



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 01:40 am BST

PDB ID : 4DNQ  
Title : Crystal Structure of DAD2 S96A mutant  
Authors : Hamiaux, C.  
Deposited on : 2012-02-08  
Resolution : 2.80 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

<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

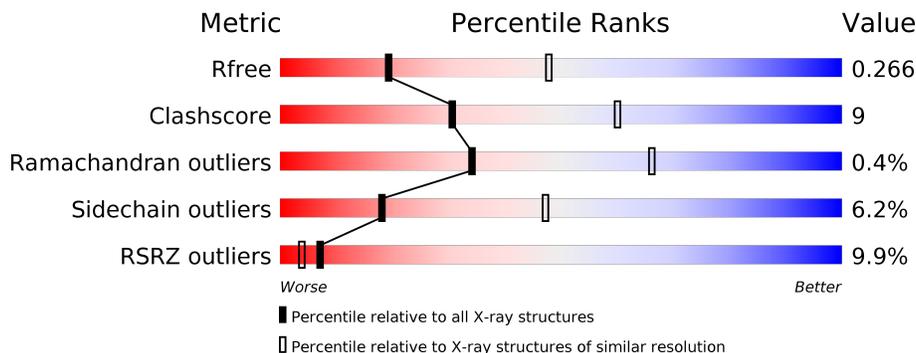
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	269	
1	B	269	
1	C	269	
1	D	269	
1	E	269	
1	F	269	

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	G	269	
1	H	269	
1	I	269	
1	J	269	
1	K	269	
1	L	269	

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 25042 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 DAD2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	262	Total 2054	C 1320	N 359	O 367	S 8	0	0	0
1	B	262	Total 2054	C 1320	N 359	O 367	S 8	0	0	0
1	C	263	Total 2063	C 1325	N 361	O 369	S 8	0	0	0
1	D	263	Total 2063	C 1325	N 361	O 369	S 8	0	0	0
1	E	262	Total 2054	C 1320	N 359	O 367	S 8	0	0	0
1	F	263	Total 2063	C 1325	N 361	O 369	S 8	0	0	0
1	G	263	Total 2063	C 1325	N 361	O 369	S 8	0	0	0
1	H	262	Total 2054	C 1320	N 359	O 367	S 8	0	0	0
1	I	264	Total 2067	C 1327	N 362	O 370	S 8	0	0	0
1	J	263	Total 2063	C 1325	N 361	O 369	S 8	0	0	0
1	K	262	Total 2054	C 1320	N 359	O 367	S 8	0	0	0
1	L	264	Total 2067	C 1327	N 362	O 370	S 8	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	33	Total 33	O 33	0	0
2	B	27	Total 27	O 27	0	0

*Continued on next page...*

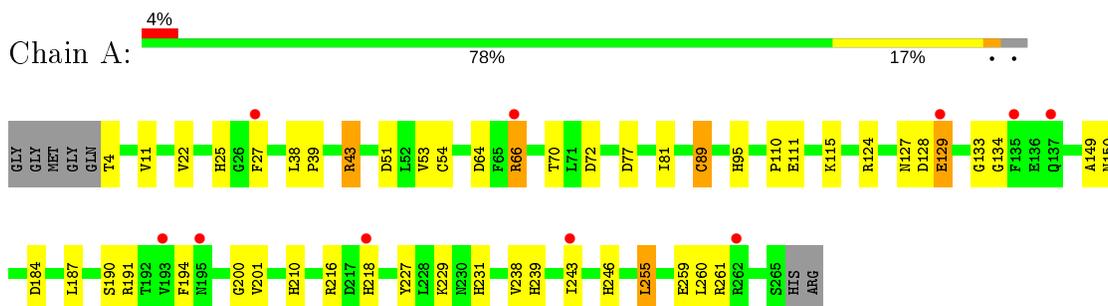
*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
2	C	35	Total 35	O 35	0	0
2	D	32	Total 32	O 32	0	0
2	E	27	Total 27	O 27	0	0
2	F	27	Total 27	O 27	0	0
2	G	31	Total 31	O 31	0	0
2	H	24	Total 24	O 24	0	0
2	I	21	Total 21	O 21	0	0
2	J	12	Total 12	O 12	0	0
2	K	22	Total 22	O 22	0	0
2	L	32	Total 32	O 32	0	0

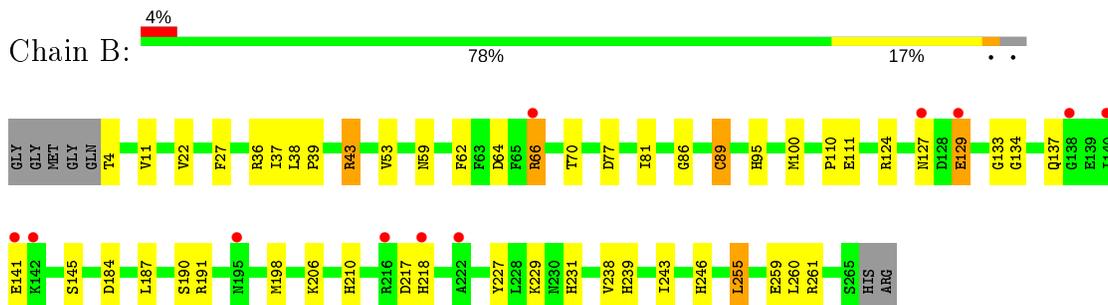
### 3 Residue-property plots [i](#)

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.

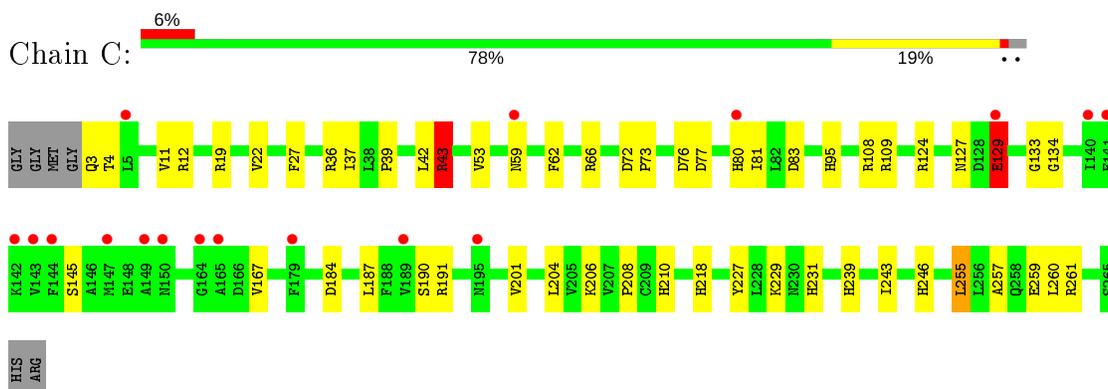
- Molecule 1: DAD2



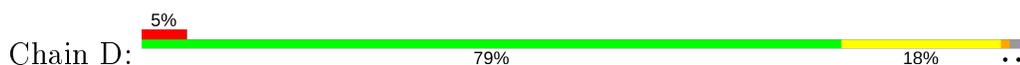
- Molecule 1: DAD2

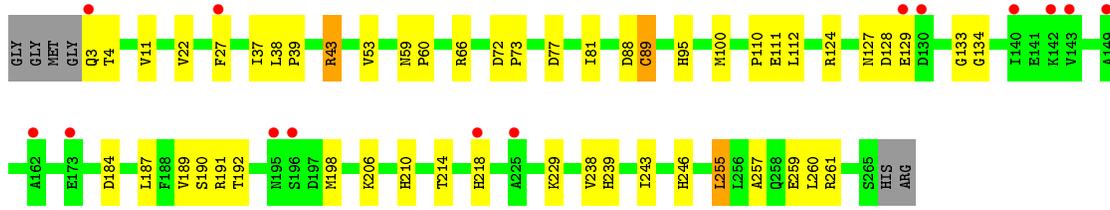


- Molecule 1: DAD2

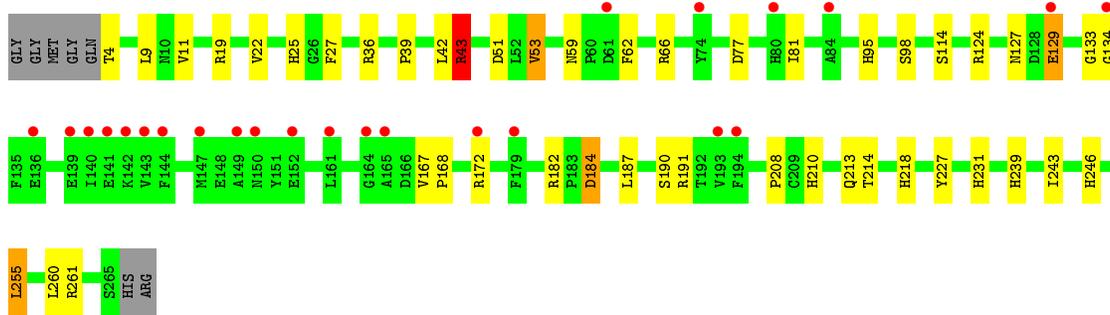
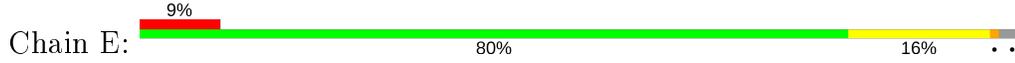


- Molecule 1: DAD2

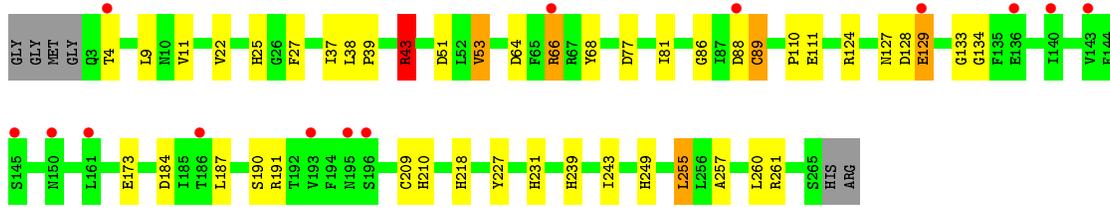
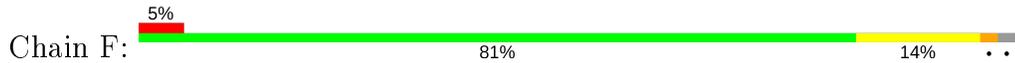




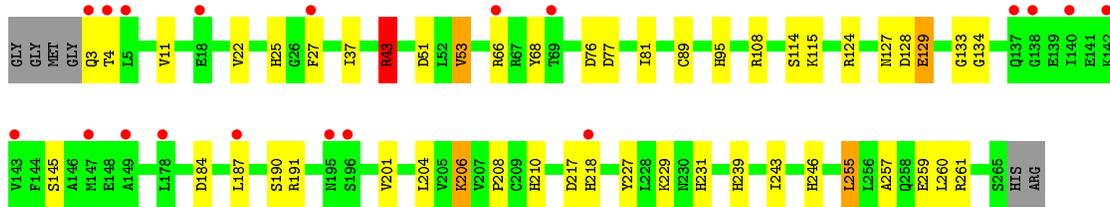
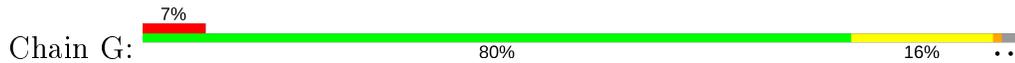
• Molecule 1: DAD2



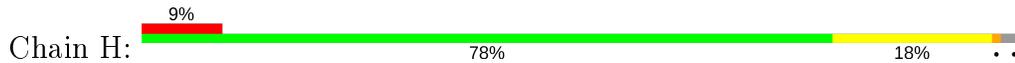
• Molecule 1: DAD2

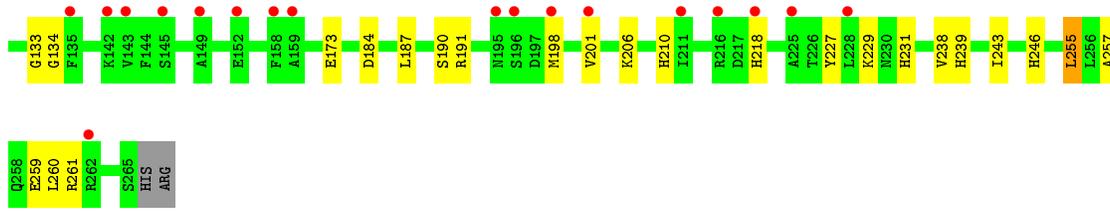


• Molecule 1: DAD2

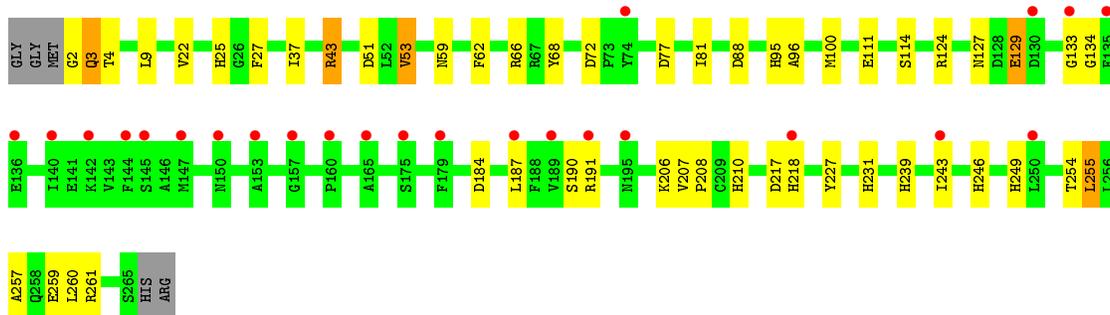
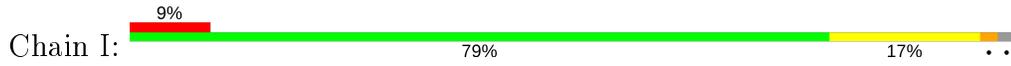


• Molecule 1: DAD2





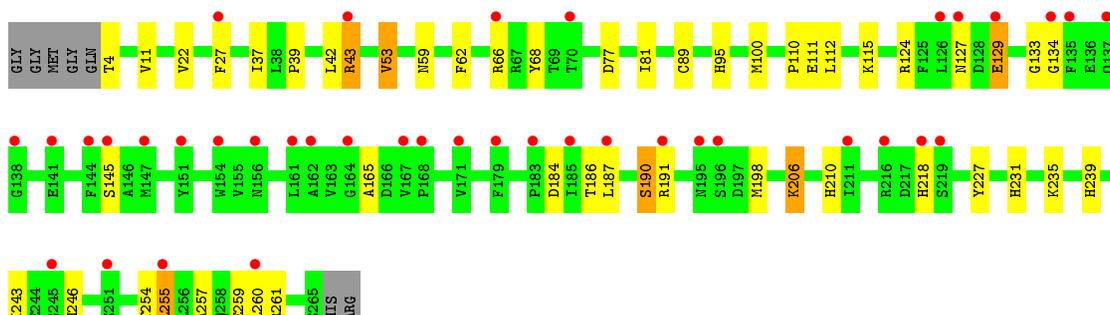
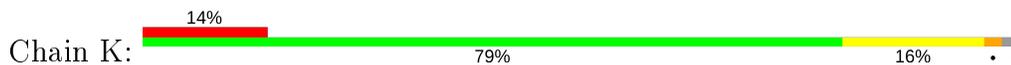
● Molecule 1: DAD2



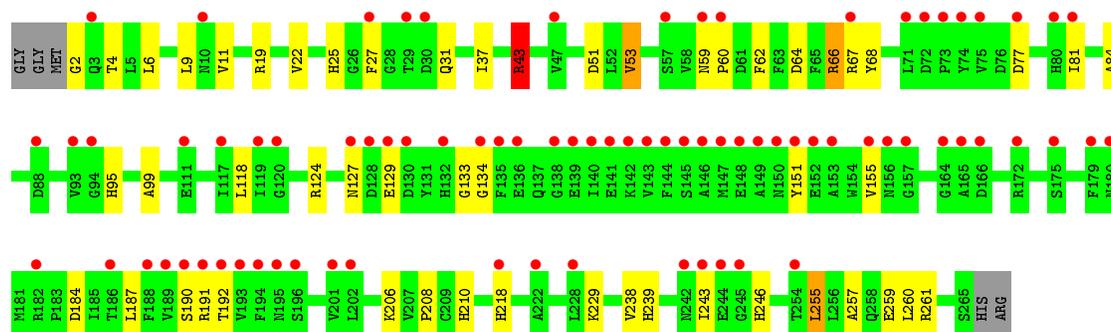
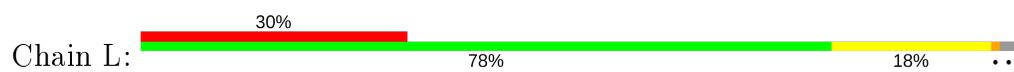
● Molecule 1: DAD2



● Molecule 1: DAD2



● Molecule 1: DAD2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	176.66Å 176.66Å 107.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	62.36 – 2.80 62.36 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (62.36-2.80) 100.0 (62.36-2.80)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	0.18	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.86 (at 2.81Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.242 , 0.272 0.240 , 0.266	Depositor DCC
$R_{free}$ test set	4632 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.9	Xtrriage
Anisotropy	0.074	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 46.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	0.015 for -h,-k,l 0.046 for h,-h-k,-l 0.017 for -k,-h,-l	Xtrriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	25042	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.05 % 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 3.3633e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.333 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	0.82	2/2105 (0.1%)	0.77	1/2865 (0.0%)
1	B	0.69	0/2105	0.75	1/2865 (0.0%)
1	C	0.70	0/2114	0.75	2/2877 (0.1%)
1	D	0.75	1/2114 (0.0%)	0.77	1/2877 (0.0%)
1	E	0.74	0/2105	0.71	1/2865 (0.0%)
1	F	0.76	2/2114 (0.1%)	0.75	2/2877 (0.1%)
1	G	0.72	1/2114 (0.0%)	0.74	1/2877 (0.0%)
1	H	0.75	2/2105 (0.1%)	0.75	1/2865 (0.0%)
1	I	0.70	0/2118	0.73	0/2882
1	J	0.64	0/2114	0.70	1/2877 (0.0%)
1	K	0.70	1/2105 (0.0%)	0.74	2/2865 (0.1%)
1	L	0.68	0/2118	0.70	1/2882 (0.0%)
All	All	0.72	9/25331 (0.0%)	0.74	14/34474 (0.0%)

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	89	CYS	CB-SG	-12.59	1.60	1.82
1	A	89	CYS	CB-SG	-11.16	1.63	1.82
1	A	54	CYS	CB-SG	-7.76	1.69	1.82
1	G	89	CYS	CB-SG	-6.99	1.70	1.82
1	H	89	CYS	CB-SG	-6.10	1.71	1.82

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	89	CYS	CA-CB-SG	12.01	135.62	114.00
1	B	89	CYS	CA-CB-SG	10.46	132.83	114.00
1	D	89	CYS	CA-CB-SG	8.35	129.03	114.00
1	K	89	CYS	CB-CA-C	-8.08	94.23	110.40
1	H	89	CYS	CA-CB-SG	7.47	127.45	114.00

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	2054	0	2034	44	0
1	B	2054	0	2034	43	0
1	C	2063	0	2042	61	0
1	D	2063	0	2042	35	0
1	E	2054	0	2034	54	0
1	F	2063	0	2042	40	0
1	G	2063	0	2042	43	0
1	H	2054	0	2034	44	0
1	I	2067	0	2045	36	1
1	J	2063	0	2042	36	0
1	K	2054	0	2034	44	0
1	L	2067	0	2045	49	0
2	A	33	0	0	3	0
2	B	27	0	0	2	0
2	C	35	0	0	1	0
2	D	32	0	0	2	0
2	E	27	0	0	5	0
2	F	27	0	0	1	0
2	G	31	0	0	0	0
2	H	24	0	0	4	0
2	I	21	0	0	3	0
2	J	12	0	0	1	0
2	K	22	0	0	2	0
2	L	32	0	0	4	0
All	All	25042	0	24470	451	1

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 9.

The worst 5 of 451 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:88:ASP:OD2	1:L:19:ARG:HD3	1.42	1.16

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:36:ARG:HG3	2:E:306:HOH:O	1.50	1.09
1:C:36:ARG:HD3	1:G:108:ARG:HD2	1.34	1.09
1:A:70:THR:HG21	1:K:254:THR:HG21	1.10	1.05
1:F:88:ASP:OD2	1:L:19:ARG:CD	2.06	1.03

All (1) 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)
1:I:72:ASP:OD2	1:I:254:THR:OG1[2_545]	1.83	0.37

## 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	260/269 (97%)	251 (96%)	8 (3%)	1 (0%)	34	66
1	B	260/269 (97%)	247 (95%)	12 (5%)	1 (0%)	34	66
1	C	261/269 (97%)	252 (97%)	7 (3%)	2 (1%)	19	49
1	D	261/269 (97%)	250 (96%)	10 (4%)	1 (0%)	34	66
1	E	260/269 (97%)	249 (96%)	10 (4%)	1 (0%)	34	66
1	F	261/269 (97%)	251 (96%)	9 (3%)	1 (0%)	34	66
1	G	261/269 (97%)	251 (96%)	9 (3%)	1 (0%)	34	66
1	H	260/269 (97%)	252 (97%)	7 (3%)	1 (0%)	34	66
1	I	262/269 (97%)	252 (96%)	8 (3%)	2 (1%)	19	49
1	J	261/269 (97%)	253 (97%)	7 (3%)	1 (0%)	34	66
1	K	260/269 (97%)	250 (96%)	9 (4%)	1 (0%)	34	66
1	L	262/269 (97%)	250 (95%)	11 (4%)	1 (0%)	34	66
All	All	3129/3228 (97%)	3008 (96%)	107 (3%)	14 (0%)	34	66

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	3	GLN
1	I	184	ASP
1	A	184	ASP
1	B	184	ASP
1	C	184	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	219/223 (98%)	206 (94%)	13 (6%)	19	49
1	B	219/223 (98%)	204 (93%)	15 (7%)	16	42
1	C	220/223 (99%)	206 (94%)	14 (6%)	17	45
1	D	220/223 (99%)	206 (94%)	14 (6%)	17	45
1	E	219/223 (98%)	206 (94%)	13 (6%)	19	49
1	F	220/223 (99%)	206 (94%)	14 (6%)	17	45
1	G	220/223 (99%)	203 (92%)	17 (8%)	13	35
1	H	219/223 (98%)	207 (94%)	12 (6%)	21	52
1	I	220/223 (99%)	205 (93%)	15 (7%)	16	42
1	J	220/223 (99%)	208 (94%)	12 (6%)	21	52
1	K	219/223 (98%)	205 (94%)	14 (6%)	17	45
1	L	220/223 (99%)	209 (95%)	11 (5%)	24	56
All	All	2635/2676 (98%)	2471 (94%)	164 (6%)	18	47

5 of 164 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	66	ARG
1	G	129	GLU
1	K	255	LEU
1	F	124	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	G	3	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 90 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	95	HIS
1	G	210	HIS
1	L	95	HIS
1	F	127	ASN
1	F	236	ASN

### 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 [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	262/269 (97%)	0.04	10 (3%) 40 30	21, 38, 63, 76	0
1	B	262/269 (97%)	0.19	11 (4%) 36 26	29, 47, 72, 89	0
1	C	263/269 (97%)	0.24	17 (6%) 18 11	24, 45, 75, 93	0
1	D	263/269 (97%)	0.19	14 (5%) 26 17	25, 43, 70, 85	0
1	E	262/269 (97%)	0.38	24 (9%) 9 5	27, 47, 80, 105	0
1	F	263/269 (97%)	0.32	14 (5%) 26 17	29, 43, 68, 80	0
1	G	263/269 (97%)	0.32	19 (7%) 15 8	27, 45, 69, 87	0
1	H	262/269 (97%)	0.43	25 (9%) 8 4	27, 46, 72, 97	0
1	I	264/269 (98%)	0.42	24 (9%) 9 5	26, 53, 119, 145	0
1	J	263/269 (97%)	0.62	34 (12%) 3 2	32, 60, 132, 168	0
1	K	262/269 (97%)	0.74	39 (14%) 2 1	35, 59, 102, 124	0
1	L	264/269 (98%)	1.40	80 (30%) 0 0	42, 58, 86, 100	0
All	All	3153/3228 (97%)	0.44	311 (9%) 7 4	21, 49, 90, 168	0

The worst 5 of 311 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	138	GLY	11.5
1	J	147	MET	9.0
1	L	189	VAL	7.7
1	L	80	HIS	7.7
1	J	179	PHE	7.3

### 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.