



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 06:50 pm GMT

PDB ID : 2F00
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

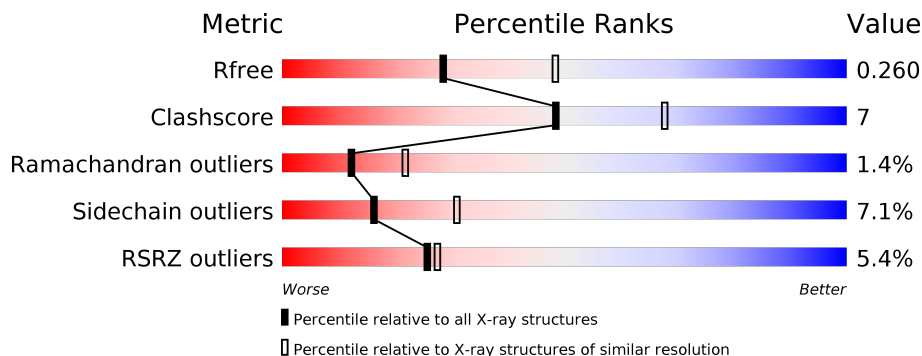
1 Overall quality at a glance

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}	111664	4155 (2.50-2.50)
Clashscore	122126	4827 (2.50-2.50)
Ramachandran outliers	120053	4735 (2.50-2.50)
Sidechain outliers	120020	4737 (2.50-2.50)
RSRZ outliers	108989	4058 (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 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	491	 9% 78% 17% ..
1	B	491	 % 83% 13% ..

2 Entry composition [i](#)

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
			Total	C	N	O	S	Se			
1	A	476	3524	2213	626	668	2	15	0	0	0
1	B	483	3678	2312	660	689	2	15	0	0	0

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 (three-letter code: MG) (formula: Mg).

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

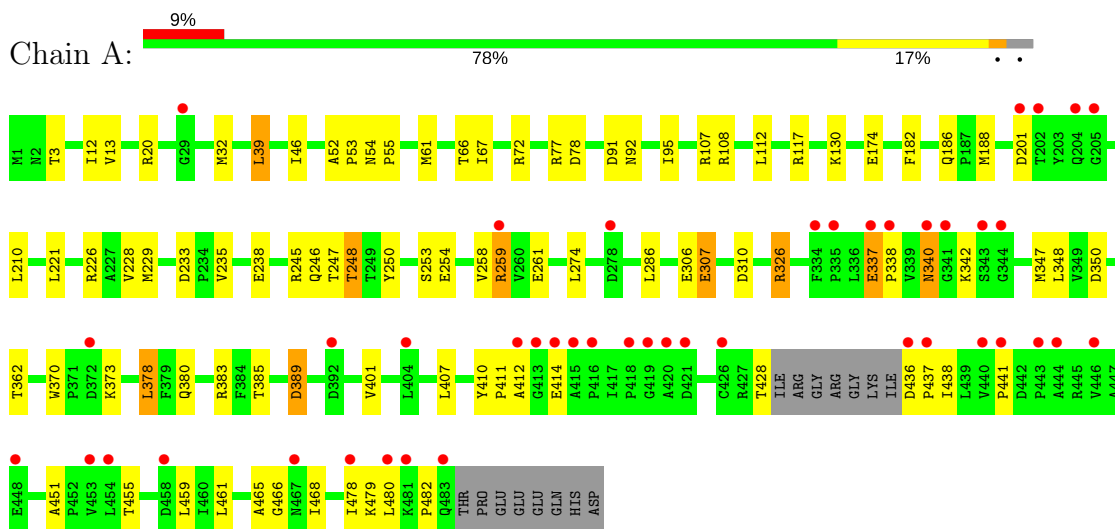
- Molecule 3 is water.

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

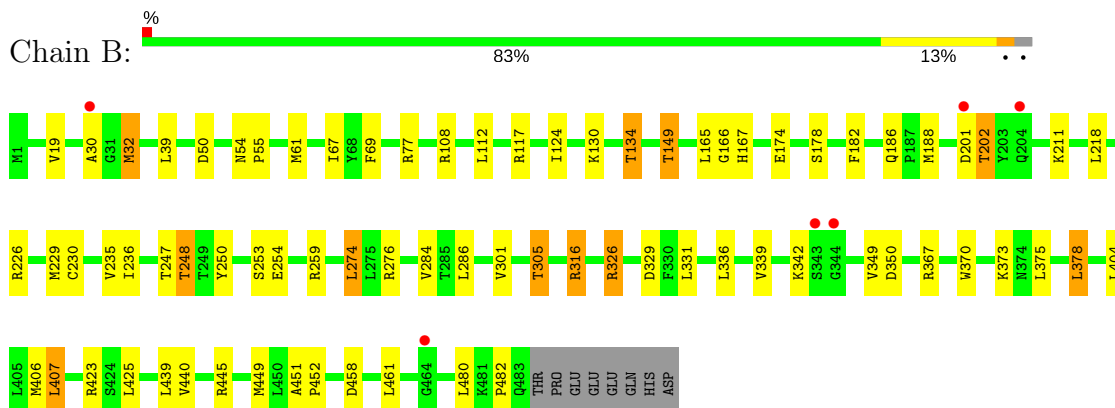
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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 29.68 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (88.00-2.50) 99.3 (29.68-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.51Å)	Xtrriage
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 (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtrriage
Anisotropy	0.372	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage'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	0.39	0/3580	0.54	1/4854 (0.0%)
1	B	0.39	0/3736	0.57	1/5053 (0.0%)
All	All	0.39	0/7316	0.55	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	5.61	110.03	103.30
1	B	482	PRO	N-CA-CB	5.16	109.49	103.30

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:HH11	1:A:259:ARG:HG2	1.10	1.15
1:A:259:ARG:NH1	1:A:259:ARG:HG2	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:CG	1:A:259:ARG:HH11	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:HG2	1:B:316:ARG:HH11	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:A:479:LYS:O	1:A:480:LEU:HB2	2.04	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:307:GLU:HG2	3:A:983:HOH:O	2.05	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:77:ARG:O	1:A:78:ASP:HB2	2.09	0.53
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:12:ILE:HG22	1:A:13:VAL:HG23	1.91	0.51
1:B:254:GLU:H	1:B:254:GLU:CD	2.14	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:N	1:A:389:ASP:OD1	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:N	1:A:91:ASP:OD1	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:316:ARG:CG	1:B:316:ARG:HH11	2.25	0.42
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: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%)	7	11
1	B	481/491 (98%)	466 (97%)	13 (3%)	2 (0%)	36	57
All	All	953/982 (97%)	911 (96%)	29 (3%)	13 (1%)	12	22

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%)	325 (93%)	26 (7%)	15	29
1	B	382/381 (100%)	356 (93%)	26 (7%)	17	33
All	All	733/762 (96%)	681 (93%)	52 (7%)	16	31

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	39	LEU
1	A	66	THR
1	A	72	ARG
1	A	107	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	428	THR
1	A	438	ILE
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	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

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	457	ASN
1	B	62	ASN
1	B	186	GLN
1	B	269	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 carbohydrates 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 [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	461/491 (93%)	0.42	44 (9%) 8 8	18, 42, 99, 100	0
1	B	468/491 (95%)	-0.00	6 (1%) 77 78	17, 34, 56, 64	0
All	All	929/982 (94%)	0.21	50 (5%) 26 27	17, 37, 83, 100	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	415	ALA	6.6
1	A	444	ALA	5.4
1	A	344	GLY	5.2
1	A	418	PRO	4.9
1	A	204	GLN	4.2
1	A	343	SER	3.9
1	A	338	PRO	3.8
1	A	440	VAL	3.8
1	A	413	GLY	3.7
1	B	30	ALA	3.7
1	A	334	PHE	3.7
1	A	467	ASN	3.6
1	A	202	THR	3.5
1	A	426	CYS	3.4
1	A	483	GLN	3.4
1	A	443	PRO	3.3
1	A	201	ASP	3.3
1	A	414	GLU	3.2
1	A	259	ARG	3.0
1	A	420	ALA	3.0
1	B	344	GLY	2.9
1	A	341	GLY	2.8
1	A	436	ASP	2.8
1	A	412	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	421	ASP	2.7
1	A	372	ASP	2.7
1	A	29	GLY	2.6
1	A	453	VAL	2.6
1	A	458	ASP	2.6
1	A	481	LYS	2.6
1	A	419	GLY	2.6
1	B	343	SER	2.6
1	A	446	VAL	2.5
1	A	480	LEU	2.5
1	A	478	ILE	2.5
1	A	335	PRO	2.5
1	A	278	ASP	2.5
1	A	340	ASN	2.4
1	A	337	GLU	2.4
1	A	437	PRO	2.3
1	A	416	PRO	2.2
1	A	448	GLU	2.2
1	A	205	GLY	2.2
1	A	404	LEU	2.2
1	B	204	GLN	2.1
1	A	441	PRO	2.1
1	A	392	ASP	2.1
1	B	201	ASP	2.1
1	B	464	GLY	2.1
1	A	454	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 carbohydrates 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(\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.