



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:28 PM UTC

PDB ID : 7ZFD / pdb_00007zfd
Title : SARS-CoV-2 Omicron RBD in complex with Omi-25 Fab
Authors : Zhou, D.; Huo, J.; Ren, J.; Stuart, D.I.
Deposited on : 2022-04-01
Resolution : 3.39 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:

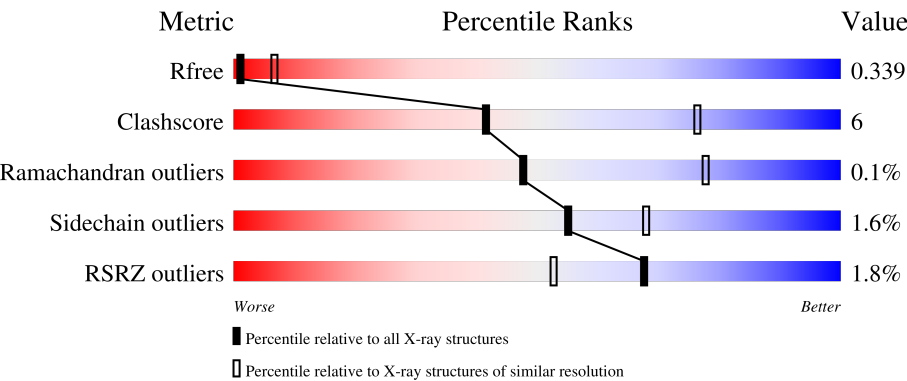
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.39 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	228	0% 80% 18% 2%
1	H	228	2% 77% 21% 2%
2	E	209	2% 77% 11% 11%
2	F	209	3% 77% 9% 13%
3	B	214	87% 12%

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Mol	Chain	Length	Quality of chain
3	L	214	% 83% 16%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9597 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 Omi-25 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	224	Total	C	N	O	S	0	1	0
			1681	1062	286	326	7			
1	A	224	Total	C	N	O	S	0	1	0
			1679	1062	284	326	7			

- Molecule 2 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	185	Total	C	N	O	S	0	0	0
			1496	965	251	272	8			
2	F	182	Total	C	N	O	S	0	0	0
			1465	940	248	269	8			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	339	ASP	GLY	variant	UNP P0DTC2
E	371	LEU	SER	variant	UNP P0DTC2
E	373	PRO	SER	variant	UNP P0DTC2
E	375	PHE	SER	variant	UNP P0DTC2
E	417	ASN	LYS	variant	UNP P0DTC2
E	440	LYS	ASN	variant	UNP P0DTC2
E	446	SER	GLY	variant	UNP P0DTC2
E	477	ASN	SER	variant	UNP P0DTC2
E	478	LYS	THR	variant	UNP P0DTC2
E	484	ALA	GLU	variant	UNP P0DTC2
E	493	ARG	GLN	variant	UNP P0DTC2
E	496	SER	GLY	variant	UNP P0DTC2
E	498	ARG	GLN	conflict	UNP P0DTC2
E	501	TYR	ASN	variant	UNP P0DTC2
E	505	HIS	TYR	variant	UNP P0DTC2
E	533	HIS	-	expression tag	UNP P0DTC2

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Chain	Residue	Modelled	Actual	Comment	Reference
E	534	HIS	-	expression tag	UNP P0DTC2
E	535	HIS	-	expression tag	UNP P0DTC2
E	536	HIS	-	expression tag	UNP P0DTC2
E	537	HIS	-	expression tag	UNP P0DTC2
E	538	HIS	-	expression tag	UNP P0DTC2
F	339	ASP	GLY	variant	UNP P0DTC2
F	371	LEU	SER	variant	UNP P0DTC2
F	373	PRO	SER	variant	UNP P0DTC2
F	375	PHE	SER	variant	UNP P0DTC2
F	417	ASN	LYS	variant	UNP P0DTC2
F	440	LYS	ASN	variant	UNP P0DTC2
F	446	SER	GLY	variant	UNP P0DTC2
F	477	ASN	SER	variant	UNP P0DTC2
F	478	LYS	THR	variant	UNP P0DTC2
F	484	ALA	GLU	variant	UNP P0DTC2
F	493	ARG	GLN	variant	UNP P0DTC2
F	496	SER	GLY	variant	UNP P0DTC2
F	498	ARG	GLN	conflict	UNP P0DTC2
F	501	TYR	ASN	variant	UNP P0DTC2
F	505	HIS	TYR	variant	UNP P0DTC2
F	533	HIS	-	expression tag	UNP P0DTC2
F	534	HIS	-	expression tag	UNP P0DTC2
F	535	HIS	-	expression tag	UNP P0DTC2
F	536	HIS	-	expression tag	UNP P0DTC2
F	537	HIS	-	expression tag	UNP P0DTC2
F	538	HIS	-	expression tag	UNP P0DTC2

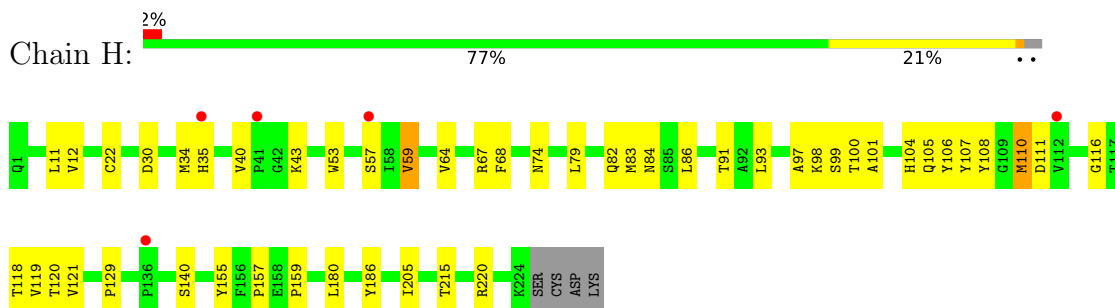
- Molecule 3 is a protein called Omi-25 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	0	0
			1638	1027	273	333	5			
3	B	213	Total	C	N	O	S	0	0	0
			1638	1027	273	333	5			

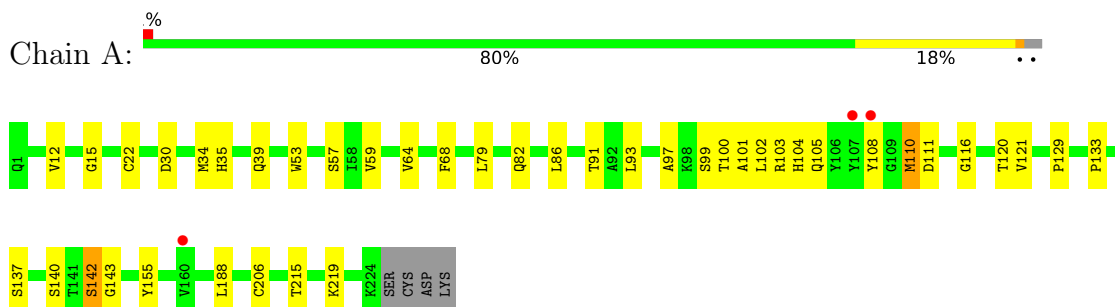
3 Residue-property plots [i](#)

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.

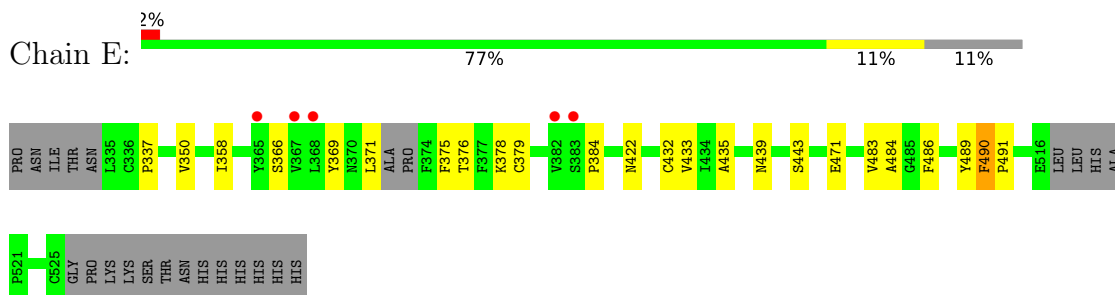
- Molecule 1: Omi-25 heavy chain



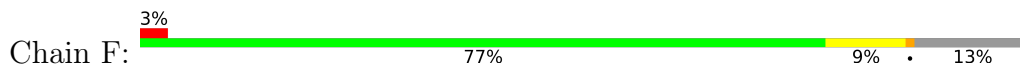
- Molecule 1: Omi-25 heavy chain

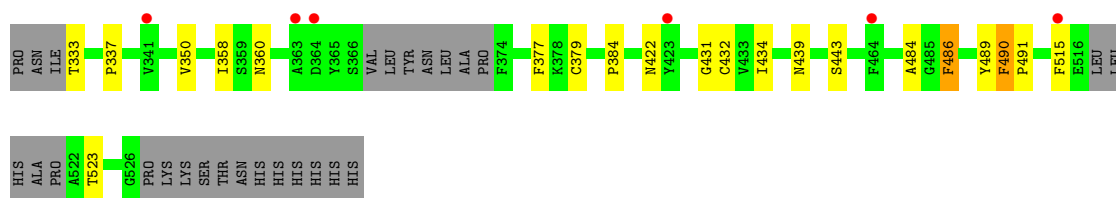


- Molecule 2: Spike protein S1

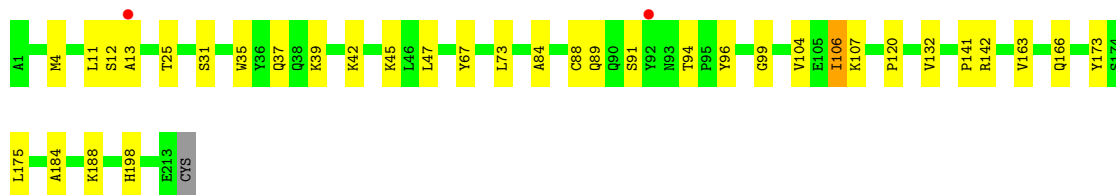
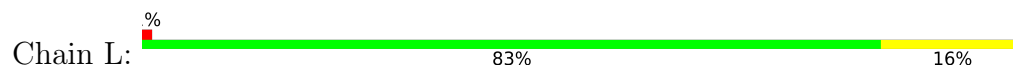


- Molecule 2: Spike protein S1

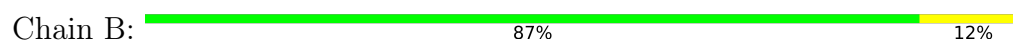




• Molecule 3: Omi-25 light chain



• Molecule 3: Omi-25 light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	123.05Å 123.05Å 223.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.67 – 3.39 68.67 – 3.39	Depositor EDS
% Data completeness (in resolution range)	99.7 (68.67-3.39) 99.7 (68.67-3.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.00 (at 3.41Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.286 , 0.338 0.286 , 0.339	Depositor DCC
R_{free} test set	1218 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	85.8	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 89.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.32$, $\langle L^2 \rangle = 0.16$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	9597	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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 [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	0.12	0/1724	0.36	0/2351
1	H	0.13	0/1726	0.34	0/2354
2	E	0.09	0/1538	0.25	0/2087
2	F	0.08	0/1505	0.23	0/2041
3	B	0.09	0/1674	0.28	0/2274
3	L	0.09	0/1674	0.29	0/2274
All	All	0.10	0/9841	0.30	0/13381

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 [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	1679	0	1629	27	0
1	H	1681	0	1629	35	0
2	E	1496	0	1424	16	0
2	F	1465	0	1386	11	0
3	B	1638	0	1588	17	0
3	L	1638	0	1588	23	0
All	All	9597	0	9244	120	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 6.

All (120) 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:104:HIS:ND1	2:F:484:ALA:O	2.17	0.78
1:H:205:ILE:HG12	1:H:220:ARG:HG2	1.68	0.75
1:A:64:VAL:HG11	1:A:68:PHE:HB2	1.68	0.74
1:H:59:VAL:HG11	3:L:94:THR:HG21	1.70	0.74
1:H:104:HIS:ND1	2:E:484:ALA:O	2.22	0.73
1:A:101:ALA:HB1	1:A:105:GLN:HB3	1.70	0.73
2:E:490:PHE:HD1	2:E:491:PRO:HD2	1.54	0.72
3:L:12:SER:HB3	3:L:107:LYS:HD2	1.70	0.72
1:H:101:ALA:HB2	1:H:108:TYR:HD2	1.58	0.68
1:H:129:PRO:HB3	1:H:155:TYR:HB3	1.76	0.68
1:A:57:SER:HB3	2:F:486:PHE:HB2	1.76	0.68
1:A:129:PRO:HB3	1:A:155:TYR:HB3	1.78	0.65
1:A:93:LEU:HD21	1:A:116:GLY:HA3	1.79	0.64
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.80	0.63
3:B:37:GLN:HB2	3:B:47:LEU:HD11	1.81	0.62
1:H:64:VAL:HG13	1:H:68:PHE:HB2	1.83	0.61
1:H:93:LEU:HD21	1:H:116:GLY:HA3	1.82	0.60
1:A:39:GLN:OE1	3:B:38:GLN:NE2	2.30	0.60
1:H:35:HIS:HE2	1:H:99:SER:HG	1.50	0.58
2:F:439:ASN:O	2:F:443:SER:OG	2.20	0.58
1:H:35:HIS:NE2	1:H:99:SER:OG	2.32	0.58
2:E:439:ASN:O	2:E:443:SER:OG	2.21	0.58
3:L:120:PRO:HD3	3:L:132:VAL:HG22	1.86	0.57
3:B:120:PRO:HD3	3:B:132:VAL:HG22	1.87	0.57
2:E:371:LEU:HD13	2:E:375:PHE:HB3	1.87	0.57
1:H:30:ASP:O	1:H:53:TRP:HB2	2.05	0.56
1:A:103:ARG:HG2	1:A:104:HIS:H	1.72	0.55
1:H:129:PRO:HD2	1:H:215:THR:HG21	1.89	0.54
3:L:31:SER:O	3:L:31:SER:OG	2.25	0.54
1:H:57:SER:HB3	2:E:486:PHE:HD2	1.73	0.54
1:A:91:THR:HG23	1:A:120:THR:HA	1.89	0.53
1:H:40:VAL:HB	1:H:43:LYS:HB2	1.92	0.52
3:L:35:TRP:CE2	3:L:73:LEU:HB2	2.45	0.52
3:L:142:ARG:HD2	3:L:173:TYR:CE1	2.45	0.52
1:A:35:HIS:HE2	1:A:99:SER:HB2	1.74	0.51
1:H:101:ALA:HB2	1:H:108:TYR:CD2	2.44	0.51
1:A:129:PRO:HD2	1:A:215:THR:HG21	1.92	0.51
1:H:34:MET:HB3	1:H:79:LEU:HD22	1.90	0.51
1:A:34:MET:HB3	1:A:79:LEU:HD22	1.92	0.51
3:B:39:LYS:HD3	3:B:84:ALA:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:39:LYS:HB2	3:B:42:LYS:HB2	1.92	0.51
1:H:118:THR:HB	1:H:159:PRO:HD3	1.92	0.51
1:A:22:CYS:HB3	1:A:79:LEU:HB3	1.92	0.51
1:A:97:ALA:HB1	1:A:110:MET:HB3	1.93	0.51
1:H:12:VAL:HG11	1:H:86:LEU:HD13	1.92	0.50
1:H:100:THR:HG22	1:H:111:ASP:HB3	1.92	0.50
1:H:22:CYS:HB3	1:H:79:LEU:HB3	1.93	0.50
1:H:11:LEU:HG	1:H:157:PRO:HD3	1.93	0.50
1:H:107:TYR:HA	3:L:91:SER:HB3	1.94	0.50
3:L:39:LYS:HB2	3:L:42:LYS:HB2	1.93	0.50
3:B:35:TRP:CE2	3:B:73:LEU:HB2	2.47	0.49
1:H:83:MET:HE1	1:H:119:VAL:HG21	1.93	0.49
2:E:371:LEU:HD22	2:E:375:PHE:CG	2.48	0.49
3:B:4:MET:HE1	3:B:25:THR:HG22	1.94	0.49
3:B:31:SER:HB3	3:B:67:TYR:CE2	2.48	0.49
2:E:379:CYS:HA	2:E:432:CYS:HA	1.94	0.49
1:H:12:VAL:HG13	1:H:121:VAL:HG22	1.94	0.49
1:A:100:THR:HG22	1:A:111:ASP:HB3	1.95	0.48
1:A:102:LEU:O	1:A:105:GLN:HB2	2.14	0.48
2:E:350:VAL:HG22	2:E:422:ASN:HB3	1.96	0.47
1:H:97:ALA:HB1	1:H:110:MET:HB3	1.96	0.47
1:H:91:THR:HG23	1:H:120:THR:HA	1.95	0.47
3:B:31:SER:O	3:B:31:SER:OG	2.26	0.47
3:L:4:MET:HE1	3:L:25:THR:HG22	1.96	0.47
3:B:94:THR:HG22	3:B:95:PRO:HA	1.97	0.47
3:L:89:GLN:HE21	3:L:96:TYR:HB3	1.79	0.46
3:L:89:GLN:NE2	3:L:96:TYR:HB3	2.29	0.46
1:H:64:VAL:CG1	1:H:68:PHE:HB2	2.44	0.46
2:F:379:CYS:HA	2:F:432:CYS:HA	1.98	0.46
1:A:206:CYS:SG	1:A:219:LYS:HB3	2.56	0.46
2:F:379:CYS:SG	2:F:384:PRO:HG3	2.56	0.46
2:F:490:PHE:CD1	2:F:491:PRO:HD2	2.52	0.45
2:E:379:CYS:SG	2:E:384:PRO:HG3	2.56	0.45
2:F:350:VAL:HG22	2:F:422:ASN:HB3	1.99	0.45
3:L:106:ILE:O	3:L:166:GLN:NE2	2.49	0.45
3:L:141:PRO:HD2	3:L:198:HIS:CE1	2.52	0.45
1:A:12:VAL:HG11	1:A:86:LEU:HD13	1.99	0.45
3:L:31:SER:HB3	3:L:67:TYR:CE2	2.51	0.45
2:E:337:PRO:HD2	2:E:358:ILE:HG23	1.98	0.45
3:L:13:ALA:O	3:L:106:ILE:HA	2.16	0.45
1:A:137:SER:H	1:A:140:SER:HB3	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:39:LYS:HD3	3:L:84:ALA:HB2	1.99	0.44
1:H:101:ALA:HB3	1:H:105:GLN:O	2.17	0.44
1:H:35:HIS:HE1	3:L:96:TYR:HE2	1.65	0.44
2:E:366:SER:HA	2:E:369:TYR:HB3	2.00	0.43
1:A:30:ASP:O	1:A:53:TRP:HB2	2.18	0.43
1:H:140:SER:O	1:H:140:SER:OG	2.28	0.43
2:E:371:LEU:HD23	2:E:371:LEU:HA	1.83	0.43
1:A:12:VAL:HG13	1:A:121:VAL:HG22	2.00	0.42
1:A:101:ALA:HB2	1:A:108:TYR:CD2	2.54	0.42
3:B:25:THR:OG1	3:B:27:GLN:O	2.35	0.42
1:A:133:PRO:HD3	1:A:219:LYS:HE3	2.01	0.42
1:H:30:ASP:CG	1:H:74:ASN:HB3	2.45	0.42
1:H:118:THR:CB	1:H:159:PRO:HD3	2.49	0.42
1:H:180:LEU:HD13	1:H:186:TYR:CE1	2.55	0.42
2:F:337:PRO:HD2	2:F:358:ILE:HG23	2.01	0.42
3:B:89:GLN:HE21	3:B:96:TYR:HB3	1.84	0.42
2:E:376:THR:OG1	2:E:435:ALA:O	2.29	0.42
3:L:11:LEU:HD11	3:L:104:VAL:HG22	2.02	0.42
1:A:142:SER:OG	1:A:143:GLY:N	2.53	0.42
3:B:88:CYS:O	3:B:99:GLY:N	2.47	0.42
1:H:67:ARG:HB3	1:H:84:ASN:O	2.19	0.41
3:L:35:TRP:CD2	3:L:73:LEU:HB2	2.56	0.41
2:F:360:ASN:HA	2:F:523:THR:HB	2.01	0.41
2:F:377:PHE:CD1	2:F:434:ILE:HG12	2.56	0.41
3:B:89:GLN:NE2	3:B:96:TYR:HB3	2.35	0.41
1:H:98:LYS:O	1:H:110:MET:HA	2.21	0.41
3:L:163:VAL:HG22	3:L:175:LEU:HD12	2.03	0.41
1:H:68:PHE:HA	1:H:82:GLN:O	2.20	0.41
2:E:483:VAL:HG21	1:A:15:GLY:C	2.46	0.41
3:B:89:GLN:HB2	3:B:98:PHE:CD2	2.56	0.41
1:A:129:PRO:CB	1:A:155:TYR:HB3	2.50	0.40
3:B:35:TRP:CD2	3:B:73:LEU:HB2	2.56	0.40
3:B:186:TYR:O	3:B:192:TYR:OH	2.36	0.40
2:E:378:LYS:O	2:E:433:VAL:N	2.41	0.40
2:E:471:GLU:O	2:E:491:PRO:HG3	2.22	0.40
3:L:88:CYS:O	3:L:99:GLY:N	2.46	0.40
3:L:184:ALA:O	3:L:188:LYS:HG3	2.20	0.40
2:F:431:GLY:HA2	2:F:515:PHE:HD2	1.86	0.40
1:A:68:PHE:HA	1:A:82:GLN:O	2.21	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	223/228 (98%)	210 (94%)	12 (5%)	1 (0%)	30	59
1	H	223/228 (98%)	212 (95%)	11 (5%)	0	100	100
2	E	179/209 (86%)	167 (93%)	12 (7%)	0	100	100
2	F	176/209 (84%)	163 (93%)	13 (7%)	0	100	100
3	B	211/214 (99%)	204 (97%)	7 (3%)	0	100	100
3	L	211/214 (99%)	204 (97%)	7 (3%)	0	100	100
All	All	1223/1302 (94%)	1160 (95%)	62 (5%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	142	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	185/191 (97%)	182 (98%)	3 (2%)	55	68
1	H	185/191 (97%)	182 (98%)	3 (2%)	55	68
2	E	163/184 (89%)	161 (99%)	2 (1%)	63	72
2	F	159/184 (86%)	155 (98%)	4 (2%)	42	62
3	B	187/189 (99%)	184 (98%)	3 (2%)	55	68
3	L	187/189 (99%)	185 (99%)	2 (1%)	65	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1066/1128 (94%)	1049 (98%)	17 (2%)	55 68

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	59	VAL
1	H	106	TYR
1	H	110	MET
2	E	489	TYR
2	E	490	PHE
3	L	45	LYS
3	L	106	ILE
1	A	59	VAL
1	A	110	MET
1	A	188	LEU
2	F	333	THR
2	F	486	PHE
2	F	489	TYR
2	F	490	PHE
3	B	45	LYS
3	B	94	THR
3	B	106	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	209	ASN
3	L	34	ASN
3	L	89	GLN
3	L	166	GLN
1	A	181	GLN
3	B	34	ASN
3	B	89	GLN

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	224/228 (98%)	-0.05	3 (1%) 75 61	57, 88, 131, 176	1 (0%)
1	H	224/228 (98%)	0.08	5 (2%) 62 47	47, 72, 120, 214	1 (0%)
2	E	185/209 (88%)	0.31	5 (2%) 56 42	70, 105, 192, 300	0
2	F	182/209 (87%)	0.26	6 (3%) 49 36	60, 116, 198, 263	0
3	B	213/214 (99%)	-0.07	1 (0%) 87 78	62, 93, 132, 159	0
3	L	213/214 (99%)	0.05	2 (0%) 81 68	56, 85, 126, 168	0
All	All	1241/1302 (95%)	0.08	22 (1%) 67 53	47, 90, 166, 300	2 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	368	LEU	4.9
2	E	382	VAL	4.4
2	F	423	TYR	3.7
2	E	365	TYR	3.5
2	E	367	VAL	3.1
2	E	383	SER	3.1
1	H	136	PRO	2.9
2	F	515	PHE	2.8
2	F	364	ASP	2.7
1	A	107	TYR	2.7
3	L	92	TYR	2.7
1	H	57	SER	2.6
2	F	341	VAL	2.6
1	A	160	VAL	2.5
1	A	108	TYR	2.4
1	H	112	VAL	2.2
1	H	41	PRO	2.2
3	L	13	ALA	2.2
2	F	464	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	35	HIS	2.2
2	F	363	ALA	2.1
3	B	47	LEU	2.1

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.