



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:18 PM UTC

PDB ID : 9HN8 / pdb_00009hn8
Title : Apo Structure of Truncated 1-deoxy-D-xylulose 5-phosphate synthase (DXPS)
from Mycobacterium tuberculosis
Authors : Gawriljuk, V.O.; Groves, M.R.
Deposited on : 2024-12-10
Resolution : 2.65 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

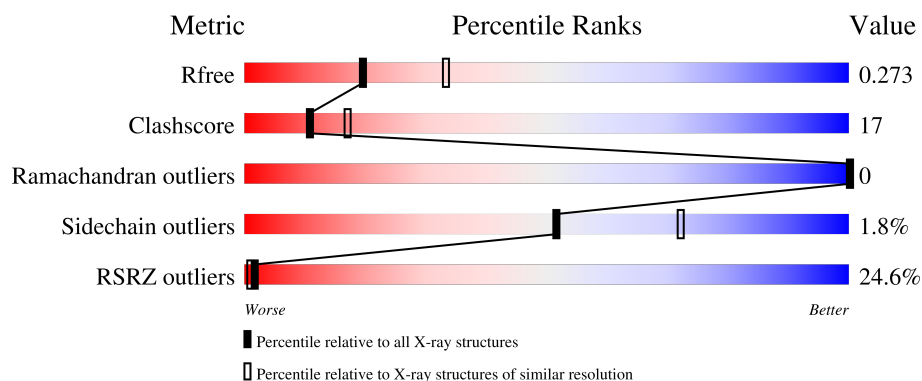
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	604	19% 56% 24% • 20%
1	B	604	20% 55% 22% 22%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7205 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 1-deoxy-D-xylulose-5-phosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	484	Total	C	N	O	S	0	0	0
			3634	2291	664	665	14			
1	B	470	Total	C	N	O	S	0	2	0
			3551	2242	652	643	14			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P9WNS3
A	-2	ALA	-	expression tag	UNP P9WNS3
A	-1	MET	-	expression tag	UNP P9WNS3
A	0	GLY	-	expression tag	UNP P9WNS3
A	228	GLY	-	linker	UNP P9WNS3
A	229	GLY	-	linker	UNP P9WNS3
A	230	GLY	-	linker	UNP P9WNS3
A	231	GLY	-	linker	UNP P9WNS3
A	232	GLY	-	linker	UNP P9WNS3
A	233	GLY	-	linker	UNP P9WNS3
A	234	GLY	-	linker	UNP P9WNS3
A	314	VAL	PRO	conflict	UNP P9WNS3
A	593	VAL	TYR	conflict	UNP P9WNS3
B	-3	GLY	-	expression tag	UNP P9WNS3
B	-2	ALA	-	expression tag	UNP P9WNS3
B	-1	MET	-	expression tag	UNP P9WNS3
B	0	GLY	-	expression tag	UNP P9WNS3
B	228	GLY	-	linker	UNP P9WNS3
B	229	GLY	-	linker	UNP P9WNS3
B	230	GLY	-	linker	UNP P9WNS3
B	231	GLY	-	linker	UNP P9WNS3
B	232	GLY	-	linker	UNP P9WNS3
B	233	GLY	-	linker	UNP P9WNS3
B	234	GLY	-	linker	UNP P9WNS3
B	314	VAL	PRO	conflict	UNP P9WNS3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	593	VAL	TYR	conflict	UNP P9WNS3

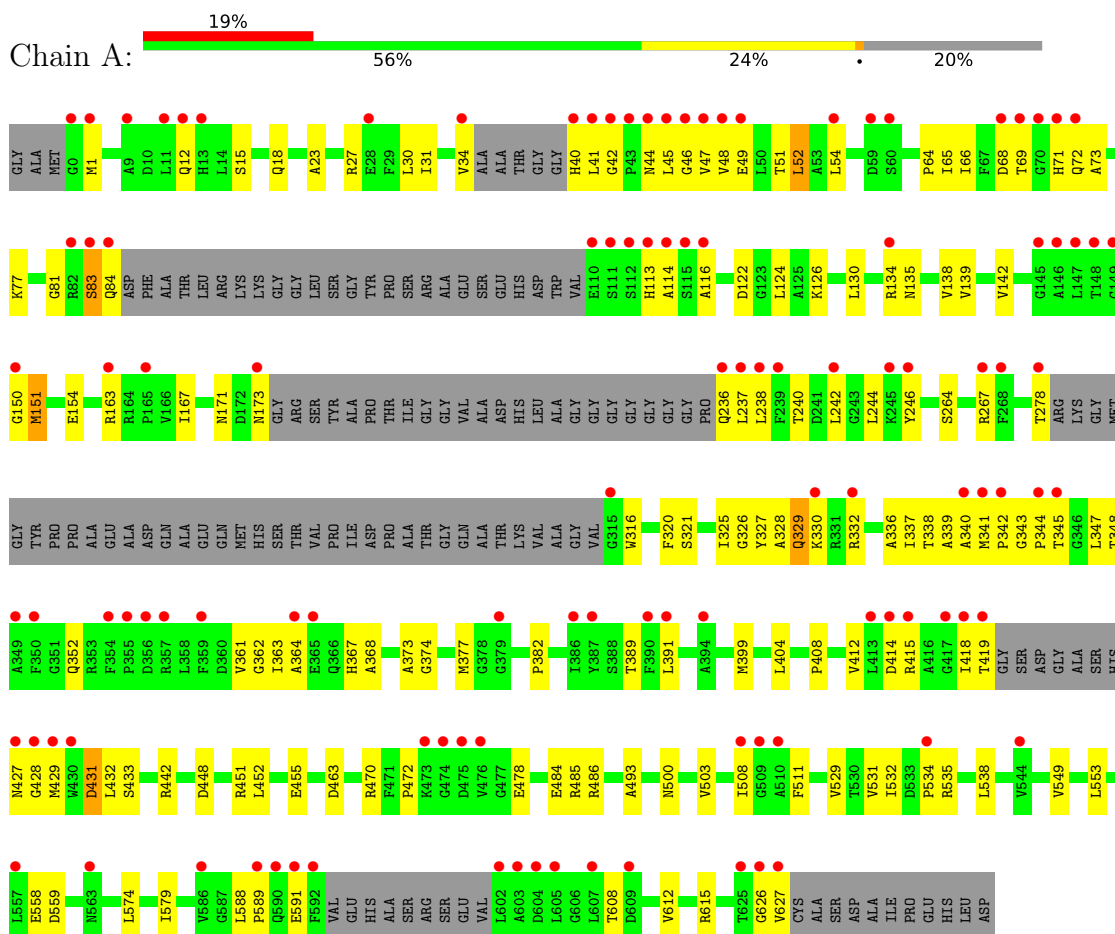
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total 6	O 6	0	0
2	B	14	Total 14	O 14	0	0

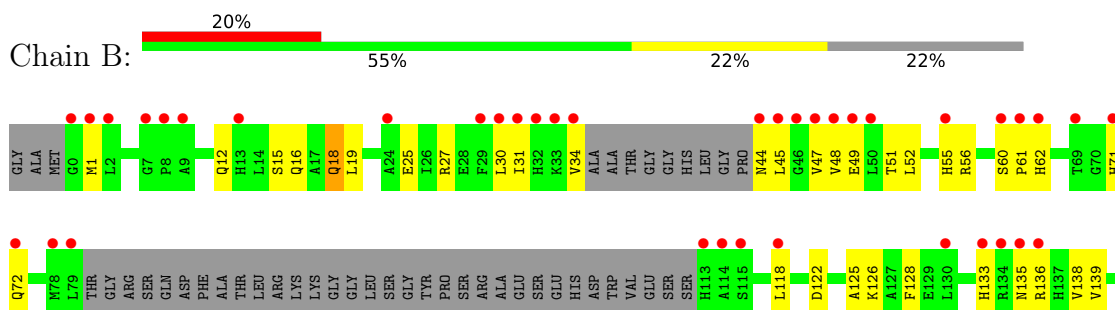
3 Residue-property plots [i](#)

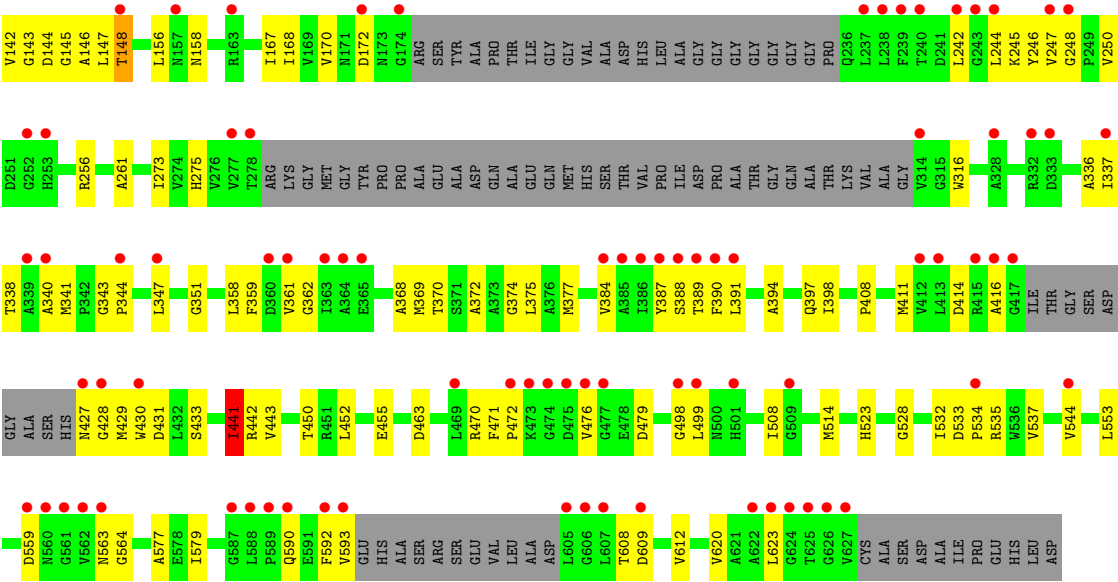
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 1-deoxy-D-xylulose-5-phosphate synthase



- Molecule 1: 1-deoxy-D-xylulose-5-phosphate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.18Å 126.77Å 80.55Å 90.00° 107.70° 90.00°	Depositor
Resolution (Å)	48.87 – 2.65 48.87 – 2.65	Depositor EDS
% Data completeness (in resolution range)	97.4 (48.87-2.65) 97.6 (48.87-2.65)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.247 , 0.273 0.247 , 0.273	Depositor DCC
R_{free} test set	1789 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	59.5	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7205	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.90% 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	0.74	3/3697 (0.1%)	0.97	11/5020 (0.2%)
1	B	0.64	0/3612	0.90	8/4903 (0.2%)
All	All	0.69	3/7309 (0.0%)	0.94	19/9923 (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	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	534	PRO	CA-C	-7.56	1.45	1.52
1	A	433	SER	C-O	5.77	1.30	1.24
1	A	42	GLY	C-O	5.61	1.31	1.24

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	327	TYR	CA-C-O	-9.40	110.58	120.55
1	B	146	ALA	N-CA-C	-8.89	102.61	113.19
1	B	534	PRO	N-CA-C	-8.62	96.97	110.50
1	A	534	PRO	N-CA-C	-8.15	96.84	110.21
1	A	382	PRO	N-CA-CB	-7.11	96.72	103.33
1	A	12	GLN	N-CA-C	7.08	119.86	111.71
1	A	329	GLN	N-CA-C	-6.59	104.34	112.38
1	B	533	ASP	CB-CA-C	-6.43	103.77	110.33
1	A	327	TYR	O-C-N	5.91	128.38	122.12
1	A	431	ASP	CA-CB-CG	5.90	118.50	112.60
1	A	163	ARG	N-CA-C	-5.88	104.25	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	148	THR	N-CA-C	-5.41	106.53	113.23
1	A	52	LEU	N-CA-C	-5.38	104.83	111.33
1	B	471	PHE	O-C-N	-5.36	118.28	121.71
1	B	18	GLN	N-CA-C	-5.29	105.69	111.82
1	A	327	TYR	CB-CA-C	5.25	119.50	110.79
1	B	498	GLY	CA-C-N	-5.22	113.14	122.07
1	B	498	GLY	C-N-CA	-5.22	113.14	122.07
1	A	151	MET	N-CA-C	-5.16	105.73	111.36

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	441	ILE	Mainchain
1	B	442[A]	ARG	Mainchain,Sidechain
1	B	442[B]	ARG	Mainchain

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	3634	0	3659	133	0
1	B	3551	0	3586	111	0
2	A	6	0	0	0	0
2	B	14	0	0	1	0
All	All	7205	0	7245	240	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 17.

All (240) 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:264:SER:O	1:A:267:ARG:HG2	1.35	1.19
1:A:47:VAL:CG2	1:A:72:GLN:HG2	1.73	1.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:VAL:HG21	1:A:72:GLN:HG2	1.25	1.15
1:B:341:MET:HE3	1:B:344:PRO:HG3	1.24	1.15
1:B:341:MET:HE3	1:B:344:PRO:CG	1.83	1.08
1:A:47:VAL:HG21	1:A:72:GLN:CG	1.91	1.00
1:A:242:LEU:HD22	1:A:244:LEU:CD1	1.92	0.99
1:A:418:ILE:HD12	1:A:588:LEU:CD2	1.97	0.94
1:A:418:ILE:HD12	1:A:588:LEU:HD22	1.48	0.94
1:B:341:MET:CE	1:B:344:PRO:HG3	1.98	0.93
1:B:133:HIS:HD2	1:B:136:ARG:CD	1.81	0.92
1:A:264:SER:HA	1:A:267:ARG:HE	1.38	0.89
1:A:34:VAL:HG21	1:A:41:LEU:HD21	1.55	0.89
1:A:77:LYS:HE2	1:A:83:SER:HB3	1.55	0.87
1:A:64:PRO:HG2	1:A:138:VAL:HG22	1.58	0.83
1:A:242:LEU:HD22	1:A:244:LEU:HD12	1.59	0.82
1:B:361:VAL:HG21	1:B:368:ALA:HB2	1.60	0.81
1:A:73:ALA:O	1:A:77:LYS:HG3	1.81	0.80
1:B:133:HIS:HD2	1:B:136:ARG:HD2	1.47	0.79
1:A:47:VAL:HG11	1:A:72:GLN:O	1.82	0.79
1:A:418:ILE:CD1	1:A:588:LEU:HD22	2.14	0.77
1:A:30:LEU:HD13	1:A:44:ASN:HB3	1.66	0.77
1:A:242:LEU:HD22	1:A:244:LEU:HD11	1.67	0.76
1:A:374:GLY:HA2	1:A:377:MET:HE2	1.68	0.76
1:B:145:GLY:O	1:B:148:THR:HG22	1.88	0.74
1:B:71:HIS:CD2	1:B:72:GLN:HG3	2.23	0.73
1:B:337:ILE:HD11	1:B:375:LEU:CD1	2.18	0.72
1:A:408:PRO:HB3	1:A:463:ASP:HA	1.71	0.72
1:B:71:HIS:HD2	1:B:72:GLN:HG3	1.54	0.72
1:A:414:ASP:OD1	1:A:415:ARG:N	2.23	0.71
1:A:451:ARG:O	1:A:455:GLU:HG3	1.90	0.71
1:A:64:PRO:CG	1:A:138:VAL:HG22	2.20	0.71
1:A:484:GLU:OE2	1:A:486:ARG:HD2	1.91	0.70
1:B:374:GLY:HA2	1:B:377:MET:HE2	1.73	0.69
1:A:77:LYS:HE2	1:A:83:SER:CB	2.22	0.67
1:A:47:VAL:HG21	1:A:72:GLN:CB	2.24	0.67
1:A:470:ARG:NH2	1:A:559:ASP:O	2.28	0.67
1:A:415:ARG:HH11	1:A:427:ASN:HB2	1.59	0.66
1:A:418:ILE:HG22	1:A:418:ILE:O	1.96	0.66
1:B:514:MET:HE1	1:B:612:VAL:HG21	1.78	0.66
1:B:340:ALA:HB2	1:B:362:GLY:HA2	1.78	0.65
1:A:242:LEU:O	1:A:242:LEU:HD23	1.97	0.65
1:B:133:HIS:CD2	1:B:136:ARG:CD	2.73	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:HIS:CD2	1:B:136:ARG:HD2	2.30	0.64
1:A:77:LYS:CE	1:A:83:SER:HB3	2.27	0.64
1:A:626:GLY:O	1:A:627:VAL:HG23	1.96	0.64
1:A:415:ARG:O	1:A:415:ARG:HG3	1.97	0.64
1:A:236:GLN:OE1	1:A:236:GLN:N	2.31	0.64
1:B:133:HIS:HD2	1:B:136:ARG:HD3	1.60	0.64
1:A:51:THR:HA	1:A:54:LEU:HG	1.80	0.63
1:A:316:TRP:NE1	1:A:472:PRO:HG2	2.14	0.63
1:A:418:ILE:O	1:A:419:THR:C	2.42	0.63
1:A:418:ILE:HD12	1:A:588:LEU:HD21	1.81	0.62
1:A:589:PRO:HB2	1:A:591:GLU:HG2	1.82	0.62
1:A:151:MET:HE2	1:B:397:GLN:HE21	1.65	0.62
1:A:626:GLY:O	1:A:627:VAL:CG2	2.48	0.61
1:B:384:VAL:HG22	1:B:411:MET:HA	1.82	0.61
1:A:342:PRO:HA	1:A:345:THR:HG22	1.82	0.60
1:B:433:SER:HA	1:B:564:GLY:O	2.00	0.60
1:A:389:THR:HB	1:A:428:GLY:H	1.66	0.59
1:A:116:ALA:HB1	1:B:370:THR:HG21	1.84	0.59
1:A:264:SER:HA	1:A:267:ARG:NE	2.12	0.59
1:A:415:ARG:NH1	1:A:427:ASN:HB2	2.18	0.59
1:B:145:GLY:C	1:B:147:LEU:H	2.11	0.58
1:A:478:GLU:N	1:A:478:GLU:OE1	2.35	0.57
1:B:71:HIS:HD2	1:B:72:GLN:H	1.51	0.57
1:A:338:THR:HG21	1:A:341:MET:C	2.29	0.57
1:B:168:ILE:HB	1:B:273:ILE:HD13	1.86	0.57
1:B:133:HIS:CD2	1:B:136:ARG:HD3	2.37	0.57
1:B:30:LEU:O	1:B:34:VAL:HG23	2.05	0.56
1:B:44:ASN:OD1	1:B:71:HIS:HB2	2.05	0.56
1:A:341:MET:HB2	1:A:344:PRO:HG2	1.87	0.56
1:B:145:GLY:C	1:B:147:LEU:N	2.59	0.55
1:A:242:LEU:CD2	1:A:244:LEU:CD1	2.78	0.55
1:B:391:LEU:HD11	1:B:398:ILE:CD1	2.37	0.55
1:B:343:GLY:N	1:B:344:PRO:HD2	2.22	0.55
1:A:574:LEU:HB2	1:A:579:ILE:HB	1.88	0.55
1:A:264:SER:O	1:A:267:ARG:CG	2.30	0.55
1:A:242:LEU:CD2	1:A:244:LEU:HD11	2.36	0.55
1:B:391:LEU:HD11	1:B:398:ILE:HD13	1.89	0.55
1:A:500:ASN:HB3	1:A:627:VAL:CG2	2.37	0.54
1:B:27:ARG:O	1:B:31:ILE:HG12	2.08	0.54
1:A:139:VAL:HG22	1:A:167:ILE:HB	1.90	0.54
1:A:626:GLY:C	1:A:627:VAL:HG23	2.32	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:TYR:CE1	1:B:248:GLY:HA2	2.43	0.54
1:B:499:LEU:HD11	1:B:523:HIS:ND1	2.23	0.53
1:B:351:GLY:HA3	1:B:358:LEU:HD22	1.90	0.53
1:B:470:ARG:HH22	1:B:559:ASP:HB3	1.74	0.53
1:A:46:GLY:HA3	1:A:278:THR:HG21	1.89	0.53
1:A:171:ASN:OD1	1:A:278:THR:OG1	2.23	0.53
1:B:388:SER:HB2	1:B:428:GLY:HA2	1.90	0.53
1:B:338:THR:HG22	1:B:347:LEU:HD12	1.90	0.53
1:A:361:VAL:HG21	1:A:368:ALA:HB2	1.91	0.53
1:B:336:ALA:HB1	1:B:347:LEU:HD13	1.89	0.53
1:A:69:THR:C	1:A:71:HIS:H	2.17	0.53
1:A:532:ILE:O	1:A:532:ILE:HG13	2.08	0.53
1:B:343:GLY:H	1:B:344:PRO:HD2	1.73	0.53
1:B:391:LEU:CD1	1:B:398:ILE:CD1	2.86	0.52
1:B:441:ILE:HG23	1:B:443:VAL:HG23	1.91	0.52
1:B:499:LEU:HB3	1:B:528:GLY:CA	2.39	0.52
1:B:563:ASN:HB2	1:B:590:GLN:CD	2.34	0.52
1:A:336:ALA:C	1:A:337:ILE:HD12	2.34	0.52
1:B:620:VAL:HA	1:B:623:LEU:HG	1.90	0.52
1:B:247:VAL:HG11	1:B:261:ALA:HB1	1.91	0.52
1:B:416:ALA:CB	1:B:470:ARG:HG2	2.39	0.52
1:A:237:LEU:HD23	1:A:237:LEU:H	1.75	0.52
1:A:238:LEU:HD12	1:A:238:LEU:H	1.74	0.52
1:A:316:TRP:CE3	1:A:452:LEU:HD12	2.45	0.51
1:B:44:ASN:O	1:B:47:VAL:HG22	2.11	0.51
1:A:500:ASN:HB3	1:A:627:VAL:HG22	1.92	0.51
1:A:608:THR:O	1:A:612:VAL:HG23	2.11	0.51
1:B:62:HIS:CD2	1:B:136:ARG:HG2	2.45	0.51
1:A:389:THR:HB	1:A:427:ASN:HB3	1.91	0.51
1:A:493:ALA:HB3	1:A:531:VAL:HB	1.93	0.51
1:B:30:LEU:HD23	1:B:45:LEU:HG	1.92	0.51
1:A:122:ASP:O	1:A:126:LYS:HG3	2.11	0.51
1:A:130:LEU:HD11	1:B:359:PHE:HZ	1.76	0.51
1:B:414:ASP:HB2	2:B:705:HOH:O	2.11	0.51
1:B:341:MET:HE3	1:B:344:PRO:CB	2.40	0.50
1:A:431:ASP:CG	1:A:470:ARG:NH1	2.68	0.50
1:A:503:VAL:HG23	1:A:529:VAL:HG13	1.93	0.50
1:A:328:ALA:C	1:A:330:LYS:N	2.67	0.50
1:B:337:ILE:HD11	1:B:375:LEU:HD11	1.93	0.50
1:B:15:SER:H	1:B:18:GLN:HB2	1.76	0.50
1:A:418:ILE:O	1:A:418:ILE:CG2	2.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:VAL:HG12	1:A:553:LEU:HD23	1.94	0.50
1:A:41:LEU:HD22	1:A:41:LEU:H	1.78	0.49
1:B:316:TRP:CZ2	1:B:476:VAL:HG23	2.47	0.49
1:A:419:THR:HG21	1:A:472:PRO:HB2	1.95	0.49
1:A:332:ARG:HG3	1:A:332:ARG:HH11	1.78	0.49
1:A:442:ARG:HA	1:A:538:LEU:O	2.13	0.49
1:B:245:LYS:NZ	1:B:246:TYR:H	2.11	0.49
1:B:553:LEU:HD13	1:B:623:LEU:HD21	1.94	0.49
1:A:236:GLN:HB3	1:A:246:TYR:CZ	2.48	0.49
1:B:48:VAL:O	1:B:52:LEU:HG	2.13	0.48
1:B:416:ALA:HB1	1:B:470:ARG:HG2	1.94	0.48
1:A:66:ILE:HD11	1:A:124:LEU:HD12	1.95	0.48
1:A:508:ILE:HB	1:A:558:GLU:HA	1.95	0.48
1:B:429:MET:HG3	1:B:430:TRP:CE2	2.49	0.48
1:A:404:LEU:HD12	1:B:592:PHE:CG	2.49	0.48
1:A:431:ASP:OD2	1:A:470:ARG:NH1	2.47	0.48
1:A:364:ALA:HB1	1:A:367:HIS:HB3	1.95	0.48
1:B:369:MET:CE	1:B:384:VAL:HG21	2.44	0.48
1:B:118:LEU:CD2	1:B:142:VAL:HG21	2.43	0.47
1:A:27:ARG:O	1:A:31:ILE:HG12	2.14	0.47
1:A:150:GLY:O	1:A:154:GLU:HG3	2.14	0.47
1:A:338:THR:HG22	1:A:341:MET:H	1.80	0.47
1:A:373:ALA:O	1:A:377:MET:HG3	2.14	0.47
1:B:391:LEU:CD1	1:B:398:ILE:HD11	2.45	0.47
1:B:52:LEU:O	1:B:56:ARG:HG3	2.14	0.47
1:A:65:ILE:HD12	1:A:139:VAL:HB	1.97	0.47
1:A:432:LEU:HG	1:A:470:ARG:NH1	2.30	0.47
1:A:485:ARG:HH21	1:A:535:ARG:CD	2.29	0.46
1:B:391:LEU:HD12	1:B:398:ILE:HD11	1.96	0.46
1:B:139:VAL:HG22	1:B:167:ILE:HB	1.97	0.46
1:B:144:ASP:HB3	1:B:172:ASP:HA	1.97	0.46
1:B:156:LEU:HB3	1:B:242:LEU:HD21	1.96	0.46
1:A:414:ASP:CG	1:A:415:ARG:H	2.21	0.46
1:B:372:ALA:HA	1:B:375:LEU:HD12	1.98	0.46
1:A:448:ASP:CG	1:A:451:ARG:HH21	2.24	0.46
1:B:47:VAL:HG11	1:B:72:GLN:HG2	1.98	0.46
1:A:27:ARG:HH22	1:A:49:GLU:CD	2.24	0.46
1:B:125:ALA:HB2	1:B:138:VAL:HG11	1.98	0.46
1:A:429:MET:HG2	1:A:588:LEU:HD13	1.98	0.45
1:B:31:ILE:O	1:B:34:VAL:HB	2.15	0.45
1:A:47:VAL:HG12	1:A:51:THR:OG1	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:ASN:HA	1:A:278:THR:O	2.17	0.45
1:A:23:ALA:O	1:A:27:ARG:HG3	2.15	0.45
1:B:429:MET:HE3	1:B:429:MET:HB2	1.75	0.45
1:B:347:LEU:HD23	1:B:347:LEU:HA	1.79	0.45
1:B:122:ASP:O	1:B:126:LYS:HG3	2.16	0.45
1:B:142:VAL:CG1	1:B:143:GLY:N	2.79	0.45
1:A:236:GLN:O	1:A:240:THR:OG1	2.29	0.44
1:A:328:ALA:C	1:A:330:LYS:H	2.25	0.44
1:A:431:ASP:OD1	1:A:470:ARG:NH1	2.50	0.44
1:A:52:LEU:HD23	1:A:52:LEU:HA	1.80	0.44
1:B:388:SER:CB	1:B:428:GLY:HA2	2.46	0.44
1:B:499:LEU:HB3	1:B:528:GLY:HA3	1.99	0.44
1:B:12:GLN:HA	1:B:56:ARG:HH11	1.81	0.44
1:B:608:THR:O	1:B:609:ASP:C	2.60	0.44
1:A:338:THR:HG21	1:A:342:PRO:N	2.32	0.44
1:B:71:HIS:CD2	1:B:72:GLN:H	2.32	0.44
1:A:338:THR:OG1	1:A:347:LEU:HD12	2.18	0.44
1:B:455:GLU:OE2	1:B:535:ARG:HB3	2.18	0.44
1:B:532:ILE:CD1	1:B:544:VAL:HG22	2.48	0.44
1:A:77:LYS:O	1:A:81:GLY:N	2.50	0.43
1:A:83:SER:HB2	1:A:84:GLN:OE1	2.17	0.43
1:B:316:TRP:NE1	1:B:472:PRO:HG2	2.33	0.43
1:A:320:PHE:CZ	1:A:412:VAL:HG21	2.54	0.43
1:B:337:ILE:HG12	1:B:359:PHE:HB2	2.00	0.43
1:B:508:ILE:HG23	1:B:537:VAL:HG21	2.01	0.43
1:B:27:ARG:NH2	1:B:49:GLU:OE2	2.52	0.43
1:B:147:LEU:N	1:B:147:LEU:HD22	2.33	0.43
1:A:113:HIS:HD2	1:A:114:ALA:N	2.16	0.43
1:A:326:GLY:O	1:A:329:GLN:HG2	2.18	0.43
1:B:142:VAL:HG12	1:B:143:GLY:N	2.33	0.43
1:B:248:GLY:O	1:B:250:VAL:HG23	2.19	0.43
1:B:341:MET:HB3	1:B:344:PRO:HG2	2.01	0.43
1:A:68:ASP:HB2	1:A:142:VAL:HG12	2.01	0.42
1:B:1:MET:H	1:B:25:GLU:CD	2.25	0.42
1:B:51:THR:O	1:B:55:HIS:HD2	2.02	0.42
1:B:128:PHE:HB3	1:B:133:HIS:O	2.18	0.42
1:A:343:GLY:N	1:A:344:PRO:HD2	2.34	0.42
1:A:418:ILE:CD1	1:A:588:LEU:HD13	2.50	0.42
1:A:429:MET:HG2	1:A:588:LEU:CD1	2.50	0.42
1:B:408:PRO:HB3	1:B:463:ASP:HA	2.00	0.42
1:A:549:VAL:CG2	1:A:579:ILE:HG12	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:TRP:HE3	1:A:452:LEU:HD12	1.85	0.42
1:A:414:ASP:CG	1:A:415:ARG:N	2.76	0.42
1:A:264:SER:C	1:A:267:ARG:HG2	2.31	0.42
1:A:41:LEU:HA	1:A:44:ASN:HB2	2.01	0.42
1:A:419:THR:HG21	1:A:472:PRO:CB	2.50	0.42
1:B:71:HIS:HD2	1:B:72:GLN:N	2.16	0.42
1:B:369:MET:HG3	1:B:394:ALA:HB1	2.02	0.42
1:B:316:TRP:CE3	1:B:452:LEU:HD12	2.55	0.41
1:B:338:THR:CG2	1:B:347:LEU:HD12	2.49	0.41
1:A:340:ALA:HB2	1:A:362:GLY:HA2	2.01	0.41
1:A:321:SER:O	1:A:325:ILE:HG12	2.20	0.41
1:A:418:ILE:HD11	1:A:588:LEU:HD13	2.02	0.41
1:B:19:LEU:HA	1:B:19:LEU:HD12	1.82	0.41
1:B:389:THR:CG2	1:B:427:ASN:HA	2.50	0.41
1:A:47:VAL:CG1	1:A:51:THR:OG1	2.69	0.41
1:A:485:ARG:HH21	1:A:535:ARG:HD3	1.86	0.41
1:B:170:VAL:HG13	1:B:275:HIS:HA	2.03	0.41
1:B:156:LEU:HD22	1:B:244:LEU:CD1	2.51	0.41
1:B:387:TYR:HB2	1:B:390:PHE:CD2	2.56	0.41
1:B:62:HIS:NE2	1:B:136:ARG:HG2	2.35	0.41
1:A:363:ILE:H	1:A:363:ILE:HD12	1.85	0.40
1:A:399:MET:HE3	1:A:399:MET:HB3	1.89	0.40
1:B:49:GLU:H	1:B:49:GLU:CD	2.28	0.40
1:B:450:THR:OG1	1:B:479:ASP:OD2	2.34	0.40
1:A:44:ASN:O	1:A:48:VAL:HG23	2.20	0.40
1:A:343:GLY:HA2	1:A:348:THR:HG23	2.03	0.40
1:B:60:SER:N	1:B:61:PRO:HD2	2.36	0.40
1:A:338:THR:HG22	1:A:339:ALA:N	2.35	0.40
1:A:615:ARG:HH11	1:A:615:ARG:HG2	1.86	0.40
1:B:16:GLN:HG2	1:B:256[A]:ARG:CZ	2.52	0.40
1:B:577:ALA:HB3	1:B:579:ILE:HD12	2.04	0.40
1:A:1:MET:HE3	1:A:1:MET:HB2	1.78	0.40
1:A:15:SER:HB3	1:A:18:GLN:HG3	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	470/604 (78%)	462 (98%)	8 (2%)	0	100	100
1	B	458/604 (76%)	447 (98%)	11 (2%)	0	100	100
All	All	928/1208 (77%)	909 (98%)	19 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	373/453 (82%)	365 (98%)	8 (2%)	47	69
1	B	363/453 (80%)	358 (99%)	5 (1%)	59	76
All	All	736/906 (81%)	723 (98%)	13 (2%)	51	72

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	HIS
1	A	45	LEU
1	A	83	SER
1	A	134	ARG
1	A	135	ASN
1	A	352	GLN
1	A	391	LEU
1	A	511	PHE

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Mol	Chain	Res	Type
1	B	135	ASN
1	B	158	ASN
1	B	431	ASP
1	B	441	ILE
1	B	593	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	HIS
1	A	72	GLN
1	A	236	GLN
1	A	329	GLN
1	B	62	HIS
1	B	71	HIS
1	B	72	GLN
1	B	133	HIS
1	B	405	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	484/604 (80%)	1.36	114 (23%)	2 1	40, 66, 114, 177	0
1	B	470/604 (77%)	1.38	121 (25%)	1 1	25, 67, 119, 152	2 (0%)
All	All	954/1208 (78%)	1.37	235 (24%)	2 1	25, 66, 118, 177	2 (0%)

All (235) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	115	SER	7.6
1	A	509	GLY	7.1
1	B	113	HIS	6.9
1	B	174	GLY	6.7
1	A	428	GLY	6.6
1	B	625	THR	6.6
1	B	8	PRO	6.6
1	A	114	ALA	5.9
1	B	61	PRO	5.7
1	A	113	HIS	5.5
1	A	591	GLU	5.4
1	A	592	PHE	5.4
1	A	627	VAL	5.4
1	B	133	HIS	5.2
1	A	0	GLY	5.1
1	A	145	GLY	5.1
1	A	341	MET	5.0
1	A	429	MET	4.9
1	A	45	LEU	4.8
1	A	242	LEU	4.7
1	B	417	GLY	4.6
1	A	43	PRO	4.5
1	B	363	ILE	4.4
1	B	242	LEU	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	589	PRO	4.4
1	B	624	GLY	4.3
1	A	34	VAL	4.3
1	B	593	VAL	4.3
1	B	607	LEU	4.2
1	B	337	ILE	4.2
1	B	47	VAL	4.2
1	B	30	LEU	4.2
1	A	268	PHE	4.1
1	B	623	LEU	4.1
1	A	418	ILE	4.1
1	B	389	THR	4.1
1	B	314	VAL	4.1
1	A	473	LYS	4.0
1	A	345	THR	4.0
1	B	386	ILE	4.0
1	B	50	LEU	4.0
1	B	562	VAL	4.0
1	B	365	GLU	3.9
1	B	605	LEU	3.9
1	A	48	VAL	3.9
1	B	588	LEU	3.9
1	A	41	LEU	3.8
1	B	498	GLY	3.8
1	B	29	PHE	3.8
1	A	47	VAL	3.8
1	A	69	THR	3.8
1	B	135	ASN	3.8
1	A	238	LEU	3.8
1	B	45	LEU	3.7
1	B	589	PRO	3.7
1	B	60	SER	3.7
1	B	114	ALA	3.7
1	B	592	PHE	3.7
1	B	44	ASN	3.7
1	A	84	GLN	3.7
1	B	388	SER	3.6
1	B	248	GLY	3.6
1	B	627	VAL	3.5
1	A	112	SER	3.5
1	A	427	ASN	3.5
1	A	46	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	54	LEU	3.5
1	A	356	ASP	3.5
1	A	607	LEU	3.5
1	A	419	THR	3.4
1	A	625	THR	3.4
1	B	332	ARG	3.4
1	B	134	ARG	3.4
1	B	559	ASP	3.4
1	B	278	THR	3.3
1	B	244	LEU	3.3
1	B	238	LEU	3.3
1	B	413	LEU	3.3
1	A	237	LEU	3.3
1	A	110	GLU	3.3
1	B	428	GLY	3.3
1	A	510	ALA	3.3
1	B	247	VAL	3.3
1	B	136	ARG	3.2
1	B	115	SER	3.2
1	A	146	ALA	3.2
1	B	416	ALA	3.2
1	B	590	GLN	3.2
1	B	34	VAL	3.2
1	B	49	GLU	3.1
1	A	590	GLN	3.1
1	B	72	GLN	3.1
1	B	427	ASN	3.1
1	B	430	TRP	3.1
1	A	344	PRO	3.1
1	B	509	GLY	3.1
1	B	415	ARG	3.1
1	B	390	PHE	3.1
1	B	347	LEU	3.0
1	A	44	ASN	3.0
1	B	253	HIS	3.0
1	B	148	THR	3.0
1	A	150	GLY	3.0
1	B	1	MET	3.0
1	A	602	LEU	3.0
1	A	349	ALA	3.0
1	B	364	ALA	3.0
1	A	355	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	605	LEU	2.9
1	B	474	GLY	2.9
1	B	391	LEU	2.9
1	A	148	THR	2.9
1	A	49	GLU	2.9
1	A	350	PHE	2.9
1	A	390	PHE	2.9
1	A	415	ARG	2.8
1	B	163	ARG	2.8
1	B	239	PHE	2.8
1	A	394	ALA	2.8
1	A	163	ARG	2.8
1	B	534	PRO	2.8
1	B	606	GLY	2.8
1	A	68	ASP	2.8
1	A	387	TYR	2.8
1	B	622	ALA	2.8
1	A	40	HIS	2.8
1	A	342	PRO	2.7
1	A	534	PRO	2.7
1	A	13	HIS	2.7
1	A	70	GLY	2.7
1	A	71	HIS	2.7
1	B	475	ASP	2.7
1	B	387	TYR	2.7
1	A	359	PHE	2.7
1	B	361	VAL	2.7
1	B	31	ILE	2.7
1	A	315	GLY	2.7
1	B	544	VAL	2.7
1	A	330	LYS	2.6
1	B	243	GLY	2.6
1	A	476	VAL	2.6
1	B	328	ALA	2.6
1	B	333	ASP	2.6
1	A	626	GLY	2.6
1	A	508	ILE	2.5
1	A	603	ALA	2.5
1	B	9	ALA	2.5
1	B	157	ASN	2.5
1	A	386	ILE	2.5
1	A	134	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	147	LEU	2.5
1	B	252	GLY	2.5
1	B	561	GLY	2.5
1	A	1	MET	2.5
1	B	69	THR	2.5
1	B	33	LYS	2.5
1	A	72	GLN	2.5
1	B	412	VAL	2.5
1	A	82	ARG	2.4
1	A	391	LEU	2.4
1	A	83	SER	2.4
1	A	42	GLY	2.4
1	A	165	PRO	2.4
1	A	173	ASN	2.4
1	A	60	SER	2.4
1	A	417	GLY	2.4
1	A	544	VAL	2.4
1	A	116	ALA	2.4
1	A	239	PHE	2.4
1	B	13	HIS	2.3
1	A	475	ASP	2.3
1	A	609	ASP	2.3
1	B	172	ASP	2.3
1	B	360	ASP	2.3
1	A	12	GLN	2.3
1	A	340	ALA	2.3
1	B	469	LEU	2.3
1	B	32	HIS	2.3
1	A	430	TRP	2.3
1	B	344	PRO	2.3
1	A	149	GLY	2.3
1	B	587	GLY	2.3
1	A	364	ALA	2.3
1	A	557	LEU	2.3
1	B	340	ALA	2.3
1	A	278	THR	2.3
1	B	240	THR	2.3
1	A	28	GLU	2.3
1	B	563	ASN	2.3
1	A	111	SER	2.3
1	A	604	ASP	2.3
1	B	384	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	2.3
1	B	560	ASN	2.3
1	B	472	PRO	2.3
1	B	609	ASP	2.3
1	B	0	GLY	2.3
1	B	477	GLY	2.3
1	B	48	VAL	2.3
1	B	130	LEU	2.3
1	B	55	HIS	2.2
1	A	414	ASP	2.2
1	A	474	GLY	2.2
1	B	118	LEU	2.2
1	B	78	MET	2.2
1	A	563	ASN	2.2
1	A	379	GLY	2.2
1	B	237	LEU	2.2
1	B	626	GLY	2.2
1	B	24	ALA	2.2
1	A	365	GLU	2.2
1	B	46	GLY	2.2
1	B	79	LEU	2.2
1	B	62	HIS	2.2
1	B	277	VAL	2.2
1	A	357	ARG	2.1
1	A	245	LYS	2.1
1	A	246	TYR	2.1
1	B	2	LEU	2.1
1	B	385	ALA	2.1
1	A	586	VAL	2.1
1	A	354	PHE	2.1
1	A	9	ALA	2.1
1	B	339	ALA	2.1
1	B	476	VAL	2.1
1	A	267	ARG	2.0
1	A	332	ARG	2.0
1	B	473	LYS	2.0
1	A	236	GLN	2.0
1	B	499	LEU	2.0
1	B	71	HIS	2.0
1	A	59	ASP	2.0
1	A	413	LEU	2.0
1	B	501	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	7	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.