



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:51 PM GMT

PDB ID : 3SCI  
Title : Crystal structure of spike protein receptor-binding domain from a predicted SARS coronavirus human strain complexed with human receptor ACE2  
Authors : Wu, K.; Peng, G.; Wilken, M.; Geraghty, R.; Li, F.  
Deposited on : 2011-06-07  
Resolution : 2.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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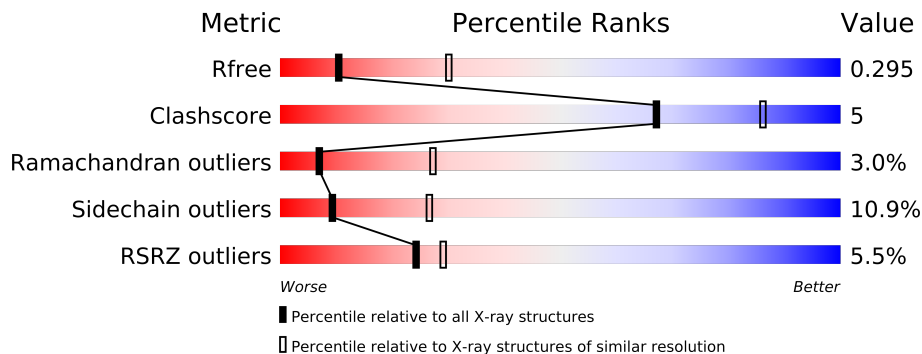
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.90 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))
$R_{free}$	66092	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	603	
1	B	603	
2	E	228	
2	F	228	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	ZN	B	901	-	X
4	CL	A	902	-	X
4	CL	B	902	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12552 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Angiotensin-convertingenzyme 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	597	Total	C	N	O	S	0	0	0
			4870	3115	806	920	29			
1	B	597	Total	C	N	O	S	0	0	0
			4870	3115	806	920	29			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	616	HIS	-	EXPRESSION TAG	UNP Q9BYF1
A	617	HIS	-	EXPRESSION TAG	UNP Q9BYF1
A	618	HIS	-	EXPRESSION TAG	UNP Q9BYF1
A	619	HIS	-	EXPRESSION TAG	UNP Q9BYF1
A	620	HIS	-	EXPRESSION TAG	UNP Q9BYF1
A	621	HIS	-	EXPRESSION TAG	UNP Q9BYF1
B	616	HIS	-	EXPRESSION TAG	UNP Q9BYF1
B	617	HIS	-	EXPRESSION TAG	UNP Q9BYF1
B	618	HIS	-	EXPRESSION TAG	UNP Q9BYF1
B	619	HIS	-	EXPRESSION TAG	UNP Q9BYF1
B	620	HIS	-	EXPRESSION TAG	UNP Q9BYF1
B	621	HIS	-	EXPRESSION TAG	UNP Q9BYF1

- Molecule 2 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	174	Total	C	N	O	S	0	0	0
			1404	912	228	257	7			
2	F	174	Total	C	N	O	S	0	0	0
			1404	912	228	257	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	442	PHE	TYR	CONFLICT	UNP P59594
E	472	PHE	LEU	CONFLICT	UNP P59594
E	528	HIS	-	EXPRESSION TAG	UNP P59594
E	529	HIS	-	EXPRESSION TAG	UNP P59594
E	530	HIS	-	EXPRESSION TAG	UNP P59594
E	531	HIS	-	EXPRESSION TAG	UNP P59594
E	532	HIS	-	EXPRESSION TAG	UNP P59594
E	533	HIS	-	EXPRESSION TAG	UNP P59594
F	442	PHE	TYR	CONFLICT	UNP P59594
F	472	PHE	LEU	CONFLICT	UNP P59594
F	528	HIS	-	EXPRESSION TAG	UNP P59594
F	529	HIS	-	EXPRESSION TAG	UNP P59594
F	530	HIS	-	EXPRESSION TAG	UNP P59594
F	531	HIS	-	EXPRESSION TAG	UNP P59594
F	532	HIS	-	EXPRESSION TAG	UNP P59594
F	533	HIS	-	EXPRESSION TAG	UNP P59594

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

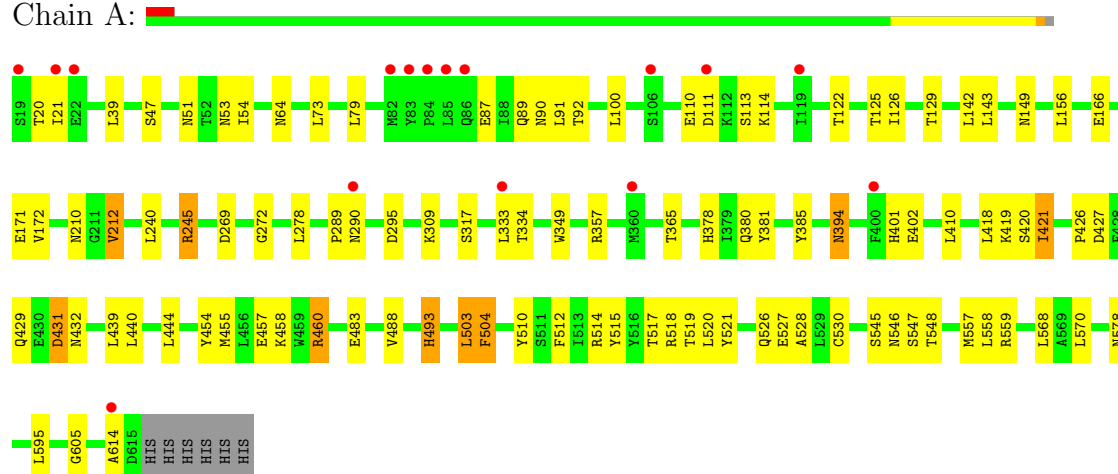
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0
4	A	1	Total Cl 1 1	0	0

### 3 Residue-property plots

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 ( $RSRZ > 2$ ). 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.

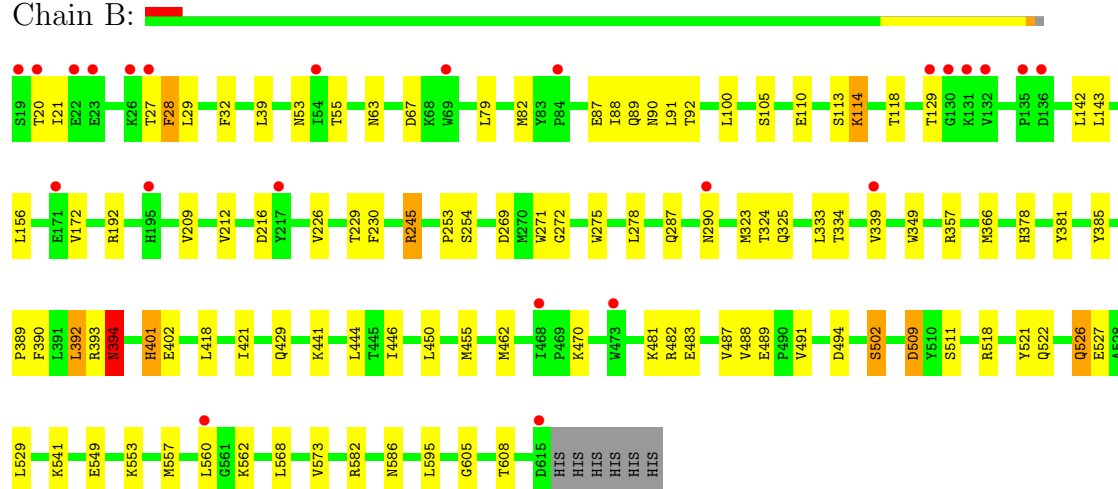
#### • Molecule 1: Angiotensin-convertingenzyme 2

Chain A:



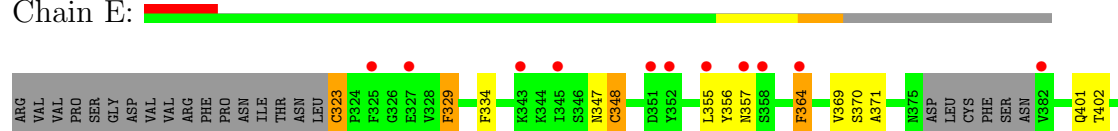
#### • Molecule 1: Angiotensin-convertingenzyme 2

Chain B:



#### • Molecule 2: Spike glycoprotein

Chain E:





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.36Å 118.33Å 111.94Å 90.00° 93.06° 90.00°	Depositor
Resolution (Å)	47.24 – 2.90 47.24 – 2.90	Depositor EDS
% Data completeness (in resolution range)	85.5 (47.24-2.90) 84.9 (47.24-2.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.54 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.226 , 0.283 0.241 , 0.295	Depositor DCC
$R_{free}$ test set	2003 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	77.7	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 30.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 46704 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	12552	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.50	0/5007	0.62	0/6803
1	B	0.52	1/5007 (0.0%)	0.62	0/6803
2	E	0.52	0/1450	0.66	0/1975
2	F	0.49	0/1450	0.61	0/1975
All	All	0.51	1/12914 (0.0%)	0.62	0/17556

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	105	SER	CB-OG	7.02	1.51	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	290	ASN	Peptide



## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4870	0	0	20	0
1	B	4870	0	0	27	0
2	E	1404	0	0	9	0
2	F	1404	0	0	7	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
All	All	12552	0	0	63	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (63) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:457:GLU:OE2	1:A:512:PHE:N	2.26	0.69
1:B:245:ARG:NH1	1:B:605:GLY:O	2.27	0.68
1:A:269:ASP:OD2	1:A:272:GLY:N	2.31	0.63
1:B:271:TRP:NE1	1:B:502:SER:O	2.32	0.62
1:B:489:GLU:OE1	1:B:489:GLU:N	2.34	0.61
1:A:378:HIS:CE1	1:A:402:GLU:OE1	2.55	0.60
1:B:90:ASN:O	1:B:92:THR:N	2.35	0.59
1:A:460:ARG:NH2	1:A:510:TYR:O	2.37	0.57
2:E:472:PHE:O	2:E:473:ASN:CB	2.54	0.55
2:F:471:ALA:O	2:F:473:ASN:N	2.39	0.55
1:B:393:ARG:O	1:B:394:ASN:CB	2.55	0.55
2:E:484:TYR:CD1	2:E:484:TYR:N	2.75	0.54
1:B:389:PRO:O	1:B:392:LEU:N	2.42	0.53
2:F:471:ALA:O	2:F:474:CYS:N	2.42	0.53
1:A:125:THR:O	1:A:129:THR:N	2.43	0.52
1:A:90:ASN:O	1:A:92:THR:N	2.43	0.51
1:A:166:GLU:OE1	1:A:493:HIS:CE1	2.63	0.51
1:B:509:ASP:C	1:B:509:ASP:OD1	2.49	0.51
1:A:454:TYR:OH	1:A:458:LYS:NZ	2.44	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:435:ASN:OD1	2:F:436:TYR:N	2.44	0.51
1:A:419:LYS:O	1:A:421:ILE:N	2.44	0.50
1:B:349:TRP:N	1:B:357:ARG:O	2.44	0.50
2:F:323:CYS:N	2:F:348:CYS:SG	2.84	0.50
1:A:526:GLN:O	1:A:530:CYS:SG	2.71	0.48
2:F:472:PHE:O	2:F:473:ASN:CB	2.62	0.48
2:E:364:PHE:CD2	2:E:364:PHE:O	2.66	0.48
1:A:245:ARG:NH1	1:A:605:GLY:O	2.48	0.47
1:B:378:HIS:CE1	1:B:402:GLU:OE1	2.68	0.47
1:B:389:PRO:O	1:B:390:PHE:C	2.54	0.47
1:B:481:LYS:O	1:B:487:VAL:N	2.48	0.46
2:E:423:TRP:N	2:E:423:TRP:CD1	2.84	0.46
1:B:253:PRO:O	1:B:254:SER:OG	2.34	0.46
1:B:269:ASP:OD2	1:B:272:GLY:N	2.49	0.46
1:A:111:ASP:O	1:A:113:SER:N	2.49	0.46
1:A:503:LEU:O	1:A:504:PHE:C	2.54	0.46
1:B:494:ASP:OD2	1:B:494:ASP:N	2.47	0.46
1:B:226:VAL:O	1:B:230:PHE:N	2.50	0.45
1:A:545:SER:O	1:A:547:SER:N	2.50	0.44
1:A:517:THR:O	1:A:518:ARG:C	2.56	0.44
1:B:114:LYS:O	1:B:118:THR:OG1	2.35	0.44
1:B:526:GLN:O	1:B:527:GLU:C	2.55	0.43
1:B:323:MET:CA	1:B:323:MET:CE	2.96	0.43
2:E:416:PHE:O	2:E:417:MET:CB	2.66	0.43
1:B:275:TRP:N	1:B:275:TRP:CD1	2.84	0.43
2:F:488:GLY:O	2:F:491:TYR:N	2.52	0.43
1:B:245:ARG:CG	1:B:245:ARG:NH2	2.81	0.43
1:A:514:ARG:CG	1:A:515:TYR:N	2.81	0.42
1:B:366:MET:CE	1:B:441:LYS:CE	2.98	0.42
1:B:378:HIS:ND1	1:B:401:HIS:ND1	2.67	0.42
2:E:329:PHE:N	2:E:329:PHE:CD1	2.87	0.42
2:E:323:CYS:N	2:E:348:CYS:SG	2.93	0.42
2:F:325:PHE:C	2:F:327:GLU:N	2.73	0.42
1:A:527:GLU:O	1:A:528:ALA:C	2.57	0.42
1:A:431:ASP:N	1:A:431:ASP:OD1	2.52	0.42
1:B:63:ASN:O	1:B:67:ASP:N	2.52	0.41
1:B:324:THR:O	1:B:325:GLN:C	2.58	0.41
1:A:349:TRP:N	1:A:357:ARG:O	2.53	0.41
2:E:487:THR:O	2:E:492:GLN:NE2	2.53	0.41
1:B:28:PHE:O	1:B:32:PHE:N	2.53	0.41
1:B:549:GLU:CD	1:B:549:GLU:N	2.74	0.41
1:B:521:TYR:O	1:B:522:GLN:C	2.60	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:471:ALA:O	2:E:472:PHE:C	2.60	0.40
1:A:520:LEU:O	1:A:521:TYR:C	2.60	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	595/603 (99%)	506 (85%)	75 (13%)	14 (2%)	9	35
1	B	595/603 (99%)	516 (87%)	68 (11%)	11 (2%)	13	44
2	E	170/228 (75%)	131 (77%)	28 (16%)	11 (6%)	2	5
2	F	170/228 (75%)	135 (79%)	25 (15%)	10 (6%)	2	7
All	All	1530/1662 (92%)	1288 (84%)	196 (13%)	46 (3%)	7	27

All (46) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	LEU
1	A	212	VAL
1	B	91	LEU
1	B	394	ASN
1	B	421	ILE
2	E	402	THR
2	E	415	ASP
2	E	416	PHE
2	E	471	ALA
2	E	472	PHE
2	E	473	ASN
2	F	415	ASP
2	F	472	PHE
2	F	473	ASN
1	A	54	ILE
1	A	289	PRO

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Mol	Chain	Res	Type
1	A	394	ASN
1	A	421	ILE
1	A	504	PHE
1	A	546	ASN
1	B	470	LYS
1	B	509	ASP
2	E	370	SER
2	E	371	ALA
2	F	374	LEU
1	A	420	SER
1	A	614	ALA
1	B	28	PHE
1	B	339	VAL
2	E	463	ASP
2	F	332	THR
2	F	401	GLN
1	A	51	ASN
1	A	64	ASN
1	A	171	GLU
1	B	113	SER
2	E	347	ASN
2	E	401	GLN
2	F	370	SER
2	F	402	THR
1	B	446	ILE
2	F	326	GLY
1	B	89	GLN
1	A	426	PRO
1	B	212	VAL
2	F	470	PRO

### 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	527/533 (99%)	468 (89%)	59 (11%)	9	25
1	B	527/533 (99%)	467 (89%)	60 (11%)	8	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	152/202 (75%)	134 (88%)	18 (12%)	8	22
2	F	152/202 (75%)	141 (93%)	11 (7%)	21	51
All	All	1358/1470 (92%)	1210 (89%)	148 (11%)	9	26

All (148) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	THR
1	A	21	ILE
1	A	39	LEU
1	A	47	SER
1	A	53	ASN
1	A	73	LEU
1	A	79	LEU
1	A	87	GLU
1	A	89	GLN
1	A	100	LEU
1	A	110	GLU
1	A	114	LYS
1	A	122	THR
1	A	126	ILE
1	A	142	LEU
1	A	143	LEU
1	A	149	ASN
1	A	156	LEU
1	A	172	VAL
1	A	210	ASN
1	A	212	VAL
1	A	240	LEU
1	A	245	ARG
1	A	278	LEU
1	A	295	ASP
1	A	309	LYS
1	A	317	SER
1	A	333	LEU
1	A	334	THR
1	A	365	THR
1	A	380	GLN
1	A	381	TYR
1	A	385	TYR
1	A	394	ASN

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Mol	Chain	Res	Type
1	A	401	HIS
1	A	410	LEU
1	A	418	LEU
1	A	427	ASP
1	A	429	GLN
1	A	431	ASP
1	A	432	ASN
1	A	439	LEU
1	A	440	LEU
1	A	444	LEU
1	A	455	MET
1	A	460	ARG
1	A	483	GLU
1	A	488	VAL
1	A	493	HIS
1	A	503	LEU
1	A	519	THR
1	A	548	THR
1	A	557	MET
1	A	558	LEU
1	A	559	ARG
1	A	568	LEU
1	A	570	LEU
1	A	578	ASN
1	A	595	LEU
1	B	20	THR
1	B	21	ILE
1	B	27	THR
1	B	29	LEU
1	B	39	LEU
1	B	53	ASN
1	B	55	THR
1	B	79	LEU
1	B	82	MET
1	B	87	GLU
1	B	88	ILE
1	B	100	LEU
1	B	110	GLU
1	B	114	LYS
1	B	129	THR
1	B	142	LEU
1	B	143	LEU

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Mol	Chain	Res	Type
1	B	156	LEU
1	B	172	VAL
1	B	192	ARG
1	B	209	VAL
1	B	216	ASP
1	B	229	THR
1	B	245	ARG
1	B	278	LEU
1	B	287	GLN
1	B	290	ASN
1	B	333	LEU
1	B	334	THR
1	B	381	TYR
1	B	385	TYR
1	B	392	LEU
1	B	394	ASN
1	B	401	HIS
1	B	418	LEU
1	B	429	GLN
1	B	444	LEU
1	B	450	LEU
1	B	455	MET
1	B	462	MET
1	B	482	ARG
1	B	483	GLU
1	B	488	VAL
1	B	491	VAL
1	B	502	SER
1	B	511	SER
1	B	518	ARG
1	B	526	GLN
1	B	529	LEU
1	B	541	LYS
1	B	553	LYS
1	B	557	MET
1	B	560	LEU
1	B	562	LYS
1	B	568	LEU
1	B	573	VAL
1	B	582	ARG
1	B	586	ASN
1	B	595	LEU

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Mol	Chain	Res	Type
1	B	608	THR
2	E	323	CYS
2	E	329	PHE
2	E	334	PHE
2	E	348	CYS
2	E	355	LEU
2	E	356	TYR
2	E	357	ASN
2	E	364	PHE
2	E	369	VAL
2	E	417	MET
2	E	423	TRP
2	E	426	ARG
2	E	427	ASN
2	E	442	PHE
2	E	467	CYS
2	E	468	THR
2	E	484	TYR
2	E	500	SER
2	F	323	CYS
2	F	328	VAL
2	F	348	CYS
2	F	355	LEU
2	F	360	PHE
2	F	364	PHE
2	F	423	TRP
2	F	427	ASN
2	F	442	PHE
2	F	468	THR
2	F	481	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains 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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	597/603 (99%)	0.20	16 (2%) 52 61	39, 70, 116, 136	0
1	B	597/603 (99%)	0.32	24 (4%) 36 43	38, 73, 119, 138	0
2	E	174/228 (76%)	0.59	19 (10%) 6 8	64, 86, 128, 137	0
2	F	174/228 (76%)	0.55	23 (13%) 4 6	64, 88, 129, 138	0
All	All	1542/1662 (92%)	0.33	82 (5%) 24 31	38, 75, 124, 138	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	442	PHE	7.0
2	E	501	PHE	6.2
2	F	443	LEU	5.7
2	F	364	PHE	5.3
1	B	136	ASP	5.2
2	E	355	LEU	5.0
2	E	352	TYR	4.9
2	E	421	LEU	4.9
1	A	86	GLN	4.9
1	B	615	ASP	4.7
2	E	382	VAL	4.6
2	E	472	PHE	4.6
1	A	85	LEU	3.8
1	B	473	TRP	3.8
1	A	19	SER	3.6
1	B	339	VAL	3.6
2	E	364	PHE	3.6
1	B	130	GLY	3.5
1	A	360	MET	3.4
2	F	418	GLY	3.4
1	A	83	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
2	F	325	PHE	3.4
2	F	354	VAL	3.3
2	E	499	LEU	3.3
2	F	419	CYS	3.2
2	E	357	ASN	3.2
1	B	26	LYS	3.2
2	F	352	TYR	3.1
2	E	443	LEU	3.0
2	F	470	PRO	3.0
1	B	135	PRO	3.0
2	F	501	PHE	2.9
1	B	27	THR	2.9
1	B	129	THR	2.9
2	E	343	LYS	2.9
2	F	365	LYS	2.9
2	F	350	ALA	2.9
1	B	22	GLU	2.8
2	F	356	TYR	2.8
2	E	410	TYR	2.8
2	F	478	LEU	2.8
1	A	111	ASP	2.7
2	E	442	PHE	2.7
1	B	19	SER	2.7
1	B	171	GLU	2.7
2	F	382	VAL	2.6
1	A	21	ILE	2.6
2	F	329	PHE	2.6
1	A	82	MET	2.6
2	E	351	ASP	2.6
1	A	84	PRO	2.6
1	A	106	SER	2.5
1	A	614	ALA	2.5
2	E	345	ILE	2.5
2	E	325	PHE	2.5
2	F	477	PRO	2.5
1	B	131	LYS	2.5
1	B	290	ASN	2.4
1	A	22	GLU	2.4
1	B	69	TRP	2.4
1	B	23	GLU	2.4
2	F	368	GLY	2.4
2	E	500	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	20	THR	2.3
2	F	359	THR	2.3
1	B	217	TYR	2.3
2	F	423	TRP	2.3
2	E	327	GLU	2.3
1	B	560	LEU	2.3
1	B	468	ILE	2.2
1	A	333	LEU	2.2
1	B	84	PRO	2.2
2	E	358	SER	2.2
1	A	290	ASN	2.2
2	F	468	THR	2.2
2	F	421	LEU	2.2
1	A	400	PHE	2.1
1	B	195	HIS	2.1
1	B	54	ILE	2.0
1	B	132	VAL	2.0
1	A	119	ILE	2.0
2	F	361	PHE	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	CL	A	902	1/1	0.40	7.58	92,92,92,92	0
4	CL	B	902	1/1	0.43	7.30	107,107,107,107	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	B	901	1/1	0.45	3.86	90,90,90,90	0
3	ZN	A	901	1/1	0.36	1.95	87,87,87,87	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.