



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:36 PM UTC

PDB ID : 7D3C / pdb_00007d3c
Title : The newly emerged SARS-like coronavirus HCoV-EMC also has an "Achilles' heel": current effective inhibitor targeting a 3C-like protease
Authors : Yan, L.; Ren, Z.
Deposited on : 2020-09-18
Resolution : 2.20 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

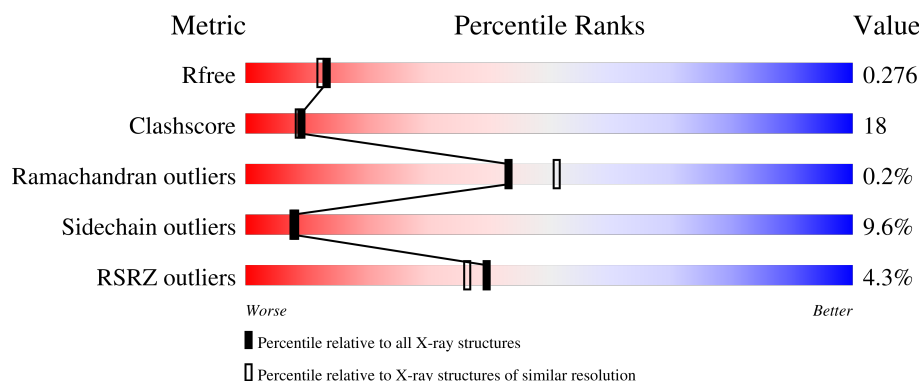
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	306	5% 70% 25% • •
1	B	306	4% 69% 24% 5% •
2	C	6	50% 50%
2	E	6	83% 17%

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	02J	C	1	-	X	-	-
2	PJE	C	5	X	-	-	-
2	02J	E	1	-	X	-	-
2	PJE	E	5	X	X	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4939 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 3C-like proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	299	Total	C	N	O	S	0	0	0
			2281	1448	377	431	25			
1	A	303	Total	C	N	O	S	0	0	0
			2305	1464	381	435	25			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	34	ASP	ASN	conflict	UNP A0A0D3MU45
A	34	ASP	ASN	conflict	UNP A0A0D3MU45

- Molecule 2 is a protein called N-[(5-METHYLISOXAZOL-3-YL)CARBONYL]ALANYL-L-VALYL-N 1 -((1R,2Z)-4-(BENZYLOXY)-4-OXO-1-{[(3R)-2-OXOPYRROLIDIN-3-YL]METHYL}BUT-2-ENYL)-L-LEUCINAMIDE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	6	Total	C	N	O	0	0	0
			49	35	6	8			
2	E	6	Total	C	N	O	0	0	0
			49	35	6	8			

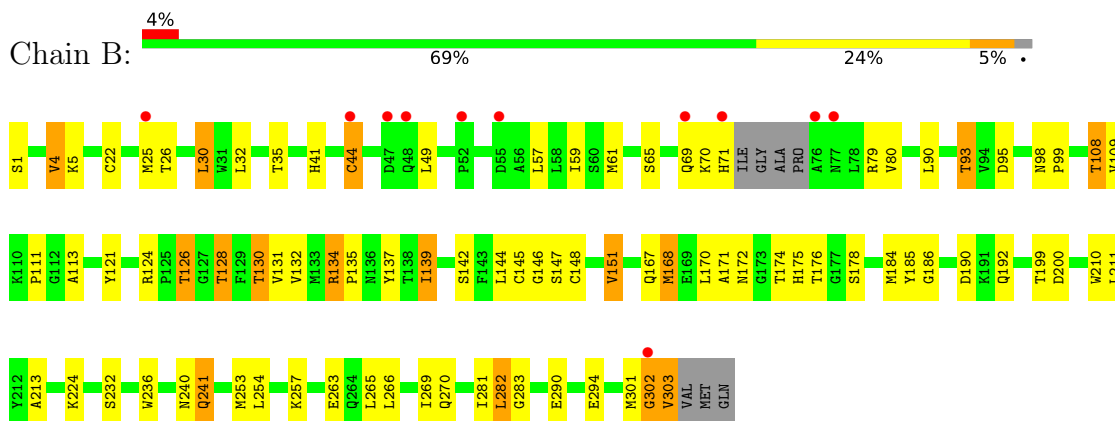
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	159	Total	O	0	0
			159	159		
3	A	94	Total	O	0	0
			94	94		
3	C	1	Total	O	0	0
			1	1		
3	E	1	Total	O	0	0
			1	1		

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: 3C-like proteinase



● Molecule 2: N-[(5-METHYLISOXAZOL-3-YL)CARBONYL]ALANYL-L-VALYL-N 1 -((1R,2Z)-4-(BENZYLOXY)-4-OXO-1-{[(3R)-2-OXOPYRROLIDIN-3-YL]METHYL}BUT-2-ENYL)-L-L EUCINAMIDE

Chain E:  83% 17%

02J1
A2
V3
L4
PJE5
0106

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.62Å 93.22Å 103.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.22 – 2.20 32.22 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (32.22-2.20) 99.7 (32.22-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.19 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, R_{free}	0.222 , 0.278 0.217 , 0.276	Depositor DCC
R_{free} test set	5116 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	31.5	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4939	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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: PJE, 010, 02J

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	0.45	0/2359	0.83	2/3211 (0.1%)
1	B	0.53	0/2333	0.88	2/3173 (0.1%)
2	C	3.28	2/19 (10.5%)	2.52	2/25 (8.0%)
2	E	3.42	2/19 (10.5%)	1.65	0/25
All	All	0.58	4/4730 (0.1%)	0.87	6/6434 (0.1%)

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	1
1	B	0	1
2	C	1	0
2	E	1	0
All	All	2	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	VAL	C-N	10.58	1.48	1.33
2	C	3	VAL	C-N	9.67	1.46	1.33
2	C	2	ALA	C-N	9.40	1.45	1.33
2	E	2	ALA	C-N	9.11	1.45	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	282	LEU	N-CA-C	7.28	121.57	111.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	168	MET	N-CA-C	5.78	117.32	108.42
1	A	168	MET	N-CA-C	5.77	117.62	108.79
2	C	2	ALA	CA-C-N	-5.71	116.11	123.19
2	C	2	ALA	C-N-CA	-5.71	116.11	123.19
1	A	283	GLY	N-CA-C	-5.50	97.34	113.30

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	5	PJE	C26
2	E	5	PJE	C26

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	282	LEU	Peptide
1	B	185	TYR	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	2305	0	2251	75	0
1	B	2281	0	2224	93	0
2	C	49	0	41	6	0
2	E	49	0	41	8	0
3	A	94	0	0	6	0
3	B	159	0	0	15	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
All	All	4939	0	4557	168	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 18.

All (168) 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:148:CYS:SG	2:C:5:PJE:C20	2.29	1.20
1:B:134:ARG:HH11	1:B:134:ARG:HG3	1.03	1.13
1:B:1:SER:N	3:B:402:HOH:O	1.86	1.08
1:B:1:SER:N	3:B:401:HOH:O	1.78	1.02
1:A:134:ARG:HH11	1:A:134:ARG:HG3	1.28	0.98
1:B:176:THR:HG23	1:B:184:MET:HE3	1.46	0.96
1:B:302:GLY:HA3	1:B:303:VAL:HG23	1.47	0.93
1:A:79:ARG:HH21	1:A:79:ARG:HG2	1.36	0.91
1:A:227:ARG:HB2	1:A:263:GLU:HB2	1.55	0.87
1:B:148:CYS:SG	2:E:5:PJE:C20	2.64	0.86
1:A:72:ILE:HD11	1:A:124:ARG:HB3	1.59	0.85
1:B:69:GLN:HG2	1:B:71:HIS:CE1	2.11	0.84
1:B:144:LEU:O	1:B:147:SER:HB2	1.78	0.82
1:B:168:MET:HE1	1:B:190:ASP:HA	1.63	0.81
1:A:134:ARG:HG3	1:A:134:ARG:NH1	1.93	0.81
1:B:253:MET:HE1	1:B:257:LYS:HE3	1.63	0.79
1:B:199:THR:HG22	3:B:456:HOH:O	1.80	0.79
1:A:303:VAL:O	3:A:401:HOH:O	2.00	0.78
1:B:134:ARG:HG3	1:B:134:ARG:NH1	1.81	0.76
1:A:22:CYS:SG	1:A:61:MET:HE1	2.29	0.73
1:A:133:MET:HE2	1:A:185:TYR:CD1	2.24	0.72
1:B:146:GLY:H	2:E:5:PJE:C	2.01	0.71
1:B:301:MET:O	1:B:302:GLY:C	2.33	0.71
1:A:22:CYS:SG	1:A:61:MET:CE	2.79	0.71
1:B:224:LYS:NZ	3:B:404:HOH:O	2.22	0.71
1:A:134:ARG:HH11	1:A:134:ARG:CG	2.04	0.71
1:B:61:MET:HE1	1:B:65:SER:HB3	1.73	0.70
1:B:302:GLY:CA	1:B:303:VAL:HG23	2.20	0.70
1:A:168:MET:HE2	2:C:2:ALA:HB1	1.73	0.70
1:A:86:GLN:NE2	3:A:402:HOH:O	2.12	0.70
1:A:133:MET:HE3	1:A:137:TYR:HA	1.75	0.69
1:B:253:MET:SD	3:B:543:HOH:O	2.51	0.69
1:A:93:THR:HG23	3:A:426:HOH:O	1.94	0.68
1:A:250:SER:HB2	1:A:297:ASN:ND2	2.08	0.68
1:B:302:GLY:HA3	1:B:303:VAL:CG2	2.24	0.67
1:A:95:ASP:OD1	1:A:96:VAL:HG23	1.93	0.67
1:A:79:ARG:HG2	1:A:79:ARG:NH2	1.99	0.66
1:B:128:THR:HG21	3:B:403:HOH:O	1.93	0.66
1:A:72:ILE:HD11	1:A:124:ARG:CB	2.25	0.66
1:A:135:PRO:HG2	1:A:201:LYS:O	1.95	0.66
1:A:262:ILE:O	1:A:266:LEU:HD13	1.95	0.66
1:B:22:CYS:SG	1:B:61:MET:HE1	2.36	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:MET:HG3	1:A:191:LYS:HG2	1.77	0.65
1:B:131:VAL:HG12	1:B:132:VAL:N	2.12	0.65
1:B:211:LEU:HD11	1:B:269:ILE:CD1	2.27	0.64
1:A:6:MET:HE1	1:A:298:MET:HE2	1.78	0.64
1:B:61:MET:CE	1:B:65:SER:HB3	2.28	0.63
1:B:170:LEU:HD12	1:B:170:LEU:N	2.12	0.62
1:A:117:VAL:O	1:A:128:THR:HG22	1.99	0.62
1:B:30:LEU:HG	1:B:151:VAL:HG21	1.82	0.62
1:B:93:THR:CG2	3:B:431:HOH:O	2.47	0.62
1:A:146:GLY:H	2:C:5:PJE:C	2.14	0.61
1:A:223:VAL:C	1:A:224:LYS:HD3	2.26	0.60
1:B:26:THR:O	2:E:6:010:H5	2.01	0.60
1:B:109:VAL:HG23	1:B:113:ALA:HB3	1.84	0.59
1:B:124:ARG:HB3	1:B:124:ARG:NH1	2.17	0.59
1:A:228:THR:O	1:A:263:GLU:HB3	2.03	0.59
1:B:171:ALA:O	1:B:174:THR:HG22	2.02	0.59
1:B:170:LEU:HD11	1:B:176:THR:HG22	1.85	0.58
1:A:6:MET:HE1	1:A:298:MET:CE	2.33	0.58
1:A:30:LEU:HD13	1:A:32:LEU:HG	1.86	0.57
1:A:294:GLU:O	1:A:298:MET:HB2	2.04	0.57
1:B:301:MET:O	1:B:302:GLY:O	2.22	0.57
1:A:242:PHE:HE1	1:A:273:TYR:CD1	2.23	0.56
1:B:211:LEU:HD11	1:B:269:ILE:HD11	1.86	0.56
1:B:71:HIS:CE1	3:B:448:HOH:O	2.58	0.56
1:A:290:GLU:HG3	3:A:458:HOH:O	2.05	0.55
1:B:109:VAL:CG2	1:B:113:ALA:HB3	2.36	0.55
1:A:167:GLN:HB2	1:A:177:GLY:HA2	1.87	0.55
1:A:200:ASP:C	1:A:201:LYS:HG2	2.31	0.55
2:C:5:PJE:O	2:C:6:010:H5	2.06	0.55
1:B:168:MET:HE3	1:B:184:MET:HE2	1.89	0.55
1:B:167:GLN:HB2	1:B:184:MET:HE1	1.89	0.55
1:B:130:THR:HG21	3:B:405:HOH:O	2.06	0.54
1:B:26:THR:O	2:E:6:010:C5	2.56	0.54
1:A:6:MET:CE	1:A:298:MET:HE2	2.37	0.54
1:B:22:CYS:SG	1:B:61:MET:CE	2.96	0.53
1:B:282:LEU:N	1:B:283:GLY:HA2	2.23	0.53
1:A:22:CYS:SG	1:A:61:MET:HE2	2.47	0.53
1:A:133:MET:CE	1:A:137:TYR:HA	2.38	0.53
1:B:148:CYS:SG	2:E:5:PJE:C21	2.97	0.53
1:B:93:THR:HG22	3:B:431:HOH:O	2.08	0.53
1:B:124:ARG:HB3	1:B:124:ARG:CZ	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:MET:C	1:A:298:MET:HE3	2.34	0.53
1:A:61:MET:CE	1:A:65:SER:HB3	2.38	0.53
1:B:142:SER:HB2	1:A:4:VAL:HG22	1.92	0.52
1:B:139:ILE:HG22	1:B:175:HIS:HB2	1.91	0.52
1:B:131:VAL:HG11	1:B:139:ILE:HG13	1.91	0.52
1:B:61:MET:CE	1:B:65:SER:CB	2.87	0.52
1:B:131:VAL:CG1	3:B:450:HOH:O	2.58	0.52
1:B:108:THR:HG23	3:B:489:HOH:O	2.08	0.52
1:B:168:MET:HE3	1:B:184:MET:CE	2.39	0.52
1:B:44:CYS:SG	1:B:49:LEU:HD23	2.50	0.52
1:B:41:HIS:CD2	2:E:4:LEU:HD22	2.44	0.52
1:A:121:TYR:CE1	1:A:147:SER:HB3	2.45	0.52
1:B:240:ASN:C	1:B:241:GLN:HG2	2.36	0.51
1:B:93:THR:HG23	3:B:431:HOH:O	2.10	0.51
1:A:211:LEU:HD11	1:A:269:ILE:HD11	1.93	0.51
1:B:30:LEU:HD13	1:B:32:LEU:HG	1.91	0.51
1:B:131:VAL:HG11	1:B:139:ILE:CG1	2.41	0.50
1:A:211:LEU:HD11	1:A:269:ILE:CD1	2.42	0.50
1:A:298:MET:HE3	1:A:299:GLN:N	2.26	0.50
1:B:137:TYR:HB3	1:B:186:GLY:H	1.76	0.50
1:B:211:LEU:HD11	1:B:269:ILE:HD13	1.92	0.50
1:B:61:MET:HE3	1:B:65:SER:CB	2.42	0.50
1:A:40:ARG:HD3	1:A:88:THR:HA	1.93	0.49
1:B:121:TYR:HD1	1:B:126:THR:HG21	1.77	0.49
1:A:133:MET:HE1	1:A:185:TYR:HB3	1.94	0.49
1:B:61:MET:HE3	1:B:65:SER:HB2	1.94	0.49
1:B:111:PRO:HB3	1:B:135:PRO:HA	1.94	0.49
1:A:98:ASN:HB3	1:A:101:THR:OG1	2.13	0.49
1:A:265:LEU:O	1:A:269:ILE:HG12	2.12	0.49
1:A:134:ARG:HB2	1:A:138:THR:O	2.13	0.48
1:A:215:ILE:HD13	1:A:260:VAL:HG21	1.94	0.48
1:A:290:GLU:OE1	3:A:403:HOH:O	2.20	0.48
1:A:109:VAL:HG22	1:A:113:ALA:HB3	1.96	0.47
1:B:170:LEU:N	1:B:170:LEU:CD1	2.77	0.47
1:B:265:LEU:O	1:B:269:ILE:HG12	2.14	0.47
1:B:211:LEU:CD1	1:B:269:ILE:HD11	2.44	0.47
1:B:254:LEU:HD23	1:B:301:MET:HE2	1.96	0.47
1:B:44:CYS:CB	1:B:57:LEU:HD11	2.45	0.47
1:B:108:THR:HG22	3:B:544:HOH:O	2.13	0.47
1:B:240:ASN:O	1:B:241:GLN:HG2	2.15	0.47
1:A:49:LEU:HB3	1:A:192:GLN:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ARG:O	1:A:126:THR:HG23	2.15	0.47
1:B:131:VAL:CG1	1:B:132:VAL:N	2.78	0.46
1:A:109:VAL:CG2	1:A:113:ALA:HB3	2.45	0.46
1:B:171:ALA:O	1:B:172:ASN:C	2.58	0.46
1:A:223:VAL:O	1:A:224:LYS:HD3	2.16	0.46
1:A:211:LEU:CD1	1:A:269:ILE:HD11	2.47	0.45
1:A:33:ASP:O	1:A:97:ALA:HA	2.16	0.45
1:B:210:TRP:O	1:B:213:ALA:HB3	2.17	0.45
1:A:61:MET:HE3	1:A:65:SER:CB	2.46	0.45
1:B:57:LEU:O	1:B:61:MET:HG2	2.16	0.45
1:B:167:GLN:CB	1:B:184:MET:HE1	2.47	0.45
1:B:148:CYS:SG	2:E:5:PJE:CA	3.06	0.44
1:A:297:ASN:HA	1:A:301:MET:HE2	1.99	0.44
1:B:236:TRP:CG	1:B:270:GLN:HE21	2.35	0.44
1:A:266:LEU:CD1	1:A:266:LEU:N	2.81	0.44
1:B:79:ARG:HD2	1:B:95:ASP:OD2	2.18	0.44
1:B:25:MET:HE1	2:E:6:010:HA	2.00	0.43
1:B:176:THR:CG2	1:B:184:MET:HE3	2.33	0.43
1:A:227:ARG:NH1	1:A:263:GLU:HG3	2.33	0.43
1:A:242:PHE:CE1	1:A:273:TYR:CD1	3.06	0.43
1:B:131:VAL:CG1	1:B:139:ILE:HG13	2.47	0.43
1:A:294:GLU:H	1:A:294:GLU:CD	2.25	0.43
1:A:111:PRO:HB3	1:A:135:PRO:HA	2.01	0.43
2:C:4:LEU:HD12	2:C:4:LEU:HA	1.89	0.42
1:B:49:LEU:HB3	1:B:192:GLN:HG3	1.99	0.42
1:B:290:GLU:HG3	3:B:519:HOH:O	2.19	0.42
1:B:134:ARG:NH1	1:B:200:ASP:OD1	2.52	0.42
1:B:281:ILE:C	1:B:283:GLY:HA2	2.44	0.42
1:B:59:ILE:HD13	1:B:59:ILE:HA	1.88	0.42
1:B:134:ARG:N	1:B:134:ARG:HD2	2.34	0.42
1:B:4:VAL:HG11	1:A:129:PHE:CG	2.54	0.42
1:A:250:SER:HB2	1:A:297:ASN:HD22	1.81	0.42
1:A:214:ALA:HA	1:A:282:LEU:HD11	2.01	0.41
1:A:224:LYS:HD3	1:A:224:LYS:N	2.30	0.41
1:B:98:ASN:HA	1:B:99:PRO:HD2	1.85	0.41
1:B:254:LEU:HD23	1:B:301:MET:CE	2.51	0.41
1:A:169:GLU:HB2	2:C:5:PJE:C29	2.51	0.41
1:A:227:ARG:HH11	1:A:263:GLU:HG3	1.86	0.41
1:B:61:MET:HE3	1:B:61:MET:HB3	1.82	0.41
1:A:44:CYS:SG	1:A:52:PRO:HB3	2.61	0.41
1:A:231:VAL:HG13	3:A:410:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:LEU:HD12	1:A:90:LEU:HA	1.77	0.41
1:A:79:ARG:HH21	1:A:79:ARG:CG	2.14	0.40
1:B:168:MET:CE	1:B:190:ASP:HA	2.43	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	301/306 (98%)	295 (98%)	6 (2%)	0	100	100
1	B	295/306 (96%)	282 (96%)	12 (4%)	1 (0%)	36	42
2	C	3/6 (50%)	3 (100%)	0	0	100	100
2	E	3/6 (50%)	3 (100%)	0	0	100	100
All	All	602/624 (96%)	583 (97%)	18 (3%)	1 (0%)	43	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	302	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	A	254/257 (99%)	230 (91%)	24 (9%)	8	9
1	B	252/257 (98%)	228 (90%)	24 (10%)	8	8
2	C	2/2 (100%)	1 (50%)	1 (50%)	0	0
2	E	2/2 (100%)	2 (100%)	0	100	100
All	All	510/518 (98%)	461 (90%)	49 (10%)	8	8

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	4	VAL
1	B	5	LYS
1	B	30	LEU
1	B	35	THR
1	B	44	CYS
1	B	70	LYS
1	B	80	VAL
1	B	90	LEU
1	B	93	THR
1	B	108	THR
1	B	126	THR
1	B	128	THR
1	B	130	THR
1	B	134	ARG
1	B	139	ILE
1	B	145	CYS
1	B	151	VAL
1	B	178	SER
1	B	232	SER
1	B	241	GLN
1	B	263	GLU
1	B	266	LEU
1	B	294	GLU
1	B	303	VAL
1	A	27	LEU
1	A	30	LEU
1	A	35	THR
1	A	59	ILE
1	A	64	HIS
1	A	67	SER
1	A	69	GLN
1	A	72	ILE
1	A	80	VAL

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Mol	Chain	Res	Type
1	A	90	LEU
1	A	128	THR
1	A	130	THR
1	A	134	ARG
1	A	174	THR
1	A	224	LYS
1	A	244	GLU
1	A	253	MET
1	A	257	LYS
1	A	263	GLU
1	A	270	GLN
1	A	271	GLN
1	A	294	GLU
1	A	298	MET
1	A	303	VAL
2	C	4	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	19	GLN
1	B	77	ASN
1	B	86	GLN
1	B	161	ASN
1	B	271	GLN
1	B	297	ASN
1	A	48	GLN
1	A	240	ASN
1	A	271	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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	PJE	C	5	2	11,13,14	7.01	7 (63%)	11,16,18	5.91	6 (54%)
2	02J	E	1	2	8,8,9	2.10	3 (37%)	9,10,12	4.13	6 (66%)
2	02J	C	1	2	8,8,9	2.09	4 (50%)	9,10,12	3.95	6 (66%)
2	PJE	E	5	2	11,13,14	7.23	8 (72%)	11,16,18	5.97	6 (54%)

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	PJE	C	5	2	1/1/3/7	6/7/18/19	0/1/1/1
2	02J	E	1	2	-	1/1/2/4	0/1/1/1
2	02J	C	1	2	-	1/1/2/4	0/1/1/1
2	PJE	E	5	2	1/1/3/7	7/7/18/19	0/1/1/1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	5	PJE	C29-N6	21.45	1.56	1.33
2	C	5	PJE	C29-N6	20.72	1.55	1.33
2	C	5	PJE	O-C	5.74	1.39	1.22
2	E	5	PJE	O-C	5.65	1.39	1.22
2	E	5	PJE	C21-C20	4.83	1.51	1.33
2	C	5	PJE	O8-C29	4.79	1.32	1.23
2	E	5	PJE	O8-C29	4.76	1.32	1.23
2	C	5	PJE	C21-C20	4.66	1.50	1.33
2	E	1	02J	CA-N	4.24	1.37	1.31
2	E	5	PJE	C26-C29	-4.18	1.41	1.52
2	C	1	02J	CA-N	4.09	1.37	1.31
2	C	5	PJE	C26-C29	-3.94	1.42	1.52
2	E	1	02J	C4-C5	2.93	1.40	1.34
2	C	1	02J	C4-C5	2.85	1.39	1.34
2	E	5	PJE	C27-C28	-2.74	1.49	1.53
2	C	5	PJE	C27-C28	-2.73	1.49	1.53
2	E	5	PJE	C21-C	2.41	1.51	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	5	PJE	C27-C26	-2.39	1.47	1.54
2	C	5	PJE	C27-C26	-2.36	1.47	1.54
2	C	1	02J	O1-N	-2.30	1.38	1.42
2	E	1	02J	O1-N	-2.10	1.38	1.42
2	C	1	02J	C-CA	2.07	1.53	1.45

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	5	PJE	O8-C29-C26	-13.25	110.89	126.21
2	C	5	PJE	O8-C29-N6	-11.75	109.33	125.76
2	E	5	PJE	O8-C29-N6	-10.81	110.65	125.76
2	C	5	PJE	O8-C29-C26	-10.62	113.93	126.21
2	E	1	02J	O1-C5-C6	8.33	127.35	115.89
2	C	1	02J	O1-C5-C6	8.06	126.97	115.89
2	C	5	PJE	C20-CA-N	-6.59	96.77	110.24
2	E	5	PJE	CA-C20-C21	-5.64	115.78	124.42
2	E	5	PJE	O-C-C21	-5.60	107.76	125.61
2	C	5	PJE	C25-CA-C20	5.27	119.05	110.99
2	C	5	PJE	O-C-C21	-5.24	108.88	125.61
2	C	5	PJE	C28-N6-C29	-5.21	102.73	113.79
2	E	1	02J	C6-C5-C4	-4.67	125.98	134.05
2	E	1	02J	C-CA-N	4.58	126.92	118.78
2	C	1	02J	C6-C5-C4	-4.57	126.15	134.05
2	E	5	PJE	C28-N6-C29	-4.43	104.39	113.79
2	C	1	02J	C-CA-N	4.21	126.25	118.78
2	E	1	02J	C4-CA-N	-3.95	107.58	111.19
2	C	1	02J	C4-CA-N	-3.74	107.77	111.19
2	E	5	PJE	C25-CA-C20	3.54	116.41	110.99
2	E	1	02J	C5-O1-N	3.19	111.72	108.47
2	C	1	02J	C5-O1-N	3.10	111.63	108.47
2	E	1	02J	O1-C5-C4	-2.86	106.67	109.88
2	C	1	02J	O1-C5-C4	-2.68	106.87	109.88

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	5	PJE	C26
2	E	5	PJE	C26

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	02J	O-C-CA-C4
2	E	1	02J	O-C-CA-C4
2	C	5	PJE	C26-C25-CA-N
2	C	5	PJE	CA-C20-C21-C
2	C	5	PJE	O-C-C21-C20
2	C	5	PJE	CA-C25-C26-C27
2	C	5	PJE	CA-C25-C26-C29
2	E	5	PJE	C26-C25-CA-C20
2	E	5	PJE	C26-C25-CA-N
2	E	5	PJE	O-C-C21-C20
2	E	5	PJE	CA-C25-C26-C27
2	E	5	PJE	CA-C25-C26-C29
2	E	5	PJE	CA-C20-C21-C
2	E	5	PJE	C21-C20-CA-C25
2	C	5	PJE	C26-C25-CA-C20

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	5	PJE	4	0
2	E	5	PJE	4	0

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	303/306 (99%)	0.48	15 (4%) 34 31	24, 39, 55, 79	0
1	B	299/306 (97%)	0.17	11 (3%) 45 42	21, 31, 55, 74	0
2	C	3/6 (50%)	0.11	0 100 100	34, 34, 35, 41	0
2	E	3/6 (50%)	0.74	0 100 100	40, 40, 41, 49	0
All	All	608/624 (97%)	0.33	26 (4%) 40 36	21, 36, 55, 79	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	76	ALA	6.2
1	B	77	ASN	3.5
1	A	76	ALA	3.4
1	A	72	ILE	3.3
1	A	303	VAL	3.2
1	B	71	HIS	3.0
1	A	273	TYR	3.0
1	A	267	TYR	2.9
1	A	302	GLY	2.8
1	A	64	HIS	2.6
1	A	25	MET	2.5
1	B	55	ASP	2.5
1	B	44	CYS	2.4
1	A	75	PRO	2.4
1	B	302	GLY	2.4
1	B	69	GLN	2.4
1	A	225	PRO	2.3
1	B	25	MET	2.3
1	B	52	PRO	2.2
1	B	47	ASP	2.2
1	A	1	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	74	ALA	2.1
1	B	48	GLN	2.0
1	A	199	THR	2.0
1	A	79	ARG	2.0
1	A	257	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	02J	E	1	8/9	0.51	0.20	55,64,74,75	0
2	02J	C	1	8/9	0.74	0.18	51,55,63,65	0
2	PJE	E	5	13/14	0.81	0.16	34,39,50,54	0
2	PJE	C	5	13/14	0.87	0.13	27,32,45,47	0

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.