



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:46 AM UTC

PDB ID : 3MQ3 / pdb_00003mq3
Title : Crystal structure of native bovine PDP1c
Authors : Guo, Y.; Ernst, S.R.; Carroll, D.W.; Hackert, M.L.
Deposited on : 2010-04-27
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
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

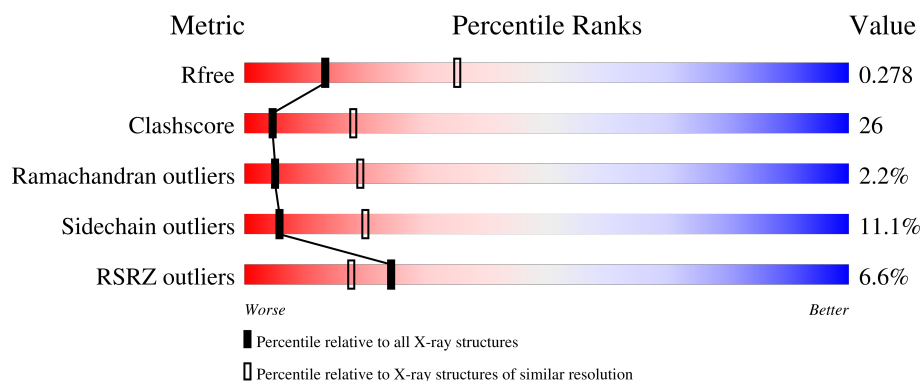
1 Overall quality at a glance

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	467	5% 53% 21% 6% 19%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3068 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 Pyruvate dehydrogenase phosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	0	0	0
			2967	1885	516	556	10			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		

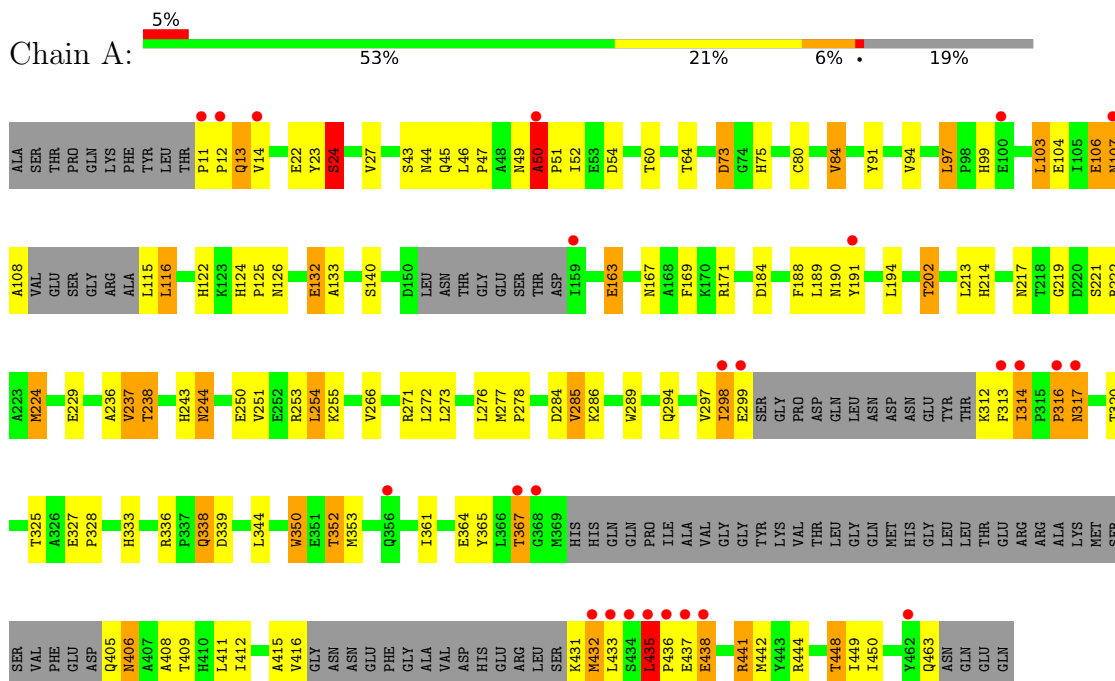
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	99	Total	O	0	0
			99	99		

3 Residue-property plots

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: Pyruvate dehydrogenase phosphatase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	75.27Å 75.27Å 172.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.00 – 2.80 27.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.0 (27.00-2.80) 97.0 (27.00-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.43 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.224 , 0.284 0.226 , 0.278	Depositor DCC
R_{free} test set	716 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.088	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 71.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3068	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% 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

Bond lengths and bond angles in the following residue types are not validated in this section: MN

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	1.09	2/3031 (0.1%)	1.21	13/4119 (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	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	450	ILE	CA-CB	-5.69	1.47	1.54
1	A	236	ALA	CA-CB	-5.28	1.46	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	314	ILE	CB-CA-C	7.63	117.53	110.13
1	A	352	THR	N-CA-C	6.96	118.95	111.36
1	A	217	ASN	N-CA-C	6.57	119.61	108.90
1	A	277	MET	CA-C-N	-5.70	114.06	119.76
1	A	277	MET	C-N-CA	-5.70	114.06	119.76
1	A	237	VAL	N-CA-C	5.67	116.03	107.80
1	A	64	THR	N-CA-C	-5.24	101.45	109.41
1	A	314	ILE	N-CA-C	5.23	113.83	107.61
1	A	350	TRP	N-CA-C	5.19	116.94	111.28
1	A	73	ASP	CA-C-N	5.16	125.24	120.34
1	A	73	ASP	C-N-CA	5.16	125.24	120.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	PHE	CA-C-N	-5.10	118.89	123.33
1	A	313	PHE	C-N-CA	-5.10	118.89	123.33

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	312	LYS	Peptide
1	A	317	ASN	Peptide
1	A	435	LEU	Peptide
1	A	50	ALA	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	2967	0	2858	151	0
2	A	2	0	0	0	0
3	A	99	0	0	10	0
All	All	3068	0	2858	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 26.

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:435:LEU:CD1	1:A:436:PRO:HD3	1.39	1.49
1:A:435:LEU:HD22	1:A:436:PRO:N	1.36	1.35
1:A:435:LEU:HD13	1:A:436:PRO:CD	1.57	1.32
1:A:435:LEU:HD22	1:A:436:PRO:CA	1.66	1.25
1:A:50:ALA:CB	1:A:51:PRO:HD2	1.64	1.24
1:A:224:MET:HE2	1:A:238:THR:HA	1.26	1.15
1:A:50:ALA:HB3	1:A:51:PRO:CD	1.79	1.12
1:A:50:ALA:CB	1:A:51:PRO:CD	2.30	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:LEU:HD22	1:A:436:PRO:CD	1.82	1.09
1:A:438:GLU:OE2	1:A:438:GLU:HA	1.58	1.01
1:A:435:LEU:HD21	1:A:436:PRO:HG3	1.44	0.99
1:A:294:GLN:O	1:A:298:ILE:HG12	1.63	0.98
1:A:11:PRO:HA	3:A:548:HOH:O	1.66	0.95
1:A:54:ASP:OD2	1:A:448:THR:HG22	1.67	0.94
1:A:435:LEU:CD2	1:A:436:PRO:N	2.30	0.94
1:A:435:LEU:CD2	1:A:436:PRO:CG	2.48	0.92
1:A:435:LEU:CD2	1:A:436:PRO:CD	2.47	0.92
1:A:124:HIS:HD2	1:A:126:ASN:H	1.11	0.92
1:A:435:LEU:CD2	1:A:436:PRO:CA	2.49	0.89
1:A:435:LEU:CD2	1:A:436:PRO:HG3	2.02	0.89
1:A:435:LEU:HD22	1:A:436:PRO:CG	2.04	0.87
1:A:435:LEU:CG	1:A:436:PRO:HD3	2.05	0.87
1:A:442:MET:HA	1:A:442:MET:HE2	1.55	0.87
1:A:50:ALA:HB3	1:A:51:PRO:HD2	0.89	0.85
1:A:435:LEU:CD1	1:A:436:PRO:CD	2.30	0.85
1:A:224:MET:CE	1:A:238:THR:HA	2.07	0.84
1:A:11:PRO:HB2	1:A:14:VAL:H	1.42	0.83
1:A:124:HIS:CD2	1:A:126:ASN:H	1.96	0.83
1:A:431:LYS:HD3	1:A:432:MET:HB2	1.62	0.81
1:A:435:LEU:H	1:A:436:PRO:CD	1.97	0.78
1:A:325:THR:HG21	3:A:498:HOH:O	1.84	0.77
1:A:250:GLU:CD	1:A:253:ARG:NH1	2.43	0.77
1:A:435:LEU:HD22	1:A:436:PRO:HA	1.65	0.76
1:A:435:LEU:HD13	1:A:435:LEU:H	1.54	0.72
1:A:184:ASP:H	1:A:190:ASN:HD22	1.37	0.72
1:A:75:HIS:HE1	1:A:284:ASP:OD1	1.72	0.71
1:A:75:HIS:CE1	1:A:284:ASP:OD1	2.43	0.71
1:A:276:LEU:HD22	1:A:278:PRO:HD2	1.72	0.71
1:A:365:TYR:HE1	1:A:405:GLN:HG3	1.55	0.70
1:A:50:ALA:HB1	1:A:51:PRO:CD	2.21	0.70
1:A:435:LEU:H	1:A:436:PRO:HD3	1.55	0.70
1:A:435:LEU:CD2	1:A:436:PRO:HA	2.19	0.70
1:A:435:LEU:HD22	1:A:435:LEU:C	2.13	0.69
1:A:250:GLU:CD	1:A:253:ARG:HH12	2.00	0.68
1:A:214:HIS:HD2	3:A:469:HOH:O	1.75	0.68
1:A:11:PRO:HB2	1:A:13:GLN:H	1.58	0.67
1:A:94:VAL:HA	1:A:97:LEU:HD22	1.77	0.66
1:A:406:ASN:HD22	1:A:409:THR:H	1.45	0.65
1:A:432:MET:SD	3:A:563:HOH:O	2.54	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:ASP:HB3	1:A:448:THR:HG21	1.78	0.65
1:A:435:LEU:HD22	1:A:436:PRO:CB	2.28	0.64
1:A:75:HIS:HD2	3:A:470:HOH:O	1.82	0.63
1:A:435:LEU:CG	1:A:436:PRO:CD	2.72	0.62
1:A:352:THR:HG22	1:A:353:MET:HG3	1.81	0.62
1:A:73:ASP:H	1:A:202:THR:HG22	1.65	0.62
1:A:238:THR:O	1:A:238:THR:HG22	2.00	0.61
1:A:124:HIS:HD2	1:A:126:ASN:N	1.92	0.61
1:A:353:MET:HE1	1:A:361:ILE:HD12	1.81	0.61
1:A:438:GLU:OE2	1:A:438:GLU:CA	2.42	0.61
1:A:435:LEU:HD13	1:A:435:LEU:N	2.16	0.60
1:A:273:LEU:HD13	1:A:316:PRO:HB2	1.84	0.60
1:A:229:GLU:HG3	3:A:506:HOH:O	2.00	0.60
1:A:435:LEU:HD13	1:A:436:PRO:HD3	0.65	0.60
1:A:317:ASN:N	3:A:550:HOH:O	2.34	0.60
1:A:45:GLN:HE21	1:A:448:THR:HB	1.67	0.60
1:A:132:GLU:HG3	1:A:188:PHE:HE1	1.67	0.59
1:A:406:ASN:ND2	1:A:409:THR:H	2.01	0.59
1:A:251:VAL:HG12	1:A:255:LYS:HE2	1.85	0.59
1:A:333:HIS:HE1	1:A:339:ASP:OD2	1.85	0.59
1:A:99:HIS:CE1	1:A:103:LEU:CD2	2.85	0.59
1:A:132:GLU:HG3	1:A:188:PHE:CE1	2.37	0.58
1:A:202:THR:HG21	1:A:221:SER:OG	2.03	0.58
1:A:75:HIS:HE1	1:A:284:ASP:CG	2.10	0.58
1:A:298:ILE:O	1:A:299:GLU:OE2	2.21	0.58
1:A:11:PRO:HB2	1:A:14:VAL:N	2.16	0.57
1:A:432:MET:C	3:A:563:HOH:O	2.47	0.57
1:A:244:ASN:C	1:A:244:ASN:HD22	2.13	0.57
1:A:298:ILE:O	1:A:299:GLU:CD	2.48	0.56
1:A:297:VAL:O	1:A:299:GLU:N	2.38	0.56
1:A:80:CYS:O	1:A:84:VAL:HG13	2.06	0.55
1:A:276:LEU:CD2	1:A:278:PRO:HD2	2.36	0.54
1:A:99:HIS:CE1	1:A:103:LEU:HD21	2.42	0.54
1:A:43:SER:HA	1:A:449:ILE:O	2.09	0.53
1:A:250:GLU:OE2	1:A:253:ARG:NH1	2.42	0.53
1:A:435:LEU:N	1:A:436:PRO:CD	2.65	0.53
1:A:353:MET:CE	1:A:361:ILE:HD12	2.39	0.53
1:A:435:LEU:CD1	1:A:436:PRO:CG	2.87	0.53
1:A:124:HIS:CD2	1:A:125:PRO:HD2	2.45	0.52
1:A:106:GLU:HA	1:A:106:GLU:OE1	2.08	0.52
1:A:116:LEU:H	1:A:116:LEU:CD2	2.21	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:MET:HE2	1:A:238:THR:CA	2.18	0.52
1:A:365:TYR:HE1	1:A:405:GLN:CG	2.22	0.52
1:A:133:ALA:HA	1:A:189:LEU:HD21	1.93	0.51
1:A:23:TYR:CG	1:A:24:SER:N	2.79	0.51
1:A:47:PRO:HA	1:A:52:ILE:HD12	1.93	0.51
1:A:184:ASP:N	1:A:190:ASN:HD22	2.07	0.51
1:A:433:LEU:O	1:A:435:LEU:HD13	2.11	0.51
1:A:285:VAL:HG13	1:A:289:TRP:CD2	2.46	0.50
1:A:47:PRO:CA	1:A:52:ILE:HD12	2.41	0.50
1:A:317:ASN:ND2	1:A:317:ASN:C	2.70	0.50
1:A:435:LEU:CD2	1:A:435:LEU:C	2.82	0.49
1:A:107:ASN:O	1:A:108:ALA:C	2.55	0.49
1:A:116:LEU:CD2	1:A:116:LEU:N	2.77	0.48
1:A:435:LEU:CD1	1:A:435:LEU:N	2.77	0.48
1:A:433:LEU:O	1:A:435:LEU:CD1	2.62	0.47
1:A:435:LEU:N	1:A:436:PRO:HD3	2.19	0.47
1:A:406:ASN:ND2	1:A:408:ALA:H	2.12	0.47
1:A:169:PHE:CG	1:A:328:PRO:HG3	2.49	0.47
1:A:222:ARG:HB3	1:A:350:TRP:NE1	2.30	0.45
1:A:336:ARG:HB3	1:A:338:GLN:OE1	2.17	0.45
1:A:435:LEU:CD2	1:A:436:PRO:CB	2.92	0.45
1:A:49:ASN:O	1:A:50:ALA:O	2.33	0.45
1:A:132:GLU:H	1:A:132:GLU:CD	2.23	0.45
1:A:116:LEU:N	1:A:116:LEU:HD22	2.31	0.45
1:A:22:GLU:O	1:A:22:GLU:HG3	2.17	0.45
1:A:22:GLU:O	1:A:22:GLU:CG	2.64	0.45
1:A:367:THR:O	1:A:367:THR:HG23	2.18	0.44
1:A:352:THR:HG21	1:A:415:ALA:HA	1.98	0.44
1:A:416:VAL:O	1:A:416:VAL:HG12	2.11	0.44
1:A:75:HIS:HE1	1:A:284:ASP:OD2	2.00	0.44
1:A:435:LEU:HD21	1:A:436:PRO:CG	2.19	0.44
1:A:442:MET:HE2	1:A:442:MET:CA	2.39	0.44
1:A:325:THR:HB	1:A:327:GLU:H	1.82	0.43
1:A:54:ASP:HB3	1:A:448:THR:CG2	2.46	0.43
1:A:255:LYS:HE3	3:A:505:HOH:O	2.18	0.43
1:A:411:LEU:HD23	1:A:411:LEU:HA	1.92	0.43
1:A:364:GLU:O	1:A:365:TYR:C	2.62	0.43
1:A:23:TYR:O	1:A:24:SER:CB	2.67	0.43
1:A:406:ASN:HD22	1:A:406:ASN:C	2.26	0.43
1:A:219:GLY:H	1:A:243:HIS:HD2	1.67	0.43
1:A:99:HIS:HE1	1:A:103:LEU:HD21	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:ARG:HA	1:A:444:ARG:O	2.19	0.42
1:A:271:ARG:HD2	3:A:512:HOH:O	2.19	0.42
1:A:73:ASP:N	1:A:202:THR:HG22	2.34	0.41
1:A:12:PRO:O	1:A:13:GLN:CB	2.68	0.41
1:A:163:GLU:O	1:A:167:ASN:ND2	2.53	0.41
1:A:406:ASN:ND2	1:A:406:ASN:C	2.78	0.41
1:A:49:ASN:HD21	1:A:441:ARG:HH21	1.69	0.41
1:A:91:TYR:CE2	1:A:171:ARG:HG3	2.55	0.41
1:A:365:TYR:CE1	1:A:405:GLN:HG3	2.44	0.41
1:A:251:VAL:O	1:A:255:LYS:HG3	2.21	0.41
1:A:412:ILE:HG13	1:A:449:ILE:HD12	2.03	0.41
1:A:50:ALA:HB1	1:A:51:PRO:HD3	2.01	0.41
1:A:191:TYR:HD1	1:A:191:TYR:HA	1.73	0.41
1:A:254:LEU:HB3	1:A:266:VAL:HG11	2.02	0.40
1:A:49:ASN:O	1:A:50:ALA:C	2.64	0.40
1:A:317:ASN:C	1:A:317:ASN:HD22	2.29	0.40
1:A:60:THR:OG1	1:A:122:HIS:HE1	2.04	0.40
1:A:103:LEU:O	1:A:104:GLU:C	2.64	0.40
1:A:44:ASN:HB3	1:A:412:ILE:HD12	2.02	0.40
1:A:442:MET:HA	1:A:442:MET:CE	2.40	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	366/467 (78%)	342 (93%)	16 (4%)	8 (2%)	5 19

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	13	GLN

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Mol	Chain	Res	Type
1	A	50	ALA
1	A	298	ILE
1	A	24	SER
1	A	107	ASN
1	A	435	LEU
1	A	316	PRO
1	A	437	GLU

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	314/404 (78%)	279 (89%)	35 (11%)	6	20

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	SER
1	A	27	VAL
1	A	46	LEU
1	A	84	VAL
1	A	97	LEU
1	A	103	LEU
1	A	106	GLU
1	A	115	LEU
1	A	116	LEU
1	A	132	GLU
1	A	140	SER
1	A	163	GLU
1	A	194	LEU
1	A	202	THR
1	A	213	LEU
1	A	224	MET
1	A	237	VAL
1	A	238	THR
1	A	244	ASN

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Mol	Chain	Res	Type
1	A	254	LEU
1	A	272	LEU
1	A	285	VAL
1	A	286	LYS
1	A	314	ILE
1	A	320	THR
1	A	338	GLN
1	A	344	LEU
1	A	367	THR
1	A	406	ASN
1	A	432	MET
1	A	435	LEU
1	A	438	GLU
1	A	441	ARG
1	A	448	THR
1	A	463	GLN

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	45	GLN
1	A	49	ASN
1	A	75	HIS
1	A	99	HIS
1	A	122	HIS
1	A	124	HIS
1	A	126	ASN
1	A	167	ASN
1	A	190	ASN
1	A	214	HIS
1	A	228	GLN
1	A	243	HIS
1	A	244	ASN
1	A	317	ASN
1	A	333	HIS
1	A	356	GLN
1	A	405	GLN
1	A	406	ASN
1	A	410	HIS
1	A	463	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](#)

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	378/467 (80%)	0.12	25 (6%) 24 18	20, 38, 72, 91	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	368	GLY	6.6
1	A	435	LEU	5.1
1	A	11	PRO	4.5
1	A	100	GLU	4.1
1	A	434	SER	4.1
1	A	317	ASN	3.9
1	A	438	GLU	3.4
1	A	432	MET	3.2
1	A	462	TYR	3.1
1	A	433	LEU	3.0
1	A	356	GLN	2.9
1	A	436	PRO	2.9
1	A	367	THR	2.8
1	A	437	GLU	2.7
1	A	14	VAL	2.6
1	A	50	ALA	2.4
1	A	316	PRO	2.2
1	A	107	ASN	2.2
1	A	12	PRO	2.2
1	A	159	ILE	2.1
1	A	298	ILE	2.1
1	A	299	GLU	2.1
1	A	314	ILE	2.1
1	A	191	TYR	2.1
1	A	313	PHE	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](#)

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($\text{\AA}^2$)	Q<0.9
2	MN	A	502	1/1	0.96	0.06	32,32,32,32	0
2	MN	A	501	1/1	0.97	0.08	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.