



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 09:18 AM GMT

PDB ID : 3HU1  
Title : Structure of p97 N-D1 R95G mutant in complex with ATPgS  
Authors : Tang, W.-K.  
Deposited on : 2009-06-12  
Resolution : 2.81 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](mailto: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.

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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

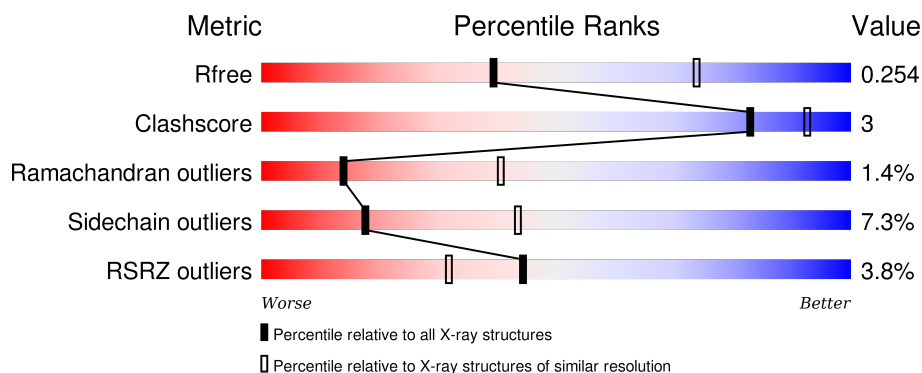
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.81 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.



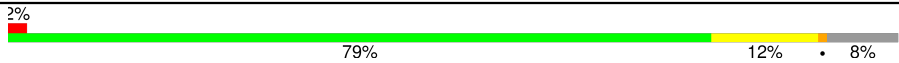
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	2676 (2.84-2.80)
Clashscore	102246	3124 (2.84-2.80)
Ramachandran outliers	100387	3072 (2.84-2.80)
Sidechain outliers	100360	3074 (2.84-2.80)
RSRZ outliers	91569	2690 (2.84-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	489	<div> <div>4%</div> <div>80% 11% • 8%</div> </div>
1	B	489	<div> <div>2%</div> <div>78% 13% • 8%</div> </div>
1	C	489	<div> <div>2%</div> <div>79% 12% • 8%</div> </div>
1	D	489	<div> <div>6%</div> <div>80% 11% • 8%</div> </div>
1	E	489	<div> <div>6%</div> <div>79% 11% • 8%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	489	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '2%', a large green segment labeled '79%', a yellow segment labeled '12%', and a small grey segment at the end labeled '8%'. The segments are separated by thin white lines.

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21445 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 Transitional endoplasmic reticulum ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	S	0	0	0
			3522	2211	623	670	18			
1	B	451	Total	C	N	O	S	0	0	0
			3522	2211	623	670	18			
1	C	451	Total	C	N	O	S	0	0	0
			3522	2211	623	670	18			
1	D	451	Total	C	N	O	S	0	0	0
			3522	2211	623	670	18			
1	E	451	Total	C	N	O	S	0	0	0
			3522	2211	623	670	18			
1	F	451	Total	C	N	O	S	0	0	0
			3522	2211	623	670	18			

There are 54 discrepancies between the modelled and reference sequences:

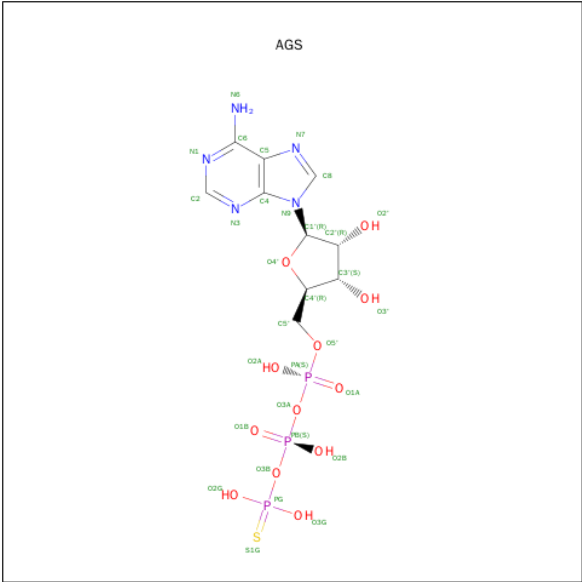
Chain	Residue	Modelled	Actual	Comment	Reference
A	95	GLY	ARG	ENGINEERED MUTATION	UNP P55072
A	482	ARG	-	EXPRESSION TAG	UNP P55072
A	483	SER	-	EXPRESSION TAG	UNP P55072
A	484	HIS	-	EXPRESSION TAG	UNP P55072
A	485	HIS	-	EXPRESSION TAG	UNP P55072
A	486	HIS	-	EXPRESSION TAG	UNP P55072
A	487	HIS	-	EXPRESSION TAG	UNP P55072
A	488	HIS	-	EXPRESSION TAG	UNP P55072
A	489	HIS	-	EXPRESSION TAG	UNP P55072
B	95	GLY	ARG	ENGINEERED MUTATION	UNP P55072
B	482	ARG	-	EXPRESSION TAG	UNP P55072
B	483	SER	-	EXPRESSION TAG	UNP P55072
B	484	HIS	-	EXPRESSION TAG	UNP P55072
B	485	HIS	-	EXPRESSION TAG	UNP P55072
B	486	HIS	-	EXPRESSION TAG	UNP P55072
B	487	HIS	-	EXPRESSION TAG	UNP P55072
B	488	HIS	-	EXPRESSION TAG	UNP P55072

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Chain	Residue	Modelled	Actual	Comment	Reference
B	489	HIS	-	EXPRESSION TAG	UNP P55072
C	95	GLY	ARG	ENGINEERED MUTATION	UNP P55072
C	482	ARG	-	EXPRESSION TAG	UNP P55072
C	483	SER	-	EXPRESSION TAG	UNP P55072
C	484	HIS	-	EXPRESSION TAG	UNP P55072
C	485	HIS	-	EXPRESSION TAG	UNP P55072
C	486	HIS	-	EXPRESSION TAG	UNP P55072
C	487	HIS	-	EXPRESSION TAG	UNP P55072
C	488	HIS	-	EXPRESSION TAG	UNP P55072
C	489	HIS	-	EXPRESSION TAG	UNP P55072
D	95	GLY	ARG	ENGINEERED MUTATION	UNP P55072
D	482	ARG	-	EXPRESSION TAG	UNP P55072
D	483	SER	-	EXPRESSION TAG	UNP P55072
D	484	HIS	-	EXPRESSION TAG	UNP P55072
D	485	HIS	-	EXPRESSION TAG	UNP P55072
D	486	HIS	-	EXPRESSION TAG	UNP P55072
D	487	HIS	-	EXPRESSION TAG	UNP P55072
D	488	HIS	-	EXPRESSION TAG	UNP P55072
D	489	HIS	-	EXPRESSION TAG	UNP P55072
E	95	GLY	ARG	ENGINEERED MUTATION	UNP P55072
E	482	ARG	-	EXPRESSION TAG	UNP P55072
E	483	SER	-	EXPRESSION TAG	UNP P55072
E	484	HIS	-	EXPRESSION TAG	UNP P55072
E	485	HIS	-	EXPRESSION TAG	UNP P55072
E	486	HIS	-	EXPRESSION TAG	UNP P55072
E	487	HIS	-	EXPRESSION TAG	UNP P55072
E	488	HIS	-	EXPRESSION TAG	UNP P55072
E	489	HIS	-	EXPRESSION TAG	UNP P55072
F	95	GLY	ARG	ENGINEERED MUTATION	UNP P55072
F	482	ARG	-	EXPRESSION TAG	UNP P55072
F	483	SER	-	EXPRESSION TAG	UNP P55072
F	484	HIS	-	EXPRESSION TAG	UNP P55072
F	485	HIS	-	EXPRESSION TAG	UNP P55072
F	486	HIS	-	EXPRESSION TAG	UNP P55072
F	487	HIS	-	EXPRESSION TAG	UNP P55072
F	488	HIS	-	EXPRESSION TAG	UNP P55072
F	489	HIS	-	EXPRESSION TAG	UNP P55072

- Molecule 2 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (three-letter code: AGS) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>12</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	C	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	D	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	E	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		
2	F	1	Total	C	N	O	P	S	0	0
			31	10	5	12	3	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

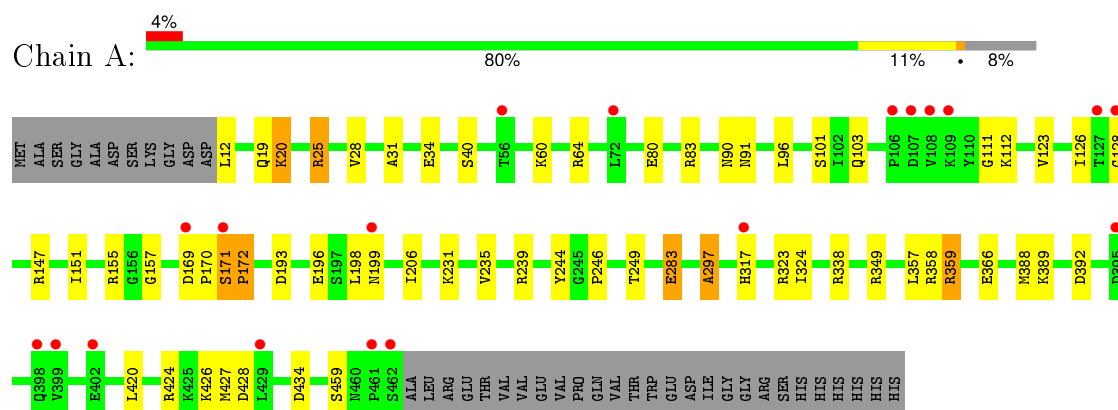
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total 27	O 27	0	0
4	B	22	Total 22	O 22	0	0
4	C	22	Total 22	O 22	0	0
4	D	16	Total 16	O 16	0	0
4	E	19	Total 19	O 19	0	0
4	F	15	Total 15	O 15	0	0

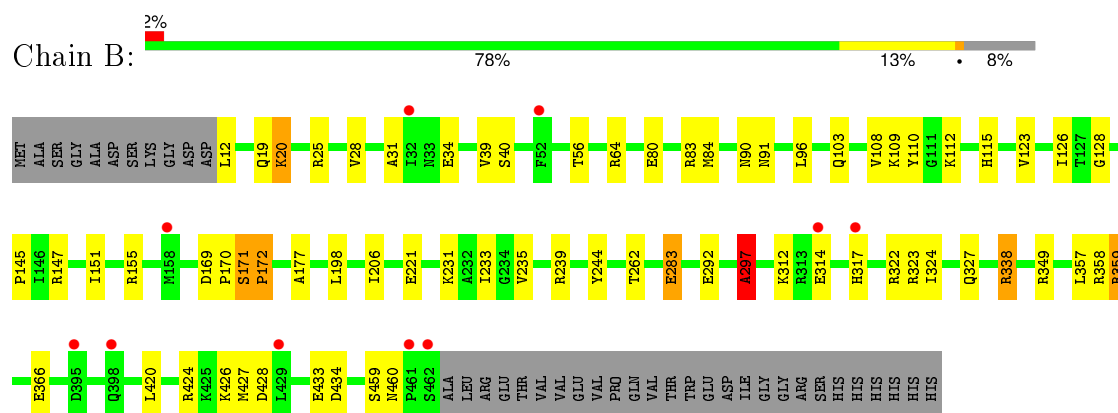
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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.

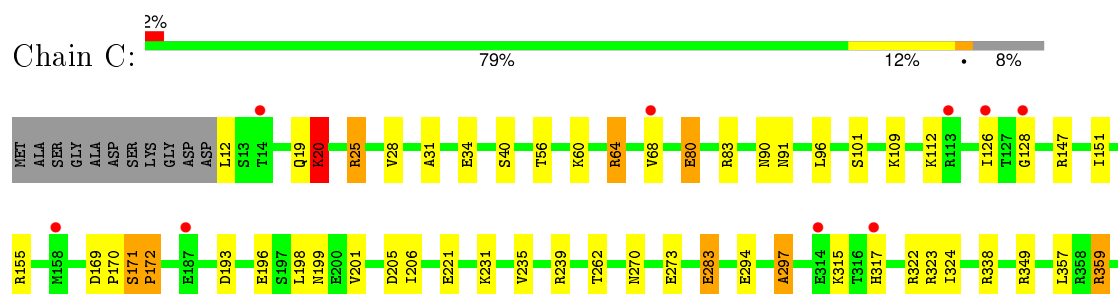
- Molecule 1: Transitional endoplasmic reticulum ATPase



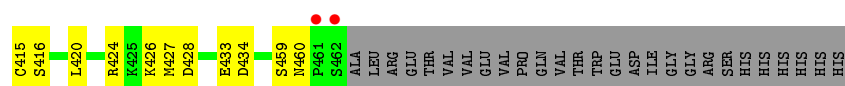
- Molecule 1: Transitional endoplasmic reticulum ATPase



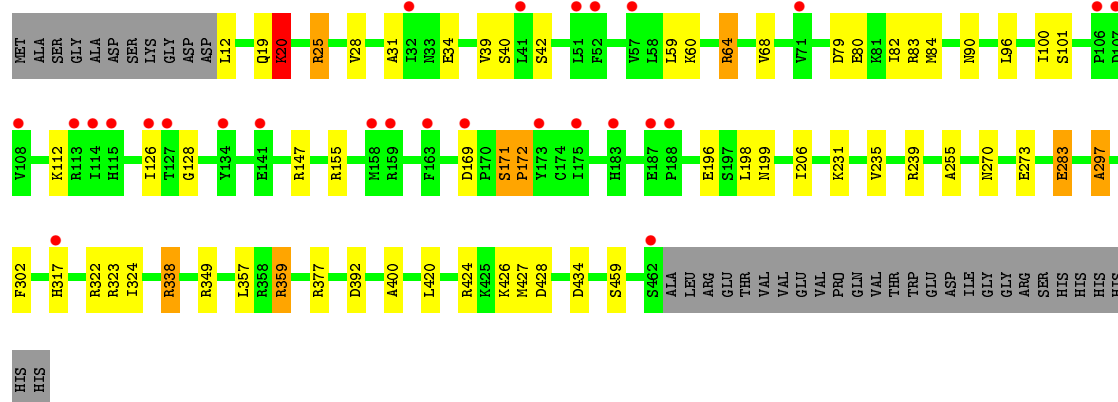
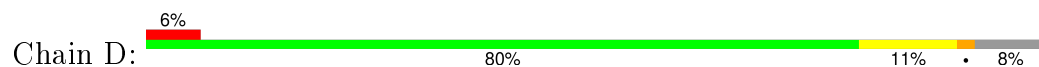
- Molecule 1: Transitional endoplasmic reticulum ATPase



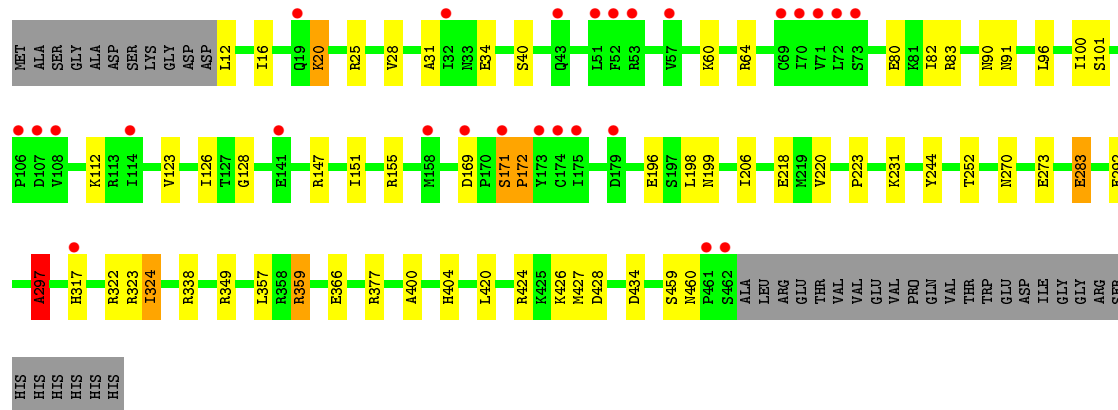
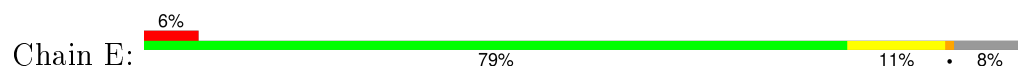




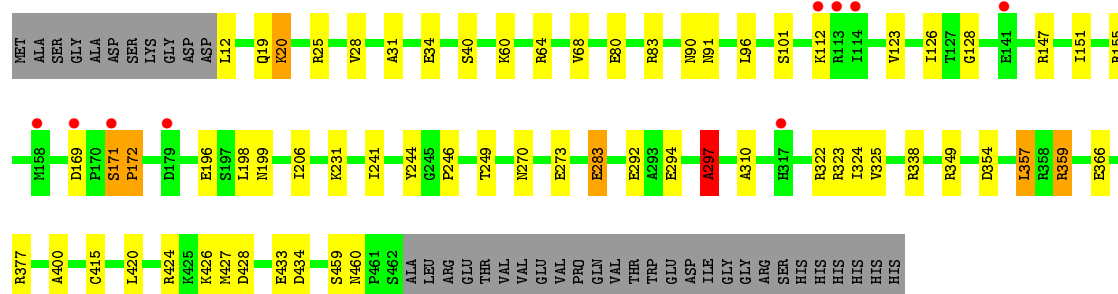
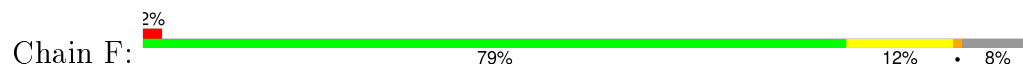
• Molecule 1: Transitional endoplasmic reticulum ATPase



• Molecule 1: Transitional endoplasmic reticulum ATPase



• Molecule 1: Transitional endoplasmic reticulum ATPase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.76Å 103.28Å 107.69Å 97.66° 91.88° 89.75°	Depositor
Resolution (Å)	25.00 – 2.81 43.92 – 2.81	Depositor EDS
% Data completeness (in resolution range)	90.2 (25.00-2.81) 87.1 (43.92-2.81)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.31 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.240 , 0.271 0.220 , 0.254	Depositor DCC
$R_{free}$ test set	4322 reflections (5.22%)	DCC
Wilson B-factor (Å <sup>2</sup> )	78.3	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 54.3	EDS
Estimated twinning fraction	0.022 for h,-k,-l 0.016 for -h,-l,-k 0.019 for -h,l,k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 87228 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	21445	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, AGS

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	0.95	0/3576	0.73	0/4832
1	B	0.97	1/3576 (0.0%)	0.73	0/4832
1	C	0.97	1/3576 (0.0%)	0.73	0/4832
1	D	0.95	0/3576	0.73	0/4832
1	E	0.99	1/3576 (0.0%)	0.74	0/4832
1	F	0.98	2/3576 (0.1%)	0.73	1/4832 (0.0%)
All	All	0.97	5/21456 (0.0%)	0.73	1/28992 (0.0%)

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	1
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	415	CYS	CB-SG	-7.06	1.70	1.82
1	F	415	CYS	CB-SG	-6.36	1.71	1.82
1	E	297	ALA	CA-CB	6.34	1.65	1.52
1	F	297	ALA	CA-CB	6.18	1.65	1.52
1	B	297	ALA	CA-CB	5.94	1.65	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	241	ILE	CG1-CB-CG2	-5.60	99.09	111.40

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	297	ALA	Peptide
1	C	297	ALA	Peptide
1	D	297	ALA	Peptide
1	E	297	ALA	Peptide
1	F	297	ALA	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	3522	0	3583	24	0
1	B	3522	0	3583	29	0
1	C	3522	0	3583	25	0
1	D	3522	0	3583	26	0
1	E	3522	0	3584	25	0
1	F	3522	0	3583	21	0
2	A	31	0	12	1	0
2	B	31	0	12	1	0
2	C	31	0	12	1	0
2	D	31	0	12	1	0
2	E	31	0	12	1	0
2	F	31	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	27	0	0	1	0
4	B	22	0	0	3	0
4	C	22	0	0	1	0
4	D	16	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	19	0	0	1	0
4	F	15	0	0	0	0
All	All	21445	0	21571	132	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 3.

The worst 5 of 132 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:F:424:ARG:HA	1:F:427:MET:HG2	1.69	0.74
1:E:424:ARG:HA	1:E:427:MET:HG2	1.73	0.71
2:B:800:AGS:S1G	1:C:359:ARG:HG2	2.31	0.71
1:D:424:ARG:HA	1:D:427:MET:HG2	1.71	0.70
2:C:800:AGS:S1G	1:D:359:ARG:HG2	2.31	0.69

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	449/489 (92%)	419 (93%)	23 (5%)	7 (2%)	12	36
1	B	449/489 (92%)	422 (94%)	21 (5%)	6 (1%)	15	42
1	C	449/489 (92%)	421 (94%)	22 (5%)	6 (1%)	15	42
1	D	449/489 (92%)	420 (94%)	23 (5%)	6 (1%)	15	42
1	E	449/489 (92%)	420 (94%)	23 (5%)	6 (1%)	15	42
1	F	449/489 (92%)	419 (93%)	24 (5%)	6 (1%)	15	42
All	All	2694/2934 (92%)	2521 (94%)	136 (5%)	37 (1%)	14	40

5 of 37 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	SER
1	B	171	SER
1	C	171	SER
1	D	171	SER
1	E	171	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	386/417 (93%)	358 (93%)	28 (7%)	17	43
1	B	386/417 (93%)	356 (92%)	30 (8%)	16	40
1	C	386/417 (93%)	355 (92%)	31 (8%)	15	39
1	D	386/417 (93%)	359 (93%)	27 (7%)	19	46
1	E	386/417 (93%)	361 (94%)	25 (6%)	21	50
1	F	386/417 (93%)	358 (93%)	28 (7%)	17	43
All	All	2316/2502 (93%)	2147 (93%)	169 (7%)	17	43

5 of 169 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	294	GLU
1	D	64	ARG
1	F	292	GLU
1	C	324	ILE
1	C	426	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	327	GLN
1	D	260	ASN

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Mol	Chain	Res	Type
1	F	285	ASN
1	D	91	ASN
1	D	285	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	AGS	A	800	3	24,33,33	1.50	2 (8%)	28,52,52	2.06	5 (17%)
2	AGS	B	800	3	24,33,33	1.64	3 (12%)	28,52,52	2.02	4 (14%)
2	AGS	C	800	3	24,33,33	1.61	4 (16%)	28,52,52	1.98	5 (17%)
2	AGS	D	800	3	24,33,33	1.53	2 (8%)	28,52,52	2.00	4 (14%)
2	AGS	E	800	3	24,33,33	1.43	3 (12%)	28,52,52	1.91	4 (14%)
2	AGS	F	800	3	24,33,33	1.62	2 (8%)	28,52,52	1.90	5 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical

component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AGS	A	800	3	-	0/15/38/38	0/3/3/3
2	AGS	B	800	3	-	0/15/38/38	0/3/3/3
2	AGS	C	800	3	-	0/15/38/38	0/3/3/3
2	AGS	D	800	3	-	0/15/38/38	0/3/3/3
2	AGS	E	800	3	-	0/15/38/38	0/3/3/3
2	AGS	F	800	3	-	0/15/38/38	0/3/3/3

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	800	AGS	C5-N7	-2.24	1.31	1.39
2	B	800	AGS	C5-N7	-2.15	1.32	1.39
2	C	800	AGS	C5-N7	-2.07	1.32	1.39
2	C	800	AGS	C5-C4	2.01	1.45	1.40
2	B	800	AGS	C5-C4	2.14	1.45	1.40

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	AGS	N3-C2-N1	-8.57	122.33	128.89
2	D	800	AGS	N3-C2-N1	-8.30	122.54	128.89
2	B	800	AGS	N3-C2-N1	-8.22	122.60	128.89
2	F	800	AGS	N3-C2-N1	-7.91	122.84	128.89
2	E	800	AGS	N3-C2-N1	-7.66	123.03	128.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	800	AGS	1	0
2	B	800	AGS	1	0
2	C	800	AGS	1	0
2	D	800	AGS	1	0
2	E	800	AGS	1	0
2	F	800	AGS	1	0



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	451/489 (92%)	0.28	19 (4%) 40 28	62, 65, 67, 72	0
1	B	451/489 (92%)	0.19	10 (2%) 65 54	62, 65, 67, 71	0
1	C	451/489 (92%)	0.21	11 (2%) 62 50	62, 65, 67, 70	0
1	D	451/489 (92%)	0.38	27 (5%) 25 15	62, 65, 67, 71	0
1	E	451/489 (92%)	0.31	27 (5%) 25 15	62, 65, 67, 77	0
1	F	451/489 (92%)	0.23	9 (1%) 68 58	62, 65, 67, 70	0
All	All	2706/2934 (92%)	0.27	103 (3%) 44 32	62, 65, 67, 77	0

The worst 5 of 103 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	158	MET	4.5
1	E	106	PRO	4.3
1	D	169	ASP	4.1
1	E	158	MET	3.9
1	D	114	ILE	3.8

### 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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	AGS	A	800	31/31	0.97	0.17	-0.99	53,60,62,64	0
2	AGS	B	800	31/31	0.98	0.14	-1.47	51,60,63,65	0
2	AGS	F	800	31/31	0.98	0.13	-1.64	51,60,62,64	0
2	AGS	D	800	31/31	0.98	0.12	-1.66	53,60,62,65	0
2	AGS	C	800	31/31	0.98	0.14	-1.88	53,61,62,65	0
2	AGS	E	800	31/31	0.99	0.15	-1.88	51,60,62,64	0
3	MG	D	801	1/1	0.96	0.12	-	55,55,55,55	0
3	MG	B	801	1/1	0.97	0.10	-	55,55,55,55	0
3	MG	E	801	1/1	0.99	0.14	-	54,54,54,54	0
3	MG	F	801	1/1	0.97	0.09	-	55,55,55,55	0
3	MG	C	801	1/1	0.98	0.12	-	54,54,54,54	0
3	MG	A	801	1/1	0.96	0.18	-	55,55,55,55	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.