



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 20, 2021 – 10:03 AM EDT

PDB ID : 6VZN  
Title : Crystal structure of human PPARgamma ligand binding domain Y473E mutant  
Authors : Shang, J.; Kojetin, D.J.  
Deposited on : 2020-02-28  
Resolution : 2.30 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 : 2.23.1  
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) : 2.23.1

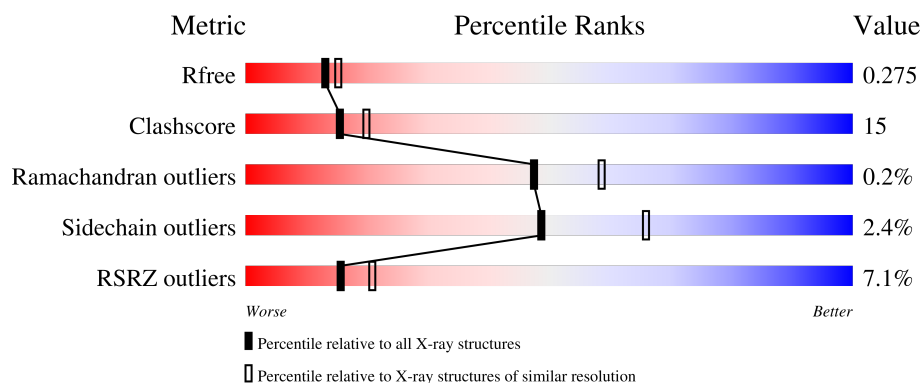
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 of 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	275	<div> <div>5%</div> <div>69%</div> <div>22%</div> <div>8%</div> </div>
1	B	275	<div> <div>8%</div> <div>68%</div> <div>22%</div> <div>9%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4172 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 Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	254	Total	C	N	O	S	0	0	0
			2037	1316	332	379	10			
1	B	251	Total	C	N	O	S	0	0	0
			2013	1303	331	370	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	473	GLU	TYR	engineered mutation	UNP P37231
B	473	GLU	TYR	engineered mutation	UNP P37231

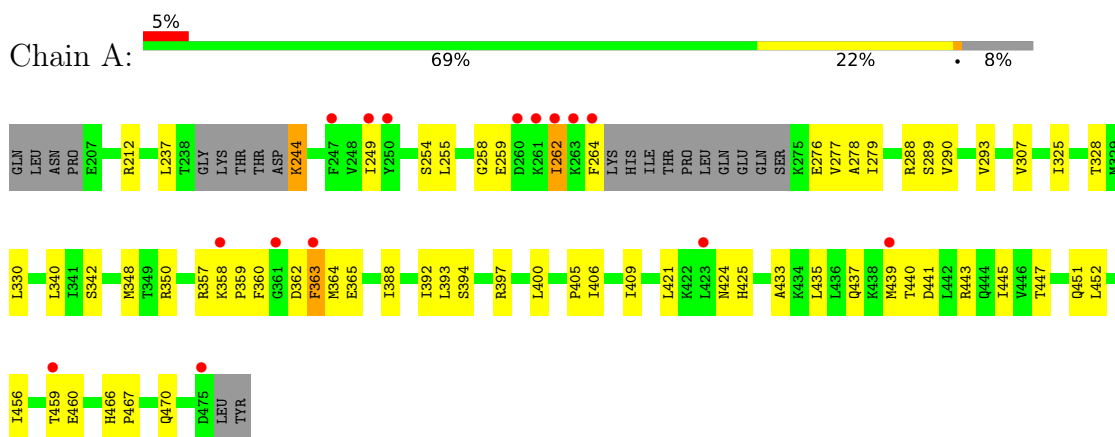
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	55	Total	O	0	0
			55	55		
2	B	67	Total	O	0	0
			67	67		

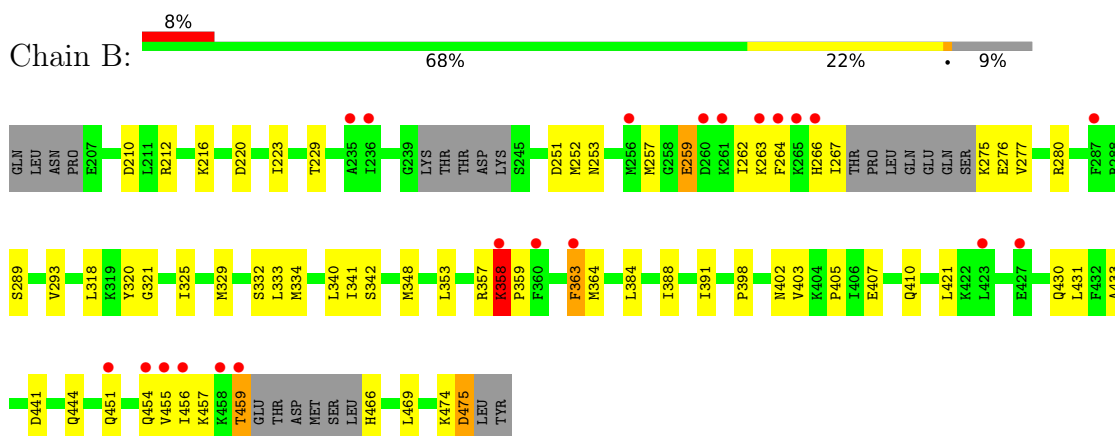
### 3 Residue-property plots [i](#)

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: Peroxisome proliferator-activated receptor gamma



- Molecule 1: Peroxisome proliferator-activated receptor gamma



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.64Å 61.75Å 118.78Å 90.00° 101.93° 90.00°	Depositor
Resolution (Å)	36.30 – 2.30 36.31 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.9 (36.30-2.30) 97.9 (36.31-2.30)	Depositor EDS
$R_{merge}$	0.02	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, $R_{free}$	0.223 , 0.278 0.224 , 0.275	Depositor DCC
$R_{free}$ test set	1993 reflections (6.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.6	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 44.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4172	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.97% 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	0.43	0/2069	0.58	0/2783
1	B	0.53	1/2045 (0.0%)	0.63	0/2749
All	All	0.48	1/4114 (0.0%)	0.61	0/5532

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	358	LYS	C-N	9.01	1.51	1.34

There are no bond angle outliers.

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	2037	0	2099	53	0
1	B	2013	0	2077	69	0
2	A	55	0	0	2	0
2	B	67	0	0	3	0
All	All	4172	0	4176	122	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 15.

All (122) 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:B:456:ILE:HA	1:B:459:THR:HG21	1.25	1.12
1:B:358:LYS:HB3	1:B:359:PRO:HD3	1.16	1.09
1:B:456:ILE:HA	1:B:459:THR:CG2	1.83	1.08
1:B:358:LYS:HB3	1:B:359:PRO:CD	1.88	1.02
1:B:358:LYS:CB	1:B:359:PRO:HD3	2.02	0.89
1:B:359:PRO:CG	1:B:456:ILE:HD11	2.02	0.89
1:B:252:MET:HE1	1:B:277:VAL:HG21	1.56	0.88
1:B:357:ARG:HG3	1:B:357:ARG:HH21	1.40	0.86
1:B:454:GLN:HA	1:B:457:LYS:HE3	1.57	0.86
1:B:359:PRO:CB	1:B:456:ILE:HD11	2.10	0.82
1:B:359:PRO:HB2	1:B:456:ILE:HD11	1.61	0.82
1:B:325:ILE:HD12	1:B:388:ILE:HB	1.63	0.80
1:B:456:ILE:C	1:B:459:THR:HG23	2.04	0.77
1:A:249:ILE:HD13	1:A:255:LEU:N	2.02	0.74
1:A:244:LYS:N	2:A:501:HOH:O	2.21	0.74
1:B:456:ILE:CA	1:B:459:THR:CG2	2.65	0.74
1:B:359:PRO:HG2	1:B:456:ILE:HD11	1.70	0.72
1:B:444:GLN:NE2	2:B:501:HOH:O	2.23	0.72
1:A:456:ILE:HA	1:A:459:THR:HG22	1.71	0.71
1:A:258:GLY:O	1:A:262:ILE:HG13	1.90	0.71
1:A:363:PHE:CE1	1:A:456:ILE:HD11	2.27	0.69
1:B:275:LYS:HB3	1:B:280:ARG:HD3	1.74	0.69
1:B:454:GLN:HA	1:B:457:LYS:CE	2.23	0.69
1:B:456:ILE:HA	1:B:459:THR:HG23	1.74	0.68
1:B:259:GLU:OE2	1:B:280:ARG:NH1	2.26	0.68
1:B:456:ILE:HD13	1:B:459:THR:HG21	1.76	0.68
1:A:249:ILE:HD13	1:A:255:LEU:CA	2.23	0.67
1:A:363:PHE:CZ	1:A:452:LEU:HB3	2.31	0.66
1:A:437:GLN:O	1:A:440:THR:HG22	1.96	0.65
1:B:454:GLN:HA	1:B:457:LYS:HG3	1.78	0.65
1:B:357:ARG:HG3	1:B:357:ARG:NH2	2.11	0.65
1:B:456:ILE:CA	1:B:459:THR:HG23	2.26	0.65
1:B:359:PRO:CG	1:B:456:ILE:CD1	2.75	0.63
1:A:249:ILE:HG23	1:A:348:MET:HA	1.81	0.62
1:A:467:PRO:HA	1:A:470:GLN:HB2	1.81	0.62
1:B:454:GLN:HA	1:B:457:LYS:CD	2.30	0.62
1:B:276:GLU:OE1	1:B:357:ARG:NE	2.31	0.61
1:B:456:ILE:CD1	1:B:459:THR:HG21	2.31	0.61
1:A:276:GLU:OE1	1:A:357:ARG:NH2	2.33	0.59
1:B:216:LYS:HD3	1:B:220:ASP:OD2	2.03	0.58
1:B:474:LYS:O	1:B:475:ASP:HB2	2.05	0.57
1:B:333:LEU:HB3	1:B:340:LEU:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:ARG:HG2	1:A:358:LYS:N	2.21	0.55
1:B:454:GLN:CA	1:B:457:LYS:HE3	2.33	0.55
1:A:259:GLU:HA	1:A:262:ILE:O	2.07	0.55
1:B:410:GLN:NE2	2:B:502:HOH:O	2.39	0.55
1:A:357:ARG:HG2	1:A:358:LYS:H	1.72	0.55
1:A:357:ARG:NH1	1:A:460:GLU:OE2	2.35	0.55
1:A:325:ILE:HD11	1:A:392:ILE:HG13	1.89	0.54
1:B:212:ARG:HD2	2:B:533:HOH:O	2.07	0.54
1:B:321:GLY:O	1:B:325:ILE:HG12	2.08	0.54
1:B:454:GLN:HA	1:B:457:LYS:CG	2.38	0.54
1:B:259:GLU:CD	1:B:280:ARG:HH22	2.10	0.53
1:B:357:ARG:HH21	1:B:357:ARG:CG	2.16	0.53
1:A:459:THR:HG23	1:A:460:GLU:HG2	1.90	0.53
1:B:320:TYR:CZ	1:B:398:PRO:HG2	2.43	0.53
1:B:454:GLN:O	1:B:457:LYS:HB2	2.09	0.52
1:B:466:HIS:HB3	1:B:469:LEU:HB2	1.91	0.52
1:A:394:SER:O	1:A:397:ARG:HG2	2.10	0.52
1:A:392:ILE:HG22	1:A:393:LEU:HD12	1.92	0.52
1:A:262:ILE:HD12	1:A:264:PHE:HZ	1.75	0.51
1:B:325:ILE:HD11	1:B:391:ILE:HB	1.93	0.51
1:A:357:ARG:HH12	1:A:460:GLU:CD	2.14	0.50
1:B:363:PHE:CE2	1:B:364:MET:HG2	2.46	0.50
1:B:262:ILE:H	1:B:262:ILE:HD12	1.76	0.50
1:A:259:GLU:O	1:A:262:ILE:O	2.30	0.50
1:A:439:MET:O	1:A:443:ARG:HG3	2.12	0.50
1:A:433:ALA:O	1:A:437:GLN:HG3	2.11	0.49
1:A:249:ILE:CD1	1:A:255:LEU:N	2.75	0.49
1:A:441:ASP:O	1:A:445:ILE:HG12	2.13	0.49
1:A:325:ILE:HG12	1:A:388:ILE:HG23	1.94	0.48
1:A:424:ASN:HB3	1:A:425:HIS:HD2	1.78	0.48
1:A:249:ILE:HD11	1:A:255:LEU:HB2	1.94	0.48
1:A:393:LEU:HD23	1:A:409:ILE:HB	1.94	0.48
1:A:447:THR:O	1:A:451:GLN:HG3	2.13	0.48
1:B:359:PRO:HG2	1:B:456:ILE:CD1	2.38	0.48
1:A:237:LEU:HD21	1:A:340:LEU:HG	1.95	0.48
1:B:289:SER:O	1:B:293:VAL:HG23	2.14	0.48
1:A:279:ILE:HD13	1:A:360:PHE:CZ	2.49	0.48
1:B:348:MET:HG2	1:B:353:LEU:HG	1.96	0.48
1:A:421:LEU:HD11	1:A:435:LEU:HD23	1.97	0.47
1:B:266:HIS:O	1:B:267:ILE:HG22	2.15	0.47
1:A:359:PRO:HG2	1:A:360:PHE:CD1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:456:ILE:O	1:B:459:THR:HG23	2.15	0.47
1:A:249:ILE:HD13	1:A:255:LEU:HA	1.96	0.46
1:B:342:SER:O	1:B:342:SER:OG	2.29	0.46
1:B:357:ARG:NH2	1:B:357:ARG:CG	2.73	0.46
1:B:421:LEU:HD22	1:B:431:LEU:HD23	1.97	0.46
1:B:403:VAL:HG12	1:B:407:GLU:HG3	1.98	0.45
1:A:277:VAL:HG13	1:A:278:ALA:N	2.31	0.45
1:B:216:LYS:CD	1:B:220:ASP:OD2	2.65	0.45
1:A:249:ILE:CD1	1:A:255:LEU:HB2	2.46	0.45
1:B:469:LEU:HA	1:B:469:LEU:HD12	1.68	0.44
1:A:350:ARG:NH2	1:A:365:GLU:OE2	2.38	0.44
1:A:262:ILE:HD12	1:A:264:PHE:CZ	2.53	0.44
1:B:223:ILE:HD13	1:B:229:THR:HG21	1.99	0.44
1:A:452:LEU:O	1:A:456:ILE:HG13	2.19	0.43
1:A:212:ARG:HA	1:A:212:ARG:HD2	1.83	0.43
1:B:430:GLN:HG3	1:B:433:ALA:HB3	2.00	0.43
1:A:289:SER:O	1:A:293:VAL:HG23	2.19	0.43
1:A:249:ILE:HD12	1:A:254:SER:OG	2.19	0.43
1:B:259:GLU:OE2	1:B:280:ARG:NH2	2.52	0.43
1:A:288:ARG:NH2	1:A:342:SER:HA	2.34	0.43
1:A:424:ASN:HB3	1:A:425:HIS:CD2	2.53	0.43
1:B:318:LEU:HA	1:B:318:LEU:HD23	1.70	0.42
1:A:290:VAL:HG21	1:A:466:HIS:CG	2.55	0.42
1:A:405:PRO:O	1:A:409:ILE:HG13	2.20	0.42
1:B:253:ASN:O	1:B:257:MET:N	2.53	0.42
1:A:400:LEU:HD13	1:A:406:ILE:HD12	2.02	0.42
1:B:384:LEU:O	1:B:388:ILE:HG23	2.20	0.42
1:A:330:LEU:HD21	1:A:364:MET:HE1	2.02	0.41
1:B:259:GLU:OE2	1:B:280:ARG:CZ	2.68	0.41
1:A:307:VAL:HG22	2:A:512:HOH:O	2.20	0.41
1:B:263:LYS:O	1:B:264:PHE:HD2	2.04	0.41
1:A:328:THR:HG21	1:A:388:ILE:HD11	2.03	0.41
1:A:363:PHE:CE1	1:A:452:LEU:HB3	2.55	0.41
1:B:334:MET:HE3	1:B:364:MET:HE3	2.02	0.41
1:B:341:ILE:HA	1:B:341:ILE:HD13	1.86	0.40
1:B:329:MET:O	1:B:332:SER:HB2	2.21	0.40
1:B:402:ASN:O	1:B:405:PRO:HD2	2.20	0.40
1:B:451:GLN:O	1:B:455:VAL:HG23	2.21	0.40
1:B:456:ILE:O	1:B:456:ILE:CG2	2.70	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	248/275 (90%)	241 (97%)	7 (3%)	0	100	100
1	B	243/275 (88%)	239 (98%)	3 (1%)	1 (0%)	34	42
All	All	491/550 (89%)	480 (98%)	10 (2%)	1 (0%)	47	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	358	LYS

### 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	228/248 (92%)	224 (98%)	4 (2%)	59	75
1	B	224/248 (90%)	217 (97%)	7 (3%)	40	55
All	All	452/496 (91%)	441 (98%)	11 (2%)	49	66

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	244	LYS
1	A	262	ILE
1	A	362	ASP
1	A	363	PHE
1	B	210	ASP

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Mol	Chain	Res	Type
1	B	251	ASP
1	B	259	GLU
1	B	363	PHE
1	B	441	ASP
1	B	459	THR
1	B	475	ASP

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

Mol	Chain	Res	Type
1	A	430	GLN
1	A	451	GLN
1	A	470	GLN

### 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	254/275 (92%)	0.34	15 (5%) 22 28	33, 53, 78, 109	0
1	B	251/275 (91%)	0.34	21 (8%) 11 15	34, 50, 85, 96	0
All	All	505/550 (91%)	0.34	36 (7%) 16 21	33, 51, 84, 109	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	456	ILE	7.2
1	A	261	LYS	5.3
1	A	363	PHE	5.0
1	B	264	PHE	4.7
1	B	455	VAL	4.6
1	B	458	LYS	4.0
1	A	260	ASP	3.7
1	B	263	LYS	3.7
1	A	358	LYS	3.7
1	B	358	LYS	3.5
1	B	454	GLN	3.4
1	A	262	ILE	3.4
1	B	235	ALA	3.4
1	B	459	THR	3.4
1	A	249	ILE	3.2
1	B	260	ASP	3.0
1	B	256	MET	2.9
1	A	263	LYS	2.8
1	B	265	LYS	2.5
1	A	247	PHE	2.4
1	A	475	ASP	2.3
1	A	361	GLY	2.3
1	B	266	HIS	2.3
1	B	427	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	423	LEU	2.3
1	A	264	PHE	2.3
1	A	250	TYR	2.2
1	A	439	MET	2.2
1	B	360	PHE	2.2
1	B	287	PHE	2.1
1	B	363	PHE	2.1
1	A	423	LEU	2.1
1	B	261	LYS	2.1
1	B	236	ILE	2.0
1	B	451	GLN	2.0
1	A	459	THR	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.