



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:55 AM GMT

PDB ID : 3S5M
Title : Crystal structures of falcilysin, a M16 metalloprotease from the malaria parasite *Plasmodium falciparum*
Authors : Morgunova, E.; Ponpuak, M.; Istvan, E.; Popov, A.; Goldberg, D.; Eneqvist, T.
Deposited on : 2011-05-23
Resolution : 1.55 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

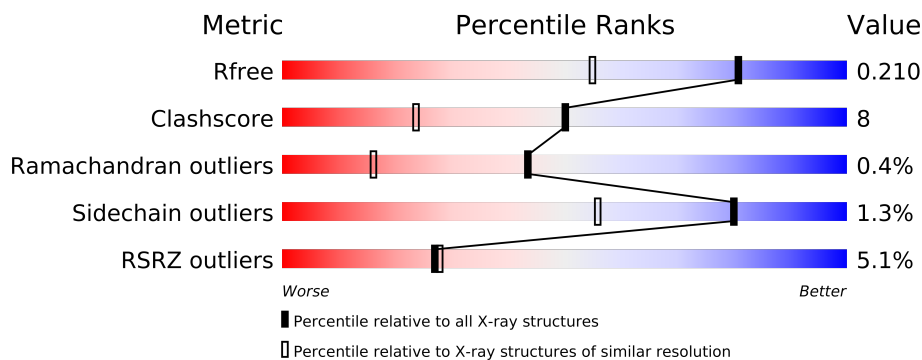
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.55 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	1117 (1.58-1.54)
Clashscore	79885	1249 (1.58-1.54)
Ramachandran outliers	78287	1212 (1.58-1.54)
Sidechain outliers	78261	1210 (1.58-1.54)
RSRZ outliers	66119	1117 (1.58-1.54)

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 ≥ 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	1193	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19225 atoms, of which 9022 are hydrogens 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 Falcilysin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	1094	Total	C	H	N	O	S	0	24	0
			18119	5852	9022	1477	1734	34			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is water.

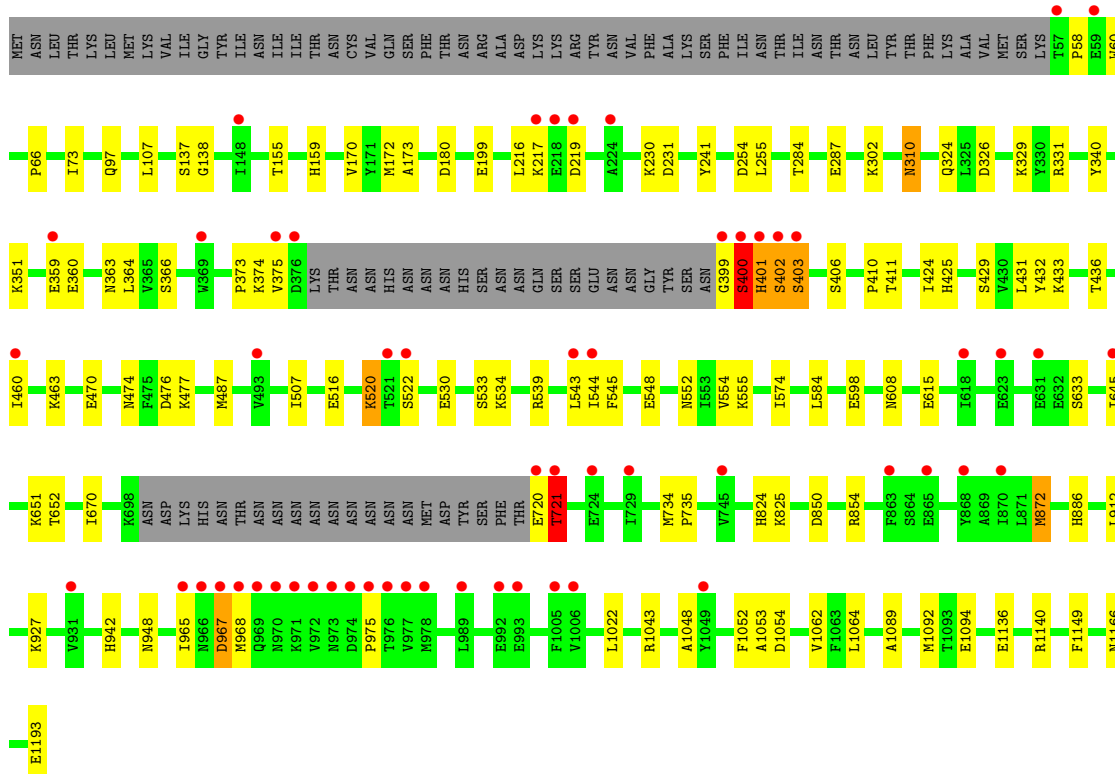
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1103	Total	O	0	0
			1103	1103		

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: Falcilysin

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.32Å 106.70Å 128.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.20 – 1.55 29.53 – 1.55	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.20-1.55) 99.4 (29.53-1.55)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 1.55Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.179 , 0.214 0.176 , 0.210	Depositor DCC
R_{free} test set	3715 reflections (2.00%)	DCC
Wilson B-factor (Å ²)	18.1	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 185784 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19225	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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, MG

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	0.70	2/9351 (0.0%)	0.78	5/12594 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	400	SER	CA-CB	-8.26	1.40	1.52
1	A	401	HIS	CA-CB	-8.11	1.36	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	539	ARG	NE-CZ-NH2	-9.56	115.52	120.30
1	A	854	ARG	NE-CZ-NH2	-8.90	115.85	120.30
1	A	975	PRO	N-CA-C	5.79	127.14	112.10
1	A	854	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	A	721	THR	N-CA-C	5.19	125.02	111.00

There are no chirality outliers.

There are no planarity outliers.

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	9097	9022	81	137	0
2	A	1	0	0	0	0
3	A	2	0	0	0	0
4	A	1103	0	0	43	0
All	All	10203	9022	81	137	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 8.

All (137) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:374:LYS:HA	1:A:402:SER:CA	1.84	1.07
1:A:374:LYS:CA	1:A:402:SER:HA	1.87	1.05
1:A:533:SER:C	4:A:2179:HOH:O	2.00	0.99
1:A:324[A]:GLN:HG3	4:A:2194:HOH:O	1.63	0.97
1:A:425:HIS:O	1:A:429[B]:SER:OG	1.88	0.92
1:A:329:LYS:HG2	4:A:2211:HOH:O	1.72	0.89
1:A:402:SER:CB	4:A:1993:HOH:O	2.22	0.87
1:A:374:LYS:HA	1:A:402:SER:HA	0.92	0.87
1:A:436[B]:THR:HG21	4:A:1309:HOH:O	1.74	0.87
1:A:137:SER:C	4:A:2176:HOH:O	2.13	0.85
1:A:608:ASN:CG	4:A:2275:HOH:O	2.15	0.83
1:A:254:ASP:HB2	4:A:2087:HOH:O	1.80	0.82
1:A:410:PRO:HB3	1:A:543:LEU:HD11	1.62	0.81
1:A:180:ASP:OD1	4:A:2208:HOH:O	1.98	0.81
1:A:399:GLY:O	1:A:400:SER:CB	2.22	0.80
1:A:544:ILE:HB	4:A:1638:HOH:O	1.83	0.79
1:A:364[B]:LEU:HD22	1:A:463:LYS:CB	2.12	0.79
1:A:410:PRO:HB3	1:A:543:LEU:CD1	2.14	0.78
1:A:172[B]:MET:HG3	1:A:173:ALA:N	1.96	0.78
1:A:375:VAL:CG1	1:A:403:SER:CB	2.61	0.78
1:A:401:HIS:O	1:A:406:SER:HB3	1.83	0.77
1:A:850:ASP:OD2	4:A:2191:HOH:O	2.03	0.76
1:A:326:ASP:O	4:A:2211:HOH:O	2.04	0.75
1:A:530:GLU:HG2	1:A:544:ILE:HG12	1.70	0.74
1:A:522:SER:HA	4:A:1392:HOH:O	1.86	0.74
1:A:138:GLY:N	4:A:2176:HOH:O	2.21	0.74
1:A:375:VAL:HG11	1:A:403:SER:CB	2.18	0.73
1:A:1094:GLU:OE1	4:A:1580:HOH:O	2.05	0.73
1:A:364[B]:LEU:HD22	1:A:463:LYS:HB2	1.74	0.69
1:A:375:VAL:HG13	1:A:403:SER:CB	2.23	0.68
1:A:359:GLU:O	4:A:2205:HOH:O	2.13	0.67
1:A:544:ILE:HD12	1:A:544:ILE:C	2.16	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:824[A]:HIS:ND1	4:A:2180:HOH:O	2.29	0.66
1:A:255:LEU:HD21	1:A:366[A]:SER:HB2	1.78	0.65
1:A:216:LEU:HD12	1:A:231:ASP:HA	1.78	0.65
1:A:255:LEU:HD21	1:A:366[B]:SER:HB3	1.78	0.65
1:A:470:GLU:OE1	4:A:2268:HOH:O	2.14	0.65
1:A:552:ASN:OD1	4:A:2207:HOH:O	2.15	0.65
1:A:399:GLY:O	4:A:1315:HOH:O	2.15	0.64
1:A:254:ASP:OD2	4:A:2289:HOH:O	2.15	0.63
1:A:159:HIS:HE2	1:A:172[B]:MET:CG	2.12	0.62
1:A:363:ASN:C	1:A:364[B]:LEU:HD23	2.22	0.60
1:A:373:PRO:O	1:A:402:SER:CB	2.49	0.60
1:A:137:SER:HA	4:A:2176:HOH:O	2.03	0.59
1:A:230:LYS:HE3	1:A:615:GLU:HG3	1.84	0.59
1:A:324[A]:GLN:CG	4:A:2194:HOH:O	2.33	0.59
1:A:431:LEU:HB3	1:A:460:ILE:CD1	2.33	0.59
1:A:402:SER:C	4:A:2089:HOH:O	2.40	0.58
1:A:402:SER:CB	4:A:2089:HOH:O	2.51	0.58
1:A:137:SER:CA	4:A:2176:HOH:O	2.50	0.58
1:A:824[B]:HIS:CE1	1:A:942:HIS:CE1	2.92	0.57
1:A:824[A]:HIS:CD2	1:A:825:LYS:HG2	2.39	0.57
1:A:363:ASN:O	1:A:364[B]:LEU:HD23	2.04	0.57
1:A:734:MET:HE2	4:A:2225:HOH:O	2.05	0.57
1:A:927:LYS:HD2	1:A:965:ILE:HD12	1.87	0.56
1:A:340:TYR:H	1:A:399:GLY:N	2.04	0.56
1:A:436[B]:THR:CG2	4:A:2270:HOH:O	2.55	0.54
1:A:967:ASP:O	1:A:968:MET:HB2	2.07	0.53
1:A:436[B]:THR:HG23	4:A:2270:HOH:O	2.09	0.53
1:A:403:SER:N	4:A:2089:HOH:O	2.42	0.52
1:A:872:MET:HE1	1:A:1052[A]:PHE:CD1	2.44	0.52
1:A:364[B]:LEU:HD22	1:A:463:LYS:HB3	1.92	0.52
1:A:155:THR:HA	4:A:2208:HOH:O	2.08	0.52
1:A:255:LEU:HD22	1:A:364[A]:LEU:HD13	1.91	0.51
1:A:401:HIS:O	1:A:402:SER:C	2.48	0.51
1:A:375:VAL:HG22	1:A:402:SER:O	2.11	0.51
1:A:1089:ALA:HA	1:A:1092:MET:HE3	1.92	0.51
1:A:720:GLU:HG3	1:A:721:THR:N	2.25	0.51
1:A:872:MET:HE3	1:A:1064:LEU:HD22	1.93	0.51
1:A:170:VAL:HG12	1:A:172[A]:MET:CG	2.41	0.50
1:A:670:ILE:HD12	1:A:948:ASN:ND2	2.25	0.50
1:A:431:LEU:HB3	1:A:460:ILE:HD13	1.92	0.50
1:A:66:PRO:HG3	1:A:310[A]:ASN:ND2	2.27	0.50
1:A:720:GLU:HG3	1:A:721:THR:H	1.78	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:401:HIS:O	1:A:406:SER:CB	2.57	0.49
1:A:516:GLU:O	1:A:520:LYS:HD3	2.12	0.49
1:A:487:MET:HE1	1:A:574:ILE:HA	1.94	0.49
1:A:433:LYS:HA	1:A:436[B]:THR:HG22	1.95	0.49
1:A:159:HIS:NE2	1:A:172[B]:MET:CG	2.76	0.49
1:A:331:ARG:HD2	4:A:1974:HOH:O	2.12	0.49
1:A:411:THR:HG23	1:A:554:VAL:CG2	2.43	0.49
1:A:324[A]:GLN:CD	4:A:2194:HOH:O	2.51	0.49
1:A:402:SER:CA	4:A:2089:HOH:O	2.62	0.48
1:A:402:SER:CB	1:A:406:SER:HB3	2.43	0.48
1:A:824[A]:HIS:HD2	1:A:825:LYS:HG2	1.78	0.48
1:A:1022:LEU:C	1:A:1022:LEU:HD23	2.33	0.48
1:A:507:ILE:HG21	1:A:555:LYS:HE2	1.95	0.48
1:A:199:GLU:OE1	1:A:199:GLU:HA	2.14	0.48
1:A:720:GLU:O	1:A:721:THR:CB	2.62	0.47
1:A:255:LEU:HD22	1:A:364[A]:LEU:CD1	2.44	0.47
1:A:872:MET:HE1	1:A:1052[A]:PHE:HD1	1.79	0.47
1:A:886:HIS:HE1	4:A:2039:HOH:O	1.98	0.47
1:A:645:ILE:HG22	4:A:1414:HOH:O	2.13	0.47
1:A:402:SER:CB	1:A:406:SER:CA	2.92	0.47
1:A:824[B]:HIS:ND1	1:A:942:HIS:CE1	2.83	0.47
1:A:608:ASN:OD1	4:A:2275:HOH:O	2.17	0.47
1:A:1052[A]:PHE:CZ	1:A:1054:ASP:HB2	2.50	0.46
1:A:107:LEU:HD11	1:A:735:PRO:HG2	1.97	0.46
1:A:364[B]:LEU:CD2	1:A:463:LYS:HB2	2.42	0.46
1:A:1140:ARG:HA	1:A:1140:ARG:HD2	1.81	0.46
1:A:241:TYR:OH	1:A:598:GLU:OE2	2.25	0.45
1:A:401:HIS:O	1:A:403:SER:N	2.49	0.45
1:A:411:THR:HG23	1:A:554:VAL:HG22	1.99	0.45
1:A:720:GLU:O	1:A:721:THR:HB	2.16	0.45
1:A:431:LEU:CB	1:A:460:ILE:CD1	2.94	0.45
1:A:608:ASN:CB	4:A:2275:HOH:O	2.59	0.45
1:A:544:ILE:HD12	1:A:545:PHE:N	2.31	0.44
1:A:734:MET:CE	4:A:2225:HOH:O	2.64	0.44
1:A:651:LYS:HG2	1:A:652:THR:O	2.17	0.44
1:A:284:THR:OG1	1:A:287:GLU:HG3	2.17	0.44
1:A:351:LYS:NZ	1:A:476:ASP:OD1	2.50	0.44
1:A:97:GLN:HB2	1:A:302:LYS:HD3	2.00	0.44
1:A:1053:ALA:HA	1:A:1062:VAL:O	2.18	0.43
1:A:474:ASN:ND2	1:A:477:LYS:HG2	2.34	0.43
1:A:1092:MET:HE1	1:A:1149:PHE:CE2	2.53	0.43
1:A:375:VAL:HG13	1:A:402:SER:O	2.19	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:912:LEU:C	1:A:912:LEU:HD23	2.39	0.43
1:A:402:SER:CB	1:A:406:SER:HA	2.48	0.43
1:A:402:SER:O	1:A:403:SER:CB	2.66	0.43
1:A:530:GLU:HG2	1:A:544:ILE:CG1	2.43	0.42
1:A:534:LYS:N	4:A:2179:HOH:O	2.38	0.42
1:A:1043:ARG:HG2	1:A:1048:ALA:O	2.19	0.42
1:A:329:LYS:CB	4:A:2211:HOH:O	2.67	0.42
1:A:1043:ARG:HD3	1:A:1043:ARG:C	2.40	0.42
1:A:1136[B]:GLU:HG2	4:A:2223:HOH:O	2.20	0.41
1:A:670:ILE:HD12	1:A:948:ASN:HD21	1.86	0.41
1:A:73:ILE:HD11	1:A:73:ILE:HG21	1.77	0.41
1:A:720:GLU:O	1:A:721:THR:HG22	2.20	0.41
1:A:58:PRO:HB2	1:A:60:TRP:CD1	2.56	0.41
1:A:533:SER:HB2	4:A:2179:HOH:O	2.21	0.41
1:A:170:VAL:HG12	1:A:172[A]:MET:HG2	2.01	0.41
1:A:548:GLU:HA	1:A:548:GLU:OE1	2.20	0.41
1:A:424:ILE:HD13	1:A:460:ILE:CD1	2.51	0.40
1:A:824[B]:HIS:CE1	1:A:942:HIS:HE1	2.39	0.40
1:A:424:ILE:HD13	1:A:460:ILE:HD12	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	1112/1193 (93%)	1077 (97%)	31 (3%)	4 (0%)	43 16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	400	SER
1	A	403	SER
1	A	721	THR
1	A	402	SER

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	1029/1104 (93%)	1015 (99%)	14 (1%)	78 52

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	217	LYS
1	A	219	ASP
1	A	310[A]	ASN
1	A	310[B]	ASN
1	A	360	GLU
1	A	432	TYR
1	A	520	LYS
1	A	584	LEU
1	A	633	SER
1	A	721	THR
1	A	872	MET
1	A	967	ASP
1	A	1166	ASN
1	A	1193	GLU

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 3 ligands modelled in this entry, 3 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, 95th 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 ²)	Q<0.9
1	A	1094/1193 (91%)	0.39	56 (5%) 27 28	11, 22, 46, 78	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	402	SER	12.2
1	A	972	VAL	10.5
1	A	968	MET	9.8
1	A	969	GLN	9.0
1	A	977	VAL	8.9
1	A	965	ILE	8.9
1	A	403	SER	8.3
1	A	967	ASP	8.0
1	A	975	PRO	8.0
1	A	976	THR	7.9
1	A	966	ASN	7.2
1	A	971	LYS	6.9
1	A	399	GLY	6.7
1	A	974	ASP	6.5
1	A	375	VAL	6.3
1	A	401	HIS	6.2
1	A	721	THR	6.0
1	A	973	ASN	6.0
1	A	376	ASP	5.8
1	A	970	ASN	5.5
1	A	218	GLU	5.0
1	A	219	ASP	4.9
1	A	631	GLU	4.6
1	A	400	SER	4.3
1	A	57	THR	4.2
1	A	460	ILE	4.0
1	A	224	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	645	ILE	3.7
1	A	1049	TYR	3.5
1	A	217	LYS	3.4
1	A	1005	PHE	3.4
1	A	978	MET	3.4
1	A	720	GLU	3.4
1	A	870	ILE	3.3
1	A	493	VAL	3.1
1	A	543	LEU	3.0
1	A	992	GLU	3.0
1	A	868	TYR	2.9
1	A	989	LEU	2.8
1	A	59	GLU	2.7
1	A	148	ILE	2.7
1	A	1006	VAL	2.6
1	A	544	ILE	2.6
1	A	863	PHE	2.6
1	A	521	THR	2.5
1	A	724	GLU	2.5
1	A	993	GLU	2.4
1	A	865	GLU	2.4
1	A	369	TRP	2.4
1	A	745	VAL	2.4
1	A	359	GLU	2.4
1	A	931	VAL	2.4
1	A	618	ILE	2.3
1	A	729	ILE	2.3
1	A	522	SER	2.2
1	A	623	GLU	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, 95th 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 ²)	Q<0.9
2	ZN	A	1194	1/1	0.10	-0.42	27,27,27,27	1
3	MG	A	2000	1/1	0.05	-0.70	21,21,21,21	0
3	MG	A	2001	1/1	0.03	-2.04	31,31,31,31	0

6.5 Other polymers

There are no such residues in this entry.