



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:18 AM UTC

PDB ID : 1F89 / pdb_00001f89
Title : Crystal structure of *Saccharomyces cerevisiae* Nit3, a member of branch 10 of the nitrilase superfamily
Authors : Kumaran, D.; Eswaramoorthy, S.; Studier, F.W.; Swaminathan, S.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2000-06-29
Resolution : 2.40 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

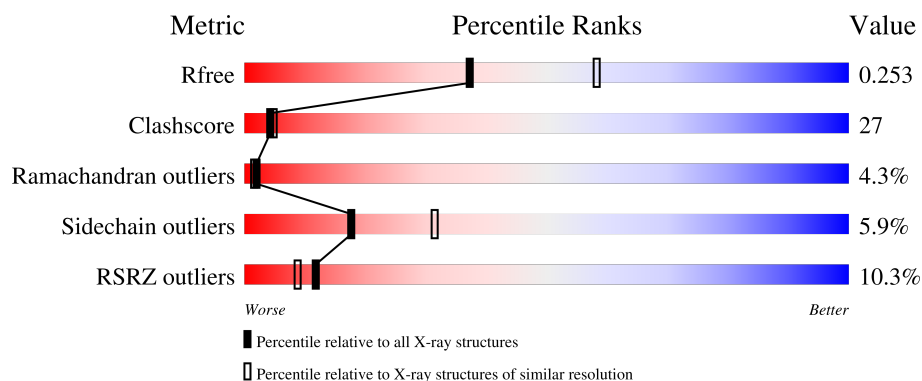
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	291	9% 47% 38% 7% 7%
1	B	291	10% 44% 41% 8% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4391 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 32.5 KDA PROTEIN YLR351C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2137	1363	366	399	9			
1	B	271	Total	C	N	O	S	0	0	0
			2137	1363	366	399	9			

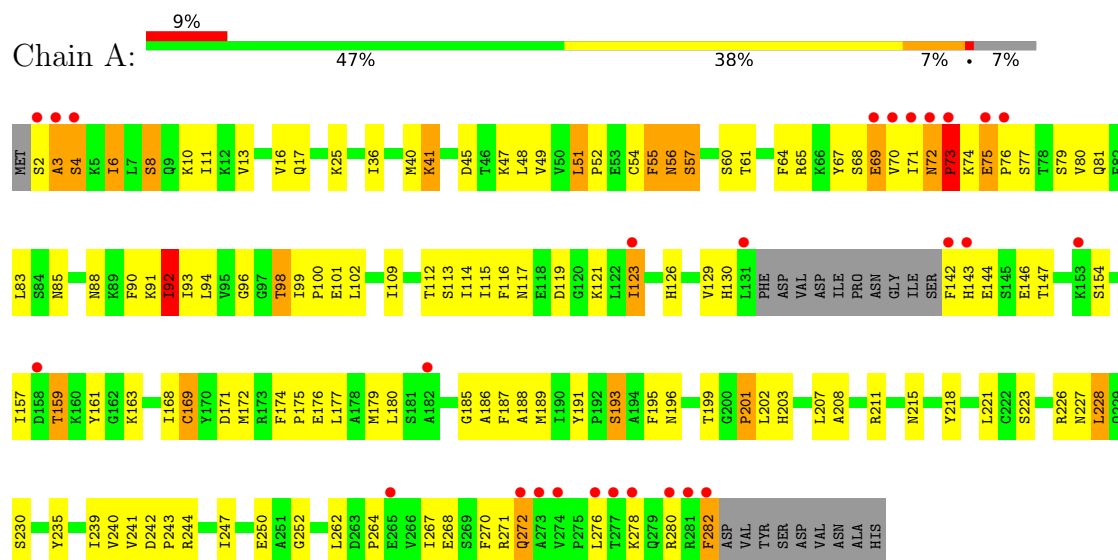
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	62	Total	O	0	0
			62	62		
2	B	55	Total	O	0	0
			55	55		

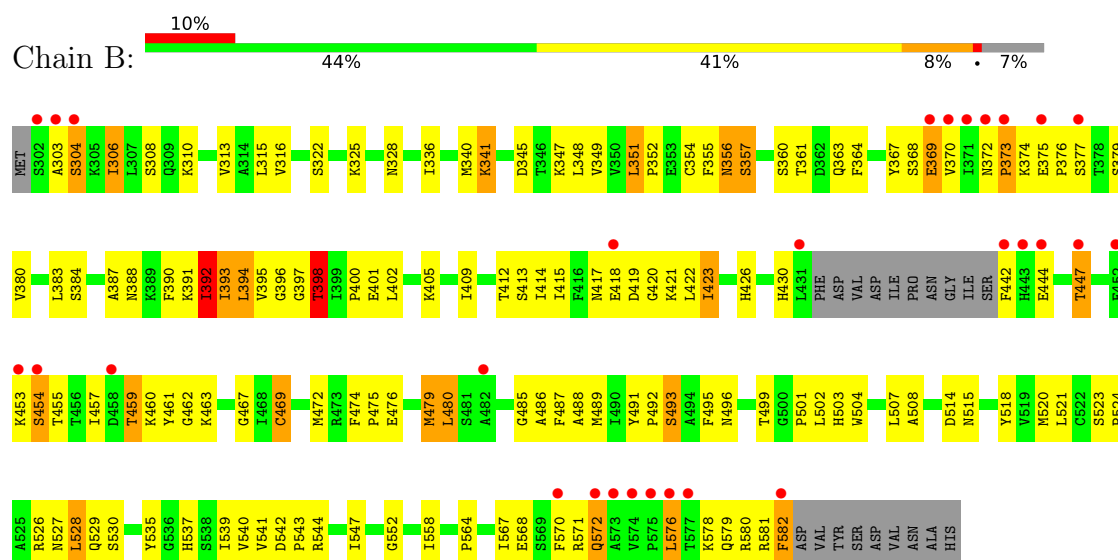
3 Residue-property plots

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: 32.5 KDA PROTEIN YLR351C



• Molecule 1: 32.5 KDA PROTEIN YLR351C



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.52Å 53.49Å 80.98Å 90.00° 112.43° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 50.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	71.7 (50.00-2.40) 85.2 (50.00-2.40)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.229 , 0.258 0.230 , 0.253	Depositor DCC
R_{free} test set	811 reflections (3.31%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtriage
Anisotropy	0.588	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4391	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 9.96% 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	0.52	0/2183	1.07	13/2950 (0.4%)
1	B	0.52	1/2183 (0.0%)	1.09	16/2950 (0.5%)
All	All	0.52	1/4366 (0.0%)	1.08	29/5900 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	479	MET	SD-CE	-5.08	1.66	1.79

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	306	ILE	N-CA-C	-8.55	104.41	112.96
1	A	93	ILE	N-CA-C	-8.11	96.14	107.99
1	B	393	ILE	N-CA-C	-8.05	96.23	107.99
1	B	392	ILE	N-CA-C	7.85	125.67	109.34
1	A	6	ILE	N-CA-C	-7.56	105.40	112.96
1	A	92	ILE	N-CA-C	7.43	124.78	109.34
1	B	503	HIS	N-CA-C	7.15	123.12	113.97
1	B	447	THR	N-CA-C	-7.11	106.51	114.62
1	A	154	SER	N-CA-C	-6.13	99.66	109.96
1	A	272	GLN	N-CA-C	-6.02	104.64	111.14
1	B	504	TRP	N-CA-C	5.99	118.33	111.02
1	B	454	SER	N-CA-C	-5.92	99.60	109.24
1	B	357	SER	N-CA-C	5.85	114.86	108.14
1	A	203	HIS	N-CA-C	5.83	121.84	114.31
1	B	394	LEU	N-CA-C	5.61	117.67	108.52
1	A	240	VAL	N-CA-C	-5.60	99.43	107.78
1	A	72	ASN	N-CA-C	5.60	118.50	109.15
1	A	57	SER	N-CA-C	5.59	114.57	108.14
1	B	552	GLY	N-CA-C	-5.58	104.65	112.13
1	A	55	PHE	N-CA-C	5.49	118.19	111.82
1	A	74	LYS	N-CA-C	-5.39	104.28	111.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	355	PHE	N-CA-C	5.36	118.04	111.82
1	B	572	GLN	N-CA-C	-5.33	105.38	111.14
1	B	374	LYS	N-CA-C	-5.33	104.37	111.24
1	B	398	THR	N-CA-C	5.15	121.77	110.80
1	A	8	SER	N-CA-C	5.14	116.88	111.28
1	A	252	GLY	N-CA-C	-5.12	105.27	112.13
1	B	423	ILE	CB-CA-C	-5.12	105.41	111.81
1	B	455	THR	N-CA-C	5.08	117.02	108.73

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2137	0	2156	112	0
1	B	2137	0	2156	130	0
2	A	62	0	0	2	0
2	B	55	0	0	4	0
All	All	4391	0	4312	233	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 27.

All (233) 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:417:ASN:HD21	1:B:421:LYS:HB3	1.04	1.14
1:A:117:ASN:HD21	1:A:121:LYS:HB3	1.00	1.08
1:A:70:VAL:HG21	1:A:99:ILE:HD12	1.36	1.07
1:A:117:ASN:ND2	1:A:121:LYS:HB3	1.78	0.97
1:B:417:ASN:ND2	1:B:421:LYS:HB3	1.81	0.95
1:B:576:LEU:O	1:B:578:LYS:HG3	1.72	0.89
1:B:541:VAL:HG12	1:B:547:ILE:HG12	1.54	0.89
1:A:159:THR:HG22	1:A:161:TYR:H	1.37	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:VAL:HG11	1:A:79:SER:HB2	1.56	0.87
1:A:169:CYS:HA	1:A:193:SER:HB3	1.57	0.86
1:B:368:SER:HA	1:B:400:PRO:HB2	1.58	0.86
1:B:306:ILE:HG22	1:B:487:PHE:HB3	1.60	0.83
1:A:175:PRO:HB2	1:A:179:MET:HE2	1.63	0.81
1:B:475:PRO:HB2	1:B:479:MET:HE2	1.62	0.80
1:B:398:THR:HA	1:B:412:THR:O	1.82	0.78
1:A:68:SER:HA	1:A:100:PRO:HB2	1.63	0.78
1:A:6:ILE:HG22	1:A:187:PHE:HB3	1.67	0.76
1:A:98:THR:HA	1:A:112:THR:O	1.86	0.76
1:A:75:GLU:O	1:A:75:GLU:HG2	1.85	0.76
1:A:68:SER:O	1:A:70:VAL:N	2.20	0.75
1:B:576:LEU:HD23	1:B:576:LEU:H	1.50	0.75
1:B:370:VAL:HG21	1:B:379:SER:HB2	1.69	0.75
1:B:368:SER:O	1:B:370:VAL:N	2.20	0.74
1:A:88:ASN:HB3	1:A:119:ASP:C	2.13	0.74
1:B:578:LYS:HB3	2:B:9:HOH:O	1.88	0.74
1:B:370:VAL:HG13	1:B:377:SER:OG	1.88	0.73
1:A:143:HIS:HA	1:A:146:GLU:HB2	1.70	0.73
1:B:417:ASN:HB3	1:B:423:ILE:HD11	1.72	0.71
1:A:239:ILE:HG22	1:A:250:GLU:HB2	1.73	0.71
1:A:282:PHE:CE2	1:B:476:GLU:HG2	2.25	0.71
1:B:341:LYS:HE3	1:B:341:LYS:O	1.91	0.70
1:B:457:ILE:HD12	1:B:457:ILE:H	1.57	0.70
1:B:388:ASN:HB3	1:B:419:ASP:C	2.16	0.69
1:B:459:THR:HG22	1:B:461:TYR:H	1.56	0.69
1:A:276:LEU:H	1:A:276:LEU:HD23	1.59	0.68
1:A:83:LEU:HD12	1:A:114:ILE:HD11	1.76	0.67
1:B:457:ILE:HD12	1:B:457:ILE:N	2.11	0.66
1:A:282:PHE:HE2	1:B:476:GLU:HG2	1.59	0.66
1:A:117:ASN:HD21	1:A:121:LYS:CB	1.93	0.66
1:A:117:ASN:HB3	1:A:123:ILE:HD11	1.77	0.66
1:B:459:THR:CG2	1:B:461:TYR:H	2.08	0.65
1:B:310:LYS:HG2	1:B:564:PRO:HD2	1.77	0.65
1:B:508:ALA:CB	1:B:539:ILE:HD11	2.27	0.64
1:B:537:HIS:O	1:B:539:ILE:HG23	1.97	0.64
1:B:348:LEU:HD23	1:B:349:VAL:N	2.13	0.64
1:B:469:CYS:O	1:B:472:MET:HG2	1.98	0.63
1:B:398:THR:H	1:B:413:SER:HA	1.63	0.63
1:B:479:MET:HA	1:B:515:ASN:ND2	2.13	0.62
1:A:157:ILE:HD12	1:A:157:ILE:H	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:SER:HA	1:B:308:SER:HA	1.82	0.62
1:A:123:ILE:HG21	1:A:157:ILE:HG23	1.81	0.62
1:A:179:MET:HA	1:A:215:ASN:ND2	2.14	0.62
1:B:423:ILE:HG21	1:B:457:ILE:HG23	1.81	0.62
1:B:469:CYS:HA	1:B:493:SER:HB3	1.82	0.61
1:A:241:VAL:HG12	1:A:247:ILE:HG12	1.82	0.61
1:A:70:VAL:HG11	1:A:79:SER:CB	2.28	0.61
1:A:196:ASN:H	1:A:199:THR:HG22	1.66	0.61
1:A:157:ILE:HD12	1:A:157:ILE:N	2.14	0.61
1:B:313:VAL:HG12	1:B:348:LEU:HB3	1.83	0.60
1:B:306:ILE:HG23	1:B:485:GLY:O	2.00	0.60
1:A:6:ILE:HG23	1:A:185:GLY:O	2.01	0.60
1:B:496:ASN:H	1:B:499:THR:HG22	1.66	0.60
1:B:372:ASN:ND2	1:B:414:ILE:HG21	2.18	0.59
1:A:227:ASN:HD22	1:A:230:SER:HB3	1.66	0.59
1:A:98:THR:H	1:A:113:SER:HA	1.68	0.59
1:B:579:GLN:N	2:B:9:HOH:O	2.35	0.59
1:B:453:LYS:HG3	1:B:454:SER:N	2.17	0.58
1:A:10:LYS:HG2	1:A:264:PRO:HD2	1.84	0.58
1:B:489:MET:HE3	1:B:491:TYR:CZ	2.39	0.58
1:A:172:MET:HE1	1:A:221:LEU:HD11	1.86	0.58
1:A:71:ILE:HG12	1:A:102:LEU:O	2.04	0.57
1:A:4:SER:HA	1:A:8:SER:HA	1.86	0.57
1:B:370:VAL:HG13	1:B:377:SER:CB	2.35	0.57
1:A:172:MET:HE2	1:A:207:LEU:HB2	1.87	0.56
1:B:377:SER:H	1:B:380:VAL:HB	1.70	0.56
1:B:361:THR:O	1:B:447:THR:HG21	2.05	0.56
1:B:444:GLU:CD	1:B:444:GLU:H	2.14	0.56
1:B:367:TYR:C	1:B:369:GLU:H	2.14	0.56
1:B:351:LEU:HD22	1:B:394:LEU:HD11	1.88	0.56
1:A:51:LEU:HD22	1:A:94:LEU:HD11	1.86	0.56
1:B:316:VAL:O	1:B:352:PRO:HD3	2.05	0.56
1:B:354:CYS:H	1:B:398:THR:HG22	1.71	0.56
1:B:356:ASN:HD22	1:B:357:SER:N	2.04	0.56
1:A:239:ILE:HG22	1:A:250:GLU:CB	2.35	0.55
1:B:417:ASN:HD21	1:B:421:LYS:CB	1.97	0.55
1:A:41:LYS:HE3	1:A:41:LYS:O	2.07	0.55
1:B:377:SER:HB3	1:B:380:VAL:HG23	1.87	0.55
1:A:282:PHE:C	1:A:282:PHE:CD2	2.85	0.55
1:A:195:PHE:HB3	1:A:199:THR:HG23	1.90	0.54
1:B:493:SER:O	1:B:523:SER:HA	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:VAL:HG21	1:A:99:ILE:CD1	2.23	0.54
1:A:91:LYS:HA	2:A:305:HOH:O	2.07	0.53
1:A:227:ASN:ND2	1:A:230:SER:HB3	2.23	0.53
1:A:189:MET:HE3	1:A:191:TYR:CZ	2.43	0.53
1:A:163:LYS:O	1:A:186:ALA:HA	2.09	0.53
1:A:211:ARG:HH22	1:B:514:ASP:CG	2.16	0.53
1:A:115:ILE:N	1:A:115:ILE:HD12	2.25	0.52
1:B:567:ILE:O	1:B:571:ARG:HG3	2.10	0.52
1:A:201:PRO:HD3	1:A:235:TYR:CD1	2.45	0.52
1:A:227:ASN:HD22	1:A:230:SER:CB	2.23	0.52
1:A:48:LEU:HD23	1:A:49:VAL:N	2.25	0.52
1:A:201:PRO:HD3	1:A:235:TYR:CE1	2.45	0.52
1:A:278:LYS:HG2	1:B:474:PHE:CE2	2.44	0.52
1:A:244:ARG:HD3	1:A:270:PHE:CE1	2.45	0.52
1:A:282:PHE:C	1:A:282:PHE:HD2	2.18	0.52
1:B:347:LYS:HD2	1:B:391:LYS:HD3	1.92	0.51
1:B:415:ILE:O	1:B:422:LEU:HD12	2.10	0.51
1:B:496:ASN:H	1:B:499:THR:CG2	2.24	0.51
1:A:243:PRO:HG2	1:A:267:ILE:HG12	1.92	0.51
1:A:61:THR:O	1:A:147:THR:HG21	2.11	0.51
1:A:36:ILE:O	1:A:40:MET:HG2	2.10	0.51
1:A:54:CYS:H	1:A:98:THR:HG22	1.76	0.50
1:B:475:PRO:C	1:B:479:MET:HE3	2.37	0.50
1:A:188:ALA:HA	1:A:218:TYR:O	2.11	0.50
1:B:415:ILE:HD12	1:B:415:ILE:N	2.27	0.50
1:A:174:PHE:CE2	1:B:578:LYS:HG2	2.46	0.50
1:B:383:LEU:HD12	1:B:414:ILE:HD11	1.94	0.50
1:B:395:VAL:C	1:B:397:GLY:H	2.20	0.50
1:A:16:VAL:O	1:A:52:PRO:HD3	2.12	0.49
1:A:242:ASP:HB2	1:A:243:PRO:HD2	1.94	0.49
1:B:579:GLN:C	1:B:581:ARG:H	2.21	0.49
1:B:520:MET:HG2	1:B:540:VAL:HG22	1.94	0.49
1:B:488:ALA:HA	1:B:518:TYR:O	2.11	0.49
1:A:77:SER:HB3	1:A:80:VAL:HG23	1.93	0.49
1:B:356:ASN:HD22	1:B:356:ASN:C	2.20	0.49
1:A:56:ASN:HD22	1:A:57:SER:N	2.11	0.49
1:A:73:PRO:HG2	1:A:75:GLU:O	2.13	0.49
1:B:377:SER:HB3	1:B:380:VAL:CG2	2.43	0.49
1:A:60:SER:O	1:A:64:PHE:HD1	1.96	0.49
1:A:196:ASN:H	1:A:199:THR:CG2	2.25	0.49
1:B:472:MET:HE1	1:B:521:LEU:HD11	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:LYS:NZ	1:A:70:VAL:HG13	2.28	0.49
1:A:282:PHE:HE1	1:B:430:HIS:CE1	2.31	0.48
1:B:582:PHE:C	1:B:582:PHE:CD2	2.91	0.48
1:A:41:LYS:HD3	1:A:41:LYS:C	2.38	0.48
1:B:582:PHE:C	1:B:582:PHE:HD2	2.22	0.48
1:B:322:SER:H	1:B:328:ASN:HD21	1.62	0.48
1:A:72:ASN:ND2	1:A:114:ILE:HG21	2.29	0.48
1:A:77:SER:HB3	1:A:80:VAL:CG2	2.44	0.48
1:B:356:ASN:C	1:B:356:ASN:ND2	2.70	0.48
1:B:368:SER:C	1:B:370:VAL:N	2.72	0.47
1:A:193:SER:O	1:A:223:SER:HA	2.14	0.47
1:B:340:MET:HG3	1:B:390:PHE:CD2	2.49	0.47
1:B:367:TYR:C	1:B:369:GLU:N	2.73	0.47
1:B:368:SER:C	1:B:370:VAL:H	2.22	0.47
1:B:402:LEU:HD13	1:B:409:ILE:CG2	2.45	0.47
1:B:463:LYS:O	1:B:486:ALA:HA	2.14	0.47
1:B:495:PHE:HB3	1:B:499:THR:HG23	1.95	0.47
1:B:542:ASP:HB2	1:B:543:PRO:HD2	1.95	0.47
1:A:67:TYR:C	1:A:69:GLU:H	2.22	0.47
1:B:543:PRO:HG2	1:B:567:ILE:HG12	1.95	0.47
1:A:268:GLU:O	1:A:272:GLN:HG2	2.15	0.47
1:A:278:LYS:HG2	1:B:474:PHE:HE2	1.80	0.47
1:B:527:ASN:HD22	1:B:530:SER:CB	2.28	0.46
1:A:2:SER:O	1:A:3:ALA:HB2	2.15	0.46
1:A:130:HIS:CE1	1:B:582:PHE:HE1	2.33	0.46
1:A:242:ASP:HB2	1:A:243:PRO:CD	2.44	0.46
1:A:90:PHE:HB2	1:A:92:ILE:HG12	1.98	0.46
1:B:467:GLY:O	1:B:492:PRO:HD2	2.15	0.46
1:A:70:VAL:HG12	1:A:77:SER:OG	2.15	0.46
1:B:472:MET:HE2	1:B:507:LEU:HB2	1.96	0.46
1:B:541:VAL:CG1	1:B:547:ILE:HG12	2.36	0.46
1:B:541:VAL:HG12	1:B:547:ILE:CG1	2.35	0.46
1:B:418:GLU:OE1	1:B:460:LYS:HD3	2.16	0.46
1:A:101:GLU:O	1:A:109:ILE:HA	2.15	0.46
1:B:578:LYS:CB	2:B:9:HOH:O	2.57	0.45
1:B:390:PHE:HB2	1:B:392:ILE:HG12	1.98	0.45
1:A:176:GLU:HA	1:A:179:MET:HE3	1.99	0.45
1:B:457:ILE:H	1:B:457:ILE:CD1	2.27	0.45
1:B:304:SER:CA	1:B:308:SER:HA	2.46	0.45
1:B:336:ILE:O	1:B:340:MET:HG2	2.17	0.45
1:B:460:LYS:HE2	1:B:461:TYR:CZ	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:395:VAL:O	1:B:397:GLY:N	2.50	0.45
1:B:304:SER:HB3	1:B:308:SER:HA	1.99	0.45
1:B:527:ASN:HD22	1:B:530:SER:HB3	1.80	0.45
1:B:354:CYS:H	1:B:398:THR:CG2	2.29	0.45
1:B:542:ASP:HB2	1:B:543:PRO:CD	2.47	0.45
1:A:169:CYS:O	1:A:172:MET:HG2	2.17	0.44
1:A:65:ARG:HB3	2:A:313:HOH:O	2.18	0.44
1:B:568:GLU:O	1:B:572:GLN:HG2	2.17	0.44
1:A:11:ILE:HD11	1:A:262:LEU:HD12	2.00	0.44
1:B:375:GLU:H	1:B:375:GLU:HG2	1.58	0.44
1:A:56:ASN:ND2	1:A:56:ASN:C	2.74	0.43
1:B:370:VAL:HG21	1:B:379:SER:CB	2.44	0.43
1:B:412:THR:HG23	1:B:426:HIS:C	2.43	0.43
1:B:401:GLU:O	1:B:409:ILE:HA	2.18	0.43
1:A:47:LYS:HD2	1:A:91:LYS:HD3	2.01	0.43
1:A:98:THR:OG1	1:A:168:ILE:HD13	2.18	0.43
1:A:81:GLN:HG3	1:A:85:ASN:HD21	1.84	0.43
1:B:315:LEU:HD12	1:B:558:ILE:HD11	2.00	0.43
1:B:325:LYS:NZ	1:B:370:VAL:HG23	2.34	0.43
1:A:112:THR:HG23	1:A:126:HIS:C	2.44	0.43
1:A:271:ARG:HD3	1:A:276:LEU:HD22	2.00	0.43
1:A:144:GLU:CD	1:A:144:GLU:H	2.26	0.42
1:B:348:LEU:HD23	1:B:348:LEU:C	2.44	0.42
1:A:67:TYR:C	1:A:69:GLU:N	2.76	0.42
1:A:13:VAL:HG12	1:A:48:LEU:HB3	2.01	0.42
1:B:367:TYR:HA	1:B:369:GLU:HG3	2.02	0.42
1:B:526:ARG:HG2	1:B:528:LEU:HD13	2.01	0.42
1:B:372:ASN:HD21	1:B:414:ILE:HG21	1.81	0.42
1:A:68:SER:C	1:A:70:VAL:N	2.76	0.42
1:A:60:SER:O	1:A:64:PHE:CD1	2.71	0.42
1:B:480:LEU:HD12	1:B:480:LEU:HA	1.91	0.42
1:A:172:MET:HE2	1:A:207:LEU:CB	2.50	0.42
1:B:363:GLN:OE1	1:B:363:GLN:HA	2.19	0.42
1:B:493:SER:O	1:B:524:PRO:HD3	2.20	0.42
1:B:544:ARG:HD3	1:B:570:PHE:CE1	2.55	0.42
1:A:25:LYS:HZ1	1:A:70:VAL:HG13	1.85	0.42
1:A:171:ASP:HB3	1:A:177:LEU:CD2	2.49	0.41
1:A:55:PHE:CD1	1:A:55:PHE:C	2.98	0.41
1:A:226:ARG:HG2	1:A:228:LEU:HD13	2.02	0.41
1:A:17:GLN:HG3	1:A:223:SER:O	2.20	0.41
1:A:174:PHE:HE2	1:B:578:LYS:HG2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:SER:CB	1:B:308:SER:HA	2.51	0.41
1:A:56:ASN:HD22	1:A:56:ASN:C	2.28	0.41
1:A:102:LEU:HD13	1:A:109:ILE:CG2	2.50	0.41
1:B:387:ALA:HB3	1:B:420:GLY:HA2	2.02	0.41
1:A:68:SER:C	1:A:70:VAL:H	2.28	0.41
1:B:384:SER:O	1:B:387:ALA:HB3	2.21	0.41
1:B:501:PRO:HG3	1:B:535:TYR:CE1	2.56	0.41
1:B:579:GLN:C	1:B:581:ARG:N	2.78	0.41
1:A:54:CYS:H	1:A:98:THR:CG2	2.33	0.41
1:A:72:ASN:HD21	1:A:114:ILE:HG21	1.86	0.41
1:A:114:ILE:HD12	1:A:116:PHE:CZ	2.56	0.41
1:B:360:SER:O	1:B:364:PHE:HD1	2.03	0.41
1:B:459:THR:HG22	1:B:462:GLY:H	1.85	0.41
1:B:579:GLN:O	1:B:581:ARG:N	2.54	0.41
1:B:393:ILE:HD13	1:B:459:THR:OG1	2.20	0.40
1:B:527:ASN:ND2	1:B:530:SER:HB3	2.36	0.40
1:B:361:THR:HA	1:B:364:PHE:CD1	2.56	0.40
1:B:479:MET:HA	1:B:515:ASN:HD21	1.83	0.40
1:B:582:PHE:CZ	2:B:9:HOH:O	2.57	0.40
1:A:208:ALA:HB2	1:A:221:LEU:HD11	2.04	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	267/291 (92%)	239 (90%)	16 (6%)	12 (4%)	2	1
1	B	267/291 (92%)	243 (91%)	13 (5%)	11 (4%)	2	2
All	All	534/582 (92%)	482 (90%)	29 (5%)	23 (4%)	2	1

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	69	GLU
1	A	73	PRO
1	A	76	PRO
1	A	92	ILE
1	B	369	GLU
1	B	373	PRO
1	B	392	ILE
1	A	3	ALA
1	A	98	THR
1	B	303	ALA
1	B	396	GLY
1	B	398	THR
1	B	469	CYS
1	A	4	SER
1	A	169	CYS
1	B	304	SER
1	B	580	ARG
1	A	280	ARG
1	A	96	GLY
1	A	123	ILE
1	B	376	PRO
1	B	405	LYS
1	A	129	VAL

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	237/255 (93%)	223 (94%)	14 (6%)	18	31
1	B	237/255 (93%)	223 (94%)	14 (6%)	18	31
All	All	474/510 (93%)	446 (94%)	28 (6%)	18	31

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LYS

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Mol	Chain	Res	Type
1	A	45	ASP
1	A	51	LEU
1	A	56	ASN
1	A	73	PRO
1	A	75	GLU
1	A	142	PHE
1	A	159	THR
1	A	180	LEU
1	A	193	SER
1	A	201	PRO
1	A	202	LEU
1	A	228	LEU
1	A	282	PHE
1	B	341	LYS
1	B	345	ASP
1	B	351	LEU
1	B	356	ASN
1	B	373	PRO
1	B	442	PHE
1	B	459	THR
1	B	480	LEU
1	B	493	SER
1	B	502	LEU
1	B	528	LEU
1	B	529	GLN
1	B	576	LEU
1	B	582	PHE

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	28	ASN
1	A	56	ASN
1	A	72	ASN
1	A	81	GLN
1	A	85	ASN
1	A	227	ASN
1	A	272	GLN
1	B	328	ASN
1	B	356	ASN
1	B	372	ASN
1	B	381	GLN

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Mol	Chain	Res	Type
1	B	385	ASN
1	B	430	HIS
1	B	515	ASN
1	B	527	ASN
1	B	572	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 [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	271/291 (93%)	0.75	27 (9%) 12 9	13, 34, 58, 68	0
1	B	271/291 (93%)	0.76	29 (10%) 11 8	12, 33, 58, 69	0
All	All	542/582 (93%)	0.75	56 (10%) 12 9	12, 34, 58, 69	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	369	GLU	5.4
1	A	70	VAL	5.4
1	B	370	VAL	5.2
1	B	443	HIS	4.9
1	B	442	PHE	4.5
1	B	573	ALA	4.3
1	A	69	GLU	4.2
1	B	304	SER	4.2
1	A	2	SER	4.2
1	A	4	SER	4.1
1	A	142	PHE	4.1
1	B	576	LEU	4.0
1	B	582	PHE	3.8
1	A	143	HIS	3.8
1	A	282	PHE	3.7
1	A	277	THR	3.5
1	B	373	PRO	3.5
1	B	482	ALA	3.5
1	A	273	ALA	3.3
1	B	577	THR	3.3
1	A	3	ALA	3.3
1	B	302	SER	3.2
1	B	458	ASP	3.2
1	A	73	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	76	PRO	3.0
1	A	131	LEU	3.0
1	A	281	ARG	3.0
1	A	274	VAL	2.8
1	B	303	ALA	2.8
1	B	372	ASN	2.8
1	A	276	LEU	2.8
1	A	158	ASP	2.7
1	B	574	VAL	2.6
1	B	444	GLU	2.6
1	A	72	ASN	2.6
1	A	75	GLU	2.6
1	A	182	ALA	2.5
1	A	71	ILE	2.5
1	B	454	SER	2.5
1	B	575	PRO	2.5
1	B	452	GLU	2.4
1	A	272	GLN	2.4
1	A	153	LYS	2.3
1	A	123	ILE	2.3
1	B	371	ILE	2.3
1	B	418	GLU	2.3
1	B	453	LYS	2.3
1	B	572	GLN	2.2
1	A	280	ARG	2.2
1	B	431	LEU	2.2
1	B	447	THR	2.2
1	B	570	PHE	2.1
1	A	265	GLU	2.1
1	B	375	GLU	2.1
1	A	278	LYS	2.0
1	B	377	SER	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.