



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 04:54 PM BST

PDB ID : 6BOP  
Title : Crystal structure of 2-methylcitrate synthase from *Aspergillus fumigatus*  
Authors : Schlachter, C.; Chruszcz, M.  
Deposited on : 2017-11-20  
Resolution : 2.71 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](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.

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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
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
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)	:	2.13.1

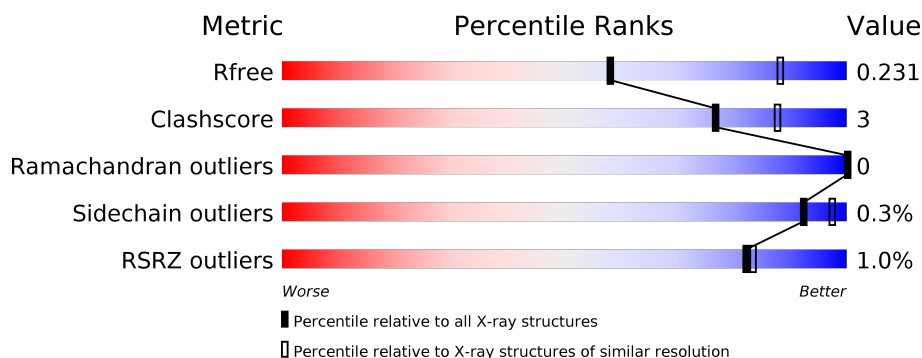
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.71 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))
$R_{free}$	130704	3359 (2.74-2.70)
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)
RSRZ outliers	127900	3276 (2.74-2.70)

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  $\geq 3$ , 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	441	<div> <div>91%</div> <div>7%</div> <div>•</div> </div>
1	B	441	<div> <div>90%</div> <div>8%</div> <div>•</div> </div>
1	C	441	<div> <div>90%</div> <div>8%</div> <div>•</div> </div>
1	D	441	<div> <div>3%</div> <div>89%</div> <div>10%</div> <div>•</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13670 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 2-methylcitrate synthase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	434	Total	C	N	O	S	0	0	0
			3383	2169	585	620	9			
1	A	434	Total	C	N	O	S	0	1	0
			3380	2170	581	620	9			
1	C	434	Total	C	N	O	S	0	0	0
			3371	2160	582	620	9			
1	D	434	Total	C	N	O	S	0	1	0
			3366	2156	581	620	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	25	SER	-	expression tag	UNP Q50I20
B	26	GLY	-	expression tag	UNP Q50I20
B	27	SER	-	expression tag	UNP Q50I20
B	28	GLY	-	expression tag	UNP Q50I20
B	352	ALA	GLY	engineered mutation	UNP Q50I20
A	25	SER	-	expression tag	UNP Q50I20
A	26	GLY	-	expression tag	UNP Q50I20
A	27	SER	-	expression tag	UNP Q50I20
A	28	GLY	-	expression tag	UNP Q50I20
A	352	ALA	GLY	engineered mutation	UNP Q50I20
C	25	SER	-	expression tag	UNP Q50I20
C	26	GLY	-	expression tag	UNP Q50I20
C	27	SER	-	expression tag	UNP Q50I20
C	28	GLY	-	expression tag	UNP Q50I20
C	352	ALA	GLY	engineered mutation	UNP Q50I20
D	25	SER	-	expression tag	UNP Q50I20
D	26	GLY	-	expression tag	UNP Q50I20
D	27	SER	-	expression tag	UNP Q50I20
D	28	GLY	-	expression tag	UNP Q50I20
D	352	ALA	GLY	engineered mutation	UNP Q50I20

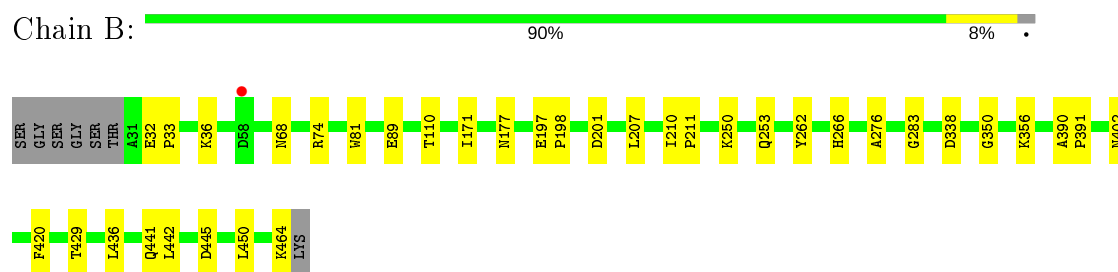
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	65	Total 65	O 65	0	0
2	A	34	Total 34	O 34	0	0
2	C	30	Total 30	O 30	0	0
2	D	41	Total 41	O 41	0	0

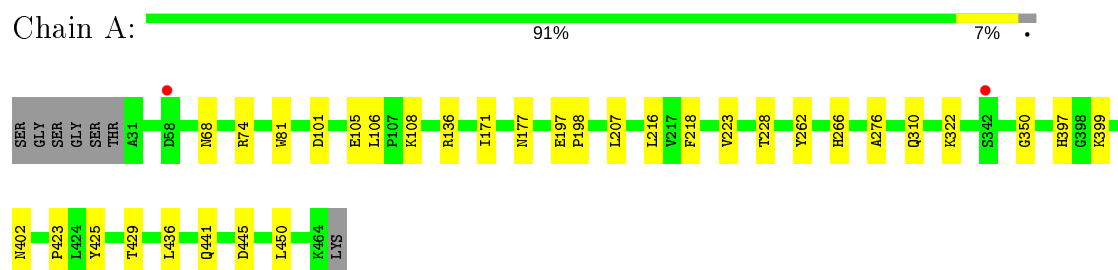
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

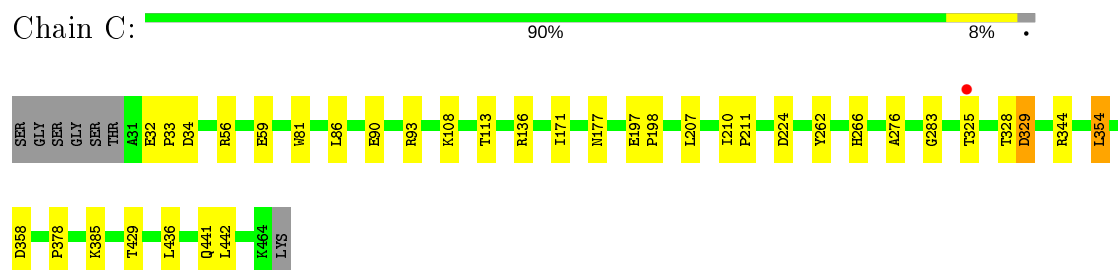
- Molecule 1: 2-methylcitrate synthase, mitochondrial



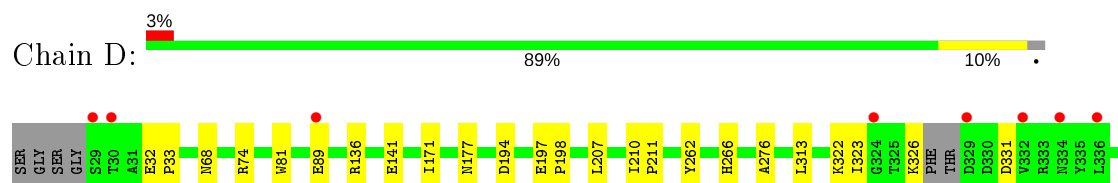
- Molecule 1: 2-methylcitrate synthase, mitochondrial

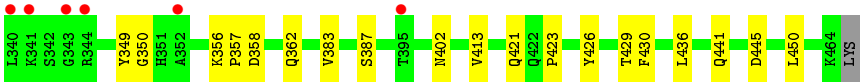


- Molecule 1: 2-methylcitrate synthase, mitochondrial



- Molecule 1: 2-methylcitrate synthase, mitochondrial





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.65Å 100.45Å 209.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.71 36.41 – 2.71	Depositor EDS
% Data completeness (in resolution range)	91.0 (40.00-2.71) 91.0 (36.41-2.71)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.10	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.87 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.187 , 0.233 0.189 , 0.231	Depositor DCC
$R_{free}$ test set	2113 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.6	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 26.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13670	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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, 0.333 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	0.65	0/3466	0.80	1/4705 (0.0%)
1	B	0.68	0/3465	0.83	2/4701 (0.0%)
1	C	0.62	0/3453	0.79	6/4689 (0.1%)
1	D	0.66	0/3449	0.79	4/4682 (0.1%)
All	All	0.65	0/13833	0.80	13/18777 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	136	ARG	NE-CZ-NH2	6.89	123.74	120.30
1	B	420	PHE	CB-CG-CD2	-6.78	116.06	120.80
1	A	136	ARG	NE-CZ-NH2	6.21	123.41	120.30
1	C	224	ASP	CB-CG-OD2	-5.83	113.06	118.30
1	D	194	ASP	CB-CG-OD1	5.75	123.48	118.30
1	C	34	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	B	338	ASP	CB-CG-OD1	5.37	123.13	118.30
1	C	344	ARG	NE-CZ-NH1	5.35	122.97	120.30
1	C	354	LEU	CB-CG-CD1	-5.18	102.20	111.00
1	D	358	ASP	CB-CG-OD1	5.16	122.95	118.30
1	D	430	PHE	CB-CG-CD1	5.13	124.39	120.80
1	D	136	ARG	NE-CZ-NH2	5.09	122.84	120.30
1	C	358	ASP	CB-CG-OD1	5.08	122.88	118.30

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	3380	0	3384	18	0
1	B	3383	0	3402	20	0
1	C	3371	0	3369	18	0
1	D	3366	0	3368	29	0
2	A	34	0	0	0	0
2	B	65	0	0	0	0
2	C	30	0	0	0	0
2	D	41	0	0	0	0
All	All	13670	0	13523	82	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.

All (82) 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:D:326:LYS:O	1:D:331:ASP:OD2	2.02	0.78
1:A:310:GLN:HG3	1:A:423:PRO:HB2	1.68	0.74
1:D:313:LEU:HD22	1:D:426:TYR:HE2	1.64	0.63
1:C:90:GLU:OE1	1:C:93:ARG:NH1	2.36	0.58
1:A:397:HIS:NE2	1:A:399:LYS:HD2	2.20	0.57
1:D:323:ILE:HG22	1:D:323:ILE:O	2.04	0.56
1:C:329:ASP:OD1	1:C:385:LYS:NZ	2.35	0.55
1:A:171:ILE:HG23	1:C:171:ILE:HG23	1.88	0.55
1:B:171:ILE:HG23	1:D:171:ILE:HG23	1.89	0.55
1:C:108:LYS:NZ	1:C:113:THR:O	2.32	0.55
1:D:313:LEU:HD22	1:D:426:TYR:CE2	2.42	0.54
1:B:207:LEU:HD11	1:B:436:LEU:HD23	1.89	0.54
1:D:383:VAL:HG21	1:D:413:VAL:HG21	1.88	0.54
1:B:197:GLU:HB2	1:B:198:PRO:HD3	1.90	0.53
1:C:197:GLU:HB2	1:C:198:PRO:HD3	1.92	0.51
1:D:197:GLU:HB2	1:D:198:PRO:HD3	1.92	0.51
1:A:262:TYR:CE2	1:A:429:THR:HG22	2.45	0.51
1:B:36:LYS:NZ	1:B:201:ASP:OD1	2.44	0.50
1:D:89:GLU:OE1	1:D:356:LYS:HE3	2.11	0.50
1:C:56:ARG:HB3	1:C:59:GLU:HG3	1.93	0.50
1:D:262:TYR:CE2	1:D:429:THR:HG22	2.47	0.49
1:D:383:VAL:CG2	1:D:413:VAL:HG21	2.43	0.49
1:B:89:GLU:OE1	1:B:356:LYS:HE3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:328:THR:HG22	1:C:329:ASP:N	2.29	0.48
1:B:262:TYR:CE2	1:B:429:THR:HG22	2.49	0.48
1:C:210:ILE:HG22	1:C:211:PRO:HD3	1.95	0.47
1:D:357:PRO:HB2	1:D:362[B]:GLN:HE21	1.79	0.47
1:C:262:TYR:CE2	1:C:429:THR:HG22	2.48	0.47
1:A:197:GLU:HB2	1:A:198:PRO:HD3	1.97	0.47
1:B:210:ILE:HG22	1:B:211:PRO:HD3	1.96	0.47
1:D:313:LEU:CD2	1:D:426:TYR:HE2	2.28	0.47
1:D:141:GLU:HA	1:D:141:GLU:OE2	2.15	0.47
1:B:210:ILE:CG2	1:B:211:PRO:HD3	2.45	0.46
1:D:323:ILE:O	1:D:323:ILE:CG2	2.64	0.46
1:D:81:TRP:CE3	1:D:441:GLN:HG2	2.50	0.46
1:D:210:ILE:HG22	1:D:211:PRO:HD3	1.97	0.46
1:C:210:ILE:CG2	1:C:211:PRO:HD3	2.46	0.45
1:C:81:TRP:CE3	1:C:441:GLN:HG2	2.51	0.45
1:D:210:ILE:CG2	1:D:211:PRO:HD3	2.46	0.45
1:D:421:GLN:O	1:D:423:PRO:HD3	2.17	0.45
1:B:283:GLY:HA3	1:B:442:LEU:HD21	1.98	0.45
1:A:81:TRP:CE3	1:A:441:GLN:HG2	2.52	0.44
1:D:445:ASP:OD1	1:D:450:LEU:HD12	2.18	0.44
1:B:110:THR:HB	1:A:228:THR:O	2.18	0.44
1:B:81:TRP:CE3	1:B:441:GLN:HG2	2.52	0.44
1:B:250:LYS:O	1:B:253:GLN:HG3	2.19	0.43
1:C:32:GLU:HB3	1:C:33:PRO:HD2	1.99	0.43
1:A:101:ASP:O	1:A:105:GLU:HG3	2.19	0.43
1:B:445:ASP:OD1	1:B:450:LEU:HD12	2.19	0.43
1:D:262:TYR:CE1	1:D:266:HIS:CE1	3.07	0.43
1:B:262:TYR:CE1	1:B:266:HIS:CE1	3.06	0.43
1:A:262:TYR:CE1	1:A:266:HIS:CE1	3.07	0.43
1:D:356:LYS:HB3	1:D:357:PRO:HD2	2.00	0.43
1:B:464:LYS:HE3	1:B:464:LYS:HB2	1.70	0.42
1:D:32:GLU:HA	1:D:33:PRO:HD3	1.79	0.42
1:A:216:LEU:HD22	1:A:223:VAL:HA	2.01	0.42
1:A:218[A]:PHE:CZ	1:A:425:TYR:HB3	2.55	0.42
1:B:68:ASN:ND2	1:B:74:ARG:O	2.53	0.42
1:D:322:LYS:HA	1:D:322:LYS:HD3	1.87	0.42
1:A:445:ASP:OD1	1:A:450:LEU:HD12	2.20	0.42
1:C:262:TYR:CE1	1:C:266:HIS:CE1	3.08	0.42
1:A:350:GLY:HA2	1:A:402:ASN:O	2.20	0.41
1:D:81:TRP:CH2	1:D:276:ALA:HB1	2.55	0.41
1:B:81:TRP:CH2	1:B:276:ALA:HB1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:LEU:O	1:A:108:LYS:NZ	2.53	0.41
1:A:68:ASN:ND2	1:A:74:ARG:O	2.52	0.41
1:A:81:TRP:CH2	1:A:276:ALA:HB1	2.55	0.41
1:C:207:LEU:HD11	1:C:436:LEU:HD23	2.02	0.41
1:C:325:THR:HG23	1:C:378:PRO:CG	2.51	0.41
1:D:68:ASN:ND2	1:D:74:ARG:O	2.53	0.41
1:A:207:LEU:HD11	1:A:436:LEU:HD23	2.02	0.41
1:D:356:LYS:HB3	1:D:357:PRO:CD	2.51	0.41
1:D:350:GLY:HA2	1:D:402:ASN:O	2.20	0.41
1:D:207:LEU:HD11	1:D:436:LEU:HD23	2.02	0.41
1:B:350:GLY:HA2	1:B:402:ASN:O	2.21	0.41
1:D:349:TYR:OH	1:D:387:SER:HA	2.21	0.41
1:B:32:GLU:HA	1:B:33:PRO:HD3	1.79	0.41
1:B:390:ALA:N	1:B:391:PRO:CD	2.85	0.40
1:C:283:GLY:HA3	1:C:442:LEU:HD11	2.03	0.40
1:C:81:TRP:CH2	1:C:276:ALA:HB1	2.56	0.40
1:A:322:LYS:HD2	1:A:322:LYS:HA	1.90	0.40
1:C:86:LEU:HD11	1:C:354:LEU:HD11	2.04	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	433/441 (98%)	425 (98%)	8 (2%)	0	100	100
1	B	432/441 (98%)	426 (99%)	6 (1%)	0	100	100
1	C	432/441 (98%)	428 (99%)	4 (1%)	0	100	100
1	D	431/441 (98%)	424 (98%)	7 (2%)	0	100	100
All	All	1728/1764 (98%)	1703 (99%)	25 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	358/365 (98%)	357 (100%)	1 (0%)	92	97
1	B	360/365 (99%)	359 (100%)	1 (0%)	92	97
1	C	357/365 (98%)	355 (99%)	2 (1%)	86	94
1	D	357/365 (98%)	356 (100%)	1 (0%)	92	97
All	All	1432/1460 (98%)	1427 (100%)	5 (0%)	92	97

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	177	ASN
1	A	177	ASN
1	C	177	ASN
1	C	329	ASP
1	D	177	ASN

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

Mol	Chain	Res	Type
1	A	266	HIS
1	D	266	HIS

### 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 monosaccharides 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, 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	434/441 (98%)	-0.38	2 (0%) 91 92	21, 34, 57, 76	0
1	B	434/441 (98%)	-0.50	1 (0%) 95 96	19, 30, 49, 65	0
1	C	434/441 (98%)	-0.28	1 (0%) 95 96	19, 38, 71, 85	0
1	D	434/441 (98%)	-0.23	14 (3%) 47 48	17, 33, 98, 127	0
All	All	1736/1764 (98%)	-0.35	18 (1%) 82 83	17, 33, 70, 127	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	30	THR	4.0
1	D	29	SER	3.2
1	D	343	GLY	3.1
1	D	324	GLY	2.8
1	D	395	THR	2.7
1	D	344	ARG	2.6
1	C	325	THR	2.6
1	D	332	VAL	2.5
1	D	334	ASN	2.4
1	D	89	GLU	2.4
1	D	340	LEU	2.4
1	D	352	ALA	2.3
1	B	58	ASP	2.3
1	A	342	SER	2.1
1	D	336	LEU	2.1
1	D	329	ASP	2.1
1	A	58	ASP	2.1
1	D	341	LYS	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 monosaccharides 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.