12	1:A:282:ILE:O	2.07	0.55
1:B:233:ASP:OD1	1:B:271:ARG:HD2	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:381:LYS:O	1:A:384:LEU:HD23	2.07	0.54
1:B:458:GLN:O	1:B:462:GLU:HG2	2.08	0.54
1:B:329:ARG:HD3	2:B:61:HOH:O	2.08	0.53
1:A:327:GLU:H	1:A:327:GLU:CD	2.11	0.53
1:A:357:GLN:N	1:A:357:GLN:HE21	2.02	0.53
1:A:333:THR:OG1	1:A:430:HIS:HE1	1.92	0.53
1:A:224:LYS:HB2	1:A:227:GLU:HG3	1.89	0.53
1:A:410:THR:OG1	1:A:430:HIS:HD2	1.92	0.53
1:A:200:GLY:O	1:A:224:LYS:HA	2.09	0.52
1:B:402:ASN:ND2	1:B:441:VAL:HA	2.23	0.52
1:A:406:ILE:HG13	1:A:408:LEU:CD1	2.40	0.52
1:A:199:SER:HB3	1:A:201:LYS:CG	2.39	0.52
1:B:372:ASP:CA	1:B:375:LEU:HD13	2.39	0.51
1:B:274:VAL:HG12	1:B:275:GLY:N	2.24	0.51
1:B:318:GLY:HA2	2:B:60:HOH:O	2.10	0.51
1:A:415:PHE:CZ	1:A:427:ILE:HD13	2.46	0.51
1:B:360:ILE:O	1:B:361:GLU:HB2	2.10	0.51
1:A:410:THR:HG21	1:A:412:TRP:CE2	2.46	0.50
1:A:447:CYS:O	1:A:449:GLU:N	2.44	0.50
1:B:440:ASP:OD2	1:B:443:ALA:HB2	2.12	0.50
1:A:419:TRP:CG	1:B:429:ARG:HD2	2.47	0.50
1:A:429:ARG:HD2	1:B:419:TRP:CG	2.46	0.50
1:A:471:PRO:O	1:A:473:PRO:HD3	2.11	0.50
1:B:394:SER:C	2:B:67:HOH:O	2.50	0.50
1:A:372:ASP:HA	1:A:375:LEU:HD13	1.94	0.50
1:B:321:ALA:CB	1:B:322:PRO:HD3	2.31	0.50
1:A:274:VAL:HG11	2:A:483:HOH:O	2.12	0.50
1:B:357:GLN:NE2	1:B:357:GLN:H	2.08	0.49
1:B:372:ASP:OD2	1:B:397:ARG:HB2	2.13	0.49
1:A:305:GLN:O	1:A:305:GLN:HG3	2.12	0.49
1:A:207:ASN:O	1:A:208:GLN:HB3	2.11	0.49
1:A:243:PRO:C	1:A:245:LYS:H	2.16	0.49
1:A:172:SER:CB	1:A:296:VAL:HG13	2.42	0.49
1:A:189:ASP:HB3	1:A:314:ARG:HB3	1.93	0.49
1:B:365:LEU:HB2	2:B:138:HOH:O	2.12	0.49
1:B:192:ALA:O	1:B:193:VAL:HG13	2.11	0.49
1:B:393:TYR:O	1:B:394:SER:O	2.31	0.49
1:B:246:LEU:HD22	1:B:291:PRO:HB2	1.95	0.49
1:B:431:LYS:HD3	1:B:432:VAL:N	2.28	0.49
1:B:365:LEU:HD11	1:B:376:MET:HG2	1.94	0.49
1:A:307:CYS:SG	1:A:308:CYS:N	2.86	0.48
1:B:210:ALA:O	1:B:211:SER:HB2	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:432:VAL:O	1:A:432:VAL:HG22	2.12	0.48
1:B:320:GLU:O	1:B:323:ARG:N	2.46	0.48
1:A:381:LYS:NZ	2:A:129:HOH:O	2.46	0.48
1:B:192:ALA:HB3	1:B:205:ILE:CD1	2.44	0.48
1:A:464:ILE:O	1:A:468:LEU:HD22	2.14	0.48
1:A:354:TYR:CD2	1:A:401:ARG:HD3	2.49	0.48
1:B:372:ASP:O	1:B:375:LEU:HD13	2.14	0.47
1:A:454:HIS:N	1:A:455:PRO:HD3	2.29	0.47
1:A:195:VAL:O	1:A:307:CYS:HB2	2.15	0.47
1:A:189:ASP:HB3	1:A:314:ARG:CB	2.44	0.47
1:A:404:GLU:OE2	1:A:437:LEU:HB3	2.14	0.47
1:B:344:VAL:CG1	1:B:348:ALA:HB3	2.45	0.47
1:A:338:ASN:ND2	1:A:340:LEU:HD22	2.30	0.46
1:A:377:LEU:O	1:A:381:LYS:HG3	2.15	0.46
1:B:269:PHE:HB3	1:B:355:LEU:HD12	1.98	0.46
1:B:456:SER:HA	1:B:459:GLU:CG	2.46	0.46
1:A:464:ILE:HG22	1:A:468:LEU:CD2	2.46	0.45
1:A:379:ILE:O	1:A:383:ILE:HG13	2.17	0.45
1:A:197:LEU:HB2	1:A:307:CYS:HA	1.99	0.45
1:B:356:PRO:O	1:B:360:ILE:HG13	2.16	0.45
1:A:173:MET:O	1:A:176:VAL:HG22	2.16	0.45
1:A:410:THR:HG22	1:A:411:SER:N	2.31	0.45
1:B:321:ALA:HB3	1:B:322:PRO:CD	2.35	0.45
1:B:271:ARG:HH22	1:B:358:ASP:CG	2.19	0.45
1:B:432:VAL:CG1	2:B:25:HOH:O	2.65	0.45
1:A:360:ILE:O	1:A:361:GLU:HB2	2.16	0.45
1:A:401:ARG:HB3	1:A:441:VAL:O	2.16	0.45
1:A:402:ASN:C	1:A:402:ASN:ND2	2.60	0.45
1:A:452:THR:HG23	1:B:243:PRO:HB2	1.99	0.44
1:B:355:LEU:HD22	1:B:355:LEU:N	2.32	0.44
1:A:295:LYS:HA	1:A:306:LEU:HA	1.99	0.44
1:A:209:VAL:HG23	1:A:210:ALA:H	1.82	0.44
1:A:172:SER:HB3	1:A:296:VAL:CA	2.44	0.44
1:B:293:LEU:HD11	1:B:306:LEU:CD1	2.46	0.44
1:B:344:VAL:HG13	1:B:348:ALA:HB3	1.99	0.44
1:B:232:HIS:N	1:B:232:HIS:CD2	2.86	0.44
1:A:404:GLU:OE2	1:A:438:ASN:N	2.51	0.43
1:B:432:VAL:HG13	2:B:25:HOH:O	2.18	0.43
1:A:247:PRO:HB3	1:A:266:LYS:HZ2	1.81	0.43
1:A:172:SER:HB2	1:A:296:VAL:HG13	2.00	0.43
1:B:384:LEU:HD12	1:B:467:LEU:CD2	2.48	0.43
1:A:376:MET:CE	1:A:380:HIS:HE1	2.31	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:292:TYR:HE2	1:A:311:LEU:HD12	1.84	0.43
1:A:399:ARG:HB2	1:A:405:TYR:CE1	2.53	0.43
1:B:379:ILE:O	1:B:383:ILE:HG13	2.18	0.43
1:A:196:SER:HB3	1:A:199:SER:HB2	2.01	0.43
1:B:422:LYS:HD3	2:B:77:HOH:O	2.18	0.43
1:A:270:CYS:HB2	1:A:289:MET:HE2	2.00	0.43
1:A:241:THR:CA	1:A:246:LEU:HD21	2.41	0.43
1:A:330:ILE:HD13	1:A:429:ARG:NH1	2.34	0.43
1:B:372:ASP:C	1:B:375:LEU:HD13	2.38	0.43
1:A:246:LEU:HD23	1:A:291:PRO:HB2	2.01	0.42
1:A:330:ILE:HD13	1:A:429:ARG:HH12	1.84	0.42
1:B:456:SER:CA	1:B:459:GLU:HG2	2.47	0.42
1:A:422:LYS:HZ3	1:A:471:PRO:HG3	1.85	0.42
1:B:407:THR:O	2:B:25:HOH:O	2.22	0.42
1:B:326:PRO:O	1:B:431:LYS:HE2	2.20	0.42
1:B:369:HIS:HB2	1:B:397:ARG:HB3	2.01	0.42
1:A:274:VAL:O	1:A:274:VAL:HG12	2.20	0.42
1:A:344:VAL:HG12	1:A:348:ALA:HB3	2.00	0.42
1:B:365:LEU:CD1	1:B:376:MET:HG2	2.49	0.41
1:B:332:THR:HG23	1:B:347:ARG:NH2	2.35	0.41
1:A:402:ASN:ND2	1:A:404:GLU:N	2.62	0.41
1:B:192:ALA:HB3	1:B:205:ILE:HD11	2.01	0.41
1:B:395:PRO:N	2:B:67:HOH:O	2.53	0.41
1:A:172:SER:HB3	1:A:296:VAL:HG13	2.02	0.41
1:B:224:LYS:HB2	1:B:227:GLU:HG3	2.02	0.41
1:A:243:PRO:O	1:A:245:LYS:N	2.54	0.41
1:B:289:MET:HA	1:B:311:LEU:O	2.21	0.41
1:B:456:SER:O	1:B:460:LEU:HB2	2.21	0.41
1:A:320:GLU:O	1:A:321:ALA:C	2.60	0.40
1:B:203:LEU:HD22	1:B:204:TYR:H	1.86	0.40
1:A:384:LEU:HD21	2:A:490:HOH:O	2.21	0.40
1:A:410:THR:CG2	1:A:411:SER:N	2.84	0.40
1:A:325:PRO:HA	1:A:326:PRO:HD3	1.98	0.40
1:B:417:ASN:HA	1:B:418:PRO:HD3	1.93	0.40
1:B:318:GLY:O	1:B:323:ARG:HA	2.22	0.40
1:B:246:LEU:CB	1:B:293:LEU:HB2	2.52	0.40
1:B:239:SER:O	1:B:245:LYS:HD2	2.21	0.40
1:B:426:ILE:HD12	1:B:426:ILE:N	2.37	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	257/309 (83%)	224 (87%)	25 (10%)	8 (3%)	7	5
1	B	236/309 (76%)	203 (86%)	29 (12%)	4 (2%)	14	17
All	All	493/618 (80%)	427 (87%)	54 (11%)	12 (2%)	9	9

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	449	GLU
1	A	452	THR
1	B	211	SER
1	B	394	SER
1	A	172	SER
1	A	448	PRO
1	A	453	PRO
1	A	178	GLY
1	A	244	TYR
1	B	208	GLN
1	A	247	PRO
1	B	274	VAL

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	242/279 (87%)	230 (95%)	12 (5%)	34	51
1	B	213/279 (76%)	201 (94%)	12 (6%)	30	45
All	All	455/558 (82%)	431 (95%)	24 (5%)	32	48

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	205	ILE
1	A	229	LEU
1	A	246	LEU
1	A	340	LEU
1	A	342	GLN
1	A	357	GLN
1	A	395	PRO
1	A	402	ASN
1	A	404	GLU
1	A	432	VAL
1	A	460	LEU
1	A	468	LEU
1	B	229	LEU
1	B	306	LEU
1	B	314	ARG
1	B	342	GLN
1	B	357	GLN
1	B	365	LEU
1	B	384	LEU
1	B	397	ARG
1	B	433	ARG
1	B	437	LEU
1	B	460	LEU
1	B	466	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	GLN
1	A	342	GLN
1	A	357	GLN
1	A	369	HIS
1	A	385	GLN
1	A	402	ASN
1	A	430	HIS
1	A	438	ASN
1	A	458	GLN
1	A	463	GLN
1	B	232	HIS
1	B	305	GLN
1	B	316	HIS

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Mol	Chain	Res	Type
1	B	342	GLN
1	B	357	GLN
1	B	367	GLN
1	B	369	HIS
1	B	385	GLN

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	269/309 (87%)	0.33	23 (8%) 11 9	33, 61, 125, 137	0
1	B	250/309 (80%)	0.41	18 (7%) 15 13	35, 62, 115, 129	0
All	All	519/618 (83%)	0.37	41 (7%) 13 11	33, 61, 117, 137	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	209	VAL	5.8
1	B	263	MET	5.4
1	B	208	GLN	5.4
1	B	210	ALA	4.2
1	B	181	SER	4.0
1	B	212	ILE	3.7
1	B	304	SER	3.6
1	A	178	GLY	3.4
1	B	223	ALA	3.3
1	B	448	PRO	3.3
1	A	247	PRO	3.3
1	B	211	SER	3.2
1	A	450	GLU	3.2
1	A	208	GLN	3.2
1	A	188	ALA	3.1
1	A	250	SER	3.0
1	A	171	TYR	3.0
1	A	306	LEU	3.0
1	A	168	GLY	2.9
1	A	180	THR	2.9
1	B	185	VAL	2.8
1	B	205	ILE	2.8
1	A	449	GLU	2.8
1	A	447	CYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	275	GLY	2.7
1	A	321	ALA	2.6
1	A	172	SER	2.5
1	A	173	MET	2.5
1	A	204	TYR	2.5
1	A	473	PRO	2.4
1	A	446	PRO	2.4
1	A	309	LEU	2.4
1	A	448	PRO	2.2
1	A	466	ARG	2.2
1	A	307	CYS	2.2
1	B	270	CYS	2.2
1	B	247	PRO	2.2
1	B	183	TYR	2.2
1	B	274	VAL	2.1
1	B	303	GLU	2.1
1	A	312	ALA	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.