



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 7, 2021 – 12:02 PM EDT

PDB ID : 7KY2
Title : Botulism Neurotoxin Light Chain A app form
Authors : Ortega, M.E.; Salazameda, N.T.
Deposited on : 2020-12-07
Resolution : 2.78 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
Xtriage (Phenix) : 1.13
EDS : 2.18
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) : 2.18

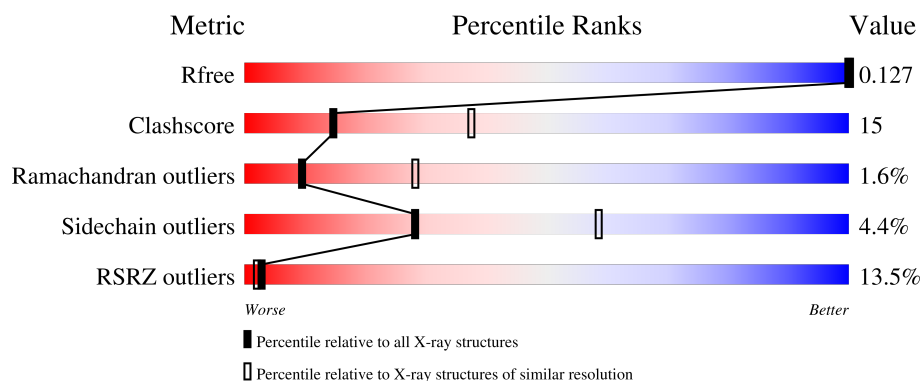
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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	4107 (2.80-2.76)
Clashscore	141614	4575 (2.80-2.76)
Ramachandran outliers	138981	4487 (2.80-2.76)
Sidechain outliers	138945	4489 (2.80-2.76)
RSRZ outliers	127900	4027 (2.80-2.76)

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	426	<div> <div>10%</div> <div>62%</div> <div>27%</div> <div>8%</div> </div>
1	B	426	<div> <div>15%</div> <div>57%</div> <div>32%</div> <div>8%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6419 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 Bont/A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			3199	2067	526	598	8			
1	B	393	Total	C	N	O	S	0	0	0
			3199	2067	526	598	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP C6K838
A	-4	HIS	-	expression tag	UNP C6K838
A	-3	HIS	-	expression tag	UNP C6K838
A	-2	HIS	-	expression tag	UNP C6K838
A	-1	HIS	-	expression tag	UNP C6K838
A	0	HIS	-	expression tag	UNP C6K838
A	1	MET	-	expression tag	UNP C6K838
A	2	GLN	-	expression tag	UNP C6K838
B	-5	HIS	-	expression tag	UNP C6K838
B	-4	HIS	-	expression tag	UNP C6K838
B	-3	HIS	-	expression tag	UNP C6K838
B	-2	HIS	-	expression tag	UNP C6K838
B	-1	HIS	-	expression tag	UNP C6K838
B	0	HIS	-	expression tag	UNP C6K838
B	1	MET	-	expression tag	UNP C6K838
B	2	GLN	-	expression tag	UNP C6K838

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

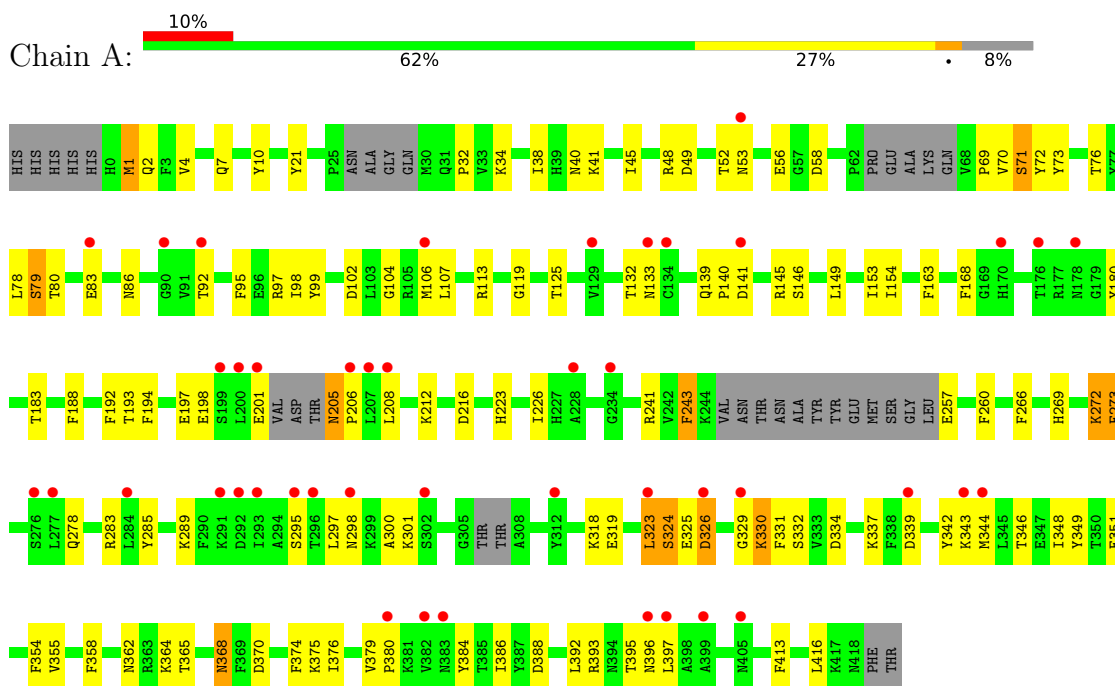
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total 8	O 8	0	0
3	B	11	Total 11	O 11	0	0

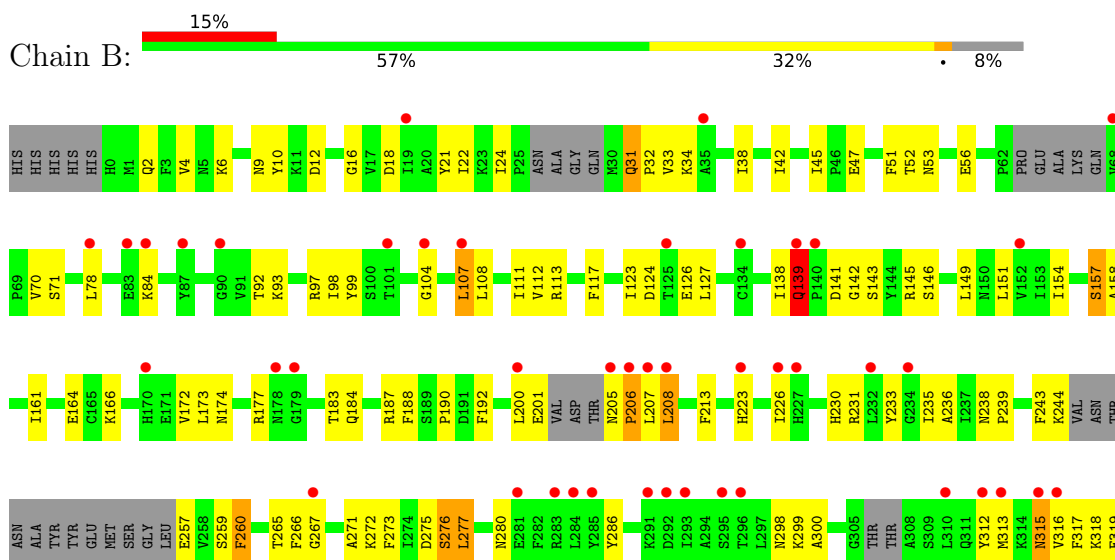
3 Residue-property plots

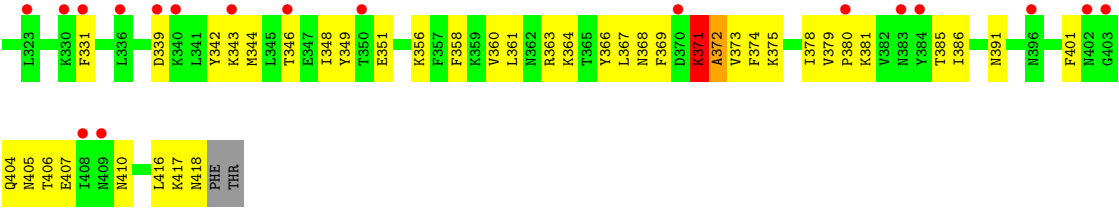
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Bont/A1



• Molecule 1: Bont/A1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.03Å 56.98Å 192.14Å 90.00° 90.26° 90.00°	Depositor
Resolution (Å)	49.01 – 2.78 49.01 – 2.78	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.01-2.78) 99.3 (49.01-2.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	50.45 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.128 , 0.170 0.112 , 0.127	Depositor DCC
R_{free} test set	693 reflections (3.23%)	wwPDB-VP
Wilson B-factor (Å ²)	66.0	Xtriage
Anisotropy	0.108	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.48 , 999.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.147 for h,-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	6419	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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: ZN

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	0.73	0/3271	0.89	1/4416 (0.0%)
1	B	0.70	0/3271	0.88	0/4416
All	All	0.72	0/6542	0.89	1/8832 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	326	ASP	CB-CA-C	5.51	121.42	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	139	GLN	Peptide

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	3199	0	3157	89	0
1	B	3199	0	3157	109	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	8	0	0	5	0
3	B	11	0	0	8	0
All	All	6419	0	6314	194	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 15.

All (194) 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:318:LYS:HG3	1:A:323:LEU:HB2	1.46	0.94
1:B:244:LYS:HB3	1:B:257:GLU:HB2	1.48	0.92
1:A:318:LYS:CG	1:A:323:LEU:HB2	2.03	0.87
1:A:358:PHE:CZ	3:A:608:HOH:O	2.30	0.85
1:A:243:PHE:CZ	1:A:273:PHE:O	2.29	0.84
1:A:21:TYR:CZ	1:B:141:ASP:HB2	2.15	0.82
1:A:325:GLU:HA	1:A:330:LYS:HG2	1.66	0.77
1:A:318:LYS:HG3	1:A:323:LEU:CB	2.14	0.77
1:B:201:GLU:HB3	1:B:205:ASN:O	1.85	0.77
1:A:205:ASN:N	1:A:205:ASN:OD1	2.21	0.74
1:B:206:PRO:HB2	1:B:207:LEU:HD12	1.70	0.73
1:B:257:GLU:O	1:B:257:GLU:HG2	1.86	0.73
1:B:31:GLN:HB2	1:B:32:PRO:HD2	1.71	0.72
1:A:201:GLU:C	1:A:205:ASN:O	2.29	0.70
1:A:4:VAL:HG21	1:A:92:THR:HG23	1.74	0.70
1:B:339:ASP:O	1:B:343:LYS:HG2	1.94	0.68
1:A:140:PRO:O	1:B:139:GLN:OE1	2.12	0.67
1:B:70:VAL:HG21	1:B:371:LYS:HA	1.76	0.67
1:B:313:MET:CE	3:B:606:HOH:O	2.41	0.67
1:A:10:TYR:O	1:A:34:LYS:NZ	2.28	0.67
1:A:71:SER:HB3	1:A:73:TYR:HE1	1.62	0.65
1:B:267:GLY:HA2	1:B:271:ALA:HB2	1.78	0.65
1:B:205:ASN:N	1:B:206:PRO:HD3	2.12	0.64
1:B:174:ASN:OD1	1:B:177:ARG:NH2	2.31	0.63
1:A:21:TYR:CE1	1:B:141:ASP:HB2	2.32	0.63
1:B:93:LYS:NZ	1:B:379:VAL:O	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:313:MET:HE1	3:B:606:HOH:O	1.98	0.63
1:B:139:GLN:HG3	1:B:142:GLY:H	1.64	0.62
1:B:344:MET:HA	1:B:348:ILE:HD12	1.82	0.62
1:A:107:LEU:HD13	1:A:349:TYR:CD2	2.35	0.61
1:A:119:GLY:HA3	1:A:180:TYR:CE2	2.36	0.61
1:A:45:ILE:HB	1:A:154:ILE:HG13	1.83	0.60
1:A:388:ASP:OD2	1:A:393:ARG:HG3	2.02	0.60
1:A:97:ARG:HA	1:A:386:ILE:HG23	1.83	0.60
1:B:172:VAL:HG13	1:B:173:LEU:HG	1.83	0.60
1:B:243:PHE:HE1	1:B:260:PHE:CE1	2.20	0.60
1:B:124:ASP:OD2	1:B:299:LYS:NZ	2.29	0.60
1:A:205:ASN:N	1:A:206:PRO:HD3	2.17	0.60
1:A:283:ARG:HD3	1:A:343:LYS:HZ2	1.66	0.59
1:A:283:ARG:HD3	1:A:343:LYS:NZ	2.17	0.59
1:A:223:HIS:CG	3:A:603:HOH:O	2.57	0.58
1:B:70:VAL:CG2	1:B:371:LYS:HA	2.33	0.57
1:A:194:PHE:HB2	1:A:370:ASP:HB3	1.87	0.57
1:A:298:ASN:OD1	1:A:331:PHE:CD2	2.59	0.56
1:A:49:ASP:OD2	1:A:52:THR:OG1	2.21	0.56
1:A:197:GLU:OE2	1:A:197:GLU:HA	2.05	0.56
1:A:98:ILE:O	1:A:104:GLY:HA3	2.06	0.56
1:B:276:SER:O	1:B:280:ASN:HB2	2.05	0.56
1:A:133:ASN:HA	1:A:183:THR:OG1	2.05	0.56
1:A:388:ASP:OD2	1:A:393:ARG:NH1	2.39	0.56
1:B:243:PHE:CE1	1:B:260:PHE:CE1	2.93	0.56
1:B:266:PHE:HA	1:B:351:GLU:HB3	1.88	0.56
1:A:21:TYR:HB3	1:A:32:PRO:HB2	1.88	0.56
1:B:70:VAL:HG12	1:B:161:ILE:HD11	1.88	0.55
1:A:297:LEU:O	1:A:300:ALA:HB3	2.07	0.55
1:B:53:ASN:HB3	1:B:56:GLU:HB2	1.87	0.55
1:B:375:LYS:HB2	1:B:416:LEU:HD11	1.89	0.54
1:A:283:ARG:NH1	1:A:343:LYS:HE3	2.22	0.54
1:A:243:PHE:O	1:A:257:GLU:HA	2.07	0.54
1:B:108:LEU:O	1:B:112:VAL:HG23	2.07	0.54
1:B:45:ILE:HB	1:B:154:ILE:HD13	1.90	0.54
1:A:339:ASP:O	1:A:343:LYS:HD3	2.07	0.54
1:B:98:ILE:O	1:B:104:GLY:HA3	2.06	0.53
1:B:78:LEU:O	1:B:84:LYS:NZ	2.41	0.53
1:A:53:ASN:HB3	1:A:56:GLU:HB2	1.90	0.53
1:A:392:LEU:HB2	1:A:395:THR:HG21	1.91	0.53
1:B:10:TYR:CE2	1:B:84:LYS:HD2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:LEU:HD13	1:B:300:ALA:CB	2.38	0.53
1:A:107:LEU:CD1	1:A:349:TYR:CD2	2.91	0.53
1:A:342:TYR:O	1:A:346:THR:HG23	2.09	0.53
1:A:243:PHE:HD1	1:A:260:PHE:HE1	1.56	0.53
1:B:138:ILE:HA	1:B:143:SER:O	2.09	0.53
1:B:97:ARG:HA	1:B:386:ILE:HG23	1.90	0.52
1:B:70:VAL:HG21	1:B:371:LYS:CA	2.38	0.52
1:A:334:ASP:HB3	1:A:337:LYS:HB2	1.91	0.52
1:B:113:ARG:NH1	1:B:319:GLU:O	2.42	0.52
1:B:356:LYS:HE2	3:B:605:HOH:O	2.09	0.52
1:B:417:LYS:HD3	1:B:418:ASN:O	2.10	0.52
1:B:9:ASN:O	1:B:12:ASP:HB2	2.10	0.52
1:A:243:PHE:HZ	1:A:273:PHE:O	1.90	0.52
1:A:113:ARG:HH21	1:A:319:GLU:HG2	1.75	0.52
1:B:298:ASN:HD21	1:B:331:PHE:H	1.58	0.52
1:B:244:LYS:HB3	1:B:257:GLU:CB	2.32	0.51
1:A:201:GLU:HA	1:A:201:GLU:OE1	2.10	0.51
1:A:48:ARG:HG2	1:A:78:LEU:HB3	1.93	0.51
1:A:285:TYR:CZ	1:A:289:LYS:HE3	2.46	0.51
1:B:361:LEU:HB2	1:B:401:PHE:CG	2.46	0.51
1:B:33:VAL:HG11	1:B:51:PHE:CZ	2.46	0.50
1:B:184:GLN:NE2	1:B:231:ARG:HB3	2.26	0.50
1:B:188:PHE:CD1	3:B:610:HOH:O	2.55	0.50
1:A:243:PHE:CE1	1:A:273:PHE:O	2.64	0.50
1:A:375:LYS:HB2	1:A:416:LEU:HD11	1.94	0.50
1:B:31:GLN:CB	1:B:32:PRO:HD2	2.40	0.50
1:B:364:LYS:CE	3:B:603:HOH:O	2.60	0.50
1:B:318:LYS:HG3	1:B:331:PHE:CE1	2.47	0.50
1:A:354:PHE:CE1	3:A:608:HOH:O	2.65	0.49
1:B:52:THR:OG1	1:B:166:LYS:HE3	2.13	0.49
1:A:193:THR:HB	1:A:376:ILE:HD13	1.93	0.49
1:A:198:GLU:OE1	1:A:362:ASN:O	2.30	0.49
1:A:216:ASP:OD1	1:A:384:TYR:OH	2.20	0.49
1:B:275:ASP:O	1:B:277:LEU:N	2.45	0.49
1:B:188:PHE:HD1	3:B:610:HOH:O	1.93	0.49
1:A:4:VAL:HG22	1:A:38:ILE:HB	1.94	0.49
1:B:4:VAL:HG21	1:B:92:THR:HG23	1.94	0.49
1:B:10:TYR:OH	1:B:47:GLU:OE2	2.30	0.48
1:A:86:ASN:OD1	1:A:379:VAL:HG21	2.14	0.48
1:B:208:LEU:HA	1:B:208:LEU:HD23	1.69	0.48
1:B:192:PHE:CE2	1:B:375:LYS:HD3	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:HIS:NE2	1:A:365:THR:HB	2.28	0.48
1:A:141:ASP:HB2	1:B:21:TYR:CZ	2.48	0.47
1:A:226:ILE:HG12	1:A:349:TYR:HB3	1.96	0.47
1:B:358:PHE:HB3	1:B:360:VAL:HG13	1.96	0.47
1:A:1:MET:SD	1:A:7:GLN:HG2	2.55	0.47
1:B:230:HIS:HE1	1:B:265:THR:OG1	1.97	0.47
1:A:354:PHE:CD1	3:A:608:HOH:O	2.55	0.47
1:A:95:PHE:HE2	1:A:153:ILE:HG13	1.80	0.47
1:B:123:ILE:HB	1:B:126:GLU:HB3	1.96	0.47
1:B:24:ILE:HD12	1:B:51:PHE:HB2	1.97	0.47
1:B:244:LYS:HE2	1:B:257:GLU:HB2	1.96	0.47
1:B:373:VAL:HG23	1:B:416:LEU:HB2	1.96	0.47
1:A:298:ASN:OD1	1:A:331:PHE:CB	2.63	0.47
1:B:10:TYR:CZ	1:B:84:LYS:HD2	2.50	0.47
1:B:117:PHE:HA	1:B:317:PHE:CZ	2.50	0.46
1:B:16:GLY:CA	1:B:139:GLN:O	2.63	0.46
1:B:363:ARG:HE	1:B:366:TYR:HA	1.81	0.46
1:B:200:LEU:HD21	1:B:361:LEU:HD21	1.97	0.46
1:B:107:LEU:O	1:B:111:ILE:HG13	2.15	0.46
1:A:379:VAL:HB	1:A:380:PRO:HD3	1.98	0.46
1:A:344:MET:HA	1:A:348:ILE:HD12	1.97	0.46
1:B:226:ILE:HG22	1:B:230:HIS:CE1	2.51	0.46
1:A:139:GLN:HG3	1:A:145:ARG:HG2	1.97	0.46
1:B:12:ASP:O	1:B:34:LYS:NZ	2.36	0.46
1:B:298:ASN:ND2	1:B:331:PHE:H	2.12	0.46
1:A:343:LYS:HD2	1:A:343:LYS:N	2.30	0.45
1:B:405:ASN:OD1	1:B:407:GLU:HG2	2.16	0.45
1:B:233:TYR:O	1:B:235:ILE:HG23	2.17	0.45
1:B:364:LYS:HE2	3:B:603:HOH:O	2.17	0.45
1:A:243:PHE:CD1	1:A:260:PHE:HE1	2.34	0.45
1:B:379:VAL:HB	1:B:380:PRO:HD3	1.96	0.45
1:B:391:ASN:OD1	1:B:404:GLN:HG3	2.17	0.45
1:B:364:LYS:HE3	3:B:603:HOH:O	2.17	0.45
1:B:70:VAL:HG21	1:B:371:LYS:O	2.16	0.44
1:A:351:GLU:O	1:A:355:VAL:HG23	2.17	0.44
1:B:372:ALA:HB1	1:B:374:PHE:CE2	2.52	0.44
1:A:40:ASN:OD1	1:A:41:LYS:NZ	2.43	0.44
1:A:241:ARG:NH2	1:A:278:GLN:OE1	2.41	0.44
1:B:157:SER:OG	1:B:158:ALA:N	2.47	0.44
1:B:315:ASN:HA	1:B:318:LYS:CB	2.48	0.44
1:B:164:GLU:HG2	1:B:187:ARG:HD2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:THR:HG23	1:A:83:GLU:H	1.83	0.43
1:A:80:THR:HG22	1:A:83:GLU:OE1	2.18	0.43
1:A:318:LYS:HD2	1:A:330:LYS:HD3	1.98	0.43
1:A:324:SER:OG	1:A:332:SER:O	2.35	0.43
1:B:344:MET:HG2	1:B:349:TYR:CZ	2.53	0.43
1:B:4:VAL:HA	1:B:38:ILE:O	2.19	0.43
1:B:177:ARG:HG2	1:B:236:ALA:O	2.18	0.43
1:B:238:ASN:HA	1:B:239:PRO:HD2	1.85	0.43
1:B:205:ASN:N	1:B:206:PRO:CD	2.81	0.43
1:A:132:THR:HB	1:A:168:PHE:HB2	2.00	0.43
1:B:381:LYS:HG2	1:B:385:THR:HG22	2.01	0.43
1:B:2:GLN:HG2	1:B:99:TYR:HE2	1.84	0.43
1:B:71:SER:H	1:B:418:ASN:ND2	2.17	0.42
1:B:363:ARG:HH11	1:B:366:TYR:HD1	1.66	0.42
1:A:395:THR:OG1	1:A:396:ASN:N	2.52	0.42
1:B:373:VAL:HG23	1:B:373:VAL:O	2.20	0.42
1:B:213:PHE:CD2	1:B:361:LEU:HD23	2.55	0.42
1:B:235:ILE:HA	1:B:286:TYR:CE1	2.55	0.42
1:A:354:PHE:HD1	3:A:608:HOH:O	2.00	0.42
1:B:6:LYS:NZ	1:B:18:ASP:OD2	2.43	0.42
1:B:149:LEU:HD21	1:B:183:THR:HG21	2.02	0.42
1:A:72:TYR:CE2	1:A:416:LEU:HD22	2.55	0.41
1:B:223:HIS:HA	1:B:226:ILE:HD12	2.02	0.41
1:B:374:PHE:CE2	1:B:406:THR:HG21	2.56	0.41
1:B:22:ILE:HD11	1:B:45:ILE:HD11	2.02	0.41
1:A:102:ASP:O	1:A:106:MET:HG3	2.20	0.41
1:A:374:PHE:CD1	1:A:413:PHE:HB3	2.56	0.41
1:A:163:PHE:CD1	1:A:188:PHE:HA	2.55	0.41
1:A:272:LYS:HD2	1:A:272:LYS:O	2.20	0.41
1:A:2:GLN:HB3	1:A:99:TYR:HE2	1.86	0.41
1:A:125:THR:OG1	1:A:301:LYS:HB2	2.21	0.41
1:A:266:PHE:HA	1:A:351:GLU:HB3	2.03	0.41
1:A:192:PHE:CE2	1:A:375:LYS:HD3	2.56	0.41
1:B:190:PRO:HB2	1:B:378:ILE:HD11	2.03	0.41
1:B:342:TYR:O	1:B:346:THR:HG23	2.21	0.41
1:A:76:THR:O	1:A:79:SER:HB2	2.20	0.40
1:B:93:LYS:HD3	1:B:378:ILE:O	2.21	0.40
1:B:127:LEU:HD13	1:B:300:ALA:HB1	2.02	0.40
1:A:41:LYS:HA	1:A:41:LYS:HD3	1.89	0.40
1:B:42:ILE:HG12	1:B:151:LEU:HB3	2.02	0.40
1:B:272:LYS:HG3	1:B:273:PHE:N	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:HIS:CD2	1:A:364:LYS:O	2.74	0.40
1:A:395:THR:HG21	1:A:397:LEU:HD12	2.03	0.40
1:B:16:GLY:HA3	1:B:139:GLN:O	2.21	0.40
1:B:312:TYR:O	1:B:316:VAL:HG23	2.21	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	381/426 (89%)	353 (93%)	24 (6%)	4 (1%)	15	41
1	B	381/426 (89%)	351 (92%)	22 (6%)	8 (2%)	7	21
All	All	762/852 (89%)	704 (92%)	46 (6%)	12 (2%)	9	28

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	330	LYS
1	A	368	ASN
1	B	276	SER
1	B	371	LYS
1	A	329	GLY
1	B	206	PRO
1	B	367	LEU
1	B	372	ALA
1	B	277	LEU
1	B	157	SER
1	B	139	GLN
1	A	69	PRO

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	354/382 (93%)	336 (95%)	18 (5%)	24	53
1	B	354/382 (93%)	341 (96%)	13 (4%)	34	65
All	All	708/764 (93%)	677 (96%)	31 (4%)	28	58

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	58	ASP
1	A	70	VAL
1	A	71	SER
1	A	79	SER
1	A	146	SER
1	A	149	LEU
1	A	205	ASN
1	A	208	LEU
1	A	212	LYS
1	A	243	PHE
1	A	272	LYS
1	A	273	PHE
1	A	295	SER
1	A	323	LEU
1	A	324	SER
1	A	326	ASP
1	A	368	ASN
1	B	31	GLN
1	B	107	LEU
1	B	139	GLN
1	B	145	ARG
1	B	146	SER
1	B	208	LEU
1	B	259	SER
1	B	260	PHE
1	B	315	ASN

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Mol	Chain	Res	Type
1	B	368	ASN
1	B	369	PHE
1	B	371	LYS
1	B	410	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	205	ASN
1	A	311	GLN
1	A	315	ASN
1	B	394	ASN
1	B	418	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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 [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	393/426 (92%)	0.99	44 (11%) 5 3	23, 44, 78, 107	0
1	B	393/426 (92%)	1.11	62 (15%) 2 1	26, 54, 84, 104	0
All	All	786/852 (92%)	1.05	106 (13%) 3 2	23, 49, 81, 107	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	GLN	6.2
1	A	329	GLY	6.0
1	B	200	LEU	4.9
1	B	178	ASN	4.5
1	B	330	LYS	3.9
1	A	343	LYS	3.7
1	B	346	THR	3.6
1	B	315	ASN	3.6
1	A	170	HIS	3.5
1	A	207	LEU	3.5
1	B	403	GLY	3.5
1	B	207	LEU	3.4
1	B	408	ILE	3.3
1	B	293	ILE	3.3
1	A	276	SER	3.2
1	A	293	ILE	3.1
1	B	296	THR	3.0
1	B	383	ASN	3.0
1	A	208	LEU	3.0
1	A	206	PRO	2.9
1	A	383	ASN	2.9
1	A	129	VAL	2.9
1	B	291	LYS	2.9
1	B	227	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	200	LEU	2.9
1	B	234	GLY	2.8
1	A	106	MET	2.8
1	A	90	GLY	2.8
1	B	206	PRO	2.8
1	B	125	THR	2.8
1	A	295	SER	2.8
1	B	384	TYR	2.8
1	B	331	PHE	2.8
1	A	296	THR	2.7
1	B	285	TYR	2.7
1	B	78	LEU	2.7
1	B	409	ASN	2.6
1	A	141	ASP	2.6
1	B	284	LEU	2.6
1	B	402	ASN	2.6
1	B	107	LEU	2.6
1	B	339	ASP	2.6
1	A	397	LEU	2.5
1	B	267	GLY	2.5
1	B	310	LEU	2.5
1	A	380	PRO	2.5
1	A	291	LYS	2.5
1	B	226	ILE	2.5
1	A	302	SER	2.5
1	B	292	ASP	2.5
1	B	312	TYR	2.5
1	A	53	ASN	2.4
1	B	84	LYS	2.4
1	B	343	LYS	2.4
1	A	234	GLY	2.4
1	B	350	THR	2.4
1	A	405	ASN	2.4
1	B	179	GLY	2.4
1	B	396	ASN	2.4
1	A	199	SER	2.4
1	B	87	TYR	2.3
1	B	134	CYS	2.3
1	A	277	LEU	2.3
1	B	19	ILE	2.3
1	A	344	MET	2.3
1	B	223	HIS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	382	VAL	2.3
1	B	170	HIS	2.3
1	B	83	GLU	2.3
1	A	292	ASP	2.3
1	B	208	LEU	2.2
1	A	399	ALA	2.2
1	B	281	GLU	2.2
1	B	232	LEU	2.2
1	A	228	ALA	2.2
1	B	90	GLY	2.2
1	A	83	GLU	2.2
1	B	340	LYS	2.2
1	B	35	ALA	2.2
1	B	101	THR	2.2
1	B	205	ASN	2.2
1	A	326	ASP	2.2
1	A	339	ASP	2.2
1	B	380	PRO	2.1
1	A	134	CYS	2.1
1	B	295	SER	2.1
1	B	140	PRO	2.1
1	A	312	TYR	2.1
1	A	323	LEU	2.1
1	B	283	ARG	2.1
1	B	104	GLY	2.1
1	A	92	THR	2.1
1	A	176	THR	2.1
1	A	178	ASN	2.1
1	B	316	VAL	2.1
1	A	284	LEU	2.0
1	B	336	LEU	2.0
1	B	68	VAL	2.0
1	B	152	VAL	2.0
1	A	201	GLU	2.0
1	A	396	ASN	2.0
1	B	313	MET	2.0
1	B	323	LEU	2.0
1	A	133	ASN	2.0
1	A	298	ASN	2.0
1	B	370	ASP	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(\AA^2)	Q<0.9
2	ZN	A	500	1/1	0.84	0.86	162,162,162,162	0
2	ZN	B	500	1/1	0.98	0.85	197,197,197,197	0

6.5 Other polymers [i](#)

There are no such residues in this entry.