



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 21, 2025 – 01:16 PM EDT

PDB ID : 5RLZ / pdb_00005rlz
Title : PanDDA analysis group deposition – Crystal Structure of SARS-CoV-2 heli-
case in complex with Z2293643386
Authors : Newman, J.A.; Yosaatmadja, Y.; Douangamath, A.; Aimon, A.; Powell, A.J.;
Dias, A.; Fearon, D.; Dunnett, L.; Brandao-Neto, J.; Krojer, T.; Skyner, R.;
Gorrie-Stone, T.; Thompson, W.; von Delft, F.; Arrowsmith, C.H.; Edwards,
A.; Bountra, C.; Gileadi, O.
Deposited on : 2020-09-16
Resolution : 1.97 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)

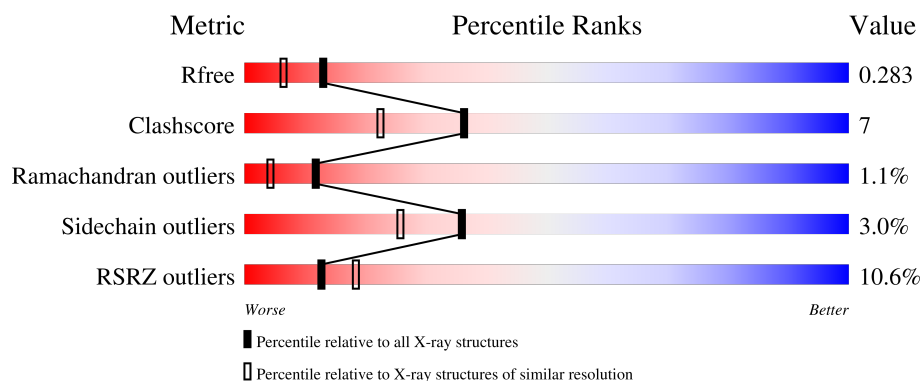
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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}	164625	1356 (1.98-1.98)
Clashscore	180529	1437 (1.98-1.98)
Ramachandran outliers	177936	1426 (1.98-1.98)
Sidechain outliers	177891	1426 (1.98-1.98)
RSRZ outliers	164620	1356 (1.98-1.98)

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	601	<div> <div>14%</div> <div>78%</div> <div>16%</div> <div>• 5%</div> </div>
1	B	601	<div> <div>7%</div> <div>79%</div> <div>17%</div> <div>• •</div> </div>

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.42

2 Entry composition [i](#)

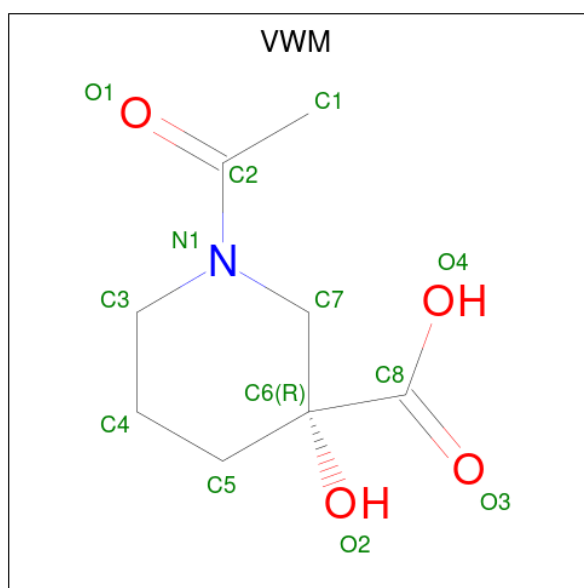
There are 5 unique types of molecules in this entry. The entry contains 9418 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	585	Total	C	N	O	S	0	1	0
			4508	2875	750	848	35			
1	A	572	Total	C	N	O	S	0	0	0
			4417	2816	737	832	32			

- Molecule 2 is (3R)-1-acetyl-3-hydroxypiperidine-3-carboxylic acid (CCD ID: VWM) (formula: $C_8H_{13}NO_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			13	8	1	4		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

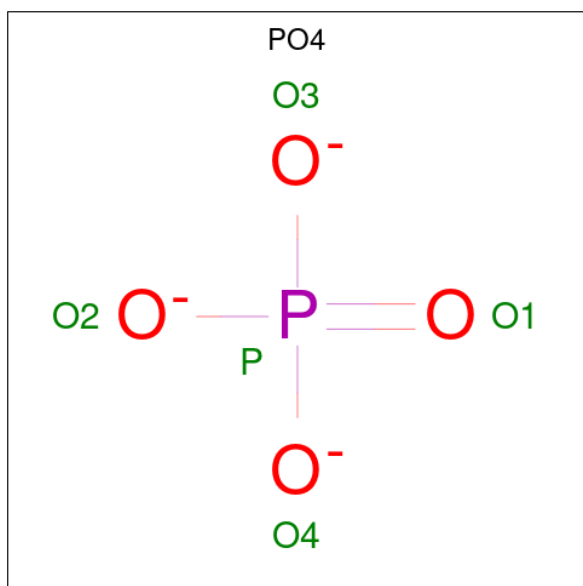
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Zn	0	0
			3	3		

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

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	250	Total	O	0	0
			250	250		
5	A	204	Total	O	0	0
			204	204		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.21Å 70.39Å 85.35Å 102.74° 95.87° 112.42°	Depositor
Resolution (Å)	81.46 – 1.97 81.45 – 1.97	Depositor EDS
% Data completeness (in resolution range)	91.0 (81.46-1.97) 91.0 (81.45-1.97)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.222 , 0.273 0.235 , 0.283	Depositor DCC
R_{free} test set	4372 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9418	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.81% 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: VWM, ZN, 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	A	0.70	0/4517	0.79	0/6156
1	B	0.70	0/4610	0.82	0/6283
All	All	0.70	0/9127	0.81	0/12439

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4417	0	4322	57	0
1	B	4508	0	4425	69	0
2	B	13	0	0	0	0
3	A	3	0	0	0	0
3	B	3	0	0	1	0
4	A	10	0	0	0	0
4	B	10	0	0	1	0
5	A	204	0	0	7	0
5	B	250	0	0	9	0
All	All	9418	0	8747	126	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 7.

All (126) 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:5:CYS:HG	3:B:703:ZN:ZN	0.66	0.98
1:B:27:CYS:SG	5:B:855:HOH:O	2.33	0.84
1:B:510:VAL:HG21	1:B:541:TYR:CD1	2.16	0.80
1:B:12:THR:HG21	1:B:25:LEU:O	1.81	0.79
1:B:48:TYR:O	5:B:801:HOH:O	2.01	0.77
1:A:60:VAL:HB	5:A:928:HOH:O	1.89	0.71
1:B:48:TYR:OH	1:B:90:PHE:O	2.05	0.71
1:A:368:ALA:O	1:A:393:ALA:HA	1.92	0.69
1:A:519:ASN:HB3	1:A:530:THR:HG23	1.77	0.66
1:B:8:CYS:SG	1:B:99:GLY:N	2.67	0.66
1:A:13:SER:OG	1:A:44:SER:HB2	1.94	0.66
1:A:510:VAL:HG21	1:A:541:TYR:CD1	2.30	0.65
1:A:275:GLN:O	1:A:395:HIS:ND1	2.30	0.64
1:B:51:ASN:OD1	5:B:802:HOH:O	2.15	0.64
1:B:271:LYS:NZ	5:B:803:HOH:O	2.16	0.63
1:A:367:THR:HA	1:A:392:ARG:O	1.99	0.62
1:A:519:ASN:HB3	1:A:530:THR:CG2	2.31	0.61
1:B:7:LEU:HD12	1:B:103:VAL:HG22	1.81	0.61
1:A:21:ARG:NH1	5:A:809:HOH:O	2.34	0.60
1:B:275:GLN:NE2	1:B:435:ASP:OD2	2.30	0.59
1:B:508:LYS:HD3	5:B:992:HOH:O	2.01	0.59
1:B:445:PRO:HB3	1:B:468:SER:HB3	1.85	0.59
1:B:554:HIS:ND1	5:B:810:HOH:O	2.32	0.57
1:B:376:ILE:HG22	1:B:400:GLY:HA3	1.85	0.57
1:A:5:CYS:SG	1:A:26:CYS:HB3	2.45	0.57
1:A:269:TYR:OH	1:A:294:GLY:HA3	2.05	0.57
1:A:163:LEU:HD23	1:A:211:TYR:CD2	2.40	0.56
1:A:4:ALA:O	1:A:24:PHE:HB2	2.06	0.56
1:B:13:SER:O	1:B:44:SER:HA	2.05	0.56
1:B:510:VAL:HG21	1:B:541:TYR:CG	2.41	0.55
1:B:28:LYS:O	1:B:32:ASP:OD1	2.24	0.55
1:B:12:THR:CG2	1:B:26:CYS:HA	2.37	0.55
1:A:7:LEU:HD13	1:A:103:VAL:HG22	1.88	0.54
1:B:277:TYR:HA	1:B:396:TYR:O	2.08	0.54
1:A:152:ALA:HB2	1:A:167:TRP:CZ3	2.42	0.54
1:B:279:THR:HB	1:B:429:MET:CE	2.39	0.53
1:B:220:ASN:N	1:B:220:ASN:OD1	2.41	0.53
1:B:385:SER:OG	5:B:804:HOH:O	2.18	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:THR:HG21	1:B:26:CYS:HA	1.92	0.52
1:A:16:CYS:SG	1:A:18:ALA:HB3	2.50	0.52
1:A:7:LEU:HD21	1:A:106:PHE:HB2	1.92	0.52
1:B:7:LEU:CD1	1:B:103:VAL:HG22	2.41	0.51
1:A:31:TYR:CE2	1:A:87:GLY:HA2	2.46	0.51
1:B:252:LEU:HB3	1:B:299:TYR:CD1	2.46	0.50
1:A:277:TYR:HA	1:A:396:TYR:O	2.10	0.50
1:A:350:SER:O	1:A:352:LEU:N	2.44	0.50
1:A:304:ILE:HA	1:A:370:ILE:O	2.11	0.50
1:B:220:ASN:O	1:B:223:ASP:OD2	2.30	0.50
1:B:7:LEU:HD21	1:B:106:PHE:HB2	1.95	0.49
1:A:331:SER:HB2	1:A:353:GLU:HG3	1.94	0.49
1:B:32:ASP:O	1:B:36:SER:OG	2.24	0.49
1:B:333:ILE:HB	1:B:358:CYS:SG	2.52	0.49
1:B:477:LYS:NZ	1:B:551:GLU:OE2	2.37	0.48
1:A:296:ALA:O	1:A:300:PRO:HA	2.13	0.48
1:B:474[B]:MET:HG2	1:B:590:LEU:HB2	1.95	0.48
1:B:103:VAL:CG1	1:B:103:VAL:O	2.61	0.48
1:B:533:VAL:HG11	1:B:560:ARG:O	2.13	0.48
1:A:59:ASP:OD1	1:A:61:THR:OG1	2.23	0.48
1:B:279:THR:HB	1:B:429:MET:HE2	1.96	0.48
1:A:1:ALA:N	5:A:819:HOH:O	2.47	0.48
1:B:2:VAL:N	5:B:828:HOH:O	2.47	0.47
1:A:235:LEU:HD21	1:A:382:TYR:CE2	2.49	0.47
1:B:14:LEU:HB2	1:B:25:LEU:O	2.15	0.47
1:B:115:THR:HA	1:B:411:LEU:O	2.15	0.47
1:B:16:CYS:O	1:B:22:ARG:HA	2.15	0.47
1:B:8:CYS:SG	1:B:99:GLY:O	2.71	0.47
1:A:12:THR:HG21	1:A:25:LEU:O	2.16	0.46
1:A:512:ILE:O	1:A:546:PHE:HA	2.15	0.46
1:B:15:ARG:HG3	1:B:24:PHE:CD2	2.50	0.46
1:B:378:MET:O	1:B:407:ALA:HB2	2.15	0.46
1:A:4:ALA:O	1:A:24:PHE:CB	2.63	0.46
1:A:318:CYS:HB3	1:A:343:PHE:CD2	2.51	0.45
1:A:351:THR:HG23	1:A:364:PRO:HG3	1.98	0.45
1:A:376:ILE:HG12	1:A:425:VAL:HG11	1.97	0.45
1:A:64:TYR:O	1:A:70:TYR:HA	2.16	0.45
1:A:585:LEU:HD22	1:A:587:PHE:CE2	2.52	0.45
1:A:34:VAL:HA	1:A:39:HIS:O	2.17	0.45
1:A:512:ILE:HA	1:A:531:GLN:O	2.17	0.45
1:A:308:ALA:O	1:A:359:THR:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:CYS:SG	1:B:99:GLY:CA	3.05	0.45
1:A:480:ILE:N	1:A:480:ILE:HD12	2.33	0.44
1:A:334:ILE:HD12	1:A:348:VAL:HG13	1.99	0.44
1:A:304:ILE:HG12	1:A:370:ILE:HB	1.98	0.44
1:B:184:GLY:C	1:B:195:ILE:HG22	2.38	0.44
1:B:245:HIS:HE1	5:B:930:HOH:O	2.00	0.44
1:B:284:PRO:HG2	1:B:566:THR:HG21	2.00	0.44
1:B:511:PHE:O	1:B:530:THR:HA	2.17	0.44
1:B:124:ASN:OD1	1:B:381:ASN:ND2	2.43	0.44
1:A:33:HIS:HA	1:A:107:ASN:OD1	2.18	0.44
1:B:14:LEU:O	1:B:24:PHE:HA	2.18	0.43
1:B:315:ASP:O	1:B:318:CYS:HB2	2.18	0.43
1:B:152:ALA:HB2	1:B:167:TRP:CZ3	2.52	0.43
1:A:591:GLU:O	1:A:592:ILE:HB	2.18	0.43
1:B:488:ILE:HD11	1:B:517:SER:HB3	2.00	0.43
1:A:490:ARG:HB2	1:A:491:PRO:HD3	2.00	0.43
1:B:254:PRO:HB3	1:B:298:TYR:CE2	2.54	0.43
1:B:282:GLY:O	1:B:288:LYS:HD3	2.19	0.43
1:B:59:ASP:OD1	1:B:59:ASP:C	2.57	0.43
1:B:372:VAL:CG1	1:B:399:ILE:HD12	2.49	0.43
1:B:152:ALA:HB2	1:B:167:TRP:CH2	2.55	0.42
1:B:386:VAL:O	1:B:390:ARG:HG2	2.18	0.42
1:B:377:SER:O	1:B:406:PRO:HA	2.19	0.42
1:A:72:CYS:SG	1:A:74:SER:HB2	2.60	0.42
1:B:5:CYS:O	1:B:9:ASN:N	2.45	0.42
1:B:452:VAL:CG1	1:B:565:ILE:HD11	2.50	0.42
1:B:544:VAL:O	1:B:572:ILE:HA	2.19	0.42
1:A:561:PHE:CD2	1:A:581:LEU:HD22	2.55	0.42
1:A:552:THR:C	5:A:892:HOH:O	2.58	0.42
1:B:31:TYR:HB2	1:B:89:VAL:HG23	2.02	0.41
1:A:280:LEU:HD11	1:A:438:LEU:HG	2.01	0.41
1:A:241:VAL:HG22	5:A:867:HOH:O	2.19	0.41
1:B:26:CYS:O	1:B:29:CYS:N	2.54	0.41
1:B:542:ASP:HA	1:B:569:LYS:HB2	2.02	0.41
1:A:152:ALA:HB2	1:A:167:TRP:CE3	2.56	0.41
1:A:584:LYS:NZ	5:A:818:HOH:O	2.44	0.41
1:A:585:LEU:HD22	1:A:587:PHE:CZ	2.55	0.41
1:B:404:GLN:NE2	4:B:706:PO4:O2	2.42	0.41
1:A:31:TYR:CD2	1:A:87:GLY:HA2	2.55	0.41
1:B:376:ILE:HD12	1:B:376:ILE:HA	1.96	0.41
1:B:504:PRO:HB3	1:B:507:ARG:NH2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:LYS:O	1:A:395:HIS:HA	2.21	0.41
1:B:84:CYS:SG	1:B:84:CYS:O	2.79	0.41
1:A:44:SER:O	1:A:45:VAL:C	2.59	0.41
1:A:377:SER:O	1:A:406:PRO:HA	2.21	0.41
1:A:151:ILE:HG12	1:A:226:VAL:HG22	2.03	0.41
1:A:359:THR:OG1	5:A:801:HOH:O	2.22	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	564/601 (94%)	529 (94%)	27 (5%)	8 (1%)	9	2
1	B	580/601 (96%)	539 (93%)	36 (6%)	5 (1%)	14	6
All	All	1144/1202 (95%)	1068 (93%)	63 (6%)	13 (1%)	12	4

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	351	THR
1	B	48	TYR
1	B	97	CYS
1	A	195	ILE
1	A	218	LYS
1	B	10	SER
1	A	283	PRO
1	A	484	VAL
1	B	9	ASN
1	A	249	ILE
1	A	45	VAL
1	B	195	ILE

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Mol	Chain	Res	Type
1	A	6	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	485/523 (93%)	472 (97%)	13 (3%)	40	30
1	B	498/523 (95%)	482 (97%)	16 (3%)	34	24
All	All	983/1046 (94%)	954 (97%)	29 (3%)	36	26

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	26	CYS
1	B	69	SER
1	B	96	THR
1	B	103	VAL
1	B	155	ARG
1	B	188	THR
1	B	191	SER
1	B	192	LYS
1	B	215	THR
1	B	220	ASN
1	B	229	SER
1	B	231	THR
1	B	259	SER
1	B	278	SER
1	B	373	PHE
1	B	458	ASP
1	A	12	THR
1	A	26	CYS
1	A	69	SER
1	A	76	LYS
1	A	156	GLU
1	A	183	THR

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Mol	Chain	Res	Type
1	A	255	THR
1	A	259	SER
1	A	373	PHE
1	A	502	ARG
1	A	530	THR
1	A	589	SER
1	A	591	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	51	ASN
1	B	268	ASN
1	B	470	GLN
1	A	179	ASN

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
4	PO4	A	705	-	4,4,4	1.12	0	6,6,6	0.71	0
2	VWM	B	701	-	13,13,13	0.87	1 (7%)	15,19,19	1.11	1 (6%)
4	PO4	B	705	-	4,4,4	0.77	0	6,6,6	0.45	0
4	PO4	A	704	-	4,4,4	0.82	0	6,6,6	0.60	0
4	PO4	B	706	-	4,4,4	1.29	0	6,6,6	0.59	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	VWM	B	701	-	-	2/10/22/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	VWM	O4-C8	-2.57	1.21	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	VWM	O4-C8-C6	2.27	117.50	113.14

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	701	VWM	C7-C6-C8-O3
2	B	701	VWM	C7-C6-C8-O4

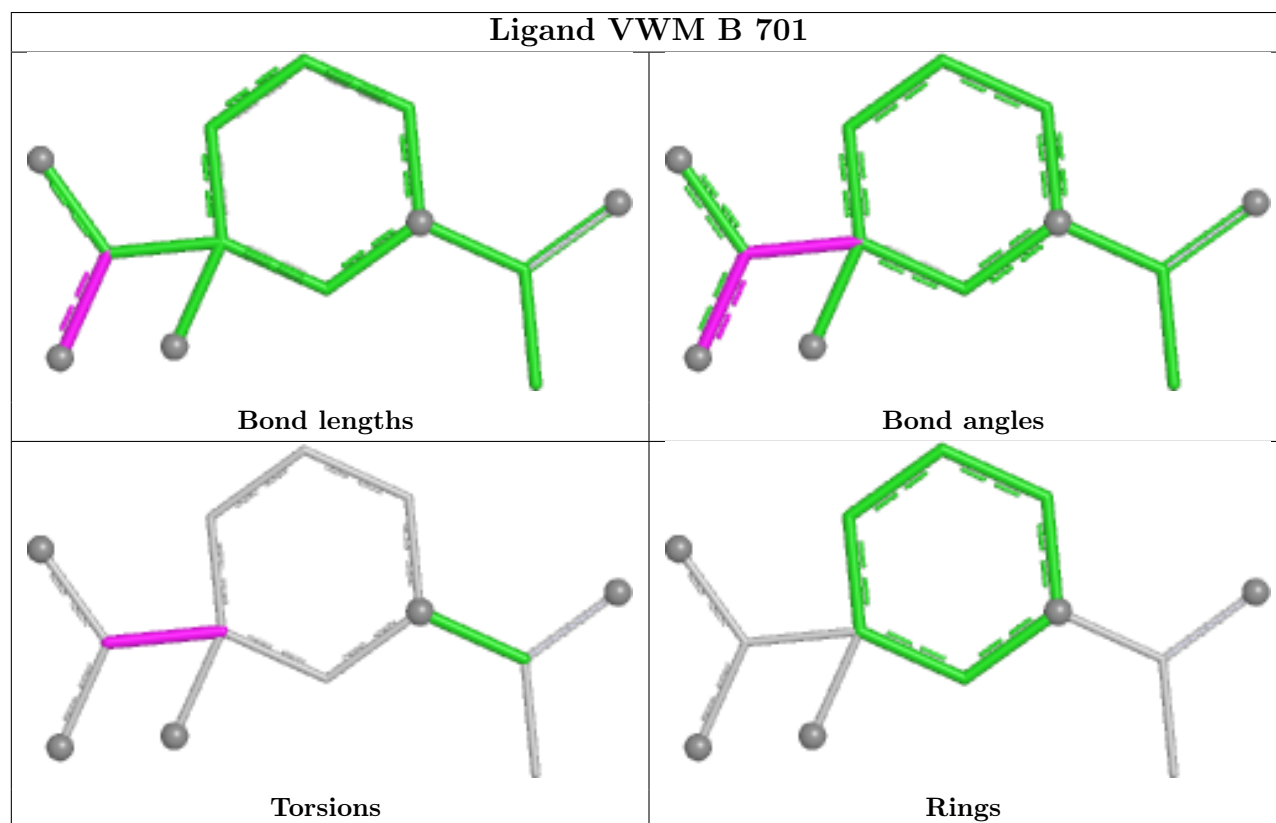
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	706	PO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	572/601 (95%)	0.89	83 (14%)	7 11	27, 57, 123, 166	0
1	B	585/601 (97%)	0.49	40 (6%)	25 34	8, 46, 96, 132	1 (0%)
All	All	1157/1202 (96%)	0.69	123 (10%)	13 18	8, 51, 112, 166	1 (0%)

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	193	VAL	6.1
1	B	474[A]	MET	5.9
1	A	592	ILE	5.5
1	A	151	ILE	5.2
1	A	219	LEU	5.1
1	A	169	VAL	4.5
1	A	81	PHE	4.5
1	A	6	VAL	4.4
1	B	214	THR	4.3
1	B	45	VAL	4.2
1	A	226	VAL	4.0
1	A	176	LEU	3.8
1	A	185	TYR	3.8
1	B	7	LEU	3.8
1	B	98	VAL	3.8
1	B	203	GLY	3.7
1	A	157	VAL	3.7
1	A	152	ALA	3.7
1	B	336	ALA	3.6
1	A	167	TRP	3.6
1	B	187	VAL	3.6
1	A	195	ILE	3.6
1	A	221	VAL	3.6
1	A	247	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	149	TYR	3.6
1	A	7	LEU	3.5
1	A	485	SER	3.5
1	A	94	LYS	3.5
1	A	158	LEU	3.5
1	A	225	PHE	3.5
1	B	2	VAL	3.5
1	A	224	TYR	3.4
1	A	181	VAL	3.4
1	A	227	LEU	3.4
1	A	170	GLY	3.3
1	B	217	TYR	3.3
1	A	154	VAL	3.3
1	A	210	VAL	3.3
1	B	14	LEU	3.3
1	B	227	LEU	3.3
1	B	12	THR	3.3
1	A	90	PHE	3.2
1	B	92	LEU	3.2
1	A	228	THR	3.2
1	A	484	VAL	3.2
1	A	256	LEU	3.2
1	A	217	TYR	3.2
1	A	200	PHE	3.2
1	A	336	ALA	3.2
1	A	92	LEU	3.1
1	B	592	ILE	3.1
1	A	182	PHE	3.1
1	A	153	THR	3.1
1	A	202	LYS	3.0
1	B	208	ALA	2.9
1	A	222	GLY	2.9
1	A	43	LEU	2.9
1	B	340	VAL	2.8
1	A	45	VAL	2.8
1	B	91	GLY	2.8
1	A	352	LEU	2.8
1	A	180	TYR	2.7
1	A	49	VAL	2.7
1	A	215	THR	2.7
1	A	208	ALA	2.6
1	B	97	CYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	413	THR	2.6
1	B	198	TYR	2.6
1	A	199	THR	2.6
1	A	209	VAL	2.6
1	A	253	TYR	2.5
1	B	44	SER	2.5
1	B	163	LEU	2.5
1	B	176	LEU	2.5
1	B	215	THR	2.5
1	B	195	ILE	2.5
1	A	165	LEU	2.5
1	B	81	PHE	2.5
1	A	103	VAL	2.5
1	B	6	VAL	2.4
1	B	213	GLY	2.4
1	A	255	THR	2.4
1	A	229	SER	2.4
1	B	593	PRO	2.4
1	A	334	ILE	2.3
1	B	4	ALA	2.3
1	B	78	PRO	2.3
1	B	102	ASN	2.3
1	A	12	THR	2.3
1	A	179	ASN	2.3
1	B	96	THR	2.3
1	A	396	TYR	2.3
1	B	219	LEU	2.3
1	A	130	LEU	2.3
1	A	147	LEU	2.3
1	A	79	ILE	2.3
1	A	46	ASN	2.3
1	A	328	ASP	2.3
1	A	241	VAL	2.2
1	B	27	CYS	2.2
1	A	302	ALA	2.2
1	A	145	PHE	2.2
1	A	14	LEU	2.2
1	A	198	TYR	2.2
1	A	335	PRO	2.2
1	A	155	ARG	2.2
1	A	13	SER	2.2
1	A	348	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	26	CYS	2.2
1	A	258	ILE	2.2
1	A	44	SER	2.1
1	A	42	VAL	2.1
1	A	24	PHE	2.1
1	A	333	ILE	2.1
1	B	43	LEU	2.1
1	A	31	TYR	2.1
1	A	70	TYR	2.1
1	B	256	LEU	2.1
1	A	166	SER	2.1
1	A	507	ARG	2.1
1	A	184	GLY	2.1
1	A	223	ASP	2.0
1	A	25	LEU	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 monosaccharides 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.

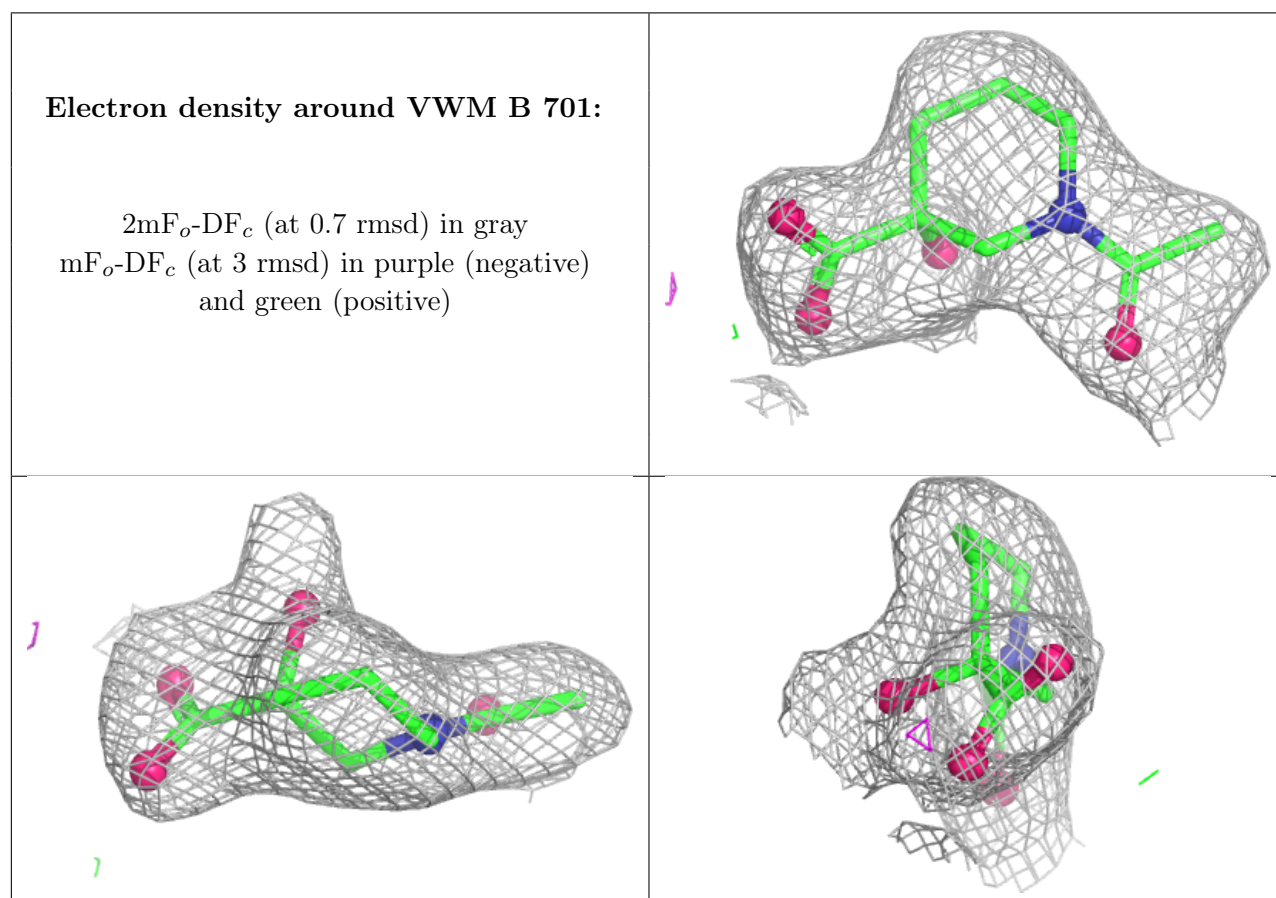
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	VWM	B	701	13/13	0.90	0.10	39,49,52,53	0
4	PO4	B	705	5/5	0.96	0.07	41,42,50,55	0
3	ZN	A	703	1/1	0.97	0.05	89,89,89,89	0
3	ZN	B	703	1/1	0.97	0.05	72,72,72,72	0
4	PO4	B	706	5/5	0.97	0.06	38,40,42,49	0
4	PO4	A	705	5/5	0.97	0.06	37,41,42,46	0
4	PO4	A	704	5/5	0.98	0.04	34,37,38,43	0
3	ZN	A	702	1/1	0.98	0.04	56,56,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ZN	B	704	1/1	0.99	0.02	74,74,74,74	0
3	ZN	A	701	1/1	0.99	0.03	61,61,61,61	0
3	ZN	B	702	1/1	1.00	0.02	43,43,43,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.