



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2021 – 06:19 PM EDT

PDB ID : 7N1B
Title : SARS-CoV-2 RLQ peptide binds to HLA-A2
Authors : Wu, D.; Mariuzza, R.A.
Deposited on : 2021-05-27
Resolution : 2.81 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 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

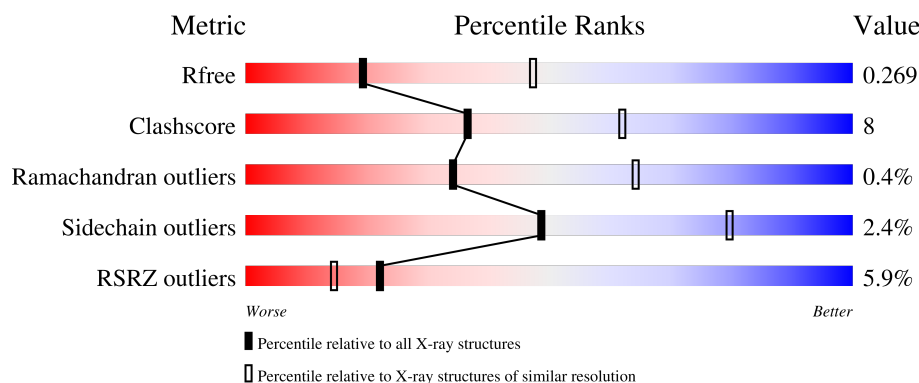
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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	275	<div> <div>4%</div> <div>83%</div> <div>16%</div> <div>.</div> </div>
1	D	275	<div> <div>11%</div> <div>79%</div> <div>17%</div> <div>..</div> </div>
2	B	100	<div> <div>%</div> <div>89%</div> <div>9%</div> <div>.</div> </div>
2	E	100	<div> <div>2%</div> <div>87%</div> <div>12%</div> <div>.</div> </div>
3	C	9	<div> <div>67%</div> <div>33%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	9	 89% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6261 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 MHC class I antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2233	1396	407	421	9			
1	D	272	Total	C	N	O	S	0	0	0
			2206	1378	399	420	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			833	530	140	159	4			
2	E	100	Total	C	N	O	S	0	0	0
			833	530	140	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
E	0	MET	-	initiating methionine	UNP P61769

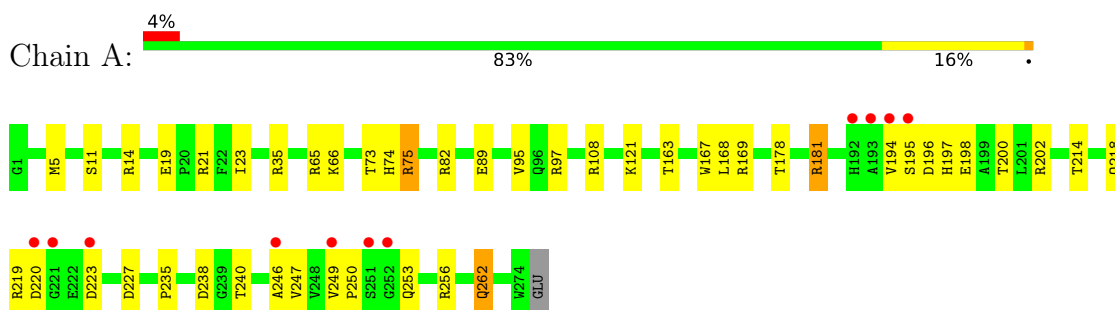
- Molecule 3 is a protein called Spike protein S2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			78	49	14	15			
3	F	9	Total	C	N	O	0	0	0
			78	49	14	15			

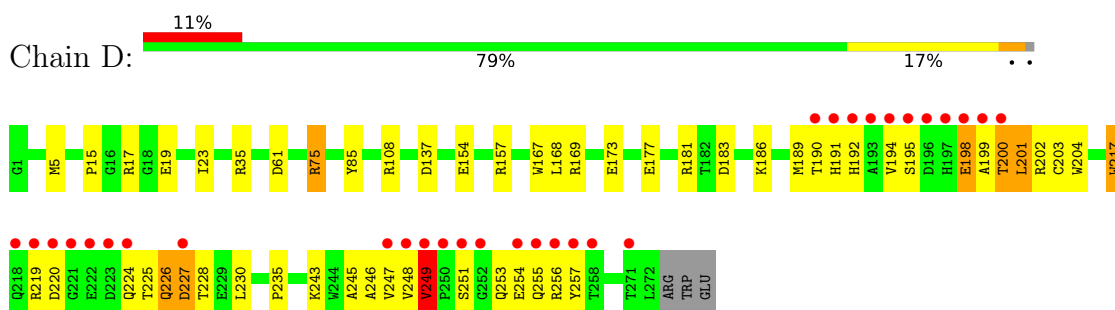
3 Residue-property plots

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.

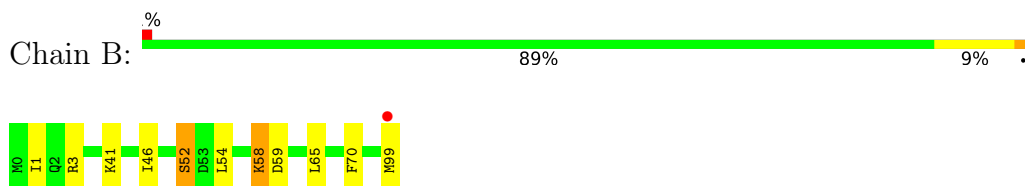
- Molecule 1: MHC class I antigen, A-2 alpha chain



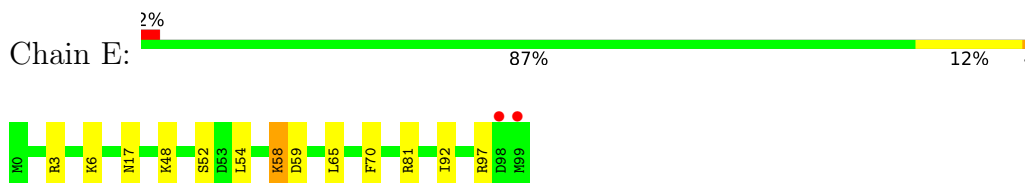
- Molecule 1: MHC class I antigen, A-2 alpha chain



- Molecule 2: Beta-2-microglobulin

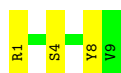


- Molecule 2: Beta-2-microglobulin




- Molecule 3: Spike protein S2

Chain C:  67% 33%



● Molecule 3: Spike protein S2

Chain F:  89% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.20Å 49.26Å 117.09Å 91.89° 92.39° 118.59°	Depositor
Resolution (Å)	43.17 – 2.81 43.17 – 2.81	Depositor EDS
% Data completeness (in resolution range)	96.0 (43.17-2.81) 96.0 (43.17-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.85 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.211 , 0.269 0.210 , 0.269	Depositor DCC
R_{free} test set	1074 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	25.6	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.042 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	6261	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.70% 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	0.26	0/2298	0.47	0/3120
1	D	0.32	0/2269	0.60	3/3081 (0.1%)
2	B	0.25	0/856	0.45	0/1158
2	E	0.31	0/856	0.47	0/1158
3	C	0.23	0/78	0.47	0/103
3	F	0.21	0/78	0.41	0/103
All	All	0.28	0/6435	0.52	3/8723 (0.0%)

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	D	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	198	GLU	CA-CB-CG	6.40	127.48	113.40
1	D	249	VAL	CG1-CB-CG2	-5.91	101.45	110.90
1	D	201	LEU	CA-CB-CG	5.15	127.14	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	200	THR	Peptide
1	D	249	VAL	Peptide

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	2233	0	2084	32	1
1	D	2206	0	2056	54	1
2	B	833	0	792	8	0
2	E	833	0	792	8	0
3	C	78	0	83	5	0
3	F	78	0	83	1	0
All	All	6261	0	5890	96	1

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 8.

All (96) 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:D:192:HIS:HB2	1:D:200:THR:CG2	1.89	1.02
1:D:226:GLN:OE1	1:D:227:ASP:HB2	1.59	1.01
1:D:192:HIS:HB2	1:D:200:THR:HG21	1.46	0.97
1:D:226:GLN:OE1	1:D:227:ASP:N	2.01	0.92
1:D:201:LEU:HG	1:D:257:TYR:HE1	1.34	0.92
1:D:201:LEU:HG	1:D:257:TYR:CE1	2.14	0.83
1:A:194:VAL:HG12	1:A:195:SER:HB2	1.66	0.78
1:D:219:ARG:NH2	1:D:253:GLN:OE1	2.11	0.78
1:A:195:SER:O	1:A:197:HIS:N	2.17	0.76
1:D:192:HIS:HB2	1:D:200:THR:HG23	1.68	0.76
1:D:198:GLU:OE2	1:D:249:VAL:O	2.04	0.75
1:D:191:HIS:HA	1:D:200:THR:OG1	1.90	0.71
2:E:17:ASN:OD1	2:E:97:ARG:NH2	2.26	0.68
1:D:226:GLN:OE1	1:D:227:ASP:CB	2.38	0.68
1:A:82:ARG:HH21	1:A:89:GLU:HG2	1.59	0.68
1:A:121:LYS:HE3	2:B:1:ILE:HG21	1.76	0.68
1:A:66:LYS:HD2	3:C:4:SER:HA	1.79	0.65
1:D:255:GLN:HA	1:D:257:TYR:CD2	2.31	0.64
1:D:199:ALA:H	1:D:249:VAL:HG22	1.63	0.64
1:D:219:ARG:HH21	1:D:256:ARG:HG3	1.62	0.64
2:B:52:SER:HB3	2:B:65:LEU:H	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:52:SER:HB3	2:E:65:LEU:H	1.62	0.63
1:D:201:LEU:HD22	1:D:247:VAL:HG23	1.82	0.62
1:D:202:ARG:HE	1:D:246:ALA:HB2	1.65	0.61
1:D:226:GLN:CD	1:D:227:ASP:HB2	2.23	0.59
1:A:227:ASP:O	1:A:247:VAL:HA	2.04	0.58
1:D:226:GLN:CD	1:D:227:ASP:N	2.57	0.58
1:A:214:THR:HB	1:A:262:GLN:HB2	1.85	0.57
1:D:200:THR:HA	1:D:248:VAL:HG12	1.87	0.56
1:A:218:GLN:HA	1:A:223:ASP:HA	1.89	0.54
1:D:15:PRO:O	1:D:17:ARG:NH2	2.41	0.54
1:D:154:GLU:HG3	1:D:157:ARG:HH22	1.72	0.54
1:D:219:ARG:HE	1:D:256:ARG:HB3	1.74	0.53
1:D:23:ILE:HG21	2:E:54:LEU:HB3	1.92	0.52
1:D:201:LEU:HB2	1:D:247:VAL:O	2.10	0.51
1:D:191:HIS:HA	1:D:200:THR:HG1	1.76	0.51
1:A:167:TRP:CE2	3:C:1:ARG:HD3	2.46	0.51
1:D:169:ARG:O	1:D:173:GLU:HG3	2.11	0.51
1:A:178:THR:O	1:A:181:ARG:HG3	2.11	0.50
2:E:3:ARG:NH1	2:E:59:ASP:OD2	2.45	0.50
1:A:253:GLN:NE2	1:A:256:ARG:HH12	2.10	0.49
2:E:48:LYS:HB2	2:E:48:LYS:HE2	1.59	0.49
1:D:254:GLU:O	1:D:257:TYR:HE2	1.96	0.49
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.95	0.48
1:A:198:GLU:HA	1:A:249:VAL:O	2.12	0.48
1:D:226:GLN:OE1	1:D:227:ASP:CA	2.61	0.48
1:D:5:MET:HB2	1:D:168:LEU:HD13	1.95	0.48
2:B:41:LYS:HB2	2:B:46:ILE:HG12	1.94	0.48
2:B:58:LYS:HD2	2:B:58:LYS:H	1.79	0.47
1:A:73:THR:HG22	3:C:8:TYR:CZ	2.49	0.47
1:D:85:TYR:OH	1:D:137:ASP:OD2	2.25	0.47
1:A:108:ARG:HA	1:A:169:ARG:HH11	1.79	0.47
1:D:189:MET:HE1	1:D:257:TYR:CE1	2.50	0.47
1:D:230:LEU:HD11	1:D:243:LYS:HE3	1.96	0.47
1:D:203:CYS:HB2	1:D:217:TRP:CE2	2.50	0.47
1:D:225:THR:HA	1:D:228:THR:OG1	2.15	0.47
1:D:190:THR:HB	1:D:192:HIS:CD2	2.50	0.47
1:A:11:SER:HA	1:A:21:ARG:O	2.16	0.46
1:A:23:ILE:HG21	2:B:54:LEU:HB3	1.98	0.46
1:A:74:HIS:CE1	1:A:97:ARG:HE	2.34	0.45
1:A:235:PRO:HG2	2:B:65:LEU:HD13	1.98	0.45
1:D:19:GLU:HG2	1:D:75:ARG:NH1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:58:LYS:H	2:E:58:LYS:HD2	1.82	0.45
1:A:163:THR:HG21	3:C:1:ARG:NH2	2.32	0.45
1:D:108:ARG:HA	1:D:169:ARG:HH11	1.81	0.45
2:B:3:ARG:NH2	2:B:59:ASP:O	2.48	0.45
1:D:235:PRO:HG2	2:E:65:LEU:HD13	1.99	0.45
1:D:201:LEU:HB3	1:D:217:TRP:CH2	2.53	0.44
1:D:201:LEU:HB3	1:D:217:TRP:CZ2	2.53	0.44
1:A:238:ASP:OD1	1:A:240:THR:HG22	2.18	0.44
1:A:202:ARG:NH1	2:B:99:MET:HG2	2.32	0.43
1:D:201:LEU:HD23	1:D:217:TRP:CZ3	2.53	0.43
2:E:81:ARG:HG3	2:E:92:ILE:HG12	2.00	0.43
1:D:192:HIS:N	1:D:200:THR:OG1	2.52	0.43
1:D:177:GLU:H	1:D:177:GLU:CD	2.22	0.43
1:D:201:LEU:HB2	1:D:247:VAL:HG23	2.00	0.42
1:A:82:ARG:NH2	1:A:89:GLU:HG2	2.30	0.42
1:A:202:ARG:HE	1:A:246:ALA:HB2	1.83	0.42
1:D:181:ARG:NH1	1:D:183:ASP:OD2	2.52	0.42
1:D:219:ARG:HB2	1:D:224:GLN:NE2	2.34	0.42
1:D:186:LYS:HA	1:D:186:LYS:HD3	1.73	0.42
1:A:19:GLU:HG2	1:A:75:ARG:HH11	1.85	0.42
1:D:202:ARG:HG2	1:D:204:TRP:NE1	2.35	0.42
1:D:167:TRP:NE1	3:F:1:ARG:HD2	2.35	0.42
1:A:66:LYS:HB2	1:A:66:LYS:HE2	1.85	0.42
1:D:194:VAL:HG11	1:D:198:GLU:HG3	2.00	0.41
1:A:11:SER:HB3	1:A:95:VAL:HG12	2.01	0.41
1:A:66:LYS:HD2	3:C:4:SER:CA	2.49	0.41
1:A:200:THR:HA	1:A:247:VAL:O	2.20	0.41
1:D:198:GLU:HA	1:D:251:SER:HB2	2.02	0.41
1:D:217:TRP:NE1	1:D:245:ALA:O	2.50	0.41
1:A:14:ARG:NH1	1:A:19:GLU:O	2.49	0.41
1:A:197:HIS:O	1:A:250:PRO:HA	2.21	0.41
1:D:219:ARG:HG2	1:D:220:ASP:OD1	2.21	0.41
1:A:19:GLU:HG2	1:A:75:ARG:NH1	2.36	0.40
1:D:256:ARG:HD3	1:D:256:ARG:HA	1.84	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ARG:NH1	1:D:61:ASP:OD2[1_455]	2.06	0.14

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	272/275 (99%)	261 (96%)	10 (4%)	1 (0%)	34	64
1	D	270/275 (98%)	256 (95%)	12 (4%)	2 (1%)	22	51
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	E	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	7 (100%)	0	0	100	100
All	All	752/768 (98%)	723 (96%)	26 (4%)	3 (0%)	34	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	ASP
1	D	226	GLN
1	D	195	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	229/231 (99%)	223 (97%)	6 (3%)	46	78
1	D	227/231 (98%)	223 (98%)	4 (2%)	59	85
2	B	94/95 (99%)	91 (97%)	3 (3%)	39	71
2	E	94/95 (99%)	91 (97%)	3 (3%)	39	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	9/9 (100%)	9 (100%)	0	100	100
3	F	9/9 (100%)	9 (100%)	0	100	100
All	All	662/670 (99%)	646 (98%)	16 (2%)	49	80

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ARG
1	A	75	ARG
1	A	181	ARG
1	A	219	ARG
1	A	220	ASP
1	A	262	GLN
2	B	52	SER
2	B	58	LYS
2	B	70	PHE
1	D	35	ARG
1	D	75	ARG
1	D	217	TRP
1	D	227	ASP
2	E	6	LYS
2	E	58	LYS
2	E	70	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	253	GLN
1	D	224	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [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, 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	274/275 (99%)	-0.05	11 (4%) 38 28	9, 20, 84, 100	0
1	D	272/275 (98%)	0.32	31 (11%) 5 3	8, 20, 111, 132	0
2	B	100/100 (100%)	-0.09	1 (1%) 82 77	12, 26, 58, 76	0
2	E	100/100 (100%)	-0.24	2 (2%) 65 56	13, 24, 54, 78	0
3	C	9/9 (100%)	-0.37	0 100 100	10, 15, 21, 23	0
3	F	9/9 (100%)	-0.49	0 100 100	12, 15, 18, 18	0
All	All	764/768 (99%)	0.05	45 (5%) 22 14	8, 21, 86, 132	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	251	SER	13.9
1	D	257	TYR	11.4
1	D	249	VAL	9.5
1	D	199	ALA	9.2
1	D	220	ASP	8.2
1	D	194	VAL	7.7
1	D	250	PRO	6.5
1	D	252	GLY	5.8
1	D	192	HIS	5.0
1	D	196	ASP	4.9
1	D	191	HIS	4.8
1	D	200	THR	4.7
1	D	195	SER	4.6
1	A	251	SER	4.6
1	D	198	GLU	4.3
2	B	99	MET	4.3
1	A	194	VAL	4.3
1	D	247	VAL	4.1
1	A	221	GLY	4.0

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Mol	Chain	Res	Type	RSRZ
1	D	256	ARG	4.0
1	D	254	GLU	4.0
1	D	248	VAL	3.9
1	A	193	ALA	3.9
1	D	271	THR	3.9
2	E	99	MET	3.9
1	D	255	GLN	3.8
1	D	193	ALA	3.6
1	D	218	GLN	3.5
1	D	223	ASP	3.5
1	D	221	GLY	3.4
1	D	227	ASP	3.3
1	D	258	THR	3.2
1	A	195	SER	3.1
1	D	190	THR	2.9
1	A	223	ASP	2.8
1	D	224	GLN	2.8
1	D	197	HIS	2.6
1	D	219	ARG	2.6
1	A	192	HIS	2.5
1	A	252	GLY	2.3
1	D	222	GLU	2.2
2	E	98	ASP	2.2
1	A	246	ALA	2.2
1	A	249	VAL	2.1
1	A	220	ASP	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 monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.