



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:55 PM UTC

PDB ID : 7KY2 / pdb_00007ky2
Title : Botulism Neurotoxin Light Chain A app form
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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 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

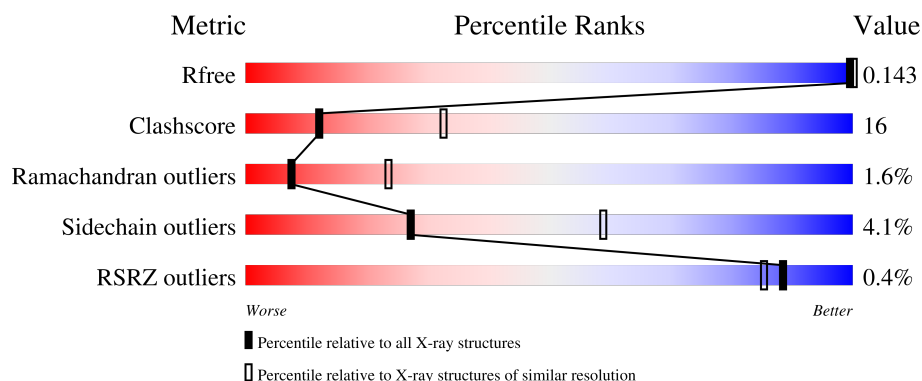
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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}	180053	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (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	 60% 29% 8%
1	B	426	 56% 33% 8%

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 (CCD ID: 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		

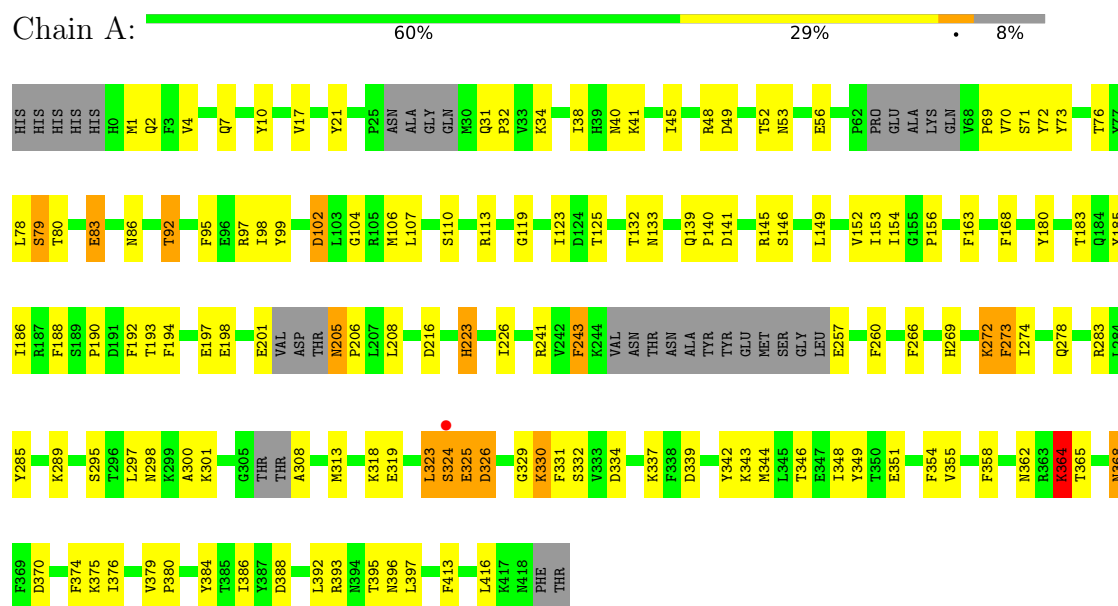
- Molecule 3 is water.

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

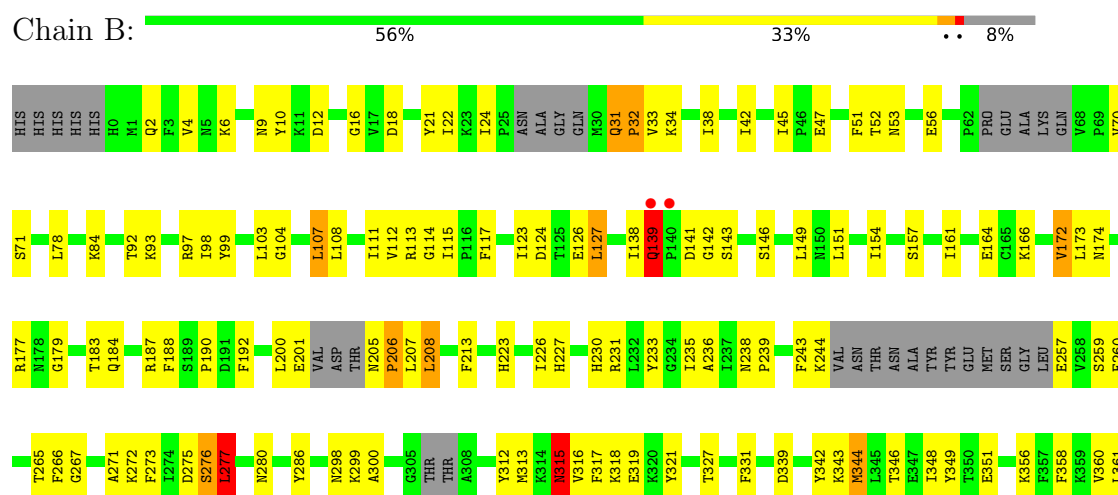
3 Residue-property plots

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



K362	R363	K364	T365	Y366	L367	N368	F369	D370	K371	A372	V373	F374	K375		I378	V379	P380	K381		T385	I386		N391		F401		Q404	N405	T406	E407		N410		L416	K417	N418	PHE	THR
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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) 99.1 (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 0.124 , 0.143	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)	0.39 , 325.3	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	0.96	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

5.1 Standard geometry

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	1.08	8/3271 (0.2%)	1.41	12/4416 (0.3%)
1	B	1.04	1/3271 (0.0%)	1.42	10/4416 (0.2%)
All	All	1.06	9/6542 (0.1%)	1.41	22/8832 (0.2%)

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

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	185	TYR	C-O	-5.98	1.17	1.24
1	A	186	ILE	C-O	-5.66	1.18	1.24
1	A	17	VAL	N-CA	5.41	1.51	1.46
1	A	83	GLU	C-O	-5.28	1.18	1.24
1	A	156	PRO	C-O	-5.23	1.17	1.23
1	A	152	VAL	C-O	-5.16	1.18	1.24
1	A	31	GLN	N-CA	5.14	1.50	1.45
1	B	227	HIS	C-O	-5.10	1.18	1.24
1	A	223	HIS	C-O	-5.08	1.18	1.24

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	326	ASP	CB-CA-C	7.27	121.42	109.72
1	A	273	PHE	N-CA-C	-6.70	103.79	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	344	MET	N-CA-C	-6.25	104.55	111.36
1	A	102	ASP	N-CA-C	-5.96	104.86	111.36
1	B	370	ASP	CA-C-N	5.91	132.83	121.54
1	B	370	ASP	C-N-CA	5.91	132.83	121.54
1	A	274	ILE	CA-C-N	5.86	129.03	120.71
1	A	274	ILE	C-N-CA	5.86	129.03	120.71
1	B	32	PRO	N-CA-C	5.69	124.20	112.47
1	B	103	LEU	CA-C-N	5.66	126.26	119.98
1	B	103	LEU	C-N-CA	5.66	126.26	119.98
1	B	277	LEU	CA-C-N	5.60	128.24	120.29
1	B	277	LEU	C-N-CA	5.60	128.24	120.29
1	A	92	THR	N-CA-C	-5.48	105.22	111.14
1	B	315	ASN	CB-CA-C	5.38	120.05	109.72
1	A	110	SER	N-CA-C	-5.22	105.59	111.28
1	A	364	LYS	N-CA-C	-5.19	106.95	113.28
1	A	326	ASP	CA-CB-CG	5.18	117.78	112.60
1	A	123	ILE	CA-C-N	5.16	128.15	120.31
1	A	123	ILE	C-N-CA	5.16	128.15	120.31
1	A	190	PRO	N-CA-C	-5.15	108.79	114.92
1	B	315	ASN	N-CA-C	-5.06	107.11	113.28

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	90	0
1	B	3199	0	3157	112	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6419	0	6314	198	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 16.

All (198) 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:B:313:MET:HE3	3:B:606:HOH:O	1.80	0.78
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:201:GLU:C	1:A:205:ASN:O	2.29	0.75
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: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
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:283:ARG:HD3	1:A:343:LYS:HZ2	1.63	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:308:ALA:HB3	1:A:313:MET:HE2	1.86	0.56
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:315:ASN:HA	1:B:318:LYS:CB	2.48	0.44
1:B:275:ASP:C	1:B:277:LEU:N	2.76	0.44
1:B:164:GLU:HG2	1:B:187:ARG:HD2	1.99	0.44
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:179:GLY:O	1:B:321:TYR:OH	2.35	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:B:275:ASP:C	1:B:277:LEU:H	2.29	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
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:114:GLY:O	1:B:115:ILE:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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%)	12	35
1	B	381/426 (89%)	351 (92%)	22 (6%)	8 (2%)	5	17
All	All	762/852 (89%)	704 (92%)	46 (6%)	12 (2%)	7	23

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%)	339 (96%)	15 (4%)	26	58
1	B	354/382 (93%)	340 (96%)	14 (4%)	28	60
All	All	708/764 (93%)	679 (96%)	29 (4%)	27	59

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

Mol	Chain	Res	Type
1	A	70	VAL
1	A	79	SER
1	A	146	SER
1	A	149	LEU
1	A	205	ASN
1	A	208	LEU
1	A	243	PHE
1	A	272	LYS
1	A	295	SER
1	A	323	LEU
1	A	324	SER
1	A	325	GLU
1	A	326	ASP
1	A	364	LYS
1	A	368	ASN
1	B	31	GLN
1	B	107	LEU
1	B	127	LEU
1	B	139	GLN
1	B	146	SER
1	B	172	VAL
1	B	208	LEU
1	B	259	SER
1	B	315	ASN
1	B	327	THR
1	B	368	ASN
1	B	369	PHE

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Mol	Chain	Res	Type
1	B	371	LYS
1	B	410	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	9	ASN
1	A	205	ASN
1	A	227	HIS
1	B	53	ASN
1	B	368	ASN
1	B	394	ASN
1	B	418	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 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 [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	393/426 (92%)	-0.73	1 (0%) 90 88	23, 44, 78, 107	0
1	B	393/426 (92%)	-0.62	2 (0%) 87 83	26, 54, 84, 104	0
All	All	786/852 (92%)	-0.68	3 (0%) 88 85	23, 49, 81, 107	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	GLN	5.3
1	B	140	PRO	2.5
1	A	324	SER	2.2

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
2	ZN	A	500	1/1	0.91	0.27	162,162,162,162	0
2	ZN	B	500	1/1	0.95	0.25	197,197,197,197	0

6.5 Other polymers [i](#)

There are no such residues in this entry.