



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 19, 2016 – 09:36 PM GMT

PDB ID : 5A9V
Title : Structure of apo BipA
Authors : Kumar, V.; Chen, Y.; Ero, R.; Li, Z.; Gao, Y.
Deposited on : 2015-07-23
Resolution : 3.31 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

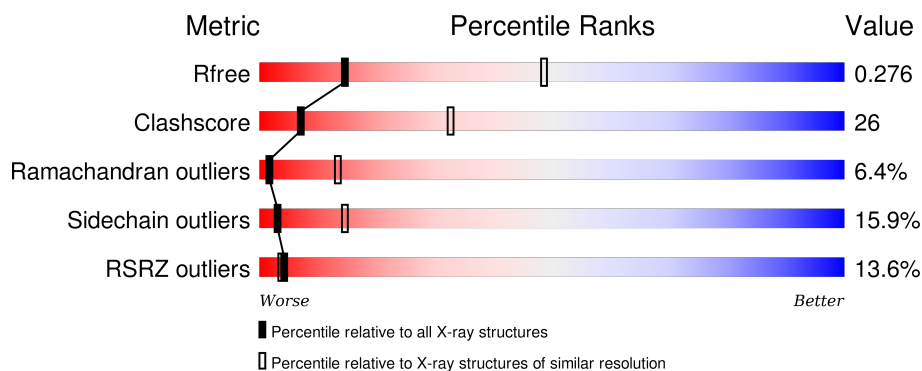
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.31 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	1198 (3.40-3.24)
Clashscore	102246	1280 (3.40-3.24)
Ramachandran outliers	100387	1260 (3.40-3.24)
Sidechain outliers	100360	1259 (3.40-3.24)
RSRZ outliers	91569	1203 (3.40-3.24)

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	607	<div> <div>12%</div> <div>59%</div> <div>25%</div> <div>5%</div> <div>9%</div> </div>
1	B	607	<div> <div>13%</div> <div>58%</div> <div>26%</div> <div>6%</div> <div>9%</div> </div>
1	C	607	<div> <div>14%</div> <div>59%</div> <div>25%</div> <div>6%</div> <div>9%</div> </div>
1	D	607	<div> <div>12%</div> <div>58%</div> <div>25%</div> <div>6%</div> <div>9%</div> </div>
1	E	607	<div> <div>13%</div> <div>59%</div> <div>25%</div> <div>5%</div> <div>9%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	607	<div><div></div><div>12%</div><div>58%</div><div>25%</div><div>6%</div><div>•</div><div>9%</div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 25398 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.

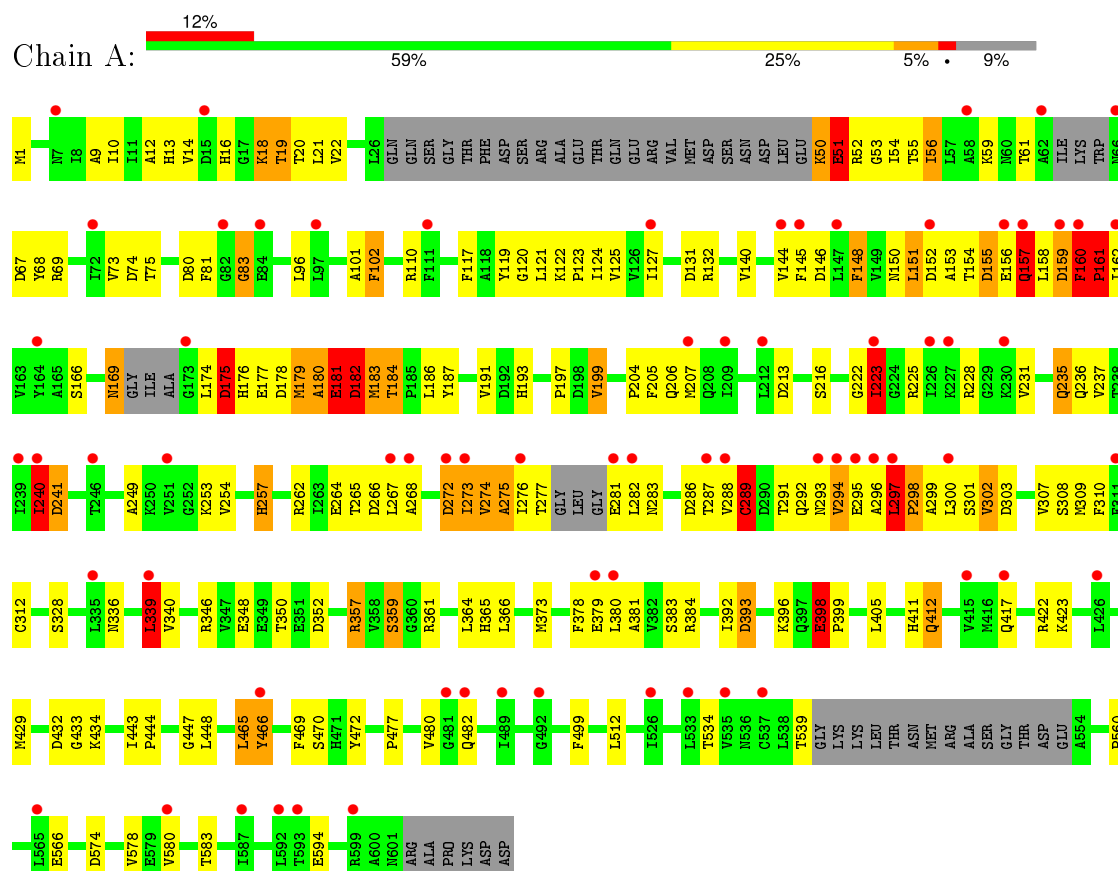
- Molecule 1 is a protein called GTP-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4233	2675	736	806	16			
1	B	555	Total	C	N	O	S	0	0	0
			4233	2675	736	806	16			
1	C	555	Total	C	N	O	S	0	0	0
			4233	2675	736	806	16			
1	D	555	Total	C	N	O	S	0	0	0
			4233	2675	736	806	16			
1	E	555	Total	C	N	O	S	0	0	0
			4233	2675	736	806	16			
1	F	555	Total	C	N	O	S	0	0	0
			4233	2675	736	806	16			

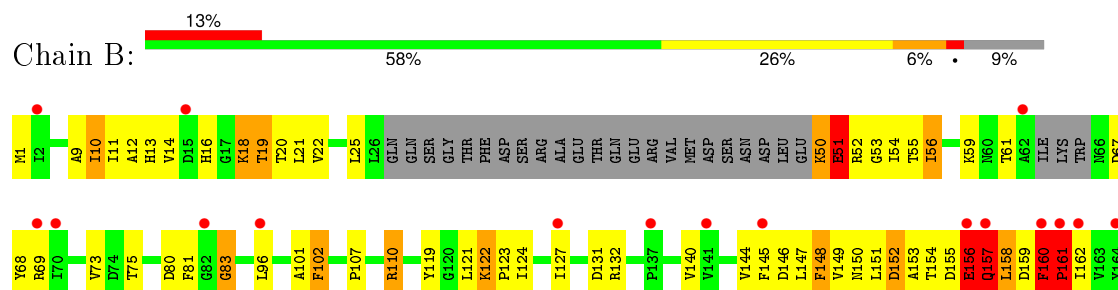
3 Residue-property plots

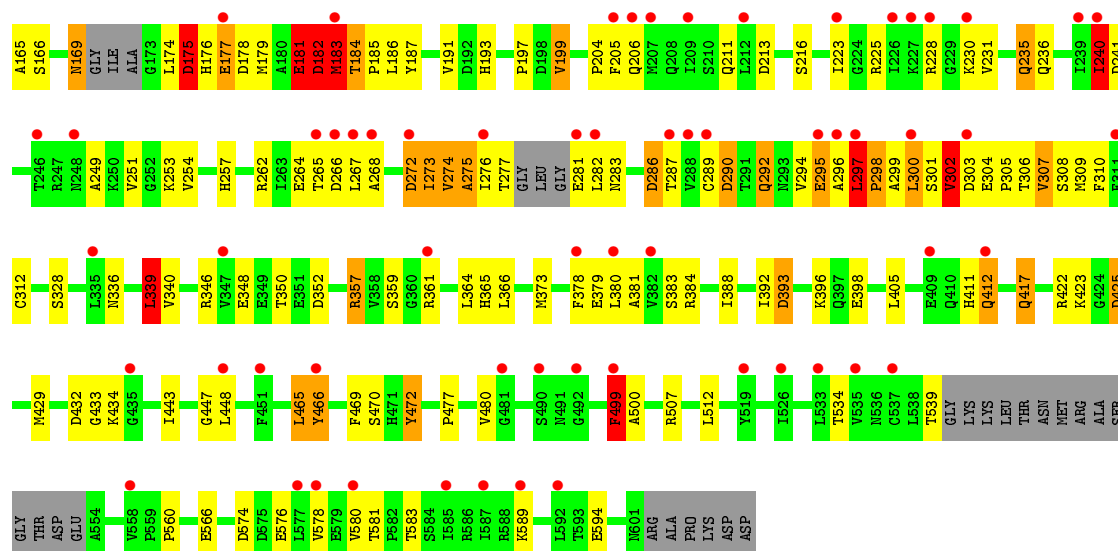
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: GTP-BINDING PROTEIN

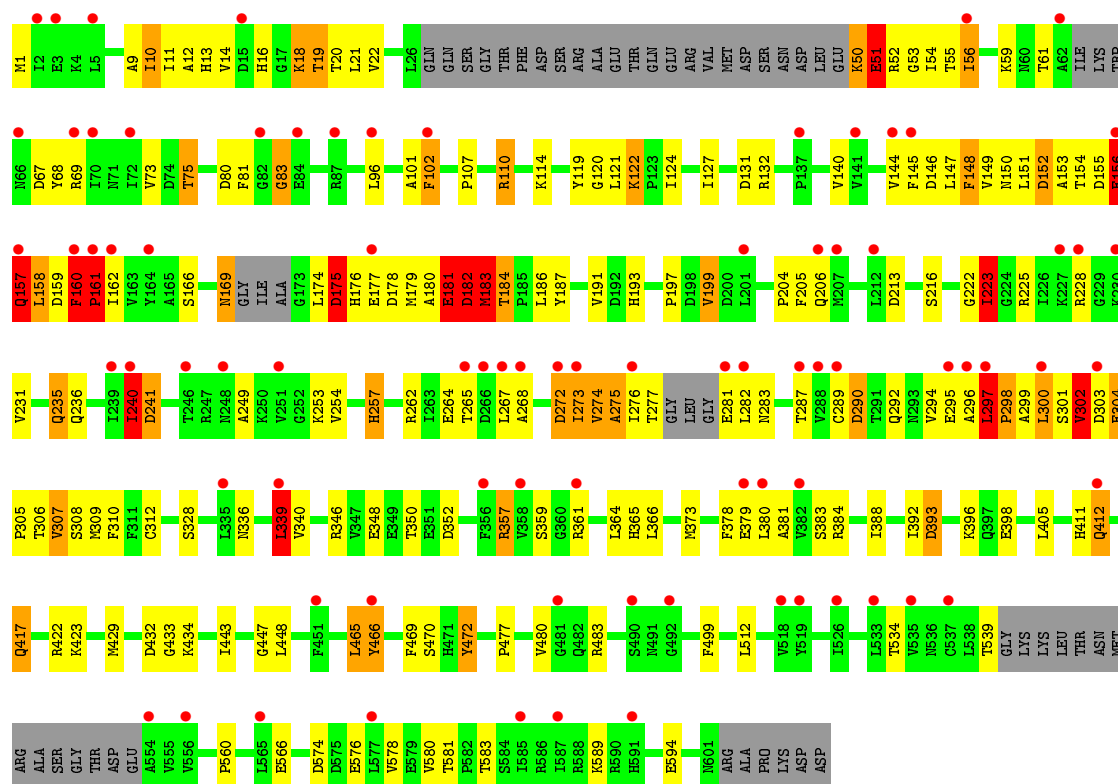


• Molecule 1: GTP-BINDING PROTEIN

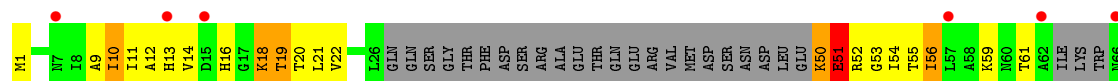


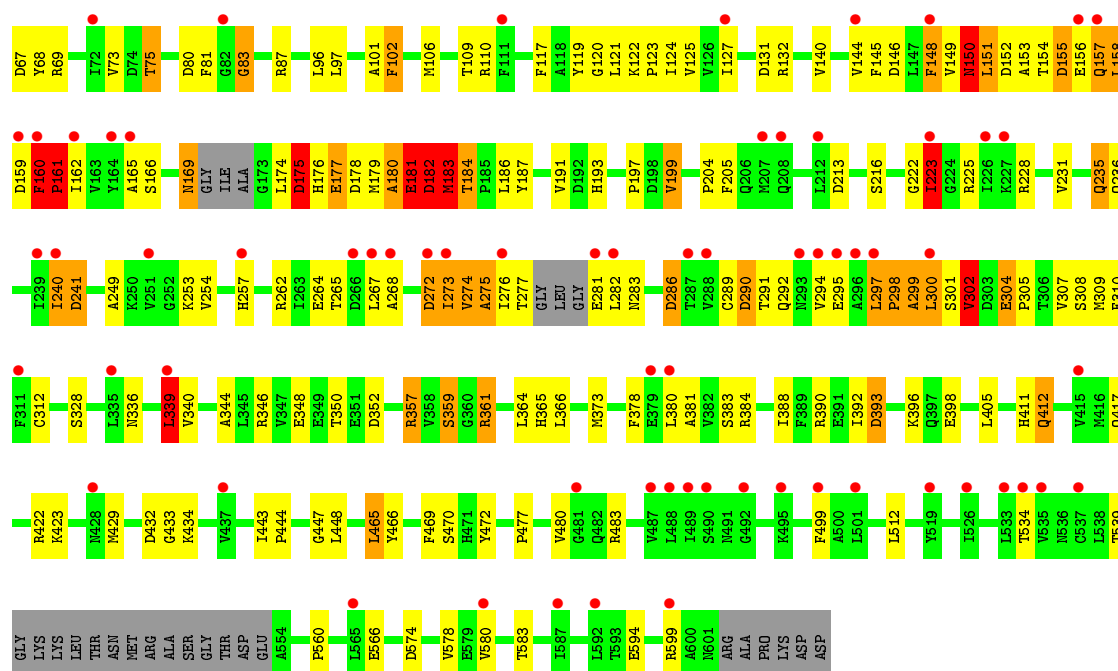


• Molecule 1: GTP-BINDING PROTEIN

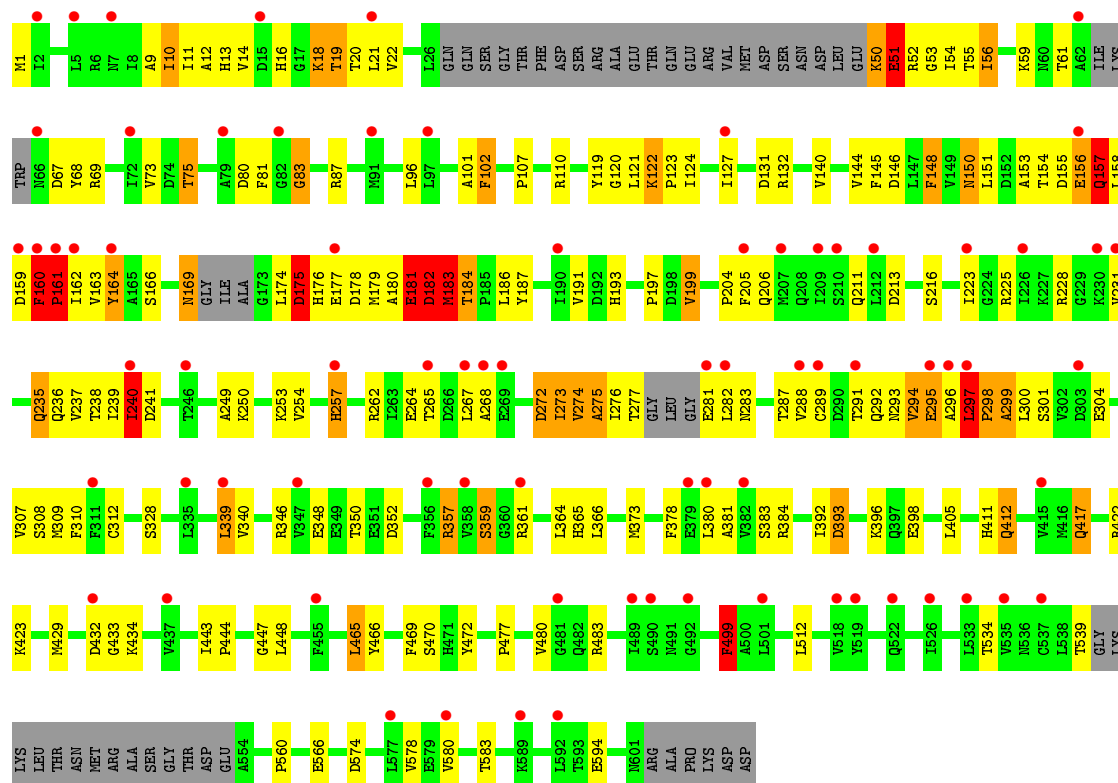


• Molecule 1: GTP-BINDING PROTEIN





• Molecule 1: GTP-BINDING PROTEIN



• Molecule 1: GTP-BINDING PROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	241.95Å 241.95Å 241.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.39 – 3.31 49.39 – 3.31	Depositor EDS
% Data completeness (in resolution range)	91.8 (49.39-3.31) 91.9 (49.39-3.31)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.263 , 0.288 0.280 , 0.276	Depositor DCC
R_{free} test set	9472 reflections (11.14%)	DCC
Wilson B-factor (Å ²)	114.3	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 132.6	EDS
Estimated twinning fraction	0.429 for l,-k,h 0.429 for -l,-k,-h 0.429 for -h,-l,-k 0.429 for -h,l,k 0.439 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	120 of 94684 reflections (0.127%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	25398	wwPDB-VP
Average B, all atoms (Å ²)	135.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.61	1/4292 (0.0%)	0.91	6/5786 (0.1%)
1	B	0.60	0/4292	0.91	8/5786 (0.1%)
1	C	0.59	0/4292	0.89	6/5786 (0.1%)
1	D	0.59	0/4292	0.89	6/5786 (0.1%)
1	E	0.58	0/4292	0.89	5/5786 (0.1%)
1	F	0.59	1/4292 (0.0%)	0.93	9/5786 (0.2%)
All	All	0.59	2/25752 (0.0%)	0.90	40/34716 (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	A	0	4
1	B	0	4
1	C	0	4
1	D	0	5
1	E	0	4
1	F	0	4
All	All	0	25

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	398	GLU	CD-OE2	-8.25	1.16	1.25
1	F	348	GLU	CD-OE2	-5.46	1.19	1.25

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	15	ASP	CB-CG-OD2	14.00	130.90	118.30
1	A	266	ASP	CB-CG-OD2	11.71	128.84	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	15	ASP	CB-CG-OD1	-11.48	107.97	118.30
1	A	398	GLU	OE1-CD-OE2	-8.39	113.23	123.30
1	F	563	MET	CA-CB-CG	-8.02	99.66	113.30

There are no chirality outliers.

5 of 25 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	ARG	Peptide
1	A	19	THR	Peptide
1	A	240	ILE	Peptide
1	A	465	LEU	Peptide
1	B	19	THR	Peptide

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	4233	0	4156	217	0
1	B	4233	0	4156	238	0
1	C	4233	0	4156	240	0
1	D	4233	0	4155	213	0
1	E	4233	0	4156	218	0
1	F	4233	0	4156	221	0
All	All	25398	0	24935	1329	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 26.

The worst 5 of 1329 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:287:THR:HB	1:A:297:LEU:CD1	1.13	1.55
1:B:169:ASN:HD21	1:B:182:ASP:N	1.09	1.48
1:A:287:THR:CB	1:A:297:LEU:CD1	1.94	1.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:THR:CB	1:C:297:LEU:HD11	1.53	1.39
1:D:169:ASN:OD1	1:D:182:ASP:CB	1.70	1.38

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	543/607 (90%)	421 (78%)	85 (16%)	37 (7%)	1	12
1	B	543/607 (90%)	420 (77%)	89 (16%)	34 (6%)	2	14
1	C	543/607 (90%)	418 (77%)	89 (16%)	36 (7%)	1	13
1	D	543/607 (90%)	419 (77%)	87 (16%)	37 (7%)	1	12
1	E	543/607 (90%)	425 (78%)	86 (16%)	32 (6%)	2	16
1	F	543/607 (90%)	424 (78%)	85 (16%)	34 (6%)	2	14
All	All	3258/3642 (90%)	2527 (78%)	521 (16%)	210 (6%)	2	13

5 of 210 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	161	PRO
1	A	181	GLU
1	A	182	ASP
1	A	223	ILE
1	A	393	ASP

5.3.2 Protein sidechains [i](#)

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	450/519 (87%)	379 (84%)	71 (16%)	3	15
1	B	450/519 (87%)	376 (84%)	74 (16%)	3	13
1	C	450/519 (87%)	381 (85%)	69 (15%)	3	16
1	D	450/519 (87%)	382 (85%)	68 (15%)	3	17
1	E	450/519 (87%)	381 (85%)	69 (15%)	3	16
1	F	450/519 (87%)	373 (83%)	77 (17%)	2	12
All	All	2700/3114 (87%)	2272 (84%)	428 (16%)	3	14

5 of 428 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	361	ARG
1	D	216	SER
1	F	297	LEU
1	C	398	GLU
1	D	50	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	482	GLN
1	D	169	ASN
1	F	176	HIS
1	D	157	GLN
1	D	193	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 [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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	555/607 (91%)	0.84	71 (12%) 5 4	79, 130, 184, 219	0
1	B	555/607 (91%)	0.84	79 (14%) 4 3	80, 129, 184, 221	0
1	C	555/607 (91%)	0.88	82 (14%) 3 3	82, 130, 183, 216	0
1	D	555/607 (91%)	0.85	73 (13%) 4 4	85, 131, 183, 218	0
1	E	555/607 (91%)	0.83	76 (13%) 4 4	83, 131, 183, 225	0
1	F	555/607 (91%)	0.84	72 (12%) 5 4	83, 131, 182, 218	0
All	All	3330/3642 (91%)	0.84	453 (13%) 4 4	79, 131, 184, 225	0

The worst 5 of 453 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	160	PHE	7.3
1	F	295	GLU	6.8
1	D	297	LEU	5.5
1	C	82	GLY	5.4
1	B	295	GLU	5.3

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.