



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:26 AM UTC

PDB ID : 4G85 / pdb_00004g85
Title : Crystal structure of human HisRS
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Deposited on : 2012-07-21
Resolution : 3.11 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

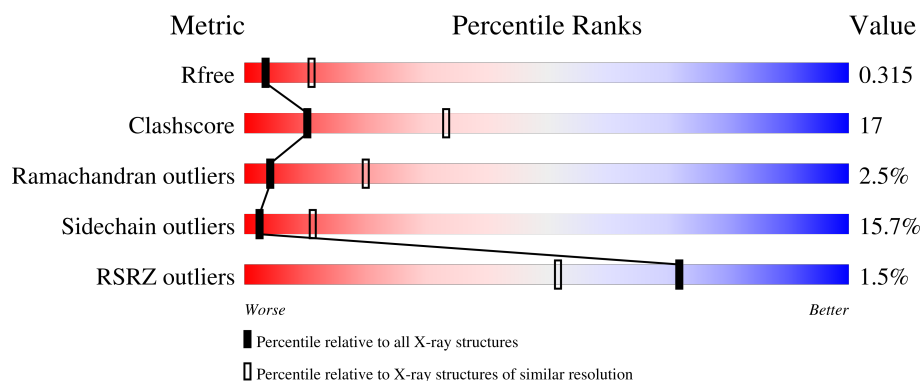
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.11 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	1816 (3.14-3.10)
Clashscore	190562	1906 (3.14-3.10)
Ramachandran outliers	187476	1802 (3.14-3.10)
Sidechain outliers	187428	1802 (3.14-3.10)
RSRZ outliers	180081	1816 (3.14-3.10)

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	517	2% 53% 22% 7% 17%
1	B	517	2% 51% 24% 6% 18%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6432 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 Histidine-tRNA ligase, cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	0	0
			3209	2051	536	608	14			
1	B	426	Total	C	N	O	S	0	0	0
			3223	2055	545	609	14			

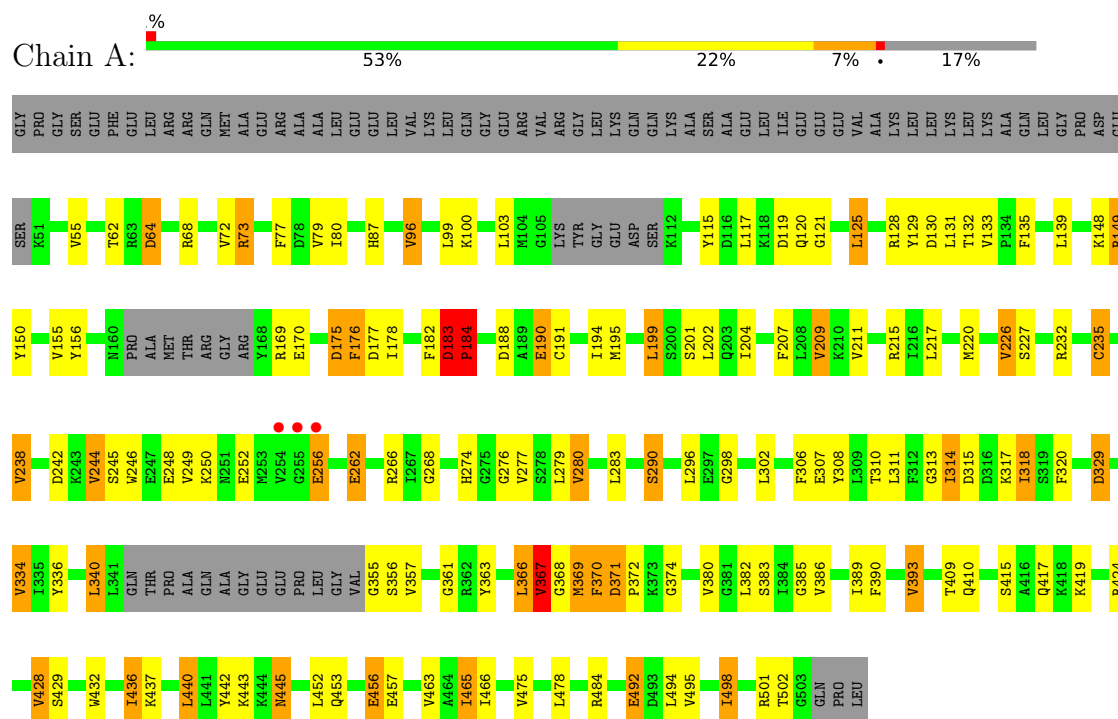
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	GLY	-	expression tag	UNP P12081
A	-9	PRO	-	expression tag	UNP P12081
A	-8	GLY	-	expression tag	UNP P12081
A	-7	SER	-	expression tag	UNP P12081
A	-6	GLU	-	expression tag	UNP P12081
A	-5	PHE	-	expression tag	UNP P12081
A	-4	GLU	-	expression tag	UNP P12081
A	-3	LEU	-	expression tag	UNP P12081
A	-2	ARG	-	expression tag	UNP P12081
A	-1	ARG	-	expression tag	UNP P12081
A	0	GLN	-	expression tag	UNP P12081
B	-10	GLY	-	expression tag	UNP P12081
B	-9	PRO	-	expression tag	UNP P12081
B	-8	GLY	-	expression tag	UNP P12081
B	-7	SER	-	expression tag	UNP P12081
B	-6	GLU	-	expression tag	UNP P12081
B	-5	PHE	-	expression tag	UNP P12081
B	-4	GLU	-	expression tag	UNP P12081
B	-3	LEU	-	expression tag	UNP P12081
B	-2	ARG	-	expression tag	UNP P12081
B	-1	ARG	-	expression tag	UNP P12081
B	0	GLN	-	expression tag	UNP P12081

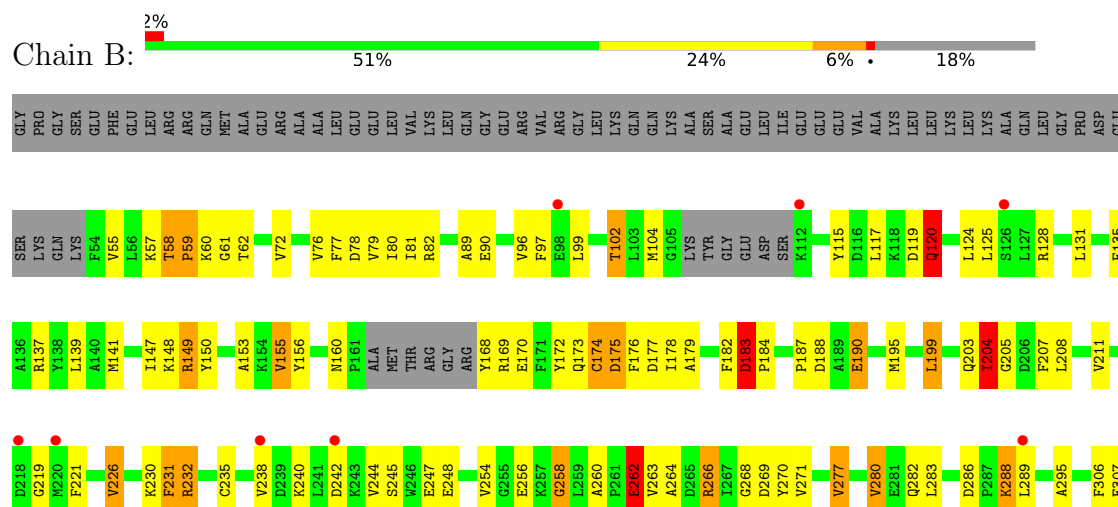
3 Residue-property plots

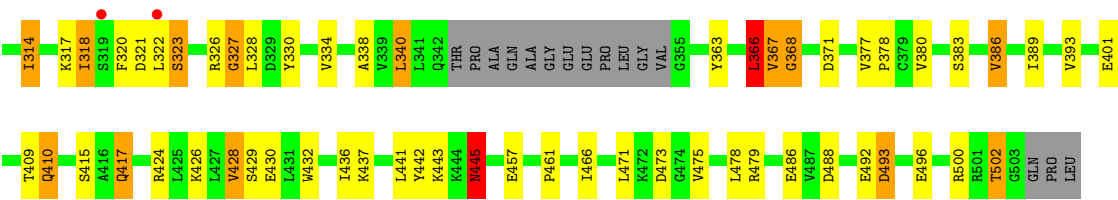
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Histidine-tRNA ligase, cytoplasmic



• Molecule 1: Histidine-tRNA ligase, cytoplasmic





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	100.45Å 100.45Å 257.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.11 50.00 – 3.11	Depositor EDS
% Data completeness (in resolution range)	97.0 (50.00-3.11) 97.2 (50.00-3.11)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.271 , 0.327 0.259 , 0.315	Depositor DCC
R_{free} test set	1213 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	94.9	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6432	wwPDB-VP
Average B, all atoms (Å ²)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.19	3/3255 (0.1%)	1.16	10/4412 (0.2%)
1	B	1.52	13/3269 (0.4%)	1.46	17/4427 (0.4%)
All	All	1.37	16/6524 (0.2%)	1.32	27/8839 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	367	VAL	CA-C	63.45	2.30	1.52
1	A	367	VAL	N-CA	35.92	1.89	1.46
1	A	367	VAL	CA-CB	34.80	1.99	1.54
1	B	367	VAL	N-CA	20.42	1.70	1.46
1	A	367	VAL	CA-C	-17.55	1.30	1.52
1	B	367	VAL	CA-CB	15.17	1.72	1.54
1	B	288	LYS	CD-CE	9.82	1.81	1.52
1	B	268	GLY	C-O	8.85	1.34	1.23
1	B	264	ALA	CA-CB	6.63	1.63	1.53
1	B	262	GLU	CD-OE2	6.55	1.37	1.25
1	B	262	GLU	CD-OE1	-6.29	1.13	1.25
1	B	258	GLY	C-N	6.01	1.41	1.33
1	B	266	ARG	NE-CZ	5.79	1.39	1.33
1	B	248	GLU	C-N	5.32	1.40	1.33
1	B	269	ASP	CG-OD2	5.12	1.35	1.25
1	B	266	ARG	CZ-NH2	5.04	1.40	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	367	VAL	CA-C-O	37.09	160.48	121.17
1	B	367	VAL	CA-CB-CG2	25.88	154.39	110.40
1	B	366	LEU	CA-C-N	-25.80	88.05	120.56
1	B	366	LEU	C-N-CA	-25.80	88.05	120.56
1	B	367	VAL	N-CA-C	21.96	131.51	110.42
1	B	367	VAL	CA-CB-CG1	-18.31	79.27	110.40
1	A	367	VAL	CA-CB-CG1	-16.05	83.12	110.40
1	A	367	VAL	N-CA-C	15.31	126.18	110.72
1	A	366	LEU	CA-C-N	-14.82	99.38	120.42
1	A	366	LEU	C-N-CA	-14.82	99.38	120.42
1	B	367	VAL	CB-CA-C	14.47	130.50	111.97
1	B	367	VAL	N-CA-CB	-11.25	97.39	110.55
1	A	367	VAL	N-CA-CB	-9.93	95.69	110.58
1	B	445	ASN	CA-C-N	-7.31	113.28	120.52
1	B	445	ASN	C-N-CA	-7.31	113.28	120.52
1	A	183	ASP	CA-C-N	7.11	128.72	119.84
1	A	183	ASP	C-N-CA	7.11	128.72	119.84
1	A	334	VAL	CB-CA-C	-6.94	100.88	111.08
1	B	58	THR	CA-C-N	5.71	126.98	119.84
1	B	58	THR	C-N-CA	5.71	126.98	119.84
1	B	183	ASP	CA-C-N	5.42	126.62	119.84
1	B	183	ASP	C-N-CA	5.42	126.62	119.84
1	A	463	VAL	N-CA-C	5.39	115.75	107.77
1	B	266	ARG	NE-CZ-NH2	-5.23	114.49	119.20
1	B	262	GLU	CG-CD-OE2	-5.12	106.62	118.40
1	B	314	ILE	CB-CA-C	5.10	116.13	110.62
1	A	465	ILE	N-CA-C	5.01	115.31	107.99

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	270	TYR	Sidechain

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	3209	0	3112	111	0
1	B	3223	0	3147	116	0
All	All	6432	0	6259	216	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 (216) 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:288:LYS:CE	1:B:288:LYS:CD	1.81	1.55
1:B:367:VAL:N	1:B:367:VAL:CA	1.70	1.52
1:A:367:VAL:CA	1:A:367:VAL:CB	1.99	1.39
1:A:367:VAL:CA	1:A:367:VAL:N	1.89	1.35
1:B:367:VAL:HA	1:B:367:VAL:C	1.48	1.35
1:B:366:LEU:C	1:B:367:VAL:CA	2.12	1.20
1:B:367:VAL:CA	1:B:368:GLY:N	2.12	1.12
1:A:367:VAL:CA	1:A:367:VAL:CG1	2.34	1.04
1:B:367:VAL:CA	1:B:367:VAL:C	2.30	1.04
1:B:366:LEU:O	1:B:367:VAL:CA	2.08	1.00
1:B:367:VAL:CA	1:B:368:GLY:H	1.71	0.96
1:A:183:ASP:OD1	1:B:443:LYS:HA	1.72	0.90
1:B:367:VAL:HA	1:B:368:GLY:N	1.77	0.90
1:B:244:VAL:HG12	1:B:245:SER:H	1.37	0.88
1:A:366:LEU:C	1:A:367:VAL:CA	2.48	0.85
1:A:195:MET:HE1	1:A:209:VAL:HG13	1.63	0.81
1:B:366:LEU:O	1:B:367:VAL:HA	1.78	0.81
1:B:155:VAL:HG21	1:B:173:GLN:HB2	1.62	0.80
1:B:204:ILE:HG12	1:B:340:LEU:HD11	1.62	0.80
1:A:195:MET:HE1	1:A:209:VAL:CG1	2.13	0.78
1:A:370:PHE:N	1:A:370:PHE:CD2	2.51	0.77
1:A:77:PHE:HE1	1:A:386:VAL:HG21	1.48	0.76
1:A:148:LYS:HG3	1:A:178:ILE:HG12	1.68	0.75
1:A:443:LYS:HA	1:B:183:ASP:OD1	1.85	0.75
1:A:370:PHE:N	1:A:370:PHE:HD2	1.84	0.73
1:A:445:ASN:N	1:A:445:ASN:HD22	1.87	0.73
1:B:288:LYS:CE	1:B:288:LYS:CG	2.67	0.73
1:A:195:MET:HG3	1:A:382:LEU:CD2	2.18	0.73
1:B:77:PHE:HE1	1:B:386:VAL:HG21	1.54	0.72
1:A:156:TYR:CE1	1:A:170:GLU:HG3	2.25	0.72
1:B:244:VAL:HG12	1:B:245:SER:N	2.06	0.70
1:B:72:VAL:O	1:B:76:VAL:HG23	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:LEU:HD12	1:A:367:VAL:HG11	1.75	0.69
1:B:149:ARG:CD	1:B:149:ARG:C	2.66	0.68
1:B:177:ASP:OD2	1:B:363:TYR:OH	2.10	0.68
1:B:367:VAL:N	1:B:367:VAL:CB	2.57	0.67
1:B:204:ILE:O	1:B:340:LEU:HD21	1.95	0.67
1:A:389:ILE:O	1:A:393:VAL:HG23	1.94	0.66
1:B:306:PHE:CZ	1:B:318:ILE:HD11	2.32	0.65
1:B:426:LYS:O	1:B:430:GLU:HG3	1.96	0.65
1:A:340:LEU:HD12	1:A:357:VAL:HG21	1.78	0.65
1:B:262:GLU:HB2	1:B:266:ARG:NH1	2.13	0.64
1:B:149:ARG:C	1:B:149:ARG:HD3	2.23	0.64
1:A:183:ASP:CB	1:A:184:PRO:HD2	2.28	0.64
1:A:492:GLU:CD	1:A:492:GLU:H	2.05	0.64
1:A:366:LEU:O	1:A:367:VAL:CA	2.46	0.63
1:B:283:LEU:HB3	1:B:289:LEU:HD12	1.79	0.63
1:A:175:ASP:HB3	1:A:383:SER:HB3	1.80	0.63
1:A:77:PHE:CE1	1:A:386:VAL:HG21	2.33	0.63
1:B:321:ASP:C	1:B:323:SER:H	2.08	0.62
1:B:417:GLN:HE21	1:B:417:GLN:H	1.48	0.61
1:A:453:GLN:HA	1:A:456:GLU:HG3	1.83	0.61
1:A:204:ILE:HD12	1:A:340:LEU:HD11	1.82	0.61
1:B:288:LYS:CD	1:B:288:LYS:NZ	2.62	0.61
1:B:492:GLU:H	1:B:492:GLU:CD	2.09	0.60
1:B:415:SER:HB2	1:B:466:ILE:O	2.02	0.60
1:B:62:THR:HG22	1:B:170:GLU:HB3	1.85	0.59
1:A:313:GLY:O	1:A:314:ILE:HD13	2.03	0.59
1:B:139:LEU:HD21	1:B:147:ILE:HG21	1.84	0.59
1:B:479:ARG:HG3	1:B:486:GLU:HG3	1.85	0.58
1:B:139:LEU:HD21	1:B:147:ILE:CG2	2.33	0.58
1:A:235:CYS:HA	1:A:238:VAL:CG2	2.33	0.58
1:B:155:VAL:CG2	1:B:173:GLN:HB2	2.33	0.58
1:A:176:PHE:C	1:A:176:PHE:CD2	2.82	0.58
1:B:232:ARG:HA	1:B:235:CYS:SG	2.43	0.57
1:B:175:ASP:HB3	1:B:383:SER:HB3	1.86	0.57
1:B:221:PHE:HB3	1:B:226:VAL:CG2	2.34	0.57
1:B:492:GLU:OE2	1:B:492:GLU:N	2.30	0.57
1:A:195:MET:HG3	1:A:382:LEU:HD23	1.85	0.57
1:A:453:GLN:HA	1:A:456:GLU:CG	2.35	0.57
1:B:471:LEU:C	1:B:473:ASP:H	2.12	0.57
1:A:367:VAL:CA	1:A:367:VAL:HG12	2.33	0.56
1:A:117:LEU:HA	1:B:117:LEU:HA	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:TYR:CE1	1:B:170:GLU:HG3	2.41	0.56
1:A:367:VAL:CA	1:A:367:VAL:HG13	2.33	0.56
1:A:199:LEU:HB3	1:A:207:PHE:CE2	2.40	0.56
1:B:424:ARG:O	1:B:428:VAL:HG22	2.05	0.55
1:A:129:TYR:CD1	1:A:130:ASP:HB2	2.41	0.55
1:B:280:VAL:HG12	1:B:320:PHE:CD2	2.42	0.55
1:A:368:GLY:O	1:A:370:PHE:N	2.40	0.55
1:B:219:GLY:HA3	1:B:295:ALA:HB2	1.89	0.55
1:B:410:GLN:HG3	1:B:437:LYS:HB2	1.88	0.55
1:B:174:CYS:O	1:B:174:CYS:SG	2.64	0.55
1:A:445:ASN:N	1:A:445:ASN:ND2	2.55	0.54
1:A:177:ASP:OD2	1:A:363:TYR:OH	2.22	0.54
1:A:367:VAL:HG12	1:A:367:VAL:C	2.32	0.54
1:B:321:ASP:OD1	1:B:323:SER:HB3	2.08	0.54
1:A:235:CYS:HA	1:A:238:VAL:HG22	1.90	0.54
1:A:410:GLN:HG3	1:A:437:LYS:HB2	1.90	0.54
1:B:203:GLN:O	1:B:205:GLY:N	2.33	0.54
1:B:271:VAL:HG13	1:B:323:SER:HB2	1.90	0.53
1:A:183:ASP:HB3	1:A:184:PRO:HD2	1.90	0.53
1:A:246:TRP:CZ2	1:A:268:GLY:HA3	2.44	0.53
1:B:445:ASN:N	1:B:445:ASN:HD22	2.06	0.53
1:A:195:MET:HE1	1:A:209:VAL:HG11	1.91	0.53
1:B:283:LEU:HA	1:B:286:ASP:HB2	1.91	0.52
1:B:156:TYR:CZ	1:B:170:GLU:HG3	2.45	0.52
1:B:389:ILE:O	1:B:393:VAL:HG23	2.10	0.52
1:B:475:VAL:CG1	1:B:488:ASP:HB2	2.39	0.52
1:B:147:ILE:HG22	1:B:179:ALA:HB3	1.92	0.52
1:B:172:TYR:HB3	1:B:386:VAL:HG11	1.92	0.52
1:B:128:ARG:NH2	1:B:153:ALA:HB1	2.25	0.52
1:A:424:ARG:O	1:A:428:VAL:HG22	2.11	0.51
1:A:195:MET:HG3	1:A:382:LEU:HD22	1.90	0.51
1:A:367:VAL:CG1	1:A:367:VAL:C	2.82	0.51
1:A:133:VAL:HG13	1:A:370:PHE:HE1	1.76	0.51
1:A:280:VAL:HG12	1:A:320:PHE:CD2	2.46	0.51
1:B:277:VAL:HG22	1:B:306:PHE:CD1	2.46	0.51
1:B:148:LYS:HG3	1:B:178:ILE:HG12	1.93	0.51
1:A:246:TRP:CZ2	1:A:268:GLY:CA	2.94	0.50
1:A:215:ARG:HH11	1:A:298:GLY:HA2	1.76	0.50
1:A:492:GLU:OE2	1:A:492:GLU:N	2.44	0.50
1:A:424:ARG:HB3	1:A:440:LEU:HD11	1.94	0.50
1:B:221:PHE:HB3	1:B:226:VAL:HG22	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:PRO:HA	1:B:190:GLU:HB3	1.94	0.50
1:B:244:VAL:CG1	1:B:245:SER:H	2.16	0.50
1:A:149:ARG:C	1:A:149:ARG:HD3	2.36	0.50
1:A:308:TYR:HA	1:A:311:LEU:HD12	1.93	0.50
1:A:79:VAL:HG13	1:A:201:SER:HB3	1.94	0.50
1:A:371:ASP:OD1	1:A:374:GLY:N	2.44	0.50
1:B:135:PHE:HB2	1:B:149:ARG:NH1	2.27	0.50
1:A:87:HIS:ND1	1:A:190:GLU:OE1	2.39	0.49
1:B:176:PHE:C	1:B:176:PHE:CD2	2.90	0.49
1:A:252:GLU:O	1:A:256:GLU:N	2.37	0.48
1:B:195:MET:HE3	1:B:338:ALA:HB2	1.94	0.48
1:A:442:TYR:OH	1:B:190:GLU:OE2	2.32	0.48
1:A:232:ARG:HH22	1:A:329:ASP:HB3	1.77	0.48
1:B:330:TYR:O	1:B:363:TYR:HB2	2.13	0.48
1:A:79:VAL:CG1	1:A:201:SER:HB3	2.43	0.47
1:A:244:VAL:CG1	1:A:248:GLU:OE2	2.62	0.47
1:A:429:SER:HA	1:A:432:TRP:CE3	2.49	0.47
1:B:254:VAL:HA	1:B:258:GLY:HA2	1.96	0.47
1:A:274:HIS:HA	1:A:279:LEU:HD22	1.95	0.47
1:A:386:VAL:HG22	1:A:390:PHE:CZ	2.50	0.47
1:A:99:LEU:O	1:A:100:LYS:C	2.58	0.47
1:B:115:TYR:OH	1:B:169:ARG:HD3	2.15	0.47
1:A:290:SER:HA	1:A:296:LEU:HD13	1.97	0.46
1:B:445:ASN:N	1:B:445:ASN:ND2	2.63	0.46
1:B:244:VAL:CG1	1:B:245:SER:N	2.77	0.46
1:B:479:ARG:HB2	1:B:486:GLU:HG2	1.98	0.46
1:B:230:LYS:O	1:B:231:PHE:C	2.58	0.46
1:A:492:GLU:CD	1:A:492:GLU:N	2.72	0.46
1:A:452:LEU:O	1:A:456:GLU:HG2	2.17	0.45
1:A:442:TYR:O	1:B:182:PHE:HB3	2.17	0.45
1:A:190:GLU:OE2	1:B:442:TYR:OH	2.34	0.45
1:B:493:ASP:OD1	1:B:493:ASP:N	2.50	0.45
1:A:62:THR:HG21	1:B:97:PHE:HE1	1.81	0.44
1:A:115:TYR:OH	1:A:169:ARG:HD3	2.16	0.44
1:A:182:PHE:CD2	1:B:442:TYR:HB2	2.52	0.44
1:A:306:PHE:CZ	1:A:318:ILE:HD11	2.51	0.44
1:B:221:PHE:HB3	1:B:226:VAL:HG21	1.99	0.44
1:A:415:SER:OG	1:A:424:ARG:NH1	2.51	0.44
1:B:149:ARG:HD3	1:B:150:TYR:CA	2.47	0.44
1:A:244:VAL:HG12	1:A:248:GLU:OE2	2.17	0.44
1:B:178:ILE:HD11	1:B:190:GLU:HG2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LEU:CD1	1:A:389:ILE:HD13	2.47	0.44
1:A:132:THR:O	1:A:135:PHE:HB3	2.18	0.44
1:A:183:ASP:OD1	1:B:442:TYR:O	2.36	0.44
1:A:279:LEU:O	1:A:283:LEU:HG	2.18	0.44
1:B:89:ALA:HA	1:B:148:LYS:O	2.18	0.44
1:B:417:GLN:HE21	1:B:417:GLN:N	2.15	0.44
1:B:283:LEU:O	1:B:286:ASP:HB3	2.18	0.43
1:A:436:ILE:CD1	1:A:495:VAL:HG13	2.48	0.43
1:B:424:ARG:HG2	1:B:466:ILE:HD12	2.00	0.43
1:B:429:SER:HA	1:B:432:TRP:CE3	2.53	0.43
1:B:99:LEU:O	1:B:102:THR:HG23	2.18	0.43
1:A:103:LEU:HB3	1:A:129:TYR:HE2	1.83	0.43
1:A:68:ARG:O	1:A:72:VAL:HG23	2.19	0.43
1:A:436:ILE:HD12	1:A:495:VAL:HG13	2.00	0.43
1:B:326:ARG:O	1:B:328:LEU:N	2.52	0.43
1:B:81:ILE:O	1:B:82:ARG:C	2.62	0.43
1:B:471:LEU:C	1:B:473:ASP:N	2.77	0.43
1:A:442:TYR:HB2	1:B:182:PHE:CD2	2.53	0.42
1:A:494:LEU:O	1:A:498:ILE:HB	2.19	0.42
1:A:361:GLY:C	1:A:380:VAL:HG23	2.44	0.42
1:B:475:VAL:HG13	1:B:488:ASP:HB2	2.01	0.42
1:A:355:GLY:C	1:A:357:VAL:H	2.27	0.42
1:A:424:ARG:HB3	1:A:440:LEU:CD1	2.49	0.42
1:B:119:ASP:C	1:B:120:GLN:HG2	2.44	0.42
1:B:410:GLN:O	1:B:461:PRO:HD2	2.18	0.42
1:A:249:VAL:O	1:A:250:LYS:C	2.62	0.42
1:A:262:GLU:O	1:A:266:ARG:HG3	2.19	0.42
1:A:119:ASP:C	1:A:121:GLY:H	2.28	0.42
1:B:61:GLY:O	1:B:168:TYR:HD1	2.02	0.42
1:B:377:VAL:HA	1:B:378:PRO:HD3	1.83	0.42
1:A:149:ARG:HG2	1:A:150:TYR:N	2.35	0.42
1:A:276:GLY:O	1:A:277:VAL:C	2.62	0.42
1:A:369:MET:C	1:A:370:PHE:CD2	2.98	0.41
1:B:188:ASP:HA	1:B:380:VAL:HG21	2.02	0.41
1:B:211:VAL:O	1:B:320:PHE:HD1	2.03	0.41
1:A:96:VAL:O	1:A:128:ARG:HG2	2.19	0.41
1:A:226:VAL:HG23	1:A:227:SER:O	2.19	0.41
1:B:90:GLU:O	1:B:149:ARG:HA	2.21	0.41
1:A:246:TRP:CH2	1:A:268:GLY:HA2	2.55	0.41
1:B:57:LYS:HE2	1:B:60:LYS:N	2.35	0.41
1:A:217:LEU:HA	1:A:220:MET:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:MET:C	1:A:370:PHE:HD2	2.29	0.41
1:A:190:GLU:O	1:A:194:ILE:HG13	2.21	0.41
1:A:73:ARG:HA	1:A:390:PHE:CE1	2.55	0.41
1:A:424:ARG:HA	1:A:466:ILE:HD12	2.03	0.41
1:A:131:LEU:HD12	1:A:131:LEU:HA	1.85	0.41
1:B:321:ASP:C	1:B:323:SER:N	2.77	0.41
1:B:367:VAL:C	1:B:367:VAL:CG1	2.94	0.41
1:B:199:LEU:HB3	1:B:207:PHE:CE2	2.56	0.40
1:B:496:GLU:O	1:B:500:ARG:HB2	2.21	0.40
1:A:211:VAL:HG22	1:A:336:TYR:HB3	2.03	0.40
1:B:410:GLN:CG	1:B:437:LYS:HB2	2.49	0.40
1:A:64:ASP:OD1	1:B:137:ARG:NH2	2.55	0.40
1:A:117:LEU:HD13	1:A:125:LEU:HB2	2.03	0.40
1:B:57:LYS:HG2	1:B:58:THR:H	1.87	0.40
1:A:244:VAL:HG12	1:A:245:SER:H	1.86	0.40
1:A:302:LEU:HD23	1:A:302:LEU:HA	1.95	0.40
1:B:78:ASP:O	1:B:79:VAL:C	2.64	0.40
1:B:59:PRO:O	1:B:60:LYS:C	2.64	0.40
1:B:260:ALA:HB3	1:B:263:VAL:HG23	2.02	0.40
1:B:326:ARG:O	1:B:327:GLY:C	2.64	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	419/517 (81%)	382 (91%)	28 (7%)	9 (2%)	5	23
1	B	418/517 (81%)	372 (89%)	34 (8%)	12 (3%)	3	18
All	All	837/1034 (81%)	754 (90%)	62 (7%)	21 (2%)	4	20

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	184	PRO
1	B	59	PRO
1	B	204	ILE
1	B	502	THR
1	A	120	GLN
1	A	369	MET
1	A	419	LYS
1	B	120	GLN
1	B	184	PRO
1	B	327	GLY
1	A	256	GLU
1	A	315	ASP
1	B	104	MET
1	B	368	GLY
1	A	502	THR
1	B	322	LEU
1	A	372	PRO
1	A	385	GLY
1	B	231	PHE
1	B	256	GLU
1	B	401	GLU

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	326/444 (73%)	273 (84%)	53 (16%)	2	10
1	B	330/444 (74%)	280 (85%)	50 (15%)	3	12
All	All	656/888 (74%)	553 (84%)	103 (16%)	2	11

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

Mol	Chain	Res	Type
1	A	55	VAL
1	A	64	ASP
1	A	73	ARG

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Mol	Chain	Res	Type
1	A	80	ILE
1	A	96	VAL
1	A	125	LEU
1	A	149	ARG
1	A	155	VAL
1	A	175	ASP
1	A	176	PHE
1	A	183	ASP
1	A	184	PRO
1	A	188	ASP
1	A	190	GLU
1	A	191	CYS
1	A	199	LEU
1	A	209	VAL
1	A	226	VAL
1	A	235	CYS
1	A	238	VAL
1	A	242	ASP
1	A	244	VAL
1	A	262	GLU
1	A	280	VAL
1	A	290	SER
1	A	307	GLU
1	A	310	THR
1	A	314	ILE
1	A	317	LYS
1	A	318	ILE
1	A	329	ASP
1	A	334	VAL
1	A	340	LEU
1	A	356	SER
1	A	367	VAL
1	A	370	PHE
1	A	371	ASP
1	A	393	VAL
1	A	409	THR
1	A	417	GLN
1	A	428	VAL
1	A	436	ILE
1	A	440	LEU
1	A	445	ASN
1	A	456	GLU

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Mol	Chain	Res	Type
1	A	457	GLU
1	A	465	ILE
1	A	475	VAL
1	A	478	LEU
1	A	484	ARG
1	A	492	GLU
1	A	498	ILE
1	A	501	ARG
1	B	55	VAL
1	B	80	ILE
1	B	96	VAL
1	B	102	THR
1	B	120	GLN
1	B	124	LEU
1	B	125	LEU
1	B	131	LEU
1	B	141	MET
1	B	149	ARG
1	B	155	VAL
1	B	160	ASN
1	B	174	CYS
1	B	175	ASP
1	B	183	ASP
1	B	190	GLU
1	B	199	LEU
1	B	204	ILE
1	B	208	LEU
1	B	226	VAL
1	B	232	ARG
1	B	238	VAL
1	B	240	LYS
1	B	242	ASP
1	B	247	GLU
1	B	262	GLU
1	B	277	VAL
1	B	280	VAL
1	B	282	GLN
1	B	307	GLU
1	B	314	ILE
1	B	317	LYS
1	B	318	ILE
1	B	323	SER

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Mol	Chain	Res	Type
1	B	334	VAL
1	B	340	LEU
1	B	366	LEU
1	B	371	ASP
1	B	386	VAL
1	B	409	THR
1	B	410	GLN
1	B	417	GLN
1	B	428	VAL
1	B	436	ILE
1	B	441	LEU
1	B	445	ASN
1	B	457	GLU
1	B	478	LEU
1	B	493	ASP
1	B	502	THR

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	160	ASN
1	A	272	GLN
1	A	445	ASN
1	B	120	GLN
1	B	160	ASN
1	B	272	GLN
1	B	417	GLN
1	B	445	ASN
1	B	451	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [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 [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	427/517 (82%)	-0.01	3 (0%) 84 68	74, 92, 122, 157	0
1	B	426/517 (82%)	0.21	10 (2%) 61 39	76, 112, 199, 246	0
All	All	853/1034 (82%)	0.10	13 (1%) 72 52	74, 98, 177, 246	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	242	ASP	4.0
1	B	319	SER	3.6
1	B	220	MET	3.1
1	A	255	GLY	3.0
1	B	218	ASP	2.6
1	B	238	VAL	2.5
1	A	256	GLU	2.5
1	B	98	GLU	2.5
1	A	254	VAL	2.4
1	B	112	LYS	2.2
1	B	322	LEU	2.2
1	B	126	SER	2.2
1	B	289	LEU	2.1

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.