



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:05 PM UTC

PDB ID : 8A60 / pdb_00008a60
Title : Crystal structure of FhuA in complex with the superinfection exclusion lipoprotein Llp
Authors : van den Berg, B.
Deposited on : 2022-06-16
Resolution : 3.37 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

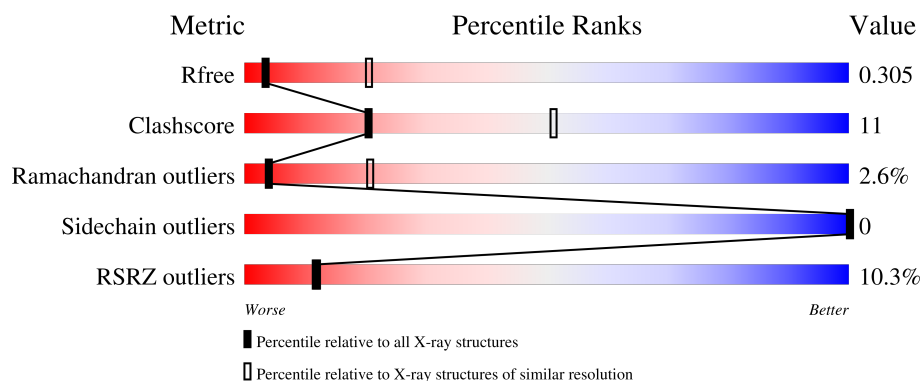
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.37 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}	180053	1047 (3.42-3.34)
Clashscore	190562	1067 (3.42-3.34)
Ramachandran outliers	187476	1056 (3.42-3.34)
Sidechain outliers	187428	1056 (3.42-3.34)
RSRZ outliers	180081	1047 (3.42-3.34)

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 ≥ 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	716	10% 70% 23% 5%
2	B	83	% 47% 28% 25%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5804 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 Ferrichrome outer membrane transporter/phage receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	677	Total	C	N	O	S	0	0	0
			5307	3347	897	1049	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP P06971
A	0	MET	-	expression tag	UNP P06971

- Molecule 2 is a protein called Lytic conversion lipoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	62	Total	C	N	O	S	0	0	0
			497	324	76	91	6			

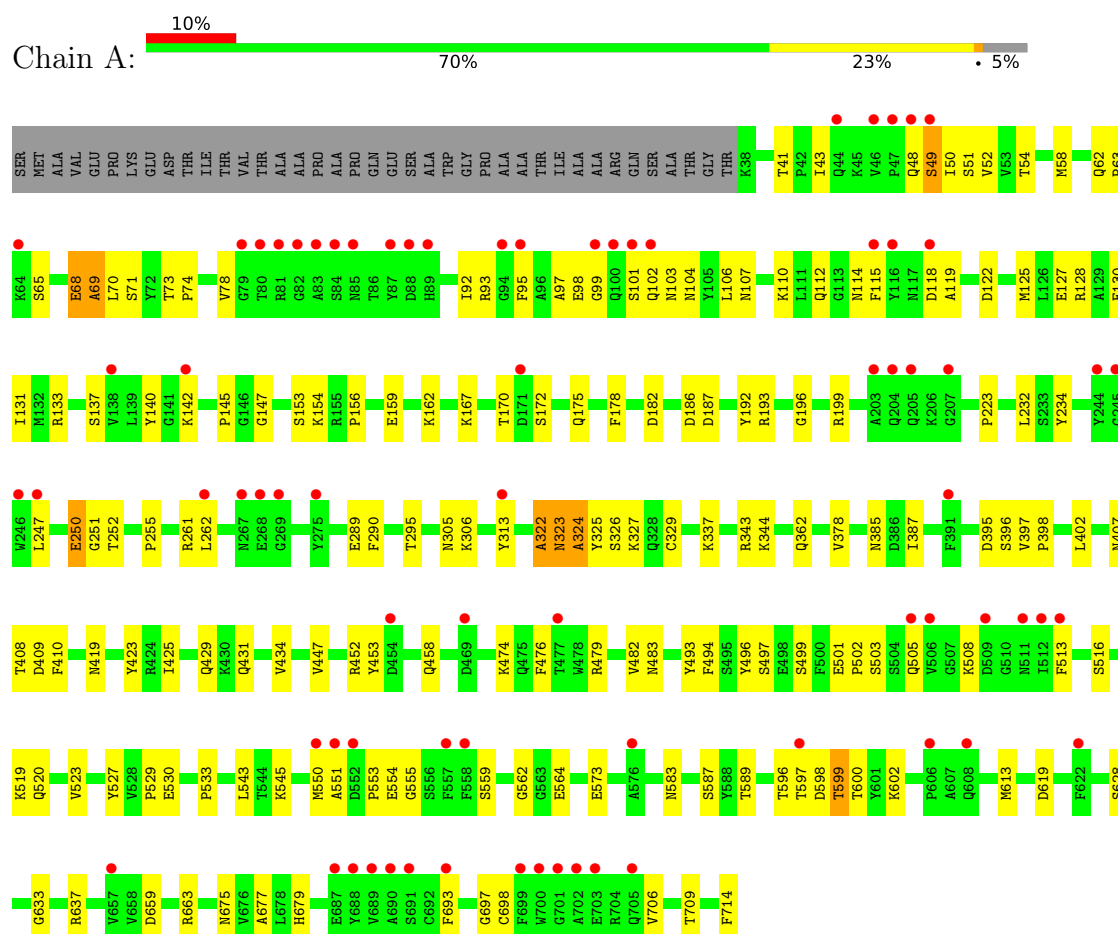
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	63	HIS	-	expression tag	UNP Q38162
B	64	HIS	-	expression tag	UNP Q38162
B	65	HIS	-	expression tag	UNP Q38162
B	66	HIS	-	expression tag	UNP Q38162
B	67	HIS	-	expression tag	UNP Q38162
B	68	HIS	-	expression tag	UNP Q38162

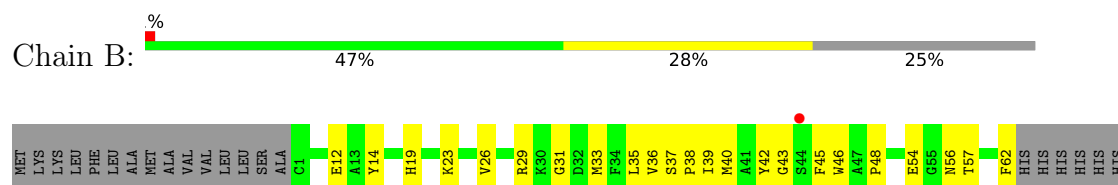
3 Residue-property plots [i](#)

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: Ferrichrome outer membrane transporter/phage receptor



- Molecule 2: Lytic conversion lipoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.29Å 98.29Å 236.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	57.85 – 3.37 57.85 – 3.37	Depositor EDS
% Data completeness (in resolution range)	77.5 (57.85-3.37) 78.1 (57.85-3.37)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 3.41Å)	Xtriage
Refinement program	PHENIX (1.20rc3_4406: ???)	Depositor
R, R_{free}	0.253 , 0.307 0.254 , 0.305	Depositor DCC
R_{free} test set	747 reflections (3.86%)	wwPDB-VP
Wilson B-factor (Å ²)	95.0	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 55.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	5804	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	0.11	0/5438	0.30	0/7387
2	B	0.10	0/513	0.30	0/692
All	All	0.11	0/5951	0.30	0/8079

There are no bond length outliers.

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	5307	0	5036	119	0
2	B	497	0	465	14	0
All	All	5804	0	5501	126	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 11.

All (126) 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:A:564:GLU:HG3	1:A:596:THR:HB	1.69	0.74
1:A:429:GLN:HG2	1:A:458:GLN:HG2	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ASP:OD2	1:A:199:ARG:NH2	2.24	0.71
1:A:378:VAL:HG12	1:A:434:VAL:HG12	1.74	0.70
1:A:186:ASP:OD1	1:A:187:ASP:N	2.25	0.69
1:A:323:ASN:O	1:A:325:TYR:N	2.28	0.66
1:A:431:GLN:HE22	1:A:502:PRO:HG3	1.61	0.66
1:A:262:LEU:HD21	1:A:402:LEU:HD12	1.78	0.66
1:A:551:ALA:HA	1:A:559:SER:HA	1.79	0.65
1:A:93:ARG:NH2	1:A:520:GLN:OE1	2.30	0.65
1:A:62:GLN:HE21	1:A:709:THR:HG21	1.62	0.64
1:A:178:PHE:CE1	1:A:196:GLY:HA3	2.33	0.64
1:A:329:CYS:O	1:A:337:LYS:NZ	2.30	0.64
1:A:52:VAL:HG22	1:A:130:GLU:HB3	1.81	0.62
1:A:62:GLN:OE1	1:A:679:HIS:NE2	2.31	0.62
1:A:98:GLU:HB2	1:A:103:ASN:HD22	1.66	0.61
1:A:520:GLN:HB2	1:A:543:LEU:HD13	1.83	0.60
1:A:74:PRO:HB3	1:A:587:SER:HB2	1.82	0.59
1:A:599:THR:H	1:A:602:LYS:HD3	1.68	0.59
1:A:598:ASP:O	1:A:600:THR:N	2.36	0.58
1:A:128:ARG:NH1	2:B:42:TYR:O	2.36	0.58
1:A:51:SER:HB3	1:A:133:ARG:HH12	1.69	0.58
1:A:494:PHE:HB3	1:A:523:VAL:HG12	1.85	0.57
1:A:597:THR:HA	1:A:602:LYS:HD2	1.87	0.57
1:A:159:GLU:N	1:A:159:GLU:OE1	2.37	0.57
1:A:496:TYR:OH	1:A:519:LYS:NZ	2.39	0.56
1:A:98:GLU:OE1	1:A:501:GLU:HB3	2.07	0.55
1:A:637:ARG:NH1	1:A:659:ASP:OD2	2.39	0.55
2:B:14:TYR:HB2	2:B:54:GLU:HG3	1.90	0.54
1:A:98:GLU:H	1:A:103:ASN:N	2.05	0.54
1:A:497:SER:HB2	1:A:520:GLN:HB3	1.90	0.54
1:A:137:SER:O	1:A:479:ARG:NH2	2.41	0.53
1:A:289:GLU:HB2	2:B:19:HIS:ND1	2.24	0.53
1:A:110:LYS:HZ1	1:A:142:LYS:HD3	1.74	0.53
1:A:118:ASP:OD1	1:A:118:ASP:N	2.37	0.53
1:A:589:THR:HB	1:A:613:MET:HE2	1.91	0.53
1:A:114:ASN:HD21	1:A:385:ASN:HB2	1.74	0.52
1:A:125:MET:HG3	1:A:234:TYR:HE1	1.75	0.51
1:A:110:LYS:HZ1	1:A:142:LYS:HB3	1.75	0.51
1:A:326:SER:HB3	1:A:329:CYS:HB2	1.91	0.51
1:A:74:PRO:HG3	1:A:573:GLU:HB2	1.93	0.51
1:A:69:ALA:C	1:A:71:SER:H	2.19	0.51
1:A:295:THR:HB	1:A:362:GLN:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:ARG:HB2	1:A:397:VAL:HG21	1.93	0.51
1:A:63:PRO:HG2	1:A:69:ALA:HB2	1.92	0.50
2:B:26:VAL:HG22	2:B:36:VAL:HG12	1.92	0.50
1:A:68:GLU:O	1:A:70:LEU:N	2.45	0.50
1:A:583:ASN:HD21	2:B:46:TRP:CD1	2.28	0.50
1:A:553:PRO:O	1:A:555:GLY:N	2.45	0.50
2:B:39:ILE:HD11	2:B:45:PHE:CE1	2.47	0.49
1:A:92:ILE:HG21	1:A:147:GLY:HA3	1.93	0.49
1:A:599:THR:N	1:A:602:LYS:HB2	2.28	0.49
1:A:447:VAL:HG12	1:A:482:VAL:HG13	1.95	0.48
1:A:95:PHE:CD1	1:A:545:LYS:HE3	2.49	0.48
1:A:327:LYS:N	1:A:395:ASP:OD2	2.47	0.48
1:A:48:GLN:H	2:B:40:MET:HG2	1.79	0.47
1:A:51:SER:OG	1:A:131:ILE:O	2.19	0.47
1:A:423:TYR:CZ	1:A:425:ILE:HG13	2.49	0.47
1:A:98:GLU:HB2	1:A:103:ASN:HA	1.95	0.47
1:A:693:PHE:CZ	1:A:697:GLY:HA3	2.49	0.47
2:B:12:GLU:O	2:B:57:THR:OG1	2.32	0.47
1:A:106:LEU:HG	1:A:107:ASN:HD22	1.79	0.47
1:A:396:SER:HA	1:A:419:ASN:HD22	1.79	0.47
1:A:162:LYS:HB2	1:A:714:PHE:HB2	1.96	0.47
1:A:95:PHE:CG	1:A:545:LYS:HE3	2.50	0.47
1:A:599:THR:OG1	1:A:600:THR:N	2.47	0.47
1:A:71:SER:OG	1:A:637:ARG:HD3	2.16	0.46
1:A:97:ALA:HB1	1:A:102:GLN:HA	1.97	0.46
1:A:533:PRO:HB3	2:B:37:SER:OG	2.16	0.46
1:A:127:GLU:HB2	1:A:154:LYS:HA	1.97	0.46
1:A:101:SER:OG	1:A:102:GLN:N	2.47	0.46
1:A:102:GLN:OE1	1:A:104:ASN:ND2	2.49	0.46
1:A:110:LYS:NZ	1:A:142:LYS:HB3	2.30	0.45
1:A:677:ALA:O	1:A:709:THR:N	2.42	0.45
1:A:99:GLY:O	1:A:505:GLN:HG2	2.16	0.45
1:A:145:PRO:HG3	1:A:499:SER:HB3	1.99	0.45
1:A:63:PRO:HB3	1:A:68:GLU:HB2	1.99	0.45
1:A:170:THR:HG22	1:A:706:VAL:H	1.82	0.45
1:A:145:PRO:HD3	1:A:501:GLU:OE1	2.17	0.45
1:A:73:THR:HG21	1:A:131:ILE:HG21	1.98	0.45
1:A:43:ILE:HD13	1:A:530:GLU:HG3	1.99	0.45
2:B:29:ARG:HD3	2:B:62:PHE:CZ	2.52	0.44
1:A:65:SER:N	1:A:68:GLU:OE1	2.48	0.44
1:A:106:LEU:HG	1:A:107:ASN:ND2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:GLN:O	1:A:50:ILE:HG23	2.18	0.44
1:A:112:GLN:HE21	1:A:119:ALA:HB2	1.83	0.44
1:A:145:PRO:O	1:A:520:GLN:NE2	2.51	0.44
1:A:255:PRO:HD3	1:A:261:ARG:HH21	1.82	0.44
1:A:503:SER:OG	1:A:513:PHE:HB2	2.17	0.44
1:A:156:PRO:HA	1:A:193:ARG:HB2	1.99	0.43
1:A:247:LEU:O	1:A:698:CYS:N	2.47	0.43
1:A:479:ARG:HD3	1:A:497:SER:HA	2.00	0.43
1:A:397:VAL:HG12	1:A:410:PHE:CE1	2.52	0.43
1:A:172:SER:O	1:A:172:SER:OG	2.26	0.43
1:A:43:ILE:HG22	2:B:23:LYS:HB3	2.00	0.43
2:B:35:LEU:HD13	2:B:46:TRP:CE2	2.53	0.43
1:A:70:LEU:O	1:A:73:THR:HG22	2.19	0.43
1:A:453:TYR:CD1	1:A:476:PHE:HB2	2.54	0.43
1:A:516:SER:OG	1:A:545:LYS:HG2	2.19	0.43
1:A:49:SER:HB2	2:B:43:GLY:HA2	2.01	0.42
1:A:501:GLU:H	1:A:501:GLU:HG3	1.61	0.42
1:A:118:ASP:O	1:A:305:ASN:ND2	2.52	0.42
1:A:182:ASP:HB3	1:A:192:TYR:HE1	1.85	0.42
1:A:619:ASP:HA	1:A:633:GLY:HA2	2.02	0.42
1:A:453:TYR:CE2	1:A:474:LYS:HG2	2.54	0.42
1:A:52:VAL:HG22	1:A:130:GLU:CB	2.50	0.42
1:A:527:TYR:CZ	1:A:529:PRO:HG3	2.55	0.42
1:A:250:GLU:O	1:A:252:THR:N	2.53	0.42
1:A:73:THR:OG1	1:A:74:PRO:HD2	2.20	0.42
1:A:140:TYR:O	1:A:479:ARG:NH1	2.52	0.41
1:A:153:SER:HB2	1:A:232:LEU:HD13	2.01	0.41
2:B:33:MET:HA	2:B:48:PRO:HA	2.02	0.41
1:A:483:ASN:HB3	1:A:493:TYR:CB	2.51	0.41
1:A:550:MET:HE1	1:A:562:GLY:HA3	2.01	0.41
1:A:115:PHE:HD2	1:A:387:ILE:HG12	1.85	0.41
1:A:167:LYS:HD3	1:A:175:GLN:CD	2.46	0.41
1:A:408:THR:HB	1:A:409:ASP:H	1.75	0.41
1:A:313:TYR:CZ	1:A:344:LYS:HB2	2.56	0.41
1:A:98:GLU:N	1:A:103:ASN:HB2	2.37	0.40
1:A:305:ASN:OD1	1:A:306:LYS:N	2.54	0.40
1:A:142:LYS:HZ3	1:A:452:ARG:NH2	2.19	0.40
1:A:54:THR:O	1:A:58:MET:N	2.48	0.40
1:A:508:LYS:H	1:A:508:LYS:HG2	1.69	0.40
1:A:663:ARG:HE	1:A:675:ASN:ND2	2.19	0.40
1:A:70:LEU:CD1	1:A:78:VAL:HG11	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:ALA:O	1:A:324:ALA:N	2.54	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	675/716 (94%)	592 (88%)	67 (10%)	16 (2%)	4	22
2	B	60/83 (72%)	46 (77%)	11 (18%)	3 (5%)	1	10
All	All	735/799 (92%)	638 (87%)	78 (11%)	19 (3%)	4	21

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	250	GLU
1	A	49	SER
1	A	69	ALA
1	A	324	ALA
1	A	554	GLU
2	B	56	ASN
1	A	41	THR
1	A	68	GLU
1	A	323	ASN
1	A	407	ASN
1	A	599	THR
1	A	628	SER
1	A	290	PHE
1	A	223	PRO
2	B	31	GLY
1	A	322	ALA
1	A	251	GLY

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Mol	Chain	Res	Type
1	A	398	PRO
2	B	38	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	566/593 (95%)	566 (100%)	0	100	100
2	B	53/71 (75%)	53 (100%)	0	100	100
All	All	619/664 (93%)	619 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	103	ASN
1	A	104	ASN
1	A	114	ASN
1	A	278	ASN
1	A	323	ASN
1	A	353	GLN
1	A	385	ASN
1	A	404	ASN
1	A	431	GLN
1	A	441	GLN
1	A	604	ASN
2	B	17	GLN
2	B	56	ASN

5.3.3 RNA ⓘ

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 oligosaccharides 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, 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	677/716 (94%)	0.55	75 (11%) 10 11	67, 109, 157, 223	0
2	B	62/83 (74%)	0.54	1 (1%) 70 55	121, 159, 206, 235	0
All	All	739/799 (92%)	0.55	76 (10%) 12 12	67, 112, 171, 235	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	82	GLY	10.5
1	A	244	TYR	9.6
1	A	81	ARG	8.6
1	A	83	ALA	8.6
1	A	702	ALA	7.7
1	A	116	TYR	7.7
1	A	557	PHE	7.6
1	A	701	GLY	7.1
1	A	79	GLY	6.9
1	A	246	TRP	6.9
1	A	550	MET	6.8
1	A	700	TRP	6.7
1	A	699	PHE	5.4
1	A	705	GLN	5.3
1	A	85	ASN	5.1
1	A	80	THR	5.1
1	A	89	HIS	4.8
1	A	115	PHE	4.8
1	A	64	LYS	4.7
1	A	552	ASP	4.7
1	A	87	TYR	4.6
1	A	688	TYR	4.6
1	A	268	GLU	4.5
1	A	84	SER	4.3

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Mol	Chain	Res	Type	RSRZ
1	A	100	GLN	4.3
1	A	513	PHE	4.2
1	A	511	ASN	4.1
1	A	204	GLN	3.9
1	A	690	ALA	3.8
1	A	95	PHE	3.8
1	A	703	GLU	3.6
1	A	48	GLN	3.5
1	A	101	SER	3.5
1	A	505	GLN	3.4
1	A	205	GLN	3.4
1	A	558	PHE	3.3
1	A	606	PRO	3.2
1	A	693	PHE	3.2
1	A	94	GLY	3.1
1	A	454	ASP	3.0
1	A	551	ALA	3.0
1	A	46	VAL	3.0
1	A	608	GLN	3.0
1	A	247	LEU	2.9
1	A	691	SER	2.9
1	A	262	LEU	2.9
1	A	275	TYR	2.9
1	A	47	PRO	2.9
1	A	687	GLU	2.9
1	A	267	ASN	2.8
1	A	171	ASP	2.8
1	A	203	ALA	2.7
1	A	597	THR	2.7
1	A	102	GLN	2.5
1	A	509	ASP	2.5
1	A	313	TYR	2.5
1	A	88	ASP	2.5
1	A	118	ASP	2.5
1	A	391	PHE	2.4
1	A	99	GLY	2.4
1	A	207	GLY	2.4
1	A	506	VAL	2.4
1	A	245	GLY	2.2
1	A	49	SER	2.2
1	A	44	GLN	2.2
1	A	269	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	576	ALA	2.1
1	A	138	VAL	2.1
1	A	469	ASP	2.1
1	A	142	LYS	2.0
1	A	512	ILE	2.0
1	A	657	VAL	2.0
1	A	689	VAL	2.0
2	B	44	SER	2.0
1	A	477	THR	2.0
1	A	622	PHE	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 oligosaccharides 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.