



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2017 – 08:38 PM EDT

PDB ID : 3DAX  
Title : Crystal structure of human CYP7A1  
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A.; Park, H.; Structural Genomics Consortium (SGC)  
Deposited on : unknown  
Resolution : 2.15 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

<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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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-20030345

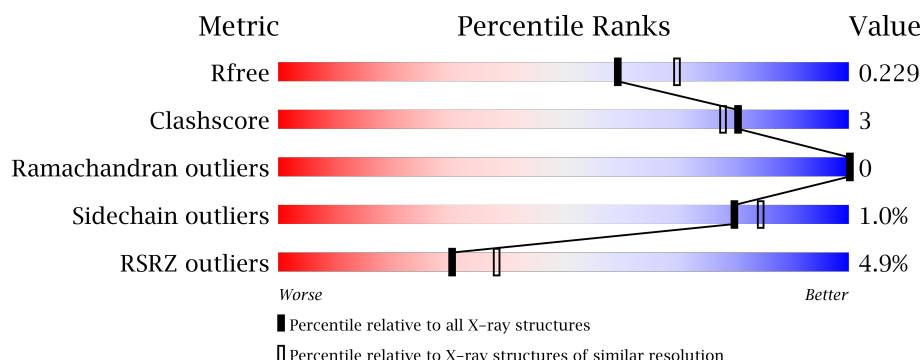
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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))
$R_{free}$	100719	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (2.16-2.16)

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	491	<div> <div>5%</div> <div>89%</div> <div>7%</div> <div>•</div> </div>
1	B	491	<div> <div>4%</div> <div>89%</div> <div>6%</div> <div>•</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7737 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 Cytochrome P450 7A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	8	0
			3773	2438	639	672	24			
1	B	470	Total	C	N	O	S	0	5	0
			3746	2416	631	675	24			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	expression tag	UNP P22680
A	19	ALA	-	expression tag	UNP P22680
A	20	LYS	-	expression tag	UNP P22680
A	21	LYS	-	expression tag	UNP P22680
A	22	THR	-	expression tag	UNP P22680
A	23	SER	-	expression tag	UNP P22680
A	24	SER	-	expression tag	UNP P22680
A	504	HIS	-	expression tag	UNP P22680
A	505	HIS	-	expression tag	UNP P22680
A	506	HIS	-	expression tag	UNP P22680
A	507	HIS	-	expression tag	UNP P22680
A	508	HIS	-	expression tag	UNP P22680
B	18	MET	-	expression tag	UNP P22680
B	19	ALA	-	expression tag	UNP P22680
B	20	LYS	-	expression tag	UNP P22680
B	21	LYS	-	expression tag	UNP P22680
B	22	THR	-	expression tag	UNP P22680
B	23	SER	-	expression tag	UNP P22680
B	24	SER	-	expression tag	UNP P22680
B	504	HIS	-	expression tag	UNP P22680
B	505	HIS	-	expression tag	UNP P22680
B	506	HIS	-	expression tag	UNP P22680
B	507	HIS	-	expression tag	UNP P22680
B	508	HIS	-	expression tag	UNP P22680

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total X 1 1	0	0
3	A	1	Total X 1 1	0	0

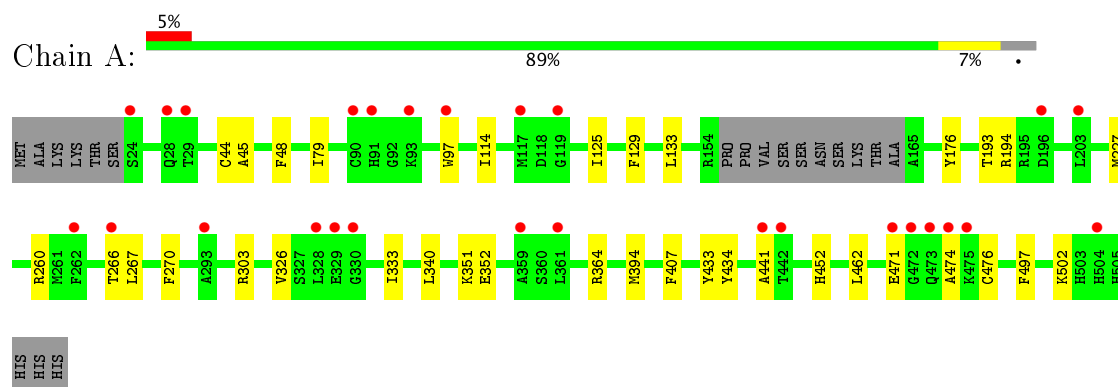
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	62	Total O 62 62	0	0
4	B	68	Total O 68 68	0	0

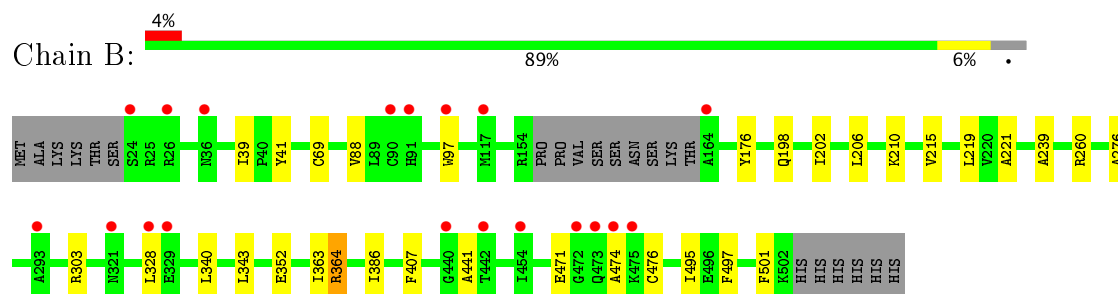
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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: Cytochrome P450 7A1



#### • Molecule 1: Cytochrome P450 7A1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.34Å 80.16Å 84.94Å 64.36° 75.23° 72.17°	Depositor
Resolution (Å)	30.00 – 2.15 29.03 – 2.15	Depositor EDS
% Data completeness (in resolution range)	93.4 (30.00-2.15) 83.7 (29.03-2.15)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.02 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, $R_{free}$	0.189 , 0.228 0.192 , 0.229	Depositor DCC
$R_{free}$ test set	2018 reflections (3.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	34.3	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 57.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7737	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 5.65% 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

### 5.1 Standard geometry

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

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	0.79	0/3891	0.63	0/5270
1	B	0.80	1/3852 (0.0%)	0.65	0/5215
All	All	0.79	1/7743 (0.0%)	0.64	0/10485

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	69	CYS	CB-SG	-5.03	1.73	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	3773	0	3693	21	0
1	B	3746	0	3669	21	0
2	A	43	0	30	2	0
2	B	43	0	30	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	62	0	0	0	0
4	B	68	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7737	0	7422	46	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 (46) 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)
2:B:601:HEM:HHD	2:B:601:HEM:HBC2	1.68	0.75
1:B:471:GLU:HB3	1:B:474:ALA:HB2	1.72	0.72
1:A:471:GLU:HB3	1:A:474:ALA:HB2	1.70	0.72
1:A:326:VAL:HG22	1:A:333:ILE:HD11	1.74	0.69
1:A:227:MET:HE1	1:B:39:ILE:HG12	1.75	0.69
1:A:114:ILE:HG22	1:A:270:PHE:CZ	2.30	0.66
1:B:97:TRP:CH2	1:B:441:ALA:HB1	2.35	0.61
1:A:114:ILE:HG22	1:A:270:PHE:HZ	1.65	0.61
1:B:206:LEU:HD21	1:B:210:LYS:HE3	1.85	0.59
1:B:303:ARG:CZ	1:B:476[A]:CYS:SG	2.90	0.58
1:A:351:LYS:HG2	1:A:434:TYR:CZ	2.39	0.58
2:A:601:HEM:HMB2	2:A:601:HEM:HBB2	1.83	0.58
1:B:471:GLU:CB	1:B:474:ALA:HB2	2.37	0.55
1:B:215[A]:VAL:HG12	1:B:219:LEU:HG	1.90	0.54
1:A:471:GLU:CB	1:A:474:ALA:HB2	2.36	0.53
2:A:601:HEM:CMB	2:A:601:HEM:HBB2	2.38	0.53
1:B:206:LEU:C	1:B:206:LEU:HD23	2.31	0.51
1:B:352:GLU:HG3	1:B:407:PHE:CD1	2.45	0.51
1:B:41:TYR:HB3	1:B:221:ALA:HB2	1.93	0.50
1:B:340:LEU:O	1:B:343:LEU:HD12	2.11	0.49
1:A:79:ILE:N	1:A:79:ILE:HD12	2.28	0.49
1:A:352:GLU:HG3	1:A:407:PHE:CD1	2.48	0.48
1:B:198:GLN:HG2	1:B:202:ILE:HD12	1.95	0.48
2:B:601:HEM:HBB2	2:B:601:HEM:HMB2	1.96	0.48
1:A:462:LEU:O	1:A:502:LYS:NZ	2.47	0.46
1:A:266:THR:HG22	1:A:267:LEU:HD23	1.97	0.45
1:B:328:LEU:HD11	1:B:501:PHE:CD2	2.52	0.44
1:B:363:ILE:HG23	1:B:386:ILE:HG23	1.98	0.44
2:B:601:HEM:CHD	2:B:601:HEM:HBC2	2.42	0.43
1:A:114:ILE:HD12	1:A:125:ILE:HD13	2.01	0.43
1:B:88:VAL:HG12	1:B:364:ARG:CZ	2.49	0.43
2:B:601:HEM:HBB2	2:B:601:HEM:CMB	2.50	0.42
1:A:193:THR:OG1	1:A:194:ARG:CD	2.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:495:ILE:C	1:B:495:ILE:HD12	2.40	0.41
1:A:303:ARG:CZ	1:A:476[A]:CYS:SG	3.08	0.41
1:A:97:TRP:CH2	1:A:441:ALA:HB1	2.55	0.41
1:A:394:MET:SD	1:A:433:TYR:HA	2.60	0.41
1:A:44:CYS:O	1:A:45:ALA:HB3	2.21	0.41
1:A:48:PHE:HB3	1:B:215[A]:VAL:HG22	2.03	0.40
1:B:239:ALA:O	1:B:276:ALA:HB1	2.21	0.40
1:A:129:PHE:O	1:A:133:LEU:HB2	2.20	0.40
1:A:340:LEU:HD23	1:A:452:HIS:CD2	2.57	0.40
1:A:303:ARG:NE	1:A:476[A]:CYS:SG	2.94	0.40
1:B:206:LEU:CD2	1:B:210:LYS:HE3	2.50	0.40

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	476/491 (97%)	470 (99%)	6 (1%)	0	100	100
1	B	471/491 (96%)	464 (98%)	7 (2%)	0	100	100
All	All	947/982 (96%)	934 (99%)	13 (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	395/433 (91%)	391 (99%)	4 (1%)	80	84
1	B	394/433 (91%)	390 (99%)	4 (1%)	80	84
All	All	789/866 (91%)	781 (99%)	8 (1%)	80	84

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	176	TYR
1	A	260	ARG
1	A	364	ARG
1	A	497	PHE
1	B	176	TYR
1	B	260	ARG
1	B	364	ARG
1	B	497	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	205	ASN
1	B	205	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](#)

Of 4 ligands modelled in this entry, 2 are unknown - leaving 2 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	HEM	A	601	1	28,50,50	2.10	6 (21%)	17,82,82	1.42	1 (5%)
2	HEM	B	601	1	28,50,50	2.25	10 (35%)	17,82,82	1.36	2 (11%)

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	HEM	A	601	1	-	0/6/54/54	0/0/8/8
2	HEM	B	601	1	-	0/6/54/54	0/0/8/8

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HEM	C3C-C2C	-5.33	1.33	1.40
2	B	601	HEM	C3C-C2C	-4.76	1.34	1.40
2	B	601	HEM	C3B-C2B	-4.69	1.34	1.40
2	A	601	HEM	C3B-C2B	-4.11	1.34	1.40
2	B	601	HEM	CMC-C2C	2.07	1.56	1.51
2	B	601	HEM	CMD-C2D	2.16	1.56	1.51
2	A	601	HEM	CAD-C3D	2.32	1.56	1.52
2	B	601	HEM	C1D-ND	2.38	1.41	1.36
2	B	601	HEM	C4A-NA	2.38	1.41	1.36
2	A	601	HEM	C3B-CAB	2.76	1.53	1.47
2	B	601	HEM	C3B-CAB	3.07	1.54	1.47
2	B	601	HEM	C3C-CAC	3.10	1.53	1.47
2	A	601	HEM	C3C-CAC	3.23	1.54	1.47
2	B	601	HEM	C4D-ND	3.26	1.40	1.36
2	A	601	HEM	C3D-C2D	5.07	1.52	1.37
2	B	601	HEM	C3D-C2D	5.20	1.53	1.37

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	A	601	HEM	CAA-CBA-CGA	-3.90	105.99	112.66
2	B	601	HEM	CBD-CAD-C3D	-2.23	108.21	112.47
2	B	601	HEM	CBA-CAA-C2A	-2.07	108.53	112.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	2	0
2	B	601	HEM	4	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

### 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	472/491 (96%)	0.11	27 (5%) 24 32	19, 36, 62, 95	0
1	B	470/491 (95%)	0.07	19 (4%) 39 46	21, 35, 63, 92	0
All	All	942/982 (95%)	0.09	46 (4%) 30 38	19, 35, 63, 95	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	474	ALA	9.2
1	A	473	GLN	7.6
1	A	97	TRP	6.6
1	B	473	GLN	5.7
1	A	474	ALA	5.2
1	B	472	GLY	5.2
1	A	91	HIS	5.1
1	B	329	GLU	4.9
1	A	472	GLY	4.7
1	A	475	LYS	4.7
1	B	91	HIS	4.5
1	B	328	LEU	4.4
1	A	329	GLU	4.1
1	A	29	THR	3.9
1	A	24	SER	3.8
1	B	97	TRP	3.5
1	A	90	CYS	3.4
1	B	90	CYS	3.4
1	B	117	MET	3.3
1	A	117	MET	3.3
1	B	442	THR	3.2
1	B	26	ARG	3.2
1	A	28	GLN	3.0
1	A	330	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	164	ALA	2.5
1	A	504	HIS	2.4
1	B	36	ASN	2.4
1	A	293	ALA	2.4
1	B	24	SER	2.3
1	A	119	GLY	2.3
1	A	361	LEU	2.3
1	A	328	LEU	2.3
1	A	266	THR	2.3
1	A	262	PHE	2.2
1	B	454	ILE	2.2
1	A	196	ASP	2.2
1	A	471	GLU	2.2
1	A	93	LYS	2.2
1	B	293	ALA	2.2
1	A	359	ALA	2.1
1	B	321	ASN	2.1
1	A	441	ALA	2.1
1	A	442	THR	2.1
1	B	440	GLY	2.1
1	B	475	LYS	2.1
1	A	203	LEU	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 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( $\text{\AA}^2$ )	Q<0.9
2	HEM	A	601	43/43	0.97	0.15	0.11	20,25,51,55	0
2	HEM	B	601	43/43	0.97	0.15	-0.06	17,25,46,50	0
3	UNX	A	1	1/1	0.66	1.99	-	2,2,2,2	1
3	UNX	B	2	1/1	0.59	2.19	-	2,2,2,2	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.