



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:41 PM UTC

PDB ID : 4ZPW / pdb_00004zpw
Title : Structure of unbound MERS-CoV spike receptor-binding domain (England1 strain).
Authors : Joyce, M.G.; Mascola, J.R.; Graham, B.S.; Kwong, P.D.
Deposited on : 2015-05-08
Resolution : 3.02 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
Mogul : 2022.3.0, CSD as543be (2022)
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

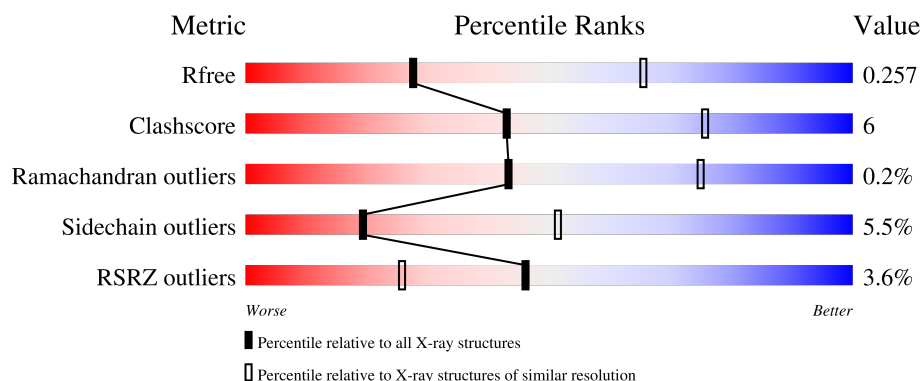
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.02 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	R	208	85% 14% .
1	S	208	6% 80% 16% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3388 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 Spike glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	208	Total	C	N	O	S	0	0	0
			1611	1029	256	315	11			
1	S	208	Total	C	N	O	S	0	0	0
			1611	1029	256	315	11			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	R	1	Total	C	N	O	0	0
			14	8	1	5		
2	R	1	Total	C	N	O	0	0
			14	8	1	5		
2	S	1	Total	C	N	O	0	0
			14	8	1	5		
2	S	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	S	1	Total	O	P	0	0
			5	4	1		

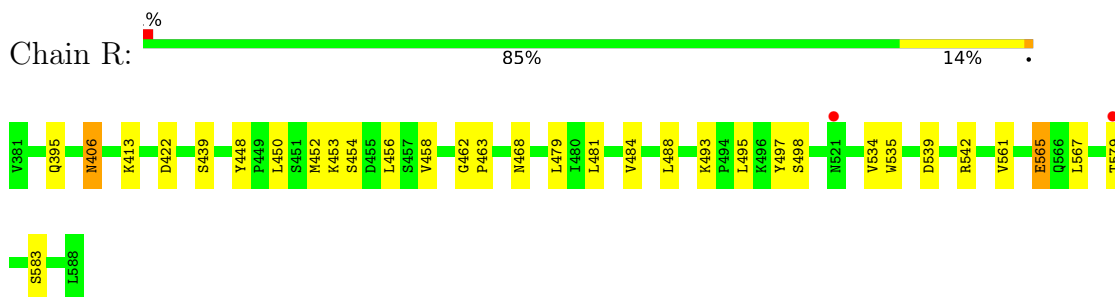
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	R	65	Total	O	0	0
			65	65		
4	S	40	Total	O	0	0
			40	40		

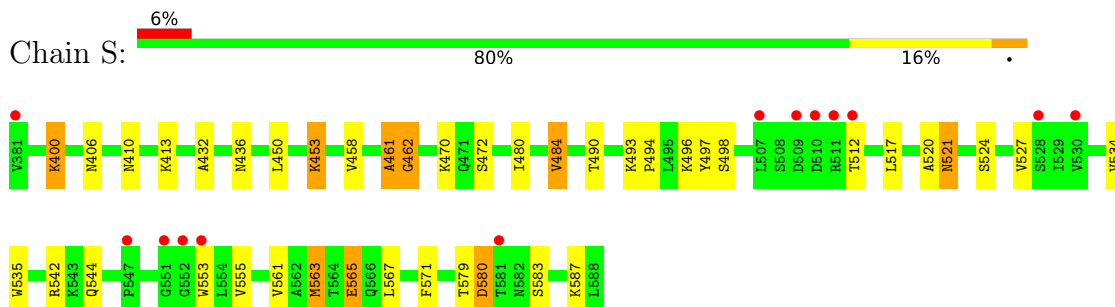
3 Residue-property plots

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: Spike glycoprotein



- Molecule 1: Spike glycoprotein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.67Å 109.89Å 125.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.73 – 3.02 43.73 – 3.02	Depositor EDS
% Data completeness (in resolution range)	90.5 (43.73-3.02) 90.5 (43.73-3.02)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.217 , 0.259 0.222 , 0.257	Depositor DCC
R_{free} test set	609 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	57.7	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 65.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3388	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, PO4

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	R	0.29	0/1651	0.71	2/2254 (0.1%)
1	S	0.30	0/1651	0.78	1/2254 (0.0%)
All	All	0.30	0/3302	0.75	3/4508 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	S	461	ALA	N-CA-C	6.31	120.98	113.16
1	R	462	GLY	CA-C-N	6.29	125.97	119.56
1	R	462	GLY	C-N-CA	6.29	125.97	119.56

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	R	1611	0	1570	17	0
1	S	1611	0	1570	28	0
2	R	28	0	26	1	0
2	S	28	0	26	1	0
3	S	5	0	0	0	0
4	R	65	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	S	40	0	0	0	0
All	All	3388	0	3192	40	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 (40) 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:S:461:ALA:N	1:S:462:GLY:HA3	2.08	0.68
1:S:542:ARG:HG2	1:S:555:VAL:HG22	1.79	0.64
1:R:422:ASP:HB3	1:R:481:LEU:HB2	1.80	0.63
1:R:456:LEU:HD13	1:R:479:LEU:HD21	1.84	0.60
1:S:480:ILE:HB	1:S:571:PHE:HB2	1.86	0.58
1:S:496:LYS:NZ	1:S:535:TRP:O	2.34	0.57
1:R:395:GLN:NE2	4:R:701:HOH:O	2.38	0.57
1:R:539:ASP:OD1	1:S:470:LYS:NZ	2.39	0.56
1:S:458:VAL:HB	1:S:461:ALA:HB3	1.87	0.56
1:R:406:ASN:HA	1:R:583:SER:HB2	1.89	0.55
1:S:472:SER:HA	1:S:520:ALA:HB1	1.88	0.54
1:R:535:TRP:CE2	1:S:517:LEU:HB2	2.43	0.53
1:S:410:ASN:HB2	1:S:587:LYS:HE3	1.91	0.53
1:S:432:ALA:O	1:S:436:ASN:ND2	2.39	0.52
1:S:494:PRO:HD2	1:S:563:MET:HE3	1.92	0.51
1:R:448:TYR:OH	1:R:452:MET:O	2.26	0.51
1:S:498:SER:HB3	1:S:534:VAL:HG23	1.93	0.50
1:R:493:LYS:NZ	1:R:565:GLU:O	2.42	0.50
1:R:463:PRO:O	1:R:468:ASN:ND2	2.44	0.50
1:S:400:LYS:HE3	1:S:400:LYS:HA	1.93	0.49
1:R:498:SER:HB3	1:R:534:VAL:HG23	1.94	0.49
1:S:406:ASN:HA	1:S:583:SER:HB2	1.95	0.49
1:R:413:LYS:HE3	2:R:602:NAG:H82	1.95	0.48
1:R:542:ARG:HB2	1:S:458:VAL:HA	1.95	0.47
1:S:453:LYS:HD3	1:S:453:LYS:C	2.40	0.47
1:S:493:LYS:NZ	1:S:565:GLU:O	2.37	0.47
1:S:410:ASN:HB3	1:S:413:LYS:HB2	1.98	0.46
1:S:497:TYR:HB2	1:S:561:VAL:HB	1.98	0.46
1:R:493:LYS:HG2	1:R:567:LEU:HB2	1.97	0.46
1:S:472:SER:HB3	1:S:521:ASN:OD1	2.16	0.45
1:S:453:LYS:HD3	1:S:453:LYS:O	2.17	0.45
1:R:535:TRP:HB2	1:S:517:LEU:HD12	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:535:TRP:CD2	1:S:517:LEU:HB2	2.52	0.43
1:R:453:LYS:HD3	1:R:481:LEU:HD22	1.99	0.43
1:S:484:VAL:HG13	1:S:567:LEU:O	2.18	0.43
1:S:524:SER:HB3	1:S:527:VAL:HG13	2.00	0.43
1:S:587:LYS:HD2	2:S:603:NAG:H82	2.02	0.42
1:R:497:TYR:HB2	1:R:561:VAL:HB	2.03	0.41
1:S:580:ASP:OD1	1:S:580:ASP:N	2.53	0.41
1:S:544:GLN:HB2	1:S:553:TRP:CH2	2.56	0.41

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	R	206/208 (99%)	201 (98%)	5 (2%)	0	100	100
1	S	206/208 (99%)	197 (96%)	8 (4%)	1 (0%)	24	59
All	All	412/416 (99%)	398 (97%)	13 (3%)	1 (0%)	43	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	S	462	GLY

5.3.2 Protein sidechains [i](#)

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	R	190/190 (100%)	180 (95%)	10 (5%)	20	52
1	S	190/190 (100%)	179 (94%)	11 (6%)	18	49
All	All	380/380 (100%)	359 (94%)	21 (6%)	19	51

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	406	ASN
1	R	439	SER
1	R	450	LEU
1	R	454	SER
1	R	458	VAL
1	R	484	VAL
1	R	488	LEU
1	R	495	LEU
1	R	565	GLU
1	R	579	THR
1	S	400	LYS
1	S	450	LEU
1	S	453	LYS
1	S	484	VAL
1	S	490	THR
1	S	512	THR
1	S	521	ASN
1	S	563	MET
1	S	565	GLU
1	S	579	THR
1	S	580	ASP

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

Mol	Chain	Res	Type
1	R	398	ASN
1	R	516	GLN
1	R	566	GLN
1	S	398	ASN
1	S	421	ASN
1	S	427	GLN
1	S	566	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NAG	R	601	1	14,14,15	0.48	0	17,19,21	0.52	0
3	PO4	S	601	-	4,4,4	0.97	0	6,6,6	0.47	0
2	NAG	R	602	1	14,14,15	0.76	1 (7%)	17,19,21	0.76	1 (5%)
2	NAG	S	603	1	14,14,15	0.17	0	17,19,21	0.43	0
2	NAG	S	602	1	14,14,15	0.35	0	17,19,21	0.52	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	R	601	1	-	0/6/23/26	0/1/1/1
2	NAG	S	603	1	-	1/6/23/26	0/1/1/1
2	NAG	S	602	1	-	3/6/23/26	0/1/1/1
2	NAG	R	602	1	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	R	602	NAG	C1-C2	2.64	1.55	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	602	NAG	C1-O5-C5	2.36	115.35	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	S	603	NAG	O5-C5-C6-O6
2	S	602	NAG	O5-C5-C6-O6
2	R	602	NAG	C3-C2-N2-C7
2	S	602	NAG	C3-C2-N2-C7
2	S	602	NAG	C1-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	R	602	NAG	1	0
2	S	603	NAG	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	R	208/208 (100%)	0.09	2 (0%) 79 59	29, 44, 82, 152	0
1	S	208/208 (100%)	0.49	13 (6%) 26 13	34, 61, 153, 216	0
All	All	416/416 (100%)	0.29	15 (3%) 46 26	29, 51, 112, 216	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	512	THR	4.0
1	S	552	GLY	3.5
1	S	509	ASP	3.3
1	S	507	LEU	3.1
1	S	528	SER	2.7
1	S	551	GLY	2.5
1	R	521	ASN	2.5
1	S	381	VAL	2.4
1	S	530	VAL	2.3
1	R	579	THR	2.3
1	S	510	ASP	2.2
1	S	581	THR	2.2
1	S	553	TRP	2.1
1	S	511	ARG	2.1
1	S	547	PRO	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](#)

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. 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	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	R	602	14/15	0.42	0.25	126,136,138,139	0
2	NAG	R	601	14/15	0.44	0.23	145,161,170,175	0
2	NAG	S	602	14/15	0.56	0.20	99,120,127,128	0
2	NAG	S	603	14/15	0.66	0.20	146,166,170,171	0
3	PO4	S	601	5/5	0.88	0.16	157,158,159,159	0

6.5 Other polymers [i](#)

There are no such residues in this entry.