



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 10:32 pm BST

PDB ID : 5GTM
Title : Modified human MxA, nucleotide-free form
Authors : Chen, Y.; Gao, S.
Deposited on : 2016-08-22
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

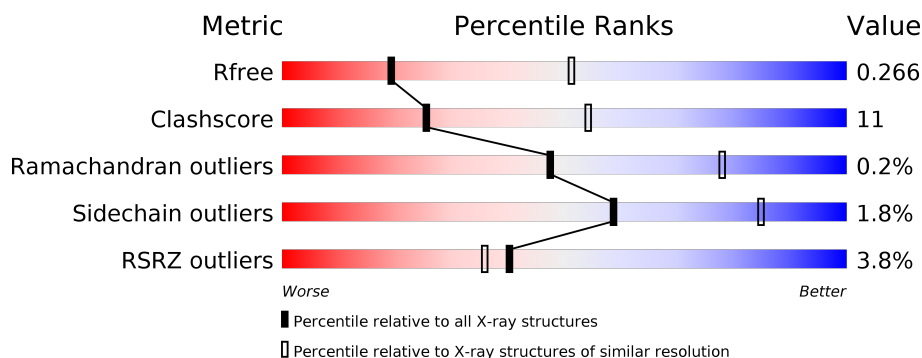
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on 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	601	<div> <div>2%</div> <div> <div></div> <div>64%</div> <div>23%</div> <div>13%</div> </div> </div>
1	B	601	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>20%</div> <div>13%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8454 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 Interferon-induced GTP-binding protein Mx1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	520	Total	C	N	O	S	0	0	0
			4229	2677	728	810	14			
1	B	522	Total	C	N	O	S	0	0	0
			4222	2666	730	812	14			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	376	SER	ILE	engineered mutation	UNP P20591
A	379	ILE	VAL	variant	UNP P20591
A	440	ALA	TYR	engineered mutation	UNP P20591
A	441	ALA	ARG	engineered mutation	UNP P20591
A	442	ALA	GLY	engineered mutation	UNP P20591
A	443	ALA	ARG	engineered mutation	UNP P20591
A	527	ASP	MET	engineered mutation	UNP P20591
A	?	-	CYS	deletion	UNP P20591
A	?	-	GLN	deletion	UNP P20591
A	?	-	ASP	deletion	UNP P20591
A	?	-	GLN	deletion	UNP P20591
A	?	-	VAL	deletion	UNP P20591
A	?	-	TYR	deletion	UNP P20591
A	?	-	ARG	deletion	UNP P20591
A	?	-	GLY	deletion	UNP P20591
A	?	-	ALA	deletion	UNP P20591
A	?	-	LEU	deletion	UNP P20591
A	?	-	GLN	deletion	UNP P20591
A	?	-	LYS	deletion	UNP P20591
A	?	-	VAL	deletion	UNP P20591
A	?	-	ARG	deletion	UNP P20591
A	?	-	GLU	deletion	UNP P20591
A	?	-	LYS	deletion	UNP P20591
A	?	-	GLU	deletion	UNP P20591
A	?	-	LEU	deletion	UNP P20591

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLU	deletion	UNP P20591
A	?	-	GLU	deletion	UNP P20591
A	?	-	GLU	deletion	UNP P20591
A	?	-	LYS	deletion	UNP P20591
A	?	-	LYS	deletion	UNP P20591
A	?	-	LYS	deletion	UNP P20591
A	?	-	LYS	deletion	UNP P20591
A	?	-	SER	deletion	UNP P20591
A	?	-	TRP	deletion	UNP P20591
A	?	-	ASP	deletion	UNP P20591
A	?	-	PHE	deletion	UNP P20591
A	614	SER	LYS	engineered mutation	UNP P20591
A	617	SER	LEU	engineered mutation	UNP P20591
A	620	SER	LEU	engineered mutation	UNP P20591
B	376	SER	ILE	engineered mutation	UNP P20591
B	379	ILE	VAL	variant	UNP P20591
B	440	ALA	TYR	engineered mutation	UNP P20591
B	441	ALA	ARG	engineered mutation	UNP P20591
B	442	ALA	GLY	engineered mutation	UNP P20591
B	443	ALA	ARG	engineered mutation	UNP P20591
B	527	ASP	MET	engineered mutation	UNP P20591
B	?	-	CYS	deletion	UNP P20591
B	?	-	GLN	deletion	UNP P20591
B	?	-	ASP	deletion	UNP P20591
B	?	-	GLN	deletion	UNP P20591
B	?	-	VAL	deletion	UNP P20591
B	?	-	TYR	deletion	UNP P20591
B	?	-	ARG	deletion	UNP P20591
B	?	-	GLY	deletion	UNP P20591
B	?	-	ALA	deletion	UNP P20591
B	?	-	LEU	deletion	UNP P20591
B	?	-	GLN	deletion	UNP P20591
B	?	-	LYS	deletion	UNP P20591
B	?	-	VAL	deletion	UNP P20591
B	?	-	ARG	deletion	UNP P20591
B	?	-	GLU	deletion	UNP P20591
B	?	-	LYS	deletion	UNP P20591
B	?	-	GLU	deletion	UNP P20591
B	?	-	LEU	deletion	UNP P20591
B	?	-	GLU	deletion	UNP P20591
B	?	-	GLU	deletion	UNP P20591
B	?	-	GLU	deletion	UNP P20591

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	LYS	deletion	UNP P20591
B	?	-	LYS	deletion	UNP P20591
B	?	-	LYS	deletion	UNP P20591
B	?	-	LYS	deletion	UNP P20591
B	?	-	SER	deletion	UNP P20591
B	?	-	TRP	deletion	UNP P20591
B	?	-	ASP	deletion	UNP P20591
B	?	-	PHE	deletion	UNP P20591
B	614	SER	LYS	engineered mutation	UNP P20591
B	617	SER	LEU	engineered mutation	UNP P20591
B	620	SER	LEU	engineered mutation	UNP P20591

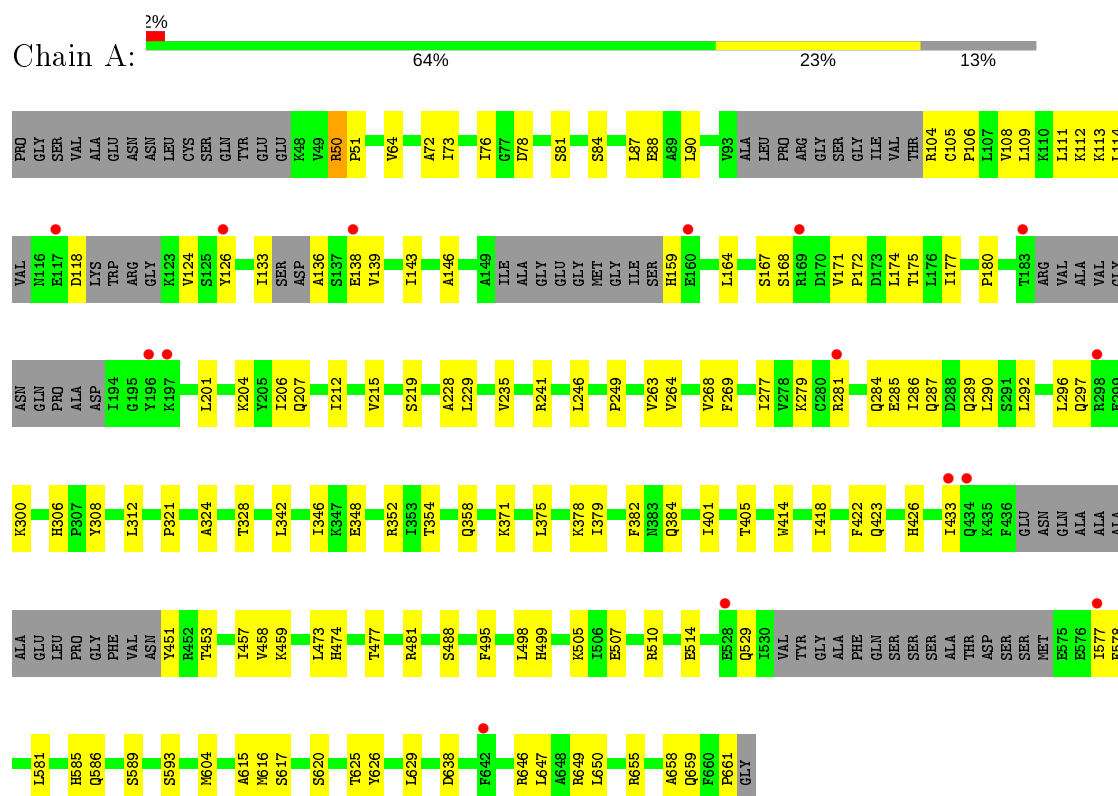
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O 1 1	0	0
2	B	2	Total O 2 2	0	0

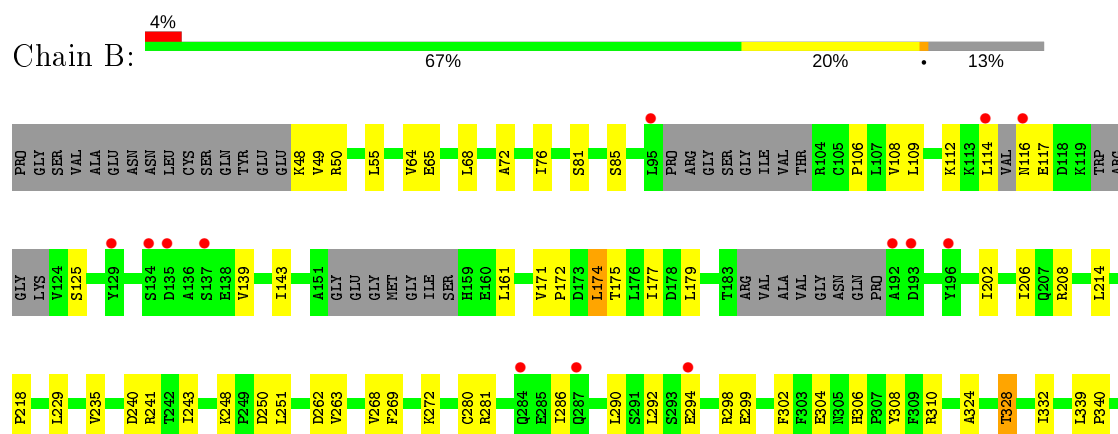
3 Residue-property plots [i](#)

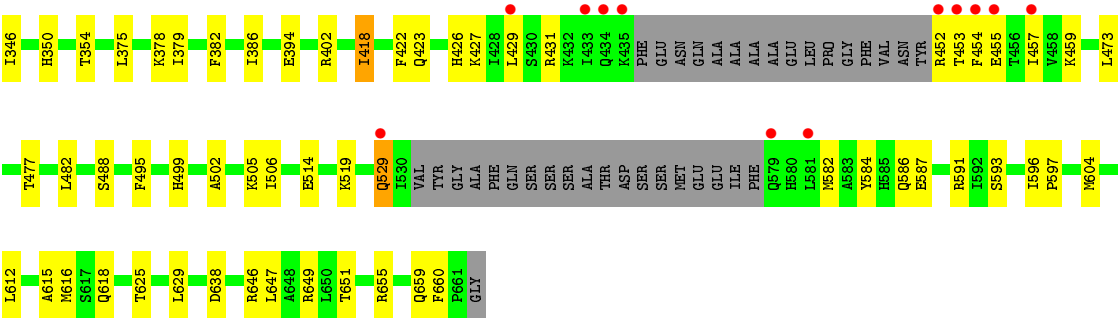
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Interferon-induced GTP-binding protein Mx1



- Molecule 1: Interferon-induced GTP-binding protein Mx1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	104.18Å 60.16Å 153.92Å 90.00° 98.62° 90.00°	Depositor
Resolution (Å)	39.20 – 2.90 39.20 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (39.20-2.90) 98.8 (39.20-2.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.204 , 0.263 0.209 , 0.266	Depositor DCC
R_{free} test set	2070 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	56.0	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 72.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8454	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.94% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.49	0/4286	0.66	0/5759
1	B	0.49	0/4277	0.68	1/5750 (0.0%)
All	All	0.49	0/8563	0.67	1/11509 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	174	LEU	CA-CB-CG	5.43	127.80	115.30

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	4229	0	4291	102	0
1	B	4222	0	4292	78	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
All	All	8454	0	8583	180	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 11.

All (180) 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:50:ARG:H	1:A:51:PRO:HD2	1.33	0.90
1:A:114:LEU:HD13	1:A:118:ASP:HA	1.60	0.82
1:B:174:LEU:HB2	1:B:328:THR:HG22	1.60	0.82
1:A:477:THR:HG22	1:A:604:MET:HG3	1.64	0.80
1:A:105:CYS:HG	1:A:159:HIS:N	1.79	0.80
1:A:422:PHE:HA	1:A:585:HIS:HE1	1.48	0.79
1:B:268:VAL:HG23	1:B:269:PHE:H	1.49	0.78
1:A:78:ASP:HB3	1:A:81:SER:HB3	1.66	0.77
1:B:206:ILE:HG22	1:B:235:VAL:HG11	1.65	0.77
1:B:625:THR:HG22	1:B:629:LEU:HG	1.67	0.77
1:A:249:PRO:HB3	1:A:277:ILE:HD11	1.66	0.76
1:A:174:LEU:HB2	1:A:328:THR:HG22	1.71	0.72
1:A:263:VAL:HG22	1:A:268:VAL:HG21	1.73	0.69
1:A:219:SER:HB3	1:A:246:LEU:HB3	1.74	0.69
1:A:268:VAL:HG23	1:A:269:PHE:H	1.58	0.68
1:B:646:ARG:HG2	1:B:649:ARG:HH21	1.58	0.68
1:A:281:ARG:HE	1:A:292:LEU:HD23	1.59	0.68
1:A:422:PHE:HA	1:A:585:HIS:CE1	2.28	0.67
1:A:206:ILE:HG22	1:A:235:VAL:HG11	1.77	0.66
1:A:281:ARG:HH21	1:A:292:LEU:HA	1.62	0.65
1:B:263:VAL:HA	1:B:268:VAL:HG22	1.79	0.64
1:A:50:ARG:N	1:A:51:PRO:HD2	2.10	0.64
1:B:281:ARG:NH2	1:B:290:LEU:O	2.21	0.64
1:B:68:LEU:HD21	1:B:346:ILE:HG12	1.80	0.63
1:B:477:THR:HG22	1:B:604:MET:HG3	1.79	0.63
1:A:458:VAL:HG13	1:A:585:HIS:HD2	1.64	0.62
1:B:114:LEU:HD22	1:B:117:GLU:HB3	1.82	0.61
1:B:250:ASP:HB2	1:B:251:LEU:HD12	1.82	0.61
1:A:505:LYS:HD3	1:A:615:ALA:HB1	1.82	0.60
1:B:263:VAL:HG22	1:B:268:VAL:HG21	1.83	0.60
1:B:473:LEU:O	1:B:477:THR:HG23	2.03	0.59
1:A:382:PHE:CE2	1:A:616:MET:HG3	2.38	0.58
1:A:422:PHE:HD1	1:A:585:HIS:HD1	1.51	0.58
1:A:281:ARG:NH2	1:A:290:LEU:O	2.37	0.58
1:B:324:ALA:O	1:B:328:THR:HG23	2.04	0.58
1:B:81:SER:HA	1:B:218:PRO:HD3	1.85	0.57
1:A:297:GLN:HA	1:A:300:LYS:HE3	1.87	0.57
1:A:378:LYS:HE3	1:A:495:PHE:CE1	2.40	0.57
1:A:105:CYS:SG	1:A:159:HIS:N	2.79	0.56
1:A:625:THR:HG22	1:A:629:LEU:HG	1.85	0.56
1:B:72:ALA:HB1	1:B:177:ILE:HD13	1.87	0.56
1:A:473:LEU:O	1:A:477:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:379:ILE:HD13	1:B:616:MET:O	2.05	0.56
1:B:429:LEU:HD23	1:B:582:MET:SD	2.46	0.56
1:A:124:VAL:HG12	1:A:164:LEU:HD13	1.88	0.56
1:B:477:THR:HG21	1:B:514:GLU:HG3	1.88	0.55
1:A:139:VAL:O	1:A:143:ILE:HG13	2.07	0.54
1:A:263:VAL:HA	1:A:268:VAL:HG22	1.88	0.54
1:A:375:LEU:O	1:A:379:ILE:HG13	2.07	0.54
1:B:55:LEU:HD12	1:B:660:PHE:CG	2.42	0.54
1:A:481:ARG:HD3	1:A:510:ARG:HD3	1.90	0.54
1:A:50:ARG:HD2	1:A:112:LYS:HD2	1.89	0.54
1:B:529:GLN:O	1:B:584:TYR:OH	2.24	0.53
1:A:458:VAL:HG13	1:A:585:HIS:CD2	2.43	0.53
1:A:133:ILE:HG21	1:A:136:ALA:N	2.23	0.53
1:A:133:ILE:HD11	1:A:138:GLU:OE1	2.10	0.53
1:B:502:ALA:O	1:B:506:ILE:HG13	2.09	0.53
1:A:109:LEU:O	1:A:175:THR:HA	2.08	0.52
1:B:354:THR:HG22	1:B:647:LEU:HD21	1.90	0.52
1:A:286:ILE:O	1:A:289:GLN:HG2	2.11	0.50
1:B:505:LYS:HD2	1:B:615:ALA:HB1	1.93	0.50
1:A:284:GLN:O	1:A:287:GLN:HG2	2.11	0.50
1:B:306:HIS:CE1	1:B:308:TYR:HB2	2.47	0.50
1:B:49:VAL:HG13	1:B:50:ARG:H	1.76	0.49
1:B:382:PHE:O	1:B:386:ILE:HG12	2.13	0.49
1:B:339:LEU:HB2	1:B:340:PRO:HD3	1.95	0.48
1:B:646:ARG:HG2	1:B:649:ARG:NH2	2.25	0.48
1:A:50:ARG:CD	1:A:112:LYS:HD2	2.42	0.48
1:A:84:SER:O	1:A:88:GLU:HG3	2.13	0.48
1:B:378:LYS:HE3	1:B:495:PHE:CE1	2.48	0.48
1:A:126:TYR:HB3	1:A:146:ALA:HB1	1.96	0.48
1:B:179:LEU:HD13	1:B:202:ILE:HG12	1.96	0.48
1:A:159:HIS:O	1:A:159:HIS:ND1	2.45	0.48
1:A:201:LEU:O	1:A:204:LYS:HB3	2.14	0.48
1:B:109:LEU:O	1:B:175:THR:HA	2.13	0.48
1:B:229:LEU:HA	1:B:229:LEU:HD23	1.67	0.48
1:B:418:ILE:HG12	1:B:593:SER:HB2	1.96	0.48
1:A:324:ALA:O	1:A:328:THR:HG23	2.14	0.47
1:B:429:LEU:HD11	1:B:454:PHE:HB3	1.96	0.47
1:B:235:VAL:CG1	1:B:241:ARG:HH12	2.27	0.47
1:A:620:SER:O	1:A:626:TYR:OH	2.27	0.47
1:A:458:VAL:HG21	1:A:581:LEU:HB3	1.96	0.47
1:A:474:HIS:ND1	1:A:514:GLU:OE2	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:ALA:HB1	1:A:177:ILE:HD13	1.95	0.47
1:B:281:ARG:HD3	1:B:286:ILE:HG12	1.96	0.46
1:A:306:HIS:HE1	1:A:308:TYR:HB2	1.79	0.46
1:B:459:LYS:HE2	1:B:529:GLN:HB2	1.96	0.46
1:B:179:LEU:HD13	1:B:202:ILE:HA	1.97	0.46
1:B:268:VAL:HG23	1:B:269:PHE:N	2.24	0.46
1:B:375:LEU:O	1:B:379:ILE:HG13	2.14	0.46
1:A:481:ARG:HD2	1:A:507:GLU:OE2	2.15	0.46
1:B:427:LYS:HE3	1:B:431:ARG:HD2	1.98	0.46
1:A:50:ARG:H	1:A:51:PRO:CD	2.18	0.46
1:A:249:PRO:HD2	1:A:279:LYS:HD2	1.98	0.46
1:B:214:LEU:HD12	1:B:243:ILE:HB	1.98	0.46
1:B:304:GLU:HA	1:B:310:ARG:HD3	1.98	0.46
1:A:354:THR:O	1:A:358:GLN:HG2	2.16	0.45
1:B:85:SER:OG	1:B:280:CYS:HB3	2.17	0.45
1:A:306:HIS:CE1	1:A:308:TYR:HB2	2.51	0.45
1:B:112:LYS:HE3	1:B:112:LYS:HB3	1.85	0.45
1:A:401:ILE:HD12	1:A:405:THR:HG22	1.98	0.45
1:B:240:ASP:OD1	1:B:272:LYS:HG2	2.17	0.45
1:A:235:VAL:CG1	1:A:241:ARG:HH22	2.29	0.45
1:B:453:THR:O	1:B:457:ILE:HD13	2.17	0.45
1:B:587:GLU:OE2	1:B:591:ARG:NH1	2.38	0.45
1:A:646:ARG:HG2	1:A:649:ARG:HH21	1.82	0.45
1:A:354:THR:HG22	1:A:647:LEU:HD21	1.99	0.45
1:A:112:LYS:O	1:A:167:SER:HA	2.17	0.44
1:A:646:ARG:HA	1:A:649:ARG:HH21	1.82	0.44
1:B:394:GLU:O	1:B:402:ARG:NE	2.50	0.44
1:B:423:GLN:HA	1:B:426:HIS:HB3	1.99	0.44
1:A:229:LEU:HA	1:A:229:LEU:HD23	1.75	0.44
1:A:76:ILE:HD11	1:A:206:ILE:CD1	2.47	0.44
1:A:113:LYS:HE3	1:A:171:VAL:O	2.18	0.44
1:A:477:THR:HG21	1:A:514:GLU:HG3	2.00	0.44
1:B:106:PRO:HB2	1:B:161:LEU:HD23	1.99	0.44
1:B:422:PHE:CE1	1:B:586:GLN:HG3	2.52	0.44
1:A:285:GLU:OE1	1:A:285:GLU:HA	2.18	0.44
1:A:215:VAL:HG13	1:A:228:ALA:HB1	1.99	0.44
1:A:112:LYS:HD3	1:A:112:LYS:HA	1.67	0.43
1:A:646:ARG:HG2	1:A:649:ARG:NH2	2.33	0.43
1:B:76:ILE:HD11	1:B:206:ILE:HD11	1.99	0.43
1:A:658:ALA:O	1:A:661:PRO:HD2	2.19	0.43
1:A:73:ILE:HD13	1:A:212:ILE:HB	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:ARG:HH21	1:B:292:LEU:HD22	1.83	0.43
1:B:647:LEU:O	1:B:651:THR:HG23	2.19	0.43
1:A:87:LEU:HA	1:A:90:LEU:HD12	2.00	0.43
1:A:589:SER:O	1:A:593:SER:N	2.40	0.43
1:B:294:GLU:O	1:B:298:ARG:N	2.49	0.43
1:A:108:VAL:HG22	1:A:177:ILE:HG13	2.01	0.42
1:A:296:LEU:HD11	1:A:321:PRO:HG3	2.01	0.42
1:A:453:THR:O	1:A:457:ILE:HG12	2.18	0.42
1:A:459:LYS:NZ	1:A:529:GLN:HG3	2.34	0.42
1:A:581:LEU:HA	1:A:581:LEU:HD23	1.75	0.42
1:B:427:LYS:O	1:B:431:ARG:HB2	2.20	0.42
1:A:109:LEU:HD22	1:A:111:LEU:HD21	1.99	0.42
1:A:104:ARG:O	1:A:180:PRO:HG3	2.20	0.42
1:A:207:GLN:HA	1:A:235:VAL:CG1	2.50	0.42
1:A:414:TRP:CZ2	1:A:418:ILE:HD11	2.54	0.42
1:B:612:LEU:O	1:B:616:MET:HG2	2.19	0.42
1:A:114:LEU:HD12	1:A:168:SER:HA	2.01	0.42
1:A:433:ILE:HD12	1:A:578:PHE:HD1	1.84	0.42
1:B:76:ILE:HD11	1:B:206:ILE:CD1	2.50	0.42
1:A:264:VAL:HG11	1:A:312:LEU:HD13	2.02	0.42
1:A:498:LEU:HD12	1:A:498:LEU:HA	1.90	0.42
1:A:268:VAL:HG23	1:A:269:PHE:N	2.31	0.42
1:A:379:ILE:HG21	1:A:617:SER:HA	2.01	0.41
1:B:139:VAL:O	1:B:143:ILE:HG13	2.20	0.41
1:A:342:LEU:O	1:A:346:ILE:HG13	2.20	0.41
1:B:65:GLU:HG2	1:B:208:ARG:HH12	1.85	0.41
1:B:350:HIS:NE2	1:B:651:THR:HG22	2.36	0.41
1:A:655:ARG:O	1:A:659:GLN:HG3	2.20	0.41
1:A:159:HIS:CD2	1:A:201:LEU:HD13	2.55	0.41
1:A:650:LEU:HD23	1:A:650:LEU:HA	1.85	0.41
1:B:55:LEU:HD12	1:B:660:PHE:CD1	2.55	0.41
1:B:596:ILE:HB	1:B:597:PRO:HD3	2.03	0.41
1:B:125:SER:O	1:B:125:SER:OG	2.38	0.41
1:B:50:ARG:HA	1:B:50:ARG:HD2	1.77	0.41
1:A:171:VAL:HG13	1:A:172:PRO:HD2	2.03	0.41
1:A:235:VAL:HG12	1:A:241:ARG:HH22	1.86	0.41
1:A:379:ILE:HD13	1:A:616:MET:O	2.20	0.41
1:B:171:VAL:HG13	1:B:172:PRO:HD2	2.03	0.41
1:B:49:VAL:HG23	1:B:332:ILE:HD11	2.03	0.41
1:A:207:GLN:HG2	1:A:235:VAL:HG22	2.03	0.41
1:A:50:ARG:N	1:A:51:PRO:CD	2.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:655:ARG:HE	1:B:659:GLN:NE2	2.19	0.41
1:A:418:ILE:CG2	1:A:589:SER:HB3	2.51	0.41
1:A:423:GLN:O	1:A:426:HIS:HB3	2.21	0.41
1:B:299:GLU:O	1:B:302:PHE:N	2.53	0.41
1:B:48:LYS:HG2	1:B:49:VAL:N	2.36	0.41
1:B:306:HIS:HE1	1:B:308:TYR:HB2	1.85	0.41
1:B:452:ARG:HA	1:B:455:GLU:HB3	2.03	0.40
1:B:625:THR:HG22	1:B:625:THR:O	2.20	0.40
1:A:488:SER:OG	1:A:499:HIS:ND1	2.44	0.40
1:A:50:ARG:NE	1:A:112:LYS:HZ3	2.20	0.40
1:A:76:ILE:HD11	1:A:206:ILE:HD11	2.03	0.40
1:B:218:PRO:HB3	1:B:248:LYS:HD2	2.04	0.40
1:B:108:VAL:HG22	1:B:177:ILE:HG13	2.04	0.40
1:A:348:GLU:O	1:A:352:ARG:HG2	2.21	0.40
1:B:488:SER:OG	1:B:499:HIS:ND1	2.45	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	502/601 (84%)	485 (97%)	15 (3%)	2 (0%)	34	66
1	B	506/601 (84%)	488 (96%)	18 (4%)	0	100	100
All	All	1008/1202 (84%)	973 (96%)	33 (3%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	ARG
1	A	106	PRO

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	475/534 (89%)	468 (98%)	7 (2%)	65	87
1	B	474/534 (89%)	464 (98%)	10 (2%)	53	81
All	All	949/1068 (89%)	932 (98%)	17 (2%)	59	85

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	VAL
1	A	371	LYS
1	A	384	GLN
1	A	451	TYR
1	A	577	ILE
1	A	586	GLN
1	A	638	ASP
1	B	64	VAL
1	B	116	ASN
1	B	262	ASP
1	B	328	THR
1	B	418	ILE
1	B	482	LEU
1	B	519	LYS
1	B	529	GLN
1	B	618	GLN
1	B	638	ASP

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

Mol	Chain	Res	Type
1	B	426	HIS
1	B	659	GLN

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 carbohydrates 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	520/601 (86%)	-0.01	15 (2%)	51 47	18, 89, 164, 204	0
1	B	522/601 (86%)	0.08	25 (4%)	30 27	22, 77, 154, 195	0
All	All	1042/1202 (86%)	0.04	40 (3%)	40 36	18, 83, 158, 204	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	192	ALA	4.8
1	B	134	SER	4.7
1	A	577	ILE	3.5
1	B	454	PHE	3.5
1	B	529	GLN	3.4
1	B	429	LEU	3.3
1	B	193	ASP	3.1
1	A	196	TYR	3.1
1	B	435	LYS	3.0
1	A	281	ARG	3.0
1	B	581	LEU	3.0
1	B	116	ASN	2.9
1	A	434	GLN	2.8
1	B	452	ARG	2.7
1	B	95	LEU	2.7
1	B	137	SER	2.5
1	B	294	GLU	2.5
1	A	117	GLU	2.4
1	A	298	ARG	2.4
1	A	433	ILE	2.4
1	B	135	ASP	2.4
1	B	579	GLN	2.4
1	B	453	THR	2.3
1	A	169	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	126	TYR	2.3
1	B	114	LEU	2.3
1	B	129	TYR	2.3
1	B	433	ILE	2.3
1	B	287	GLN	2.2
1	B	434	GLN	2.2
1	A	183	THR	2.2
1	A	138	GLU	2.2
1	A	160	GLU	2.2
1	B	455	GLU	2.2
1	A	197	LYS	2.1
1	B	284	GLN	2.1
1	B	457	ILE	2.1
1	A	528	GLU	2.0
1	A	642	PHE	2.0
1	B	196	TYR	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 carbohydrates 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.