



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:16 AM UTC

PDB ID : 9I9S / pdb_00009i9s
Title : Room temperature structure of PBP2a at the LCLS
Authors : Grieco, A.; Botha, S.; Martin-Garcia, J.M.
Deposited on : 2025-02-07
Resolution : 2.80 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

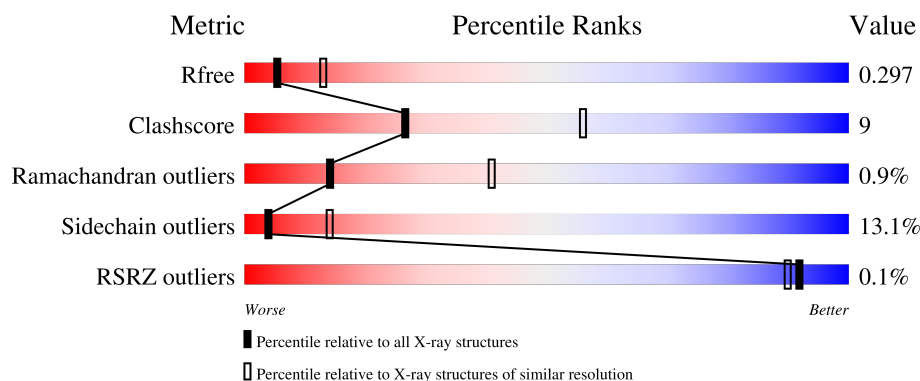
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	643	 69% 27% .
1	B	643	 71% 25% . .

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
3	CL	B	704	-	-	X	-
3	CL	B	706	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10400 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 MecA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	643	Total	C	N	O	S	0	1	0
			5169	3257	873	1023	16			
1	B	638	Total	C	N	O	S	0	1	0
			5135	3237	866	1017	15			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	246	GLU	GLY	conflict	UNP B3VQ68
A	664	TYR	-	expression tag	UNP B3VQ68
A	665	ASP	-	expression tag	UNP B3VQ68
A	666	ILE	-	expression tag	UNP B3VQ68
A	667	ASP	-	expression tag	UNP B3VQ68
A	668	GLU	-	expression tag	UNP B3VQ68
B	246	GLU	GLY	conflict	UNP B3VQ68
B	664	TYR	-	expression tag	UNP B3VQ68
B	665	ASP	-	expression tag	UNP B3VQ68
B	666	ILE	-	expression tag	UNP B3VQ68
B	667	ASP	-	expression tag	UNP B3VQ68
B	668	GLU	-	expression tag	UNP B3VQ68

- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Cd	0	0
			4	4		
2	B	3	Total	Cd	0	0
			3	3		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	0
			2	2		
3	B	5	Total	Cl	0	0
			5	5		

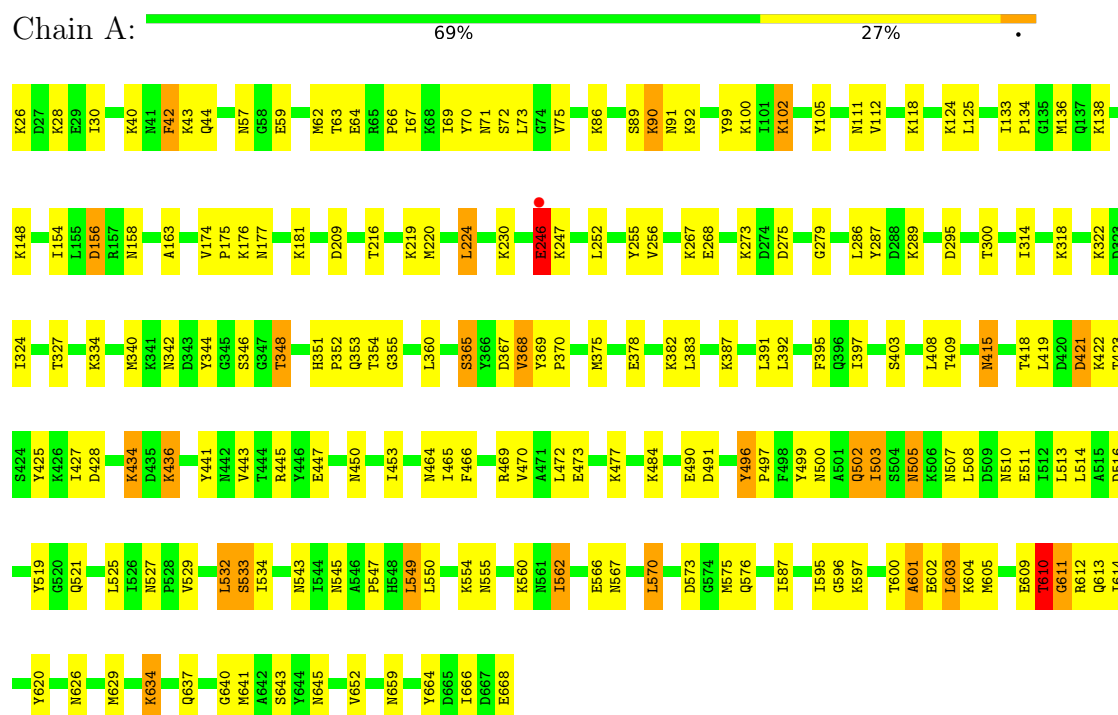
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	43	Total	O	0	0
			43	43		
4	B	39	Total	O	0	0
			39	39		

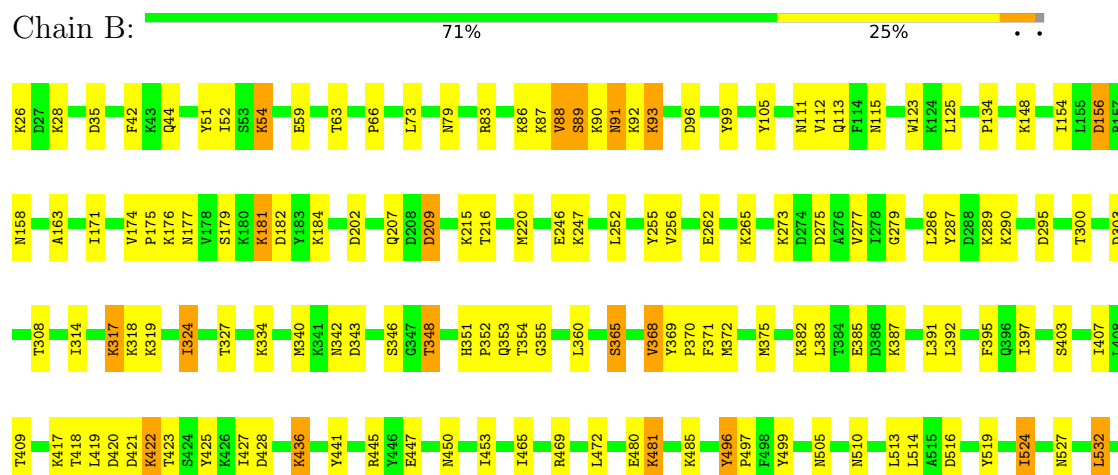
3 Residue-property plots

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: MecA



• Molecule 1: MecA





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.40Å 107.70Å 185.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.27 – 2.80 35.27 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.0 (35.27-2.80) 99.0 (35.27-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.220 , 0.290 0.230 , 0.297	Depositor DCC
R_{free} test set	2040 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	78.7	Xtriage
Anisotropy	0.150	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 78.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10400	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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: CD, CL

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	0.49	0/5255	1.06	2/7065 (0.0%)
1	B	0.49	0/5220	1.06	6/7018 (0.1%)
All	All	0.49	0/10475	1.06	8/14083 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	ASP	CA-CB-CG	6.38	118.98	112.60
1	A	573	ASP	CA-CB-CG	6.13	118.73	112.60
1	B	209	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	436	LYS	CB-CA-C	5.19	119.69	109.72
1	B	308	THR	CA-CB-OG1	-5.11	101.94	109.60
1	A	209	ASP	CA-CB-CG	5.08	117.69	112.60
1	B	202	ASP	CA-CB-CG	5.02	117.62	112.60
1	B	420	ASP	CA-CB-CG	5.00	117.60	112.60

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	5169	0	5165	97	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	5135	0	5131	94	0
2	A	4	0	0	0	0
2	B	3	0	0	0	0
3	A	2	0	0	1	0
3	B	5	0	0	9	0
4	A	43	0	0	9	0
4	B	39	0	0	9	0
All	All	10400	0	10296	191	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 9.

All (191) 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:30:ILE:HA	4:A:808:HOH:O	1.38	1.18
1:B:35:ASP:HB3	4:B:801:HOH:O	1.54	1.07
1:B:548:HIS:CE1	3:B:704:CL:CL	2.61	0.91
1:A:268:GLU:OE2	4:A:801:HOH:O	1.94	0.85
1:A:219:LYS:HG2	1:A:220:MET:H	1.45	0.82
1:A:562:ILE:HB	4:A:818:HOH:O	1.88	0.74
1:A:219:LYS:HG2	1:A:220:MET:N	2.02	0.74
1:A:133:ILE:HB	1:A:136:MET:HE2	1.68	0.73
1:B:220:MET:HA	1:B:220:MET:HE2	1.71	0.72
1:B:275:ASP:HB2	4:B:806:HOH:O	1.91	0.70
1:B:327:THR:OG1	1:B:549:LEU:HA	1.93	0.69
1:A:71:ASN:OD1	1:A:72:SER:N	2.26	0.68
1:A:219:LYS:CG	1:A:220:MET:H	2.06	0.68
1:A:327:THR:OG1	1:A:549:LEU:HA	1.93	0.67
1:A:352:PRO:HD2	1:A:353:GLN:OE1	1.94	0.67
1:A:441:TYR:OH	1:A:516:ASP:OD1	2.13	0.67
1:B:441:TYR:OH	1:B:516:ASP:OD1	2.12	0.67
1:A:434:LYS:NZ	1:A:511:GLU:OE1	2.28	0.66
1:B:255:TYR:HB3	1:B:368:VAL:HG12	1.78	0.66
1:B:352:PRO:HD2	1:B:353:GLN:OE1	1.96	0.66
1:B:524:ILE:H	1:B:524:ILE:HD13	1.61	0.66
1:B:303:ASP:HA	4:B:804:HOH:O	1.97	0.65
1:A:30:ILE:HG12	4:A:808:HOH:O	1.96	0.64
1:B:627:MET:CE	1:B:629:MET:HB2	2.28	0.64
1:B:51:TYR:HA	1:B:54:LYS:HD2	1.79	0.63
1:B:88:VAL:O	1:B:89:SER:C	2.42	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:554:LYS:C	3:B:704:CL:CL	2.78	0.63
1:A:445:ARG:HA	1:A:464:ASN:HD22	1.63	0.63
1:A:425:TYR:OH	1:A:473:GLU:HG3	1.99	0.63
1:B:627:MET:HE1	1:B:629:MET:HB2	1.81	0.62
1:A:255:TYR:HB3	1:A:368:VAL:HG12	1.80	0.62
1:B:154:ILE:HB	1:B:163:ALA:HB3	1.81	0.62
1:A:30:ILE:CG1	4:A:808:HOH:O	2.48	0.61
1:B:555:ASN:N	3:B:704:CL:CL	2.71	0.61
1:A:154:ILE:HB	1:A:163:ALA:HB3	1.83	0.60
1:A:567:ASN:HA	1:A:570:LEU:HD23	1.84	0.60
1:B:171:ILE:HD11	1:B:220:MET:HE1	1.85	0.59
1:B:88:VAL:O	1:B:89:SER:O	2.20	0.59
1:B:553:THR:O	3:B:704:CL:CL	2.59	0.57
1:B:369:TYR:HB2	1:B:370:PRO:HD3	1.85	0.57
1:A:42:PHE:HB2	1:A:63:THR:HG22	1.86	0.57
1:A:224:LEU:C	1:A:224:LEU:HD12	2.29	0.57
1:B:99:TYR:CD2	1:B:134:PRO:HG3	2.40	0.56
1:A:99:TYR:CD2	1:A:134:PRO:HG3	2.41	0.56
1:B:427:ILE:HD12	1:B:453:ILE:HG13	1.88	0.55
1:A:369:TYR:HB2	1:A:370:PRO:HD3	1.88	0.54
1:A:422:LYS:O	1:A:423:THR:C	2.51	0.54
1:B:422:LYS:O	1:B:423:THR:C	2.51	0.54
1:A:516:ASP:HA	1:A:519:TYR:CE1	2.43	0.54
1:A:89:SER:OG	1:A:90:LYS:N	2.41	0.53
1:B:35:ASP:CB	4:B:801:HOH:O	2.28	0.53
1:A:427:ILE:HD12	1:A:453:ILE:HG13	1.90	0.53
1:B:580:ASN:ND2	4:B:803:HOH:O	2.42	0.53
1:B:548:HIS:CD2	3:B:704:CL:CL	2.93	0.52
1:A:428:ASP:HA	1:A:450:ASN:OD1	2.10	0.52
1:B:428:ASP:HA	1:B:450:ASN:OD1	2.10	0.52
1:B:371:PHE:CE1	1:B:375:MET:HE1	2.45	0.51
1:B:516:ASP:HA	1:B:519:TYR:CE1	2.45	0.51
1:B:555:ASN:HA	3:B:704:CL:CL	2.48	0.51
1:B:481:LYS:O	1:B:481:LYS:HD3	2.10	0.51
1:A:603:LEU:HD23	1:A:612:ARG:HB2	1.93	0.51
1:A:397:ILE:HA	1:A:499:TYR:HD2	1.76	0.50
1:B:207:GLN:HG3	3:B:706:CL:CL	2.49	0.50
1:A:89:SER:HB3	1:A:92:LYS:HB2	1.94	0.50
1:A:102:LYS:HG3	4:A:827:HOH:O	2.12	0.50
1:B:469:ARG:HG2	4:B:809:HOH:O	2.12	0.49
1:A:105:TYR:O	1:A:314:ILE:HG21	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:LEU:HD22	1:A:395:PHE:CD1	2.47	0.49
1:B:397:ILE:HA	1:B:499:TYR:HD2	1.78	0.49
1:B:587:ILE:CD1	1:B:617:PHE:CE1	2.96	0.48
1:A:30:ILE:CA	4:A:808:HOH:O	2.20	0.48
1:A:75:VAL:HG23	1:A:75:VAL:O	2.13	0.48
1:B:90:LYS:HG2	4:B:833:HOH:O	2.12	0.48
1:A:600:THR:O	1:A:601:ALA:HB2	2.14	0.48
1:A:286:LEU:HD21	1:A:497:PRO:HD3	1.96	0.48
1:B:286:LEU:HD21	1:B:497:PRO:HD3	1.96	0.48
1:B:369:TYR:HD1	1:B:372:MET:HE2	1.79	0.48
1:B:317:LYS:HA	1:B:317:LYS:HE3	1.96	0.47
1:B:360:LEU:HD22	1:B:395:PHE:CD1	2.49	0.47
1:A:156:ASP:OD1	1:A:158:ASN:N	2.41	0.47
1:A:344:TYR:CE1	1:A:634:LYS:HE3	2.48	0.47
1:A:370:PRO:HB2	1:A:375:MET:HE2	1.96	0.47
1:A:496:TYR:CD1	1:A:497:PRO:HD2	2.50	0.47
1:A:555:ASN:CG	3:A:705:CL:CL	2.83	0.47
1:A:340:MET:HE1	1:A:346:SER:O	2.15	0.47
1:B:600:THR:O	1:B:601:ALA:HB2	2.13	0.47
1:A:640:GLY:O	1:A:643:SER:N	2.46	0.47
1:B:496:TYR:CD1	1:B:497:PRO:HD2	2.50	0.46
1:A:351:HIS:HD2	1:A:354:THR:OG1	1.97	0.46
1:A:355:GLY:O	1:A:547:PRO:HA	2.14	0.46
1:B:617:PHE:HB2	1:B:645:ASN:HB3	1.95	0.46
1:B:355:GLY:O	1:B:547:PRO:HA	2.13	0.46
1:A:247:LYS:O	1:A:365:SER:OG	2.34	0.46
1:B:35:ASP:OD1	1:B:83:ARG:NH1	2.46	0.46
1:A:464:ASN:OD1	1:A:519:TYR:CD2	2.68	0.46
1:B:255:TYR:HB3	1:B:368:VAL:CG1	2.44	0.46
1:A:491:ASP:HA	1:A:500:ASN:OD1	2.16	0.46
1:B:351:HIS:HD2	1:B:354:THR:OG1	1.98	0.46
1:B:407:ILE:HG12	1:B:597:LYS:HD3	1.97	0.46
1:B:445:ARG:HD3	1:B:465:ILE:HD12	1.98	0.46
1:A:464:ASN:OD1	1:A:519:TYR:HD2	1.99	0.46
1:B:613:GLN:N	1:B:637:GLN:OE1	2.48	0.46
1:B:105:TYR:O	1:B:314:ILE:HG21	2.16	0.46
1:B:207:GLN:CG	3:B:706:CL:CL	3.01	0.46
1:B:256:VAL:HA	1:B:279:GLY:HA2	1.98	0.46
1:A:502:GLN:HE22	1:A:525:LEU:HD12	1.80	0.45
1:A:640:GLY:O	1:A:641:MET:C	2.59	0.45
1:A:256:VAL:HA	1:A:279:GLY:HA2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:LEU:HD22	1:A:534:ILE:HG21	1.97	0.45
1:B:383:LEU:HB3	1:B:391:LEU:HD13	1.98	0.45
1:A:421:ASP:OD1	1:A:421:ASP:N	2.50	0.45
1:A:610:THR:O	1:A:611:GLY:C	2.59	0.45
1:A:613:GLN:N	1:A:637:GLN:OE1	2.50	0.45
1:B:575:MET:HE3	1:B:596:GLY:HA2	1.98	0.45
1:A:348:THR:HB	1:A:532:LEU:HD12	1.99	0.45
1:A:443:VAL:O	1:A:465:ILE:CD1	2.65	0.45
1:A:322:LYS:HA	1:A:322:LYS:HE3	1.99	0.45
1:B:156:ASP:OD1	1:B:158:ASN:N	2.41	0.45
1:A:287:TYR:CZ	1:A:550:LEU:HD11	2.52	0.45
1:A:246:GLU:OE2	1:A:367:ASP:OD1	2.35	0.44
1:A:469:ARG:O	1:A:472:LEU:N	2.50	0.44
1:A:496:TYR:CD1	1:A:496:TYR:C	2.95	0.44
1:B:247:LYS:O	1:B:365:SER:OG	2.34	0.44
1:B:469:ARG:O	1:B:472:LEU:N	2.50	0.44
1:B:640:GLY:O	1:B:643:SER:N	2.46	0.44
1:A:543:ASN:OD1	1:A:560:LYS:HA	2.18	0.44
1:B:640:GLY:O	1:B:641:MET:C	2.60	0.44
1:B:59:GLU:O	1:B:63:THR:OG1	2.33	0.44
1:B:148:LYS:HA	1:B:295:ASP:OD1	2.18	0.44
1:A:575:MET:HE3	1:A:596:GLY:HA2	1.99	0.44
1:A:255:TYR:HB3	1:A:368:VAL:CG1	2.46	0.44
1:B:496:TYR:CD1	1:B:496:TYR:C	2.95	0.44
1:B:90:LYS:HG3	1:B:91:ASN:OD1	2.18	0.43
1:B:343:ASP:HB3	1:B:633:VAL:HG13	2.00	0.43
1:A:71:ASN:OD1	1:A:71:ASN:C	2.61	0.43
1:A:91:ASN:ND2	1:A:118:LYS:HB2	2.33	0.43
1:A:224:LEU:HD12	1:A:224:LEU:O	2.18	0.43
1:B:176:LYS:CE	1:B:209:ASP:HB3	2.49	0.43
1:A:59:GLU:O	1:A:63:THR:OG1	2.33	0.43
1:B:96:ASP:OD1	1:B:113:GLN:HG2	2.18	0.43
1:B:287:TYR:CZ	1:B:550:LEU:HD11	2.52	0.43
1:B:348:THR:HB	1:B:532:LEU:HD12	2.00	0.43
1:A:629:MET:SD	1:A:652:VAL:HG11	2.58	0.43
1:A:273:LYS:HB3	1:A:275:ASP:OD1	2.17	0.43
1:A:415:ASN:C	1:A:415:ASN:HD22	2.26	0.43
1:B:543:ASN:OD1	1:B:560:LYS:HA	2.19	0.43
1:B:262:GLU:HA	1:B:265:LYS:HD2	2.01	0.43
1:A:503:ILE:H	1:A:503:ILE:HD13	1.84	0.43
1:A:612:ARG:C	1:A:637:GLN:OE1	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:659:ASN:ND2	4:A:805:HOH:O	2.52	0.42
1:B:340:MET:HE1	1:B:346:SER:O	2.18	0.42
1:A:529:VAL:O	1:A:533:SER:OG	2.36	0.42
1:B:42:PHE:CE1	1:B:66:PRO:HB3	2.54	0.42
1:B:273:LYS:HB3	1:B:275:ASP:OD1	2.19	0.42
1:A:555:ASN:N	4:A:803:HOH:O	2.52	0.42
1:B:371:PHE:CD1	1:B:375:MET:HE1	2.55	0.42
1:A:70:TYR:CD2	1:A:75:VAL:HG21	2.54	0.42
1:A:436:LYS:H	1:A:436:LYS:HD3	1.85	0.42
1:B:93:LYS:HD2	1:B:123:TRP:CZ2	2.55	0.42
1:B:99:TYR:HD2	1:B:112:VAL:HG11	1.85	0.42
1:A:148:LYS:HA	1:A:295:ASP:OD1	2.19	0.42
1:B:421:ASP:OD1	1:B:421:ASP:N	2.52	0.42
1:B:573:ASP:O	1:B:576:GLN:HB3	2.19	0.42
1:A:645:ASN:HD22	1:A:645:ASN:H	1.68	0.42
1:B:626:ASN:ND2	1:B:664:TYR:O	2.53	0.42
1:B:35:ASP:OD2	4:B:801:HOH:O	2.22	0.42
1:B:174:VAL:O	1:B:175:PRO:C	2.62	0.42
1:B:369:TYR:HD1	1:B:372:MET:CE	2.32	0.42
1:B:629:MET:SD	1:B:652:VAL:HG11	2.60	0.42
1:A:99:TYR:HD2	1:A:112:VAL:HG11	1.83	0.42
1:A:383:LEU:HB3	1:A:391:LEU:HD13	2.03	0.41
1:A:603:LEU:HD11	1:A:614:ILE:HD11	2.01	0.41
1:A:42:PHE:CZ	1:A:67:ILE:HG12	2.55	0.41
1:A:626:ASN:ND2	1:A:664:TYR:O	2.53	0.41
1:B:351:HIS:CD2	1:B:354:THR:OG1	2.73	0.41
1:B:425:TYR:CB	1:B:469:ARG:NH2	2.83	0.41
1:B:580:ASN:CB	4:B:803:HOH:O	2.69	0.41
1:A:521:GLN:HE22	1:A:602:GLU:HB2	1.85	0.41
1:A:505:ASN:C	1:A:507:ASN:N	2.78	0.41
1:A:466:PHE:O	1:A:470:VAL:HG23	2.20	0.41
1:B:179:SER:OG	1:B:181:LYS:O	2.36	0.41
1:B:553:THR:C	3:B:704:CL:CL	3.01	0.41
1:B:619:SER:O	1:B:620:TYR:HB3	2.20	0.41
1:B:327:THR:HG23	1:B:549:LEU:C	2.46	0.41
1:A:327:THR:HG23	1:A:549:LEU:C	2.46	0.40
1:A:174:VAL:O	1:A:175:PRO:C	2.64	0.40
1:B:154:ILE:HG12	1:B:324:ILE:HG22	2.03	0.40
1:A:57:ASN:HB2	1:A:62:MET:HE2	2.03	0.40
1:A:613:GLN:CD	1:A:641:MET:HE1	2.47	0.40
1:B:360:LEU:O	1:B:395:PHE:CE2	2.75	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:VAL:O	1:A:176:LYS:N	2.55	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	642/643 (100%)	569 (89%)	67 (10%)	6 (1%)	14	41
1	B	635/643 (99%)	571 (90%)	58 (9%)	6 (1%)	14	41
All	All	1277/1286 (99%)	1140 (89%)	125 (10%)	12 (1%)	14	41

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	89	SER
1	A	605	MET
1	A	610	THR
1	A	246	GLU
1	B	246	GLU
1	A	601	ALA
1	A	620	TYR
1	B	601	ALA
1	B	617	PHE
1	B	620	TYR
1	A	611	GLY
1	B	648	ILE

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	574/574 (100%)	494 (86%)	80 (14%)	3	12
1	B	571/574 (100%)	501 (88%)	70 (12%)	4	16
All	All	1145/1148 (100%)	995 (87%)	150 (13%)	4	14

All (150) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LYS
1	A	28	LYS
1	A	40	LYS
1	A	42	PHE
1	A	43	LYS
1	A	44	GLN
1	A	64	GLU
1	A	66	PRO
1	A	69	ILE
1	A	73	LEU
1	A	86	LYS
1	A	90	LYS
1	A	100	LYS
1	A	102	LYS
1	A	111	ASN
1	A	124	LYS
1	A	125	LEU
1	A	138	LYS
1	A	156	ASP
1	A	177	ASN
1	A	181	LYS
1	A	216	THR
1	A	224	LEU
1	A	230	LYS
1	A	246	GLU
1	A	252	LEU
1	A	267	LYS

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Mol	Chain	Res	Type
1	A	289	LYS
1	A	300	THR
1	A	318	LYS
1	A	324	ILE
1	A	334	LYS
1	A	342	ASN
1	A	348	THR
1	A	365	SER
1	A	368	VAL
1	A	378	GLU
1	A	382	LYS
1	A	387	LYS
1	A	392	LEU
1	A	403	SER
1	A	409	THR
1	A	415	ASN
1	A	418	THR
1	A	419	LEU
1	A	421	ASP
1	A	434	LYS
1	A	436	LYS
1	A	447	GLU
1	A	477	LYS
1	A	484	LYS
1	A	490	GLU
1	A	496	TYR
1	A	502	GLN
1	A	503	ILE
1	A	505	ASN
1	A	508	LEU
1	A	510	ASN
1	A	513	LEU
1	A	514	LEU
1	A	527	ASN
1	A	532	LEU
1	A	533	SER
1	A	545	ASN
1	A	549	LEU
1	A	554	LYS
1	A	562	ILE
1	A	566	GLU
1	A	570	LEU

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Mol	Chain	Res	Type
1	A	576	GLN
1	A	587	ILE
1	A	595	ILE
1	A	597	LYS
1	A	603	LEU
1	A	604	LYS
1	A	609	GLU
1	A	610	THR
1	A	634	LYS
1	A	666	ILE
1	A	668	GLU
1	B	26	LYS
1	B	28	LYS
1	B	44	GLN
1	B	52	ILE
1	B	54	LYS
1	B	73	LEU
1	B	79	ASN
1	B	86	LYS
1	B	87	LYS
1	B	88	VAL
1	B	91	ASN
1	B	92	LYS
1	B	93	LYS
1	B	111	ASN
1	B	115	ASN
1	B	125	LEU
1	B	156	ASP
1	B	177	ASN
1	B	181	LYS
1	B	184	LYS
1	B	215	LYS
1	B	216	THR
1	B	252	LEU
1	B	277	VAL
1	B	289	LYS
1	B	290	LYS
1	B	300	THR
1	B	317	LYS
1	B	318	LYS
1	B	319	LYS
1	B	324	ILE

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Mol	Chain	Res	Type
1	B	334	LYS
1	B	342	ASN
1	B	348	THR
1	B	365	SER
1	B	368	VAL
1	B	382	LYS
1	B	385	GLU
1	B	387	LYS
1	B	392	LEU
1	B	403	SER
1	B	409	THR
1	B	417	LYS
1	B	418	THR
1	B	419	LEU
1	B	422	LYS
1	B	436	LYS
1	B	447	GLU
1	B	480	GLU
1	B	481	LYS
1	B	485	LYS
1	B	496	TYR
1	B	505	ASN
1	B	510	ASN
1	B	513	LEU
1	B	514	LEU
1	B	524	ILE
1	B	527	ASN
1	B	532	LEU
1	B	545	ASN
1	B	549	LEU
1	B	556	LYS
1	B	559	LYS
1	B	560	LYS
1	B	565	LYS
1	B	573	ASP
1	B	595	ILE
1	B	622	LYS
1	B	651	LYS
1	B	666	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	79	ASN
1	A	91	ASN
1	A	129	HIS
1	A	342	ASN
1	A	351	HIS
1	A	393	ASN
1	A	442	ASN
1	A	464	ASN
1	A	502	GLN
1	A	505	ASN
1	A	507	ASN
1	A	521	GLN
1	A	567	ASN
1	A	576	GLN
1	A	577	GLN
1	A	613	GLN
1	A	645	ASN
1	B	32	ASN
1	B	79	ASN
1	B	342	ASN
1	B	351	HIS
1	B	393	ASN
1	B	502	GLN
1	B	521	GLN
1	B	567	ASN
1	B	645	ASN

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

Of 14 ligands modelled in this entry, 14 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	643/643 (100%)	-0.67	1 (0%) 91 88	39, 98, 162, 237	1 (0%)
1	B	638/643 (99%)	-0.70	0 100 100	42, 102, 156, 210	1 (0%)
All	All	1281/1286 (99%)	-0.68	1 (0%) 92 90	39, 100, 160, 237	2 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	246	GLU	2.4

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(Å ²)	Q<0.9
3	CL	A	705	1/1	0.81	0.10	247,247,247,247	0
3	CL	B	704	1/1	0.84	0.10	294,294,294,294	0
3	CL	B	705	1/1	0.87	0.05	96,96,96,96	0
2	CD	A	704	1/1	0.91	0.11	198,198,198,198	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CL	A	706	1/1	0.98	0.14	37,37,37,37	0
2	CD	B	701	1/1	0.98	0.06	154,154,154,154	0
2	CD	A	701	1/1	0.98	0.08	178,178,178,178	0
3	CL	B	708	1/1	0.98	0.05	39,39,39,39	0
3	CL	B	707	1/1	0.99	0.08	54,54,54,54	0
3	CL	B	706	1/1	0.99	0.03	61,61,61,61	0
2	CD	B	703	1/1	1.00	0.02	65,65,65,65	0
2	CD	A	702	1/1	1.00	0.04	81,81,81,81	0
2	CD	A	703	1/1	1.00	0.01	63,63,63,63	0
2	CD	B	702	1/1	1.00	0.02	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.