



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:39 AM UTC

PDB ID : 4W4R / pdb_00004w4r
Title : Crystal structure of ent-kaurene synthase BJKS from bradyrhizobium japonicum
Authors : Liu, W.; Zheng, Y.; Huang, C.H.; Ko, T.P.; Guo, R.T.
Deposited on : 2014-08-15
Resolution : 1.92 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

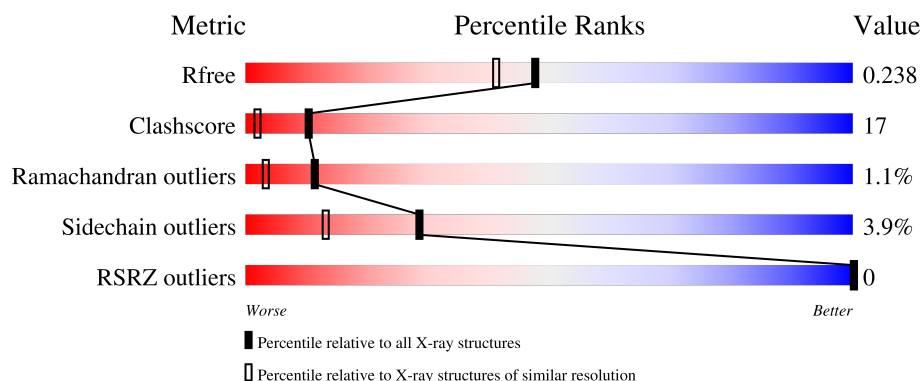
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	300	 67% 20% • 10%
1	B	300	 68% 20% •• 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4934 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 Uncharacterized protein blr2150.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2130	1349	385	384	12			
1	B	276	Total	C	N	O	S	0	0	0
			2180	1376	397	396	11			

- Molecule 2 is water.

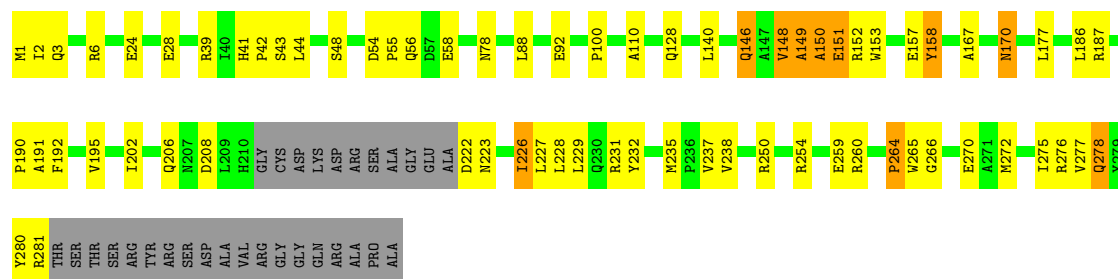
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	314	Total	O	0	0
			314	314		
2	B	310	Total	O	0	0
			310	310		

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.

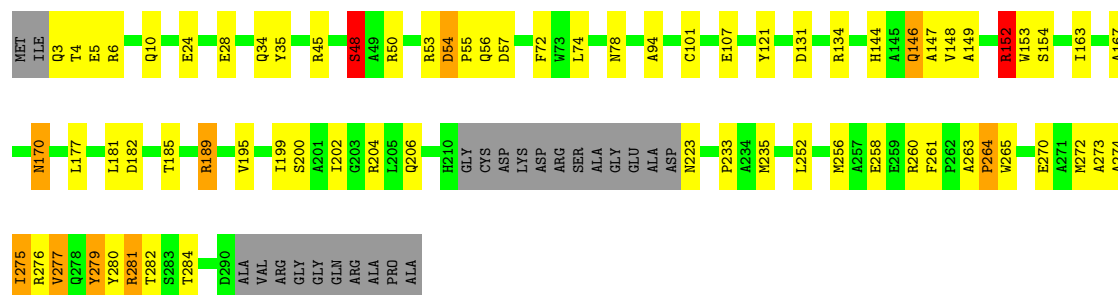
- Molecule 1: Uncharacterized protein blr2150

Chain A: 



- Molecule 1: Uncharacterized protein blr2150

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	65.80Å 65.80Å 135.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.92 25.00 – 1.92	Depositor EDS
% Data completeness (in resolution range)	94.8 (25.00-1.92) 97.6 (25.00-1.92)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.09 (at 1.92Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.191 , 0.232 0.192 , 0.238	Depositor DCC
R_{free} test set	2138 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.092 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4934	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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	0.58	0/2175	0.97	6/2956 (0.2%)
1	B	0.51	0/2226	0.98	11/3025 (0.4%)
All	All	0.55	0/4401	0.98	17/5981 (0.3%)

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 (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	281	ARG	N-CA-C	9.97	124.51	111.75
1	A	48	SER	N-CA-C	7.58	119.54	111.28
1	B	265	TRP	N-CA-C	7.43	119.38	111.28
1	B	152	ARG	N-CA-C	7.26	119.90	110.53
1	B	48	SER	N-CA-C	6.23	118.59	111.11
1	B	54	ASP	CA-C-N	6.22	126.42	119.32
1	B	54	ASP	C-N-CA	6.22	126.42	119.32
1	A	167	ALA	N-CA-C	5.86	120.01	112.86
1	B	279	TYR	N-CA-C	-5.85	106.19	113.38
1	A	265	TRP	N-CA-C	5.81	117.61	111.28
1	B	167	ALA	N-CA-C	5.71	119.83	112.86
1	B	5	GLU	N-CA-C	5.71	118.28	111.71
1	B	147	ALA	N-CA-C	-5.40	106.66	113.20
1	A	150	ALA	N-CA-C	5.37	118.59	108.65
1	B	154	SER	N-CA-C	-5.30	102.68	110.52
1	A	151	GLU	N-CA-C	-5.28	98.60	107.32
1	A	146	GLN	N-CA-C	5.21	117.38	111.02

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	158	TYR	Sidechain

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	2130	0	2108	71	0
1	B	2180	0	2149	78	0
2	A	314	0	0	7	0
2	B	310	0	0	15	0
All	All	4934	0	4257	146	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 17.

All (146) 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:229:LEU:HG	1:A:237:VAL:HG11	1.42	1.02
1:B:277:VAL:HG23	1:B:280:TYR:CE1	2.07	0.88
1:B:152:ARG:HH11	1:B:152:ARG:HG3	1.40	0.85
1:B:279:TYR:HB3	2:B:1005:HOH:O	1.76	0.84
1:A:148:VAL:HG12	1:A:149:ALA:N	1.96	0.80
1:A:266:GLY:O	1:A:270:GLU:HG3	1.83	0.79
1:B:199:ILE:HD11	1:B:273:ALA:HA	1.65	0.79
1:B:263:ALA:HA	2:B:730:HOH:O	1.82	0.79
1:A:223:ASN:HB3	1:A:226:ILE:CG2	2.13	0.78
1:A:190:PRO:HD2	2:A:551:HOH:O	1.83	0.78
1:B:280:TYR:HD1	1:B:281:ARG:H	1.29	0.76
1:A:223:ASN:HB3	1:A:226:ILE:HG22	1.67	0.76
1:A:3:GLN:H	1:A:3:GLN:CD	1.93	0.76
1:B:170:ASN:HD22	1:B:170:ASN:H	1.34	0.76
1:B:276:ARG:HB2	1:B:279:TYR:CZ	2.21	0.75
1:B:280:TYR:HD1	1:B:281:ARG:N	1.85	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:GLU:HG3	2:B:982:HOH:O	1.86	0.74
1:A:28:GLU:OE2	1:A:281:ARG:HD3	1.88	0.73
1:B:195:VAL:HG11	1:B:256:MET:CE	2.19	0.72
1:A:237:VAL:HG13	1:A:238:VAL:N	2.05	0.71
1:B:199:ILE:HD11	1:B:273:ALA:CA	2.20	0.71
1:B:280:TYR:CD1	1:B:281:ARG:N	2.58	0.71
1:A:152:ARG:HB2	1:A:231:ARG:HD3	1.73	0.70
1:A:237:VAL:HG13	1:A:238:VAL:H	1.54	0.70
1:B:72:PHE:CE1	1:B:281:ARG:HB2	2.27	0.70
1:B:260:ARG:HH11	1:B:260:ARG:HG2	1.57	0.69
1:A:43:SER:HB3	1:A:264:PRO:HB3	1.72	0.69
1:A:24:GLU:O	1:A:28:GLU:HG3	1.92	0.69
1:B:281:ARG:HD2	2:B:796:HOH:O	1.95	0.67
1:A:229:LEU:CG	1:A:237:VAL:HG11	2.24	0.66
1:B:277:VAL:HG21	1:B:282:THR:HB	1.78	0.65
1:B:202:ILE:O	1:B:206:GLN:HG3	1.97	0.65
1:A:2:ILE:HD12	1:A:58:GLU:HG3	1.79	0.65
1:A:158:TYR:OH	1:A:208:ASP:OD2	2.07	0.65
1:B:279:TYR:HB2	2:B:1010:HOH:O	1.96	0.65
1:B:24:GLU:O	1:B:28:GLU:HG3	1.97	0.64
1:A:3:GLN:CD	1:A:3:GLN:N	2.56	0.63
1:A:148:VAL:O	1:A:150:ALA:N	2.28	0.63
1:B:274:ALA:O	1:B:275:ILE:C	2.42	0.62
1:B:281:ARG:HH11	1:B:281:ARG:HG2	1.64	0.61
1:B:256:MET:HA	1:B:256:MET:HE2	1.81	0.60
1:B:277:VAL:HG22	1:B:282:THR:HG22	1.84	0.60
1:B:277:VAL:CG2	1:B:282:THR:HB	2.31	0.60
1:A:146:GLN:OE1	1:A:152:ARG:HA	2.02	0.60
1:B:233:PRO:HD3	2:B:701:HOH:O	2.01	0.59
1:B:3:GLN:HG3	2:B:726:HOH:O	2.02	0.59
1:A:187:ARG:NH2	2:A:654:HOH:O	2.14	0.59
1:B:204:ARG:O	1:B:204:ARG:HD3	2.03	0.58
1:A:170:ASN:HD22	1:A:170:ASN:H	1.50	0.57
1:B:206:GLN:HG2	1:B:276:ARG:NH1	2.19	0.57
1:A:223:ASN:HB3	1:A:226:ILE:HG21	1.85	0.57
1:A:43:SER:CB	1:A:264:PRO:HB3	2.35	0.57
1:B:256:MET:HE1	1:B:261:PHE:HE2	1.70	0.57
1:B:195:VAL:HG11	1:B:256:MET:HE3	1.86	0.56
1:B:260:ARG:HG2	1:B:260:ARG:NH1	2.18	0.56
1:A:152:ARG:HB3	1:A:227:LEU:CD1	2.35	0.56
1:A:146:GLN:HG3	1:A:153:TRP:CE2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ALA:HB2	2:A:452:HOH:O	2.06	0.55
1:A:192:PHE:O	1:A:195:VAL:HG22	2.06	0.55
1:B:35:TYR:CE2	1:B:275:ILE:HB	2.41	0.55
1:B:195:VAL:HG11	1:B:256:MET:HE1	1.87	0.55
1:A:153:TRP:HB2	1:A:157:GLU:OE1	2.05	0.55
1:A:235:MET:HE2	1:B:235:MET:HG3	1.89	0.55
1:B:170:ASN:H	1:B:170:ASN:ND2	2.04	0.55
1:B:277:VAL:HA	1:B:280:TYR:CZ	2.41	0.55
1:A:1:MET:HE2	2:A:589:HOH:O	2.06	0.54
1:A:152:ARG:HB3	1:A:227:LEU:HD11	1.89	0.54
1:B:277:VAL:HA	1:B:280:TYR:CE2	2.42	0.54
1:A:148:VAL:C	1:A:150:ALA:H	2.13	0.54
1:A:228:LEU:HB2	1:A:237:VAL:HG23	1.90	0.53
1:B:146:GLN:HG3	1:B:153:TRP:CD2	2.43	0.53
1:B:281:ARG:HG2	1:B:281:ARG:NH1	2.23	0.53
1:B:223:ASN:N	2:B:988:HOH:O	2.42	0.53
1:A:232:TYR:O	1:A:235:MET:HB3	2.09	0.53
1:A:128:GLN:HG2	1:A:187:ARG:HH21	1.74	0.52
1:A:235:MET:CE	1:B:235:MET:HG3	2.39	0.52
1:A:235:MET:HG3	1:B:235:MET:HE2	1.92	0.52
1:B:199:ILE:CD1	1:B:273:ALA:HA	2.39	0.51
1:B:152:ARG:HG3	1:B:152:ARG:NH1	2.15	0.51
1:B:121:TYR:HB3	2:B:902:HOH:O	2.11	0.51
1:A:158:TYR:CD1	1:A:222:ASP:OD2	2.64	0.50
1:A:272:MET:O	1:A:275:ILE:HG22	2.11	0.50
1:B:281:ARG:HG2	1:B:281:ARG:O	2.10	0.50
1:B:146:GLN:HG3	1:B:153:TRP:CE2	2.45	0.50
1:A:229:LEU:HG	1:A:237:VAL:CG1	2.28	0.50
1:A:41:HIS:HB3	1:A:42:PRO:HD3	1.93	0.50
1:B:280:TYR:CD2	2:B:1010:HOH:O	2.55	0.49
1:B:252:LEU:O	1:B:256:MET:HG2	2.11	0.49
1:B:34:GLN:HB3	2:B:958:HOH:O	2.13	0.49
1:B:181:LEU:HD22	1:B:264:PRO:CD	2.43	0.49
1:B:189:ARG:NH2	1:B:260:ARG:O	2.43	0.49
1:B:277:VAL:HG22	1:B:282:THR:CG2	2.43	0.49
1:A:223:ASN:CB	1:A:226:ILE:HG22	2.42	0.48
1:A:250:ARG:O	1:A:254:ARG:HG3	2.14	0.48
1:A:237:VAL:CG1	1:A:238:VAL:H	2.25	0.48
1:A:1:MET:HB2	1:A:56:GLN:NE2	2.29	0.48
1:A:177:LEU:C	1:A:177:LEU:HD23	2.40	0.47
1:A:78:ASN:ND2	1:A:140:LEU:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:PHE:CD1	1:B:281:ARG:HG3	2.50	0.47
1:A:191:ALA:HB2	1:A:259:GLU:OE2	2.14	0.47
1:A:152:ARG:HD2	1:A:227:LEU:HD13	1.96	0.47
1:B:146:GLN:C	1:B:148:VAL:H	2.22	0.47
1:B:275:ILE:HG13	2:B:1010:HOH:O	2.14	0.47
1:B:275:ILE:HD11	1:B:280:TYR:CD2	2.50	0.47
1:A:270:GLU:HG2	2:A:605:HOH:O	2.14	0.46
1:B:256:MET:HE2	1:B:256:MET:CA	2.44	0.46
1:A:277:VAL:HA	1:A:280:TYR:CZ	2.51	0.46
1:B:199:ILE:HD11	1:B:273:ALA:N	2.30	0.46
1:A:39:ARG:NH2	2:A:401:HOH:O	2.45	0.45
1:B:280:TYR:HD2	2:B:1010:HOH:O	1.94	0.45
1:A:88:LEU:O	1:A:92:GLU:HG3	2.16	0.45
1:B:152:ARG:NH1	1:B:152:ARG:CG	2.76	0.45
1:A:237:VAL:CG1	1:A:238:VAL:N	2.76	0.45
1:A:254:ARG:HE	1:A:254:ARG:HB3	1.62	0.45
1:B:6:ARG:O	1:B:10:GLN:HG3	2.16	0.45
1:A:148:VAL:C	1:A:150:ALA:N	2.74	0.45
1:A:228:LEU:HB2	1:A:237:VAL:CG2	2.47	0.45
1:B:94:ALA:HB2	1:B:101:CYS:HA	1.98	0.44
1:B:199:ILE:HD12	1:B:199:ILE:HA	1.84	0.44
1:B:177:LEU:C	1:B:177:LEU:HD23	2.43	0.44
1:A:276:ARG:HD3	1:A:276:ARG:HA	1.73	0.43
1:B:28:GLU:HG2	1:B:284:THR:HB	2.00	0.43
1:A:44:LEU:HD23	1:A:44:LEU:HA	1.83	0.43
1:B:78:ASN:HD21	1:B:144:HIS:HB2	1.82	0.43
1:A:100:PRO:HA	1:A:110:ALA:HB2	2.00	0.43
1:A:148:VAL:HG12	1:A:149:ALA:H	1.79	0.43
1:B:163:ILE:HG23	1:B:200:SER:HB2	1.99	0.43
1:A:186:LEU:HD23	1:A:186:LEU:HA	1.89	0.43
1:B:45:ARG:NH1	2:B:947:HOH:O	2.50	0.43
1:A:202:ILE:O	1:A:206:GLN:HG3	2.20	0.42
1:A:6:ARG:NH2	1:A:58:GLU:OE1	2.53	0.42
1:A:226:ILE:HG23	1:A:227:LEU:N	2.35	0.42
1:B:182:ASP:OD1	1:B:185:THR:HG23	2.19	0.42
1:B:48:SER:OG	1:B:55:PRO:HD3	2.20	0.41
1:B:170:ASN:ND2	1:B:170:ASN:N	2.67	0.41
1:A:24:GLU:HB2	2:A:519:HOH:O	2.20	0.41
1:B:134:ARG:HD2	2:B:944:HOH:O	2.21	0.41
1:B:275:ILE:HG12	1:B:280:TYR:HE2	1.86	0.41
1:A:158:TYR:CE1	1:A:222:ASP:CG	2.98	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ASN:H	1:A:170:ASN:ND2	2.16	0.41
1:B:54:ASP:O	1:B:57:ASP:HB2	2.21	0.41
1:A:54:ASP:HA	1:A:55:PRO:HD3	1.95	0.40
1:B:272:MET:O	1:B:275:ILE:HG22	2.21	0.40
1:A:150:ALA:HB1	1:A:151:GLU:H	1.68	0.40
1:A:277:VAL:HG13	1:A:278:GLN:N	2.37	0.40
1:B:53:ARG:HA	1:B:53:ARG:HD3	1.77	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	266/300 (89%)	259 (97%)	5 (2%)	2 (1%)	16	6
1	B	272/300 (91%)	260 (96%)	8 (3%)	4 (2%)	8	2
All	All	538/600 (90%)	519 (96%)	13 (2%)	6 (1%)	11	3

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	ALA
1	B	149	ALA
1	B	275	ILE
1	B	4	THR
1	A	148	VAL
1	B	277	VAL

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	216/237 (91%)	211 (98%)	5 (2%)	44	30
1	B	222/237 (94%)	210 (95%)	12 (5%)	20	7
All	All	438/474 (92%)	421 (96%)	17 (4%)	28	13

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	ASN
1	A	226	ILE
1	A	260	ARG
1	A	264	PRO
1	A	278	GLN
1	B	48	SER
1	B	50	ARG
1	B	56	GLN
1	B	74	LEU
1	B	131	ASP
1	B	146	GLN
1	B	152	ARG
1	B	170	ASN
1	B	189	ARG
1	B	258	GLU
1	B	264	PRO
1	B	270	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	34	GLN
1	A	56	GLN
1	A	160	HIS
1	A	170	ASN

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Mol	Chain	Res	Type
1	A	194	GLN
1	A	206	GLN
1	A	207	ASN
1	A	223	ASN
1	B	78	ASN
1	B	170	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	270/300 (90%)	-1.41	0 100 100	18, 31, 57, 73	0
1	B	276/300 (92%)	-1.32	0 100 100	22, 37, 69, 82	0
All	All	546/600 (91%)	-1.36	0 100 100	18, 34, 64, 82	0

There are no RSRZ outliers to report.

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.