



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:31 PM UTC

PDB ID : 6IOE / pdb_00006ioe
Title : Crystal structure of the CysR-CTLD2 fragment of human MR at basic pH (pH 8.5)
Authors : Hu, Z.; He, Y.
Deposited on : 2018-10-30
Resolution : 2.90 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

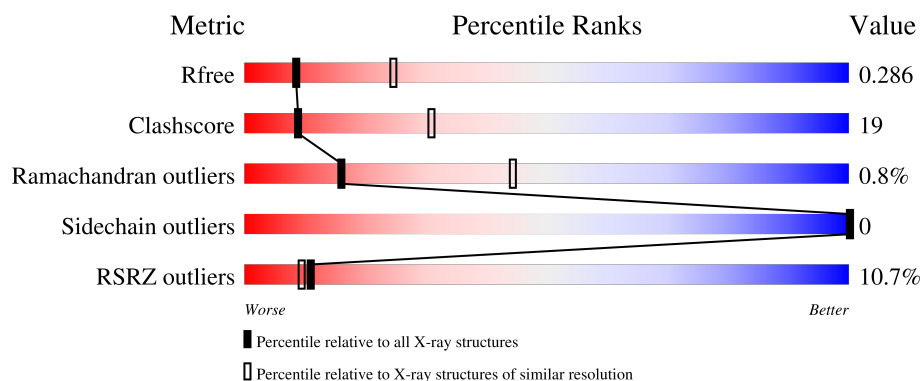
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	475	7% 67% 27% • •
1	B	475	14% 59% 33% • 6%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7242 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 Macrophage mannose receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	454	Total	C	N	O	S	0	0	0
			3654	2317	618	693	26			
1	B	446	Total	C	N	O	S	0	0	0
			3588	2282	603	677	26			

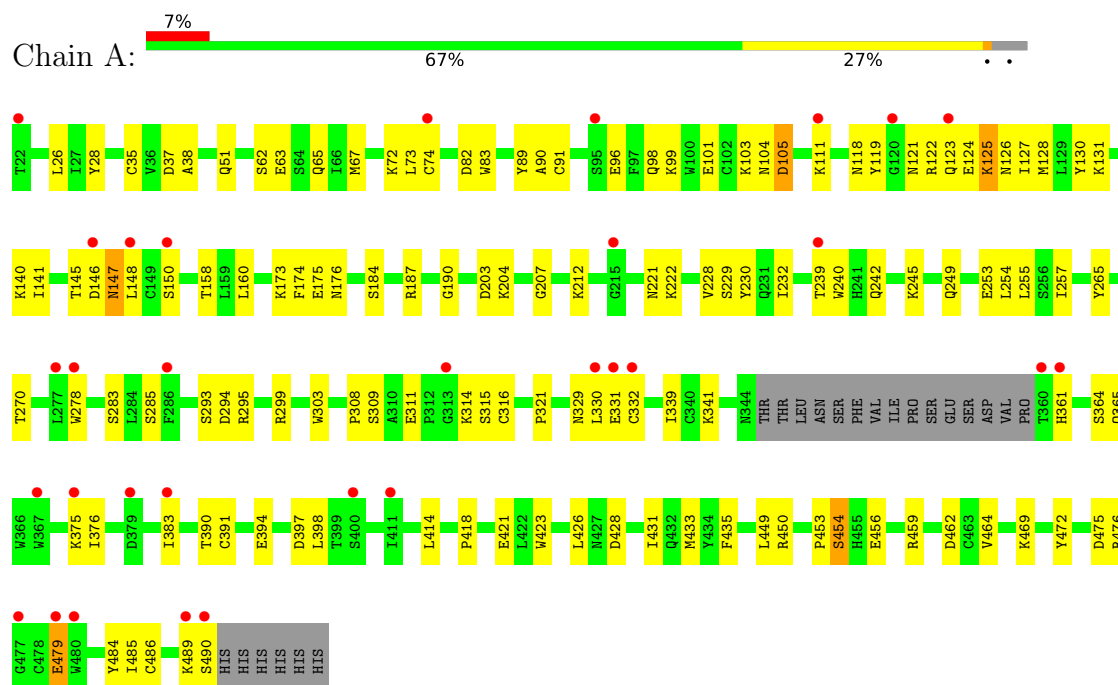
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	491	HIS	-	expression tag	UNP P22897
A	492	HIS	-	expression tag	UNP P22897
A	493	HIS	-	expression tag	UNP P22897
A	494	HIS	-	expression tag	UNP P22897
A	495	HIS	-	expression tag	UNP P22897
A	496	HIS	-	expression tag	UNP P22897
B	491	HIS	-	expression tag	UNP P22897
B	492	HIS	-	expression tag	UNP P22897
B	493	HIS	-	expression tag	UNP P22897
B	494	HIS	-	expression tag	UNP P22897
B	495	HIS	-	expression tag	UNP P22897
B	496	HIS	-	expression tag	UNP P22897

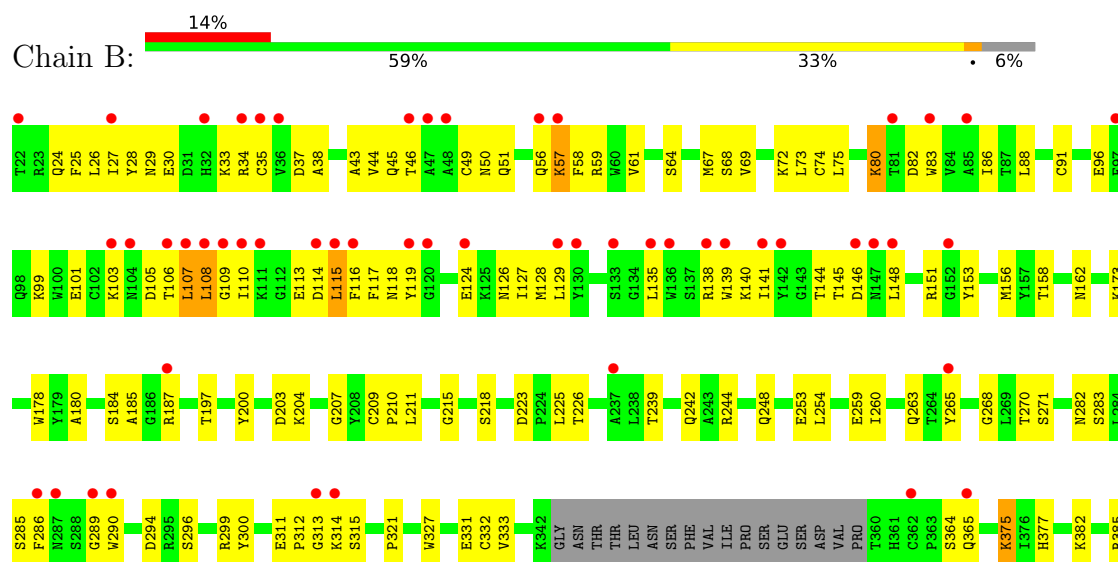
3 Residue-property plots

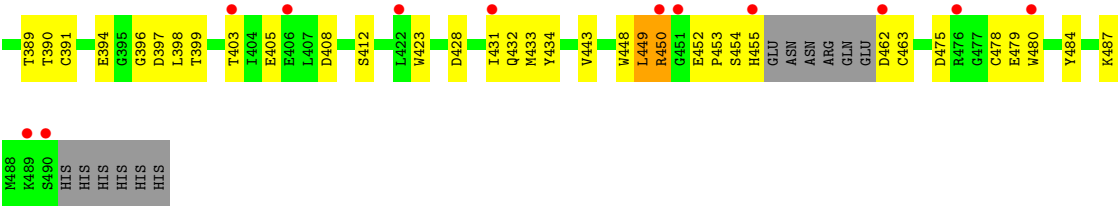
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Macrophage mannose receptor 1



• Molecule 1: Macrophage mannose receptor 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	241.43Å 72.24Å 109.53Å 90.00° 114.51° 90.00°	Depositor
Resolution (Å)	29.85 – 2.90 29.85 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.0 (29.85-2.90) 97.0 (29.85-2.90)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.236 , 0.284 0.237 , 0.286	Depositor DCC
R_{free} test set	2012 reflections (5.23%)	wwPDB-VP
Wilson B-factor (Å ²)	55.5	Xtriage
Anisotropy	1.123	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.025 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7242	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.85% 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	0.25	0/3756	0.60	4/5090 (0.1%)
1	B	0.37	2/3689 (0.1%)	0.73	11/4999 (0.2%)
All	All	0.32	2/7445 (0.0%)	0.67	15/10089 (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	B	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	225	LEU	C-O	6.08	1.30	1.24
1	B	80	LYS	CE-NZ	5.66	1.66	1.49

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	108	LEU	CB-CG-CD1	-9.63	81.82	110.70
1	B	375	LYS	CG-CD-CE	-8.74	91.20	111.30
1	B	50	ASN	CA-C-N	-8.42	108.67	122.65
1	B	50	ASN	C-N-CA	-8.42	108.67	122.65
1	A	331	GLU	CA-CB-CG	-7.50	99.09	114.10
1	B	110	ILE	N-CA-C	6.59	118.56	109.80
1	B	108	LEU	CB-CG-CD2	6.37	129.80	110.70
1	A	96	GLU	CA-CB-CG	5.92	125.94	114.10
1	B	449	LEU	CA-C-N	-5.90	112.91	122.82
1	B	449	LEU	C-N-CA	-5.90	112.91	122.82
1	A	479	GLU	CA-C-N	5.86	134.30	122.70
1	A	479	GLU	C-N-CA	5.86	134.30	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	110	ILE	CA-C-N	5.50	128.69	120.87
1	B	110	ILE	C-N-CA	5.50	128.69	120.87
1	B	57	LYS	CD-CE-NZ	-5.09	95.60	111.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	103	LYS	Peptide
1	B	107	LEU	Peptide
1	B	115	LEU	Peptide
1	B	450	ARG	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	3654	0	3463	113	1
1	B	3588	0	3406	151	1
All	All	7242	0	6869	262	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 19.

All (262) 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:314:LYS:NZ	1:A:316:CYS:H	1.24	1.36
1:A:314:LYS:HZ2	1:A:316:CYS:N	1.46	1.13
1:B:289:GLY:HA2	1:B:450:ARG:HB3	1.31	1.10
1:B:26:LEU:HD12	1:B:51:GLN:HG2	1.24	1.09
1:A:314:LYS:HE2	1:A:330:LEU:C	1.81	1.05
1:A:140:LYS:NZ	1:A:146:ASP:O	1.90	1.04
1:A:140:LYS:NZ	1:A:146:ASP:C	2.19	1.01
1:A:140:LYS:HZ3	1:A:146:ASP:C	1.74	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:LYS:HZ3	1:A:315:SER:H	1.13	0.92
1:A:433:MET:HE1	1:A:456:GLU:HB3	1.51	0.91
1:A:314:LYS:NZ	1:A:316:CYS:N	2.08	0.91
1:A:314:LYS:NZ	1:A:315:SER:H	1.70	0.90
1:B:450:ARG:HA	1:B:452:GLU:H	1.37	0.89
1:B:26:LEU:HD12	1:B:51:GLN:CG	2.03	0.88
1:A:253:GLU:HG2	1:A:294:ASP:HB3	1.56	0.86
1:B:311:GLU:HB2	1:B:314:LYS:HG2	1.58	0.83
1:A:314:LYS:HE2	1:A:330:LEU:O	1.77	0.83
1:A:221:ASN:ND2	1:A:265:TYR:OH	2.11	0.82
1:B:61:VAL:HG21	1:B:67:MET:HE3	1.61	0.82
1:B:26:LEU:CD1	1:B:51:GLN:HG2	2.09	0.81
1:A:459:ARG:NH1	1:A:462:ASP:OD1	2.18	0.77
1:A:391:CYS:HG	1:A:486:CYS:CB	1.97	0.77
1:B:108:LEU:HD13	1:B:148:LEU:HD11	1.67	0.76
1:A:391:CYS:SG	1:A:486:CYS:HB3	2.25	0.75
1:A:314:LYS:HZ1	1:A:316:CYS:H	1.33	0.74
1:A:103:LYS:O	1:A:105:ASP:N	2.20	0.74
1:A:239:THR:OG1	1:A:332:CYS:O	2.06	0.73
1:A:140:LYS:HE2	1:A:145:THR:C	2.13	0.73
1:B:59:ARG:NH1	1:B:180:ALA:O	2.22	0.72
1:B:197:THR:HG21	1:B:203:ASP:HB2	1.71	0.72
1:B:45:GLN:HA	1:B:127:ILE:HG12	1.72	0.71
1:A:314:LYS:NZ	1:A:315:SER:N	2.39	0.71
1:A:314:LYS:HZ2	1:A:316:CYS:H	0.71	0.71
1:B:283:SER:HB3	1:B:450:ARG:HE	1.56	0.70
1:A:221:ASN:O	1:A:230:TYR:N	2.27	0.67
1:B:399:THR:HA	1:B:487:LYS:HB2	1.76	0.67
1:B:449:LEU:O	1:B:452:GLU:HB2	1.94	0.67
1:B:299:ARG:HG3	1:B:443:VAL:HG11	1.77	0.67
1:A:314:LYS:HE3	1:A:329:ASN:HB2	1.78	0.66
1:A:423:TRP:HB2	1:A:484:TYR:HB3	1.77	0.66
1:B:408:ASP:O	1:B:412:SER:OG	2.13	0.65
1:A:426:LEU:HD23	1:A:464:VAL:HG21	1.78	0.65
1:A:391:CYS:SG	1:A:486:CYS:CB	2.83	0.65
1:B:223:ASP:HB3	1:B:265:TYR:CZ	2.31	0.64
1:A:140:LYS:HZ3	1:A:147:ASN:CA	2.11	0.64
1:B:239:THR:HG23	1:B:242:GLN:H	1.63	0.63
1:A:295:ARG:HG2	1:A:295:ARG:HH11	1.63	0.63
1:B:390:THR:O	1:B:394:GLU:HG3	1.99	0.62
1:B:223:ASP:OD1	1:B:226:THR:OG1	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:428:ASP:OD1	1:B:462:ASP:HB3	2.00	0.62
1:A:83:TRP:CD1	1:A:128:MET:HE1	2.35	0.62
1:B:364:SER:O	1:B:365:GLN:HB2	2.00	0.62
1:B:73:LEU:HB2	1:B:88:LEU:HB3	1.82	0.61
1:A:314:LYS:HZ3	1:A:315:SER:N	1.93	0.61
1:B:140:LYS:HE2	1:B:146:ASP:C	2.26	0.61
1:B:480:TRP:N	1:B:480:TRP:CD1	2.69	0.61
1:B:144:THR:HG22	1:B:145:THR:H	1.66	0.61
1:A:140:LYS:HZ3	1:A:147:ASN:N	1.98	0.60
1:B:428:ASP:HB2	1:B:431:ILE:O	2.01	0.60
1:B:162:ASN:ND2	1:B:204:LYS:O	2.34	0.60
1:B:223:ASP:HB3	1:B:265:TYR:CE1	2.35	0.60
1:B:215:GLY:O	1:B:218:SER:OG	2.19	0.60
1:B:289:GLY:HA2	1:B:450:ARG:CB	2.20	0.60
1:B:45:GLN:HB3	1:B:126:ASN:HA	1.83	0.60
1:B:184:SER:C	1:B:187:ARG:HG3	2.27	0.60
1:A:160:LEU:HD13	1:B:311:GLU:HG2	1.84	0.59
1:A:239:THR:HG23	1:A:242:GLN:H	1.68	0.59
1:A:82:ASP:OD1	1:A:83:TRP:N	2.36	0.59
1:A:140:LYS:HZ1	1:A:146:ASP:C	1.96	0.59
1:B:259:GLU:OE2	1:B:299:ARG:NH1	2.35	0.59
1:B:140:LYS:HZ3	1:B:141:ILE:H	1.51	0.59
1:B:268:GLY:O	1:B:271:SER:OG	2.21	0.58
1:B:331:GLU:CD	1:B:331:GLU:H	2.11	0.58
1:A:126:ASN:ND2	1:A:127:ILE:O	2.36	0.58
1:B:423:TRP:HB2	1:B:484:TYR:HB3	1.86	0.57
1:B:140:LYS:HZ3	1:B:141:ILE:N	2.03	0.57
1:A:454:SER:OG	1:A:475:ASP:O	2.21	0.57
1:A:222:LYS:HA	1:A:228:VAL:O	2.04	0.57
1:B:286:PHE:HA	1:B:450:ARG:HH22	1.69	0.57
1:A:418:PRO:HB3	1:A:469:LYS:HA	1.85	0.57
1:B:44:VAL:HG11	1:B:88:LEU:HG	1.86	0.57
1:B:28:TYR:O	1:B:140:LYS:N	2.38	0.56
1:B:67:MET:HA	1:B:74:CYS:HA	1.86	0.56
1:B:107:LEU:HD11	1:B:116:PHE:HB3	1.87	0.56
1:A:212:LYS:HD2	1:A:212:LYS:N	2.22	0.55
1:A:140:LYS:HE2	1:A:146:ASP:N	2.22	0.55
1:B:253:GLU:HG2	1:B:294:ASP:HB3	1.88	0.55
1:B:244:ARG:O	1:B:248:GLN:HG3	2.07	0.55
1:B:115:LEU:HB2	1:B:129:LEU:HD21	1.89	0.55
1:B:314:LYS:HD2	1:B:331:GLU:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:433:MET:HE1	1:B:455:HIS:CD2	2.43	0.54
1:B:107:LEU:HG	1:B:108:LEU:C	2.33	0.54
1:A:283:SER:O	1:A:450:ARG:NH1	2.36	0.54
1:B:450:ARG:CA	1:B:452:GLU:H	2.15	0.54
1:A:270:THR:HB	1:A:321:PRO:HB3	1.91	0.53
1:A:67:MET:SD	1:A:72:LYS:HA	2.48	0.53
1:A:184:SER:O	1:A:184:SER:OG	2.27	0.53
1:A:390:THR:O	1:A:394:GLU:HG3	2.08	0.53
1:B:28:TYR:CE2	1:B:49:CYS:SG	3.02	0.53
1:B:118:ASN:OD1	1:B:119:TYR:N	2.41	0.53
1:B:454:SER:N	1:B:475:ASP:OD1	2.41	0.52
1:A:397:ASP:OD1	1:A:398:LEU:N	2.43	0.52
1:B:209:CYS:HB2	1:B:211:LEU:HD12	1.91	0.52
1:B:108:LEU:CD1	1:B:148:LEU:HD11	2.39	0.52
1:B:200:TYR:O	1:B:204:LYS:HD3	2.10	0.52
1:A:383:ILE:HD12	1:A:383:ILE:H	1.75	0.52
1:B:141:ILE:HD11	1:B:148:LEU:HA	1.92	0.52
1:B:432:GLN:HG2	1:B:433:MET:HE3	1.92	0.52
1:A:67:MET:HA	1:A:74:CYS:HA	1.92	0.52
1:B:117:PHE:HD2	1:B:139:TRP:CZ3	2.28	0.52
1:A:255:LEU:HD21	1:A:257:ILE:HD11	1.92	0.51
1:B:397:ASP:OD1	1:B:398:LEU:N	2.42	0.51
1:A:376:ILE:HG12	1:A:485:ILE:HG12	1.92	0.51
1:B:49:CYS:HA	1:B:56:GLN:HE22	1.76	0.51
1:B:290:TRP:HD1	1:B:453:PRO:HD2	1.75	0.51
1:B:382:LYS:NZ	1:B:484:TYR:OH	2.43	0.51
1:B:289:GLY:HA3	1:B:448:TRP:HB2	1.92	0.51
1:A:73:LEU:HD23	1:A:90:ALA:HA	1.92	0.51
1:A:245:LYS:O	1:A:249:GLN:HG2	2.11	0.51
1:A:314:LYS:HZ2	1:A:315:SER:N	2.09	0.50
1:B:101:GLU:N	1:B:109:GLY:O	2.44	0.50
1:B:375:LYS:HD3	1:B:377:HIS:CD2	2.46	0.50
1:A:173:LYS:HD3	1:A:176:ASN:HA	1.94	0.50
1:B:480:TRP:N	1:B:480:TRP:HD1	2.09	0.50
1:B:294:ASP:OD1	1:B:296:SER:OG	2.30	0.50
1:A:428:ASP:HB2	1:A:431:ILE:O	2.12	0.50
1:B:283:SER:HB3	1:B:450:ARG:NE	2.26	0.50
1:A:26:LEU:HD12	1:A:51:GLN:HG2	1.94	0.49
1:A:253:GLU:HG3	1:A:254:LEU:N	2.27	0.49
1:A:454:SER:N	1:A:475:ASP:OD1	2.44	0.49
1:A:158:THR:HB	1:A:207:GLY:HA3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:ILE:CD1	1:B:58:PHE:HB2	2.43	0.49
1:B:26:LEU:HD23	1:B:26:LEU:HA	1.54	0.49
1:B:37:ASP:OD2	1:B:38:ALA:N	2.45	0.49
1:B:72:LYS:HD3	1:B:178:TRP:HB2	1.93	0.49
1:A:314:LYS:HG3	1:A:329:ASN:HB2	1.95	0.48
1:A:376:ILE:HG13	1:A:485:ILE:HG23	1.95	0.48
1:B:239:THR:OG1	1:B:332:CYS:O	2.32	0.48
1:A:140:LYS:HZ3	1:A:147:ASN:HA	1.76	0.48
1:A:364:SER:O	1:A:365:GLN:HB2	2.14	0.48
1:B:24:GLN:O	1:B:151:ARG:NH1	2.47	0.48
1:B:253:GLU:HG3	1:B:254:LEU:N	2.27	0.48
1:A:285:SER:O	1:A:450:ARG:NH1	2.47	0.48
1:A:309:SER:OG	1:A:311:GLU:HG3	2.14	0.48
1:A:124:GLU:O	1:A:126:ASN:N	2.47	0.48
1:A:278:TRP:HH2	1:A:330:LEU:O	1.97	0.48
1:B:28:TYR:N	1:B:139:TRP:HB3	2.28	0.48
1:A:65:GLN:HE21	1:A:99:LYS:NZ	2.12	0.48
1:A:28:TYR:HD1	1:A:35:CYS:SG	2.37	0.47
1:B:184:SER:CA	1:B:187:ARG:HG3	2.44	0.47
1:B:30:GLU:HA	1:B:33:LYS:HD3	1.97	0.47
1:B:107:LEU:HB2	1:B:138:ARG:NH1	2.28	0.47
1:A:454:SER:OG	1:A:475:ASP:OD1	2.31	0.47
1:B:108:LEU:O	1:B:108:LEU:HG	2.14	0.47
1:A:449:LEU:HD12	1:A:472:TYR:HB2	1.95	0.47
1:B:67:MET:HG2	1:B:68:SER:N	2.29	0.47
1:B:187:ARG:HA	1:B:187:ARG:NE	2.29	0.47
1:B:28:TYR:HA	1:B:139:TRP:HD1	1.80	0.47
1:B:29:ASN:ND2	1:B:135:LEU:O	2.46	0.47
1:A:314:LYS:CE	1:A:330:LEU:N	2.78	0.46
1:B:43:ALA:HA	1:B:126:ASN:HD21	1.80	0.46
1:B:80:LYS:NZ	1:B:96:GLU:O	2.48	0.46
1:B:244:ARG:HA	1:B:254:LEU:HG	1.95	0.46
1:B:452:GLU:OE1	1:B:454:SER:OG	2.32	0.46
1:B:283:SER:O	1:B:450:ARG:NH2	2.48	0.46
1:B:58:PHE:CE2	1:B:68:SER:HB2	2.49	0.46
1:A:83:TRP:HZ2	1:A:121:ASN:HB3	1.80	0.46
1:B:58:PHE:CE2	1:B:75:LEU:HD21	2.49	0.46
1:B:259:GLU:OE2	1:B:434:TYR:OH	2.33	0.46
1:B:35:CYS:O	1:B:46:THR:HA	2.16	0.46
1:B:148:LEU:HD23	1:B:148:LEU:H	1.80	0.46
1:A:221:ASN:O	1:A:229:SER:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:VAL:HG23	1:B:180:ALA:HB3	1.97	0.46
1:B:187:ARG:HA	1:B:187:ARG:HE	1.80	0.46
1:B:64:SER:HB3	1:B:99:LYS:HB3	1.98	0.46
1:B:239:THR:OG1	1:B:333:VAL:HA	2.16	0.46
1:B:397:ASP:O	1:B:487:LYS:N	2.49	0.46
1:A:453:PRO:HB2	1:A:456:GLU:OE2	2.16	0.46
1:A:421:GLU:OE2	1:A:476:ARG:NH1	2.49	0.45
1:B:108:LEU:HD13	1:B:148:LEU:CD1	2.43	0.45
1:B:270:THR:HB	1:B:321:PRO:HB3	1.98	0.45
1:B:285:SER:O	1:B:450:ARG:NH1	2.49	0.45
1:A:62:SER:OG	1:A:63:GLU:N	2.48	0.45
1:A:118:ASN:ND2	1:A:119:TYR:O	2.48	0.45
1:B:144:THR:HG22	1:B:145:THR:N	2.30	0.45
1:B:185:ALA:C	1:B:187:ARG:HG2	2.42	0.45
1:A:489:LYS:HD2	1:A:490:SER:H	1.81	0.45
1:B:74:CYS:N	1:B:91:CYS:SG	2.90	0.45
1:B:86:ILE:HG12	1:B:128:MET:HA	1.98	0.45
1:B:107:LEU:HG	1:B:108:LEU:N	2.32	0.45
1:A:37:ASP:OD2	1:A:38:ALA:N	2.50	0.45
1:A:299:ARG:NH2	1:A:435:PHE:O	2.46	0.45
1:B:82:ASP:OD1	1:B:83:TRP:N	2.50	0.45
1:B:185:ALA:O	1:B:187:ARG:HG2	2.17	0.45
1:A:376:ILE:HD12	1:A:414:LEU:HB2	1.99	0.44
1:B:51:GLN:O	1:B:57:LYS:NZ	2.48	0.44
1:B:286:PHE:C	1:B:450:ARG:HH12	2.24	0.44
1:A:89:TYR:HB2	1:A:98:GLN:NE2	2.33	0.44
1:A:303:TRP:CE2	1:A:308:PRO:HG3	2.52	0.44
1:B:105:ASP:HB3	1:B:138:ARG:HH12	1.83	0.44
1:B:403:THR:HG22	1:B:405:GLU:H	1.82	0.44
1:A:184:SER:HB2	1:A:190:GLY:HA2	2.00	0.43
1:A:314:LYS:HE3	1:A:329:ASN:C	2.42	0.43
1:A:82:ASP:OD1	1:A:130:TYR:HA	2.18	0.43
1:B:67:MET:SD	1:B:72:LYS:HA	2.58	0.43
1:B:452:GLU:HA	1:B:453:PRO:C	2.42	0.43
1:B:463:CYS:N	1:B:478:CYS:SG	2.91	0.43
1:A:232:ILE:HD13	1:A:339:ILE:HG12	2.00	0.43
1:A:83:TRP:CZ2	1:A:121:ASN:HB3	2.54	0.43
1:A:212:LYS:HD2	1:A:212:LYS:H	1.83	0.43
1:B:146:ASP:OD1	1:B:146:ASP:N	2.51	0.43
1:B:158:THR:HB	1:B:207:GLY:HA3	2.00	0.43
1:A:314:LYS:HE3	1:A:316:CYS:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:LYS:HE2	1:A:330:LEU:N	2.34	0.43
1:B:106:THR:HG23	1:B:138:ARG:HB3	2.01	0.43
1:A:174:PHE:CE1	1:A:175:GLU:HG3	2.54	0.42
1:B:28:TYR:H	1:B:140:LYS:H	1.68	0.42
1:B:260:ILE:HG13	1:B:434:TYR:CE2	2.54	0.42
1:B:30:GLU:O	1:B:33:LYS:HG2	2.19	0.42
1:B:80:LYS:HD2	1:B:113:GLU:OE1	2.19	0.42
1:B:126:ASN:O	1:B:128:MET:HE2	2.19	0.42
1:B:263:GLN:HE22	1:B:327:TRP:HE1	1.67	0.42
1:B:313:GLY:C	1:B:314:LYS:HD3	2.45	0.42
1:A:74:CYS:N	1:A:91:CYS:SG	2.92	0.42
1:B:28:TYR:H	1:B:139:TRP:HB3	1.84	0.42
1:B:140:LYS:HE3	1:B:145:THR:C	2.44	0.42
1:A:240:TRP:HD1	1:A:278:TRP:CE3	2.38	0.42
1:A:365:GLN:HB3	1:A:375:LYS:HZ1	1.85	0.42
1:A:140:LYS:HA	1:A:140:LYS:HD2	1.55	0.41
1:B:28:TYR:CZ	1:B:33:LYS:HD2	2.55	0.41
1:B:263:GLN:HE21	1:B:300:TYR:HE1	1.68	0.41
1:B:34:ARG:NH2	1:B:46:THR:O	2.53	0.41
1:B:391:CYS:O	1:B:396:GLY:N	2.51	0.41
1:A:82:ASP:OD2	1:A:131:LYS:HE3	2.20	0.41
1:B:118:ASN:HA	1:B:139:TRP:HH2	1.85	0.41
1:A:203:ASP:C	1:A:204:LYS:HG2	2.45	0.41
1:B:128:MET:HE3	1:B:128:MET:HB2	1.73	0.41
1:A:228:VAL:HG11	1:A:341:LYS:HE3	2.02	0.41
1:B:209:CYS:CB	1:B:211:LEU:HD12	2.51	0.41
1:A:141:ILE:HG13	1:A:148:LEU:HG	2.03	0.41
1:A:329:ASN:OD1	1:A:329:ASN:N	2.48	0.41
1:B:403:THR:HG22	1:B:405:GLU:N	2.36	0.41
1:B:385:ARG:O	1:B:389:THR:HG23	2.21	0.41
1:B:80:LYS:CD	1:B:113:GLU:OE1	2.69	0.41
1:B:119:TYR:HE2	1:B:124:GLU:O	2.04	0.41
1:B:282:ASN:HA	1:B:315:SER:O	2.21	0.41
1:B:25:PHE:CE2	1:B:59:ARG:HA	2.56	0.41
1:A:314:LYS:CG	1:A:329:ASN:HB2	2.51	0.40
1:B:144:THR:HG22	1:B:146:ASP:H	1.86	0.40
1:B:156:MET:HG2	1:B:210:PRO:HD2	2.03	0.40
1:A:253:GLU:OE1	1:A:293:SER:OG	2.24	0.40
1:B:450:ARG:HH11	1:B:450:ARG:HG2	1.85	0.40
1:B:479:GLU:C	1:B:480:TRP:CD1	3.00	0.40
1:A:121:ASN:ND2	1:A:123:GLN:HB2	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:ARG:NH2	1:B:312:PRO:HG2	2.37	0.40
1:A:26:LEU:HD23	1:A:26:LEU:HA	1.70	0.40
1:A:101:GLU:OE2	1:A:111:LYS:HG3	2.20	0.40
1:A:122:ARG:HH11	1:A:125:LYS:NZ	2.20	0.40
1:A:147:ASN:O	1:A:150:SER:N	2.51	0.40
1:A:140:LYS:HD2	1:A:141:ILE:H	1.87	0.40
1:B:114:ASP:C	1:B:115:LEU:HD23	2.46	0.40
1:B:173:LYS:HB2	1:B:178:TRP:CZ3	2.57	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:A:479:GLU:OE2	1:B:375:LYS:NZ[1_545]	2.19	0.01

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	450/475 (95%)	411 (91%)	33 (7%)	6 (1%)	9	32
1	B	440/475 (93%)	401 (91%)	38 (9%)	1 (0%)	43	72
All	All	890/950 (94%)	812 (91%)	71 (8%)	7 (1%)	16	44

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	105	ASP
1	A	125	LYS
1	A	454	SER
1	B	153	TYR

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Mol	Chain	Res	Type
1	A	361	HIS
1	A	147	ASN

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	397/418 (95%)	397 (100%)	0	100	100
1	B	390/418 (93%)	390 (100%)	0	100	100
All	All	787/836 (94%)	787 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	123	GLN
1	A	221	ASN
1	A	241	HIS
1	A	344	ASN
1	A	413	GLN
1	B	24	GLN
1	B	29	ASN
1	B	56	GLN
1	B	121	ASN
1	B	126	ASN
1	B	231	GLN
1	B	250	GLN
1	B	263	GLN
1	B	287	ASN
1	B	455	HIS

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	454/475 (95%)	0.68	31 (6%) 23 18	38, 59, 87, 128	0
1	B	446/475 (93%)	1.04	65 (14%) 6 4	36, 65, 125, 147	0
All	All	900/950 (94%)	0.86	96 (10%) 11 9	36, 62, 108, 147	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	108	LEU	6.2
1	B	135	LEU	5.9
1	B	455	HIS	5.6
1	B	35	CYS	5.1
1	B	480	TRP	4.9
1	B	120	GLY	4.6
1	A	330	LEU	4.5
1	B	107	LEU	4.5
1	B	106	THR	4.2
1	B	142	TYR	4.1
1	B	313	GLY	3.8
1	B	103	LYS	3.6
1	B	119	TYR	3.6
1	B	139	TRP	3.6
1	B	148	LEU	3.6
1	B	104	ASN	3.5
1	B	47	ALA	3.5
1	A	360	THR	3.5
1	B	115	LEU	3.3
1	B	490	SER	3.2
1	B	451	GLY	3.1
1	B	48	ALA	3.1
1	A	146	ASP	3.1
1	B	286	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	95	SER	3.0
1	B	146	ASP	3.0
1	B	365	GLN	2.9
1	A	331	GLU	2.9
1	B	489	LYS	2.9
1	A	400	SER	2.9
1	A	479	GLU	2.8
1	B	187	ARG	2.8
1	A	22	THR	2.8
1	B	85	ALA	2.8
1	B	116	PHE	2.8
1	A	379	ASP	2.8
1	B	290	TRP	2.8
1	B	289	GLY	2.8
1	B	138	ARG	2.7
1	A	332	CYS	2.7
1	A	490	SER	2.7
1	B	109	GLY	2.7
1	A	120	GLY	2.7
1	B	114	ASP	2.7
1	B	314	LYS	2.7
1	B	56	GLN	2.7
1	A	313	GLY	2.7
1	A	148	LEU	2.7
1	A	215	GLY	2.6
1	B	36	VAL	2.6
1	A	383	ILE	2.6
1	A	286	PHE	2.6
1	B	32	HIS	2.6
1	B	129	LEU	2.6
1	B	141	ILE	2.6
1	B	287	ASN	2.6
1	B	403	THR	2.5
1	B	136	TRP	2.5
1	A	150	SER	2.5
1	B	133	SER	2.5
1	B	27	ILE	2.5
1	B	450	ARG	2.5
1	A	367	TRP	2.4
1	B	83	TRP	2.4
1	B	22	THR	2.4
1	A	375	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	265	TYR	2.4
1	A	411	ILE	2.4
1	A	74	CYS	2.4
1	B	147	ASN	2.3
1	B	152	GLY	2.3
1	B	46	THR	2.3
1	B	130	TYR	2.3
1	A	111	LYS	2.3
1	B	431	ILE	2.3
1	B	110	ILE	2.3
1	B	81	THR	2.3
1	B	124	GLU	2.3
1	B	476	ARG	2.2
1	B	97	PHE	2.2
1	B	237	ALA	2.2
1	A	480	TRP	2.2
1	B	406	GLU	2.1
1	A	123	GLN	2.1
1	A	489	LYS	2.1
1	B	57	LYS	2.1
1	B	111	LYS	2.1
1	B	422	LEU	2.1
1	A	239	THR	2.1
1	B	362	CYS	2.1
1	A	477	GLY	2.1
1	A	277	LEU	2.0
1	B	34	ARG	2.0
1	B	462	ASP	2.0
1	A	361	HIS	2.0
1	A	278	TRP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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.