



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:02 AM UTC

PDB ID : 2F00 / pdb_00002f00
Title : Escherichia coli MurC
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Deposited on : 2005-11-10
Resolution : 2.50 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

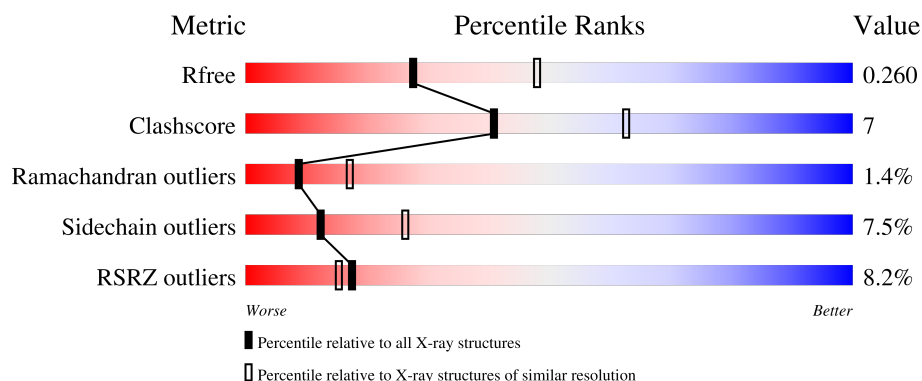
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 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	491	15% 78% 16% • •
1	B	491	83% 13% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7582 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 UDP-N-acetylmuramate--L-alanine ligase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	Se	0	0	0
			3524	2213	626	668	2	15			
1	B	483	Total	C	N	O	S	Se	0	0	0
			3678	2312	660	689	2	15			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P17952
A	16	MSE	MET	modified residue	UNP P17952
A	32	MSE	MET	modified residue	UNP P17952
A	61	MSE	MET	modified residue	UNP P17952
A	111	MSE	MET	modified residue	UNP P17952
A	116	MSE	MET	modified residue	UNP P17952
A	136	MSE	MET	modified residue	UNP P17952
A	188	MSE	MET	modified residue	UNP P17952
A	200	MSE	MET	modified residue	UNP P17952
A	229	MSE	MET	modified residue	UNP P17952
A	282	MSE	MET	modified residue	UNP P17952
A	347	MSE	MET	modified residue	UNP P17952
A	377	MSE	MET	modified residue	UNP P17952
A	406	MSE	MET	modified residue	UNP P17952
A	449	MSE	MET	modified residue	UNP P17952
B	1	MSE	MET	modified residue	UNP P17952
B	16	MSE	MET	modified residue	UNP P17952
B	32	MSE	MET	modified residue	UNP P17952
B	61	MSE	MET	modified residue	UNP P17952
B	111	MSE	MET	modified residue	UNP P17952
B	116	MSE	MET	modified residue	UNP P17952
B	136	MSE	MET	modified residue	UNP P17952
B	188	MSE	MET	modified residue	UNP P17952
B	200	MSE	MET	modified residue	UNP P17952
B	229	MSE	MET	modified residue	UNP P17952

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Chain	Residue	Modelled	Actual	Comment	Reference
B	282	MSE	MET	modified residue	UNP P17952
B	347	MSE	MET	modified residue	UNP P17952
B	377	MSE	MET	modified residue	UNP P17952
B	406	MSE	MET	modified residue	UNP P17952
B	449	MSE	MET	modified residue	UNP P17952

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	1	0
2	B	1	Total Mg 1 1	1	0

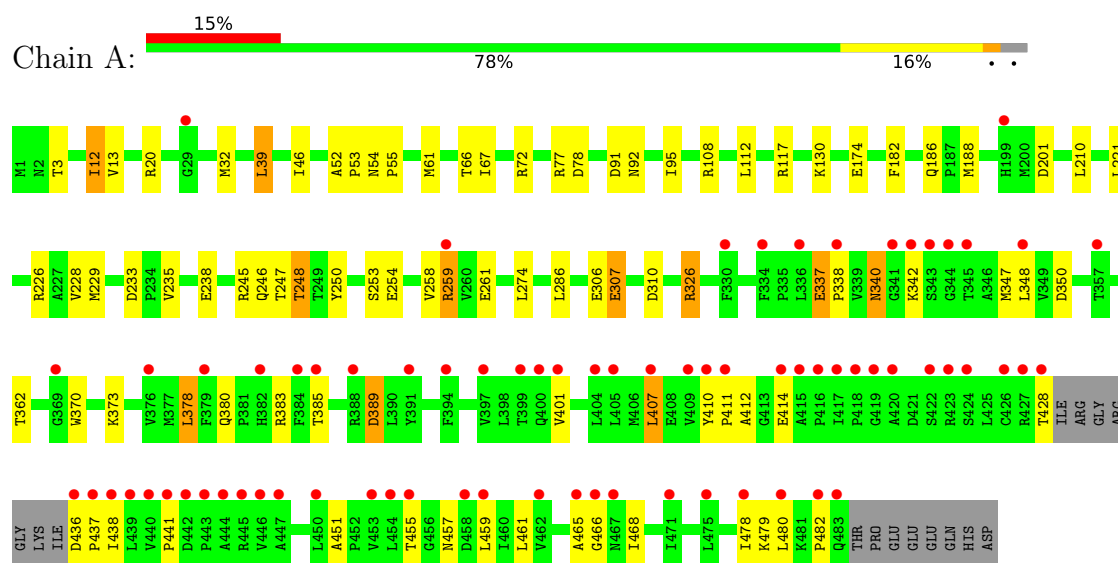
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	135	Total O 135 135	5	0
3	B	243	Total O 243 243	5	0

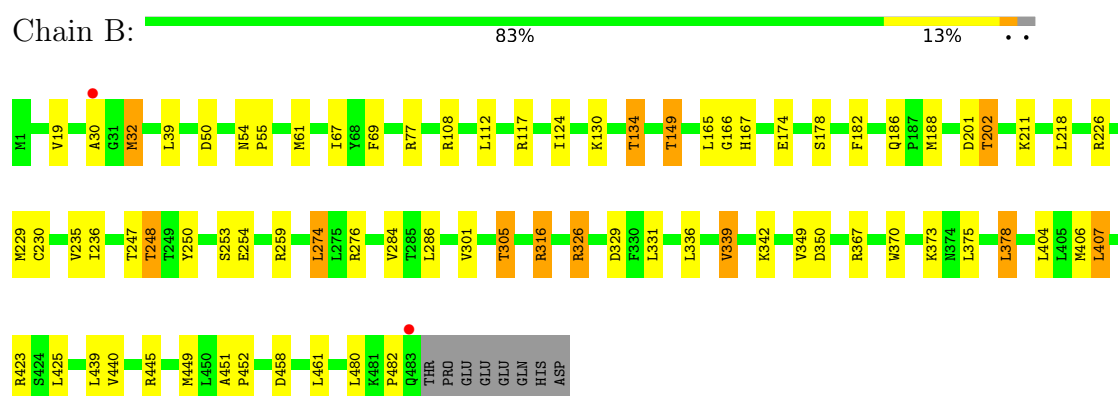
3 Residue-property plots

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: UDP-N-acetylmuramate--L-alanine ligase



- Molecule 1: UDP-N-acetylmuramate--L-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.93Å 93.13Å 176.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.00 – 2.50 88.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (88.00-2.50) 99.2 (88.00-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.217 , 0.260 0.216 , 0.260	Depositor DCC
R_{free} test set	2137 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.372	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7582	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.51	0/3580	0.81	1/4854 (0.0%)
1	B	0.52	0/3736	0.82	1/5053 (0.0%)
All	All	0.52	0/7316	0.82	2/9907 (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	A	482	PRO	N-CA-CB	6.46	110.03	103.25
1	B	482	PRO	N-CA-CB	5.94	109.49	103.25

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	3524	0	3377	49	0
1	B	3678	0	3646	48	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	135	0	0	1	0
3	B	243	0	0	3	0
All	All	7582	0	7023	93	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 (93) 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:259:ARG:HG2	1:A:259:ARG:HH11	1.10	1.15
1:A:259:ARG:HG2	1:A:259:ARG:NH1	1.78	0.94
1:A:337:GLU:HB3	1:A:338:PRO:HD3	1.55	0.86
1:B:149:THR:HG22	1:B:166:GLY:H	1.40	0.84
1:B:32:MSE:HE3	1:B:108:ARG:HG3	1.63	0.81
1:B:378:LEU:HD21	1:B:407:LEU:HD13	1.62	0.81
1:B:226:ARG:HD3	1:B:248:THR:HG23	1.64	0.78
1:B:301:VAL:O	1:B:305:THR:HG23	1.85	0.77
1:A:259:ARG:HH11	1:A:259:ARG:CG	1.96	0.76
1:B:124:ILE:HD13	1:B:134:THR:HB	1.68	0.76
1:A:221:LEU:O	1:A:245:ARG:NH1	2.20	0.74
1:B:316:ARG:HH11	1:B:316:ARG:HG2	1.53	0.72
1:B:149:THR:HG23	3:B:943:HOH:O	1.92	0.68
1:B:178:SER:O	1:B:202:THR:HG21	1.93	0.68
1:A:186:GLN:HE22	1:B:186:GLN:HE21	1.44	0.65
1:A:347:MSE:HE3	1:A:459:LEU:HB2	1.79	0.64
1:A:259:ARG:NH1	1:A:261:GLU:OE1	2.30	0.64
1:A:54:ASN:HB2	1:A:55:PRO:CD	2.29	0.63
1:B:130:LYS:O	1:B:134:THR:HG22	1.99	0.63
1:B:276:ARG:HH11	1:B:305:THR:HG21	1.64	0.63
1:B:130:LYS:O	1:B:134:THR:CG2	2.49	0.60
1:B:117:ARG:NH2	3:B:919:HOH:O	2.35	0.60
1:A:436:ASP:N	1:A:437:PRO:CD	2.65	0.59
1:A:54:ASN:HB2	1:A:55:PRO:HD2	1.84	0.59
1:A:130:LYS:HB3	1:A:174:GLU:HG3	1.85	0.58
1:B:407:LEU:HD12	1:B:440:VAL:HB	1.86	0.57
1:B:370:TRP:HB3	1:B:373:LYS:HG3	1.85	0.57
1:B:211:LYS:HD3	1:B:235:VAL:HG13	1.87	0.56
1:A:479:LYS:O	1:A:480:LEU:HB2	2.04	0.56
1:A:307:GLU:HG2	3:A:983:HOH:O	2.05	0.56
1:B:254:GLU:H	1:B:254:GLU:CD	2.14	0.56
1:B:54:ASN:HB2	1:B:55:PRO:CD	2.36	0.55
1:B:54:ASN:HB2	1:B:55:PRO:HD2	1.89	0.54
1:A:229:MSE:HE3	1:A:247:THR:CG2	2.39	0.53
1:A:20:ARG:HH22	1:A:78:ASP:HB3	1.74	0.52
1:A:233:ASP:OD1	1:A:235:VAL:HG12	2.08	0.52
1:A:117:ARG:CZ	1:B:188:MSE:HE1	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:326:ARG:HD2	1:B:350:ASP:OD2	2.09	0.52
1:B:149:THR:HG22	1:B:166:GLY:N	2.16	0.52
1:A:77:ARG:O	1:A:78:ASP:HB2	2.09	0.51
1:A:12:ILE:HG22	1:A:13:VAL:HG23	1.91	0.51
1:A:186:GLN:HE22	1:B:186:GLN:NE2	2.08	0.51
1:A:61:MSE:HE3	1:A:67:ILE:HB	1.93	0.50
1:B:445:ARG:O	1:B:449:MSE:HG2	2.12	0.50
1:A:478:ILE:HG13	1:A:480:LEU:H	1.76	0.49
1:B:226:ARG:HD3	1:B:248:THR:CG2	2.40	0.49
1:A:92:ASN:HB3	1:A:95:ILE:HD13	1.93	0.49
1:B:230:CYS:HB2	1:B:250:TYR:CZ	2.48	0.49
1:B:423:ARG:HB3	1:B:439:LEU:HD11	1.94	0.49
1:B:329:ASP:HB3	1:B:350:ASP:HB3	1.95	0.49
1:A:253:SER:O	1:A:259:ARG:HD3	2.13	0.48
1:B:149:THR:CG2	1:B:166:GLY:H	2.19	0.48
1:A:326:ARG:NH1	1:A:350:ASP:OD1	2.46	0.48
1:A:326:ARG:HD2	1:A:350:ASP:OD2	2.13	0.48
1:A:226:ARG:NH2	1:A:246:GLN:HB3	2.29	0.48
1:B:274:LEU:HD22	1:B:284:VAL:HG21	1.96	0.47
1:A:188:MSE:HE1	1:B:117:ARG:NE	2.30	0.47
1:A:229:MSE:HE3	1:A:247:THR:HG21	1.95	0.47
1:A:378:LEU:HD22	1:A:407:LEU:HD12	1.96	0.47
1:B:378:LEU:HD21	1:B:407:LEU:CD1	2.39	0.47
1:A:370:TRP:HB3	1:A:373:LYS:HG3	1.97	0.47
1:A:436:ASP:N	1:A:437:PRO:HD3	2.31	0.46
1:A:254:GLU:HA	1:A:259:ARG:CD	2.46	0.46
1:B:50:ASP:O	1:B:69:PHE:HA	2.16	0.46
1:B:349:VAL:HB	1:B:461:LEU:HD23	1.98	0.46
1:A:389:ASP:OD1	1:A:389:ASP:N	2.49	0.45
1:A:226:ARG:HH21	1:A:246:GLN:HB3	1.81	0.45
1:A:347:MSE:HB2	1:A:459:LEU:HA	1.98	0.45
1:B:253:SER:O	1:B:259:ARG:HG3	2.17	0.45
1:B:134:THR:HG21	1:B:174:GLU:HB2	1.99	0.44
1:B:378:LEU:CD2	1:B:407:LEU:HD13	2.40	0.44
1:A:32:MSE:SE	1:A:108:ARG:HG3	2.68	0.44
1:A:340:ASN:ND2	1:A:451:ALA:O	2.50	0.44
1:A:91:ASP:OD1	1:A:91:ASP:N	2.45	0.43
1:B:336:LEU:HA	1:B:339:VAL:CG1	2.48	0.43
1:A:54:ASN:CB	1:A:55:PRO:CD	2.96	0.43
1:B:276:ARG:HH11	1:B:305:THR:CG2	2.30	0.43
1:B:149:THR:HB	1:B:165:LEU:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:MSE:HA	1:B:32:MSE:CE	2.48	0.42
1:B:61:MSE:HE3	1:B:67:ILE:HB	2.00	0.42
1:B:316:ARG:HH11	1:B:316:ARG:CG	2.25	0.42
1:A:250:TYR:HA	1:A:258:VAL:O	2.19	0.42
1:A:380:GLN:HB2	1:A:468:ILE:HD12	2.01	0.42
1:A:337:GLU:HB3	1:A:338:PRO:CD	2.38	0.41
1:A:410:TYR:CE1	1:A:412:ALA:HB2	2.55	0.41
1:A:39:LEU:HB3	1:A:46:ILE:HD11	2.02	0.41
1:A:253:SER:O	1:A:259:ARG:CD	2.68	0.41
1:B:451:ALA:HB3	1:B:452:PRO:HD3	2.01	0.41
1:A:52:ALA:HA	1:A:53:PRO:HD2	1.81	0.41
1:B:167:HIS:HB2	3:B:1098:HOH:O	2.20	0.41
1:B:404:LEU:HD21	1:B:406:MSE:CE	2.51	0.40
1:A:228:VAL:HG22	1:A:248:THR:HG23	2.03	0.40
1:B:229:MSE:HE2	1:B:236:ILE:HG23	2.03	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	472/491 (96%)	445 (94%)	16 (3%)	11 (2%)	5	8
1	B	481/491 (98%)	466 (97%)	13 (3%)	2 (0%)	30	49
All	All	953/982 (97%)	911 (96%)	29 (3%)	13 (1%)	9	17

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	30	ALA
1	A	414	GLU
1	A	337	GLU

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Mol	Chain	Res	Type
1	A	342	LYS
1	A	383	ARG
1	A	455	THR
1	A	466	GLY
1	B	342	LYS
1	A	340	ASN
1	A	441	PRO
1	A	411	PRO
1	A	465	ALA
1	A	401	VAL

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	351/381 (92%)	323 (92%)	28 (8%)	11	24
1	B	382/381 (100%)	355 (93%)	27 (7%)	13	29
All	All	733/762 (96%)	678 (92%)	55 (8%)	12	26

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	12	ILE
1	A	39	LEU
1	A	66	THR
1	A	72	ARG
1	A	112	LEU
1	A	182	PHE
1	A	201	ASP
1	A	210	LEU
1	A	238	GLU
1	A	248	THR
1	A	259	ARG
1	A	274	LEU

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Mol	Chain	Res	Type
1	A	286	LEU
1	A	306	GLU
1	A	307	GLU
1	A	310	ASP
1	A	326	ARG
1	A	348	LEU
1	A	362	THR
1	A	378	LEU
1	A	385	THR
1	A	389	ASP
1	A	407	LEU
1	A	428	THR
1	A	438	ILE
1	A	457	ASN
1	A	461	LEU
1	B	19	VAL
1	B	32	MSE
1	B	39	LEU
1	B	77	ARG
1	B	112	LEU
1	B	134	THR
1	B	149	THR
1	B	182	PHE
1	B	201	ASP
1	B	202	THR
1	B	218	LEU
1	B	247	THR
1	B	248	THR
1	B	274	LEU
1	B	286	LEU
1	B	305	THR
1	B	316	ARG
1	B	326	ARG
1	B	331	LEU
1	B	339	VAL
1	B	367	ARG
1	B	375	LEU
1	B	378	LEU
1	B	407	LEU
1	B	425	LEU
1	B	458	ASP
1	B	480	LEU

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

Mol	Chain	Res	Type
1	A	62	ASN
1	A	184	HIS
1	A	186	GLN
1	A	212	GLN
1	A	265	GLN
1	A	457	ASN
1	B	62	ASN
1	B	186	GLN
1	B	322	GLN

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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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	461/491 (93%)	0.63	74 (16%) 4 4	18, 42, 99, 100	0
1	B	468/491 (95%)	-0.04	2 (0%) 88 86	17, 34, 56, 64	0
All	All	929/982 (94%)	0.29	76 (8%) 17 15	17, 37, 83, 100	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	428	THR	4.5
1	A	438	ILE	4.4
1	A	401	VAL	4.2
1	A	417	ILE	4.1
1	A	483	GLN	4.0
1	B	30	ALA	3.8
1	A	443	PRO	3.5
1	A	440	VAL	3.4
1	A	420	ALA	3.4
1	A	259	ARG	3.4
1	A	442	ASP	3.3
1	A	446	VAL	3.3
1	A	385	THR	3.3
1	A	405	LEU	3.2
1	A	454	LEU	3.2
1	A	388	ARG	3.2
1	A	450	LEU	3.0
1	A	404	LEU	3.0
1	A	453	VAL	2.9
1	A	343	SER	2.9
1	A	400	GLN	2.9
1	B	483	GLN	2.9
1	A	455	THR	2.9
1	A	418	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	345	THR	2.8
1	A	415	ALA	2.8
1	A	444	ALA	2.8
1	A	391	TYR	2.7
1	A	439	LEU	2.7
1	A	426	CYS	2.7
1	A	336	LEU	2.7
1	A	480	LEU	2.6
1	A	437	PRO	2.6
1	A	482	PRO	2.6
1	A	376	VAL	2.6
1	A	338	PRO	2.6
1	A	427	ARG	2.6
1	A	471	ILE	2.5
1	A	465	ALA	2.5
1	A	411	PRO	2.5
1	A	459	LEU	2.5
1	A	29	GLY	2.4
1	A	416	PRO	2.4
1	A	382	HIS	2.4
1	A	334	PHE	2.4
1	A	441	PRO	2.4
1	A	341	GLY	2.4
1	A	369	GLY	2.4
1	A	447	ALA	2.4
1	A	199	HIS	2.3
1	A	475	LEU	2.3
1	A	410	TYR	2.3
1	A	458	ASP	2.3
1	A	384	PHE	2.3
1	A	423	ARG	2.3
1	A	414	GLU	2.3
1	A	397	VAL	2.3
1	A	462	VAL	2.3
1	A	348	LEU	2.2
1	A	344	GLY	2.2
1	A	379	PHE	2.2
1	A	419	GLY	2.2
1	A	466	GLY	2.2
1	A	422	SER	2.2
1	A	330	PHE	2.2
1	A	394	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	467	ASN	2.1
1	A	478	ILE	2.1
1	A	436	ASP	2.1
1	A	445	ARG	2.1
1	A	342	LYS	2.1
1	A	407	LEU	2.1
1	A	357	THR	2.1
1	A	409	VAL	2.0
1	A	399	THR	2.0
1	A	424	SER	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	A	902	1/1	-	-	48,48,48,48	1
2	MG	B	901	1/1	-	-	44,44,44,44	1

6.5 Other polymers [i](#)

There are no such residues in this entry.