



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 10:31 PM GMT

PDB ID : 1TW9
Title : Glutathione Transferase-2, apo form, from the nematode Heligmosomoides polygyrus
Authors : Schuller, D.J.; Liu, Q.; Kriksunov, I.A.; Campbell, A.M.; Barrett, J.; Brophy, P.M.; Hao, Q.
Deposited on : 2004-06-30
Resolution : 1.71 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

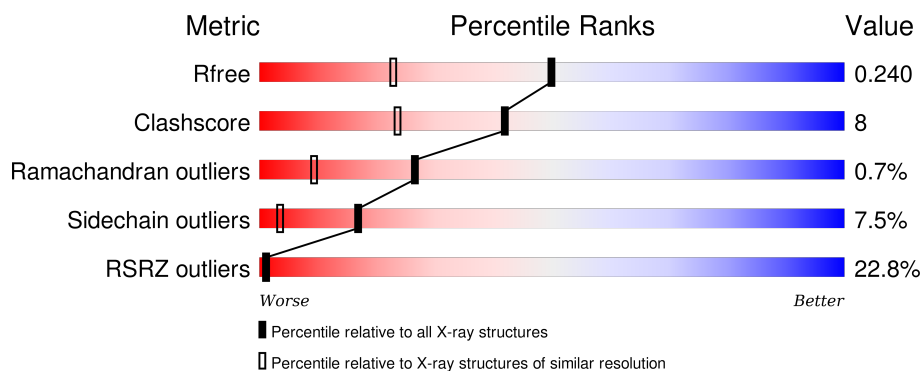
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.71 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}	91344	3998 (1.74-1.70)
Clashscore	102246	4425 (1.74-1.70)
Ramachandran outliers	100387	4360 (1.74-1.70)
Sidechain outliers	100360	4360 (1.74-1.70)
RSRZ outliers	91569	4010 (1.74-1.70)

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. 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	206	<div> <div>18%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>• 7%</div> </div> </div>
1	B	206	<div> <div>18%</div> <div> <div></div> <div>74%</div> <div>13%</div> <div>• • 9%</div> </div> </div>
1	C	206	<div> <div>17%</div> <div> <div></div> <div>80%</div> <div>13%</div> <div>• 6%</div> </div> </div>
1	D	206	<div> <div>20%</div> <div> <div></div> <div>77%</div> <div>15%</div> <div>• • 6%</div> </div> </div>
1	E	206	<div> <div>24%</div> <div> <div></div> <div>73%</div> <div>17%</div> <div>• • 6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	206	<div><div></div><div>30%</div><div></div><div>77%</div><div></div><div>18%</div><div></div><div></div></div>
1	G	206	<div><div></div><div>17%</div><div></div><div>75%</div><div></div><div>15%</div><div></div><div>8%</div></div>
1	H	206	<div><div></div><div>27%</div><div></div><div>71%</div><div></div><div>16%</div><div></div><div>5%</div><div>7%</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14254 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 glutathione S-transferase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1531	989	256	280	6			
1	B	188	Total	C	N	O	S	0	0	0
			1507	974	252	275	6			
1	C	194	Total	C	N	O	S	0	0	0
			1563	1013	260	284	6			
1	D	194	Total	C	N	O	S	0	0	0
			1559	1008	260	285	6			
1	E	193	Total	C	N	O	S	0	0	0
			1542	997	257	282	6			
1	F	205	Total	C	N	O	S	0	0	0
			1645	1066	271	301	7			
1	G	189	Total	C	N	O	S	0	0	0
			1516	980	254	276	6			
1	H	191	Total	C	N	O	S	0	0	0
			1531	989	256	280	6			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	CLONING ARTIFACT	UNP Q9NJK6
A	2	VAL	-	CLONING ARTIFACT	UNP Q9NJK6
A	21	VAL	ILE	CONFLICT	UNP Q9NJK6
A	40	VAL	ALA	CONFLICT	UNP Q9NJK6
A	91	LEU	PRO	CONFLICT	UNP Q9NJK6
A	108	THR	MET	CONFLICT	UNP Q9NJK6
A	120	PRO	LEU	CONFLICT	UNP Q9NJK6
A	140	LEU	PRO	CONFLICT	UNP Q9NJK6
B	1	MET	-	CLONING ARTIFACT	UNP Q9NJK6
B	2	VAL	-	CLONING ARTIFACT	UNP Q9NJK6
B	21	VAL	ILE	CONFLICT	UNP Q9NJK6
B	40	VAL	ALA	CONFLICT	UNP Q9NJK6
B	91	LEU	PRO	CONFLICT	UNP Q9NJK6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	108	THR	MET	CONFLICT	UNP Q9NJJQ6
B	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6
B	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6
C	1	MET	-	CLONING ARTIFACT	UNP Q9NJJQ6
C	2	VAL	-	CLONING ARTIFACT	UNP Q9NJJQ6
C	21	VAL	ILE	CONFLICT	UNP Q9NJJQ6
C	40	VAL	ALA	CONFLICT	UNP Q9NJJQ6
C	91	LEU	PRO	CONFLICT	UNP Q9NJJQ6
C	108	THR	MET	CONFLICT	UNP Q9NJJQ6
C	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6
C	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6
D	1	MET	-	CLONING ARTIFACT	UNP Q9NJJQ6
D	2	VAL	-	CLONING ARTIFACT	UNP Q9NJJQ6
D	21	VAL	ILE	CONFLICT	UNP Q9NJJQ6
D	40	VAL	ALA	CONFLICT	UNP Q9NJJQ6
D	91	LEU	PRO	CONFLICT	UNP Q9NJJQ6
D	108	THR	MET	CONFLICT	UNP Q9NJJQ6
D	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6
D	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6
E	1	MET	-	CLONING ARTIFACT	UNP Q9NJJQ6
E	2	VAL	-	CLONING ARTIFACT	UNP Q9NJJQ6
E	21	VAL	ILE	CONFLICT	UNP Q9NJJQ6
E	40	VAL	ALA	CONFLICT	UNP Q9NJJQ6
E	91	LEU	PRO	CONFLICT	UNP Q9NJJQ6
E	108	THR	MET	CONFLICT	UNP Q9NJJQ6
E	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6
E	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6
F	1	MET	-	CLONING ARTIFACT	UNP Q9NJJQ6
F	2	VAL	-	CLONING ARTIFACT	UNP Q9NJJQ6
F	21	VAL	ILE	CONFLICT	UNP Q9NJJQ6
F	40	VAL	ALA	CONFLICT	UNP Q9NJJQ6
F	91	LEU	PRO	CONFLICT	UNP Q9NJJQ6
F	108	THR	MET	CONFLICT	UNP Q9NJJQ6
F	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6
F	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6
G	1	MET	-	CLONING ARTIFACT	UNP Q9NJJQ6
G	2	VAL	-	CLONING ARTIFACT	UNP Q9NJJQ6
G	21	VAL	ILE	CONFLICT	UNP Q9NJJQ6
G	40	VAL	ALA	CONFLICT	UNP Q9NJJQ6
G	91	LEU	PRO	CONFLICT	UNP Q9NJJQ6
G	108	THR	MET	CONFLICT	UNP Q9NJJQ6
G	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6
H	1	MET	-	CLONING ARTIFACT	UNP Q9NJJQ6
H	2	VAL	-	CLONING ARTIFACT	UNP Q9NJJQ6
H	21	VAL	ILE	CONFLICT	UNP Q9NJJQ6
H	40	VAL	ALA	CONFLICT	UNP Q9NJJQ6
H	91	LEU	PRO	CONFLICT	UNP Q9NJJQ6
H	108	THR	MET	CONFLICT	UNP Q9NJJQ6
H	120	PRO	LEU	CONFLICT	UNP Q9NJJQ6
H	140	LEU	PRO	CONFLICT	UNP Q9NJJQ6

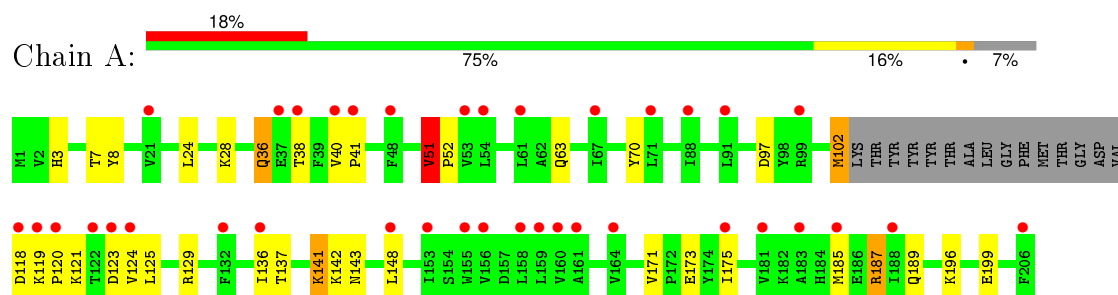
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	384	Total O 384 384	0	0
2	B	220	Total O 220 220	0	0
2	C	285	Total O 285 285	0	0
2	D	251	Total O 251 251	0	0
2	E	149	Total O 149 149	0	0
2	F	186	Total O 186 186	0	0
2	G	216	Total O 216 216	0	0
2	H	169	Total O 169 169	0	0

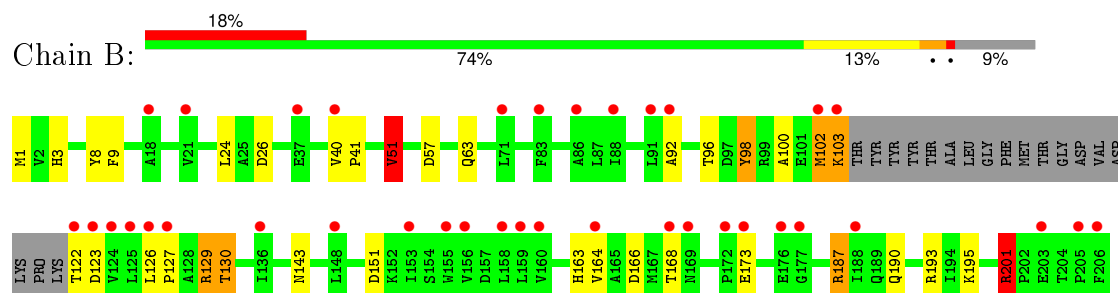
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

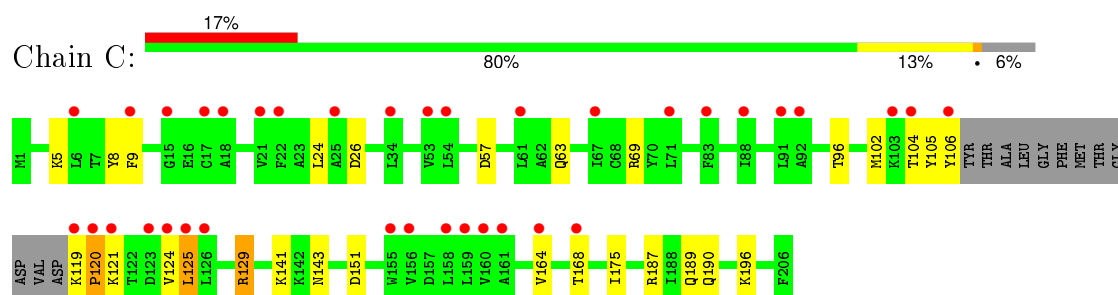
- Molecule 1: glutathione S-transferase 2



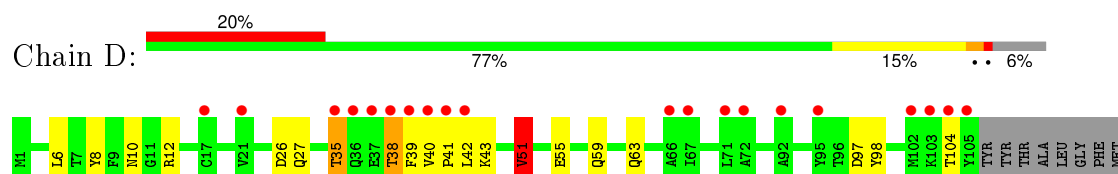
- Molecule 1: glutathione S-transferase 2

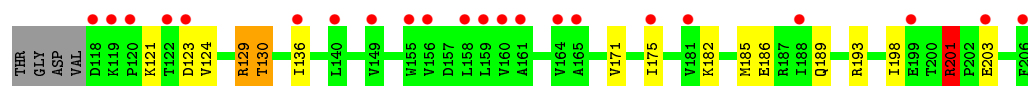


- Molecule 1: glutathione S-transferase 2

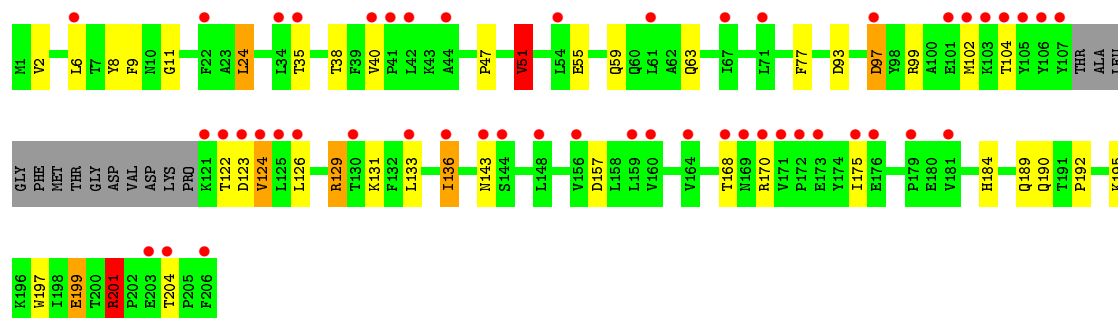


- Molecule 1: glutathione S-transferase 2

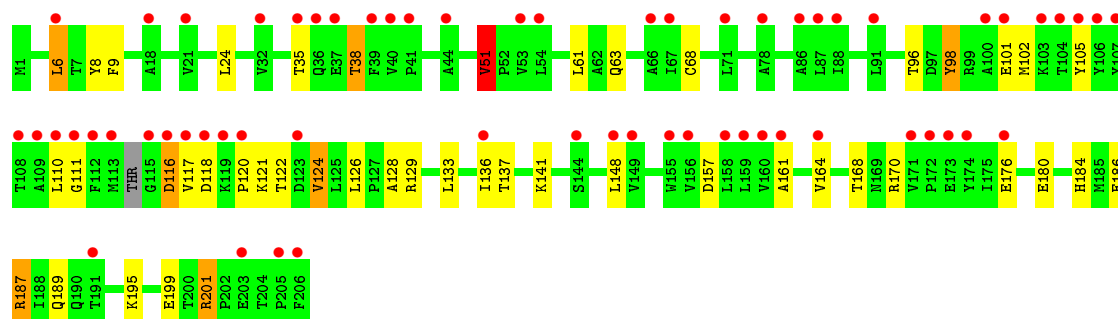
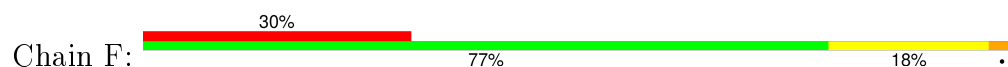




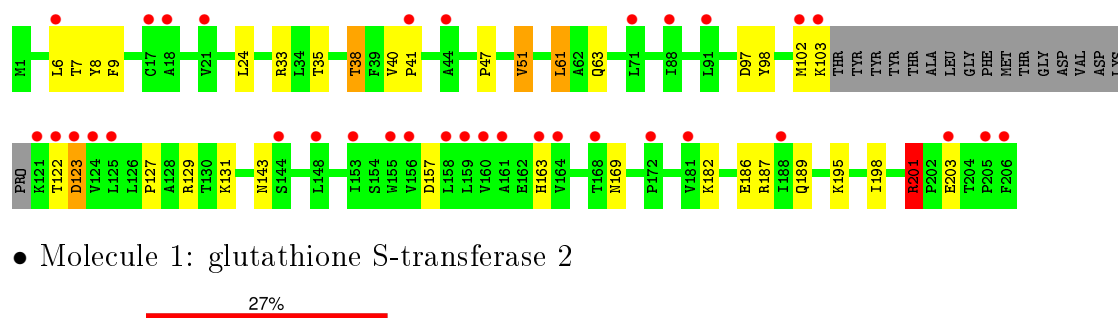
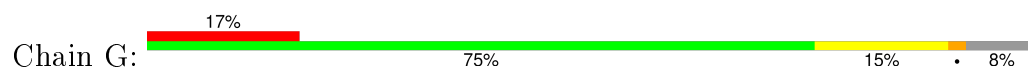
- Molecule 1: glutathione S-transferase 2



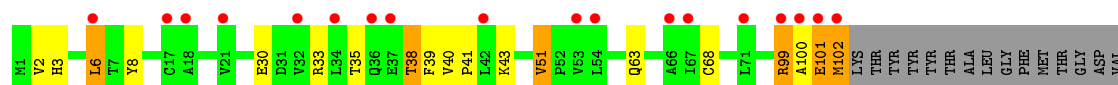
- Molecule 1: glutathione S-transferase 2

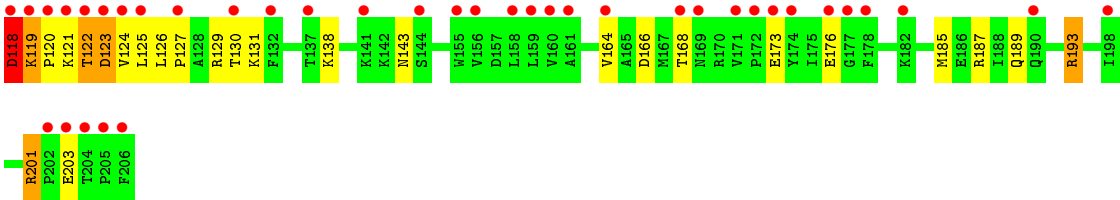


- Molecule 1: glutathione S-transferase 2



- Molecule 1: glutathione S-transferase 2





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	72.68Å 74.03Å 88.57Å 79.09° 80.08° 81.55°	Depositor
Resolution (Å)	25.00 – 1.71 24.26 – 1.71	Depositor EDS
% Data completeness (in resolution range)	95.2 (25.00-1.71) 91.0 (24.26-1.71)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 1.71Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.180 , 0.232 0.189 , 0.240	Depositor DCC
R_{free} test set	4565 reflections (2.55%)	DCC
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.9	EDS
Estimated twinning fraction	0.009 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 184047 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14254	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	1.05	2/1568 (0.1%)	1.11	6/2120 (0.3%)
1	B	0.88	0/1543	1.00	8/2086 (0.4%)
1	C	0.96	0/1602	1.01	6/2166 (0.3%)
1	D	0.95	0/1597	0.99	5/2159 (0.2%)
1	E	0.94	2/1579 (0.1%)	0.98	8/2136 (0.4%)
1	F	0.97	2/1686 (0.1%)	1.04	6/2280 (0.3%)
1	G	0.88	0/1552	1.06	8/2097 (0.4%)
1	H	0.87	1/1568 (0.1%)	1.02	7/2120 (0.3%)
All	All	0.94	7/12695 (0.1%)	1.03	54/17164 (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

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	199	GLU	CD-OE2	17.71	1.45	1.25
1	E	199	GLU	CD-OE1	8.35	1.34	1.25
1	H	68	CYS	CB-SG	-5.76	1.72	1.81
1	A	70	TYR	CD2-CE2	-5.55	1.31	1.39
1	A	70	TYR	CD1-CE1	-5.39	1.31	1.39

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	201	ARG	NE-CZ-NH1	14.00	127.30	120.30
1	C	187	ARG	NE-CZ-NH1	13.65	127.13	120.30
1	B	201	ARG	NE-CZ-NH1	11.89	126.24	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	187	ARG	NE-CZ-NH1	11.70	126.15	120.30
1	F	201	ARG	NE-CZ-NH1	11.46	126.03	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	102	MET	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	1531	0	1527	26	0
1	B	1507	0	1503	27	0
1	C	1563	0	1561	20	0
1	D	1559	0	1556	28	0
1	E	1542	0	1527	29	0
1	F	1645	0	1634	36	1
1	G	1516	0	1516	16	0
1	H	1531	0	1527	27	0
2	A	384	0	0	12	1
2	B	220	0	0	10	0
2	C	285	0	0	8	0
2	D	251	0	0	8	1
2	E	149	0	0	11	1
2	F	186	0	0	8	0
2	G	216	0	0	8	0
2	H	169	0	0	4	0
All	All	14254	0	12351	207	2

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.

The worst 5 of 207 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:G:195:LYS:NZ	2:G:268:HOH:O	1.73	1.18
1:E:2:VAL:HB	2:E:1781:HOH:O	0.96	1.12
1:C:106:TYR:C	2:C:1685:HOH:O	1.88	1.10
1:B:96:THR:HG22	2:B:337:HOH:O	1.52	1.08
1:E:143:ASN:CG	2:E:1376:HOH:O	1.95	1.03

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:240:HOH:O	2:E:1566:HOH:O[1_444]	1.76	0.44
1:F:180:GLU:OE1	2:D:367:HOH:O[1_665]	2.07	0.13

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	187/206 (91%)	184 (98%)	2 (1%)	1 (0%)	34	15
1	B	184/206 (89%)	180 (98%)	3 (2%)	1 (0%)	34	15
1	C	190/206 (92%)	186 (98%)	2 (1%)	2 (1%)	17	4
1	D	190/206 (92%)	187 (98%)	2 (1%)	1 (0%)	34	15
1	E	189/206 (92%)	185 (98%)	3 (2%)	1 (0%)	34	15
1	F	201/206 (98%)	197 (98%)	2 (1%)	2 (1%)	19	5
1	G	185/206 (90%)	181 (98%)	3 (2%)	1 (0%)	34	15
1	H	187/206 (91%)	180 (96%)	5 (3%)	2 (1%)	17	4
All	All	1513/1648 (92%)	1480 (98%)	22 (2%)	11 (1%)	26	9

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	111	GLY
1	H	63	GLN
1	B	63	GLN
1	C	63	GLN
1	E	63	GLN

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	162/174 (93%)	151 (93%)	11 (7%)	20	4
1	B	159/174 (91%)	147 (92%)	12 (8%)	17	3
1	C	165/174 (95%)	157 (95%)	8 (5%)	31	11
1	D	165/174 (95%)	153 (93%)	12 (7%)	17	4
1	E	161/174 (92%)	149 (92%)	12 (8%)	17	3
1	F	173/174 (99%)	161 (93%)	12 (7%)	19	4
1	G	160/174 (92%)	145 (91%)	15 (9%)	11	2
1	H	162/174 (93%)	146 (90%)	16 (10%)	10	2
All	All	1307/1392 (94%)	1209 (92%)	98 (8%)	17	3

5 of 98 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	24	LEU
1	E	201	ARG
1	H	122	THR
1	E	51	VAL
1	E	131	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	27	GLN

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Mol	Chain	Res	Type
1	D	190	GLN
1	H	27	GLN
1	D	63	GLN
1	E	63	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	191/206 (92%)	1.14	37 (19%) 1 1	29, 33, 37, 41	0
1	B	188/206 (91%)	1.22	37 (19%) 1 1	28, 33, 39, 42	0
1	C	194/206 (94%)	1.16	36 (18%) 2 2	29, 33, 38, 40	0
1	D	194/206 (94%)	1.25	42 (21%) 1 1	29, 33, 39, 43	0
1	E	193/206 (93%)	1.47	49 (25%) 1 1	29, 34, 39, 41	0
1	F	205/206 (99%)	1.44	61 (29%) 1 1	29, 33, 39, 43	0
1	G	189/206 (91%)	1.23	34 (17%) 2 2	29, 33, 38, 41	0
1	H	191/206 (92%)	1.45	56 (29%) 1 1	29, 34, 38, 41	0
All	All	1545/1648 (93%)	1.30	352 (22%) 1 1	28, 33, 39, 43	0

The worst 5 of 352 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	107	TYR	12.1
1	C	120	PRO	9.9
1	B	124	VAL	8.0
1	F	112	PHE	7.8
1	D	206	PHE	7.5

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.