



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:09 AM UTC

PDB ID : 1VL2 / pdb_00001vl2
Title : Crystal structure of Argininosuccinate synthase (TM1780) from *Thermotoga maritima* at 1.65 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2004-07-08
Resolution : 1.65 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

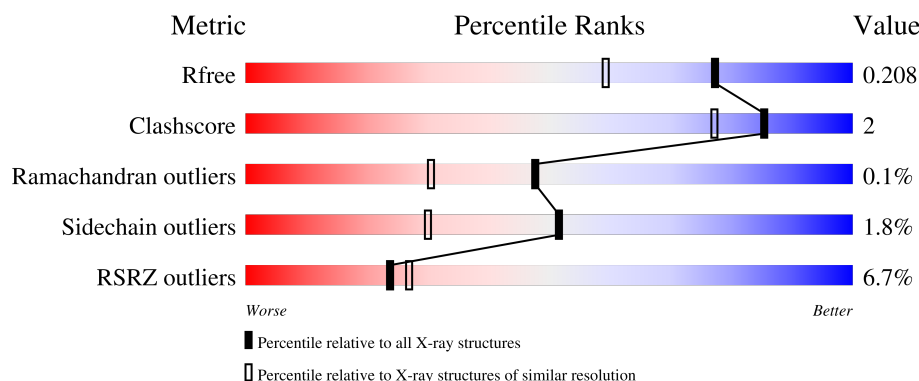
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

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	421	10% 87% 7% • 5%
1	B	421	5% 88% 5% 6%
1	C	421	5% 88% 5% 7%
1	D	421	5% 87% 5% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13531 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 Argininosuccinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	398	Total	C	N	O	S	0	2	0
			3117	2007	521	577	12			
1	B	394	Total	C	N	O	S	0	2	0
			3104	1997	516	579	12			
1	C	393	Total	C	N	O	S	0	2	0
			3089	1990	511	576	12			
1	D	389	Total	C	N	O	S	0	2	0
			3037	1961	501	563	12			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9X2A1
A	-10	GLY	-	expression tag	UNP Q9X2A1
A	-9	SER	-	expression tag	UNP Q9X2A1
A	-8	ASP	-	expression tag	UNP Q9X2A1
A	-7	LYS	-	expression tag	UNP Q9X2A1
A	-6	ILE	-	expression tag	UNP Q9X2A1
A	-5	HIS	-	expression tag	UNP Q9X2A1
A	-4	HIS	-	expression tag	UNP Q9X2A1
A	-3	HIS	-	expression tag	UNP Q9X2A1
A	-2	HIS	-	expression tag	UNP Q9X2A1
A	-1	HIS	-	expression tag	UNP Q9X2A1
A	0	HIS	-	expression tag	UNP Q9X2A1
B	-11	MET	-	expression tag	UNP Q9X2A1
B	-10	GLY	-	expression tag	UNP Q9X2A1
B	-9	SER	-	expression tag	UNP Q9X2A1
B	-8	ASP	-	expression tag	UNP Q9X2A1
B	-7	LYS	-	expression tag	UNP Q9X2A1
B	-6	ILE	-	expression tag	UNP Q9X2A1
B	-5	HIS	-	expression tag	UNP Q9X2A1
B	-4	HIS	-	expression tag	UNP Q9X2A1
B	-3	HIS	-	expression tag	UNP Q9X2A1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	expression tag	UNP Q9X2A1
B	-1	HIS	-	expression tag	UNP Q9X2A1
B	0	HIS	-	expression tag	UNP Q9X2A1
C	-11	MET	-	expression tag	UNP Q9X2A1
C	-10	GLY	-	expression tag	UNP Q9X2A1
C	-9	SER	-	expression tag	UNP Q9X2A1
C	-8	ASP	-	expression tag	UNP Q9X2A1
C	-7	LYS	-	expression tag	UNP Q9X2A1
C	-6	ILE	-	expression tag	UNP Q9X2A1
C	-5	HIS	-	expression tag	UNP Q9X2A1
C	-4	HIS	-	expression tag	UNP Q9X2A1
C	-3	HIS	-	expression tag	UNP Q9X2A1
C	-2	HIS	-	expression tag	UNP Q9X2A1
C	-1	HIS	-	expression tag	UNP Q9X2A1
C	0	HIS	-	expression tag	UNP Q9X2A1
D	-11	MET	-	expression tag	UNP Q9X2A1
D	-10	GLY	-	expression tag	UNP Q9X2A1
D	-9	SER	-	expression tag	UNP Q9X2A1
D	-8	ASP	-	expression tag	UNP Q9X2A1
D	-7	LYS	-	expression tag	UNP Q9X2A1
D	-6	ILE	-	expression tag	UNP Q9X2A1
D	-5	HIS	-	expression tag	UNP Q9X2A1
D	-4	HIS	-	expression tag	UNP Q9X2A1
D	-3	HIS	-	expression tag	UNP Q9X2A1
D	-2	HIS	-	expression tag	UNP Q9X2A1
D	-1	HIS	-	expression tag	UNP Q9X2A1
D	0	HIS	-	expression tag	UNP Q9X2A1

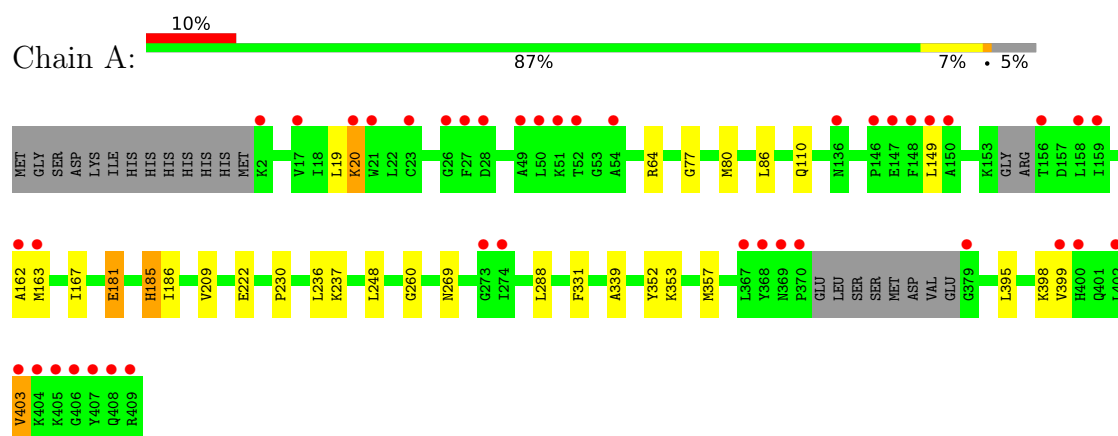
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	315	Total O 315 315	0	0
2	B	298	Total O 298 298	0	0
2	C	312	Total O 312 312	0	0
2	D	259	Total O 259 259	0	0

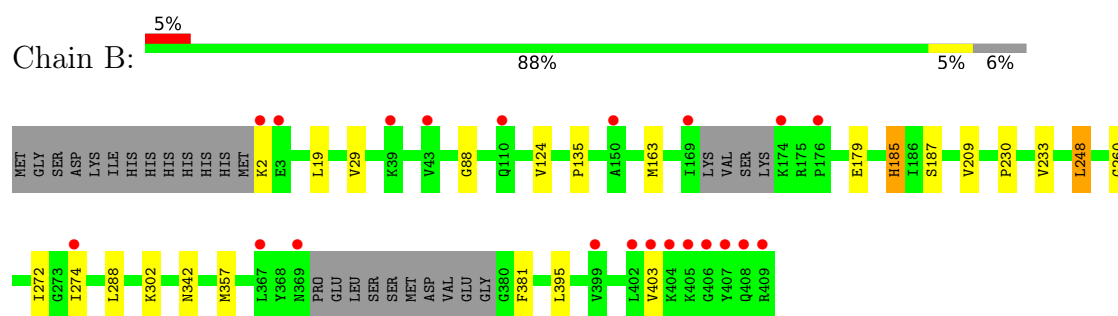
3 Residue-property plots [i](#)

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.

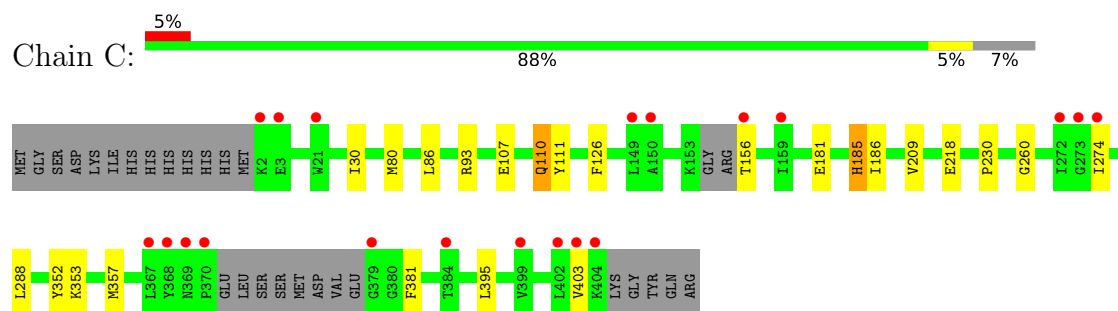
- Molecule 1: Argininosuccinate synthase



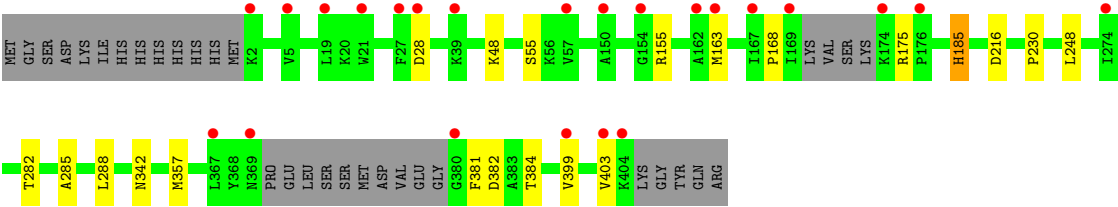
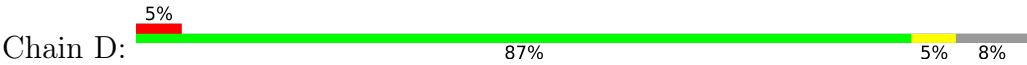
- Molecule 1: Argininosuccinate synthase



- Molecule 1: Argininosuccinate synthase



- Molecule 1: Argininosuccinate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.32Å 114.91Å 149.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.21 – 1.65 49.21 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.21-1.65) 99.7 (49.21-1.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.2.0001, CNS	Depositor
R, R_{free}	0.177 , 0.209 (Not available) , 0.208	Depositor DCC
R_{free} test set	9757 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13531	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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 49.39 % 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 7.5130e-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	1.04	3/3197 (0.1%)	0.96	0/4321
1	B	0.99	0/3184	0.95	2/4308 (0.0%)
1	C	1.02	0/3168	0.94	0/4285
1	D	1.01	0/3116	0.95	1/4225 (0.0%)
All	All	1.01	3/12665 (0.0%)	0.95	3/17139 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	77	GLY	C-O	-5.53	1.19	1.24
1	A	339	ALA	CA-CB	-5.15	1.44	1.53
1	A	331	PHE	C-O	-5.04	1.18	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	342	ASN	N-CA-C	5.90	120.29	113.38
1	B	88	GLY	N-CA-C	5.28	119.27	112.83
1	D	342	ASN	N-CA-C	5.04	119.02	112.86

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	3117	0	3063	18	0
1	B	3104	0	3043	18	0
1	C	3089	0	3035	13	0
1	D	3037	0	2962	13	0
2	A	315	0	0	6	0
2	B	298	0	0	2	0
2	C	312	0	0	2	0
2	D	259	0	0	3	0
All	All	13531	0	12103	54	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 2.

All (54) 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:C:230:PRO:HG3	1:C:288:LEU:HD11	1.63	0.80
1:B:233:VAL:HG23	1:B:248:LEU:HG	1.69	0.74
1:C:357[B]:MET:SD	2:D:532:HOH:O	2.50	0.68
1:D:230:PRO:HG3	1:D:288:LEU:HD11	1.74	0.67
1:A:230:PRO:HG3	1:A:288:LEU:HD11	1.76	0.67
2:A:571:HOH:O	1:B:357[B]:MET:SD	2.52	0.67
1:A:357[B]:MET:SD	2:B:547:HOH:O	2.52	0.66
2:C:628:HOH:O	1:D:357[B]:MET:HG2	1.95	0.66
2:A:650:HOH:O	1:B:357[B]:MET:HG2	1.97	0.65
1:A:80:MET:HG2	1:A:86:LEU:HD23	1.79	0.65
1:D:357[B]:MET:SD	2:D:441:HOH:O	2.55	0.63
1:B:274:ILE:HD13	1:D:399:VAL:HG21	1.81	0.63
1:B:230:PRO:HG3	1:B:288:LEU:HD11	1.83	0.61
1:A:237:LYS:HB2	2:A:634:HOH:O	2.03	0.58
1:B:19:LEU:HD13	1:B:29:VAL:HG11	1.86	0.58
1:B:381:PHE:CD1	1:C:395:LEU:HD11	2.39	0.56
1:C:110:GLN:HG2	1:C:111:TYR:CD2	2.42	0.55
1:C:181:GLU:HG3	1:C:186:ILE:HG12	1.91	0.53
1:A:162:ALA:HA	1:A:167:ILE:HD12	1.92	0.51
1:A:395:LEU:HD11	1:D:381:PHE:CD1	2.46	0.51
2:C:628:HOH:O	1:D:357[B]:MET:CG	2.58	0.50
1:B:2:LYS:N	2:B:646:HOH:O	2.45	0.50
1:B:233:VAL:CG2	1:B:248:LEU:HG	2.40	0.49
2:A:434:HOH:O	1:B:357[B]:MET:SD	2.60	0.49
1:C:93:ARG:HD2	1:C:126:PHE:CE2	2.48	0.49
1:A:222:GLU:HG3	1:A:236:LEU:HD21	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:ARG:HD3	2:A:544:HOH:O	2.14	0.48
1:D:382:ASP:OD1	1:D:384:THR:HB	2.15	0.47
1:C:30:ILE:HD13	1:C:107:GLU:HG3	1.98	0.46
1:B:124:VAL:CG2	1:B:272:ILE:HD12	2.46	0.46
1:B:395:LEU:HD11	1:C:381:PHE:CD1	2.50	0.46
1:A:209:VAL:O	1:A:260:GLY:HA2	2.16	0.46
1:A:20:LYS:HG3	1:A:167:ILE:HD11	1.98	0.45
1:D:216:ASP:CB	2:D:483:HOH:O	2.63	0.45
1:C:80:MET:HG2	1:C:86:LEU:HD23	1.99	0.45
1:A:398:LYS:HG2	1:B:135:PRO:HB3	2.00	0.44
1:A:237:LYS:CB	2:A:634:HOH:O	2.64	0.44
1:A:181:GLU:HB2	1:A:186:ILE:HG12	2.00	0.43
1:C:209:VAL:O	1:C:260:GLY:HA2	2.17	0.43
1:B:274:ILE:HD13	1:D:399:VAL:CG2	2.46	0.43
1:A:352:TYR:O	1:A:353:LYS:C	2.63	0.42
1:D:48:LYS:HE3	1:D:168:PRO:O	2.19	0.42
1:A:185:HIS:CD2	1:A:185:HIS:C	2.97	0.42
1:B:185:HIS:CD2	1:B:185:HIS:C	2.98	0.41
1:C:352:TYR:O	1:C:353:LYS:C	2.63	0.41
1:A:399:VAL:HG21	1:C:274:ILE:HD13	2.02	0.41
1:B:179:GLU:HA	1:B:187:SER:O	2.20	0.41
1:D:282:THR:HG22	1:D:285:ALA:HB3	2.02	0.41
1:D:185:HIS:CD2	1:D:185:HIS:C	2.98	0.41
1:B:209:VAL:O	1:B:260:GLY:HA2	2.20	0.41
1:D:28:ASP:OD1	1:D:55:SER:HB3	2.21	0.41
1:C:185:HIS:CD2	1:C:185:HIS:C	2.98	0.40
1:A:269:ASN:HB3	1:B:302:LYS:HD3	2.03	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	394/421 (94%)	380 (96%)	13 (3%)	1 (0%)	36	21
1	B	390/421 (93%)	376 (96%)	14 (4%)	0	100	100
1	C	389/421 (92%)	373 (96%)	15 (4%)	1 (0%)	36	21
1	D	385/421 (91%)	373 (97%)	12 (3%)	0	100	100
All	All	1558/1684 (92%)	1502 (96%)	54 (4%)	2 (0%)	48	30

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	403	VAL
1	C	403	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	321/356 (90%)	312 (97%)	9 (3%)	38	15
1	B	321/356 (90%)	317 (99%)	4 (1%)	63	45
1	C	320/356 (90%)	316 (99%)	4 (1%)	61	42
1	D	310/356 (87%)	304 (98%)	6 (2%)	50	28
All	All	1272/1424 (89%)	1249 (98%)	23 (2%)	51	30

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	20	LYS
1	A	110	GLN
1	A	149	LEU
1	A	163	MET
1	A	181	GLU
1	A	185	HIS
1	A	248	LEU
1	A	403	VAL

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Mol	Chain	Res	Type
1	B	163	MET
1	B	185	HIS
1	B	248	LEU
1	B	403	VAL
1	C	110	GLN
1	C	156	THR
1	C	185	HIS
1	C	218	GLU
1	D	155	ARG
1	D	163	MET
1	D	175	ARG
1	D	185	HIS
1	D	248	LEU
1	D	403	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	D	393	HIS

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers

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	398/421 (94%)	0.28	41 (10%)	12 12	8, 16, 38, 49	2 (0%)
1	B	394/421 (93%)	0.10	21 (5%)	32 36	8, 16, 33, 57	2 (0%)
1	C	393/421 (93%)	0.09	20 (5%)	33 37	8, 16, 35, 61	2 (0%)
1	D	389/421 (92%)	0.26	23 (5%)	28 32	8, 17, 35, 57	2 (0%)
All	All	1574/1684 (93%)	0.18	105 (6%)	24 27	8, 16, 36, 61	8 (0%)

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	169	ILE	6.6
1	D	367	LEU	6.4
1	C	274	ILE	6.2
1	A	403	VAL	5.8
1	B	403	VAL	5.8
1	D	403	VAL	5.4
1	C	403	VAL	5.2
1	A	405	LYS	4.9
1	C	367	LEU	4.9
1	B	406	GLY	4.7
1	B	402	LEU	4.5
1	A	407	TYR	4.5
1	D	169	ILE	4.5
1	B	407	TYR	4.4
1	D	404	LYS	4.2
1	A	150	ALA	4.1
1	B	405	LYS	4.1
1	A	406	GLY	4.0
1	B	367	LEU	3.9
1	A	159	ILE	3.8
1	A	21	TRP	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	404	LYS	3.8
1	A	402	LEU	3.7
1	A	370	PRO	3.7
1	C	2	LYS	3.7
1	D	150	ALA	3.6
1	C	399	VAL	3.6
1	D	2	LYS	3.5
1	A	367	LEU	3.5
1	C	159	ILE	3.5
1	C	273	GLY	3.4
1	C	379	GLY	3.4
1	D	274	ILE	3.3
1	A	379	GLY	3.3
1	B	274	ILE	3.3
1	D	5	VAL	3.2
1	A	50	LEU	3.1
1	C	150	ALA	3.1
1	D	369	ASN	3.1
1	A	399	VAL	3.1
1	C	272	ILE	3.0
1	A	146	PRO	3.0
1	C	21	TRP	3.0
1	C	156	THR	3.0
1	A	2	LYS	3.0
1	A	400	HIS	2.9
1	A	163	MET	2.9
1	B	399	VAL	2.8
1	D	57	VAL	2.8
1	A	369	ASN	2.8
1	D	174	LYS	2.8
1	A	149	LEU	2.8
1	B	409	ARG	2.8
1	A	52	THR	2.8
1	A	23	CYS	2.8
1	B	404	LYS	2.7
1	B	150	ALA	2.7
1	A	51	LYS	2.6
1	C	402	LEU	2.6
1	A	162	ALA	2.6
1	D	154	GLY	2.6
1	C	370	PRO	2.6
1	A	28	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	409	ARG	2.5
1	A	156	THR	2.5
1	C	149	LEU	2.5
1	A	27	PHE	2.5
1	A	404	LYS	2.4
1	D	163	MET	2.4
1	D	176	PRO	2.4
1	A	49	ALA	2.4
1	A	368	TYR	2.4
1	A	54	ALA	2.4
1	D	167	ILE	2.4
1	B	174	LYS	2.4
1	B	408	GLN	2.4
1	A	158	LEU	2.4
1	D	19	LEU	2.4
1	A	274	ILE	2.4
1	A	408	GLN	2.3
1	C	368	TYR	2.3
1	D	162	ALA	2.3
1	C	369	ASN	2.3
1	B	176	PRO	2.3
1	D	39	LYS	2.3
1	C	384	THR	2.2
1	A	148	PHE	2.2
1	D	27	PHE	2.2
1	D	380	GLY	2.2
1	B	43	VAL	2.2
1	A	17	VAL	2.2
1	A	26	GLY	2.2
1	B	2	LYS	2.1
1	A	147	GLU	2.1
1	B	369	ASN	2.1
1	D	399	VAL	2.1
1	D	28	ASP	2.1
1	A	20	LYS	2.1
1	A	136	ASN	2.0
1	B	3	GLU	2.0
1	C	3	GLU	2.0
1	A	273	GLY	2.0
1	B	110	GLN	2.0
1	D	21	TRP	2.0
1	B	39	LYS	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.