



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:02 PM UTC

PDB ID : 7B5T / pdb_00007b5t
Title : S. agalactiae BusR transcription factor
Authors : Bandera, A.M.; Witte, G.
Deposited on : 2020-12-07
Resolution : 2.80 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

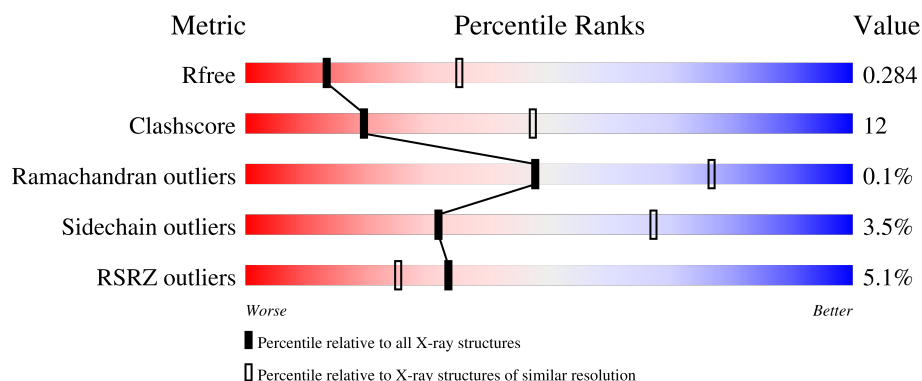
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	215	6% 73% 24% ..
1	B	215	5% 75% 19% ..
1	C	215	6% 69% 27% ..
1	D	215	3% 67% 25% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6528 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 GntR family transcriptional regulator.

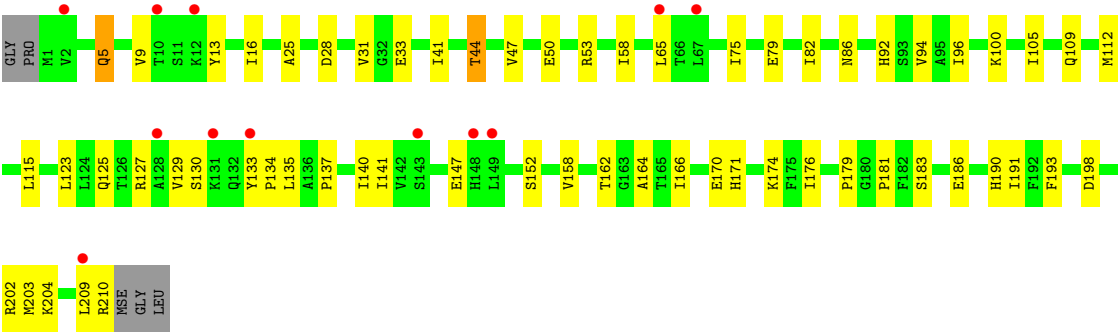
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	Se	0	0	0
			1656	1053	279	321	3			
1	B	207	Total	C	N	O	Se	0	0	0
			1628	1035	275	316	2			
1	D	200	Total	C	N	O	Se	0	0	0
			1573	1002	267	302	2			
1	C	210	Total	C	N	O	Se	0	0	0
			1649	1048	278	320	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP K0JNC6
A	0	PRO	-	expression tag	UNP K0JNC6
B	-1	GLY	-	expression tag	UNP K0JNC6
B	0	PRO	-	expression tag	UNP K0JNC6
D	-1	GLY	-	expression tag	UNP K0JNC6
D	0	PRO	-	expression tag	UNP K0JNC6
C	-1	GLY	-	expression tag	UNP K0JNC6
C	0	PRO	-	expression tag	UNP K0JNC6

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total	O	0	0
			6	6		
2	B	2	Total	O	0	0
			2	2		
2	D	5	Total	O	0	0
			5	5		
2	C	9	Total	O	0	0
			9	9		



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	114.18Å 114.18Å 240.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.22 – 2.80 48.22 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.1 (48.22-2.80) 99.1 (48.22-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.257 , 0.269 0.260 , 0.284	Depositor DCC
R_{free} test set	1996 reflections (5.30%)	wwPDB-VP
Wilson B-factor (Å ²)	104.6	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 73.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.030 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6528	wwPDB-VP
Average B, all atoms (Å ²)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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](#)

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	0.28	0/1681	0.57	0/2266
1	B	0.24	0/1653	0.56	0/2230
1	C	0.40	0/1674	0.69	1/2258 (0.0%)
1	D	0.28	0/1598	0.61	0/2155
All	All	0.31	0/6606	0.61	1/8909 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	5	GLN	N-CA-CB	-6.08	101.21	111.17

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	208	ASN	Mainchain

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	1656	0	1671	49	0
1	B	1628	0	1638	32	0
1	C	1649	0	1664	55	0
1	D	1573	0	1586	40	0
2	A	6	0	0	0	0
2	B	2	0	0	0	0
2	C	9	0	0	0	0
2	D	5	0	0	1	0
All	All	6528	0	6559	151	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 12.

All (151) 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:90:THR:HG23	1:A:92:HIS:H	1.37	0.88
1:A:147:GLU:HG2	1:A:148:HIS:H	1.42	0.85
1:C:204:LYS:HE2	1:C:209:LEU:HD23	1.65	0.77
1:C:13:TYR:HD1	1:C:58:ILE:HD11	1.49	0.77
1:C:198:ASP:O	1:C:202:ARG:HG3	1.87	0.73
1:C:50:GLU:OE2	1:C:53:ARG:NH1	2.22	0.73
1:D:89:GLU:OE1	1:C:129:VAL:HB	1.90	0.71
1:C:204:LYS:CE	1:C:209:LEU:HD23	2.20	0.71
1:D:23:ARG:NH2	1:C:137:PRO:O	2.25	0.70
1:C:170:GLU:HB3	1:C:190:HIS:HB2	1.73	0.69
1:C:13:TYR:CD1	1:C:58:ILE:HD11	2.27	0.68
1:A:105:ILE:HD13	1:B:116:ALA:HB2	1.78	0.66
1:D:85:LEU:HD22	1:C:130:SER:HB2	1.77	0.66
1:B:205:THR:HA	1:B:209:LEU:HB2	1.79	0.65
1:B:141:ILE:HG12	1:B:190:HIS:CD2	2.31	0.65
1:B:35:LEU:HD12	1:B:73:ALA:HB3	1.79	0.65
1:B:170:GLU:HB3	1:B:190:HIS:HB2	1.78	0.65
1:C:152:SER:HB2	1:C:181:PRO:O	1.97	0.65
1:D:148:HIS:HB3	1:D:156:LEU:HD11	1.79	0.63
1:C:204:LYS:CD	1:C:209:LEU:HD23	2.29	0.63
1:A:153:ILE:HD11	1:A:185:ILE:HD11	1.81	0.62
1:D:110:GLN:NE2	2:D:301:HOH:O	2.33	0.62
1:D:12:LYS:HD3	1:D:47:VAL:HG22	1.82	0.62
1:C:92:HIS:HD2	1:C:94:VAL:HG22	1.65	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:GLU:HG2	1:A:148:HIS:N	2.12	0.62
1:A:168:ALA:HB1	1:A:175:PHE:HE1	1.65	0.61
1:A:142:VAL:HA	1:A:208:ASN:OD1	2.01	0.61
1:A:90:THR:HG23	1:A:92:HIS:N	2.15	0.60
1:C:171:HIS:HE1	1:C:186:GLU:HB2	1.66	0.60
1:C:204:LYS:HE2	1:C:209:LEU:CD2	2.31	0.60
1:A:144:GLU:OE1	1:A:187:GLN:NE2	2.21	0.59
1:D:112:MSE:SE	1:C:112:MSE:HG3	2.52	0.59
1:D:158:VAL:O	1:D:162:THR:HG22	2.02	0.59
1:A:200:TYR:HB3	1:C:125:GLN:HE21	1.67	0.59
1:C:105:ILE:HG23	1:C:109:GLN:HE22	1.68	0.58
1:A:42:ALA:HB1	1:A:47:VAL:O	2.04	0.58
1:A:153:ILE:CD1	1:A:185:ILE:HD11	2.34	0.58
1:C:204:LYS:HD2	1:C:209:LEU:HD23	1.86	0.58
1:C:16:ILE:HD11	1:C:47:VAL:HG21	1.85	0.57
1:A:147:GLU:OE1	1:A:147:GLU:N	2.38	0.57
1:D:23:ARG:NH1	1:D:28:ASP:OD2	2.38	0.57
1:A:168:ALA:HB2	1:C:135:LEU:HD21	1.87	0.56
1:C:152:SER:HA	1:C:183:SER:O	2.05	0.56
1:D:16:ILE:HD11	1:D:47:VAL:HG21	1.86	0.56
1:D:34:LYS:HG2	1:D:35:LEU:O	2.05	0.56
1:A:148:HIS:HB2	1:A:151:LYS:HD2	1.88	0.56
1:D:187:GLN:HG2	1:D:187:GLN:O	2.06	0.55
1:D:197:GLU:HG3	1:C:28:ASP:OD1	2.06	0.55
1:C:171:HIS:CE1	1:C:186:GLU:HB2	2.41	0.55
1:C:140:ILE:HG23	1:C:193:PHE:HE1	1.72	0.54
1:B:195:GLY:HA3	1:B:199:VAL:HG11	1.89	0.54
1:A:16:ILE:O	1:A:20:VAL:HG12	2.07	0.54
1:A:37:SER:O	1:A:41:ILE:HG13	2.08	0.54
1:A:123:LEU:HD22	1:B:98:LYS:HD3	1.88	0.54
1:A:105:ILE:HG13	1:A:106:LYS:N	2.22	0.53
1:A:147:GLU:HB3	1:A:210:ARG:HG2	1.90	0.53
1:B:22:GLN:HA	1:B:85:LEU:HD11	1.89	0.53
1:A:6:SER:HA	1:D:173:GLY:O	2.08	0.53
1:A:147:GLU:HA	1:A:210:ARG:NH2	2.23	0.53
1:C:133:TYR:N	1:C:134:PRO:HD2	2.24	0.53
1:A:208:ASN:C	1:A:209:LEU:HD12	2.33	0.52
1:D:162:THR:HG21	1:D:203:MSE:SE	2.59	0.52
1:A:66:THR:HG23	1:A:68:LYS:H	1.74	0.52
1:B:196:ASP:O	1:B:199:VAL:HG22	2.09	0.52
1:A:208:ASN:O	1:A:209:LEU:HD12	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:50:GLU:OE2	1:D:54:LYS:NZ	2.43	0.52
1:A:54:LYS:HD3	1:D:182:PHE:O	2.10	0.51
1:C:31:VAL:O	1:C:75:ILE:HB	2.10	0.51
1:C:179:PRO:HB2	1:C:183:SER:OG	2.11	0.51
1:A:200:TYR:CD1	1:C:125:GLN:HG2	2.46	0.50
1:B:105:ILE:HG13	1:B:106:LYS:N	2.26	0.50
1:D:131:LYS:NZ	1:C:86:ASN:OD1	2.43	0.50
1:C:96:ILE:O	1:C:100:LYS:HG3	2.11	0.50
1:C:162:THR:HG23	1:C:164:ALA:H	1.76	0.50
1:D:92:HIS:CD2	1:D:94:VAL:HG12	2.47	0.49
1:A:69:HIS:HD2	1:A:74:ILE:HD12	1.76	0.49
1:D:158:VAL:HG21	1:D:166:ILE:CD1	2.41	0.49
1:C:50:GLU:OE2	1:C:53:ARG:CZ	2.60	0.49
1:D:92:HIS:HD2	1:D:94:VAL:HG12	1.76	0.49
1:B:16:ILE:O	1:B:20:VAL:HG12	2.12	0.49
1:A:142:VAL:HG12	1:A:189:ASP:O	2.12	0.49
1:B:40:THR:O	1:B:44:THR:HG23	2.13	0.49
1:D:158:VAL:HG21	1:D:166:ILE:HD11	1.94	0.49
1:A:185:ILE:HD12	1:A:185:ILE:N	2.28	0.48
1:B:132:GLN:NE2	1:C:9:VAL:HG12	2.28	0.48
1:A:90:THR:HG21	1:B:201:ALA:HB1	1.96	0.48
1:A:147:GLU:OE2	1:A:206:TYR:O	2.32	0.48
1:B:101:ILE:O	1:B:105:ILE:HG23	2.14	0.48
1:A:90:THR:HG21	1:B:201:ALA:CB	2.44	0.47
1:D:114:GLU:O	1:D:118:LEU:HD13	2.14	0.47
1:B:25:ALA:O	1:B:78:LYS:HE3	2.14	0.47
1:C:140:ILE:HG13	1:C:191:ILE:HB	1.97	0.47
1:A:144:GLU:HA	1:A:187:GLN:NE2	2.29	0.46
1:A:101:ILE:HG23	1:D:115:LEU:HD11	1.97	0.46
1:A:168:ALA:HB3	1:A:192:PHE:HB2	1.96	0.46
1:D:133:TYR:CD1	1:D:134:PRO:HD3	2.50	0.46
1:D:157:ASN:HB3	1:D:160:HIS:HB3	1.97	0.46
1:B:101:ILE:HG23	1:C:115:LEU:HD11	1.97	0.46
1:A:102:ARG:HA	1:A:105:ILE:HG12	1.98	0.46
1:A:195:GLY:HA3	1:A:199:VAL:HG11	1.98	0.46
1:D:96:ILE:O	1:D:100:LYS:HG3	2.16	0.46
1:C:65:LEU:HD23	1:C:75:ILE:HA	1.97	0.46
1:C:133:TYR:H	1:C:134:PRO:HD2	1.81	0.46
1:C:33:GLU:O	1:C:75:ILE:HG12	2.16	0.45
1:D:133:TYR:N	1:D:134:PRO:HD2	2.30	0.45
1:D:147:GLU:HB3	1:D:148:HIS:H	1.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:LYS:HB3	1:B:40:THR:OG1	2.17	0.45
1:A:79:GLU:O	1:A:82:ILE:HG12	2.17	0.45
1:A:123:LEU:CD2	1:B:98:LYS:HD3	2.45	0.45
1:A:149:LEU:H	1:A:149:LEU:HD23	1.81	0.45
1:C:65:LEU:HD21	1:C:75:ILE:HD13	1.99	0.45
1:C:158:VAL:O	1:C:162:THR:HG22	2.16	0.44
1:C:162:THR:HG21	1:C:203:MSE:SE	2.67	0.44
1:D:140:ILE:HD11	1:D:209:LEU:HD12	1.99	0.44
1:C:204:LYS:HZ3	1:C:210:ARG:HD2	1.82	0.44
1:B:130:SER:HB2	1:D:127:ARG:NH1	2.32	0.44
1:D:47:VAL:HG12	1:D:51:THR:OG1	2.18	0.44
1:C:158:VAL:HG11	1:C:166:ILE:CD1	2.48	0.44
1:D:20:VAL:HG13	1:D:59:LEU:HD21	2.00	0.43
1:B:152:SER:HB2	1:B:181:PRO:O	2.19	0.43
1:B:12:LYS:HD3	1:B:12:LYS:HA	1.87	0.43
1:C:92:HIS:O	1:C:96:ILE:HG12	2.18	0.43
1:D:140:ILE:CG2	1:D:191:ILE:HB	2.48	0.43
1:B:148:HIS:HA	1:B:151:LYS:NZ	2.33	0.43
1:C:25:ALA:HB1	1:C:82:ILE:HG12	2.01	0.43
1:A:128:ALA:O	1:A:132:GLN:HG2	2.19	0.43
1:B:10:THR:OG1	1:B:11:SER:N	2.51	0.43
1:B:59:LEU:HD22	1:B:64:ILE:HD12	2.01	0.42
1:D:198:ASP:O	1:D:202:ARG:HG3	2.19	0.42
1:B:9:VAL:HG12	1:B:10:THR:O	2.19	0.42
1:D:93:SER:HA	1:D:96:ILE:HG12	2.00	0.42
1:D:140:ILE:HD11	1:D:204:LYS:HA	2.00	0.42
1:D:65:LEU:HA	1:D:76:LEU:HG	2.02	0.42
1:A:138:TYR:HE2	1:C:127:ARG:HH21	1.67	0.42
1:A:205:THR:HB	1:B:87:GLN:HE22	1.83	0.42
1:B:210:ARG:HD3	1:B:210:ARG:HA	1.95	0.42
1:B:209:LEU:HA	1:B:209:LEU:HD23	1.74	0.42
1:C:133:TYR:N	1:C:134:PRO:CD	2.83	0.42
1:D:38:ARG:HD2	1:D:53:ARG:HB2	2.00	0.41
1:A:101:ILE:O	1:A:105:ILE:HG23	2.20	0.41
1:A:143:SER:OG	1:A:144:GLU:N	2.53	0.41
1:C:133:TYR:CD2	1:C:134:PRO:HD3	2.56	0.41
1:C:162:THR:HG23	1:C:164:ALA:N	2.36	0.41
1:C:174:LYS:O	1:C:176:ILE:HD12	2.20	0.41
1:A:112:MSE:SE	1:B:108:GLN:HB3	2.70	0.41
1:B:209:LEU:HB3	1:B:210:ARG:H	1.55	0.41
1:A:130:SER:HA	1:C:127:ARG:NH1	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:ILE:HA	1:C:44:THR:OG1	2.21	0.41
1:A:126:THR:HA	1:A:129:VAL:HG22	2.03	0.41
1:C:5:GLN:HG3	1:C:5:GLN:O	2.21	0.40
1:D:98:LYS:HG3	1:C:123:LEU:HD13	2.02	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	209/215 (97%)	196 (94%)	13 (6%)	0	100	100
1	B	205/215 (95%)	195 (95%)	9 (4%)	1 (0%)	24	55
1	C	208/215 (97%)	200 (96%)	8 (4%)	0	100	100
1	D	198/215 (92%)	189 (96%)	9 (4%)	0	100	100
All	All	820/860 (95%)	780 (95%)	39 (5%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	209	LEU

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	181/179 (101%)	175 (97%)	6 (3%)	33	69
1	B	177/179 (99%)	168 (95%)	9 (5%)	21	54
1	C	180/179 (101%)	176 (98%)	4 (2%)	45	78
1	D	170/179 (95%)	164 (96%)	6 (4%)	32	67
All	All	708/716 (99%)	683 (96%)	25 (4%)	32	67

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	66	THR
1	A	146	SER
1	A	148	HIS
1	A	152	SER
1	A	184	VAL
1	B	40	THR
1	B	67	LEU
1	B	105	ILE
1	B	108	GLN
1	B	142	VAL
1	B	144	GLU
1	B	184	VAL
1	B	209	LEU
1	B	210	ARG
1	D	75	ILE
1	D	131	LYS
1	D	145	ASP
1	D	155	GLU
1	D	156	LEU
1	D	194	VAL
1	C	44	THR
1	C	79	GLU
1	C	141	ILE
1	C	147	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	69	HIS

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Mol	Chain	Res	Type
1	A	87	GLN
1	A	110	GLN
1	A	161	GLN
1	A	190	HIS
1	B	148	HIS
1	B	157	ASN
1	B	160	HIS
1	B	171	HIS
1	D	14	GLN
1	D	110	GLN
1	D	160	HIS
1	C	14	GLN
1	C	87	GLN
1	C	104	ASN
1	C	109	GLN
1	C	125	GLN
1	C	148	HIS
1	C	187	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](#)

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

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	208/215 (96%)	0.59	13 (6%) 26 19	92, 123, 169, 226	0
1	B	205/215 (95%)	0.53	10 (4%) 35 27	73, 126, 181, 204	0
1	C	207/215 (96%)	0.62	12 (5%) 29 22	78, 118, 174, 201	0
1	D	198/215 (92%)	0.56	7 (3%) 47 38	81, 123, 184, 235	0
All	All	818/860 (95%)	0.58	42 (5%) 33 25	73, 123, 181, 235	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	131	LYS	4.1
1	C	2	VAL	4.1
1	A	32	GLY	3.7
1	C	143	SER	3.4
1	D	139	GLU	3.3
1	A	45	PHE	3.2
1	B	128	ALA	3.1
1	A	207	PHE	3.0
1	C	209	LEU	3.0
1	A	6	SER	3.0
1	D	16	ILE	2.9
1	B	129	VAL	2.9
1	B	8	ILE	2.8
1	A	31	VAL	2.6
1	B	56	LEU	2.6
1	B	153	ILE	2.6
1	A	0	PRO	2.5
1	C	67	LEU	2.5
1	A	89	GLU	2.5
1	C	149	LEU	2.4
1	D	31	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	11	SER	2.3
1	A	206	TYR	2.3
1	C	65	LEU	2.3
1	C	148	HIS	2.3
1	C	133	TYR	2.3
1	B	4	GLU	2.3
1	A	161	GLN	2.3
1	C	128	ALA	2.2
1	B	133	TYR	2.2
1	B	146	SER	2.2
1	C	10	THR	2.2
1	A	50	GLU	2.2
1	D	32	GLY	2.1
1	B	182	PHE	2.1
1	D	202	ARG	2.1
1	A	148	HIS	2.1
1	A	38	ARG	2.1
1	B	52	ALA	2.1
1	C	12	LYS	2.0
1	A	91	SER	2.0
1	D	77	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.