



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:44 PM UTC

PDB ID : 3VWA / pdb_00003vwa
Title : Crystal structure of Cex1p
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Deposited on : 2012-08-13
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
Mogul	:	2022.3.0, CSD as543be (2022)
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

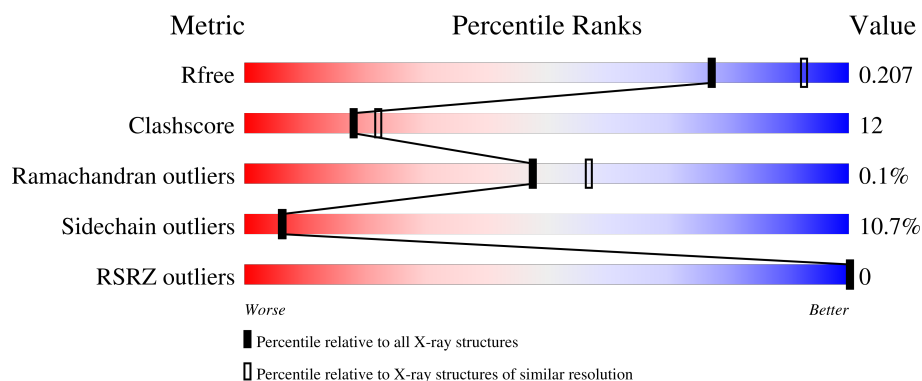
1 Overall quality at a glance

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	560	 63% 27% • 5%
1	B	560	 69% 21% • • 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8822 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 Cytoplasmic export protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	531	Total	C	N	O	S	Se	0	0	0
			4184	2702	687	780	8	7			
1	B	529	Total	C	N	O	S	Se	0	0	0
			4180	2699	689	777	8	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	PRO	-	expression tag	UNP Q12453
A	6	LEU	-	expression tag	UNP Q12453
A	7	HIS	-	expression tag	UNP Q12453
A	8	MSE	-	expression tag	UNP Q12453
B	5	PRO	-	expression tag	UNP Q12453
B	6	LEU	-	expression tag	UNP Q12453
B	7	HIS	-	expression tag	UNP Q12453
B	8	MSE	-	expression tag	UNP Q12453

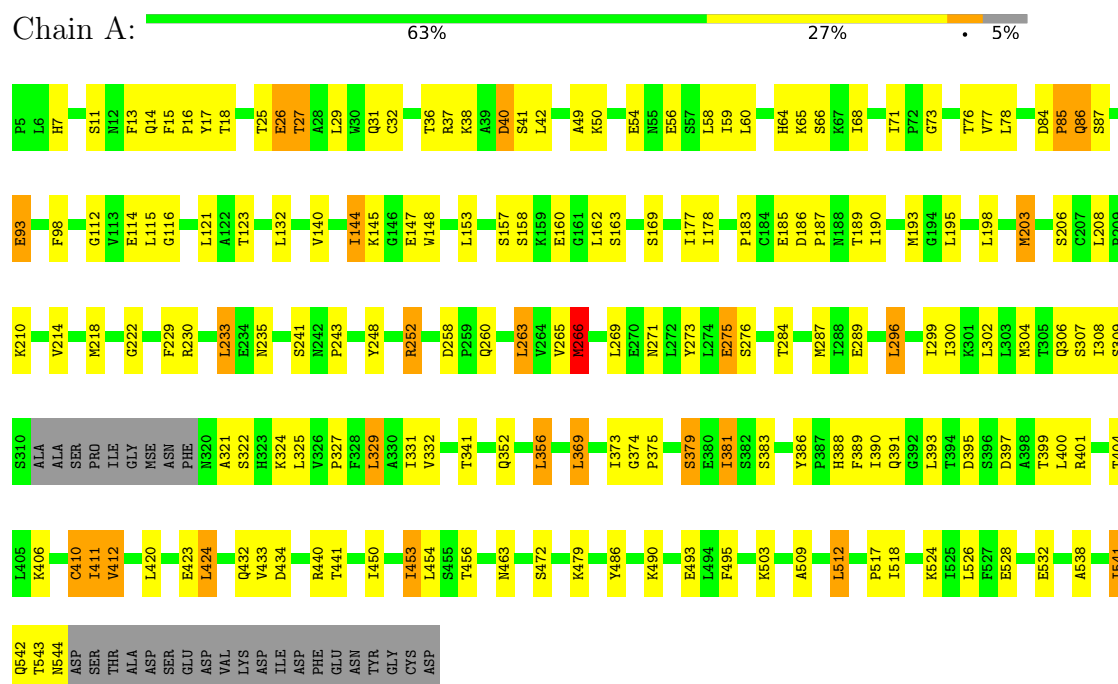
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	229	Total	O	0	0
			229	229		
2	B	229	Total	O	0	0
			229	229		

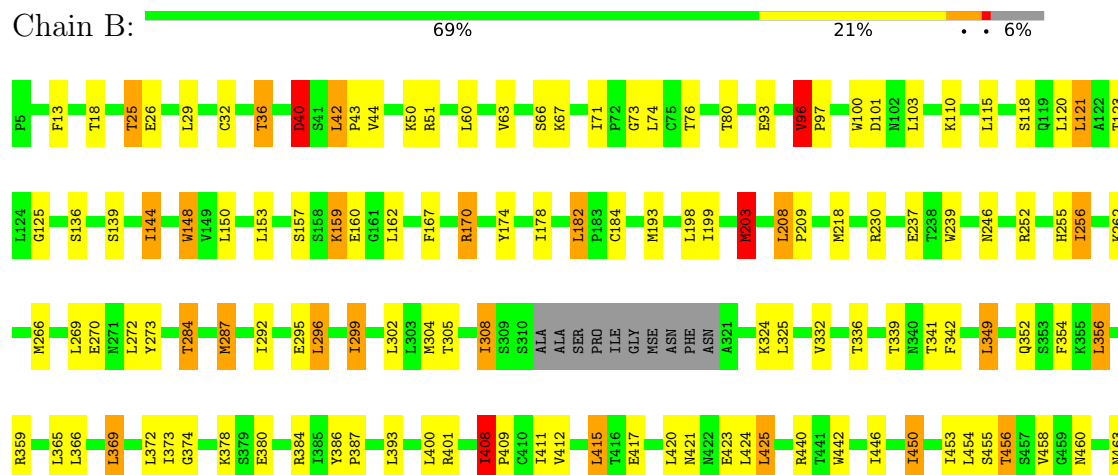
3 Residue-property plots

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: Cytoplasmic export protein 1



• Molecule 1: Cytoplasmic export protein 1



K474	D475	L488	K503	T506	L512	L513	D514	I518	V519	R520	G521	R522	K523	K524	I525	L526	I541	G542	T543	ASN	ASP	SER	THR	ALA	ASP	ASP	SER	GLU	ASP	VAL	LYS	ASP	ILE	ASP	PHE	GLU	ASN	TYR	GLY	CYS	ASP
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	34.06Å 182.22Å 112.44Å 90.00° 98.69° 90.00°	Depositor
Resolution (Å)	45.56 – 2.20 45.56 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.7 (45.56-2.20) 98.8 (45.56-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.56 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.166 , 0.207 0.166 , 0.207	Depositor DCC
R_{free} test set	6784 reflections (7.31%)	wwPDB-VP
Wilson B-factor (Å ²)	32.5	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 23.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.430 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.575 for H, K, L 0.425 for -H, -K, H+L	Depositor
Outliers	0 of 67710 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8822	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.95% 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	1.04	1/4256 (0.0%)	1.17	13/5763 (0.2%)
1	B	1.07	8/4252 (0.2%)	1.18	14/5756 (0.2%)
All	All	1.06	9/8508 (0.1%)	1.18	27/11519 (0.2%)

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	1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	96	VAL	CA-CB	8.46	1.60	1.54
1	B	118	SER	N-CA	6.31	1.53	1.46
1	B	110	LYS	C-O	-6.06	1.17	1.24
1	B	408	ILE	CA-CB	6.01	1.57	1.54
1	A	495	PHE	C-O	5.72	1.30	1.24
1	B	125	GLY	N-CA	5.38	1.52	1.45
1	B	239	TRP	CD2-CE2	5.37	1.50	1.41
1	B	148	TRP	CA-C	-5.20	1.46	1.52
1	B	100	TRP	CD2-CE2	5.14	1.50	1.41

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	356	LEU	CA-C-N	-9.70	110.94	120.83
1	A	356	LEU	C-N-CA	-9.70	110.94	120.83
1	B	218	MSE	CA-CB-CG	-7.09	99.91	114.10
1	B	96	VAL	CA-C-O	6.86	123.76	119.51
1	B	96	VAL	CB-CA-C	6.60	116.53	110.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	MSE	CG-SE-CE	6.57	113.37	98.92
1	B	203	MSE	CB-CA-C	-6.37	103.03	111.82
1	B	408	ILE	O-C-N	6.21	124.39	120.42
1	A	287	MSE	CA-CB-CG	-6.15	101.80	114.10
1	B	292	ILE	N-CA-C	6.13	116.31	110.42
1	B	193	MSE	N-CA-C	-5.92	104.77	112.23
1	B	458	VAL	N-CA-C	-5.92	104.75	110.72
1	B	423	GLU	N-CA-C	5.88	117.77	111.36
1	A	40	ASP	N-CA-C	5.87	119.63	112.23
1	A	85	PRO	N-CA-C	-5.70	102.87	111.41
1	B	287	MSE	CA-CB-CG	-5.69	102.72	114.10
1	B	374	GLY	CA-C-N	5.57	125.04	119.24
1	B	374	GLY	C-N-CA	5.57	125.04	119.24
1	A	410	CYS	CA-C-N	-5.55	114.41	122.69
1	A	410	CYS	C-N-CA	-5.55	114.41	122.69
1	A	404	THR	N-CA-C	-5.20	105.50	111.07
1	A	260	GLN	CA-C-N	5.19	125.70	119.94
1	A	260	GLN	C-N-CA	5.19	125.70	119.94
1	A	11	SER	N-CA-C	5.07	117.55	111.71
1	A	140	VAL	N-CA-C	5.02	115.33	107.99
1	B	101	ASP	N-CA-C	5.02	118.50	112.38
1	B	356	LEU	N-CA-C	5.01	116.20	109.83

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	40	ASP	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	4184	0	4315	103	0
1	B	4180	0	4323	104	1
2	A	229	0	0	9	1
2	B	229	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8822	0	8638	207	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 12.

All (207) 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:B:203:MSE:HE3	1:B:208:LEU:HD23	1.33	1.08
1:B:354:PHE:HZ	1:B:369:LEU:HD21	1.20	1.03
1:B:266:MSE:HE1	1:B:269:LEU:CD2	1.91	1.01
1:B:67:LYS:NZ	2:B:722:HOH:O	1.75	1.00
1:B:421:ASN:ND2	1:B:455:SER:OG	1.95	0.99
1:B:266:MSE:CE	1:B:269:LEU:HD23	1.94	0.97
1:A:25:THR:HG22	1:A:32:CYS:HB3	1.46	0.95
1:A:266:MSE:HA	1:A:266:MSE:HE2	1.47	0.94
1:B:266:MSE:HE1	1:B:269:LEU:HD22	1.51	0.90
1:B:203:MSE:HE3	1:B:208:LEU:CD2	2.04	0.87
1:B:304:MSE:HE3	1:B:352:GLN:CD	2.00	0.86
1:A:266:MSE:HA	1:A:266:MSE:CE	2.02	0.86
1:B:67:LYS:HE2	1:B:80:THR:OG1	1.75	0.85
1:B:284:THR:HG22	2:B:750:HOH:O	1.77	0.84
1:B:266:MSE:CE	1:B:269:LEU:CD2	2.55	0.83
1:B:412:VAL:HA	1:B:415:LEU:HD22	1.63	0.81
1:A:203:MSE:HE2	2:A:684:HOH:O	1.79	0.81
1:B:304:MSE:HE3	1:B:352:GLN:NE2	1.96	0.79
1:B:25:THR:HG22	1:B:32:CYS:HB3	1.64	0.79
1:B:354:PHE:CZ	1:B:369:LEU:HD21	2.12	0.79
1:B:13:PHE:O	2:B:746:HOH:O	2.00	0.78
1:B:541:ILE:HG23	1:B:543:THR:HG23	1.66	0.77
1:B:148:TRP:HE1	1:B:284:THR:HG21	1.50	0.75
1:A:189:THR:HB	1:A:193:MSE:CE	2.17	0.75
1:A:265:VAL:HG12	1:A:266:MSE:HE3	1.68	0.74
1:B:378:LYS:NZ	2:B:711:HOH:O	2.22	0.73
1:B:296:LEU:HD13	1:B:332:VAL:HG23	1.70	0.72
1:B:18:THR:CG2	1:B:36:THR:HB	2.20	0.71
1:A:76:THR:HB	1:A:93:GLU:HG3	1.73	0.70
1:B:266:MSE:HE3	1:B:266:MSE:HA	1.73	0.70
1:A:114:GLU:HG3	1:A:203:MSE:HE1	1.75	0.69
1:A:13:PHE:O	2:A:777:HOH:O	2.09	0.69
1:B:208:LEU:HD22	1:B:209:PRO:HD2	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:LEU:HD13	1:A:411:ILE:HG21	1.76	0.67
1:B:359:ARG:HG3	1:B:400:LEU:HD22	1.77	0.67
1:A:509:ALA:O	1:A:512:LEU:HB2	1.95	0.67
1:B:266:MSE:HE2	1:B:269:LEU:HD23	1.76	0.67
1:A:40:ASP:HB3	1:A:42:LEU:H	1.60	0.66
1:B:40:ASP:HB3	1:B:42:LEU:H	1.61	0.66
1:A:393:LEU:O	1:A:401:ARG:HD2	1.96	0.65
1:B:442:TRP:O	1:B:446:ILE:HG12	1.97	0.64
1:A:524:LYS:O	1:A:528:GLU:HG2	1.98	0.64
1:B:373:ILE:HD11	1:B:411:ILE:HD12	1.80	0.63
1:B:336:THR:HB	1:B:342:PHE:CE2	2.34	0.63
1:A:412:VAL:HG11	1:A:453:ILE:HD11	1.80	0.63
1:A:85:PRO:O	1:A:86:GLN:HG2	1.98	0.62
1:A:25:THR:CG2	1:A:32:CYS:HB3	2.25	0.62
1:A:27:THR:HG22	1:A:29:LEU:H	1.65	0.62
1:A:66:SER:O	1:A:77:VAL:HG21	1.99	0.62
1:B:121:LEU:HB3	1:B:230:ARG:HD3	1.82	0.61
1:B:463:ASN:HD22	1:B:503:LYS:HZ2	1.47	0.61
1:A:397:ASP:OD2	1:A:399:THR:OG1	2.17	0.61
1:A:185:GLU:O	1:A:187:PRO:HD3	2.01	0.61
1:A:304:MSE:HG2	1:A:356:LEU:HD11	1.82	0.61
1:B:417:GLU:O	1:B:421:ASN:HB2	2.01	0.61
1:B:308:ILE:HG21	1:B:356:LEU:HD23	1.84	0.60
1:A:186:ASP:HB3	1:A:189:THR:OG1	2.01	0.60
1:B:76:THR:HB	1:B:93:GLU:HG3	1.83	0.60
1:A:395:ASP:O	1:A:401:ARG:HD3	2.01	0.60
1:B:63:VAL:HG13	1:B:80:THR:HB	1.84	0.59
1:A:189:THR:HB	1:A:193:MSE:HE2	1.83	0.59
1:B:144:ILE:HD13	2:B:818:HOH:O	2.03	0.59
1:A:266:MSE:HE2	1:A:269:LEU:HB3	1.85	0.58
1:A:26:GLU:HG2	1:A:31:GLN:HG2	1.85	0.58
1:B:393:LEU:O	1:B:401:ARG:HD2	2.03	0.58
1:A:412:VAL:HG11	1:A:453:ILE:CD1	2.34	0.58
1:B:460:ASN:ND2	2:B:691:HOH:O	2.27	0.58
1:A:296:LEU:HD13	1:A:332:VAL:HG23	1.86	0.58
1:A:390:ILE:HG23	1:A:391:GLN:HE21	1.69	0.58
1:B:266:MSE:HE2	1:B:296:LEU:HG	1.85	0.57
1:B:349:LEU:C	1:B:349:LEU:HD23	2.28	0.57
1:A:300:ILE:CG2	1:A:304:MSE:HE3	2.34	0.57
1:B:304:MSE:HE3	1:B:352:GLN:CG	2.34	0.56
1:A:84:ASP:HB3	1:A:87:SER:HB2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:MSE:HE3	1:B:269:LEU:HB3	1.87	0.56
1:B:270:GLU:OE1	1:B:324:LYS:HE2	2.05	0.56
1:B:380:GLU:O	1:B:384:ARG:HB2	2.05	0.56
1:A:189:THR:HB	1:A:193:MSE:HE3	1.86	0.56
1:A:178:ILE:O	1:A:178:ILE:HG22	2.06	0.56
1:B:463:ASN:ND2	1:B:503:LYS:NZ	2.55	0.55
1:A:64:HIS:O	1:A:68:ILE:HG12	2.06	0.55
1:B:308:ILE:CG2	1:B:356:LEU:CD2	2.85	0.55
1:A:412:VAL:HG22	1:A:420:LEU:CD1	2.37	0.55
1:A:541:ILE:HG13	1:A:542:GLN:N	2.20	0.55
1:B:174:TYR:O	1:B:178:ILE:HG12	2.07	0.54
1:A:29:LEU:HD21	1:A:177:ILE:HG21	1.90	0.54
1:A:327:PRO:O	1:A:331:ILE:HG12	2.08	0.54
1:B:273:TYR:OH	2:B:687:HOH:O	2.18	0.53
1:B:409:PRO:O	1:B:412:VAL:HG22	2.09	0.53
1:A:324:LYS:C	1:A:327:PRO:HD2	2.33	0.53
1:B:63:VAL:O	1:B:66:SER:HB2	2.08	0.53
1:B:308:ILE:CG2	1:B:356:LEU:HD23	2.39	0.53
1:A:390:ILE:HG23	1:A:391:GLN:NE2	2.24	0.53
1:B:26:GLU:OE1	1:B:50:LYS:NZ	2.36	0.52
1:B:304:MSE:CE	1:B:352:GLN:NE2	2.71	0.52
1:B:450:ILE:HG23	1:B:454:LEU:HD21	1.91	0.52
1:B:386:TYR:HB3	1:B:387:PRO:HD3	1.91	0.52
1:A:463:ASN:OD1	1:A:503:LYS:NZ	2.39	0.52
1:B:463:ASN:HD22	1:B:503:LYS:NZ	2.07	0.52
1:B:463:ASN:ND2	1:B:503:LYS:HZ2	2.08	0.52
1:A:266:MSE:HG2	1:A:299:ILE:HD13	1.92	0.52
1:A:229:PHE:CE2	1:A:233:LEU:HD22	2.45	0.51
1:B:18:THR:HG22	1:B:36:THR:O	2.10	0.51
1:A:388:HIS:HE1	2:A:734:HOH:O	1.92	0.51
1:B:73:GLY:HA3	1:B:123:THR:OG1	2.09	0.51
1:A:266:MSE:HE2	1:A:269:LEU:CB	2.41	0.50
1:A:189:THR:CB	1:A:193:MSE:CE	2.88	0.50
1:A:248:TYR:CE2	1:A:252:ARG:HD2	2.46	0.50
1:A:433:VAL:HG22	1:A:433:VAL:O	2.12	0.50
1:B:446:ILE:O	1:B:450:ILE:HB	2.12	0.50
1:B:474:LYS:NZ	2:B:717:HOH:O	2.44	0.50
1:A:373:ILE:HG12	1:A:381:ILE:HD13	1.93	0.49
1:B:417:GLU:HG2	1:B:421:ASN:ND2	2.27	0.49
1:B:408:ILE:N	1:B:409:PRO:CD	2.77	0.48
1:B:203:MSE:CE	1:B:208:LEU:HD23	2.25	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:LYS:HD3	1:A:68:ILE:HD11	1.94	0.48
1:A:517:PRO:HG3	2:A:636:HOH:O	2.14	0.48
1:A:7:HIS:HE1	1:A:84:ASP:OD2	1.97	0.48
1:A:27:THR:O	1:A:50:LYS:NZ	2.42	0.47
1:A:121:LEU:HD21	1:A:195:LEU:HD21	1.95	0.47
1:A:189:THR:CB	1:A:193:MSE:HE3	2.44	0.47
1:A:258:ASP:C	1:A:258:ASP:OD2	2.57	0.47
1:B:199:ILE:HG23	1:B:203:MSE:HE2	1.96	0.47
1:A:183:PRO:HG2	1:A:190:ILE:HG13	1.96	0.47
1:A:144:ILE:HD12	1:A:145:LYS:HG3	1.97	0.47
1:B:159:LYS:NZ	2:B:729:HOH:O	2.48	0.46
1:B:446:ILE:HG22	1:B:450:ILE:HD13	1.96	0.46
1:A:241:SER:O	1:A:243:PRO:HD3	2.14	0.46
1:A:308:ILE:HG22	1:A:308:ILE:O	2.15	0.46
1:B:174:TYR:CE2	1:B:178:ILE:HG13	2.50	0.46
1:B:421:ASN:ND2	1:B:455:SER:HG	2.08	0.46
1:A:73:GLY:HA3	1:A:123:THR:OG1	2.14	0.46
1:A:37:ARG:NH2	1:A:40:ASP:OD1	2.47	0.46
1:B:541:ILE:CG2	1:B:543:THR:HG23	2.42	0.46
1:B:255:HIS:CE1	1:B:295:GLU:OE2	2.69	0.46
1:B:454:LEU:HD23	1:B:454:LEU:N	2.31	0.46
1:A:193:MSE:HE2	1:A:222:GLY:HA3	1.97	0.46
1:B:42:LEU:HD22	1:B:43:PRO:HD2	1.97	0.46
1:B:136:SER:O	1:B:139:SER:HB2	2.16	0.46
1:B:256:ILE:HD12	1:B:256:ILE:HA	1.68	0.46
1:A:116:GLY:HA3	1:A:148:TRP:CD2	2.51	0.45
1:A:538:ALA:O	1:A:541:ILE:HG23	2.16	0.45
1:B:324:LYS:HA	1:B:324:LYS:HD3	1.57	0.45
1:A:432:GLN:NE2	1:A:472:SER:OG	2.49	0.45
1:A:230:ARG:HD3	2:A:791:HOH:O	2.16	0.45
1:A:273:TYR:CD1	1:A:331:ILE:CD1	2.99	0.45
1:A:49:ALA:HB1	1:A:59:ILE:HD12	1.99	0.45
1:A:214:VAL:O	1:A:218:MSE:HG3	2.16	0.44
1:A:412:VAL:HG22	1:A:420:LEU:HD11	1.99	0.44
1:B:255:HIS:HE1	1:B:295:GLU:OE2	1.99	0.44
1:A:17:TYR:OH	2:A:760:HOH:O	2.20	0.44
1:A:263:LEU:HD11	1:A:321:ALA:CB	2.47	0.44
1:B:40:ASP:HB3	1:B:42:LEU:N	2.31	0.44
1:B:456:THR:HG21	1:B:460:ASN:OD1	2.17	0.44
1:A:266:MSE:CE	1:A:266:MSE:CA	2.86	0.44
1:B:408:ILE:HB	1:B:409:PRO:HD3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:MSE:HG2	1:B:325:LEU:HD11	2.01	0.43
1:B:365:LEU:O	1:B:369:LEU:HD23	2.18	0.43
1:A:271:ASN:O	1:A:275:GLU:HG2	2.18	0.43
1:A:379:SER:O	1:A:383:SER:HB3	2.18	0.43
1:B:18:THR:HG21	1:B:36:THR:HB	1.97	0.43
1:A:15:PHE:HA	1:A:16:PRO:HD3	1.89	0.43
1:B:18:THR:HG23	1:B:36:THR:HB	1.98	0.43
1:B:167:PHE:HA	1:B:170:ARG:HH11	1.83	0.43
1:B:514:ASP:O	1:B:520:ARG:HD3	2.18	0.43
1:A:112:GLY:HA3	2:A:601:HOH:O	2.17	0.43
1:A:373:ILE:HA	1:A:381:ILE:HD11	2.01	0.43
1:A:412:VAL:CG1	1:A:453:ILE:CD1	2.96	0.43
1:B:71:ILE:HB	1:B:74:LEU:HD12	1.99	0.43
1:B:339:THR:O	1:B:341:THR:HG23	2.19	0.43
1:A:486:TYR:CE1	1:A:490:LYS:HE2	2.54	0.43
1:B:440:ARG:HD2	1:B:475:ASP:OD1	2.19	0.43
1:A:147:GLU:OE2	1:A:248:TYR:OH	2.28	0.43
1:B:182:LEU:HD22	1:B:184:CYS:O	2.19	0.43
1:A:18:THR:HG23	1:A:38:LYS:HG3	1.99	0.42
1:A:178:ILE:O	1:A:178:ILE:CG2	2.66	0.42
1:A:441:THR:HG21	1:A:479:LYS:HG2	2.01	0.42
1:A:306:GLN:O	1:A:309:SER:HB2	2.19	0.42
1:A:374:GLY:O	1:A:375:PRO:C	2.60	0.42
1:A:486:TYR:CZ	1:A:490:LYS:HE2	2.54	0.42
1:A:266:MSE:CG	1:A:299:ILE:HD13	2.49	0.42
1:A:325:LEU:O	1:A:329:LEU:HB2	2.19	0.42
1:B:96:VAL:HA	1:B:97:PRO:HD3	1.93	0.42
1:A:289:GLU:HG2	1:A:341:THR:HG21	2.02	0.42
1:B:160:GLU:OE2	1:B:160:GLU:HA	2.19	0.42
1:A:15:PHE:CD2	1:A:16:PRO:HD2	2.55	0.42
1:A:158:SER:OG	1:A:160:GLU:HG2	2.19	0.42
1:B:120:LEU:HD22	1:B:150:LEU:HD21	2.02	0.42
1:B:336:THR:HB	1:B:342:PHE:CD2	2.55	0.42
1:B:488:LEU:HB2	1:B:526:LEU:HD11	2.01	0.42
1:A:434:ASP:O	1:A:440:ARG:HD3	2.20	0.41
1:B:425:LEU:HB2	2:B:806:HOH:O	2.21	0.41
1:A:37:ARG:O	1:A:41:SER:N	2.54	0.41
1:A:412:VAL:CG1	1:A:453:ILE:HD12	2.51	0.41
1:A:424:LEU:HD23	1:A:424:LEU:HA	1.78	0.41
1:B:42:LEU:CD2	1:B:43:PRO:HD2	2.50	0.41
1:B:121:LEU:HB3	1:B:230:ARG:CD	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:TYR:CE1	1:A:37:ARG:HG3	2.56	0.41
1:A:98:PHE:HE1	1:A:148:TRP:CZ2	2.39	0.41
1:A:406:LYS:NZ	2:A:774:HOH:O	2.54	0.41
1:B:262:LYS:HB3	1:B:299:ILE:HD11	2.02	0.41
1:B:518:ILE:HD11	1:B:522:ARG:HH21	1.86	0.41
1:A:386:TYR:HA	1:A:389:PHE:HB3	2.03	0.41
1:B:51:ARG:CZ	1:B:60:LEU:HD21	2.51	0.41
1:B:366:LEU:HA	1:B:369:LEU:HD23	2.03	0.41
1:B:512:LEU:CD2	1:B:524:LYS:HA	2.51	0.41
1:A:410:CYS:HB3	2:A:661:HOH:O	2.20	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:543:THR:O	2:A:636:HOH:O[2_746]	1.89	0.31

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	527/560 (94%)	511 (97%)	15 (3%)	1 (0%)	43	51
1	B	525/560 (94%)	515 (98%)	10 (2%)	0	100	100
All	All	1052/1120 (94%)	1026 (98%)	25 (2%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	86	GLN

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	480/502 (96%)	422 (88%)	58 (12%)	5	4
1	B	481/502 (96%)	436 (91%)	45 (9%)	8	9
All	All	961/1004 (96%)	858 (89%)	103 (11%)	6	6

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	26	GLU
1	A	27	THR
1	A	36	THR
1	A	54	GLU
1	A	56	GLU
1	A	58	LEU
1	A	60	LEU
1	A	71	ILE
1	A	78	LEU
1	A	93	GLU
1	A	115	LEU
1	A	132	LEU
1	A	144	ILE
1	A	153	LEU
1	A	157	SER
1	A	162	LEU
1	A	163	SER
1	A	169	SER
1	A	198	LEU
1	A	203	MSE
1	A	206	SER
1	A	208	LEU
1	A	210	LYS
1	A	233	LEU
1	A	235	ASN
1	A	252	ARG

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Mol	Chain	Res	Type
1	A	263	LEU
1	A	266	MSE
1	A	275	GLU
1	A	276	SER
1	A	284	THR
1	A	296	LEU
1	A	302	LEU
1	A	307	SER
1	A	322	SER
1	A	329	LEU
1	A	352	GLN
1	A	369	LEU
1	A	379	SER
1	A	381	ILE
1	A	400	LEU
1	A	411	ILE
1	A	412	VAL
1	A	423	GLU
1	A	424	LEU
1	A	450	ILE
1	A	453	ILE
1	A	454	LEU
1	A	456	THR
1	A	493	GLU
1	A	512	LEU
1	A	518	ILE
1	A	526	LEU
1	A	532	GLU
1	A	541	ILE
1	A	543	THR
1	A	544	ASN
1	B	25	THR
1	B	29	LEU
1	B	36	THR
1	B	40	ASP
1	B	42	LEU
1	B	44	VAL
1	B	96	VAL
1	B	103	LEU
1	B	115	LEU
1	B	121	LEU
1	B	144	ILE

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Mol	Chain	Res	Type
1	B	153	LEU
1	B	157	SER
1	B	159	LYS
1	B	162	LEU
1	B	170	ARG
1	B	182	LEU
1	B	198	LEU
1	B	203	MSE
1	B	208	LEU
1	B	237	GLU
1	B	246	ASN
1	B	252	ARG
1	B	256	ILE
1	B	272	LEU
1	B	284	THR
1	B	287	MSE
1	B	296	LEU
1	B	299	ILE
1	B	302	LEU
1	B	305	THR
1	B	308	ILE
1	B	349	LEU
1	B	369	LEU
1	B	372	LEU
1	B	408	ILE
1	B	415	LEU
1	B	420	LEU
1	B	424	LEU
1	B	425	LEU
1	B	450	ILE
1	B	453	ILE
1	B	456	THR
1	B	506	THR
1	B	526	LEU

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	7	HIS
1	A	228	ASN
1	A	246	ASN
1	A	271	ASN

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Mol	Chain	Res	Type
1	A	352	GLN
1	A	388	HIS
1	A	391	GLN
1	A	432	GLN
1	A	460	ASN
1	A	502	ASN
1	B	7	HIS
1	B	31	GLN
1	B	119	GLN
1	B	129	ASN
1	B	246	ASN
1	B	249	GLN
1	B	255	HIS
1	B	352	GLN
1	B	463	ASN
1	B	542	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	524/560 (93%)	-1.20	0 100 100	19, 35, 56, 82	0
1	B	522/560 (93%)	-1.26	0 100 100	20, 33, 50, 75	0
All	All	1046/1120 (93%)	-1.23	0 100 100	19, 34, 54, 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.