



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 03:46 pm BST

PDB ID : 4MC0
Title : Hedycaryol apo
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Deposited on : 2013-08-21
Resolution : 2.70 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	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

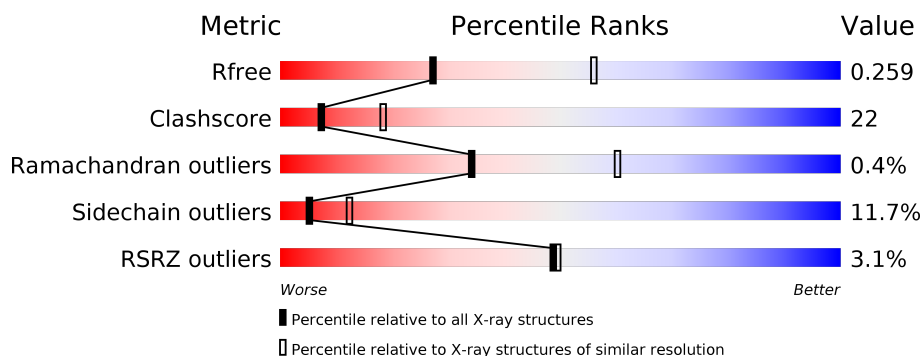
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 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	346	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> % 60% 21% • 16% </div> </div>
1	B	346	<div> <div style="width: 4%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; align-items: center;"> 4% 49% 29% 6% 16% </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4721 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 Putative sesquiterpene cyclase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			2311	1463	414	424	10			
1	B	290	Total	C	N	O	S	0	0	0
			2300	1457	410	423	10			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	LEU	-	EXPRESSION TAG	UNP E4MYY0
A	340	GLU	-	EXPRESSION TAG	UNP E4MYY0
A	341	HIS	-	EXPRESSION TAG	UNP E4MYY0
A	342	HIS	-	EXPRESSION TAG	UNP E4MYY0
A	343	HIS	-	EXPRESSION TAG	UNP E4MYY0
A	344	HIS	-	EXPRESSION TAG	UNP E4MYY0
A	345	HIS	-	EXPRESSION TAG	UNP E4MYY0
A	346	HIS	-	EXPRESSION TAG	UNP E4MYY0
B	339	LEU	-	EXPRESSION TAG	UNP E4MYY0
B	340	GLU	-	EXPRESSION TAG	UNP E4MYY0
B	341	HIS	-	EXPRESSION TAG	UNP E4MYY0
B	342	HIS	-	EXPRESSION TAG	UNP E4MYY0
B	343	HIS	-	EXPRESSION TAG	UNP E4MYY0
B	344	HIS	-	EXPRESSION TAG	UNP E4MYY0
B	345	HIS	-	EXPRESSION TAG	UNP E4MYY0
B	346	HIS	-	EXPRESSION TAG	UNP E4MYY0

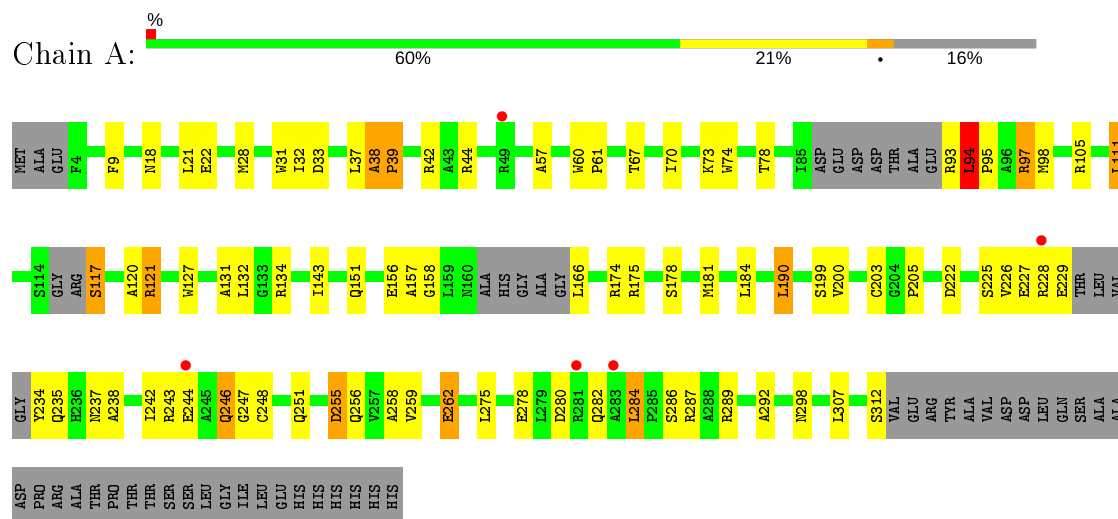
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	56	Total	O	0	0
			56	56		
2	B	54	Total	O	0	0
			54	54		

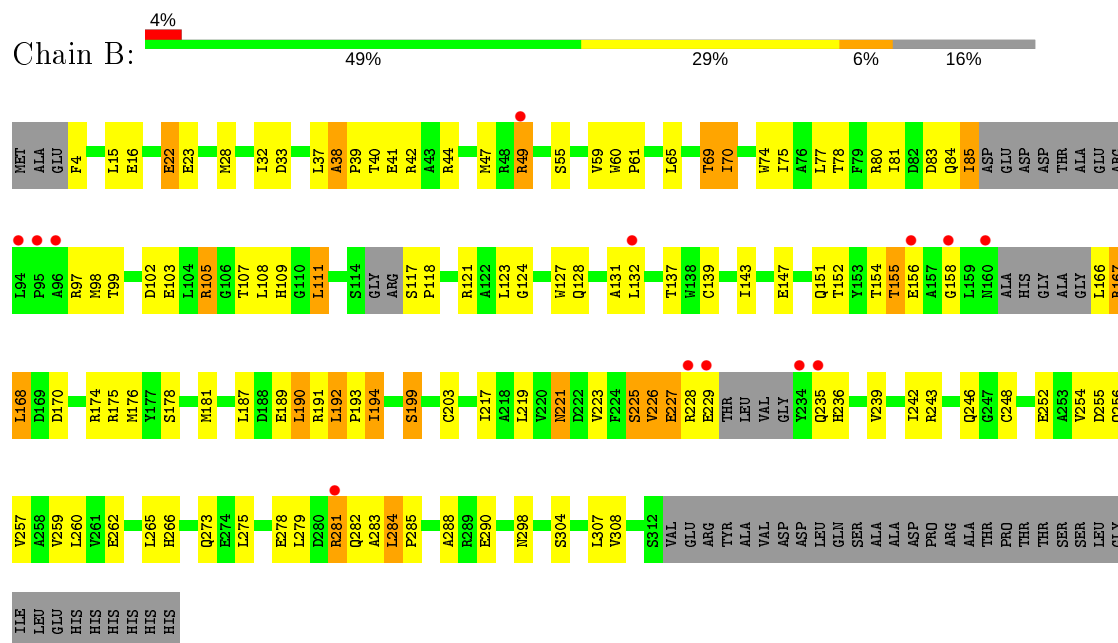
3 Residue-property plots

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: Putative sesquiterpene cyclase



• Molecule 1: Putative sesquiterpene cyclase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.38 Å 80.60 Å 97.55 Å 90.00° 94.69° 90.00°	Depositor
Resolution (Å)	15.00 – 2.70 14.97 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.3 (15.00-2.70) 95.3 (14.97-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.69 Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.207 , 0.247 0.217 , 0.259	Depositor DCC
R_{free} test set	1198 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.3	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	0.95	EDS
Total number of atoms	4721	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 51.76 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.3572e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.75	1/2362 (0.0%)	0.67	4/3214 (0.1%)
1	B	0.95	2/2351 (0.1%)	0.77	5/3200 (0.2%)
All	All	0.86	3/4713 (0.1%)	0.72	9/6414 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	39	PRO	N-CD	5.33	1.55	1.47
1	B	39	PRO	N-CD	5.19	1.55	1.47
1	B	193	PRO	N-CD	5.08	1.54	1.47

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	MET	C-N-CD	6.05	141.11	128.40
1	B	181	MET	C-N-CD	6.01	141.03	128.40
1	A	284	LEU	C-N-CD	5.83	140.64	128.40
1	A	117	SER	C-N-CD	5.82	140.62	128.40
1	B	60	TRP	C-N-CD	5.73	140.44	128.40
1	B	111	LEU	C-N-CD	5.63	140.23	128.40
1	B	38	ALA	C-N-CD	5.62	140.19	128.40
1	B	192	LEU	C-N-CD	5.53	140.01	128.40
1	A	38	ALA	C-N-CD	5.46	139.87	128.40

There are no chirality outliers.

There are no planarity outliers.

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	2311	0	2262	93	1
1	B	2300	0	2251	116	3
2	A	56	0	0	10	0
2	B	54	0	0	5	0
All	All	4721	0	4513	204	4

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 22.

All (204) 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:235:GLN:NE2	1:A:244:GLU:OE1	1.68	1.24
1:B:105:ARG:NE	1:B:147:GLU:OE1	1.76	1.18
1:B:80:ARG:CZ	1:B:84:GLN:OE1	1.96	1.14
1:A:93:ARG:HG2	1:A:94:LEU:HD23	1.40	1.01
1:A:94:LEU:HD23	1:A:94:LEU:H	1.20	1.01
1:B:227:GLU:OE2	1:B:227:GLU:N	1.94	0.99
1:A:61:PRO:HD2	1:A:298:ASN:HD21	1.24	0.97
1:B:151:GLN:O	1:B:154:THR:HG22	1.64	0.96
1:A:227:GLU:CG	2:A:420:HOH:O	2.13	0.96
1:B:246:GLN:HE22	1:B:256:GLN:HE22	1.01	0.95
1:B:152:THR:O	1:B:156:GLU:HG3	1.64	0.95
1:B:33:ASP:OD1	1:B:44:ARG:NH2	2.00	0.94
1:B:108:LEU:HD21	1:B:123:LEU:HD21	1.49	0.94
1:A:74:TRP:O	1:A:78:THR:HG22	1.69	0.92
1:A:127:TRP:CZ2	1:A:143:ILE:HD11	2.04	0.91
1:B:243:ARG:HH11	1:B:243:ARG:HG2	1.33	0.91
1:A:227:GLU:HG2	2:A:420:HOH:O	1.70	0.91
1:B:108:LEU:CD2	1:B:123:LEU:HD21	2.01	0.90
1:B:167:ARG:HG2	1:B:167:ARG:HH11	1.32	0.90
1:B:28:MET:O	1:B:32:ILE:HG13	1.76	0.85
1:B:226:VAL:HG12	1:B:227:GLU:OE2	1.76	0.85
1:B:70:ILE:CD1	1:B:190:LEU:HD23	2.07	0.85
1:A:93:ARG:CG	1:A:94:LEU:HD23	2.07	0.84
1:B:278:GLU:OE1	1:B:281:ARG:HD2	1.78	0.84
1:B:167:ARG:HD2	1:B:170:ASP:HB2	1.58	0.83
1:B:166:LEU:HD21	1:B:236:HIS:NE2	1.92	0.83
1:A:117:SER:HB3	1:A:120:ALA:HB3	1.60	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:ASN:O	1:B:225:SER:OG	1.96	0.82
1:B:70:ILE:HD11	1:B:190:LEU:HD23	1.62	0.81
1:B:166:LEU:CD1	1:B:235:GLN:HG3	2.10	0.81
1:A:243:ARG:HG2	1:A:243:ARG:HH11	1.42	0.81
1:A:235:GLN:HE22	1:A:244:GLU:CD	1.82	0.81
1:B:166:LEU:CD2	1:B:236:HIS:NE2	2.45	0.80
1:B:80:ARG:NH2	1:B:84:GLN:OE1	2.15	0.79
1:B:105:ARG:HE	1:B:147:GLU:CD	1.87	0.78
1:A:93:ARG:HG2	1:A:94:LEU:CD2	2.14	0.77
1:A:31:TRP:CD1	1:A:73:LYS:HE2	2.19	0.77
1:B:152:THR:O	1:B:156:GLU:CG	2.33	0.77
1:A:94:LEU:HD23	1:A:94:LEU:N	1.99	0.76
1:A:238:ALA:O	1:A:242:ILE:HG13	1.85	0.76
1:B:246:GLN:HE22	1:B:256:GLN:NE2	1.82	0.75
1:A:235:GLN:CD	1:A:244:GLU:OE1	2.24	0.75
1:B:167:ARG:HG2	1:B:167:ARG:NH1	1.98	0.74
1:B:80:ARG:NH1	1:B:84:GLN:OE1	2.21	0.74
1:A:280:ASP:OD1	1:A:289:ARG:NH1	2.20	0.74
1:A:127:TRP:CH2	1:A:131:ALA:HB2	2.23	0.73
1:A:227:GLU:HG3	2:A:420:HOH:O	1.80	0.73
1:B:243:ARG:NH1	1:B:243:ARG:HG2	1.96	0.73
1:A:199:SER:O	1:A:203:CYS:HB3	1.87	0.73
1:A:222:ASP:HB3	1:A:237:ASN:OD1	1.90	0.72
1:A:94:LEU:N	1:A:95:PRO:HD3	2.05	0.72
1:A:235:GLN:HG3	1:A:235:GLN:O	1.88	0.72
1:B:189:GLU:HA	1:B:192:LEU:HD12	1.73	0.71
1:A:70:ILE:HD11	1:A:190:LEU:HD23	1.73	0.70
1:A:127:TRP:CZ3	1:A:131:ALA:HB2	2.26	0.70
1:B:127:TRP:CH2	1:B:131:ALA:HB2	2.27	0.70
1:A:93:ARG:CD	1:A:94:LEU:HD23	2.21	0.69
1:B:174:ARG:HH11	1:B:236:HIS:CD2	2.10	0.69
1:A:94:LEU:N	1:A:95:PRO:CD	2.56	0.69
1:B:166:LEU:HD21	1:B:236:HIS:CE1	2.28	0.69
1:B:109:HIS:NE2	1:B:147:GLU:OE2	2.26	0.68
1:A:175:ARG:HD2	1:A:178:SER:OG	1.94	0.68
1:A:33:ASP:OD1	1:A:44:ARG:NH2	2.27	0.68
1:A:151:GLN:HG2	1:B:246:GLN:O	1.92	0.68
1:B:166:LEU:CD1	1:B:235:GLN:CG	2.73	0.67
1:A:21:LEU:HD13	1:A:57:ALA:HB2	1.76	0.67
1:A:61:PRO:HD2	1:A:298:ASN:ND2	2.04	0.67
1:A:205:PRO:HB2	1:A:275:LEU:HD12	1.74	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ARG:C	1:A:95:PRO:CD	2.63	0.67
1:A:229:GLU:HB2	1:A:234:TYR:HB3	1.78	0.66
1:A:289:ARG:HD3	2:A:412:HOH:O	1.95	0.66
2:A:447:HOH:O	1:B:168:LEU:HD23	1.94	0.66
1:B:242:ILE:O	1:B:246:GLN:HB2	1.96	0.65
1:A:117:SER:HB3	1:A:120:ALA:CB	2.28	0.64
1:A:255:ASP:N	1:A:255:ASP:OD1	2.31	0.64
1:B:166:LEU:HD11	1:B:235:GLN:CG	2.29	0.63
1:B:108:LEU:HD23	1:B:123:LEU:HD21	1.81	0.62
1:B:22:GLU:OE2	1:B:22:GLU:HA	1.99	0.62
1:A:93:ARG:HG2	1:A:94:LEU:H	1.63	0.62
1:B:166:LEU:HD11	1:B:235:GLN:HG3	1.80	0.62
1:A:31:TRP:CG	1:A:73:LYS:HE2	2.35	0.62
1:A:93:ARG:HD3	1:A:94:LEU:CD2	2.29	0.61
1:B:49:ARG:CG	1:B:49:ARG:HH21	2.13	0.61
1:B:15:LEU:HD13	1:B:308:VAL:HG21	1.83	0.61
1:A:94:LEU:H	1:A:94:LEU:CD2	2.07	0.61
1:B:77:LEU:HD21	1:B:123:LEU:HD12	1.83	0.60
1:A:93:ARG:C	1:A:95:PRO:HD2	2.22	0.60
1:B:246:GLN:NE2	1:B:256:GLN:HE22	1.86	0.60
1:A:259:VAL:O	1:A:262:GLU:OE2	2.20	0.60
1:B:194:ILE:HD11	1:B:288:ALA:HA	1.84	0.60
1:B:49:ARG:HG3	1:B:49:ARG:HH21	1.67	0.59
1:A:28:MET:O	1:A:32:ILE:HG13	2.04	0.57
1:B:15:LEU:HG	1:B:16:GLU:N	2.17	0.57
1:B:70:ILE:HD11	1:B:190:LEU:CD2	2.32	0.57
1:B:278:GLU:OE1	1:B:281:ARG:CD	2.49	0.57
1:B:199:SER:O	1:B:203:CYS:HB3	2.03	0.57
1:A:93:ARG:CG	1:A:94:LEU:H	2.17	0.57
1:B:265:LEU:O	1:B:265:LEU:HD12	2.04	0.57
1:A:38:ALA:HB1	1:A:44:ARG:HG3	1.87	0.57
1:B:227:GLU:HG2	2:B:445:HOH:O	2.05	0.57
1:A:70:ILE:CD1	1:A:190:LEU:HD23	2.35	0.56
1:A:158:GLY:HA3	1:B:256:GLN:NE2	2.21	0.56
1:A:18:ASN:HB2	1:A:61:PRO:HA	1.87	0.56
1:A:243:ARG:HG2	1:A:243:ARG:NH1	2.19	0.56
1:B:74:TRP:O	1:B:78:THR:HG22	2.06	0.55
1:A:235:GLN:OE1	1:A:244:GLU:OE1	2.24	0.55
1:A:278:GLU:O	1:A:282:GLN:HG3	2.06	0.55
1:B:282:GLN:O	1:B:283:ALA:HB3	2.06	0.55
1:B:243:ARG:CG	1:B:243:ARG:HH11	2.13	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:PRO:HD2	1:B:298:ASN:HD21	1.72	0.54
1:B:255:ASP:O	1:B:259:VAL:HG23	2.08	0.54
1:B:235:GLN:HA	1:B:235:GLN:OE1	2.08	0.54
1:B:266:HIS:HB3	2:B:402:HOH:O	2.06	0.54
2:A:447:HOH:O	1:B:168:LEU:CD2	2.55	0.54
1:A:93:ARG:O	1:A:95:PRO:HD2	2.07	0.53
1:A:93:ARG:HD3	1:A:94:LEU:HD21	1.90	0.53
1:A:243:ARG:HH11	1:A:243:ARG:CG	2.14	0.53
1:B:32:ILE:HG23	1:B:37:LEU:HB2	1.90	0.53
1:A:174:ARG:NH2	1:B:252:GLU:OE2	2.42	0.52
1:A:151:GLN:CG	1:B:246:GLN:O	2.58	0.52
1:B:187:LEU:O	1:B:190:LEU:HB2	2.09	0.52
1:A:156:GLU:CG	1:A:174:ARG:HE	2.23	0.52
1:A:9:PHE:HE1	2:A:453:HOH:O	1.93	0.52
1:B:139:CYS:O	1:B:143:ILE:HG12	2.10	0.52
1:A:31:TRP:CZ3	1:A:37:LEU:HD11	2.46	0.51
1:A:93:ARG:C	1:A:95:PRO:HD3	2.30	0.51
1:B:37:LEU:O	1:B:80:ARG:HD3	2.11	0.51
1:A:243:ARG:NH1	1:A:243:ARG:CG	2.73	0.51
1:A:93:ARG:CD	1:A:94:LEU:CD2	2.88	0.51
1:B:42:ARG:HD3	2:B:408:HOH:O	2.11	0.49
1:B:49:ARG:CG	1:B:49:ARG:NH2	2.73	0.49
1:B:117:SER:O	1:B:121:ARG:HG3	2.12	0.49
1:B:124:GLY:O	1:B:128:GLN:HB2	2.13	0.49
1:B:81:ILE:O	1:B:85:ILE:HG13	2.12	0.49
1:A:246:GLN:NE2	1:A:256:GLN:OE1	2.44	0.49
1:B:23:GLU:HB3	1:B:65:LEU:HD11	1.95	0.49
1:A:94:LEU:N	1:A:94:LEU:CD2	2.73	0.49
1:A:111:LEU:HD12	2:A:449:HOH:O	2.12	0.49
1:B:15:LEU:CD1	1:B:308:VAL:HG21	2.43	0.49
1:B:275:LEU:O	1:B:279:LEU:HG	2.13	0.48
1:A:111:LEU:CD1	2:A:449:HOH:O	2.61	0.48
1:B:248:CYS:HB2	1:B:252:GLU:OE1	2.13	0.48
1:A:158:GLY:HA3	1:B:256:GLN:HE21	1.78	0.48
1:B:243:ARG:HD2	2:B:429:HOH:O	2.13	0.48
1:A:97:ARG:HA	1:A:97:ARG:HD3	1.54	0.47
1:B:107:THR:OG1	1:B:123:LEU:HD23	2.14	0.47
1:B:108:LEU:HD23	1:B:123:LEU:CD2	2.44	0.47
1:B:41:GLU:HA	1:B:41:GLU:OE1	2.15	0.47
1:A:32:ILE:HG23	1:A:37:LEU:HB2	1.96	0.47
1:B:166:LEU:HD13	1:B:235:GLN:HG3	1.91	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:226:VAL:HG21	1:B:239:VAL:HG11	1.97	0.47
1:A:246:GLN:C	1:A:248:CYS:H	2.17	0.47
1:A:97:ARG:HG2	1:A:157:ALA:CB	2.45	0.46
1:B:167:ARG:CD	1:B:170:ASP:HB2	2.38	0.46
1:B:166:LEU:HD22	1:B:236:HIS:CD2	2.50	0.46
1:A:67:THR:O	1:A:70:ILE:HG12	2.16	0.46
1:B:166:LEU:HD11	1:B:235:GLN:HG2	1.98	0.46
1:B:38:ALA:HB1	1:B:44:ARG:HG3	1.97	0.45
1:B:105:ARG:CZ	1:B:147:GLU:OE1	2.56	0.45
1:A:184:LEU:HD23	1:A:184:LEU:HA	1.79	0.45
1:A:105:ARG:HG2	1:A:105:ARG:O	2.17	0.45
1:B:4:PHE:CD2	1:B:254:VAL:HG21	2.53	0.44
1:B:154:THR:HG23	1:B:155:THR:N	2.32	0.44
1:B:155:THR:O	1:B:158:GLY:N	2.51	0.44
1:B:226:VAL:CG1	1:B:227:GLU:OE2	2.56	0.44
1:A:251:GLN:HG3	1:A:251:GLN:O	2.17	0.44
1:B:15:LEU:HA	2:B:411:HOH:O	2.17	0.44
1:B:175:ARG:HD2	1:B:178:SER:OG	2.18	0.44
1:A:93:ARG:HG2	1:A:94:LEU:N	2.30	0.44
1:A:225:SER:O	1:A:229:GLU:HG2	2.17	0.43
1:A:93:ARG:CG	1:A:94:LEU:CD2	2.86	0.43
1:B:284:LEU:HB3	1:B:285:PRO:HD2	1.99	0.43
1:B:219:LEU:O	1:B:223:VAL:HG23	2.18	0.43
1:B:15:LEU:HD13	1:B:308:VAL:CG2	2.48	0.43
1:A:227:GLU:HG3	1:A:228:ARG:N	2.33	0.43
1:A:60:TRP:CZ2	1:A:184:LEU:HD22	2.53	0.43
1:B:273:GLN:OE1	1:B:273:GLN:HA	2.18	0.43
1:A:258:ALA:HA	2:A:453:HOH:O	2.18	0.43
1:B:123:LEU:O	1:B:123:LEU:HG	2.18	0.42
1:B:156:GLU:OE2	1:B:174:ARG:NE	2.20	0.42
1:B:65:LEU:O	1:B:69:THR:OG1	2.33	0.42
1:A:246:GLN:O	1:A:248:CYS:N	2.52	0.42
1:B:103:GLU:O	1:B:107:THR:HG23	2.20	0.42
1:A:93:ARG:HD2	1:A:95:PRO:HD3	2.01	0.42
1:B:127:TRP:HA	1:B:127:TRP:CE3	2.55	0.42
1:B:154:THR:CG2	1:B:155:THR:N	2.82	0.42
1:A:226:VAL:HG13	1:A:227:GLU:N	2.35	0.42
1:A:243:ARG:HG3	1:A:248:CYS:O	2.20	0.41
1:B:217:ILE:O	1:B:221:ASN:HB2	2.19	0.41
1:B:257:VAL:O	1:B:260:LEU:HB2	2.20	0.41
1:A:205:PRO:HB2	1:A:275:LEU:CD1	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:GLU:HB3	1:B:65:LEU:CD1	2.50	0.41
1:B:127:TRP:CE2	1:B:143:ILE:HD11	2.55	0.41
1:B:166:LEU:HD22	1:B:236:HIS:NE2	2.30	0.41
1:A:117:SER:O	1:A:121:ARG:HD2	2.20	0.41
1:B:117:SER:HA	1:B:118:PRO:HD2	1.94	0.41
1:A:278:GLU:HG3	1:A:278:GLU:O	2.21	0.41
1:B:37:LEU:HA	1:B:37:LEU:HD23	1.88	0.41
1:B:55:SER:O	1:B:59:VAL:HG23	2.21	0.41
1:A:93:ARG:CG	1:A:94:LEU:N	2.83	0.41
1:B:167:ARG:CG	1:B:167:ARG:HH11	2.13	0.41
1:B:151:GLN:HE21	1:B:154:THR:HG21	1.86	0.40
1:A:200:VAL:HG11	1:A:292:ALA:HB2	2.03	0.40
1:B:75:ILE:HA	1:B:78:THR:HG22	2.04	0.40

All (4) 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:128:GLN:NE2	1:B:266:HIS:CD2[4_748]	1.47	0.73
1:B:44:ARG:NH1	1:B:44:ARG:NH1[2_858]	1.79	0.41
1:A:44:ARG:NH1	1:A:44:ARG:NH1[2_757]	1.99	0.21
1:B:128:GLN:NE2	1:B:266:HIS:NE2[4_748]	2.08	0.12

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	281/346 (81%)	277 (99%)	2 (1%)	2 (1%)	22	46
1	B	280/346 (81%)	277 (99%)	3 (1%)	0	100	100
All	All	561/692 (81%)	554 (99%)	5 (1%)	2 (0%)	34	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	LEU
1	A	247	GLY

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	239/281 (85%)	219 (92%)	20 (8%)	11	25
1	B	238/281 (85%)	202 (85%)	36 (15%)	3	7
All	All	477/562 (85%)	421 (88%)	56 (12%)	5	12

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	39	PRO
1	A	42	ARG
1	A	94	LEU
1	A	97	ARG
1	A	98	MET
1	A	111	LEU
1	A	121	ARG
1	A	132	LEU
1	A	134	ARG
1	A	166	LEU
1	A	190	LEU
1	A	246	GLN
1	A	255	ASP
1	A	262	GLU
1	A	284	LEU
1	A	286	SER
1	A	287	ARG
1	A	307	LEU
1	A	312	SER
1	B	22	GLU

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Mol	Chain	Res	Type
1	B	40	THR
1	B	47	MET
1	B	49	ARG
1	B	69	THR
1	B	70	ILE
1	B	83	ASP
1	B	85	ILE
1	B	97	ARG
1	B	98	MET
1	B	99	THR
1	B	102	ASP
1	B	105	ARG
1	B	111	LEU
1	B	132	LEU
1	B	137	THR
1	B	155	THR
1	B	167	ARG
1	B	168	LEU
1	B	176	MET
1	B	190	LEU
1	B	191	ARG
1	B	194	ILE
1	B	199	SER
1	B	221	ASN
1	B	225	SER
1	B	226	VAL
1	B	227	GLU
1	B	228	ARG
1	B	229	GLU
1	B	262	GLU
1	B	281	ARG
1	B	284	LEU
1	B	290	GLU
1	B	304	SER
1	B	307	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	160	ASN
1	A	236	HIS

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Mol	Chain	Res	Type
1	A	246	GLN
1	A	256	GLN
1	A	298	ASN
1	B	35	ASN
1	B	256	GLN
1	B	298	ASN
1	B	302	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 carbohydrates 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	291/346 (84%)	-0.10	5 (1%) 70 72	29, 54, 89, 110	0
1	B	290/346 (83%)	-0.11	13 (4%) 33 31	29, 52, 96, 132	0
All	All	581/692 (83%)	-0.10	18 (3%) 49 49	29, 53, 93, 132	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	95	PRO	4.4
1	B	229	GLU	4.3
1	B	96	ALA	4.2
1	B	234	TYR	3.9
1	B	235	GLN	3.8
1	A	228	ARG	3.8
1	B	228	ARG	3.6
1	A	49	ARG	3.4
1	B	49	ARG	3.2
1	B	160	ASN	3.0
1	B	132	LEU	2.6
1	A	244	GLU	2.5
1	A	283	ALA	2.4
1	B	156	GLU	2.4
1	B	94	LEU	2.2
1	B	158	GLY	2.1
1	B	281	ARG	2.1
1	A	281	ARG	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.