



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 10:40 AM EDT

PDB ID : 5M98 / pdb_00005m98
Title : Crystal structure of urate oxidase from zebrafish
Authors : Zanotti, G.; Cendron, I.; Percudani, R.; Berni, R.
Deposited on : 2016-11-01
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

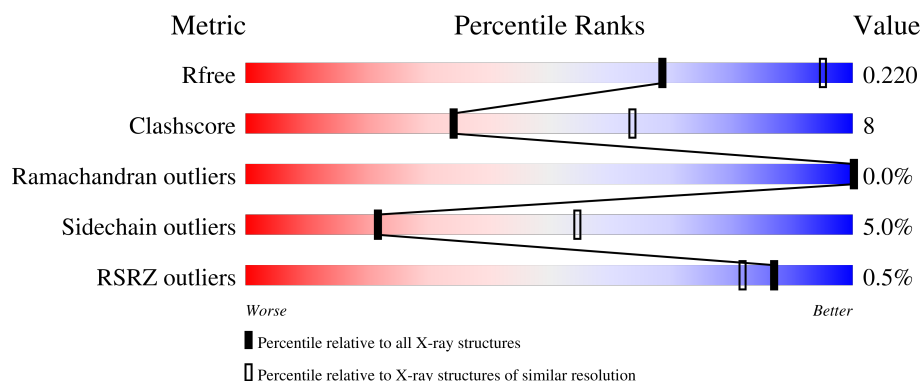
1 Overall quality at a glance

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	298	% 76% 19% • •
1	B	298	76% 18% • •
1	C	298	% 80% 15% • •
1	D	298	79% 16% • •
1	E	298	76% 19% • •

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Mol	Chain	Length	Quality of chain
1	F	298	% 73%21%••
1	G	298	% 76%18%••
1	H	298	% 78%15%••

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 18677 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 Uricase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	B	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	C	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	D	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	E	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	F	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	G	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			
1	H	287	Total	C	N	O	S	0	0	0
			2322	1482	405	426	9			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	13	Total	O	0	0
			13	13		
2	B	15	Total	O	0	0
			15	15		
2	C	23	Total	O	0	0
			23	23		
2	D	15	Total	O	0	0
			15	15		
2	E	5	Total	O	0	0
			5	5		
2	F	9	Total	O	0	0
			9	9		

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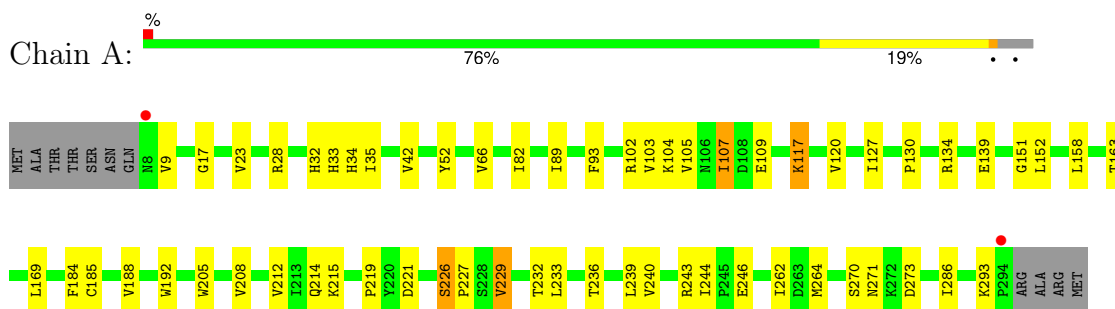
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	7	Total	O	0	0
			7	7		
2	H	14	Total	O	0	0
			14	14		

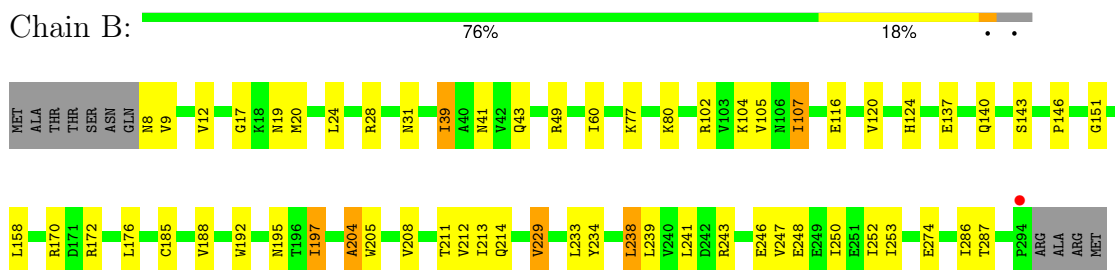
3 Residue-property plots [i](#)

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.

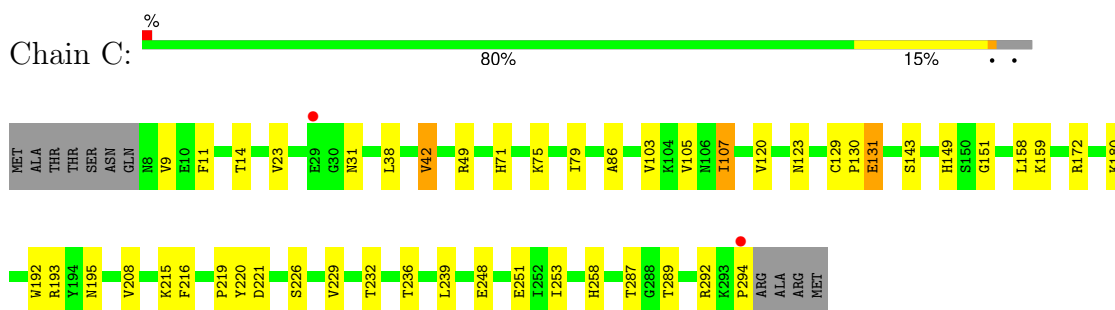
• Molecule 1: Uricase



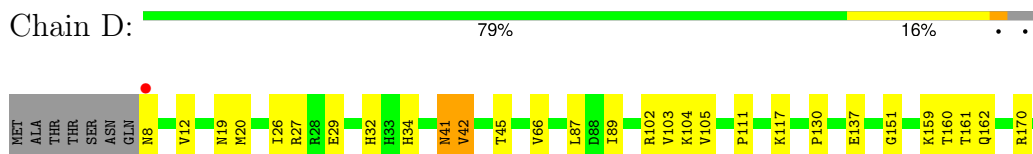
• Molecule 1: Uricase

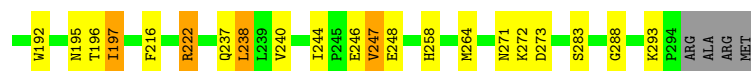


• Molecule 1: Uricase



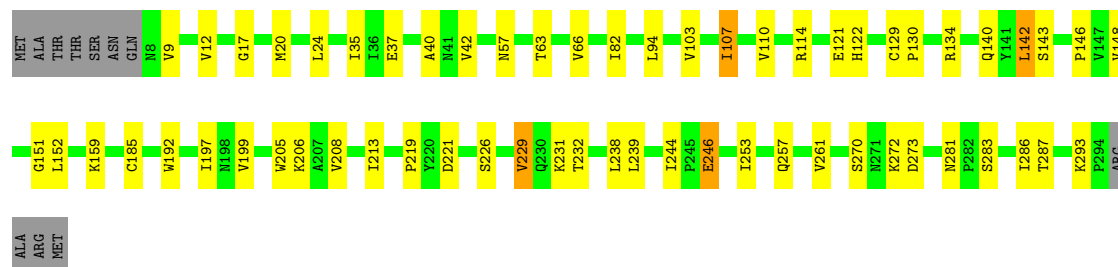
• Molecule 1: Uricase





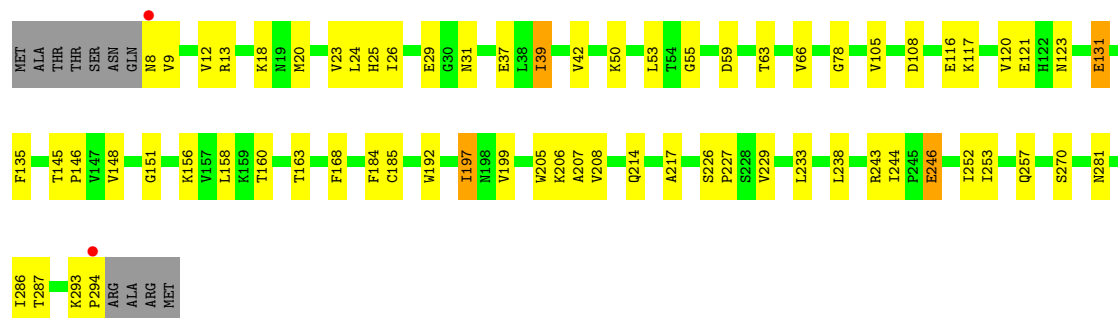
• Molecule 1: Uricase

Chain E: 76% 19%



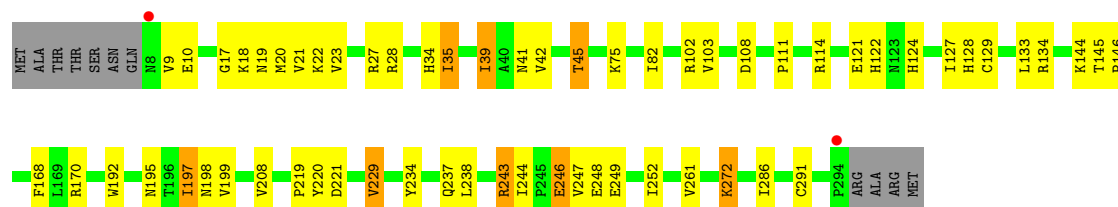
• Molecule 1: Uricase

Chain F: 73% 21%



• Molecule 1: Uricase

Chain G: 76% 18%



• Molecule 1: Uricase

Chain H: 78% 15%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.34Å 127.42Å 132.60Å 90.00° 102.21° 90.00°	Depositor
Resolution (Å)	83.20 – 2.80 83.20 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.9 (83.20-2.80) 94.9 (83.20-2.80)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.82Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.204 , 0.242 0.206 , 0.220	Depositor DCC
R_{free} test set	3240 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 27.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	18677	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 59.32 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7963e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.91	0/2374	1.12	3/3217 (0.1%)
1	B	0.92	0/2374	1.09	4/3217 (0.1%)
1	C	0.95	0/2374	1.11	3/3217 (0.1%)
1	D	0.94	0/2374	1.11	1/3217 (0.0%)
1	E	0.90	0/2374	1.10	1/3217 (0.0%)
1	F	0.86	0/2374	1.04	2/3217 (0.1%)
1	G	0.96	0/2374	1.06	0/3217
1	H	0.93	0/2374	1.10	3/3217 (0.1%)
All	All	0.92	0/18992	1.09	17/25736 (0.1%)

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	E	0	1

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	273	ASP	N-CA-C	6.72	120.68	111.54
1	F	9	VAL	N-CA-C	6.03	116.18	108.82
1	B	9	VAL	N-CA-C	5.99	117.54	108.80
1	A	226	SER	CA-C-N	5.97	127.30	119.84
1	A	226	SER	C-N-CA	5.97	127.30	119.84
1	H	250	ILE	N-CA-C	5.78	116.39	107.78
1	C	258	HIS	N-CA-C	5.76	119.30	110.20
1	B	17	GLY	N-CA-C	5.55	119.81	111.42
1	B	60	ILE	N-CA-C	5.45	116.04	108.84
1	C	226	SER	CA-C-N	5.39	126.24	120.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	226	SER	C-N-CA	5.39	126.24	120.47
1	E	148	VAL	N-CA-C	5.36	115.61	108.11
1	H	129	CYS	N-CA-CB	5.27	116.67	110.71
1	B	204	ALA	N-CA-C	-5.22	105.48	111.07
1	F	131	GLU	N-CA-C	-5.22	105.48	111.07
1	H	39	ILE	CB-CA-C	5.13	117.35	110.42
1	D	258	HIS	N-CA-C	5.02	117.74	109.76

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	142	LEU	Peptide

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	2322	0	2325	40	0
1	B	2322	0	2325	35	0
1	C	2322	0	2325	35	0
1	D	2322	0	2325	30	0
1	E	2322	0	2325	45	0
1	F	2322	0	2325	53	0
1	G	2322	0	2325	45	0
1	H	2322	0	2325	37	0
2	A	13	0	0	0	0
2	B	15	0	0	0	0
2	C	23	0	0	1	0
2	D	15	0	0	0	0
2	E	5	0	0	0	0
2	F	9	0	0	0	0
2	G	7	0	0	1	0
2	H	14	0	0	0	0
All	All	18677	0	18600	282	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 8.

All (282) 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:MET:HE3	1:A:271:ASN:HB2	1.52	0.89
1:G:9:VAL:HG21	1:H:238:LEU:HD11	1.55	0.88
1:B:229:VAL:HG21	1:B:286:ILE:HD13	1.56	0.87
1:F:229:VAL:HG21	1:F:286:ILE:HD13	1.56	0.87
1:G:20:MET:HA	1:G:39:ILE:HD11	1.61	0.82
1:D:264:MET:HE3	1:D:271:ASN:HB2	1.60	0.82
1:H:241:LEU:HD23	1:H:247:VAL:HG23	1.63	0.81
1:A:151:GLY:O	1:A:205:TRP:HZ3	1.64	0.80
1:A:215:LYS:HB3	1:A:236:THR:HG22	1.62	0.79
1:H:215:LYS:HB3	1:H:236:THR:HG22	1.64	0.79
1:D:170:ARG:HH22	1:D:177:THR:HA	1.48	0.78
1:D:20:MET:H	1:D:41:ASN:HD22	1.30	0.76
1:C:38:LEU:HD22	1:C:107:ILE:HD11	1.65	0.75
1:F:158:LEU:HD13	1:F:185:CYS:HB2	1.70	0.72
1:B:28:ARG:NH2	1:D:273:ASP:O	2.22	0.72
1:A:158:LEU:HD13	1:A:185:CYS:HB2	1.72	0.72
1:B:151:GLY:O	1:B:205:TRP:HZ3	1.71	0.72
1:B:19:ASN:HD22	1:B:41:ASN:CG	1.99	0.71
1:E:151:GLY:O	1:E:205:TRP:HZ3	1.72	0.71
1:F:253:ILE:HG12	1:F:287:THR:HG22	1.73	0.70
1:D:20:MET:H	1:D:41:ASN:ND2	1.89	0.70
1:E:20:MET:HB3	1:F:281:ASN:HD22	1.56	0.70
1:F:151:GLY:O	1:F:205:TRP:HZ3	1.73	0.70
1:A:117:LYS:HE3	1:C:220:TYR:CZ	2.27	0.69
1:E:229:VAL:CG2	1:E:286:ILE:HG13	2.25	0.67
1:C:42:VAL:HB	1:C:105:VAL:HG12	1.77	0.67
1:A:42:VAL:HG12	1:A:105:VAL:HG22	1.77	0.66
1:G:238:LEU:HD11	1:H:9:VAL:HG21	1.76	0.66
1:D:244:ILE:CG2	1:D:247:VAL:HG13	2.25	0.66
1:A:214:GLN:NE2	1:H:120:VAL:HG22	2.11	0.66
1:G:27:ARG:NH2	1:G:34:HIS:CD2	2.65	0.65
1:H:156:LYS:NZ	1:H:257:GLN:OE1	2.30	0.65
1:C:129:CYS:O	1:C:129:CYS:SG	2.55	0.64
1:G:243:ARG:HG3	1:G:243:ARG:HH11	1.63	0.64
1:E:20:MET:HE3	1:F:281:ASN:ND2	2.12	0.64
1:B:49:ARG:HG3	1:B:49:ARG:HH11	1.62	0.64
1:F:42:VAL:HG11	1:F:66:VAL:HG22	1.80	0.63
1:C:49:ARG:HG2	1:C:49:ARG:HH11	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:MET:HE3	1:A:271:ASN:CB	2.28	0.63
1:A:229:VAL:CG2	1:A:286:ILE:HG13	2.29	0.63
1:G:146:PRO:O	1:G:197:ILE:HD11	1.99	0.63
1:F:20:MET:SD	1:F:39:ILE:HD11	2.39	0.62
1:D:151:GLY:HA2	1:D:192:TRP:CZ3	2.35	0.62
1:H:42:VAL:HG22	1:H:66:VAL:HG11	1.82	0.62
1:G:27:ARG:NH2	1:G:34:HIS:HD2	1.98	0.61
1:E:229:VAL:HG21	1:E:286:ILE:HG13	1.82	0.61
1:F:31:ASN:O	1:F:123:ASN:HB3	2.01	0.60
1:H:234:TYR:CE2	1:H:238:LEU:HD22	2.37	0.59
1:F:25:HIS:CE1	1:F:78:GLY:HA2	2.38	0.59
1:A:42:VAL:HG11	1:A:89:ILE:HG21	1.85	0.59
1:C:215:LYS:HG2	1:C:236:THR:HG22	1.85	0.58
1:H:111:PRO:HG3	1:H:130:PRO:HB3	1.85	0.58
1:E:94:LEU:HD22	1:E:142:LEU:HB2	1.86	0.58
1:B:211:THR:O	1:B:214:GLN:HB3	2.03	0.58
1:B:229:VAL:CG2	1:B:286:ILE:HD13	2.30	0.58
1:B:104:LYS:HD2	1:B:137:GLU:OE2	2.04	0.57
1:A:192:TRP:CZ3	1:A:208:VAL:HG11	2.40	0.57
1:B:253:ILE:HG12	1:B:287:THR:HG22	1.87	0.57
1:C:219:PRO:C	1:C:221:ASP:H	2.13	0.57
1:E:229:VAL:HG23	1:E:286:ILE:HG21	1.87	0.57
1:G:20:MET:CA	1:G:39:ILE:HD11	2.32	0.57
1:G:21:VAL:HA	2:G:304:HOH:O	2.06	0.56
1:C:38:LEU:HD13	1:C:107:ILE:HD12	1.87	0.56
1:D:27:ARG:NH2	1:D:34:HIS:HD1	2.03	0.56
1:H:45:THR:CG2	1:H:102:ARG:HE	2.18	0.56
1:H:45:THR:HG23	1:H:102:ARG:HE	1.70	0.56
1:H:215:LYS:CB	1:H:236:THR:HG22	2.33	0.56
1:E:20:MET:HB3	1:F:281:ASN:ND2	2.21	0.56
1:E:199:VAL:HG13	1:E:246:GLU:HG3	1.87	0.56
1:F:233:LEU:HD11	1:F:252:ILE:HB	1.88	0.56
1:B:253:ILE:HG23	1:B:287:THR:HG22	1.88	0.55
1:G:170:ARG:HG3	1:G:170:ARG:HH11	1.70	0.55
1:F:50:LYS:HD2	1:F:59:ASP:OD2	2.07	0.55
1:A:229:VAL:HG21	1:A:286:ILE:HG13	1.88	0.55
1:F:146:PRO:HG2	1:F:197:ILE:HG21	1.89	0.54
1:A:102:ARG:NH1	1:A:139:GLU:OE2	2.41	0.54
1:G:39:ILE:HG23	1:G:108:ASP:HB2	1.88	0.54
1:H:20:MET:SD	1:H:39:ILE:HD11	2.48	0.54
1:B:158:LEU:HD13	1:B:185:CYS:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:42:VAL:HG11	1:F:66:VAL:CG2	2.38	0.54
1:F:160:THR:CG2	1:H:35:ILE:HD11	2.38	0.54
1:G:19:ASN:CG	1:G:41:ASN:HD22	2.16	0.53
1:G:129:CYS:O	1:G:129:CYS:SG	2.64	0.53
1:C:253:ILE:HG12	1:C:287:THR:HG22	1.89	0.53
1:F:25:HIS:HD2	1:F:26:ILE:N	2.06	0.53
1:G:243:ARG:HH11	1:G:243:ARG:CG	2.21	0.53
1:H:207:ALA:HB3	1:H:244:ILE:HD11	1.91	0.53
1:H:111:PRO:CG	1:H:130:PRO:HB3	2.38	0.53
1:B:140:GLN:NE2	1:B:146:PRO:HD3	2.24	0.53
1:D:111:PRO:HG3	1:D:130:PRO:HB3	1.90	0.52
1:G:229:VAL:HG21	1:G:286:ILE:HD13	1.88	0.52
1:F:270:SER:O	1:H:75:LYS:HE2	2.09	0.52
1:A:229:VAL:HG23	1:A:286:ILE:HG21	1.92	0.52
1:E:82:ILE:HB	1:E:107:ILE:HG13	1.91	0.52
1:A:117:LYS:HE3	1:C:220:TYR:OH	2.08	0.52
1:F:24:LEU:HD12	1:F:37:GLU:HB2	1.92	0.52
1:A:293:LYS:HZ3	1:B:8:ASN:N	2.08	0.52
1:F:207:ALA:CB	1:F:244:ILE:HD11	2.39	0.52
1:B:20:MET:SD	1:B:39:ILE:HD11	2.50	0.52
1:B:188:VAL:HG11	1:B:212:VAL:HG11	1.91	0.51
1:E:226:SER:OG	1:E:232:THR:OG1	2.24	0.51
1:A:262:ILE:O	1:A:271:ASN:ND2	2.41	0.51
1:E:17:GLY:HA3	1:E:42:VAL:O	2.09	0.51
1:E:192:TRP:CZ3	1:E:208:VAL:HG11	2.45	0.51
1:A:134:ARG:HA	1:A:205:TRP:CH2	2.44	0.51
1:G:18:LYS:C	1:G:19:ASN:HD22	2.18	0.51
1:C:195:ASN:HB3	1:C:248:GLU:HB2	1.92	0.51
1:F:156:LYS:O	1:H:127:ILE:HD12	2.10	0.51
1:H:192:TRP:CZ3	1:H:208:VAL:HG11	2.46	0.50
1:A:117:LYS:CE	1:C:220:TYR:OH	2.59	0.50
1:A:188:VAL:HG11	1:A:212:VAL:HG11	1.92	0.50
1:E:146:PRO:HG2	1:E:197:ILE:HG21	1.92	0.50
1:E:205:TRP:C	1:E:205:TRP:CD1	2.89	0.50
1:D:104:LYS:HE2	1:D:137:GLU:OE2	2.12	0.50
1:E:57:ASN:O	1:F:168:PHE:HA	2.11	0.50
1:D:186:THR:OG1	1:D:187:SER:N	2.45	0.50
1:G:22:LYS:HG2	1:G:39:ILE:HD13	1.94	0.50
1:H:222:ARG:NH1	1:H:222:ARG:HB3	2.27	0.49
1:C:38:LEU:HD21	1:C:79:ILE:HD13	1.92	0.49
1:H:35:ILE:HD12	1:H:126:PHE:CE1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:240:VAL:HG12	1:D:247:VAL:HG21	1.95	0.49
1:G:146:PRO:HG2	1:G:197:ILE:HG12	1.93	0.49
1:C:149:HIS:NE2	1:C:193:ARG:HB2	2.28	0.49
1:D:159:LYS:HG3	1:D:216:PHE:CZ	2.48	0.49
1:D:244:ILE:HG22	1:D:247:VAL:HG13	1.92	0.49
1:E:134:ARG:HG3	1:E:205:TRP:CE3	2.48	0.49
1:E:273:ASP:HB3	1:G:28:ARG:NH2	2.27	0.49
1:C:71:HIS:HD2	2:C:320:HOH:O	1.94	0.49
1:D:32:HIS:HD2	1:D:34:HIS:NE2	2.11	0.49
1:E:151:GLY:HA2	1:E:192:TRP:CZ3	2.48	0.49
1:E:159:LYS:HA	1:G:124:HIS:O	2.13	0.49
1:F:63:THR:HA	1:F:66:VAL:HG12	1.94	0.49
1:B:24:LEU:C	1:B:24:LEU:HD23	2.38	0.48
1:B:49:ARG:HG3	1:B:49:ARG:NH1	2.28	0.48
1:E:140:GLN:HB2	1:E:146:PRO:HB3	1.94	0.48
1:B:105:VAL:HG12	1:B:107:ILE:CD1	2.44	0.48
1:C:158:LEU:HD23	1:C:159:LYS:N	2.28	0.48
1:D:45:THR:HB	1:D:102:ARG:HB3	1.95	0.48
1:A:32:HIS:HD2	1:A:34:HIS:CE1	2.30	0.48
1:D:161:THR:OG1	1:D:162:GLN:N	2.47	0.48
1:E:244:ILE:HG22	1:E:246:GLU:HG2	1.95	0.48
1:H:214:GLN:HE22	1:H:243:ARG:HH22	1.61	0.48
1:C:151:GLY:HA2	1:C:192:TRP:CZ3	2.49	0.47
1:C:251:GLU:HG3	1:C:289:THR:OG1	2.13	0.47
1:E:40:ALA:HA	1:E:107:ILE:HG22	1.96	0.47
1:B:233:LEU:HD11	1:B:252:ILE:CG2	2.44	0.47
1:B:31:ASN:HD21	1:D:222:ARG:HH11	1.61	0.47
1:E:63:THR:HA	1:E:66:VAL:HG12	1.97	0.47
1:D:195:ASN:HB3	1:D:248:GLU:HB2	1.97	0.47
1:E:205:TRP:CD1	1:E:206:LYS:N	2.83	0.47
1:F:24:LEU:C	1:F:24:LEU:HD23	2.40	0.47
1:G:192:TRP:CZ3	1:G:208:VAL:HG11	2.49	0.47
1:G:219:PRO:C	1:G:221:ASP:H	2.23	0.47
1:C:11:PHE:CD2	1:C:49:ARG:HD3	2.50	0.47
1:G:168:PHE:HA	1:H:57:ASN:O	2.15	0.47
1:G:199:VAL:HG13	1:G:246:GLU:HG3	1.97	0.47
1:G:82:ILE:HD11	1:G:134:ARG:HG2	1.97	0.46
1:H:20:MET:HA	1:H:39:ILE:HD11	1.97	0.46
1:C:292:ARG:O	1:C:294:PRO:HD3	2.15	0.46
1:E:293:LYS:HD2	1:F:8:ASN:N	2.30	0.46
1:B:146:PRO:HG2	1:B:197:ILE:HG21	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:GLU:O	1:D:26:ILE:HG22	2.15	0.46
1:F:229:VAL:HG23	1:F:286:ILE:HG21	1.97	0.46
1:A:17:GLY:HA3	1:A:42:VAL:O	2.15	0.46
1:A:232:THR:O	1:A:236:THR:HG23	2.14	0.46
1:H:249:GLU:HG3	1:H:291:CYS:HB3	1.97	0.46
1:B:105:VAL:HG12	1:B:107:ILE:HD12	1.98	0.46
1:E:9:VAL:HG21	1:F:238:LEU:HD11	1.97	0.46
1:E:281:ASN:HD21	1:F:20:MET:HE3	1.81	0.46
1:G:195:ASN:HB3	1:G:248:GLU:HB2	1.97	0.46
1:E:20:MET:HE3	1:F:281:ASN:HD21	1.78	0.46
1:F:31:ASN:O	1:F:123:ASN:CG	2.59	0.46
1:G:122:HIS:HE1	1:G:124:HIS:CD2	2.34	0.46
1:H:231:LYS:HE3	1:H:235:ASP:OD2	2.16	0.45
1:A:130:PRO:HD2	1:C:130:PRO:HD2	1.97	0.45
1:A:28:ARG:HG2	1:A:33:HIS:CD2	2.51	0.45
1:B:229:VAL:HG23	1:B:286:ILE:HG21	1.98	0.45
1:D:20:MET:N	1:D:41:ASN:HD22	2.07	0.45
1:A:82:ILE:HG21	1:A:109:GLU:HB2	1.97	0.45
1:C:31:ASN:O	1:C:123:ASN:HB3	2.17	0.45
1:C:158:LEU:HD23	1:C:158:LEU:C	2.41	0.45
1:C:86:ALA:HA	1:C:105:VAL:HG21	1.98	0.45
1:E:270:SER:O	1:G:75:LYS:HE2	2.16	0.45
1:A:102:ARG:HH12	1:A:104:LYS:HD2	1.80	0.45
1:G:114:ARG:HD2	1:G:121:GLU:HB3	1.97	0.45
1:B:124:HIS:HB2	1:D:160:THR:OG1	2.17	0.45
1:F:25:HIS:HE1	1:F:78:GLY:HA2	1.80	0.45
1:A:226:SER:HA	1:A:227:PRO:HD3	1.79	0.45
1:C:14:THR:OG1	1:C:49:ARG:NH1	2.40	0.45
1:C:232:THR:O	1:C:236:THR:HG23	2.16	0.45
1:G:249:GLU:HG3	1:G:291:CYS:HB3	1.97	0.45
1:B:192:TRP:CZ3	1:B:208:VAL:HG11	2.52	0.44
1:G:272:LYS:HB2	1:G:272:LYS:HE3	1.74	0.44
1:F:117:LYS:O	1:F:120:VAL:HG12	2.17	0.44
1:F:158:LEU:O	1:H:126:PHE:N	2.48	0.44
1:G:243:ARG:C	1:G:244:ILE:HG13	2.41	0.44
1:A:151:GLY:O	1:A:205:TRP:CZ3	2.56	0.44
1:A:163:THR:HG21	1:A:184:PHE:HB2	1.99	0.44
1:C:107:ILE:HD13	1:C:107:ILE:HG21	1.72	0.44
1:C:159:LYS:HB2	1:C:216:PHE:CZ	2.52	0.44
1:E:231:LYS:HG3	1:F:55:GLY:N	2.33	0.44
1:B:195:ASN:HB3	1:B:248:GLU:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:PRO:C	1:A:221:ASP:H	2.24	0.44
1:D:244:ILE:HG21	1:D:247:VAL:HG13	2.00	0.44
1:A:127:ILE:HD13	1:C:131:GLU:HG2	2.00	0.44
1:C:9:VAL:HG21	1:D:238:LEU:HD11	1.99	0.44
1:E:122:HIS:CD2	1:G:220:TYR:HA	2.52	0.44
1:H:32:HIS:CD2	1:H:34:HIS:HE2	2.36	0.44
1:G:17:GLY:HA3	1:G:42:VAL:O	2.18	0.43
1:F:205:TRP:CD1	1:F:206:LYS:N	2.86	0.43
1:F:207:ALA:HB1	1:F:244:ILE:HD11	1.99	0.43
1:F:293:LYS:HA	1:F:294:PRO:HD2	1.81	0.43
1:G:45:THR:CG2	1:G:102:ARG:HH21	2.31	0.43
1:G:237:GLN:HG3	1:G:252:ILE:HD12	2.00	0.43
1:D:237:GLN:NE2	1:D:288:GLY:O	2.51	0.43
1:E:283:SER:C	1:F:18:LYS:HD2	2.43	0.43
1:F:151:GLY:HA2	1:F:192:TRP:CZ3	2.52	0.43
1:F:163:THR:HG21	1:F:184:PHE:HB2	2.00	0.43
1:G:192:TRP:HB2	1:G:247:VAL:HG13	2.00	0.43
1:H:129:CYS:O	1:H:129:CYS:SG	2.76	0.43
1:F:156:LYS:NZ	1:F:257:GLN:OE1	2.48	0.43
1:D:87:LEU:HD21	1:D:197:ILE:HD11	2.01	0.43
1:G:45:THR:HG23	1:G:102:ARG:HE	1.83	0.43
1:F:214:GLN:HE22	1:F:243:ARG:HH22	1.65	0.43
1:A:240:VAL:HG13	1:A:244:ILE:HD12	2.00	0.43
1:B:239:LEU:O	1:B:243:ARG:HG3	2.19	0.43
1:E:114:ARG:HD2	1:E:121:GLU:HB3	2.01	0.43
1:E:129:CYS:HA	1:E:130:PRO:HD3	1.88	0.43
1:F:199:VAL:HG22	1:F:246:GLU:HG3	1.99	0.43
1:B:19:ASN:ND2	1:B:41:ASN:CG	2.73	0.42
1:E:134:ARG:HG3	1:E:205:TRP:CZ3	2.54	0.42
1:H:244:ILE:O	1:H:247:VAL:HG22	2.19	0.42
1:B:204:ALA:O	1:B:205:TRP:C	2.61	0.42
1:E:219:PRO:C	1:E:221:ASP:H	2.27	0.42
1:A:66:VAL:HG23	1:A:93:PHE:CZ	2.54	0.42
1:C:180:LYS:HB3	1:C:180:LYS:HE3	1.95	0.42
1:E:37:GLU:HB3	1:E:110:VAL:HB	2.00	0.42
1:F:31:ASN:O	1:F:123:ASN:CB	2.66	0.42
1:B:43:GLN:OE1	1:B:102:ARG:NH2	2.53	0.42
1:F:25:HIS:CD2	1:F:26:ILE:N	2.86	0.42
1:A:233:LEU:C	1:A:233:LEU:HD23	2.44	0.42
1:C:38:LEU:HD22	1:C:107:ILE:CD1	2.42	0.42
1:D:19:ASN:OD1	1:D:41:ASN:HB3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:273:ASP:OD1	1:E:273:ASP:N	2.50	0.42
1:F:116:GLU:HG3	1:F:121:GLU:HG2	2.02	0.42
1:G:114:ARG:CD	1:G:121:GLU:HB3	2.50	0.41
1:E:20:MET:CE	1:F:281:ASN:HD21	2.33	0.41
1:F:192:TRP:CZ3	1:F:208:VAL:HG11	2.54	0.41
1:G:234:TYR:HA	1:G:237:GLN:OE1	2.20	0.41
1:E:253:ILE:HG12	1:E:287:THR:HG22	2.02	0.41
1:H:241:LEU:HD23	1:H:247:VAL:CG2	2.44	0.41
1:F:217:ALA:HB1	1:H:122:HIS:CD2	2.56	0.41
1:G:144:LYS:HA	1:G:144:LYS:HD3	1.93	0.41
1:D:272:LYS:O	1:D:272:LYS:CG	2.69	0.41
1:A:239:LEU:O	1:A:243:ARG:HG3	2.21	0.41
1:F:253:ILE:HG12	1:F:287:THR:CG2	2.48	0.41
1:G:122:HIS:CE1	1:G:124:HIS:CD2	3.09	0.41
1:H:104:LYS:HE2	1:H:137:GLU:OE2	2.21	0.41
1:E:281:ASN:ND2	1:F:20:MET:HE3	2.36	0.41
1:C:86:ALA:HB1	1:C:105:VAL:CG2	2.50	0.41
1:A:52:TYR:HB3	1:B:234:TYR:HB2	2.02	0.41
1:A:82:ILE:HB	1:A:107:ILE:HD11	2.02	0.41
1:B:170:ARG:NH2	1:B:176:LEU:O	2.54	0.41
1:E:185:CYS:HB3	1:E:257:GLN:O	2.20	0.41
1:H:239:LEU:HD12	1:H:239:LEU:HA	1.90	0.41
1:H:245:PRO:HA	1:H:292:ARG:NH2	2.36	0.41
1:B:233:LEU:HD11	1:B:252:ILE:HG22	2.02	0.40
1:D:42:VAL:HG22	1:D:66:VAL:HG11	2.02	0.40
1:E:152:LEU:HB3	1:E:205:TRP:CH2	2.56	0.40
1:F:226:SER:HA	1:F:227:PRO:HD3	1.87	0.40
1:H:195:ASN:OD1	1:H:196:THR:HG22	2.22	0.40
1:A:152:LEU:HB3	1:A:205:TRP:CH2	2.56	0.40
1:D:89:ILE:HB	1:D:105:VAL:HG21	2.04	0.40
1:G:111:PRO:HG3	1:G:133:LEU:CD2	2.50	0.40
1:H:232:THR:O	1:H:236:THR:HG23	2.21	0.40
1:C:192:TRP:CZ3	1:C:208:VAL:HG11	2.57	0.40
1:C:253:ILE:HG23	1:C:287:THR:HG22	2.04	0.40
1:F:108:ASP:OD1	1:F:135:PHE:HB3	2.22	0.40
1:A:9:VAL:HG11	1:B:238:LEU:HD11	2.04	0.40
1:E:24:LEU:HD23	1:E:24:LEU:C	2.46	0.40
1:G:35:ILE:HD12	1:G:35:ILE:HA	1.92	0.40
1:G:127:ILE:HG13	1:G:128:HIS:N	2.37	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	285/298 (96%)	277 (97%)	8 (3%)	0	100	100
1	B	285/298 (96%)	276 (97%)	9 (3%)	0	100	100
1	C	285/298 (96%)	275 (96%)	10 (4%)	0	100	100
1	D	285/298 (96%)	276 (97%)	9 (3%)	0	100	100
1	E	285/298 (96%)	275 (96%)	10 (4%)	0	100	100
1	F	285/298 (96%)	274 (96%)	11 (4%)	0	100	100
1	G	285/298 (96%)	272 (95%)	13 (5%)	0	100	100
1	H	285/298 (96%)	278 (98%)	6 (2%)	1 (0%)	30	60
All	All	2280/2384 (96%)	2203 (97%)	76 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	143	SER

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	258/267 (97%)	248 (96%)	10 (4%)	28	64
1	B	258/267 (97%)	241 (93%)	17 (7%)	15	43
1	C	258/267 (97%)	247 (96%)	11 (4%)	26	60
1	D	258/267 (97%)	243 (94%)	15 (6%)	18	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	258/267 (97%)	246 (95%)	12 (5%)	23	57
1	F	258/267 (97%)	246 (95%)	12 (5%)	23	57
1	G	258/267 (97%)	244 (95%)	14 (5%)	20	51
1	H	258/267 (97%)	245 (95%)	13 (5%)	22	54
All	All	2064/2136 (97%)	1960 (95%)	104 (5%)	22	54

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	VAL
1	A	35	ILE
1	A	103	VAL
1	A	107	ILE
1	A	117	LYS
1	A	120	VAL
1	A	169	LEU
1	A	229	VAL
1	A	246	GLU
1	A	270	SER
1	B	12	VAL
1	B	39	ILE
1	B	77	LYS
1	B	80	LYS
1	B	107	ILE
1	B	116	GLU
1	B	120	VAL
1	B	143	SER
1	B	172	ARG
1	B	197	ILE
1	B	213	ILE
1	B	229	VAL
1	B	238	LEU
1	B	241	LEU
1	B	246	GLU
1	B	247	VAL
1	B	250	ILE
1	C	23	VAL
1	C	42	VAL
1	C	75	LYS
1	C	103	VAL
1	C	107	ILE

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Mol	Chain	Res	Type
1	C	120	VAL
1	C	131	GLU
1	C	143	SER
1	C	172	ARG
1	C	229	VAL
1	C	239	LEU
1	D	8	ASN
1	D	12	VAL
1	D	29	GLU
1	D	41	ASN
1	D	42	VAL
1	D	103	VAL
1	D	117	LYS
1	D	196	THR
1	D	197	ILE
1	D	222	ARG
1	D	238	LEU
1	D	246	GLU
1	D	247	VAL
1	D	283	SER
1	D	293	LYS
1	E	12	VAL
1	E	35	ILE
1	E	103	VAL
1	E	107	ILE
1	E	143	SER
1	E	213	ILE
1	E	229	VAL
1	E	238	LEU
1	E	239	LEU
1	E	246	GLU
1	E	261	VAL
1	E	272	LYS
1	F	12	VAL
1	F	13	ARG
1	F	23	VAL
1	F	29	GLU
1	F	39	ILE
1	F	53	LEU
1	F	105	VAL
1	F	131	GLU
1	F	145	THR

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Mol	Chain	Res	Type
1	F	148	VAL
1	F	197	ILE
1	F	246	GLU
1	G	10	GLU
1	G	23	VAL
1	G	35	ILE
1	G	39	ILE
1	G	45	THR
1	G	103	VAL
1	G	145	THR
1	G	197	ILE
1	G	198	ASN
1	G	229	VAL
1	G	243	ARG
1	G	246	GLU
1	G	261	VAL
1	G	272	LYS
1	H	12	VAL
1	H	35	ILE
1	H	39	ILE
1	H	42	VAL
1	H	45	THR
1	H	76	LEU
1	H	103	VAL
1	H	120	VAL
1	H	197	ILE
1	H	238	LEU
1	H	239	LEU
1	H	266	LYS
1	H	293	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	34	HIS
1	A	106	ASN
1	A	214	GLN
1	A	281	ASN
1	B	19	ASN
1	B	31	ASN
1	B	34	HIS

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Mol	Chain	Res	Type
1	B	71	HIS
1	B	162	GLN
1	B	281	ASN
1	B	285	ASN
1	C	34	HIS
1	C	68	ASN
1	C	71	HIS
1	C	281	ASN
1	D	8	ASN
1	D	32	HIS
1	D	41	ASN
1	D	43	GLN
1	D	214	GLN
1	E	8	ASN
1	E	71	HIS
1	E	140	GLN
1	E	162	GLN
1	F	25	HIS
1	F	34	HIS
1	F	41	ASN
1	F	71	HIS
1	F	140	GLN
1	G	8	ASN
1	G	19	ASN
1	G	34	HIS
1	G	41	ASN
1	G	71	HIS
1	G	92	HIS
1	G	122	HIS
1	G	123	ASN
1	G	124	HIS
1	H	32	HIS
1	H	92	HIS
1	H	106	ASN
1	H	281	ASN

5.3.3 RNA ⓘ

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	287/298 (96%)	-0.19	2 (0%) 84 77	14, 29, 48, 78	0
1	B	287/298 (96%)	-0.35	1 (0%) 90 86	12, 25, 46, 67	0
1	C	287/298 (96%)	-0.33	2 (0%) 84 77	13, 25, 49, 63	0
1	D	287/298 (96%)	-0.43	1 (0%) 90 86	13, 24, 44, 68	0
1	E	287/298 (96%)	-0.15	0 100 100	18, 31, 51, 85	0
1	F	287/298 (96%)	-0.23	2 (0%) 84 77	17, 30, 50, 67	0
1	G	287/298 (96%)	-0.13	2 (0%) 84 77	16, 32, 48, 75	0
1	H	287/298 (96%)	-0.05	2 (0%) 84 77	17, 33, 53, 71	0
All	All	2296/2384 (96%)	-0.23	12 (0%) 87 82	12, 28, 49, 85	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	294	PRO	3.6
1	A	8	ASN	3.6
1	H	8	ASN	3.5
1	A	294	PRO	2.8
1	D	8	ASN	2.8
1	C	294	PRO	2.6
1	F	294	PRO	2.5
1	B	294	PRO	2.4
1	H	294	PRO	2.4
1	F	8	ASN	2.3
1	G	8	ASN	2.3
1	C	29	GLU	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.