



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 12:08 AM UTC

PDB ID : 5WWP / pdb_00005wwp
Title : Crystal structure of Middle East respiratory syndrome coronavirus helicase (MERS-CoV nsp13)
Authors : Hao, W.; Wojdyla, J.A.; Zhao, R.; Han, R.; Das, R.; Zlatev, I.; Manoharan, M.; Wang, M.; Cui, S.
Deposited on : 2017-01-03
Resolution : 3.00 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

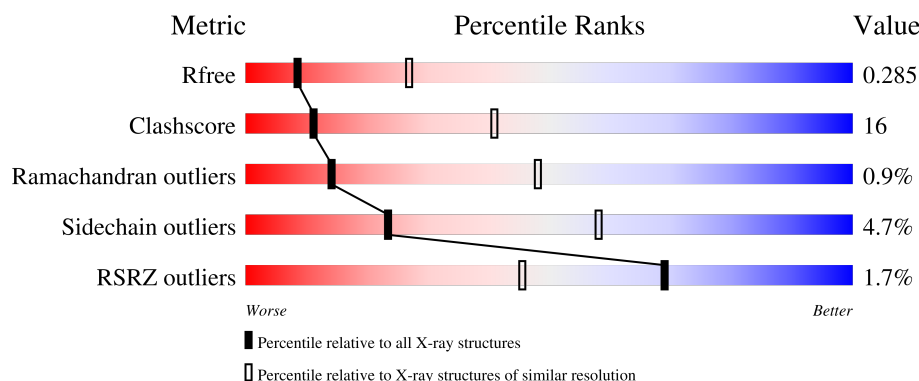
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	A	600	
1	B	600	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	A	602	-	-	X	-
2	ZN	B	602	-	-	X	-
2	ZN	B	603	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8571 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 ORF1ab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	591	Total	C	N	O	S	0	0	0
			4579	2917	767	861	34			
1	A	515	Total	C	N	O	S	0	0	0
			3961	2514	669	744	34			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	expression tag	UNP K0BWD0
B	0	PRO	-	expression tag	UNP K0BWD0
A	-1	GLY	-	expression tag	UNP K0BWD0
A	0	PRO	-	expression tag	UNP K0BWD0

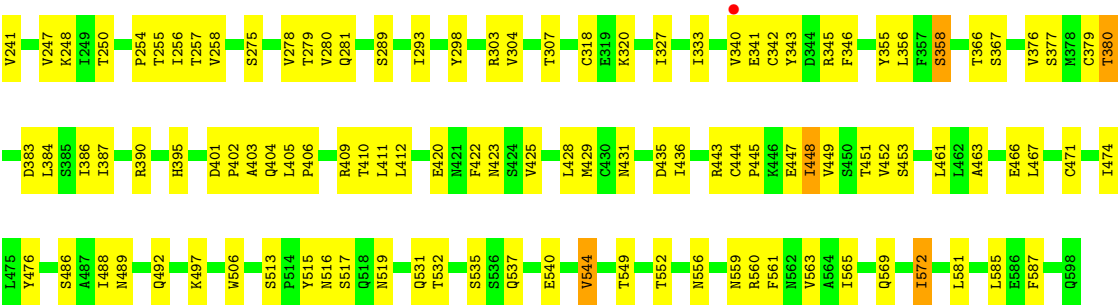
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Zn	0	0
			3	3		
2	A	3	Total	Zn	0	0
			3	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	185.68Å 185.68Å 185.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.95 – 3.00 48.95 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.95-3.00) 99.5 (48.95-3.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.232 , 0.277 0.240 , 0.285	Depositor DCC
R_{free} test set	1829 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	82.9	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.5	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	0.92	EDS
Total number of atoms	8571	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, SO4

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	0.71	0/4046	0.60	1/5504 (0.0%)
1	B	0.56	0/4679	0.54	1/6364 (0.0%)
All	All	0.63	0/8725	0.57	2/11868 (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	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	0	PRO	N-CA-CB	6.78	110.46	103.00
1	B	0	PRO	N-CA-CB	6.71	110.38	103.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	18	THR	Peptide
1	A	71	PHE	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	3961	0	3939	129	1
1	B	4579	0	4564	150	1
2	A	3	0	0	3	0
2	B	3	0	0	4	0
3	A	20	0	0	1	0
3	B	5	0	0	1	0
All	All	8571	0	8503	278	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 16.

All (278) 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:B:55:CYS:HG	2:B:603:ZN:ZN	0.73	1.00
1:B:19:CYS:HG	2:B:602:ZN:ZN	0.66	0.97
1:A:19:CYS:HG	2:A:602:ZN:ZN	0.63	0.93
1:B:534:ASP:OD2	1:B:560:ARG:NH2	2.05	0.89
1:A:94:LYS:NZ	1:A:95:ASN:OD1	2.11	0.84
1:A:513:SER:OG	1:A:519:ASN:OD1	1.97	0.80
1:A:486:SER:OG	1:A:517:SER:OG	1.97	0.78
1:B:19:CYS:SG	2:B:602:ZN:ZN	1.71	0.76
1:B:20:ILE:HG23	1:B:21:ARG:HG2	1.68	0.74
1:B:285:GLY:N	3:B:604:SO4:O1	2.17	0.74
1:A:405:LEU:HB2	1:A:537:GLN:OE1	1.88	0.72
1:A:8:CYS:SG	2:A:601:ZN:ZN	1.77	0.72
1:B:177:ASN:OD1	1:B:178:ARG:N	2.23	0.72
1:A:379:CYS:O	1:A:423:ASN:ND2	2.23	0.71
1:A:531:GLN:OE1	1:A:535:SER:OG	2.09	0.70
1:A:19:CYS:SG	2:A:602:ZN:ZN	1.77	0.70
1:A:257:THR:OG1	1:A:258:VAL:N	2.24	0.69
1:A:49:VAL:HG12	1:A:58:SER:HB3	1.75	0.69
1:B:113:ASP:OD2	1:B:115:THR:OG1	2.10	0.68
1:A:532:THR:HB	1:A:535:SER:H	1.58	0.68
1:A:8:CYS:HB2	1:A:10:SER:H	1.56	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:LYS:HG2	1:B:354:GLN:H	1.59	0.67
1:B:167:TRP:HZ2	1:B:176:LEU:HD21	1.60	0.67
1:B:513:SER:OG	1:B:519:ASN:OD1	2.11	0.66
1:B:366:THR:OG1	1:B:367:SER:N	2.27	0.66
1:A:380:THR:OG1	1:A:383:ASP:OD2	2.11	0.66
1:A:59:ASP:OD1	1:A:61:THR:N	2.28	0.66
1:B:333:ILE:HD11	1:B:356:LEU:HD13	1.78	0.65
1:A:404:GLN:O	1:A:559:ASN:ND2	2.29	0.64
1:B:451:THR:HG22	1:B:586:GLU:H	1.63	0.64
1:B:490:ARG:NH1	1:B:525:MET:SD	2.71	0.64
1:B:258:VAL:HG22	1:B:259:PRO:O	1.99	0.63
1:A:60:VAL:HA	1:A:63:LEU:HD12	1.82	0.62
1:B:201:GLU:O	1:B:210:SER:N	2.28	0.62
1:B:329:LYS:HD3	1:B:354:GLN:HB2	1.81	0.62
1:A:18:THR:O	1:A:18:THR:OG1	2.16	0.62
1:A:304:VAL:HG22	1:A:355:TYR:CD1	2.35	0.62
1:A:55:CYS:SG	1:A:56:GLY:HA3	2.40	0.61
1:B:257:THR:HG22	1:B:258:VAL:H	1.65	0.61
1:B:159:GLY:O	1:B:162:GLN:N	2.24	0.61
1:B:180:TYR:HD2	1:B:227:LEU:HD11	1.65	0.61
1:B:376:VAL:O	1:B:423:ASN:ND2	2.32	0.61
1:A:405:LEU:HD21	1:A:560:ARG:HG2	1.82	0.61
1:B:257:THR:HG22	1:B:258:VAL:HG12	1.83	0.61
1:A:18:THR:HG23	1:A:39:HIS:HB3	1.83	0.61
1:B:19:CYS:SG	1:B:39:HIS:ND1	2.74	0.60
1:B:197:GLU:N	1:B:214:SER:OG	2.32	0.60
1:B:335:PRO:HG3	1:B:362:ALA:HB2	1.82	0.60
1:A:515:TYR:O	1:A:519:ASN:ND2	2.33	0.60
1:A:448:ILE:HG22	1:A:587:PHE:CE1	2.35	0.60
1:B:519:ASN:O	1:B:522:SER:OG	2.11	0.60
1:A:279:THR:HG23	1:A:435:ASP:H	1.67	0.60
1:A:532:THR:H	1:A:535:SER:HB3	1.66	0.60
1:A:552:THR:O	1:A:556:ASN:ND2	2.34	0.60
1:B:333:ILE:HB	1:B:358:SER:HB2	1.84	0.59
1:A:333:ILE:HB	1:A:358:SER:HB2	1.85	0.59
1:A:32:ASP:HA	1:A:35:ILE:HD12	1.85	0.59
1:A:420:GLU:OE1	1:A:431:ASN:ND2	2.34	0.59
1:A:489:ASN:HB2	1:A:549:THR:HG23	1.85	0.59
1:B:161:ARG:HB2	1:B:211:TYR:HD2	1.68	0.58
1:B:180:TYR:CD2	1:B:227:LEU:HD11	2.38	0.58
1:A:516:ASN:ND2	3:A:605:SO4:S	2.74	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:THR:HB	1:A:26:CYS:HA	1.84	0.58
1:B:37:THR:HG23	1:B:39:HIS:H	1.69	0.57
1:B:308:ALA:O	1:B:360:ILE:HG13	2.05	0.57
1:B:519:ASN:HD22	1:B:530:THR:HG1	1.53	0.57
1:A:448:ILE:HG22	1:A:587:PHE:CZ	2.40	0.56
1:A:31:TYR:O	1:A:35:ILE:HG13	2.05	0.56
1:A:303:ARG:HH12	1:A:356:LEU:HD13	1.70	0.56
1:A:447:GLU:O	1:A:451:THR:OG1	2.22	0.56
1:B:117:SER:HB3	1:B:415:GLY:HA3	1.88	0.56
1:A:12:THR:OG1	1:A:13:SER:N	2.38	0.56
1:A:449:VAL:HG11	1:A:463:ALA:HB2	1.88	0.56
1:B:544:VAL:HG23	1:B:572:ILE:HG13	1.88	0.55
1:B:561:PHE:CZ	1:B:565:ILE:HD11	2.41	0.55
1:B:18:THR:HG23	1:B:41:MET:HB3	1.87	0.55
1:B:371:LEU:HB3	1:B:396:ILE:HG12	1.88	0.55
1:A:32:ASP:HB2	1:A:103:ILE:HG12	1.89	0.55
1:A:30:CYS:O	1:A:34:VAL:HG22	2.07	0.55
1:B:55:CYS:SG	2:B:603:ZN:ZN	1.86	0.54
1:A:8:CYS:SG	1:A:29:CYS:SG	3.04	0.54
1:A:247:VAL:HG23	1:A:248:LYS:HG3	1.88	0.54
1:B:90:PHE:HA	1:B:94:LYS:NZ	2.22	0.54
1:A:366:THR:OG1	1:A:367:SER:N	2.40	0.54
1:B:305:VAL:HG22	1:B:356:LEU:HB2	1.90	0.54
1:B:453:SER:O	1:B:459:ASN:N	2.31	0.54
1:A:281:GLN:HG3	1:A:402:PRO:HD2	1.89	0.54
1:B:249:ILE:HG13	1:B:274:TYR:CZ	2.43	0.53
1:B:40:LYS:O	1:B:60:VAL:HG23	2.09	0.53
1:B:406:PRO:HB3	1:B:422:PHE:CZ	2.44	0.53
1:A:71:PHE:HD1	1:A:75:HIS:HB2	1.72	0.53
1:A:19:CYS:SG	1:A:33:HIS:NE2	2.81	0.53
1:B:489:ASN:OD1	1:B:492:GLN:HG2	2.09	0.53
1:B:473:LYS:HE3	1:B:582:PHE:CE2	2.44	0.53
1:A:405:LEU:HG	1:A:563:VAL:HG21	1.91	0.53
1:A:451:THR:HG21	1:A:585:LEU:HD23	1.91	0.53
1:B:31:TYR:CE2	1:B:87:GLY:HA2	2.44	0.52
1:B:65:LEU:HD23	1:B:81:PHE:CZ	2.45	0.52
1:A:59:ASP:OD1	1:A:59:ASP:C	2.53	0.52
1:A:425:VAL:O	1:A:429:MET:HG3	2.10	0.52
1:A:561:PHE:CZ	1:A:585:LEU:HD21	2.45	0.52
1:B:44:SER:OG	1:B:45:VAL:N	2.42	0.52
1:B:78:VAL:HG13	1:B:79:CYS:H	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:THR:HG22	1:A:41:MET:HB2	1.90	0.52
1:B:215:THR:HB	1:B:216:THR:HG22	1.91	0.52
1:A:31:TYR:CD2	1:A:35:ILE:HD11	2.45	0.52
1:B:19:CYS:SG	1:B:39:HIS:CE1	3.03	0.51
1:B:585:LEU:HD23	1:B:587:PHE:CE2	2.46	0.51
1:B:182:PHE:O	1:B:198:TYR:N	2.44	0.51
1:B:279:THR:HG23	1:B:435:ASP:H	1.76	0.51
1:A:5:CYS:C	1:A:7:VAL:H	2.18	0.51
1:A:343:TYR:CE2	1:A:345:ARG:HB2	2.46	0.51
1:B:167:TRP:CZ2	1:B:176:LEU:HD21	2.41	0.51
1:B:262:PHE:HE1	1:B:290:HIS:CD2	2.29	0.51
1:A:340:VAL:HG12	1:A:341:GLU:H	1.76	0.51
1:B:451:THR:CG2	1:B:586:GLU:H	2.23	0.51
1:B:425:VAL:O	1:B:429:MET:HG3	2.11	0.51
1:B:520:ALA:O	1:B:523:ARG:HB2	2.11	0.51
1:B:467:LEU:H	1:B:467:LEU:HD12	1.76	0.51
1:A:59:ASP:OD1	1:A:61:THR:HG22	2.11	0.51
1:B:105:GLU:HG2	1:B:122:LEU:HD11	1.93	0.51
1:B:155:LYS:HB2	1:B:166:VAL:HG21	1.93	0.50
1:B:487:ALA:O	1:B:517:SER:OG	2.29	0.50
1:A:410:THR:OG1	1:A:411:LEU:N	2.44	0.50
1:A:476:TYR:O	1:A:492:GLN:NE2	2.34	0.50
1:B:161:ARG:HB2	1:B:211:TYR:CD2	2.47	0.49
1:B:491:PRO:O	1:B:494:THR:HG22	2.12	0.49
1:A:445:PRO:HD3	1:A:569:GLN:O	2.11	0.49
1:B:447:GLU:O	1:B:451:THR:HG23	2.12	0.49
1:B:198:TYR:HD2	1:B:211:TYR:HE1	1.61	0.49
1:B:371:LEU:HD13	1:B:391:ILE:HG21	1.94	0.49
1:B:252:LEU:HB3	1:B:299:TYR:CD1	2.47	0.49
1:A:34:VAL:HG12	1:A:40:LYS:HA	1.94	0.49
1:B:244:GLU:OE1	1:B:246:TYR:OH	2.25	0.49
1:B:59:ASP:OD1	1:B:61:THR:OG1	2.27	0.48
1:B:121:THR:HG22	1:B:421:ASN:HD22	1.79	0.48
1:A:254:PRO:HB3	1:A:298:TYR:CE2	2.49	0.48
1:B:160:GLU:HA	1:B:161:ARG:HA	1.52	0.48
1:A:18:THR:CG2	1:A:39:HIS:HB3	2.44	0.48
1:B:258:VAL:HG23	1:B:262:PHE:HB2	1.96	0.48
1:B:48:TYR:OH	1:B:90:PHE:O	2.26	0.48
1:B:473:LYS:HG2	1:B:582:PHE:CE1	2.48	0.48
1:B:376:VAL:HG12	1:B:400:GLY:HA3	1.95	0.47
1:B:55:CYS:SG	1:B:75:HIS:CE1	3.07	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:TYR:HA	1:B:193:VAL:O	2.14	0.47
1:B:451:THR:O	1:B:455:LEU:HB2	2.13	0.47
1:A:304:VAL:HG22	1:A:355:TYR:HD1	1.79	0.47
1:A:532:THR:N	1:A:535:SER:HB3	2.29	0.47
1:B:138:LEU:O	1:B:142:GLU:HG3	2.15	0.47
1:A:250:THR:O	1:A:395:HIS:NE2	2.32	0.47
1:A:142:GLU:C	1:A:144:ALA:H	2.23	0.47
1:B:37:THR:CG2	1:B:39:HIS:H	2.28	0.47
1:B:122:LEU:O	1:B:126:THR:OG1	2.33	0.47
1:B:271:LYS:C	1:B:273:GLY:H	2.23	0.47
1:A:444:CYS:HB2	1:A:449:VAL:HG23	1.97	0.47
1:B:35:ILE:HD13	1:A:497:LYS:HG2	1.96	0.47
1:B:262:PHE:CE1	1:B:290:HIS:NE2	2.82	0.47
1:B:90:PHE:HA	1:B:94:LYS:HZ1	1.78	0.47
1:A:444:CYS:HB3	1:A:448:ILE:CG1	2.44	0.47
1:B:304:VAL:HG22	1:B:355:TYR:CD1	2.50	0.47
1:B:16:CYS:SG	1:B:19:CYS:SG	3.13	0.46
1:B:245:ARG:NH1	1:B:275:SER:OG	2.48	0.46
1:B:496:VAL:O	1:B:500:ILE:HG22	2.15	0.46
1:A:544:VAL:HG23	1:A:572:ILE:HG13	1.96	0.46
1:B:256:ILE:HA	1:B:257:THR:HA	1.62	0.46
1:B:280:VAL:HG22	1:B:436:ILE:HB	1.97	0.46
1:A:47:PRO:O	1:A:49:VAL:HG22	2.15	0.46
1:B:248:LYS:HA	1:B:274:TYR:HE1	1.80	0.46
1:A:383:ASP:O	1:A:387:ILE:HG13	2.16	0.46
1:B:258:VAL:CG2	1:B:262:PHE:HB2	2.45	0.46
1:B:350:GLU:C	1:B:352:ASN:H	2.24	0.46
1:B:406:PRO:HB3	1:B:422:PHE:CE2	2.51	0.46
1:A:78:VAL:HG23	1:A:79:CYS:H	1.81	0.46
1:B:60:VAL:HG12	1:B:84:CYS:HB2	1.96	0.46
1:B:72:CYS:HB3	1:B:75:HIS:H	1.81	0.46
1:B:266:VAL:HG13	1:B:298:TYR:CE1	2.51	0.46
1:A:516:ASN:HA	1:A:519:ASN:HD22	1.81	0.46
1:B:535:SER:OG	1:B:535:SER:O	2.33	0.46
1:A:327:ILE:HD12	1:A:327:ILE:H	1.80	0.46
1:A:489:ASN:OD1	1:A:492:GLN:HG3	2.15	0.46
1:A:585:LEU:HD22	1:A:587:PHE:CE2	2.50	0.46
1:A:471:CYS:O	1:A:587:PHE:HB3	2.16	0.45
1:A:142:GLU:O	1:A:144:ALA:N	2.49	0.45
1:B:350:GLU:O	1:B:352:ASN:N	2.49	0.45
1:A:47:PRO:O	1:A:49:VAL:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:ASP:OD1	1:A:32:ASP:N	2.47	0.45
1:B:215:THR:HB	1:B:216:THR:HA	1.97	0.45
1:B:471:CYS:HA	1:B:572:ILE:O	2.17	0.45
1:A:443:ARG:NE	1:A:540:GLU:OE1	2.49	0.45
1:A:279:THR:HG22	1:A:435:ASP:OD2	2.17	0.45
1:B:563:VAL:O	1:B:567:ARG:NE	2.49	0.45
1:B:12:THR:HB	1:B:26:CYS:HA	1.98	0.45
1:B:304:VAL:HG12	1:B:370:ILE:HB	1.98	0.45
1:B:498:ASN:HA	1:B:501:THR:HG22	1.98	0.45
1:A:55:CYS:SG	1:A:75:HIS:HE1	2.39	0.45
1:B:546:PHE:CE2	1:B:548:GLN:HB3	2.51	0.44
1:B:376:VAL:HG21	1:B:425:VAL:HG11	1.99	0.44
1:B:227:LEU:HD23	1:B:227:LEU:HA	1.86	0.44
1:A:16:CYS:SG	1:A:19:CYS:SG	3.14	0.44
1:B:175:PRO:HG2	1:B:180:TYR:CE1	2.52	0.44
1:A:448:ILE:O	1:A:452:VAL:HG22	2.18	0.44
1:B:479:ASN:H	1:B:491:PRO:HG2	1.82	0.44
1:A:37:THR:C	1:A:39:HIS:H	2.26	0.44
1:B:150:ALA:HB1	1:B:168:GLU:HG3	2.00	0.44
1:B:358:SER:OG	1:B:359:THR:O	2.36	0.44
1:B:505:ALA:O	1:B:507:SER:N	2.50	0.44
1:A:307:THR:HA	1:A:358:SER:O	2.18	0.44
1:B:500:ILE:HD11	1:B:507:SER:HA	2.00	0.43
1:A:72:CYS:C	1:A:74:ASP:N	2.76	0.43
1:A:139:ARG:O	1:A:139:ARG:HG3	2.16	0.43
1:A:506:TRP:CD1	1:A:506:TRP:N	2.86	0.43
1:B:42:VAL:CG2	1:B:60:VAL:HG21	2.48	0.43
1:A:8:CYS:SG	1:A:26:CYS:SG	3.13	0.43
1:B:323:LYS:HG3	1:B:324:TYR:CD1	2.52	0.43
1:B:519:ASN:ND2	1:B:530:THR:OG1	2.33	0.43
1:B:71:PHE:CD2	1:B:77:PRO:HD3	2.54	0.43
1:B:154:ILE:HD13	1:B:154:ILE:HA	1.91	0.43
1:B:384:LEU:HD11	1:B:423:ASN:HB3	1.99	0.43
1:B:451:THR:HG22	1:B:586:GLU:HG3	1.99	0.43
1:B:200:PHE:HB3	1:B:209:VAL:HG21	2.01	0.43
1:B:325:LEU:HD23	1:B:325:LEU:HA	1.73	0.43
1:A:255:THR:OG1	1:A:256:ILE:N	2.51	0.43
1:A:18:THR:HG1	1:A:39:HIS:CG	2.36	0.43
1:A:142:GLU:C	1:A:144:ALA:N	2.77	0.43
1:B:1:ALA:HA	1:B:15:ARG:NH2	2.34	0.43
1:A:108:ARG:CZ	1:A:122:LEU:HD22	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:THR:O	1:A:412:LEU:HD12	2.19	0.43
1:A:531:GLN:HB3	1:A:535:SER:HG	1.82	0.43
1:B:478:GLY:C	1:B:489:ASN:HD21	2.26	0.42
1:A:289:SER:HB3	1:A:320:LYS:HG3	2.01	0.42
1:A:386:ILE:O	1:A:390:ARG:HG3	2.20	0.42
1:B:32:ASP:HB2	1:B:103:ILE:HG12	2.00	0.42
1:A:121:THR:O	1:A:125:THR:OG1	2.36	0.42
1:A:488:ILE:HG12	1:A:517:SER:HB3	2.01	0.42
1:B:198:TYR:CD2	1:B:211:TYR:HE1	2.38	0.42
1:A:453:SER:HB2	1:A:461:LEU:HB3	2.02	0.42
1:A:63:LEU:HA	1:A:71:PHE:O	2.20	0.42
1:A:544:VAL:CG2	1:A:572:ILE:HG13	2.50	0.42
1:A:21:ARG:HD2	1:A:21:ARG:HA	1.82	0.42
1:B:94:LYS:HE2	1:B:94:LYS:HB3	1.88	0.42
1:A:280:VAL:HG22	1:A:436:ILE:HB	2.02	0.42
1:A:466:GLU:O	1:A:467:LEU:C	2.63	0.42
1:B:164:LEU:HD23	1:B:164:LEU:HA	1.85	0.42
1:B:548:GLN:NE2	1:B:576:MET:SD	2.93	0.42
1:A:37:THR:O	1:A:39:HIS:N	2.53	0.42
1:A:444:CYS:HB3	1:A:448:ILE:HD11	2.02	0.42
1:A:105:GLU:OE1	1:A:126:THR:HG22	2.20	0.41
1:B:453:SER:HA	1:B:457:TYR:HB2	2.02	0.41
1:A:78:VAL:HG23	1:A:79:CYS:N	2.35	0.41
1:B:1:ALA:HA	1:B:15:ARG:HH21	1.85	0.41
1:B:158:VAL:HG23	1:B:162:GLN:HB2	2.02	0.41
1:B:423:ASN:HB2	1:B:425:VAL:H	1.84	0.41
1:A:18:THR:OG1	1:A:39:HIS:CG	2.73	0.41
1:A:143:GLU:OE1	1:A:231:SER:N	2.53	0.41
1:A:240:ILE:HG21	1:A:240:ILE:HD13	1.74	0.41
1:B:446:LYS:HA	1:B:449:VAL:HG12	2.03	0.41
1:B:548:GLN:CD	1:B:577:THR:H	2.29	0.41
1:A:318:CYS:HB3	1:A:346:PHE:CE2	2.55	0.41
1:B:272:SER:HA	1:B:278:VAL:HG21	2.01	0.41
1:B:548:GLN:OE1	1:B:577:THR:HG22	2.21	0.41
1:A:5:CYS:C	1:A:7:VAL:N	2.76	0.41
1:A:581:LEU:HD12	1:A:581:LEU:HA	1.86	0.41
1:A:401:ASP:C	1:A:403:ALA:H	2.28	0.41
1:B:429:MET:HE2	1:B:429:MET:HB3	1.73	0.41
1:A:289:SER:O	1:A:293:ILE:HG12	2.21	0.41
1:B:92:LEU:HD23	1:B:92:LEU:HA	1.71	0.41
1:A:384:LEU:HD11	1:A:423:ASN:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:LEU:HD23	1:A:428:LEU:HA	1.90	0.41
1:B:544:VAL:CG2	1:B:568:ALA:HB2	2.51	0.40
1:B:446:LYS:HA	1:B:449:VAL:CG1	2.51	0.40
1:B:449:VAL:HG11	1:B:463:ALA:HB2	2.03	0.40
1:B:505:ALA:C	1:B:507:SER:N	2.79	0.40
1:A:26:CYS:HB2	1:A:97:CYS:HB3	2.03	0.40
1:A:31:TYR:CE1	1:A:35:ILE:HG12	2.55	0.40
1:A:340:VAL:HG12	1:A:341:GLU:N	2.36	0.40
1:A:561:PHE:CE2	1:A:565:ILE:HD11	2.57	0.40
1:B:326:ASN:O	1:B:328:ALA:N	2.54	0.40
1:A:409:ARG:NH2	1:A:422:PHE:O	2.55	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:B:79:CYS:CB	1:A:79:CYS:SG[8_445]	1.40	0.80

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	511/600 (85%)	458 (90%)	48 (9%)	5 (1%)	12	45
1	B	587/600 (98%)	532 (91%)	50 (8%)	5 (1%)	14	48
All	All	1098/1200 (92%)	990 (90%)	98 (9%)	10 (1%)	14	48

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	504	PRO
1	B	215	THR

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Mol	Chain	Res	Type
1	B	123	ALA
1	B	351	THR
1	B	506	TRP
1	A	45	VAL
1	A	126	THR
1	A	406	PRO
1	A	38	PRO
1	A	73	VAL

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	443/521 (85%)	417 (94%)	26 (6%)	18	50
1	B	512/521 (98%)	493 (96%)	19 (4%)	30	64
All	All	955/1042 (92%)	910 (95%)	45 (5%)	23	58

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	27	CYS
1	B	37	THR
1	B	72	CYS
1	B	78	VAL
1	B	79	CYS
1	B	97	CYS
1	B	126	THR
1	B	148	SER
1	B	203	ILE
1	B	220	THR
1	B	227	LEU
1	B	241	VAL
1	B	257	THR
1	B	359	THR
1	B	366	THR

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Mol	Chain	Res	Type
1	B	377	SER
1	B	494	THR
1	B	524	SER
1	B	573	LEU
1	A	2	VAL
1	A	6	VAL
1	A	18	THR
1	A	20	ILE
1	A	35	ILE
1	A	41	MET
1	A	50	CYS
1	A	55	CYS
1	A	95	ASN
1	A	102	SER
1	A	125	THR
1	A	126	THR
1	A	145	SER
1	A	236	THR
1	A	241	VAL
1	A	275	SER
1	A	278	VAL
1	A	342	CYS
1	A	358	SER
1	A	376	VAL
1	A	377	SER
1	A	380	THR
1	A	448	ILE
1	A	474	ILE
1	A	544	VAL
1	A	572	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	B	421	ASN
1	B	458	ASN
1	B	489	ASN
1	A	290	HIS
1	A	354	GLN
1	A	423	ASN
1	A	569	GLN

5.3.3 RNA [i](#)

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](#)

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 for Mogul analysis.

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	$\# Z > 2$
3	SO4	A	604	-	4,4,4	0.19	0	6,6,6	0.27	0
3	SO4	A	605	-	4,4,4	0.20	0	6,6,6	0.18	0
3	SO4	A	606	-	4,4,4	0.21	0	6,6,6	0.19	0
3	SO4	B	604	-	4,4,4	0.19	0	6,6,6	0.21	0
3	SO4	A	607	-	4,4,4	0.27	0	6,6,6	0.19	0

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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	605	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	604	SO4	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	A	515/600 (85%)	-0.18	5 (0%) 79 59	33, 54, 87, 140	0
1	B	591/600 (98%)	0.13	14 (2%) 59 36	39, 74, 113, 133	0
All	All	1106/1200 (92%)	-0.01	19 (1%) 69 45	33, 62, 109, 140	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	550	ALA	3.7
1	B	516	ASN	3.4
1	B	224	ILE	2.7
1	B	340	VAL	2.7
1	B	93	TYR	2.5
1	B	338	ALA	2.4
1	B	494	THR	2.4
1	A	39	HIS	2.4
1	B	165	LEU	2.4
1	B	596	LYS	2.3
1	B	207	ASP	2.3
1	B	483	ASP	2.1
1	B	96	MET	2.1
1	B	92	LEU	2.1
1	B	485	SER	2.1
1	A	340	VAL	2.1
1	A	56	GLY	2.1
1	A	57	VAL	2.0
1	A	49	VAL	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
3	SO4	B	604	5/5	0.62	0.14	89,94,101,120	0
3	SO4	A	605	5/5	0.70	0.28	85,94,108,118	0
3	SO4	A	604	5/5	0.80	0.31	73,79,87,113	0
3	SO4	A	606	5/5	0.86	0.10	74,79,87,107	0
3	SO4	A	607	5/5	0.91	0.09	70,71,73,79	0
2	ZN	A	603	1/1	0.95	0.08	167,167,167,167	0
2	ZN	A	602	1/1	0.98	0.07	50,50,50,50	0
2	ZN	B	603	1/1	0.99	0.05	52,52,52,52	0
2	ZN	B	601	1/1	0.99	0.04	45,45,45,45	0
2	ZN	A	601	1/1	1.00	0.05	52,52,52,52	0
2	ZN	B	602	1/1	1.00	0.04	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.