



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:10 PM UTC

PDB ID : 4OAF / pdb_00004oaf
Title : Crystal structure of the cytosolic domain of mouse MiD51
Authors : Loson, O.C.; Kaiser, J.T.; Chan, D.C.
Deposited on : 2014-01-04
Resolution : 2.20 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

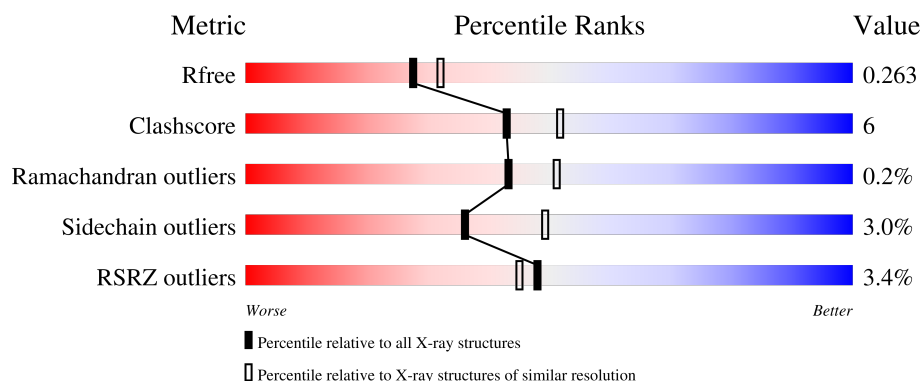
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	335	4% 86% 11% .
1	B	335	4% 83% 13% ..
1	C	335	4% 81% 14% ..
1	D	335	4% 79% 17% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10570 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 Mitochondrial dynamic protein MID51.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2532	1623	424	474	11			
1	B	327	Total	C	N	O	S	0	0	0
			2525	1614	429	471	11			
1	C	326	Total	C	N	O	S	0	0	0
			2539	1628	429	471	11			
1	D	327	Total	C	N	O	S	0	0	0
			2525	1615	424	475	11			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	GLY	-	expression tag	UNP Q8BGV8
A	130	PRO	-	expression tag	UNP Q8BGV8
A	131	LEU	-	expression tag	UNP Q8BGV8
A	132	GLY	-	expression tag	UNP Q8BGV8
A	133	SER	-	expression tag	UNP Q8BGV8
B	129	GLY	-	expression tag	UNP Q8BGV8
B	130	PRO	-	expression tag	UNP Q8BGV8
B	131	LEU	-	expression tag	UNP Q8BGV8
B	132	GLY	-	expression tag	UNP Q8BGV8
B	133	SER	-	expression tag	UNP Q8BGV8
C	129	GLY	-	expression tag	UNP Q8BGV8
C	130	PRO	-	expression tag	UNP Q8BGV8
C	131	LEU	-	expression tag	UNP Q8BGV8
C	132	GLY	-	expression tag	UNP Q8BGV8
C	133	SER	-	expression tag	UNP Q8BGV8
D	129	GLY	-	expression tag	UNP Q8BGV8
D	130	PRO	-	expression tag	UNP Q8BGV8
D	131	LEU	-	expression tag	UNP Q8BGV8
D	132	GLY	-	expression tag	UNP Q8BGV8
D	133	SER	-	expression tag	UNP Q8BGV8

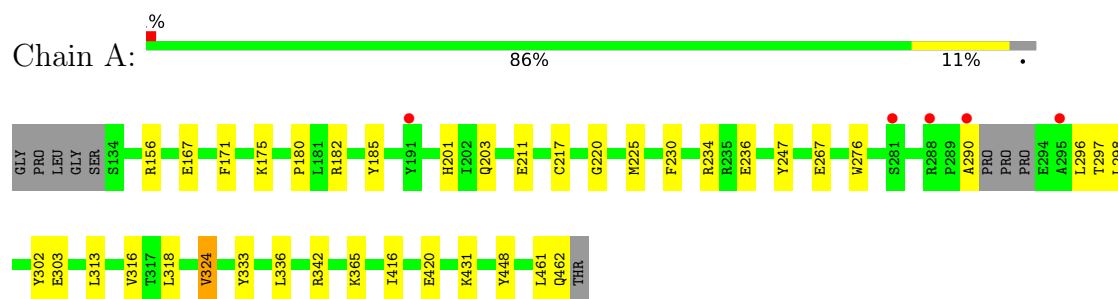
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	135	Total 135	O 135	0	0
2	B	105	Total 105	O 105	0	0
2	C	108	Total 108	O 108	0	0
2	D	101	Total 101	O 101	0	0

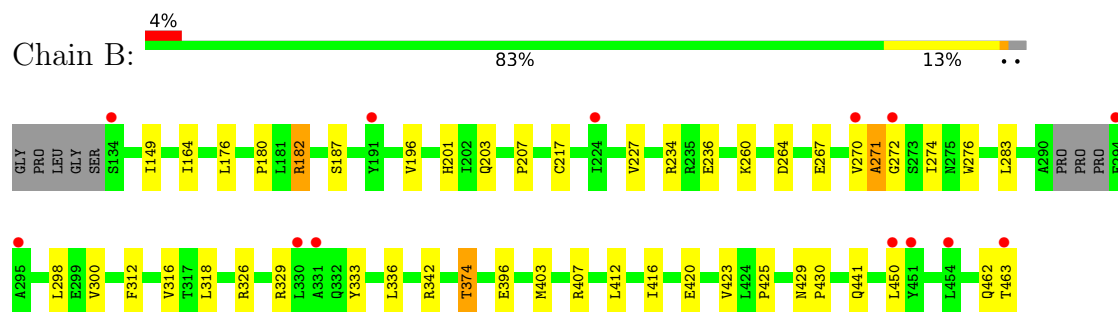
3 Residue-property plots

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.

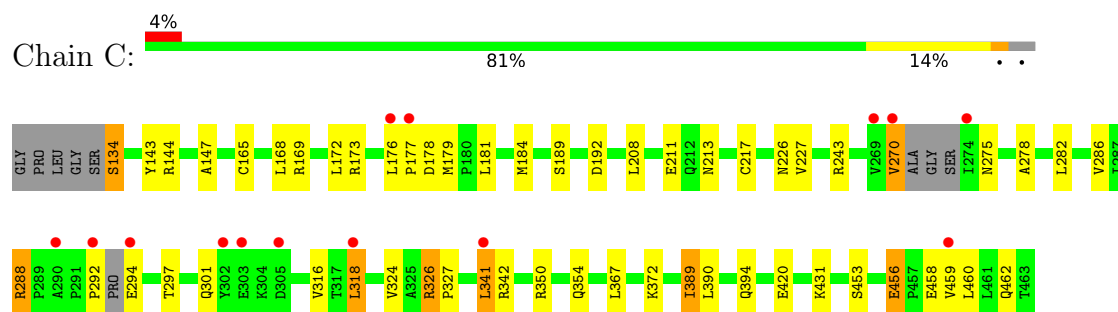
• Molecule 1: Mitochondrial dynamic protein MID51



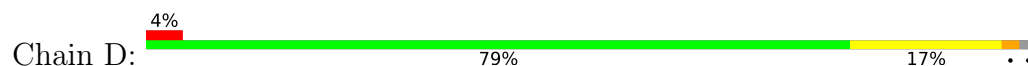
• Molecule 1: Mitochondrial dynamic protein MID51

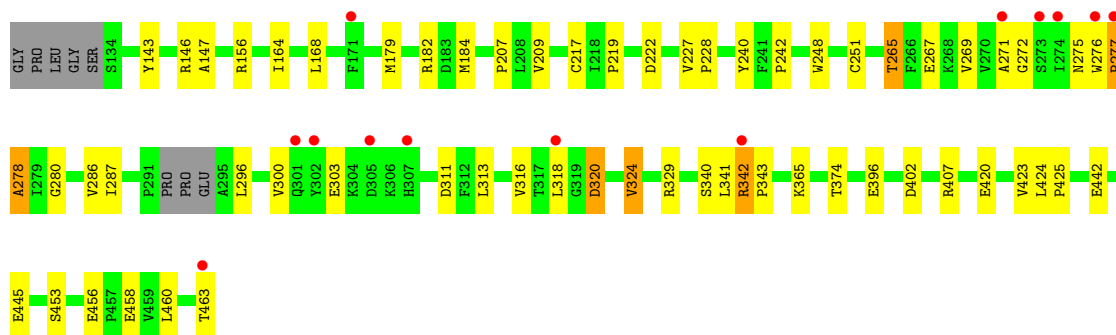


• Molecule 1: Mitochondrial dynamic protein MID51



• Molecule 1: Mitochondrial dynamic protein MID51





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.08Å 78.59Å 102.31Å 90.00° 96.61° 90.00°	Depositor
Resolution (Å)	32.01 – 2.20 32.01 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.7 (32.01-2.20) 85.9 (32.01-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.204 , 0.258 0.206 , 0.263	Depositor DCC
R_{free} test set	2000 reflections (2.86%)	wwPDB-VP
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.808	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 57.9	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	0.96	EDS
Total number of atoms	10570	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 48.54 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.4753e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.52	0/2586	0.82	0/3530
1	B	0.49	0/2577	0.85	4/3518 (0.1%)
1	C	0.50	0/2594	0.91	4/3542 (0.1%)
1	D	0.47	0/2577	0.86	1/3519 (0.0%)
All	All	0.49	0/10334	0.86	9/14109 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	271	ALA	N-CA-C	8.47	123.67	113.16
1	C	326	ARG	CA-C-N	8.43	128.49	119.90
1	C	326	ARG	C-N-CA	8.43	128.49	119.90
1	D	278	ALA	N-CA-C	-7.52	103.05	112.90
1	C	456	GLU	CA-C-N	7.13	127.29	119.87
1	C	456	GLU	C-N-CA	7.13	127.29	119.87
1	B	182	ARG	N-CA-C	-5.07	104.17	110.41
1	B	326	ARG	CA-C-N	5.06	125.06	119.90
1	B	326	ARG	C-N-CA	5.06	125.06	119.90

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	2532	0	2526	29	0
1	B	2525	0	2520	28	0
1	C	2539	0	2539	39	0
1	D	2525	0	2518	37	0
2	A	135	0	0	5	1
2	B	105	0	0	5	1
2	C	108	0	0	11	0
2	D	101	0	0	4	0
All	All	10570	0	10103	126	1

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 6.

All (126) 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:182:ARG:NH1	1:C:179:MET:O	1.96	0.99
1:A:342:ARG:NH2	2:A:542:HOH:O	1.94	0.98
1:B:182:ARG:NH1	1:D:179:MET:O	1.98	0.95
1:D:420:GLU:OE2	2:D:597:HOH:O	1.84	0.95
1:B:187:SER:OG	1:B:342:ARG:NH2	2.05	0.89
1:A:156:ARG:NH1	1:A:303:GLU:OE2	2.08	0.87
1:C:192:ASP:OD2	1:C:342:ARG:NH2	2.12	0.82
1:C:217:CYS:HB3	1:C:318:LEU:HD23	1.64	0.77
1:C:134:SER:OG	2:C:592:HOH:O	2.08	0.72
1:B:180:PRO:HA	1:D:209:VAL:HG11	1.69	0.72
1:C:144:ARG:NH2	2:C:576:HOH:O	2.21	0.72
1:C:286:VAL:HB	1:C:301:GLN:HB3	1.71	0.72
1:D:342:ARG:HG3	1:D:343:PRO:HD3	1.73	0.71
1:B:396:GLU:OE1	2:B:505:HOH:O	2.09	0.70
1:A:167:GLU:OE1	2:A:625:HOH:O	2.08	0.70
1:C:420:GLU:OE1	2:C:589:HOH:O	2.09	0.68
1:D:219:PRO:HG2	1:D:222:ASP:OD2	1.94	0.68
1:B:271:ALA:N	1:B:272:GLY:HA3	2.12	0.65
1:A:225:MET:HE2	1:A:333:TYR:HD1	1.63	0.64
1:D:396:GLU:HG2	1:D:407:ARG:CZ	2.28	0.64
1:B:234:ARG:HD2	1:B:236:GLU:OE2	1.98	0.64
1:C:189:SER:HA	1:C:342:ARG:NH1	2.15	0.62
1:D:320:ASP:N	1:D:320:ASP:OD1	2.33	0.61
1:D:311:ASP:OD1	2:D:522:HOH:O	2.16	0.61
1:B:149:ILE:HG12	1:B:196:VAL:O	2.01	0.60
1:C:372:LYS:NZ	2:C:523:HOH:O	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:156:ARG:NH1	1:D:303:GLU:OE2	2.34	0.59
1:B:260:LYS:NZ	2:B:558:HOH:O	2.36	0.58
1:C:226:ASN:ND2	2:C:526:HOH:O	2.00	0.57
1:C:431:LYS:NZ	2:C:503:HOH:O	2.38	0.56
1:B:176:LEU:HD13	1:B:270:VAL:HG22	1.87	0.56
1:A:333:TYR:HB3	1:A:336:LEU:HD12	1.88	0.56
1:C:292:PRO:HA	1:C:294:GLU:N	2.21	0.56
1:A:313:LEU:HD11	1:A:324:VAL:HG13	1.87	0.55
1:D:143:TYR:HA	1:D:147:ALA:HB3	1.88	0.55
1:B:276:TRP:CH2	1:B:298:LEU:HD21	2.40	0.55
1:C:453:SER:HB2	1:C:460:LEU:HG	1.90	0.53
1:A:302:TYR:CD2	1:A:303:GLU:HG3	2.44	0.53
1:A:234:ARG:HD2	1:A:236:GLU:OE2	2.07	0.53
1:A:247:TYR:HB2	1:C:211:GLU:OE1	2.09	0.53
1:B:283:LEU:O	2:B:597:HOH:O	2.19	0.53
1:D:240:TYR:O	1:D:242:PRO:HD3	2.09	0.53
1:A:461:LEU:O	1:A:462:GLN:HG2	2.09	0.53
1:A:416:ILE:O	1:A:420:GLU:HG3	2.10	0.52
1:D:267:GLU:HG2	1:D:296:LEU:HB3	1.92	0.52
1:C:189:SER:HA	1:C:342:ARG:HH11	1.74	0.52
1:D:271:ALA:N	1:D:272:GLY:HA3	2.24	0.52
1:D:463:THR:HG22	2:D:505:HOH:O	2.10	0.52
1:A:431:LYS:NZ	2:A:530:HOH:O	2.41	0.52
1:C:143:TYR:HA	1:C:147:ALA:HB3	1.91	0.51
1:A:217:CYS:HB3	1:A:318:LEU:HD12	1.92	0.51
1:A:276:TRP:CH2	1:A:298:LEU:HD11	2.46	0.51
1:D:324:VAL:HG21	1:D:342:ARG:HH21	1.76	0.51
1:D:329:ARG:NH2	1:D:442:GLU:OE1	2.41	0.50
1:C:270:VAL:O	2:C:596:HOH:O	2.20	0.50
1:A:365:LYS:NZ	2:A:629:HOH:O	2.40	0.50
1:D:423:VAL:HG12	1:D:425:PRO:HD3	1.93	0.50
1:B:462:GLN:HG2	2:B:567:HOH:O	2.12	0.50
1:B:333:TYR:HB3	1:B:336:LEU:HD12	1.94	0.49
1:D:179:MET:SD	1:D:269:VAL:HG21	2.52	0.49
1:D:453:SER:HB2	1:D:460:LEU:HG	1.95	0.49
1:D:168:LEU:HD13	1:D:184:MET:HE1	1.95	0.49
1:D:217:CYS:HB3	1:D:318:LEU:HD13	1.96	0.48
1:C:367:LEU:HD12	1:C:389:ILE:HD11	1.96	0.48
1:C:462:GLN:OE1	1:C:462:GLN:N	2.38	0.48
1:D:143:TYR:HE1	1:D:365:LYS:HE2	1.79	0.48
1:D:445:GLU:HB3	2:D:550:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:374:THR:OG1	1:D:463:THR:HG21	2.15	0.47
1:B:416:ILE:O	1:B:420:GLU:HG3	2.15	0.47
1:D:277:PRO:HB3	1:D:280:GLY:HA3	1.97	0.47
1:C:326:ARG:HA	1:C:327:PRO:HD3	1.68	0.47
1:A:185:TYR:CE2	1:C:173:ARG:HD2	2.50	0.46
1:D:276:TRP:HA	1:D:278:ALA:H	1.80	0.46
1:D:275:ASN:H	1:D:278:ALA:HB2	1.81	0.46
1:C:213:ASN:OD1	2:C:527:HOH:O	2.21	0.46
1:C:341:LEU:H	1:C:341:LEU:HD12	1.79	0.46
1:D:164:ILE:HD11	1:D:300:VAL:HG11	1.98	0.46
1:C:189:SER:CA	1:C:342:ARG:HH11	2.28	0.46
1:A:156:ARG:NH1	1:A:303:GLU:CD	2.73	0.46
1:D:456:GLU:HB2	1:D:458:GLU:OE2	2.15	0.46
1:A:276:TRP:CZ3	1:A:298:LEU:HD11	2.50	0.46
1:B:267:GLU:HA	1:B:271:ALA:HB3	1.98	0.46
1:A:185:TYR:CD2	1:C:173:ARG:HD2	2.51	0.46
1:B:182:ARG:HB2	1:B:207:PRO:HG2	1.98	0.45
1:C:275:ASN:HB3	1:C:278:ALA:HB3	1.98	0.45
1:C:172:LEU:HD22	1:C:184:MET:HE2	1.98	0.45
1:A:201:HIS:NE2	1:A:203:GLN:OE1	2.50	0.45
1:A:290:ALA:HB3	1:A:297:THR:HG23	1.99	0.44
1:C:350:ARG:O	1:C:354:GLN:HG3	2.17	0.44
1:B:407:ARG:NH2	2:B:505:HOH:O	2.49	0.44
1:A:267:GLU:HB2	1:A:296:LEU:HD21	1.98	0.44
1:A:171:PHE:CE1	1:A:175:LYS:HD2	2.53	0.44
1:A:220:GLY:N	1:A:230:PHE:O	2.49	0.44
1:B:423:VAL:HG12	1:B:425:PRO:HD3	2.00	0.44
1:C:168:LEU:HD13	1:C:184:MET:HE1	2.00	0.43
1:C:177:PRO:O	2:C:561:HOH:O	2.21	0.43
1:C:165:CYS:O	1:C:169:ARG:HG3	2.18	0.43
1:C:390:LEU:O	1:C:394:GLN:HG2	2.18	0.43
1:A:211:GLU:H	1:A:211:GLU:CD	2.26	0.43
1:B:412:LEU:HD22	1:B:450:LEU:HD22	2.01	0.43
1:B:201:HIS:NE2	1:B:203:GLN:OE1	2.52	0.43
1:D:182:ARG:HB2	1:D:207:PRO:HG2	2.01	0.43
1:A:180:PRO:HG3	1:A:247:TYR:HB3	2.01	0.42
1:D:251:CYS:SG	1:D:265:THR:OG1	2.72	0.42
1:C:168:LEU:HD23	1:C:168:LEU:HA	1.85	0.42
1:B:429:ASN:HA	1:B:430:PRO:HD2	1.87	0.42
1:C:176:LEU:HD13	1:C:179:MET:SD	2.59	0.42
1:C:243:ARG:NE	2:C:530:HOH:O	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:374:THR:OG1	1:B:463:THR:HG21	2.19	0.42
1:C:288:ARG:HD3	1:C:288:ARG:HA	1.80	0.42
1:B:164:ILE:HD12	1:B:300:VAL:HG21	2.02	0.42
1:B:270:VAL:O	1:B:274:ILE:HG13	2.20	0.42
1:A:182:ARG:HD2	1:C:181:LEU:O	2.21	0.41
1:D:313:LEU:HD11	1:D:324:VAL:HG13	2.03	0.41
1:C:178:ASP:OD1	2:C:537:HOH:O	2.22	0.41
1:D:179:MET:HE3	1:D:248:TRP:CH2	2.56	0.41
1:B:217:CYS:HB3	1:B:318:LEU:HD12	2.02	0.41
1:C:243:ARG:HH11	1:C:243:ARG:HG2	1.85	0.41
1:D:146:ARG:HH12	1:D:463:THR:C	2.29	0.41
1:D:217:CYS:HB3	1:D:318:LEU:CD1	2.51	0.41
1:D:227:VAL:HA	1:D:228:PRO:HD2	1.99	0.41
1:D:251:CYS:HG	1:D:265:THR:HG1	1.61	0.41
1:A:448:TYR:HB2	2:A:580:HOH:O	2.21	0.40
1:B:276:TRP:CZ3	1:B:298:LEU:HD21	2.56	0.40
1:B:298:LEU:HD11	1:B:312:PHE:CE2	2.57	0.40
1:B:403:MET:O	1:B:407:ARG:HG2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:536:HOH:O	2:B:525:HOH:O[1_655]	2.19	0.01

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	322/335 (96%)	315 (98%)	7 (2%)	0	100	100
1	B	323/335 (96%)	311 (96%)	11 (3%)	1 (0%)	36	42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	320/335 (96%)	312 (98%)	8 (2%)	0	100	100
1	D	323/335 (96%)	305 (94%)	17 (5%)	1 (0%)	36	42
All	All	1288/1340 (96%)	1243 (96%)	43 (3%)	2 (0%)	43	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	329	ARG
1	D	277	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	273/287 (95%)	271 (99%)	2 (1%)	76	87
1	B	271/287 (94%)	266 (98%)	5 (2%)	51	68
1	C	275/287 (96%)	260 (94%)	15 (6%)	19	24
1	D	272/287 (95%)	261 (96%)	11 (4%)	28	38
All	All	1091/1148 (95%)	1058 (97%)	33 (3%)	36	49

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	316	VAL
1	A	324	VAL
1	B	227	VAL
1	B	264	ASP
1	B	316	VAL
1	B	374	THR
1	B	441	GLN
1	C	134	SER
1	C	208	LEU
1	C	227	VAL

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Mol	Chain	Res	Type
1	C	270	VAL
1	C	282	LEU
1	C	288	ARG
1	C	297	THR
1	C	316	VAL
1	C	318	LEU
1	C	324	VAL
1	C	341	LEU
1	C	389	ILE
1	C	456	GLU
1	C	458	GLU
1	C	459	VAL
1	D	265	THR
1	D	286	VAL
1	D	287	ILE
1	D	316	VAL
1	D	320	ASP
1	D	324	VAL
1	D	340	SER
1	D	341	LEU
1	D	342	ARG
1	D	402	ASP
1	D	424	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	441	GLN
1	D	275	ASN
1	D	429	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 [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	326/335 (97%)	-0.09	5 (1%) 72 69	22, 35, 58, 77	0
1	B	327/335 (97%)	0.23	13 (3%) 42 39	28, 44, 74, 98	0
1	C	326/335 (97%)	0.16	14 (4%) 40 36	29, 41, 77, 96	0
1	D	327/335 (97%)	0.20	13 (3%) 42 39	26, 41, 75, 95	0
All	All	1306/1340 (97%)	0.12	45 (3%) 48 45	22, 41, 73, 98	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	274	ILE	5.5
1	D	276	TRP	4.4
1	C	269	VAL	4.3
1	C	270	VAL	4.1
1	B	454	LEU	4.0
1	B	134	SER	3.8
1	D	277	PRO	3.7
1	C	303	GLU	3.6
1	A	290	ALA	3.4
1	D	302	TYR	3.2
1	C	292	PRO	3.1
1	B	451	TYR	3.0
1	B	463	THR	2.8
1	C	294	GLU	2.8
1	B	294	GLU	2.7
1	D	305	ASP	2.6
1	D	273	SER	2.6
1	A	295	ALA	2.5
1	C	341	LEU	2.5
1	C	177	PRO	2.5
1	D	171	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	176	LEU	2.4
1	B	224	ILE	2.4
1	C	459	VAL	2.3
1	B	330	LEU	2.3
1	D	342	ARG	2.3
1	D	463	THR	2.3
1	C	290	ALA	2.2
1	C	274	ILE	2.2
1	A	288	ARG	2.2
1	D	271	ALA	2.2
1	B	270	VAL	2.2
1	D	307	HIS	2.2
1	C	305	ASP	2.1
1	B	272	GLY	2.1
1	B	331	ALA	2.1
1	C	318	LEU	2.1
1	B	191	TYR	2.1
1	C	302	TYR	2.1
1	D	301	GLN	2.1
1	A	281	SER	2.0
1	B	295	ALA	2.0
1	B	450	LEU	2.0
1	A	191	TYR	2.0
1	D	318	LEU	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.