



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 07:43 AM UTC

PDB ID : 3MGU / pdb_00003mgu
Title : Structure of *S. cerevisiae* Tpa1 protein, a proline hydroxylase modifying ribosomal protein Rps23
Authors : Henri, J.; Rispal, D.; Bayart, E.; van Tilbeurgh, H.; Seraphin, B.; Graille, M.
Deposited on : 2010-04-07
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

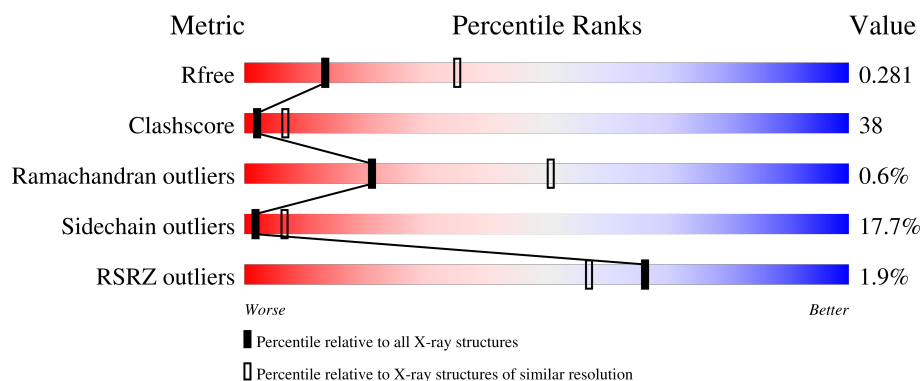
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	644	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4406 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 PKHD-type hydroxylase TPA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	533	Total	C	N	O	S	0	0	0
			4372	2813	730	816	13			

- Molecule 2 is FE (II) ION (CCD ID: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total	O	0	0
			33	33		

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	81.18Å 104.84Å 205.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.80 19.92 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.3 (19.92-2.80) 99.2 (19.92-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.239 , 0.296 0.232 , 0.281	Depositor DCC
R_{free} test set	1106 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.861	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4406	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 5.31% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FE2

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	0.40	0/4480	0.84	3/6048 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	158	THR	N-CA-C	6.17	119.41	110.28
1	A	221	LEU	CA-C-N	5.68	126.01	119.93
1	A	221	LEU	C-N-CA	5.68	126.01	119.93

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	4372	0	4307	326	0
2	A	1	0	0	0	0
3	A	33	0	0	3	0
All	All	4406	0	4307	326	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 38.

All (326) 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:202:SER:HA	1:A:409:PRO:HG3	1.28	1.15
1:A:424:ASN:HD21	1:A:426:GLN:HE21	1.04	0.97
1:A:289:ARG:HG2	1:A:289:ARG:HH11	1.26	0.96
1:A:371:LEU:H	1:A:597:ASN:HD21	1.08	0.96
1:A:175:PRO:HD2	1:A:212:PHE:HA	1.50	0.94
1:A:507:PRO:HA	1:A:615:VAL:HG12	1.50	0.93
1:A:69:ALA:HB1	1:A:112:ASN:HD21	1.37	0.89
1:A:461:ARG:HD3	1:A:461:ARG:H	1.37	0.88
1:A:535:GLU:OE1	1:A:604:ARG:HD2	1.74	0.87
1:A:395:HIS:CD2	1:A:397:LYS:HG2	2.10	0.86
1:A:200:PRO:HG2	1:A:409:PRO:HD2	1.58	0.86
1:A:479:LYS:HG3	1:A:500:ILE:HD12	1.59	0.84
1:A:31:ASN:HA	3:A:668:HOH:O	1.79	0.81
1:A:289:ARG:HH11	1:A:289:ARG:CG	1.97	0.77
1:A:337:TYR:O	1:A:340:GLN:HG2	1.85	0.76
1:A:442:ASN:ND2	1:A:445:PHE:H	1.83	0.75
1:A:285:PRO:HG3	1:A:416:TYR:CZ	2.21	0.75
1:A:87:ILE:O	1:A:157:LEU:HD22	1.86	0.75
1:A:417:ILE:HG23	1:A:502:ILE:HG12	1.69	0.74
1:A:91:ASN:HD21	1:A:151:THR:HG22	1.50	0.73
1:A:394:TYR:HD1	1:A:509:MET:SD	2.10	0.73
1:A:181:TRP:H	1:A:211:GLN:HE22	1.36	0.73
1:A:394:TYR:CD1	1:A:509:MET:SD	2.82	0.72
1:A:177:PRO:HG3	1:A:212:PHE:CE2	2.23	0.72
1:A:424:ASN:ND2	1:A:426:GLN:HE21	1.84	0.71
1:A:277:ASN:C	1:A:277:ASN:HD22	1.98	0.71
1:A:69:ALA:CB	1:A:112:ASN:HD21	2.03	0.71
1:A:27:LYS:HE3	1:A:27:LYS:N	2.06	0.71
1:A:560:ASP:O	1:A:561:ASP:HB2	1.89	0.71
1:A:138:LEU:HD23	1:A:168:ILE:HD12	1.73	0.69
1:A:356:GLN:O	1:A:360:VAL:HG23	1.92	0.69
1:A:126:ARG:HB3	1:A:138:LEU:HD13	1.73	0.69
1:A:535:GLU:OE2	1:A:627:LYS:HE2	1.92	0.69
1:A:169:SER:HB2	1:A:244:TRP:NE1	2.09	0.68
1:A:220:VAL:HG12	1:A:220:VAL:O	1.92	0.68
1:A:157:LEU:HA	1:A:228:ASP:HB2	1.76	0.67
1:A:166:ARG:HD3	1:A:244:TRP:CD1	2.29	0.67
1:A:483:PHE:CE2	1:A:487:LEU:HD23	2.28	0.67
1:A:67:LEU:HD21	1:A:172:LEU:HD21	1.77	0.66
1:A:386:LYS:HZ3	1:A:464:ASN:HA	1.59	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:413:ARG:O	1:A:510:ASP:HA	1.94	0.66
1:A:371:LEU:H	1:A:597:ASN:ND2	1.90	0.66
1:A:511:PHE:HB3	1:A:614:TYR:HA	1.78	0.66
1:A:474:LEU:O	1:A:478:HIS:HD2	1.79	0.65
1:A:156:LEU:HB3	1:A:229:VAL:HG13	1.76	0.65
1:A:162:VAL:HG22	1:A:220:VAL:HG13	1.79	0.65
1:A:51:GLN:HG2	1:A:52:PRO:HA	1.77	0.65
1:A:385:GLN:HE22	1:A:463:ASN:ND2	1.95	0.65
1:A:73:GLU:CD	1:A:111:PRO:HB2	2.22	0.64
1:A:193:PRO:HD3	1:A:205:SER:OG	1.97	0.64
1:A:395:HIS:HD2	1:A:397:LYS:HG2	1.59	0.64
1:A:522:LEU:H	1:A:522:LEU:HD12	1.62	0.64
1:A:285:PRO:HG3	1:A:416:TYR:CE1	2.32	0.64
1:A:494:CYS:HB2	1:A:633:LYS:HG2	1.79	0.64
1:A:73:GLU:OE2	1:A:111:PRO:HB2	1.98	0.63
1:A:402:PRO:HB2	1:A:417:ILE:HD13	1.80	0.63
1:A:505:PHE:O	1:A:622:SER:HB3	1.98	0.63
1:A:139:SER:HB2	1:A:248:PRO:HA	1.81	0.63
1:A:442:ASN:HD22	1:A:445:PHE:H	1.47	0.63
1:A:208:LEU:HD23	1:A:208:LEU:H	1.62	0.62
1:A:123:LYS:HE2	1:A:127:ASP:OD2	1.99	0.62
1:A:546:TRP:CE3	1:A:594:PRO:HG2	2.36	0.61
1:A:558:MET:HB3	1:A:588:VAL:HA	1.82	0.61
1:A:47:ILE:HD12	1:A:132:VAL:O	2.00	0.61
1:A:353:GLU:O	1:A:357:LYS:HD3	2.01	0.61
1:A:187:GLY:O	1:A:238:ARG:NH1	2.33	0.61
1:A:460:SER:HB2	1:A:462:GLU:HG2	1.83	0.60
1:A:353:GLU:HA	1:A:356:GLN:HE21	1.66	0.60
1:A:371:LEU:N	1:A:597:ASN:HD21	1.91	0.60
1:A:507:PRO:HA	1:A:615:VAL:CG1	2.27	0.60
1:A:91:ASN:ND2	1:A:151:THR:HG22	2.16	0.60
1:A:446:THR:O	1:A:450:ILE:HD13	2.01	0.60
1:A:447:LYS:HE3	1:A:451:LYS:NZ	2.16	0.60
1:A:36:ASP:C	1:A:36:ASP:OD1	2.44	0.59
1:A:26:ILE:HG13	1:A:27:LYS:CE	2.32	0.59
1:A:142:LYS:HD3	1:A:246:HIS:ND1	2.17	0.59
1:A:277:ASN:HD22	1:A:278:VAL:N	2.00	0.59
1:A:52:PRO:HG3	1:A:205:SER:OG	2.02	0.59
1:A:209:VAL:HG23	1:A:209:VAL:O	2.03	0.59
1:A:80:PHE:HD1	1:A:92:GLN:O	1.86	0.59
1:A:74:ILE:HG12	1:A:78:ILE:HD12	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:SER:O	1:A:413:ARG:NH1	2.36	0.58
1:A:174:LEU:N	1:A:175:PRO:HD3	2.17	0.58
1:A:67:LEU:HA	1:A:70:VAL:HG22	1.85	0.58
1:A:289:ARG:HD2	1:A:629:LYS:HB2	1.84	0.58
1:A:29:MET:HA	3:A:657:HOH:O	2.04	0.58
1:A:133:THR:HG21	1:A:168:ILE:CD1	2.34	0.58
1:A:173:TYR:C	1:A:174:LEU:HD22	2.29	0.58
1:A:418:ASP:HB3	1:A:501:LEU:CD2	2.33	0.58
1:A:395:HIS:HD2	1:A:397:LYS:H	1.52	0.57
1:A:461:ARG:HD3	1:A:461:ARG:N	2.12	0.57
1:A:162:VAL:HG12	1:A:162:VAL:O	2.04	0.57
1:A:520:ASN:HB3	1:A:523:LEU:HG	1.86	0.57
1:A:370:PHE:HD2	1:A:371:LEU:HD13	1.69	0.57
1:A:60:GLU:HA	1:A:213:ASN:O	2.04	0.56
1:A:64:ASP:O	1:A:68:ARG:HG3	2.04	0.56
1:A:36:ASP:OD1	1:A:38:THR:HG23	2.06	0.56
1:A:72:LYS:O	1:A:76:THR:HG23	2.05	0.56
1:A:74:ILE:HG23	1:A:78:ILE:HD12	1.87	0.56
1:A:303:MET:SD	1:A:485:LYS:HD3	2.45	0.56
1:A:364:SER:HA	1:A:601:LEU:O	2.05	0.56
1:A:424:ASN:C	1:A:424:ASN:HD22	2.14	0.56
1:A:121:TYR:CD2	1:A:143:THR:HG22	2.40	0.56
1:A:390:GLN:HG3	1:A:391:GLU:H	1.70	0.56
1:A:196:LEU:HB2	1:A:199:VAL:HB	1.87	0.56
1:A:26:ILE:N	1:A:27:LYS:HZ1	2.04	0.56
1:A:342:ILE:HG13	1:A:343:SER:N	2.20	0.56
1:A:386:LYS:O	1:A:390:GLN:HB3	2.06	0.56
1:A:453:ILE:O	1:A:454:SER:CB	2.53	0.56
1:A:195:ILE:O	1:A:196:LEU:HG	2.06	0.55
1:A:91:ASN:HD21	1:A:151:THR:CG2	2.19	0.55
1:A:342:ILE:HD11	1:A:346:HIS:HB2	1.89	0.55
1:A:137:LYS:HG2	1:A:254:GLY:HA3	1.88	0.55
1:A:174:LEU:HD21	1:A:241:ILE:HD12	1.87	0.55
1:A:547:GLU:HA	1:A:548:SER:OG	2.06	0.55
1:A:266:ASN:C	1:A:266:ASN:ND2	2.65	0.55
1:A:169:SER:HB2	1:A:244:TRP:HE1	1.70	0.54
1:A:522:LEU:HD12	1:A:522:LEU:N	2.23	0.54
1:A:546:TRP:O	1:A:593:PRO:HB3	2.06	0.54
1:A:173:TYR:HB3	1:A:175:PRO:HD3	1.88	0.54
1:A:513:LEU:O	1:A:515:THR:HG23	2.08	0.54
1:A:296:GLU:CD	1:A:481:THR:HG22	2.33	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:424:ASN:HD21	1:A:426:GLN:NE2	1.88	0.54
1:A:53:TYR:O	1:A:55:TRP:HD1	1.91	0.54
1:A:604:ARG:HB3	1:A:608:VAL:CG1	2.38	0.54
1:A:263:TRP:O	1:A:266:ASN:HB3	2.08	0.54
1:A:378:LEU:O	1:A:382:VAL:HG22	2.07	0.54
1:A:138:LEU:HD12	1:A:138:LEU:N	2.23	0.53
1:A:190:ARG:O	1:A:227:HIS:HA	2.08	0.53
1:A:464:ASN:O	1:A:468:GLU:CD	2.51	0.53
1:A:404:LYS:HE2	1:A:418:ASP:OD2	2.09	0.53
1:A:51:GLN:HG2	1:A:52:PRO:CA	2.38	0.53
1:A:250:VAL:HA	1:A:255:TYR:CE1	2.44	0.52
1:A:163:ILE:HD11	1:A:166:ARG:CZ	2.39	0.52
1:A:157:LEU:HD22	1:A:157:LEU:H	1.74	0.52
1:A:385:GLN:OE1	1:A:463:ASN:HB2	2.09	0.52
1:A:208:LEU:H	1:A:208:LEU:CD2	2.22	0.52
1:A:261:GLU:HB2	1:A:265:ARG:HH12	1.75	0.52
1:A:390:GLN:HG3	1:A:391:GLU:N	2.23	0.52
1:A:447:LYS:HD2	1:A:447:LYS:C	2.34	0.52
1:A:67:LEU:HD13	1:A:174:LEU:HD13	1.91	0.52
1:A:523:LEU:N	1:A:523:LEU:HD23	2.25	0.52
1:A:447:LYS:HE3	1:A:451:LYS:HZ3	1.74	0.52
1:A:197:PRO:O	1:A:198:ASN:HB2	2.10	0.52
1:A:368:GLU:O	1:A:369:SER:C	2.53	0.52
1:A:385:GLN:HE22	1:A:463:ASN:CG	2.17	0.52
1:A:474:LEU:HD22	1:A:478:HIS:CD2	2.45	0.51
1:A:507:PRO:O	1:A:615:VAL:O	2.27	0.51
1:A:385:GLN:HE21	1:A:389:GLU:HB3	1.75	0.51
1:A:399:VAL:C	1:A:400:LYS:HD2	2.36	0.51
1:A:135:ALA:HB3	1:A:168:ILE:HD11	1.92	0.51
1:A:144:ASP:O	1:A:144:ASP:OD2	2.28	0.51
1:A:171:ILE:HD13	1:A:173:TYR:CZ	2.46	0.51
1:A:503:ARG:HH11	1:A:503:ARG:CG	2.23	0.51
1:A:87:ILE:HG22	1:A:157:LEU:CD2	2.40	0.51
1:A:460:SER:HB2	1:A:462:GLU:CG	2.41	0.51
1:A:71:ARG:NH1	1:A:175:PRO:O	2.44	0.51
1:A:162:VAL:HG22	1:A:220:VAL:CG1	2.40	0.51
1:A:560:ASP:O	1:A:561:ASP:CB	2.57	0.51
1:A:53:TYR:OH	1:A:191:LEU:HD12	2.10	0.50
1:A:92:GLN:O	1:A:93:SER:HB3	2.11	0.50
1:A:92:GLN:HA	1:A:148:ASN:ND2	2.27	0.50
1:A:266:ASN:HD22	1:A:267:ASN:N	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:LEU:O	1:A:207:LYS:HA	2.10	0.50
1:A:497:SER:OG	1:A:631:ASP:HB2	2.12	0.50
1:A:247:ILE:O	1:A:249:GLN:NE2	2.44	0.50
1:A:342:ILE:HG12	1:A:347:LEU:CD2	2.42	0.50
1:A:138:LEU:HD12	1:A:138:LEU:H	1.76	0.50
1:A:142:LYS:HG2	1:A:142:LYS:O	2.11	0.49
1:A:400:LYS:HD2	1:A:400:LYS:N	2.26	0.49
1:A:276:SER:O	1:A:279:LEU:HD23	2.13	0.49
1:A:87:ILE:HG22	1:A:157:LEU:HD21	1.94	0.49
1:A:461:ARG:H	1:A:461:ARG:CD	2.18	0.49
1:A:474:LEU:HD13	1:A:626:VAL:HG21	1.95	0.49
1:A:623:ARG:NH1	1:A:625:ASP:OD2	2.46	0.49
1:A:189:LEU:HB3	1:A:208:LEU:HD21	1.95	0.49
1:A:299:HIS:CE1	1:A:485:LYS:NZ	2.81	0.49
1:A:370:PHE:CD2	1:A:371:LEU:HD13	2.47	0.49
1:A:66:LEU:O	1:A:70:VAL:HG22	2.11	0.49
1:A:115:LYS:O	1:A:119:ILE:HG23	2.12	0.49
1:A:508:GLY:HA2	1:A:617:TRP:HD1	1.78	0.49
1:A:72:LYS:HZ2	1:A:72:LYS:N	2.10	0.49
1:A:74:ILE:HD13	1:A:147:ILE:HG21	1.95	0.49
1:A:68:ARG:O	1:A:72:LYS:HB2	2.13	0.48
1:A:163:ILE:HD11	1:A:166:ARG:NH2	2.28	0.48
1:A:158:THR:HG23	1:A:409:PRO:O	2.13	0.48
1:A:70:VAL:HG21	1:A:116:LEU:HD22	1.95	0.48
1:A:494:CYS:HB2	1:A:633:LYS:CG	2.42	0.48
1:A:296:GLU:OE1	1:A:481:THR:HG22	2.13	0.48
1:A:359:PHE:HB2	1:A:365:LEU:HD12	1.95	0.48
1:A:447:LYS:NZ	1:A:473:ASP:OD1	2.44	0.48
1:A:548:SER:HA	1:A:549:GLY:HA2	1.62	0.48
1:A:511:PHE:CE2	1:A:513:LEU:HD11	2.48	0.48
1:A:414:TYR:HA	1:A:510:ASP:CG	2.38	0.48
1:A:128:PHE:O	1:A:132:VAL:HG22	2.14	0.48
1:A:382:VAL:HG12	1:A:466:ASP:HB3	1.96	0.48
1:A:402:PRO:HG3	1:A:468:GLU:HG2	1.95	0.48
1:A:157:LEU:HD23	1:A:157:LEU:C	2.39	0.47
1:A:147:ILE:O	1:A:148:ASN:ND2	2.47	0.47
1:A:524:LYS:HD3	1:A:524:LYS:HA	1.65	0.47
1:A:137:LYS:N	1:A:137:LYS:HD2	2.30	0.47
1:A:151:THR:O	1:A:154:CYS:HB2	2.15	0.47
1:A:156:LEU:CB	1:A:229:VAL:HG13	2.43	0.47
1:A:546:TRP:O	1:A:547:GLU:HG2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:523:LEU:O	1:A:526:ASN:HB3	2.14	0.47
1:A:599:PHE:CE2	1:A:601:LEU:HG	2.50	0.47
1:A:173:TYR:CD2	1:A:210:PRO:HB3	2.50	0.47
1:A:294:PHE:C	1:A:294:PHE:CD2	2.92	0.47
1:A:169:SER:HB2	1:A:244:TRP:CD1	2.49	0.47
1:A:86:ASP:N	1:A:86:ASP:OD1	2.47	0.46
1:A:122:SER:O	1:A:126:ARG:HG3	2.16	0.46
1:A:380:LYS:HB2	1:A:596:TRP:CZ2	2.50	0.46
1:A:431:ILE:O	1:A:435:LEU:HD22	2.15	0.46
1:A:30:PHE:HZ	1:A:119:ILE:O	1.98	0.46
1:A:196:LEU:HD21	1:A:428:GLU:HG3	1.98	0.46
1:A:209:VAL:O	1:A:209:VAL:CG2	2.63	0.46
1:A:177:PRO:HG3	1:A:212:PHE:CZ	2.50	0.46
1:A:560:ASP:HB3	1:A:609:LEU:HD11	1.97	0.46
1:A:586:ASP:C	1:A:586:ASP:OD1	2.58	0.46
1:A:59:HIS:O	1:A:60:GLU:HG2	2.15	0.46
1:A:166:ARG:CZ	1:A:244:TRP:CE2	2.99	0.46
1:A:352:ILE:O	1:A:356:GLN:HG3	2.15	0.46
1:A:431:ILE:HG22	1:A:435:LEU:CD2	2.45	0.46
1:A:26:ILE:HG13	1:A:27:LYS:HE3	1.96	0.46
1:A:174:LEU:N	1:A:175:PRO:CD	2.78	0.46
1:A:468:GLU:OE1	1:A:468:GLU:N	2.49	0.46
1:A:62:VAL:CG2	1:A:213:ASN:HD21	2.29	0.45
1:A:163:ILE:O	1:A:164:GLY:C	2.58	0.45
1:A:604:ARG:HB3	1:A:608:VAL:HG13	1.98	0.45
1:A:26:ILE:C	1:A:27:LYS:HE3	2.40	0.45
1:A:396:SER:HB2	1:A:405:THR:HG21	1.97	0.45
1:A:124:GLN:H	1:A:124:GLN:NE2	2.13	0.45
1:A:613:LYS:O	1:A:614:TYR:C	2.59	0.45
1:A:155:HIS:HB3	1:A:231:GLU:HG3	1.99	0.45
1:A:422:TYR:O	1:A:423:ARG:HD3	2.17	0.45
1:A:385:GLN:NE2	1:A:463:ASN:ND2	2.64	0.45
1:A:375:LYS:HA	1:A:375:LYS:HD3	1.61	0.45
1:A:443:PHE:O	1:A:447:LYS:HB2	2.16	0.45
1:A:512:THR:C	1:A:513:LEU:HD12	2.41	0.45
1:A:153:GLY:HA2	1:A:231:GLU:OE1	2.17	0.44
1:A:354:LYS:HB3	1:A:354:LYS:HE3	1.56	0.44
1:A:596:TRP:CD1	1:A:597:ASN:HB2	2.52	0.44
1:A:78:ILE:CD1	1:A:147:ILE:HD12	2.48	0.44
1:A:303:MET:HE2	1:A:303:MET:HB3	1.90	0.44
1:A:554:TYR:HA	1:A:594:PRO:HD3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:LEU:HA	1:A:70:VAL:CG2	2.48	0.44
1:A:147:ILE:HG22	1:A:239:LEU:HD11	1.99	0.44
1:A:506:ARG:NH2	1:A:509:MET:HE2	2.32	0.44
1:A:126:ARG:HB3	1:A:138:LEU:CD1	2.43	0.44
1:A:173:TYR:O	1:A:174:LEU:HD22	2.17	0.44
1:A:261:GLU:C	1:A:263:TRP:H	2.25	0.44
1:A:88:TYR:CD1	1:A:90:VAL:HG12	2.53	0.44
1:A:59:HIS:C	1:A:60:GLU:HG2	2.42	0.44
1:A:520:ASN:OD1	1:A:522:LEU:HD13	2.17	0.44
1:A:530:ILE:HD12	1:A:531:ASP:H	1.83	0.44
1:A:113:LEU:HD12	1:A:113:LEU:O	2.17	0.43
1:A:547:GLU:HA	1:A:548:SER:CB	2.48	0.43
1:A:462:GLU:HG2	1:A:462:GLU:H	1.54	0.43
1:A:547:GLU:HA	1:A:548:SER:HA	1.83	0.43
1:A:31:ASN:HB3	1:A:34:ILE:HG23	2.00	0.43
1:A:192:PHE:HA	1:A:193:PRO:HD2	1.82	0.43
1:A:413:ARG:HD2	1:A:509:MET:HB2	2.00	0.43
1:A:493:LEU:HD13	1:A:632:VAL:HG11	2.00	0.43
1:A:620:LYS:HD2	1:A:620:LYS:HA	1.74	0.43
1:A:121:TYR:CG	1:A:143:THR:HG22	2.53	0.43
1:A:549:GLY:C	1:A:551:LEU:H	2.26	0.43
1:A:220:VAL:O	1:A:222:PRO:HD3	2.18	0.43
1:A:450:ILE:CG2	1:A:469:LEU:HD21	2.49	0.43
1:A:411:LYS:O	1:A:412:ALA:HB2	2.19	0.43
1:A:503:ARG:NH1	1:A:625:ASP:OD1	2.52	0.43
1:A:279:LEU:N	1:A:279:LEU:HD22	2.34	0.42
1:A:62:VAL:CG2	1:A:67:LEU:HG	2.49	0.42
1:A:342:ILE:HG12	1:A:347:LEU:HD21	2.00	0.42
1:A:450:ILE:HG22	1:A:469:LEU:HD21	2.01	0.42
1:A:604:ARG:HB3	1:A:608:VAL:HG11	1.99	0.42
1:A:471:LEU:HB3	1:A:502:ILE:HD12	2.02	0.42
1:A:556:LEU:O	1:A:610:GLU:HA	2.19	0.42
1:A:560:ASP:HB3	1:A:609:LEU:CD1	2.48	0.42
1:A:133:THR:HG21	1:A:168:ILE:HD11	2.00	0.42
1:A:289:ARG:HG2	1:A:289:ARG:NH1	2.08	0.42
1:A:289:ARG:CG	1:A:289:ARG:NH1	2.65	0.42
1:A:610:GLU:HG3	3:A:663:HOH:O	2.20	0.42
1:A:261:GLU:HB2	1:A:265:ARG:NH1	2.34	0.42
1:A:365:LEU:HD23	1:A:367:ILE:HG13	2.01	0.42
1:A:503:ARG:CG	1:A:503:ARG:NH1	2.83	0.42
1:A:613:LYS:HE3	1:A:614:TYR:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:464:ASN:O	1:A:468:GLU:OE1	2.37	0.42
1:A:27:LYS:C	1:A:29:MET:H	2.28	0.42
1:A:447:LYS:HD3	1:A:469:LEU:HD22	2.02	0.42
1:A:382:VAL:CG1	1:A:466:ASP:HB3	2.50	0.42
1:A:503:ARG:NH1	1:A:625:ASP:OD2	2.52	0.42
1:A:299:HIS:HE1	1:A:485:LYS:HZ2	1.68	0.41
1:A:129:PHE:CD2	1:A:245:TYR:CE2	3.08	0.41
1:A:284:PHE:O	1:A:406:ALA:HA	2.18	0.41
1:A:371:LEU:HB2	1:A:597:ASN:HD21	1.86	0.41
1:A:62:VAL:HG23	1:A:67:LEU:HG	2.02	0.41
1:A:129:PHE:CE2	1:A:245:TYR:HE2	2.38	0.41
1:A:173:TYR:CE2	1:A:210:PRO:HB3	2.55	0.41
1:A:511:PHE:HE2	1:A:513:LEU:HD11	1.85	0.41
1:A:228:ASP:OD1	1:A:228:ASP:C	2.62	0.41
1:A:294:PHE:CD2	1:A:295:HIS:N	2.88	0.41
1:A:591:ASN:O	1:A:593:PRO:HD3	2.20	0.41
1:A:196:LEU:O	1:A:197:PRO:C	2.61	0.41
1:A:472:ILE:O	1:A:476:VAL:HG23	2.21	0.41
1:A:70:VAL:HG12	1:A:113:LEU:HA	2.03	0.41
1:A:158:THR:HA	1:A:228:ASP:HB3	2.03	0.41
1:A:414:TYR:HB3	1:A:510:ASP:HB3	2.02	0.41
1:A:503:ARG:NH1	1:A:625:ASP:CG	2.79	0.41
1:A:87:ILE:H	1:A:87:ILE:HG12	1.48	0.41
1:A:261:GLU:CD	1:A:261:GLU:H	2.28	0.41
1:A:386:LYS:NZ	1:A:464:ASN:HA	2.31	0.41
1:A:451:LYS:HB3	1:A:451:LYS:HE2	1.77	0.41
1:A:124:GLN:H	1:A:124:GLN:HE21	1.69	0.41
1:A:129:PHE:O	1:A:133:THR:HB	2.21	0.41
1:A:202:SER:HA	1:A:409:PRO:CG	2.21	0.41
1:A:414:TYR:HA	1:A:510:ASP:OD1	2.21	0.41
1:A:516:LYS:HG3	1:A:609:LEU:HD23	2.03	0.41
1:A:486:TYR:CZ	1:A:490:LEU:HD21	2.56	0.41
1:A:150:TYR:OH	1:A:156:LEU:HB2	2.21	0.40
1:A:364:SER:CA	1:A:601:LEU:O	2.69	0.40
1:A:173:TYR:CE1	1:A:210:PRO:HG2	2.56	0.40
1:A:208:LEU:O	1:A:210:PRO:HD3	2.21	0.40
1:A:395:HIS:CD2	1:A:397:LYS:H	2.35	0.40
1:A:133:THR:HG21	1:A:168:ILE:HD13	2.03	0.40
1:A:452:ILE:HD12	1:A:452:ILE:HA	1.95	0.40
1:A:157:LEU:CA	1:A:228:ASP:HB2	2.48	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	521/644 (81%)	472 (91%)	46 (9%)	3 (1%)	21	51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	164	GLY
1	A	212	PHE
1	A	369	SER

5.3.2 Protein sidechains [i](#)

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	487/582 (84%)	401 (82%)	86 (18%)	2	6

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LYS
1	A	34	ILE
1	A	36	ASP
1	A	38	THR
1	A	43	LEU
1	A	62	VAL
1	A	64	ASP
1	A	67	LEU

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Mol	Chain	Res	Type
1	A	72	LYS
1	A	81	THR
1	A	87	ILE
1	A	91	ASN
1	A	113	LEU
1	A	117	ARG
1	A	123	LYS
1	A	124	GLN
1	A	129	PHE
1	A	133	THR
1	A	137	LYS
1	A	138	LEU
1	A	147	ILE
1	A	151	THR
1	A	156	LEU
1	A	157	LEU
1	A	158	THR
1	A	169	SER
1	A	171	ILE
1	A	182	LYS
1	A	191	LEU
1	A	195	ILE
1	A	213	ASN
1	A	220	VAL
1	A	229	VAL
1	A	234	VAL
1	A	261	GLU
1	A	263	TRP
1	A	264	VAL
1	A	266	ASN
1	A	277	ASN
1	A	289	ARG
1	A	291	ILE
1	A	332	GLU
1	A	342	ILE
1	A	347	LEU
1	A	349	SER
1	A	354	LYS
1	A	365	LEU
1	A	371	LEU
1	A	375	LYS
1	A	382	VAL

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Mol	Chain	Res	Type
1	A	396	SER
1	A	405	THR
1	A	417	ILE
1	A	424	ASN
1	A	435	LEU
1	A	440	LEU
1	A	442	ASN
1	A	447	LYS
1	A	451	LYS
1	A	452	ILE
1	A	462	GLU
1	A	463	ASN
1	A	474	LEU
1	A	485	LYS
1	A	487	LEU
1	A	490	LEU
1	A	502	ILE
1	A	503	ARG
1	A	509	MET
1	A	512	THR
1	A	518	ARG
1	A	522	LEU
1	A	523	LEU
1	A	524	LYS
1	A	538	LEU
1	A	548	SER
1	A	586	ASP
1	A	588	VAL
1	A	604	ARG
1	A	608	VAL
1	A	610	GLU
1	A	612	VAL
1	A	623	ARG
1	A	628	MET
1	A	632	VAL
1	A	634	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	91	ASN

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Mol	Chain	Res	Type
1	A	112	ASN
1	A	148	ASN
1	A	211	GLN
1	A	213	ASN
1	A	214	GLN
1	A	242	GLN
1	A	249	GLN
1	A	266	ASN
1	A	277	ASN
1	A	299	HIS
1	A	340	GLN
1	A	356	GLN
1	A	358	GLN
1	A	366	GLN
1	A	385	GLN
1	A	390	GLN
1	A	395	HIS
1	A	424	ASN
1	A	437	ASN
1	A	442	ASN
1	A	478	HIS
1	A	499	GLN
1	A	597	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	533/644 (82%)	0.23	10 (1%) 66 57	45, 63, 100, 124	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	463	ASN	5.3
1	A	530	ILE	3.3
1	A	464	ASN	3.1
1	A	111	PRO	2.9
1	A	26	ILE	2.4
1	A	454	SER	2.4
1	A	34	ILE	2.3
1	A	553	GLY	2.1
1	A	276	SER	2.0
1	A	460	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FE2	A	1501	1/1	0.95	0.15	67,67,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.