



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:50 PM UTC

PDB ID : 2UXT / pdb_00002uxt
Title : SufI Protein from Escherichia Coli
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Deposited on : 2007-03-29
Resolution : 1.90 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

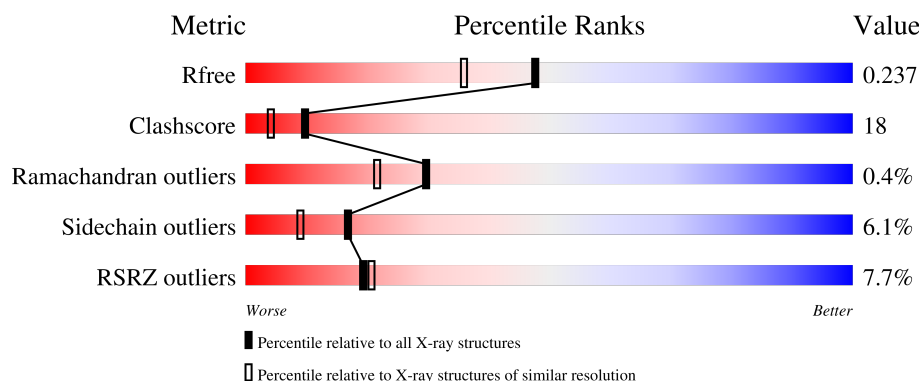
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	451	
1	B	451	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7014 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 PROTEIN SUFI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3313	2103	589	607	14			
1	B	420	Total	C	N	O	S	0	0	0
			3295	2095	580	606	14			

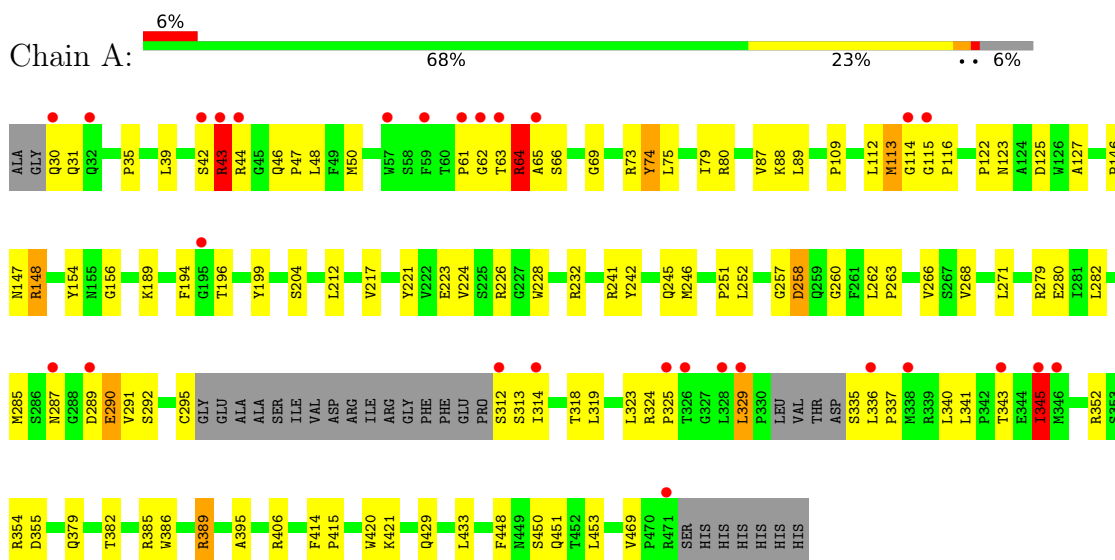
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	229	Total	O	0	0
			229	229		
2	B	177	Total	O	0	0
			177	177		

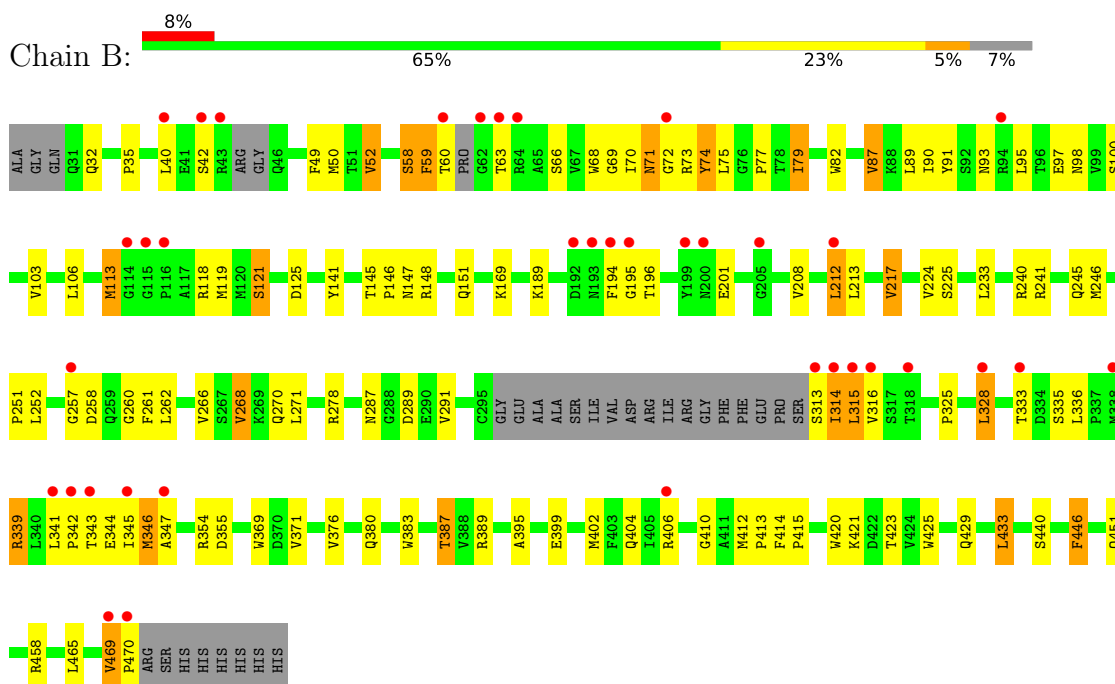
3 Residue-property plots

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: PROTEIN SUFI



• Molecule 1: PROTEIN SUFI



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.47Å 89.55Å 153.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.84 – 1.90 39.84 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (39.84-1.90) 98.0 (39.84-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 1.89Å)	Xtriage
Refinement program	TNT 5.13.1.0	Depositor
R, R_{free}	0.191 , (Not available) 0.198 , 0.237	Depositor DCC
R_{free} test set	2976 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.8	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	7014	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.41	0/3402	0.84	8/4645 (0.2%)
1	B	0.40	0/3382	0.87	7/4619 (0.2%)
All	All	0.41	0/6784	0.86	15/9264 (0.2%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	71	ASN	N-CA-C	12.47	128.21	112.24
1	B	113	MET	N-CA-C	8.59	120.42	111.14
1	A	329	LEU	C-N-CD	-8.17	91.52	125.00
1	A	113	MET	N-CA-C	7.88	119.56	110.97
1	B	74	TYR	N-CA-C	-7.73	97.99	110.20
1	B	59	PHE	N-CA-C	-7.32	104.24	113.02
1	A	74	TYR	N-CA-C	-6.91	97.99	109.46
1	A	345	ILE	N-CA-C	6.25	118.45	109.51
1	B	446	PHE	N-CA-C	-6.01	101.63	110.52
1	B	87	VAL	N-CA-C	5.38	115.85	107.99
1	B	262	LEU	N-CA-C	-5.27	102.49	110.39
1	A	448	PHE	N-CA-C	-5.18	100.59	109.24
1	A	258	ASP	CB-CA-C	-5.16	102.55	110.81
1	A	450	SER	N-CA-C	-5.14	102.67	110.28
1	A	453	LEU	N-CA-C	5.07	117.47	111.33

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3313	0	3262	109	0
1	B	3295	0	3243	125	0
2	A	229	0	0	8	0
2	B	177	0	0	6	0
All	All	7014	0	6505	231	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 18.

All (231) 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:347:ALA:HB1	1:B:433:LEU:HD23	1.20	1.12
1:A:312:SER:HA	1:A:314:ILE:H	1.25	1.00
1:B:40:LEU:HD22	1:B:77:PRO:HB2	1.46	0.95
1:A:312:SER:N	1:A:313:SER:HB3	1.82	0.93
1:B:58:SER:HB2	1:B:60:THR:HG23	1.54	0.89
1:B:402:MET:HE3	1:B:420:TRP:CD2	2.09	0.87
1:B:212:LEU:HD12	1:B:315:LEU:HD13	1.57	0.87
1:B:50:MET:HE1	1:B:79:ILE:HD11	1.56	0.87
1:A:282:LEU:HD12	1:A:336:LEU:HD22	1.54	0.86
1:B:347:ALA:HB1	1:B:433:LEU:CD2	2.06	0.85
1:B:148:ARG:HG2	1:B:148:ARG:HH11	1.41	0.84
1:A:312:SER:HA	1:A:314:ILE:N	1.93	0.84
1:B:266:VAL:HG11	1:B:415:PRO:HB2	1.60	0.84
1:A:228:TRP:CZ3	1:A:336:LEU:HD21	2.12	0.84
1:B:402:MET:HE3	1:B:420:TRP:CE2	2.14	0.83
1:A:282:LEU:CD1	1:A:336:LEU:HD22	2.08	0.82
1:B:119:MET:HE3	1:B:145:THR:HG23	1.61	0.81
1:A:287:ASN:HB3	2:A:2134:HOH:O	1.81	0.80
1:B:98:ASN:CB	1:B:119:MET:HE2	2.12	0.79
1:B:35:PRO:HD3	1:B:217:VAL:HG11	1.66	0.78
1:B:119:MET:HE3	1:B:145:THR:CG2	2.12	0.78
1:B:380:GLN:OE1	1:B:469:VAL:HG22	1.84	0.77
1:B:50:MET:HG3	1:B:70:ILE:HG21	1.68	0.76
1:A:257:GLY:HA2	2:A:2205:HOH:O	1.85	0.76
1:A:64:ARG:HD3	1:A:194:PHE:CD1	2.21	0.76
1:B:347:ALA:CB	1:B:433:LEU:HD23	2.09	0.74
1:A:65:ALA:HB3	1:A:196:THR:HG21	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:LEU:CD2	1:B:77:PRO:HB2	2.19	0.73
1:A:345:ILE:CD1	1:A:420:TRP:HE1	2.02	0.73
1:B:341:LEU:CD1	1:B:345:ILE:HD11	2.19	0.73
1:B:148:ARG:HG2	1:B:148:ARG:NH1	2.01	0.73
1:B:212:LEU:CD1	1:B:315:LEU:HD13	2.18	0.72
1:A:312:SER:HA	1:A:314:ILE:HD12	1.71	0.72
1:A:345:ILE:HD11	1:A:420:TRP:HE1	1.53	0.72
1:B:50:MET:HE1	1:B:79:ILE:CD1	2.20	0.71
1:A:379:GLN:O	1:A:382:THR:HG22	1.89	0.71
1:A:35:PRO:HD3	1:A:217:VAL:HG11	1.73	0.71
1:B:50:MET:CE	1:B:79:ILE:HD11	2.21	0.70
1:B:89:LEU:HD23	1:B:91:TYR:CE1	2.27	0.70
1:A:42:SER:O	1:A:43:ARG:HB2	1.90	0.70
1:A:345:ILE:HD11	1:A:420:TRP:NE1	2.07	0.70
1:B:40:LEU:HD22	1:B:77:PRO:CB	2.21	0.69
1:B:98:ASN:HB2	1:B:119:MET:HE2	1.73	0.69
1:A:287:ASN:O	1:A:287:ASN:ND2	2.26	0.69
1:A:312:SER:CA	1:A:314:ILE:H	2.04	0.68
1:B:245:GLN:HG2	1:B:270:GLN:HG3	1.75	0.68
1:A:226:ARG:HD3	1:A:329:LEU:CD2	2.24	0.67
1:A:257:GLY:HA3	1:A:260:GLY:O	1.95	0.66
1:B:257:GLY:HA2	2:B:2160:HOH:O	1.95	0.66
1:B:58:SER:HB2	1:B:60:THR:CG2	2.24	0.66
1:A:289:ASP:HB2	2:A:2134:HOH:O	1.94	0.65
1:B:380:GLN:CD	1:B:469:VAL:HG22	2.20	0.64
1:B:341:LEU:HD23	1:B:342:PRO:HD2	1.78	0.64
1:B:339:ARG:HG2	1:B:339:ARG:HH11	1.63	0.64
1:B:32:GLN:HB2	1:B:217:VAL:HG21	1.80	0.62
1:A:228:TRP:CE3	1:A:336:LEU:HD21	2.34	0.62
1:A:414:PHE:CE2	1:B:251:PRO:HB3	2.34	0.62
1:A:258:ASP:HB2	1:A:421:LYS:HA	1.81	0.62
1:A:312:SER:CA	1:A:313:SER:HB3	2.29	0.62
1:B:395:ALA:HB3	1:B:451:GLN:HB2	1.81	0.62
1:B:50:MET:HG3	1:B:70:ILE:CG2	2.30	0.62
1:B:148:ARG:HA	2:B:2176:HOH:O	1.98	0.62
1:B:50:MET:HE3	1:B:70:ILE:HG21	1.81	0.62
1:A:63:THR:HG21	2:A:2022:HOH:O	1.98	0.61
1:B:406:ARG:HD3	1:B:433:LEU:CD2	2.30	0.61
1:A:226:ARG:HD3	1:A:329:LEU:HD21	1.81	0.61
1:B:49:PHE:CZ	1:B:90:ILE:HD12	2.36	0.61
1:A:285:MET:HE3	1:A:325:PRO:HG3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:LEU:HB3	1:A:337:PRO:CD	2.32	0.60
1:A:406:ARG:HD2	1:A:433:LEU:HB2	1.84	0.60
1:A:289:ASP:O	1:A:325:PRO:HG2	2.02	0.60
1:A:30:GLN:HG3	2:A:2140:HOH:O	2.01	0.59
1:B:406:ARG:HH11	1:B:433:LEU:CD2	2.15	0.59
1:B:212:LEU:HD12	1:B:315:LEU:CD1	2.31	0.59
1:A:355:ASP:OD2	1:A:389:ARG:NH1	2.35	0.59
1:B:98:ASN:HB3	1:B:119:MET:HE2	1.84	0.59
1:A:287:ASN:OD1	1:B:410:GLY:HA3	2.02	0.58
1:B:406:ARG:HD3	1:B:433:LEU:HD22	1.84	0.58
1:A:406:ARG:CD	1:A:433:LEU:HB2	2.33	0.58
1:B:146:PRO:O	1:B:147:ASN:HB2	2.03	0.58
1:B:268:VAL:HG21	1:B:271:LEU:HD13	1.86	0.58
1:A:266:VAL:HG11	1:A:415:PRO:HB2	1.85	0.58
1:B:68:TRP:CE3	1:B:95:LEU:HD13	2.38	0.58
1:B:406:ARG:NH1	2:B:2149:HOH:O	2.29	0.57
1:B:339:ARG:HH11	1:B:339:ARG:CG	2.18	0.57
1:B:458:ARG:HD3	2:B:2070:HOH:O	2.05	0.57
1:A:312:SER:CB	1:A:314:ILE:HD13	2.35	0.56
1:A:395:ALA:HB3	1:A:451:GLN:HB2	1.85	0.56
1:A:285:MET:CE	1:A:325:PRO:HG3	2.35	0.56
1:A:62:GLY:N	1:A:196:THR:OG1	2.38	0.56
1:A:323:LEU:O	1:A:325:PRO:HD3	2.06	0.56
1:B:59:PHE:O	1:B:60:THR:O	2.24	0.56
1:A:148:ARG:NH1	1:A:148:ARG:HG2	2.19	0.56
1:A:325:PRO:HB2	2:A:2142:HOH:O	2.07	0.55
1:B:35:PRO:HD3	1:B:217:VAL:CG1	2.36	0.55
1:B:369:TRP:CZ2	1:B:371:VAL:HG12	2.42	0.55
1:A:39:LEU:HD21	1:A:80:ARG:CZ	2.37	0.55
1:A:46:GLN:HB3	1:A:47:PRO:HD2	1.88	0.55
1:A:212:LEU:HD21	1:A:242:TYR:HE2	1.72	0.55
1:B:224:VAL:O	1:B:325:PRO:HA	2.07	0.55
1:A:382:THR:HG21	2:A:2180:HOH:O	2.07	0.53
1:A:385:ARG:HH11	1:A:406:ARG:HH22	1.56	0.53
1:B:68:TRP:CZ3	1:B:95:LEU:HD13	2.43	0.53
1:B:341:LEU:HD13	1:B:345:ILE:HD11	1.90	0.53
1:A:312:SER:HB2	1:A:314:ILE:HD13	1.91	0.52
1:A:312:SER:CA	1:A:314:ILE:HD12	2.37	0.52
1:B:258:ASP:HB2	1:B:421:LYS:HA	1.92	0.52
1:B:241:ARG:HD2	1:B:425:TRP:CE2	2.44	0.52
1:A:148:ARG:HG2	1:A:148:ARG:HH11	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:TYR:O	1:A:189:LYS:NZ	2.43	0.52
1:A:245:GLN:HA	1:A:252:LEU:HD11	1.92	0.52
1:B:63:THR:HG21	1:B:195:GLY:HA3	1.91	0.52
1:B:257:GLY:HA3	1:B:260:GLY:O	2.10	0.52
1:A:35:PRO:CD	1:A:217:VAL:HG11	2.40	0.51
1:B:341:LEU:HD23	1:B:342:PRO:CD	2.40	0.51
1:A:352:ARG:HD3	1:A:354:ARG:NH2	2.26	0.51
1:B:49:PHE:CE2	1:B:90:ILE:HD12	2.46	0.51
1:B:69:GLY:HA3	1:B:73:ARG:O	2.11	0.51
1:A:352:ARG:HD3	1:A:354:ARG:CZ	2.40	0.51
1:A:146:PRO:O	1:A:147:ASN:HB2	2.09	0.51
1:A:336:LEU:HB3	1:A:337:PRO:HD2	1.93	0.50
1:B:74:TYR:O	1:B:75:LEU:HB2	2.10	0.50
1:B:383:TRP:CH2	1:B:420:TRP:HH2	2.30	0.50
1:A:64:ARG:HD3	1:A:194:PHE:CE1	2.47	0.50
1:B:68:TRP:CD2	1:B:95:LEU:HD13	2.46	0.50
1:A:282:LEU:HD13	1:A:336:LEU:HD22	1.93	0.50
1:A:226:ARG:NH2	1:A:329:LEU:HD21	2.26	0.50
1:A:74:TYR:O	1:A:75:LEU:HB2	2.12	0.49
1:B:344:GLU:HB3	1:B:346:MET:HE3	1.94	0.49
1:A:212:LEU:HD21	1:A:242:TYR:CE2	2.47	0.49
1:A:148:ARG:HH11	1:A:148:ARG:CG	2.26	0.49
1:B:245:GLN:NE2	1:B:270:GLN:HE21	2.11	0.49
1:B:52:VAL:HG22	1:B:93:ASN:HA	1.96	0.48
1:B:97:GLU:CD	1:B:148:ARG:HE	2.21	0.48
1:A:224:VAL:HG21	1:A:285:MET:HG3	1.94	0.48
1:B:354:ARG:O	1:B:387:THR:HG22	2.14	0.47
1:B:421:LYS:HE3	1:B:423:THR:O	2.13	0.47
1:A:379:GLN:HE21	1:A:469:VAL:CG2	2.28	0.47
1:B:208:VAL:CG1	1:B:240:ARG:NH2	2.77	0.47
1:B:355:ASP:OD1	1:B:389:ARG:NH2	2.35	0.47
1:B:404:GLN:HG2	1:B:406:ARG:HD2	1.97	0.47
1:A:69:GLY:HA3	1:A:73:ARG:O	2.14	0.46
1:A:226:ARG:HD3	1:A:329:LEU:HD22	1.96	0.46
1:A:354:ARG:HG3	1:A:386:TRP:NE1	2.29	0.46
1:A:389:ARG:HD3	1:A:429:GLN:OE1	2.15	0.46
1:B:35:PRO:CD	1:B:217:VAL:CG1	2.93	0.46
1:B:98:ASN:HB3	1:B:119:MET:HG2	1.97	0.46
1:A:263:PRO:HA	1:A:341:LEU:HB2	1.97	0.46
1:B:82:TRP:CH2	1:B:169:LYS:HD2	2.51	0.46
1:A:50:MET:HE2	1:A:79:ILE:HD11	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:ARG:CZ	1:B:376:VAL:HG13	2.46	0.46
1:B:355:ASP:HA	1:B:387:THR:HG23	1.98	0.45
1:A:335:SER:N	1:A:336:LEU:CA	2.80	0.45
1:A:385:ARG:NH1	1:A:406:ARG:HH22	2.13	0.45
1:A:389:ARG:HG2	1:A:429:GLN:HG3	1.97	0.45
1:B:194:PHE:CD1	1:B:194:PHE:N	2.84	0.45
1:B:328:LEU:N	1:B:328:LEU:HD13	2.31	0.45
1:B:339:ARG:CG	1:B:339:ARG:NH1	2.78	0.45
1:B:50:MET:HE2	1:B:141:TYR:CZ	2.51	0.45
1:B:389:ARG:HG2	1:B:429:GLN:HB2	1.98	0.45
1:A:385:ARG:HH11	1:A:406:ARG:NH2	2.13	0.45
1:B:50:MET:CE	1:B:141:TYR:CZ	3.00	0.45
1:A:226:ARG:HH21	1:A:329:LEU:HD21	1.81	0.44
1:B:246:MET:HE3	1:B:291:VAL:CG2	2.48	0.44
1:B:50:MET:O	1:B:91:TYR:HA	2.18	0.44
1:B:369:TRP:CZ2	1:B:371:VAL:CG1	3.00	0.44
1:A:223:GLU:HA	1:A:324:ARG:O	2.18	0.44
1:B:100:SER:OG	1:B:145:THR:HA	2.17	0.44
1:A:75:LEU:HA	1:A:156:GLY:O	2.17	0.44
1:B:40:LEU:H	1:B:40:LEU:HD23	1.82	0.44
1:A:35:PRO:CD	1:A:217:VAL:CG1	2.95	0.44
1:A:89:LEU:O	1:A:127:ALA:HA	2.18	0.44
1:B:71:ASN:N	1:B:72:GLY:HA2	2.33	0.44
1:B:383:TRP:HH2	1:B:420:TRP:HH2	1.65	0.44
1:B:469:VAL:O	1:B:470:PRO:O	2.36	0.44
1:A:122:PRO:O	1:A:123:ASN:HB2	2.17	0.44
1:B:68:TRP:CE3	1:B:95:LEU:CD1	3.01	0.43
1:B:103:VAL:HG12	1:B:106:LEU:HB2	2.00	0.43
1:A:312:SER:OG	1:A:314:ILE:HB	2.18	0.43
1:A:335:SER:N	1:A:336:LEU:C	2.75	0.43
1:B:261:PHE:O	1:B:341:LEU:HD12	2.19	0.43
1:B:201:GLU:CG	1:B:458:ARG:HH21	2.31	0.43
1:B:89:LEU:HD21	1:B:91:TYR:CZ	2.54	0.43
1:B:287:ASN:OD1	1:B:289:ASP:HB2	2.19	0.43
1:A:212:LEU:HD11	1:A:295:CYS:SG	2.58	0.43
1:A:271:LEU:HD21	1:A:279:ARG:HD2	2.01	0.43
1:B:339:ARG:NH2	1:B:344:GLU:OE2	2.51	0.43
1:A:251:PRO:HB3	1:B:414:PHE:CE2	2.53	0.43
1:B:343:THR:OG1	1:B:344:GLU:N	2.51	0.43
1:A:232:ARG:HG2	1:A:280:GLU:HG2	2.01	0.42
1:A:292:SER:HB3	1:A:319:LEU:CD1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:ARG:HD3	1:B:433:LEU:HD23	2.01	0.42
1:B:208:VAL:HG13	1:B:240:ARG:NH2	2.34	0.42
1:B:241:ARG:HD2	1:B:425:TRP:CZ2	2.55	0.42
1:A:312:SER:CB	1:A:314:ILE:CD1	2.98	0.42
1:A:355:ASP:CG	1:A:389:ARG:NH1	2.77	0.42
1:B:89:LEU:CD2	1:B:91:TYR:CE1	2.99	0.42
1:B:148:ARG:HH11	1:B:148:ARG:CG	2.17	0.42
1:A:113:MET:HB2	1:A:114:GLY:HA2	2.01	0.42
1:A:42:SER:OG	1:A:43:ARG:N	2.53	0.42
1:B:89:LEU:CD2	1:B:91:TYR:CZ	3.03	0.42
1:B:465:LEU:C	1:B:465:LEU:HD23	2.44	0.42
1:A:65:ALA:HB3	1:A:196:THR:CG2	2.45	0.41
1:A:246:MET:HE3	1:A:291:VAL:CG2	2.51	0.41
1:B:63:THR:CG2	1:B:195:GLY:HA3	2.50	0.41
1:A:345:ILE:HD11	1:A:420:TRP:CE2	2.54	0.41
1:A:47:PRO:HB3	1:A:88:LYS:CG	2.50	0.41
1:A:48:LEU:HB2	1:A:89:LEU:HD12	2.02	0.41
1:A:109:PRO:HD2	1:A:112:LEU:HD12	2.03	0.41
1:A:262:LEU:O	1:A:340:LEU:HB2	2.21	0.41
1:B:113:MET:HE2	1:B:399:GLU:CD	2.44	0.41
1:B:189:LYS:NZ	2:B:2067:HOH:O	2.42	0.41
1:A:50:MET:CE	1:A:79:ILE:HD11	2.49	0.41
1:A:290:GLU:HG2	1:A:291:VAL:N	2.35	0.41
1:B:68:TRP:CH2	1:B:95:LEU:HD13	2.56	0.41
1:B:98:ASN:OD1	1:B:121:SER:HA	2.21	0.41
1:B:440:SER:HB3	1:B:446:PHE:CE2	2.56	0.41
1:B:395:ALA:CB	1:B:451:GLN:HB2	2.49	0.41
1:A:154:TYR:CD2	1:A:199:TYR:HB2	2.57	0.40
1:A:292:SER:HB3	1:A:319:LEU:HD11	2.02	0.40
1:A:295:CYS:O	1:A:318:THR:OG1	2.28	0.40
1:B:233:LEU:O	1:B:278:ARG:HA	2.21	0.40
1:B:412:MET:HA	1:B:413:PRO:HD3	1.91	0.40
1:B:58:SER:CB	1:B:60:THR:HG23	2.39	0.40
1:B:314:ILE:H	1:B:314:ILE:HG12	1.60	0.40
1:B:335:SER:HA	2:B:2111:HOH:O	2.21	0.40
1:A:73:ARG:HD2	2:A:2026:HOH:O	2.20	0.40
1:A:221:TYR:CD2	1:A:221:TYR:C	2.99	0.40
1:B:58:SER:C	1:B:60:THR:H	2.27	0.40
1:B:341:LEU:CD2	1:B:343:THR:CG2	2.98	0.40
1:A:115:GLY:HA3	1:A:116:PRO:HD2	1.90	0.40

There are no symmetry-related clashes.

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	416/451 (92%)	395 (95%)	18 (4%)	3 (1%)	18	10
1	B	412/451 (91%)	393 (95%)	19 (5%)	0	100	100
All	All	828/902 (92%)	788 (95%)	37 (4%)	3 (0%)	30	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	ARG
1	A	61	PRO
1	A	64	ARG

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	363/386 (94%)	348 (96%)	15 (4%)	27	19
1	B	362/386 (94%)	333 (92%)	29 (8%)	11	5
All	All	725/772 (94%)	681 (94%)	44 (6%)	17	9

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLN
1	A	43	ARG
1	A	44	ARG

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Mol	Chain	Res	Type
1	A	64	ARG
1	A	66	SER
1	A	87	VAL
1	A	125	ASP
1	A	148	ARG
1	A	204	SER
1	A	241	ARG
1	A	268	VAL
1	A	290	GLU
1	A	343	THR
1	A	345	ILE
1	A	389	ARG
1	B	42	SER
1	B	52	VAL
1	B	58	SER
1	B	66	SER
1	B	79	ILE
1	B	87	VAL
1	B	118	ARG
1	B	121	SER
1	B	125	ASP
1	B	151	GLN
1	B	196	THR
1	B	212	LEU
1	B	213	LEU
1	B	217	VAL
1	B	225	SER
1	B	252	LEU
1	B	268	VAL
1	B	313	SER
1	B	314	ILE
1	B	315	LEU
1	B	316	VAL
1	B	328	LEU
1	B	333	THR
1	B	336	LEU
1	B	339	ARG
1	B	346	MET
1	B	387	THR
1	B	433	LEU
1	B	469	VAL

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	107	GLN
1	A	187	GLN
1	A	372	ASN
1	A	404	GLN
1	A	407	ASN
1	A	438	GLN
1	A	443	HIS
1	A	449	ASN
1	A	451	GLN
1	B	107	GLN
1	B	144	ASN
1	B	151	GLN
1	B	270	GLN
1	B	438	GLN
1	B	449	ASN
1	B	451	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](#)

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 ⓘ

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	422/451 (93%)	0.36	28 (6%) 24 26	15, 26, 50, 72	0
1	B	420/451 (93%)	0.55	37 (8%) 15 16	18, 29, 52, 78	0
All	All	842/902 (93%)	0.46	65 (7%) 19 21	15, 27, 51, 78	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	469	VAL	6.2
1	A	287	ASN	5.7
1	A	62	GLY	5.6
1	B	60	THR	5.6
1	A	336	LEU	5.4
1	A	61	PRO	5.0
1	A	345	ILE	4.9
1	B	194	PHE	4.7
1	A	63	THR	4.6
1	A	65	ALA	4.6
1	B	316	VAL	4.5
1	A	57	TRP	4.4
1	A	343	THR	4.3
1	B	315	LEU	4.3
1	B	314	ILE	4.2
1	A	114	GLY	3.8
1	B	470	PRO	3.7
1	A	471	ARG	3.7
1	B	195	GLY	3.5
1	B	115	GLY	3.5
1	B	313	SER	3.4
1	B	114	GLY	3.4
1	A	289	ASP	3.4
1	A	325	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	200	ASN	3.3
1	A	329	LEU	3.2
1	B	42	SER	3.2
1	A	43	ARG	3.2
1	A	312	SER	3.2
1	B	62	GLY	3.1
1	B	205	GLY	3.1
1	B	64	ARG	3.1
1	A	338	MET	3.0
1	B	192	ASP	2.9
1	A	195	GLY	2.9
1	B	347	ALA	2.9
1	B	343	THR	2.8
1	A	346	MET	2.8
1	B	40	LEU	2.7
1	B	341	LEU	2.7
1	A	115	GLY	2.6
1	B	43	ARG	2.6
1	B	63	THR	2.6
1	B	342	PRO	2.6
1	A	59	PHE	2.6
1	A	42	SER	2.5
1	A	328	LEU	2.5
1	A	32	GLN	2.4
1	A	30	GLN	2.4
1	A	314	ILE	2.4
1	B	193	ASN	2.3
1	B	116	PRO	2.3
1	B	345	ILE	2.2
1	B	328	LEU	2.2
1	B	338	MET	2.2
1	A	326	THR	2.2
1	B	212	LEU	2.2
1	B	318	THR	2.1
1	A	44	ARG	2.1
1	B	72	GLY	2.1
1	B	406	ARG	2.1
1	B	199	TYR	2.1
1	B	333	THR	2.1
1	B	94	ARG	2.1
1	B	257	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.