



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:15 PM GMT

PDB ID : 4OAF
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

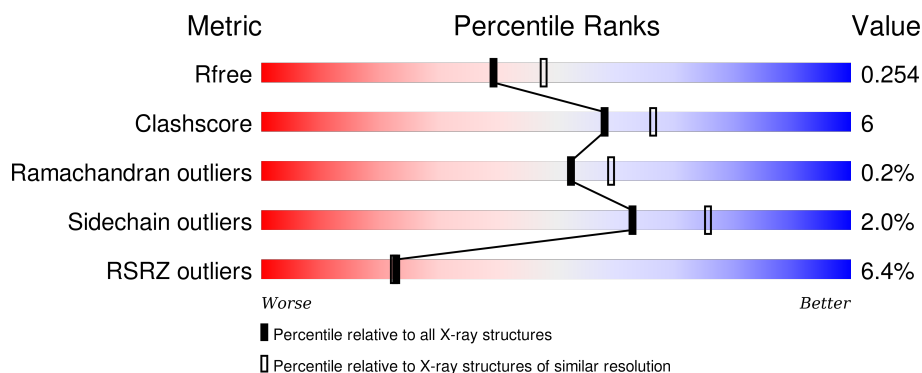
1 Overall quality at a glance

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}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>4%</div> <div>87%</div> <div>10%</div> <div>.</div> </div>
1	B	335	<div> <div>9%</div> <div>84%</div> <div>13%</div> <div>.</div> </div>
1	C	335	<div> <div>7%</div> <div>81%</div> <div>15%</div> <div>..</div> </div>
1	D	335	<div> <div>5%</div> <div>80%</div> <div>16%</div> <div>..</div> </div>

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

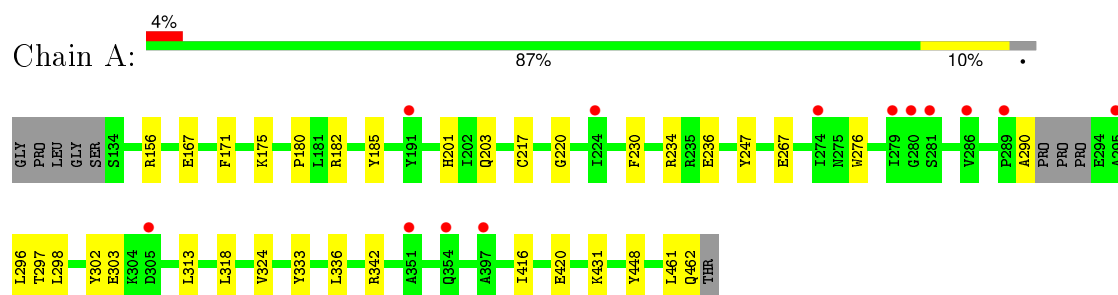
- Molecule 2 is water.

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

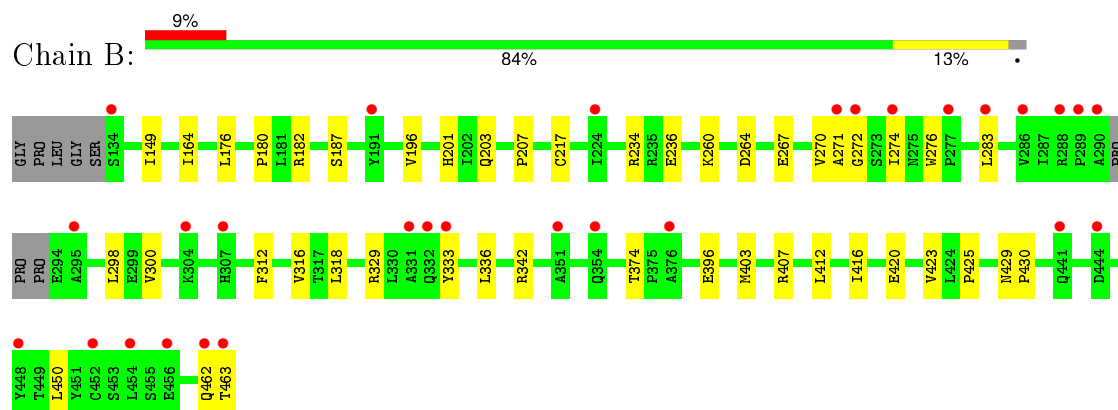
3 Residue-property plots [i](#)

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.

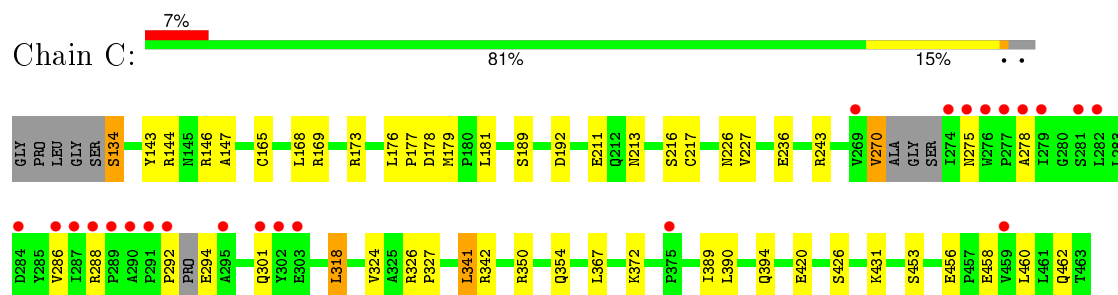
• Molecule 1: Mitochondrial dynamic protein MID51



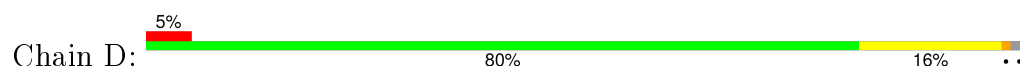
• Molecule 1: Mitochondrial dynamic protein MID51

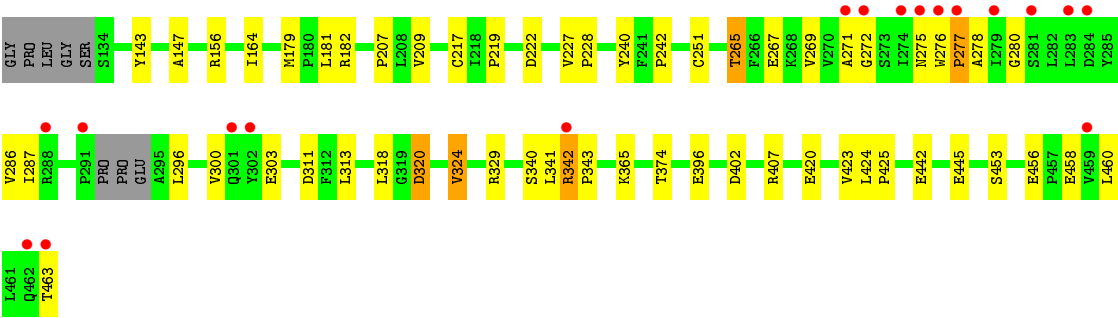


• Molecule 1: Mitochondrial dynamic protein MID51



• 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 39.30 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.7 (32.01-2.20) 86.0 (39.30-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.204 , 0.258 0.199 , 0.254	Depositor DCC
R_{free} test set	1781 reflections (2.84%)	DCC
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.790	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 58.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	7 of 70031 reflections (0.010%)	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 48.43 % 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 8.6155e-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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.39	0/2586	0.55	0/3530
1	B	0.36	0/2577	0.54	0/3518
1	C	0.38	0/2594	0.58	0/3542
1	D	0.35	0/2577	0.56	0/3519
All	All	0.37	0/10334	0.56	0/14109

There are no bond length outliers.

There are no bond angle outliers.

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	26	0
1	B	2525	0	2520	28	0
1	C	2539	0	2539	38	0
1	D	2525	0	2518	35	0
2	A	135	0	0	4	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	120	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 (120) 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:C:286:VAL:HB	1:C:301:GLN:HB3	1.71	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: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: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:OD1	1:D:320:ASP:N	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
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:D:240:TYR:O	1:D:242:PRO:HD3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:LEU:O	2:B:597:HOH:O	2.19	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:D:324:VAL:HG21	1:D:342:ARG:HH21	1.76	0.51
1:A:276:TRP:CH2	1:A:298:LEU:HD11	2.46	0.51
1:A:217:CYS:HB3	1:A:318:LEU:HD12	1.92	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: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:179:MET:HE2	1:D:181:LEU:HD11	1.95	0.48
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:D:143:TYR:HE1	1:D:365:LYS:HE2	1.79	0.48
1:C:462:GLN:OE1	1:C:462:GLN:N	2.38	0.48
1:D:445:GLU:HB3	2:D:550:HOH:O	2.14	0.47
1:D:374:THR:OG1	1:D:463:THR:HG21	2.15	0.47
1:D:277:PRO:HB3	1:D:280:GLY:HA3	1.97	0.47
1:B:416:ILE:O	1:B:420:GLU:HG3	2.15	0.47
1:C:326:ARG:HA	1:C:327:PRO:HD3	1.68	0.47
1:D:276:TRP:HA	1:D:278:ALA:H	1.80	0.46
1:A:185:TYR:CE2	1:C:173:ARG:HD2	2.50	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:D:164:ILE:HD11	1:D:300:VAL:HG11	1.98	0.46
1:C:341:LEU:H	1:C:341:LEU:HD12	1.79	0.46
1:C:189:SER:CA	1:C:342:ARG:HH11	2.28	0.46
1:D:456:GLU:HB2	1:D:458:GLU:OE2	2.15	0.46
1:B:267:GLU:HA	1:B:271:ALA:HB3	1.98	0.46
1:A:276:TRP:CZ3	1:A:298:LEU:HD11	2.50	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:275:ASN:HB3	1:C:278:ALA:HB3	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:177:PRO:O	2:C:561:HOH:O	2.21	0.43
1:C:390:LEU:O	1:C:394:GLN:HG2	2.18	0.43
1:C:165:CYS:O	1:C:169:ARG:HG3	2.18	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:HA	1:C:168:LEU:HD23	1.85	0.42
1:C:176:LEU:HD13	1:C:179:MET:SD	2.59	0.42
1:B:429:ASN:HA	1:B:430:PRO:HD2	1.87	0.42
1:C:243:ARG:NE	2:C:530:HOH:O	2.52	0.42
1:B:374:THR:OG1	1:B:463:THR:HG21	2.19	0.42
1:C:288:ARG:HA	1:C:288:ARG:HD3	1.80	0.42
1:B:270:VAL:O	1:B:274:ILE:HG13	2.20	0.42
1:C:216:SER:OG	1:C:236:GLU:OE2	2.31	0.42
1:B:164:ILE:HD12	1:B:300:VAL:HG21	2.02	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:A:156:ARG:NH1	1:A:303:GLU:CD	2.73	0.41
1:D:217:CYS:HB3	1:D:318:LEU:CD1	2.51	0.41
1:D:251:CYS:HG	1:D:265:THR:HG1	1.61	0.41
1:C:243:ARG:HG2	1:C:243:ARG:HH11	1.85	0.41
1:B:217:CYS:HB3	1:B:318:LEU:HD12	2.02	0.41
1:D:227:VAL:HA	1:D:228:PRO:HD2	1.99	0.41
1:B:298:LEU:HD11	1:B:312:PHE:CE2	2.57	0.40
1:B:276:TRP:CZ3	1:B:298:LEU:HD21	2.56	0.40
1:A:448:TYR:HB2	2:A:580:HOH:O	2.21	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 [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	322/335 (96%)	315 (98%)	7 (2%)	0	100	100
1	B	323/335 (96%)	311 (96%)	11 (3%)	1 (0%)	46	50
1	C	320/335 (96%)	312 (98%)	8 (2%)	0	100	100
1	D	323/335 (96%)	305 (94%)	17 (5%)	1 (0%)	46	50
All	All	1288/1340 (96%)	1243 (96%)	43 (3%)	2 (0%)	52	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	329	ARG
1	D	277	PRO

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	273/287 (95%)	273 (100%)	0	100	100
1	B	271/287 (94%)	269 (99%)	2 (1%)	88	94
1	C	275/287 (96%)	265 (96%)	10 (4%)	42	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	D	272/287 (95%)	262 (96%)	10 (4%)	41 50
All	All	1091/1148 (95%)	1069 (98%)	22 (2%)	63 76

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	264	ASP
1	B	316	VAL
1	C	134	SER
1	C	146	ARG
1	C	227	VAL
1	C	270	VAL
1	C	318	LEU
1	C	324	VAL
1	C	341	LEU
1	C	426	SER
1	C	456	GLU
1	C	458	GLU
1	D	265	THR
1	D	286	VAL
1	D	287	ILE
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

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates 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%)	0.08	13 (3%) 42 41	22, 35, 58, 77	0
1	B	327/335 (97%)	0.37	29 (8%) 12 11	28, 44, 74, 98	0
1	C	326/335 (97%)	0.31	23 (7%) 19 18	29, 41, 77, 96	0
1	D	327/335 (97%)	0.23	18 (5%) 29 28	26, 41, 75, 95	0
All	All	1306/1340 (97%)	0.25	83 (6%) 23 22	22, 41, 73, 98	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	274	ILE	9.5
1	D	463	THR	7.0
1	D	277	PRO	6.6
1	C	291	PRO	5.9
1	C	290	ALA	5.9
1	C	288	ARG	4.9
1	B	463	THR	4.6
1	C	286	VAL	4.5
1	B	274	ILE	4.5
1	D	275	ASN	4.4
1	D	276	TRP	4.0
1	B	288	ARG	3.8
1	D	272	GLY	3.7
1	C	276	TRP	3.7
1	C	292	PRO	3.6
1	C	275	ASN	3.6
1	B	354	GLN	3.6
1	B	286	VAL	3.6
1	A	280	GLY	3.5
1	C	302	TYR	3.5
1	B	224	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	286	VAL	3.4
1	B	454	LEU	3.3
1	C	277	PRO	3.3
1	D	291	PRO	3.3
1	D	459	VAL	3.3
1	B	295	ALA	3.3
1	B	441	GLN	3.3
1	C	295	ALA	3.2
1	C	303	GLU	3.2
1	C	279	ILE	3.1
1	D	283	LEU	3.1
1	B	277	PRO	3.1
1	B	290	ALA	3.1
1	D	271	ALA	3.0
1	B	289	PRO	2.9
1	C	459	VAL	2.9
1	B	332	GLN	2.9
1	A	295	ALA	2.8
1	A	351	ALA	2.8
1	C	289	PRO	2.8
1	A	354	GLN	2.7
1	C	278	ALA	2.7
1	D	302	TYR	2.7
1	A	397	ALA	2.7
1	A	305	ASP	2.6
1	D	281	SER	2.6
1	C	282	LEU	2.6
1	C	281	SER	2.5
1	A	281	SER	2.5
1	D	284	ASP	2.5
1	C	301	GLN	2.5
1	A	274	ILE	2.5
1	C	269	VAL	2.4
1	B	351	ALA	2.4
1	C	375	PRO	2.4
1	A	224	ILE	2.4
1	B	462	GLN	2.3
1	C	274	ILE	2.2
1	A	191	TYR	2.2
1	B	452	CYS	2.2
1	B	283	LEU	2.2
1	A	289	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	331	ALA	2.2
1	D	342	ARG	2.2
1	D	301	GLN	2.2
1	B	333	TYR	2.2
1	B	456	GLU	2.2
1	B	191	TYR	2.1
1	B	134	SER	2.1
1	B	304	LYS	2.1
1	B	376	ALA	2.1
1	B	444	ASP	2.1
1	D	288	ARG	2.1
1	B	271	ALA	2.1
1	C	287	ILE	2.1
1	D	279	ILE	2.1
1	A	279	ILE	2.1
1	C	284	ASP	2.0
1	B	272	GLY	2.0
1	B	448	TYR	2.0
1	B	307	HIS	2.0
1	D	462	GLN	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 carbohydrates 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.