



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2022 – 02:11 PM JST

PDB ID : 7DNB
Title : Crystal structure of PhoCl barrel
Authors : Wen, Y.; Lemieux, J.M.
Deposited on : 2020-12-09
Resolution : 2.81 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.26
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.26

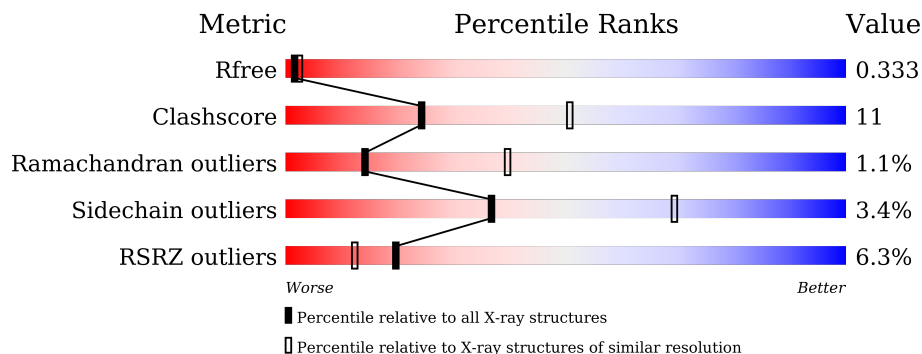
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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}	130704	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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	242	<div> <div>2%</div> <div> <div>51%</div> <div>16%</div> <div>•</div> <div>32%</div> </div> </div>
1	B	242	<div> <div>6%</div> <div> <div>46%</div> <div>19%</div> <div>•</div> <div>33%</div> </div> </div>
1	C	242	<div> <div>4%</div> <div> <div>66%</div> <div>12%</div> <div>•</div> <div>21%</div> </div> </div>
1	D	242	<div> <div>5%</div> <div> <div>48%</div> <div>17%</div> <div>35%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4999 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 PhoCl Barrel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	190	Total	C	N	O	S	0	1	0
			1429	918	237	265	9			
1	A	165	Total	C	N	O	S	1	1	0
			1247	804	203	232	8			
1	B	162	Total	C	N	O	S	0	1	0
			1186	754	200	225	7			
1	D	158	Total	C	N	O	S	0	1	0
			1136	723	191	215	7			

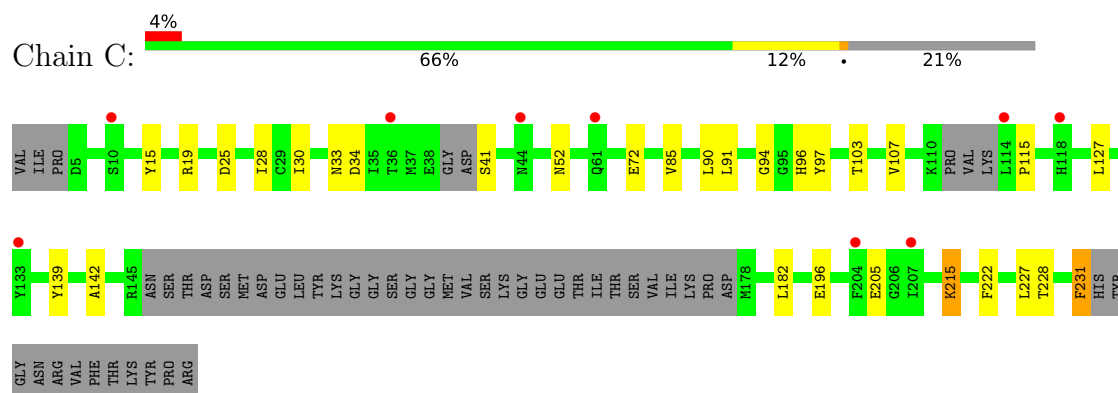
- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		

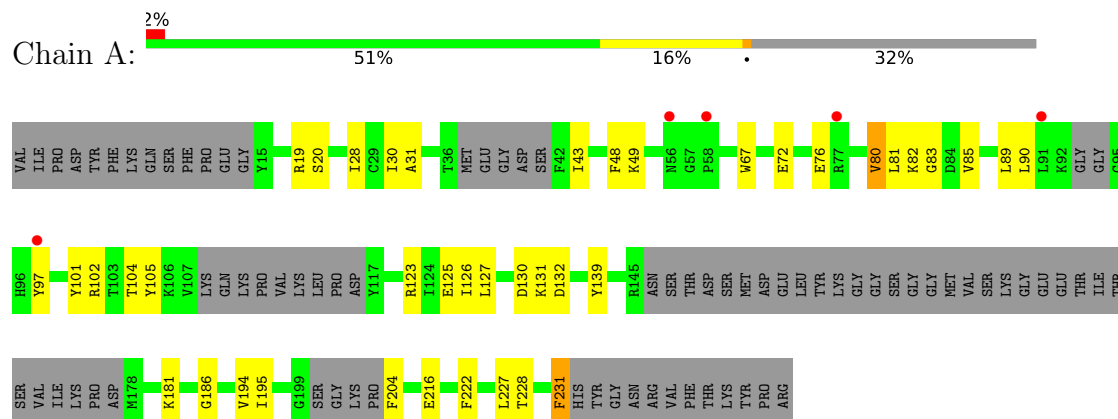
3 Residue-property plots [i](#)

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.

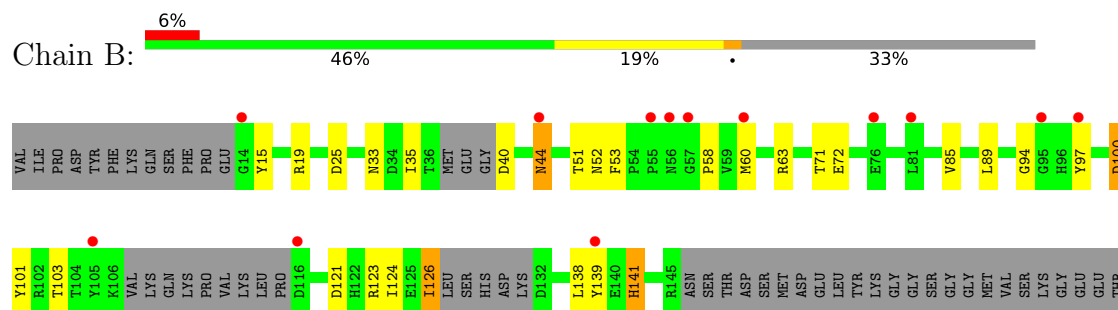
• Molecule 1: PhoCl Barrel

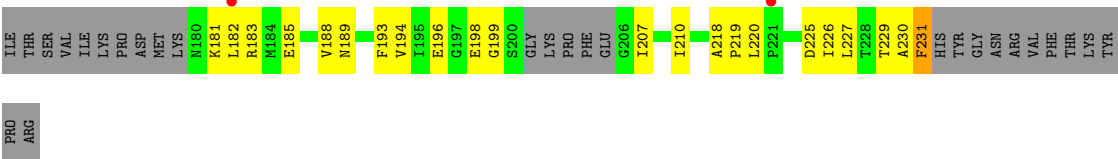


• Molecule 1: PhoCl Barrel

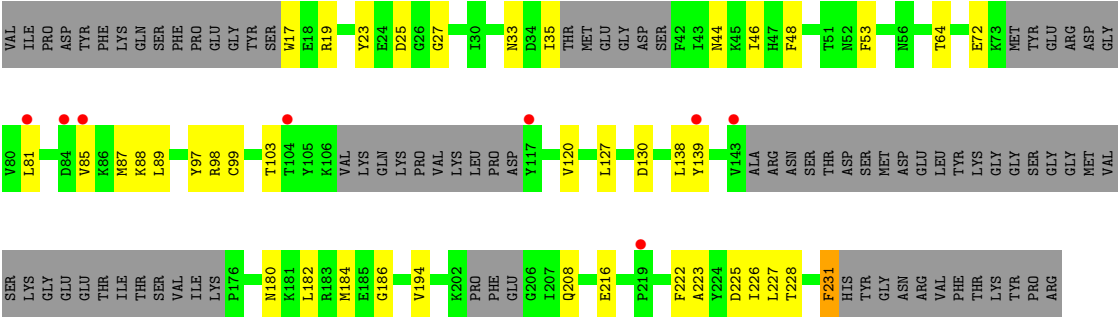


• Molecule 1: PhoCl Barrel





● Molecule 1: PhoCl Barrel



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.40Å 119.64Å 65.26Å 90.00° 107.66° 90.00°	Depositor
Resolution (Å)	43.11 – 2.81 43.11 – 2.81	Depositor EDS
% Data completeness (in resolution range)	71.3 (43.11-2.81) 71.4 (43.11-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.97 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.275 , 0.335 0.275 , 0.333	Depositor DCC
R_{free} test set	1177 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	44.7	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 61.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	4999	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.06% 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: NA, NFA

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	0.26	0/1261	0.50	0/1710
1	B	0.32	0/1195	0.50	0/1619
1	C	0.25	0/1451	0.44	0/1969
1	D	0.25	0/1146	0.44	0/1558
All	All	0.27	0/5053	0.47	0/6856

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	1247	0	1127	27	0
1	B	1186	0	1051	39	0
1	C	1429	0	1282	17	0
1	D	1136	0	990	26	0
2	B	1	0	0	0	0
All	All	4999	0	4450	104	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 11.

All (104) 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:230:ALA:H	1:B:231:NFA:HB3	1.50	0.76
1:C:196:GLU:HG3	1:C:215:LYS:HG3	1.71	0.72
1:D:19:ARG:HB2	1:D:103:THR:HG23	1.72	0.70
1:B:126:ILE:HA	1:B:138:LEU:HD13	1.72	0.69
1:D:17:TRP:N	1:D:33:ASN:O	2.27	0.68
1:B:139:TYR:HE2	1:B:207:ILE:HG23	1.59	0.67
1:A:82:LYS:NZ	1:A:104:THR:OG1	2.19	0.66
1:C:94:GLY:HA3	1:B:94:GLY:HA3	1.80	0.63
1:B:181:LYS:HA	1:B:198:GLU:HA	1.81	0.62
1:B:33:ASN:HD21	1:B:44:ASN:HB3	1.66	0.61
1:B:53:PHE:HB2	1:B:189:ASN:HD21	1.65	0.61
1:D:88:LYS:HA	1:D:98:ARG:HA	1.82	0.61
1:A:89:LEU:HD23	1:A:97:TYR:HD2	1.66	0.60
1:B:230:ALA:N	1:B:231:NFA:HB3	2.17	0.60
1:C:15:TYR:HB2	1:C:107:VAL:HA	1.84	0.60
1:D:228:THR:HA	1:D:231:NFA:HD1	1.83	0.60
1:D:72:GLU:N	1:D:120:VAL:O	2.33	0.59
1:B:89:LEU:HB3	1:B:97:TYR:HB3	1.84	0.59
1:A:195:ILE:HD11	1:A:231:NFA:HZ	1.83	0.59
1:C:72:GLU:HA	1:C:85:VAL:HB	1.86	0.58
1:B:124:ILE:HG22	1:B:124:ILE:O	2.05	0.57
1:B:19:ARG:NH1	1:B:229:THR:O	2.39	0.56
1:D:226:ILE:HG13	1:D:227:LEU:HG	1.86	0.56
1:D:194:VAL:HB	1:D:216:GLU:HG2	1.88	0.56
1:B:72:GLU:HA	1:B:85:VAL:HB	1.88	0.55
1:D:25:ASP:OD1	1:D:97:TYR:OH	2.21	0.55
1:A:123:ARG:NH1	1:A:125:GLU:OE1	2.39	0.55
1:A:130:ASP:O	1:A:132:ASP:N	2.39	0.55
1:B:58:PRO:HA	1:B:63:ARG:HE	1.71	0.55
1:A:43:ILE:HG12	1:A:181:LYS:HB2	1.88	0.55
1:A:228:THR:HA	1:A:231:NFA:HD1	1.89	0.55
1:B:126:ILE:CA	1:B:138:LEU:HD13	2.36	0.55
1:B:25:ASP:OD1	1:B:97:TYR:OH	2.21	0.54
1:A:194:VAL:HB	1:A:216:GLU:HB2	1.89	0.54
1:D:127:LEU:HD11	1:D:139:TYR:HB2	1.89	0.54
1:B:123:ARG:CB	1:B:141:HIS:HB3	2.38	0.54
1:A:72:GLU:HA	1:A:85:VAL:HB	1.90	0.53
1:B:53:PHE:HB2	1:B:189:ASN:ND2	2.23	0.53
1:C:222:PHE:HE2	1:C:227:LEU:HD11	1.73	0.53
1:D:44:ASN:ND2	1:D:180:ASN:OD1	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:PHE:HE2	1:A:227:LEU:HD11	1.74	0.53
1:B:138:LEU:HD12	1:B:139:TYR:H	1.74	0.53
1:B:19:ARG:HB2	1:B:103:THR:HG23	1.92	0.52
1:D:64:THR:HB	1:D:89:LEU:HD11	1.91	0.51
1:D:138:LEU:HD11	1:D:228:THR:HB	1.92	0.51
1:B:199:GLY:HA3	1:B:210:ILE:HA	1.91	0.51
1:A:19:ARG:HD3	1:A:101:TYR:HD1	1.76	0.51
1:A:89:LEU:HB3	1:A:97:TYR:HB3	1.92	0.50
1:A:127:LEU:HD11	1:A:139:TYR:HB2	1.93	0.50
1:D:27:GLY:HA3	1:D:53:PHE:CD1	2.46	0.50
1:B:218:ALA:O	1:B:220:LEU:N	2.45	0.50
1:B:230:ALA:H	1:B:231:NFA:CB	2.20	0.50
1:D:87:MET:O	1:D:99:CYS:N	2.45	0.50
1:A:67:TRP:CE2	1:A:89:LEU:HD13	2.47	0.49
1:C:91:LEU:HD21	1:C:97:TYR:HB2	1.93	0.49
1:B:33:ASN:ND2	1:B:44:ASN:HB3	2.27	0.49
1:C:28:ILE:HD13	1:A:28:ILE:HD11	1.95	0.49
1:C:28:ILE:HG21	1:A:28:ILE:HD11	1.95	0.49
1:A:126:ILE:HD11	1:A:228:THR:HG21	1.95	0.49
1:B:33:ASN:HD21	1:B:44:ASN:HD22	1.60	0.49
1:C:33:ASN:OD1	1:C:34:ASP:N	2.46	0.48
1:B:139:TYR:CE2	1:B:207:ILE:HG23	2.45	0.48
1:A:222:PHE:CE2	1:A:227:LEU:HD11	2.47	0.48
1:D:48:PHE:O	1:D:186:GLY:HA2	2.12	0.48
1:B:185:GLU:HA	1:B:194:VAL:HG22	1.96	0.48
1:A:81:LEU:HD23	1:A:105:TYR:HB2	1.96	0.48
1:B:188:VAL:HG23	1:B:193:PHE:HE1	1.79	0.47
1:B:123:ARG:O	1:B:124:ILE:HG13	2.13	0.47
1:D:19:ARG:NH1	1:D:103:THR:OG1	2.47	0.47
1:A:90:LEU:H	1:A:90:LEU:HD12	1.80	0.47
1:A:76:GLU:HG3	1:A:80:VAL:O	2.15	0.46
1:D:33:ASN:HB2	1:D:46:ILE:HG23	1.98	0.46
1:D:184:MET:HB3	1:D:184:MET:HE2	1.79	0.46
1:C:90:LEU:HD12	1:C:96:HIS:CD2	2.50	0.45
1:D:23:TYR:CD1	1:D:99:CYS:HB2	2.52	0.45
1:B:183:ARG:CB	1:B:196:GLU:HG2	2.47	0.45
1:B:226:ILE:HG13	1:B:227:LEU:HG	1.99	0.44
1:B:15:TYR:CZ	1:B:35:ILE:HD12	2.53	0.44
1:A:89:LEU:O	1:A:97:TYR:N	2.50	0.44
1:A:48:PHE:O	1:A:186:GLY:HA2	2.18	0.43
1:C:228:THR:HA	1:C:231:NFA:HD1	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:44:ASN:N	1:B:44:ASN:OD1	2.51	0.43
1:B:60:MET:SD	1:B:189:ASN:ND2	2.91	0.43
1:B:85:VAL:HG13	1:B:101:TYR:HB2	1.99	0.43
1:C:127:LEU:HD11	1:C:139:TYR:HB2	1.99	0.43
1:D:180:ASN:ND2	1:D:208:GLN:OE1	2.49	0.43
1:D:64:THR:OG1	1:D:225:ASP:OD2	2.29	0.43
1:C:142:ALA:HB3	1:C:205:GLU:HA	2.00	0.43
1:D:72:GLU:HB3	1:D:120:VAL:HB	2.00	0.43
1:D:35:ILE:HG12	1:D:44:ASN:OD1	2.19	0.42
1:A:31:ALA:HB2	1:A:48:PHE:CD1	2.55	0.42
1:C:52:ASN:ND2	1:A:102:ARG:O	2.53	0.42
1:B:51:THR:HG23	1:B:52:ASN:OD1	2.20	0.42
1:D:184:MET:HE1	1:D:231:NFA:CZ	2.49	0.42
1:B:124:ILE:O	1:B:124:ILE:CG2	2.68	0.41
1:C:30:ILE:HD12	1:A:30:ILE:HG12	2.02	0.41
1:D:223:ALA:O	1:D:226:ILE:HG12	2.20	0.41
1:D:231:NFA:HA	1:D:231:NFA:HD2	1.81	0.41
1:B:100:ASP:OD1	1:B:100:ASP:N	2.54	0.41
1:B:19:ARG:HD3	1:B:101:TYR:CD1	2.55	0.41
1:C:25:ASP:OD1	1:C:97:TYR:OH	2.37	0.41
1:C:19:ARG:HG3	1:C:103:THR:HG22	2.01	0.40
1:A:80:VAL:HG22	1:A:81:LEU:H	1.86	0.40
1:B:71:THR:HG23	1:B:121:ASP:OD1	2.22	0.40

There are no symmetry-related clashes.

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	153/242 (63%)	137 (90%)	13 (8%)	3 (2%)	7	23
1	B	150/242 (62%)	129 (86%)	19 (13%)	2 (1%)	12	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	182/242 (75%)	172 (94%)	9 (5%)	1 (0%)	29	59
1	D	146/242 (60%)	136 (93%)	9 (6%)	1 (1%)	22	51
All	All	631/968 (65%)	574 (91%)	50 (8%)	7 (1%)	14	39

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	182	LEU
1	A	80	VAL
1	A	131	LYS
1	B	219	PRO
1	C	115	PRO
1	A	83	GLY
1	D	85	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	119/208 (57%)	116 (98%)	3 (2%)	47	78
1	B	108/208 (52%)	102 (94%)	6 (6%)	21	49
1	C	135/208 (65%)	132 (98%)	3 (2%)	52	81
1	D	103/208 (50%)	99 (96%)	4 (4%)	32	64
All	All	465/832 (56%)	449 (97%)	16 (3%)	37	69

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	41	SER
1	C	182	LEU
1	C	215	LYS
1	A	20	SER
1	A	49	LYS

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Mol	Chain	Res	Type
1	A	204	PHE
1	B	40	ASP
1	B	44	ASN
1	B	100	ASP
1	B	126	ILE
1	B	141	HIS
1	B	225	ASP
1	D	81	LEU
1	D	130	ASP
1	D	182	LEU
1	D	222	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	B	44	ASN
1	D	187	ASN

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	$\# Z > 2$
1	NFA	A	231	1	12,12,12	1.67	2 (16%)	15,15,15	0.92	1 (6%)
1	NFA	C	231	1	12,12,12	1.66	2 (16%)	15,15,15	0.90	0
1	NFA	B	231	1	12,12,12	1.65	2 (16%)	15,15,15	1.23	2 (13%)
1	NFA	D	231	1	12,12,12	1.65	2 (16%)	15,15,15	0.97	1 (6%)

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
1	NFA	A	231	1	-	1/8/8/8	0/1/1/1
1	NFA	C	231	1	-	5/8/8/8	0/1/1/1
1	NFA	B	231	1	-	6/8/8/8	0/1/1/1
1	NFA	D	231	1	-	5/8/8/8	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	231	NFA	C-NXT	5.05	1.45	1.32
1	C	231	NFA	C-NXT	5.04	1.45	1.32
1	B	231	NFA	C-NXT	5.03	1.45	1.32
1	D	231	NFA	C-NXT	5.01	1.45	1.32
1	B	231	NFA	O-C	-2.58	1.18	1.23
1	A	231	NFA	O-C	-2.54	1.19	1.23
1	C	231	NFA	O-C	-2.53	1.19	1.23
1	D	231	NFA	O-C	-2.50	1.19	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	231	NFA	CA-C-NXT	3.32	122.38	116.68
1	B	231	NFA	O-C-NXT	-2.24	119.11	123.00
1	A	231	NFA	CA-C-NXT	2.06	120.22	116.68
1	D	231	NFA	CA-C-NXT	2.01	120.13	116.68

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	231	NFA	O-C-CA-N
1	C	231	NFA	O-C-CA-CB
1	C	231	NFA	NXT-C-CA-CB
1	A	231	NFA	NXT-C-CA-N
1	B	231	NFA	NXT-C-CA-N
1	D	231	NFA	O-C-CA-N
1	B	231	NFA	N-CA-CB-CG
1	D	231	NFA	N-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
1	D	231	NFA	CA-CB-CG-CD2
1	D	231	NFA	CA-CB-CG-CD1
1	B	231	NFA	CA-CB-CG-CD2
1	B	231	NFA	CA-CB-CG-CD1
1	B	231	NFA	O-C-CA-N
1	B	231	NFA	C-CA-CB-CG
1	D	231	NFA	C-CA-CB-CG
1	C	231	NFA	CA-CB-CG-CD2
1	C	231	NFA	CA-CB-CG-CD1

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	231	NFA	2	0
1	C	231	NFA	1	0
1	B	231	NFA	3	0
1	D	231	NFA	3	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 [i](#)

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

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	164/242 (67%)	0.50	5 (3%)	50	40	8, 35, 66, 86	1 (0%)
1	B	161/242 (66%)	0.68	15 (9%)	8	4	8, 39, 75, 91	0
1	C	189/242 (78%)	0.40	9 (4%)	30	21	3, 29, 55, 75	1 (0%)
1	D	157/242 (64%)	0.71	13 (8%)	11	6	12, 45, 73, 86	1 (0%)
All	All	671/968 (69%)	0.57	42 (6%)	20	12	3, 36, 70, 91	3 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	TYR	4.3
1	B	97	TYR	4.3
1	B	221	PRO	3.7
1	C	44	ASN	3.6
1	B	57	GLY	3.5
1	B	55	PRO	3.5
1	D	143	VAL	3.4
1	B	182	LEU	3.3
1	B	56	ASN	3.3
1	C	114	LEU	3.1
1	D	117	TYR	3.1
1	D	81	LEU	3.0
1	B	60	MET	3.0
1	D	85	VAL	3.0
1	D	51	THR	2.8
1	D	104	THR	2.8
1	C	207	ILE	2.7
1	D	84	ASP	2.7
1	B	44	ASN	2.7
1	B	95	GLY	2.6
1	B	14	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	133	TYR	2.5
1	C	36	THR	2.4
1	C	118	HIS	2.4
1	C	204	PHE	2.4
1	B	81	LEU	2.3
1	D	56	ASN	2.3
1	D	139	TYR	2.3
1	A	58	PRO	2.3
1	D	219	PRO	2.3
1	B	116	ASP	2.2
1	A	91	LEU	2.2
1	D	47	HIS	2.2
1	A	77	ARG	2.2
1	A	97	TYR	2.2
1	C	61	GLN	2.1
1	B	76	GLU	2.1
1	D	30	ILE	2.1
1	C	10	SER	2.1
1	D	52	ASN	2.1
1	B	105	TYR	2.0
1	A	56	ASN	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(\AA^2)	Q<0.9
1	NFA	D	231	12/12	0.81	0.39	44,48,53,55	0
1	NFA	B	231	12/12	0.84	0.24	36,39,44,46	0
1	NFA	A	231	12/12	0.90	0.24	26,27,33,36	0
1	NFA	C	231	12/12	0.91	0.20	19,24,28,29	0

6.3 Carbohydrates [i](#)

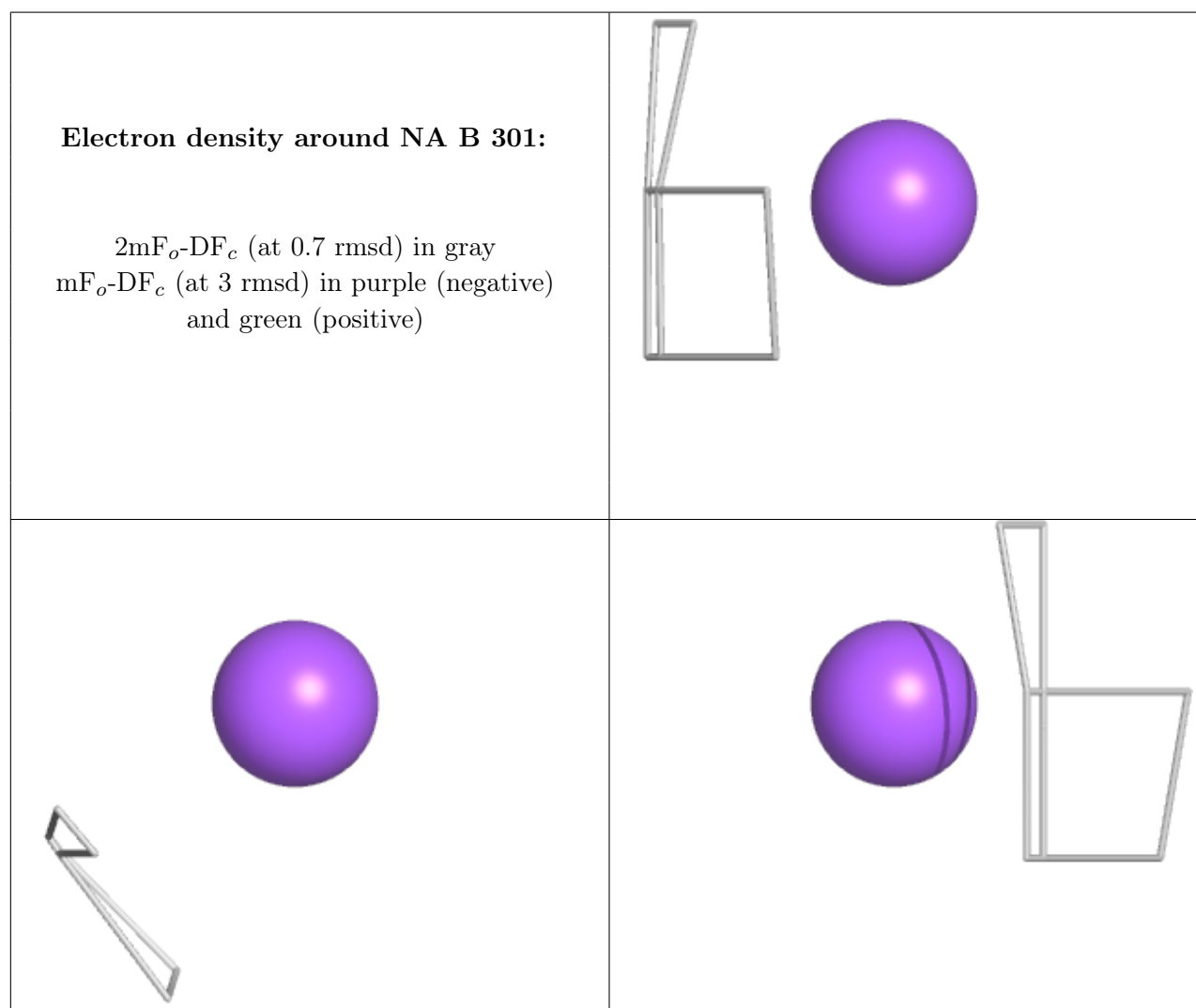
There are no monosaccharides in this entry.

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	NA	B	301	1/1	0.84	0.14	2,2,2,2	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.