



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 02:15 AM UTC

PDB ID : 4EGQ / pdb_00004egq
Title : Crystal structure of D-alanine-D-alanine ligase B from Burkholderia pseudo-mallei
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2012-03-31
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

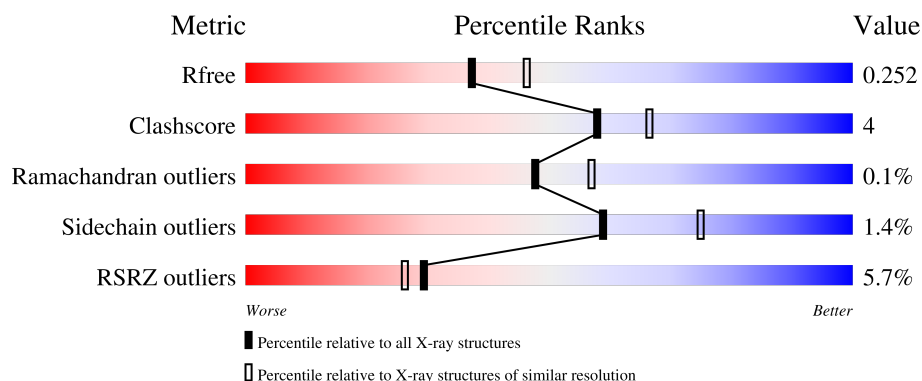
1 Overall quality at a glance

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	316	7% 83% 7% 9%
1	B	316	3% 88% 8%
1	C	316	4% 87% 8%
1	D	316	7% 83% 7% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8592 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 D-alanine–D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2030	1299	353	372	6			
1	B	302	Total	C	N	O	S	0	0	0
			2158	1382	367	403	6			
1	C	302	Total	C	N	O	S	0	0	0
			2168	1395	369	398	6			
1	D	288	Total	C	N	O	S	0	0	0
			2043	1316	353	368	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q3JNE1
A	-2	PRO	-	expression tag	UNP Q3JNE1
A	-1	GLY	-	expression tag	UNP Q3JNE1
A	0	SER	-	expression tag	UNP Q3JNE1
B	-3	GLY	-	expression tag	UNP Q3JNE1
B	-2	PRO	-	expression tag	UNP Q3JNE1
B	-1	GLY	-	expression tag	UNP Q3JNE1
B	0	SER	-	expression tag	UNP Q3JNE1
C	-3	GLY	-	expression tag	UNP Q3JNE1
C	-2	PRO	-	expression tag	UNP Q3JNE1
C	-1	GLY	-	expression tag	UNP Q3JNE1
C	0	SER	-	expression tag	UNP Q3JNE1
D	-3	GLY	-	expression tag	UNP Q3JNE1
D	-2	PRO	-	expression tag	UNP Q3JNE1
D	-1	GLY	-	expression tag	UNP Q3JNE1
D	0	SER	-	expression tag	UNP Q3JNE1

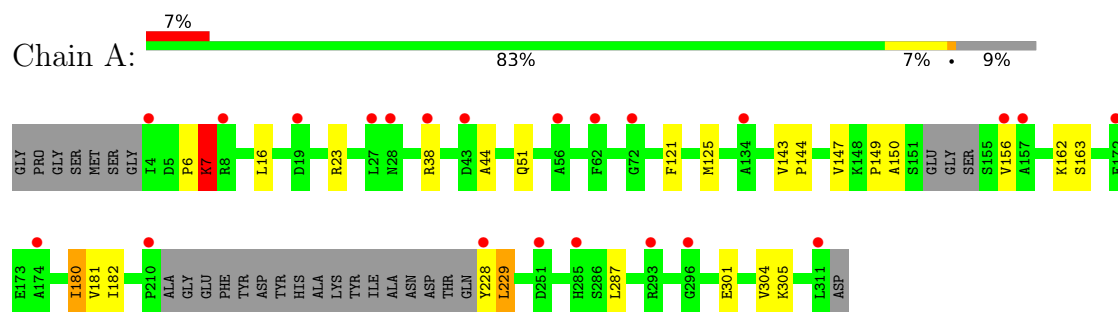
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total 19	O 19	0	0
2	B	60	Total 60	O 60	0	0
2	C	61	Total 61	O 61	0	0
2	D	53	Total 53	O 53	0	0

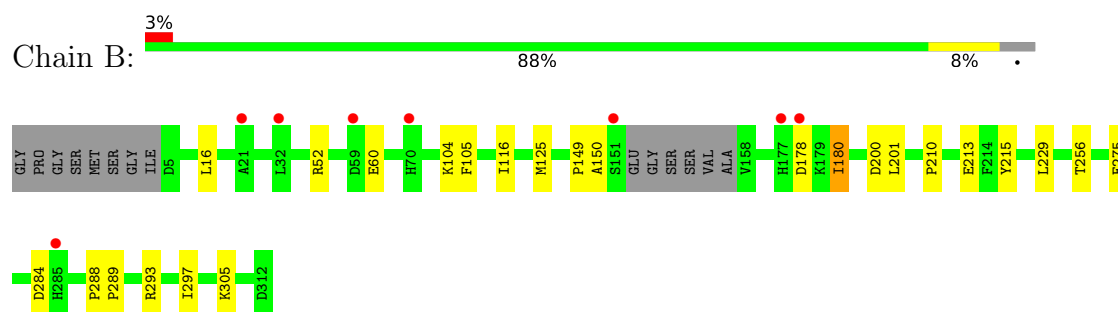
3 Residue-property plots [i](#)

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.

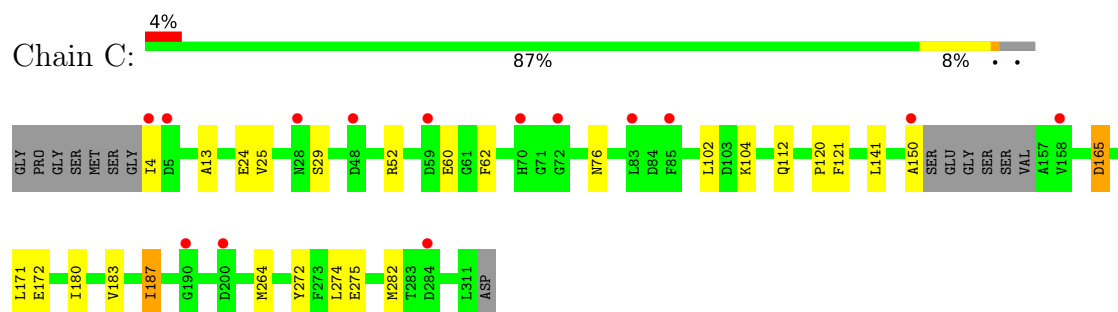
• Molecule 1: D-alanine–D-alanine ligase



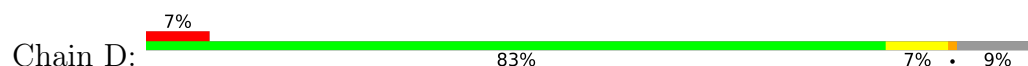
• Molecule 1: D-alanine–D-alanine ligase

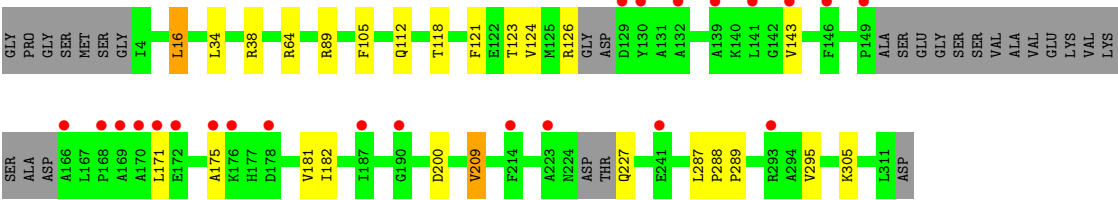


• Molecule 1: D-alanine–D-alanine ligase



• Molecule 1: D-alanine–D-alanine ligase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.39Å 66.09Å 98.56Å 84.47° 80.50° 83.59°	Depositor
Resolution (Å)	35.27 – 2.20 35.27 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.0 (35.27-2.20) 97.0 (35.27-2.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.210 , 0.250 0.213 , 0.252	Depositor DCC
R_{free} test set	2772 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.005	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8592	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.67% 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

5.1 Standard geometry

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	0.78	0/2070	0.85	3/2828 (0.1%)
1	B	0.84	0/2203	0.87	1/3008 (0.0%)
1	C	0.90	0/2216	0.90	1/3025 (0.0%)
1	D	0.81	0/2085	0.92	7/2845 (0.2%)
All	All	0.84	0/8574	0.89	12/11706 (0.1%)

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	A	0	1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	64	ARG	N-CA-CB	-6.96	100.04	111.66
1	C	172	GLU	N-CA-C	6.78	119.25	111.11
1	D	64	ARG	CA-CB-CG	6.20	126.49	114.10
1	D	287	LEU	CA-C-N	-6.05	114.14	120.38
1	D	287	LEU	C-N-CA	-6.05	114.14	120.38
1	B	116	ILE	N-CA-C	-6.02	103.45	108.63
1	A	287	LEU	CA-C-N	-5.44	114.78	120.38
1	A	287	LEU	C-N-CA	-5.44	114.78	120.38
1	D	118	THR	CA-C-N	-5.21	115.02	120.38
1	D	118	THR	C-N-CA	-5.21	115.02	120.38
1	D	143	VAL	N-CA-C	5.04	119.76	108.88
1	A	229	LEU	N-CA-C	5.01	117.14	109.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	7	LYS	Peptide

5.2 Too-close contacts

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	2030	0	1935	19	0
1	B	2158	0	2048	14	1
1	C	2168	0	2079	17	1
1	D	2043	0	1953	23	0
2	A	19	0	0	0	0
2	B	60	0	0	1	0
2	C	61	0	0	0	0
2	D	53	0	0	1	0
All	All	8592	0	8015	73	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 4.

All (73) 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:D:124:VAL:HG22	1:D:181:VAL:HG13	1.33	1.08
1:D:124:VAL:CG2	1:D:181:VAL:CG1	2.32	1.07
1:D:124:VAL:HG23	1:D:181:VAL:CG1	2.00	0.92
1:D:124:VAL:HG22	1:D:181:VAL:CG1	1.99	0.90
1:D:124:VAL:HG21	1:D:171:LEU:HD22	1.56	0.87
1:A:143:VAL:HG13	1:A:144:PRO:HA	1.59	0.84
1:D:124:VAL:HG23	1:D:181:VAL:HG12	1.58	0.82
1:B:150:ALA:HB3	1:B:180:ILE:HG23	1.62	0.79
1:A:125:MET:HG2	1:A:180:ILE:HD12	1.66	0.76
1:A:6:PRO:O	1:A:7:LYS:CB	2.35	0.73
1:D:295:VAL:HG12	1:D:295:VAL:O	1.87	0.73
1:C:52:ARG:NH2	1:C:60:GLU:OE2	2.24	0.70
1:D:124:VAL:CG2	1:D:181:VAL:HG13	2.02	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ALA:HB3	1:A:180:ILE:HG23	1.74	0.69
1:A:149:PRO:HG2	1:A:156:VAL:HG11	1.73	0.69
1:C:4:ILE:O	1:C:4:ILE:HG23	1.94	0.68
1:A:143:VAL:CG1	1:A:144:PRO:HA	2.26	0.64
1:C:24:GLU:OE2	1:C:25:VAL:HG23	1.98	0.63
1:A:150:ALA:HA	1:A:182:ILE:HD12	1.80	0.62
1:D:209:VAL:HG12	1:D:227:GLN:HA	1.83	0.61
1:A:301:GLU:O	1:A:304:VAL:HG12	2.02	0.60
1:C:120:PRO:HB2	1:C:141:LEU:HD22	1.87	0.57
1:B:149:PRO:HB3	1:B:178:ASP:HB2	1.87	0.56
1:D:200:ASP:OD1	1:D:305:LYS:HE2	2.07	0.53
1:B:200:ASP:OD1	1:B:305:LYS:HE2	2.09	0.52
1:C:112:GLN:HG3	1:C:121:PHE:CE2	2.44	0.52
1:C:264:MET:HE3	1:C:274:LEU:HD11	1.91	0.51
1:B:104:LYS:HE3	1:B:275:GLU:OE1	2.10	0.51
1:D:112:GLN:OE1	1:D:121:PHE:CZ	2.64	0.51
1:A:143:VAL:HG11	1:A:162:LYS:O	2.10	0.51
1:D:126:ARG:NH1	1:D:175:ALA:O	2.44	0.51
1:D:295:VAL:O	1:D:295:VAL:CG1	2.58	0.50
1:D:105:PHE:CD1	1:D:182:ILE:HD11	2.46	0.49
1:A:149:PRO:CG	1:A:156:VAL:HG11	2.41	0.48
1:C:171:LEU:HD21	1:C:183:VAL:HG21	1.95	0.48
1:B:215:TYR:CE1	1:B:229:LEU:HD11	2.49	0.48
1:C:150:ALA:CB	1:C:180:ILE:HG23	2.44	0.47
1:C:264:MET:HE3	1:C:274:LEU:CD1	2.44	0.47
1:C:104:LYS:HE3	1:C:275:GLU:OE1	2.15	0.47
1:D:124:VAL:CG2	1:D:181:VAL:HG11	2.35	0.46
1:B:210:PRO:HD2	1:B:213:GLU:OE2	2.16	0.46
1:A:147:VAL:HG13	1:A:181:VAL:CG2	2.46	0.45
1:C:171:LEU:HD21	1:C:183:VAL:CG2	2.46	0.45
1:B:125:MET:HG2	1:B:180:ILE:HD12	1.98	0.45
1:A:304:VAL:HG13	1:A:305:LYS:N	2.32	0.44
1:B:105:PHE:CZ	1:B:180:ILE:HG12	2.53	0.44
1:A:23:ARG:HH22	1:A:51:GLN:HE22	1.65	0.44
1:C:150:ALA:HB3	1:C:180:ILE:HG23	2.00	0.44
1:D:34:LEU:O	1:D:38:ARG:HG2	2.18	0.44
1:D:124:VAL:CG2	1:D:181:VAL:HG12	2.23	0.44
1:A:147:VAL:CG1	1:A:181:VAL:HG21	2.49	0.43
1:D:123:THR:HG23	2:D:444:HOH:O	2.18	0.43
1:D:16:LEU:C	1:D:16:LEU:CD1	2.92	0.43
1:D:288:PRO:HB2	1:D:289:PRO:HD3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:VAL:HG13	1:A:181:VAL:HG21	2.01	0.43
1:B:215:TYR:CD1	1:B:229:LEU:HD11	2.54	0.43
1:B:125:MET:HE2	1:B:125:MET:HB3	1.91	0.43
1:C:13:ALA:HB2	1:C:62:PHE:CD1	2.54	0.43
1:C:187:ILE:HD12	1:C:272:TYR:CE2	2.54	0.42
1:C:4:ILE:O	1:C:4:ILE:CG2	2.63	0.42
1:D:112:GLN:OE1	1:D:121:PHE:CE2	2.72	0.42
1:A:228:TYR:CB	1:A:229:LEU:HD12	2.50	0.42
1:A:150:ALA:HA	1:A:182:ILE:CD1	2.49	0.41
1:C:76:ASN:CB	1:C:102:LEU:HD12	2.51	0.41
1:B:256:THR:HG23	2:B:448:HOH:O	2.20	0.41
1:D:16:LEU:HD12	1:D:16:LEU:O	2.20	0.41
1:B:288:PRO:HB2	1:B:289:PRO:HD3	2.03	0.41
1:D:121:PHE:CD1	1:D:121:PHE:C	2.99	0.41
1:A:38:ARG:NH2	1:A:44:ALA:O	2.54	0.40
1:C:29:SER:OG	1:C:282:MET:HB2	2.20	0.40
1:B:201:LEU:HD22	1:B:297:ILE:HD13	2.03	0.40
1:B:284:ASP:O	1:B:293:ARG:NH2	2.45	0.40
1:A:121:PHE:CD1	1:A:121:PHE:C	3.00	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)
1:B:60:GLU:OE2	1:C:165:ASP:OD1[1_445]	2.09	0.11

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	282/316 (89%)	280 (99%)	1 (0%)	1 (0%)	30 34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	298/316 (94%)	296 (99%)	2 (1%)	0	100	100
1	C	298/316 (94%)	296 (99%)	2 (1%)	0	100	100
1	D	280/316 (89%)	276 (99%)	4 (1%)	0	100	100
All	All	1158/1264 (92%)	1148 (99%)	9 (1%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	LYS

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	184/240 (77%)	181 (98%)	3 (2%)	55	71
1	B	197/240 (82%)	194 (98%)	3 (2%)	57	73
1	C	200/240 (83%)	198 (99%)	2 (1%)	68	81
1	D	183/240 (76%)	180 (98%)	3 (2%)	55	71
All	All	764/960 (80%)	753 (99%)	11 (1%)	59	75

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	163	SER
1	A	180	ILE
1	B	16	LEU
1	B	52	ARG
1	B	180	ILE
1	C	165	ASP
1	C	187	ILE
1	D	16	LEU
1	D	89	ARG

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Mol	Chain	Res	Type
1	D	209	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	270	ASN
1	B	270	ASN
1	C	270	ASN
1	D	270	ASN

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 [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	288/316 (91%)	0.71	22 (7%)	20 17	22, 44, 71, 93	0
1	B	302/316 (95%)	0.54	8 (2%)	57 54	21, 36, 54, 66	0
1	C	302/316 (95%)	0.35	14 (4%)	37 34	20, 33, 50, 67	0
1	D	288/316 (91%)	0.60	23 (7%)	18 16	22, 36, 64, 77	0
All	All	1180/1264 (93%)	0.55	67 (5%)	29 26	20, 37, 63, 93	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	141	LEU	3.8
1	B	285	HIS	3.7
1	A	210	PRO	3.7
1	D	149	PRO	3.6
1	D	172	GLU	3.6
1	D	169	ALA	3.5
1	C	284	ASP	3.4
1	C	5	ASP	3.2
1	C	4	ILE	3.2
1	D	178	ASP	3.1
1	B	21	ALA	3.1
1	B	70	HIS	3.1
1	A	19	ASP	3.1
1	D	129	ASP	3.0
1	B	59	ASP	2.9
1	A	56	ALA	2.9
1	D	170	ALA	2.9
1	D	187	ILE	2.9
1	A	8	ARG	2.9
1	A	27	LEU	2.9
1	D	143	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	130	TYR	2.8
1	D	146	PHE	2.8
1	B	178	ASP	2.7
1	D	214	PHE	2.7
1	B	177	HIS	2.7
1	C	28	ASN	2.7
1	D	241	GLU	2.7
1	A	251	ASP	2.7
1	D	176	LYS	2.6
1	B	151	SER	2.6
1	A	296	GLY	2.6
1	A	134	ALA	2.6
1	C	190	GLY	2.5
1	D	190	GLY	2.5
1	A	157	ALA	2.5
1	C	70	HIS	2.5
1	A	72	GLY	2.5
1	A	311	LEU	2.4
1	A	156	VAL	2.4
1	A	4	ILE	2.4
1	C	85	PHE	2.4
1	A	43	ASP	2.3
1	A	228	TYR	2.3
1	D	171	LEU	2.3
1	D	166	ALA	2.3
1	A	28	ASN	2.2
1	C	150	ALA	2.2
1	D	175	ALA	2.2
1	A	293	ARG	2.2
1	C	83	LEU	2.2
1	C	200	ASP	2.1
1	A	172	GLU	2.1
1	C	72	GLY	2.1
1	D	223	ALA	2.1
1	C	59	ASP	2.1
1	A	38	ARG	2.1
1	A	174	ALA	2.1
1	D	139	ALA	2.1
1	C	48	ASP	2.1
1	C	158	VAL	2.1
1	A	62	PHE	2.1
1	A	285	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	293	ARG	2.0
1	D	132	ALA	2.0
1	B	32	LEU	2.0
1	D	168	PRO	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.