



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 9, 2026 – 02:26 am BST

PDB ID : 9QBK / pdb_00009qbk
Title : Bacillus Botulinum neurotoxin-like Protein 1 (BBP1)
Authors : Scaletti Hutchinson, E.; Krc, A.; Walse, E.; Stenmark, P.
Deposited on : 2025-03-03
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

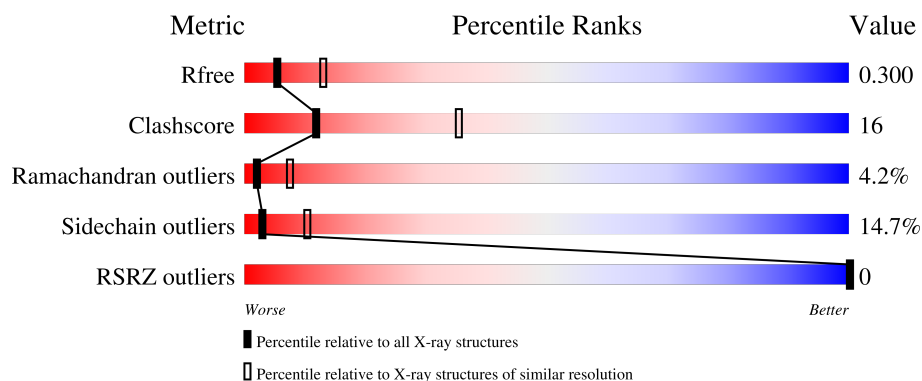
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	401	
1	B	401	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11603 atoms, of which 5602 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/Bt.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	390	Total	C	H	N	O	S	215	0	0
			5785	1916	2796	497	568	8			
1	B	390	Total	C	H	N	O	S	208	0	0
			5804	1919	2806	500	570	9			

- Molecule 2 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ni	0	0
			1	1		

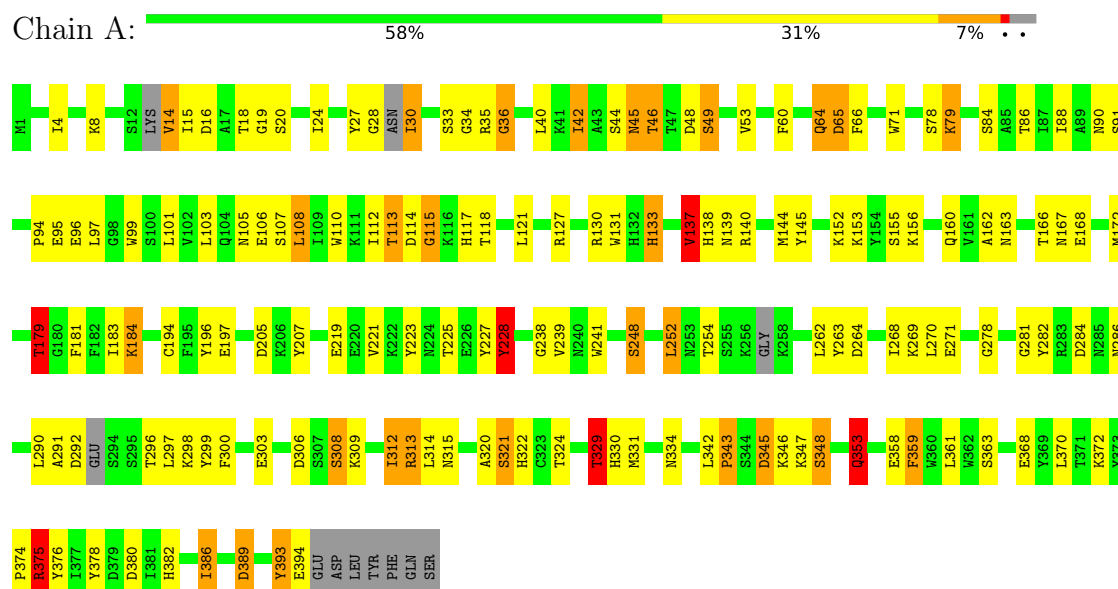
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	O	0	0
			4	4		
3	B	9	Total	O	0	0
			9	9		

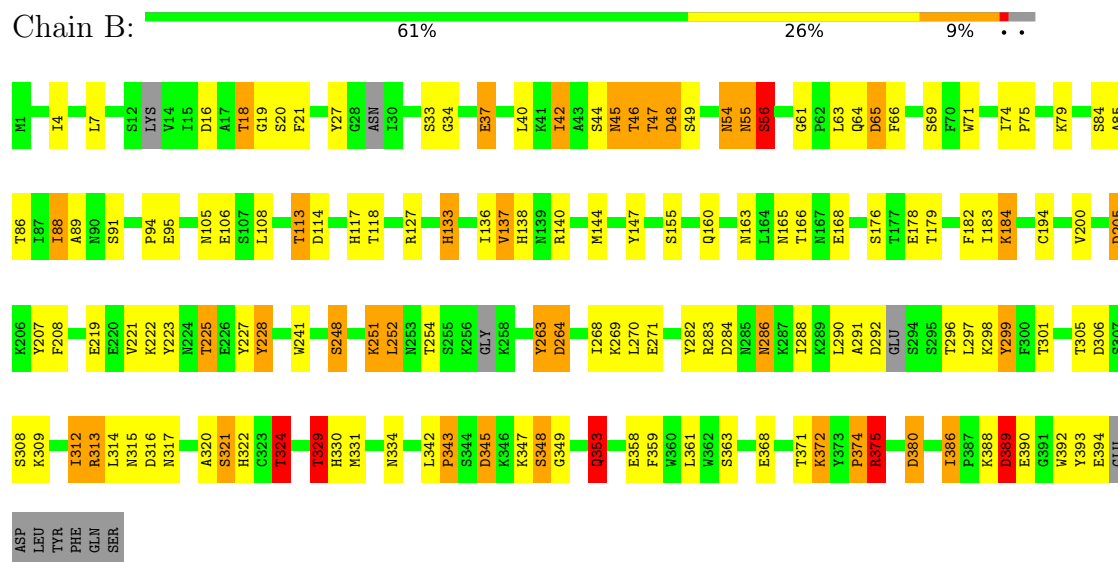
3 Residue-property plots

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/Bt



• Molecule 1: BoNT/Bt



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	243.74Å 51.45Å 104.53Å 90.00° 115.39° 90.00°	Depositor
Resolution (Å)	37.93 – 2.80 37.93 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.6 (37.93-2.80) 98.7 (37.93-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.240 , 0.301 0.241 , 0.300	Depositor DCC
R_{free} test set	1429 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	64.8	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 71.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.469 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11603	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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: NI

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	0.61	0/3065	1.28	20/4178 (0.5%)
1	B	0.60	0/3074	1.28	24/4186 (0.6%)
All	All	0.61	0/6139	1.28	44/8364 (0.5%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	329	THR	CA-CB-OG1	-12.73	90.51	109.60
1	A	329	THR	CA-CB-OG1	-11.91	91.73	109.60
1	A	65	ASP	CB-CA-C	8.78	124.53	109.51
1	B	353	GLN	N-CA-CB	8.65	123.44	110.29
1	B	86	THR	CA-CB-OG1	-8.37	97.05	109.60
1	A	114	ASP	CA-CB-CG	8.33	120.93	112.60
1	A	353	GLN	CB-CA-C	-7.70	95.29	109.46
1	A	86	THR	CA-CB-OG1	-7.57	98.24	109.60
1	B	324	THR	CA-CB-OG1	-7.40	98.50	109.60
1	B	65	ASP	CB-CA-C	7.39	123.06	109.46
1	A	343	PRO	N-CA-C	-7.29	97.44	112.47
1	B	228	TYR	N-CA-CB	-7.19	99.34	110.56
1	B	353	GLN	CB-CA-C	-7.11	96.39	109.46
1	A	353	GLN	N-CA-CB	7.10	121.08	110.29
1	B	179	THR	CA-CB-OG1	-7.05	99.02	109.60
1	B	114	ASP	CA-CB-CG	6.89	119.49	112.60
1	A	324	THR	CA-CB-OG1	-6.51	99.83	109.60
1	A	118	THR	CA-CB-OG1	-6.50	99.85	109.60
1	B	375	ARG	N-CA-CB	6.13	120.85	110.49
1	A	133	HIS	CA-CB-CG	-6.08	107.72	113.80
1	B	343	PRO	N-CA-C	-6.07	99.97	112.47
1	B	374	PRO	N-CA-CB	-6.01	96.94	103.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	359	PHE	CB-CA-C	-6.00	103.83	111.86
1	A	137	VAL	N-CA-CB	-5.89	104.31	111.21
1	B	118	THR	CA-CB-OG1	-5.85	100.82	109.60
1	A	228	TYR	N-CA-CB	-5.83	101.76	111.20
1	B	345	ASP	CA-CB-CG	5.81	118.41	112.60
1	A	359	PHE	CB-CA-C	-5.80	103.74	111.63
1	A	108	LEU	N-CA-CB	-5.71	101.74	110.65
1	B	380	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	152	LYS	N-CA-CB	5.66	118.45	109.51
1	B	228	TYR	CB-CA-C	5.56	118.97	109.80
1	A	300	PHE	CA-CB-CG	5.54	119.34	113.80
1	A	375	ARG	N-CA-CB	5.52	119.81	110.49
1	B	329	THR	OG1-CB-CG2	5.35	120.00	109.30
1	A	179	THR	CA-CB-OG1	-5.34	101.59	109.60
1	B	133	HIS	CA-CB-CG	-5.31	108.49	113.80
1	B	205	ASP	CA-CB-CG	5.29	117.89	112.60
1	B	184	LYS	CB-CA-C	5.21	118.76	109.38
1	B	389	ASP	CA-CB-CG	5.19	117.79	112.60
1	A	358	GLU	CB-CA-C	5.18	118.72	109.29
1	B	264	ASP	CA-CB-CG	5.17	117.77	112.60
1	A	184	LYS	CB-CA-C	5.09	118.54	109.38
1	B	225	THR	CA-CB-OG1	-5.01	102.08	109.60

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	2989	2796	2658	90	0
1	B	2998	2806	2676	90	1
2	A	1	0	0	0	1
3	A	4	0	0	0	0
3	B	9	0	0	1	0
All	All	6001	5602	5334	179	1

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 (179) 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:228:TYR:OH	1:B:264:ASP:O	1.71	1.06
1:B:228:TYR:OH	1:B:264:ASP:N	2.09	0.85
1:A:228:TYR:OH	1:A:264:ASP:O	1.97	0.83
1:A:228:TYR:OH	1:A:264:ASP:N	2.12	0.82
1:A:343:PRO:O	1:A:348:SER:HA	1.79	0.82
1:B:27:TYR:O	1:B:48:ASP:HA	1.82	0.80
1:A:322:HIS:O	1:A:342:LEU:O	1.99	0.79
1:B:329:THR:O	1:B:331:MET:N	2.17	0.77
1:A:329:THR:O	1:A:331:MET:N	2.19	0.75
1:B:228:TYR:HH	1:B:264:ASP:C	1.91	0.74
1:B:342:LEU:HB3	1:B:343:PRO:HD3	1.70	0.73
1:A:368:GLU:OE2	1:A:374:PRO:O	2.07	0.73
1:B:251:LYS:NZ	1:B:358:GLU:O	2.23	0.72
1:A:40:LEU:O	1:A:184:LYS:HA	1.90	0.71
1:B:228:TYR:CZ	1:B:264:ASP:O	2.44	0.70
1:B:42:ILE:HD11	1:B:49:SER:HA	1.74	0.70
1:B:368:GLU:OE2	1:B:374:PRO:O	2.09	0.70
1:B:65:ASP:HB3	1:B:140:ARG:HG2	1.74	0.69
1:B:343:PRO:O	1:B:348:SER:HA	1.93	0.69
1:B:106:GLU:OE1	1:B:127:ARG:HD3	1.93	0.69
1:B:228:TYR:OH	1:B:264:ASP:C	2.36	0.68
1:B:322:HIS:O	1:B:342:LEU:O	2.11	0.68
1:A:342:LEU:HB3	1:A:343:PRO:HD3	1.76	0.67
1:B:65:ASP:HB2	1:B:138:HIS:O	1.94	0.67
1:B:297:LEU:HA	1:B:315:ASN:O	1.94	0.67
1:A:45:ASN:O	1:A:46:THR:O	2.13	0.67
1:B:353:GLN:OE1	1:B:380:ASP:HB2	1.97	0.65
1:A:78:SER:OG	1:A:79:LYS:HD3	1.96	0.64
1:B:221:VAL:HA	1:B:227:TYR:OH	1.97	0.64
1:A:353:GLN:OE1	1:A:380:ASP:HB2	1.99	0.62
1:B:284:ASP:HB3	3:B:502:HOH:O	1.99	0.61
1:A:228:TYR:CZ	1:A:264:ASP:O	2.53	0.61
1:A:221:VAL:HA	1:A:227:TYR:OH	2.01	0.61
1:B:16:ASP:HB3	1:B:21:PHE:CB	2.32	0.59
1:A:42:ILE:HG22	1:A:183:ILE:HG22	1.85	0.59
1:B:320:ALA:O	1:B:321:SER:CB	2.51	0.58
1:A:96:GLU:HG2	1:A:115:GLY:HA2	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:PRO:O	1:A:95:GLU:C	2.47	0.57
1:B:166:THR:HG22	1:B:168:GLU:H	1.70	0.57
1:B:61:GLY:HA2	1:B:163:ASN:HD22	1.69	0.57
1:B:228:TYR:CE2	1:B:263:TYR:HA	2.40	0.57
1:A:374:PRO:O	1:A:375:ARG:HB2	2.05	0.57
1:A:376:TYR:OH	1:B:380:ASP:OD1	2.16	0.57
1:A:27:TYR:O	1:A:48:ASP:HA	2.05	0.57
1:B:40:LEU:O	1:B:184:LYS:HA	2.05	0.57
1:B:241:TRP:CE3	1:B:248:SER:HB3	2.40	0.56
1:B:371:THR:C	1:B:372:LYS:HE3	2.31	0.56
1:A:297:LEU:HA	1:A:315:ASN:O	2.05	0.56
1:A:166:THR:O	1:A:168:GLU:N	2.39	0.55
1:B:305:THR:HB	1:B:347:LYS:CB	2.35	0.55
1:B:284:ASP:O	1:B:324:THR:OG1	2.16	0.55
1:A:320:ALA:O	1:A:321:SER:CB	2.55	0.55
1:A:110:TRP:HE1	1:A:138:HIS:HE1	1.53	0.55
1:B:306:ASP:O	1:B:308:SER:N	2.40	0.55
1:B:85:ALA:HB2	1:B:176:SER:HB2	1.89	0.54
1:A:312:ILE:CD1	1:A:342:LEU:HD22	2.38	0.54
1:B:223:TYR:C	1:B:225:THR:H	2.13	0.54
1:B:65:ASP:CB	1:B:140:ARG:HG2	2.38	0.53
1:B:297:LEU:O	1:B:298:LYS:C	2.51	0.53
1:B:268:ILE:HD11	1:B:288:ILE:CG2	2.39	0.53
1:A:35:ARG:O	1:A:36:GLY:C	2.52	0.53
1:A:90:ASN:O	1:A:166:THR:OG1	2.26	0.53
1:A:65:ASP:HB2	1:A:138:HIS:O	2.08	0.53
1:A:144:MET:O	1:A:155:SER:HA	2.09	0.53
1:B:46:THR:HB	1:B:48:ASP:OD1	2.09	0.53
1:B:223:TYR:HE1	1:B:286:ASN:HB3	1.75	0.52
1:B:228:TYR:OH	1:B:264:ASP:CA	2.57	0.52
1:A:65:ASP:HB3	1:A:140:ARG:HG2	1.92	0.51
1:B:74:ILE:HD11	1:B:108:LEU:HD13	1.91	0.51
1:B:54:ASN:O	1:B:56:SER:N	2.43	0.51
1:A:238:GLY:HA3	1:A:264:ASP:O	2.11	0.51
1:A:16:ASP:OD2	1:A:20:SER:N	2.42	0.51
1:A:28:GLY:O	1:A:30:ILE:HG22	2.12	0.50
1:A:42:ILE:HD11	1:A:49:SER:HA	1.93	0.50
1:B:374:PRO:O	1:B:375:ARG:CB	2.56	0.49
1:B:144:MET:O	1:B:155:SER:HA	2.12	0.49
1:A:99:TRP:HB3	1:A:112:ILE:HG22	1.94	0.49
1:A:228:TYR:HH	1:A:264:ASP:C	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:PHE:CZ	1:B:334:ASN:HB3	2.47	0.49
1:A:66:PHE:CE2	1:A:138:HIS:CD2	3.01	0.49
1:A:374:PRO:O	1:A:375:ARG:CB	2.57	0.49
1:B:147:TYR:CD2	1:B:200:VAL:HG11	2.48	0.49
1:B:374:PRO:O	1:B:375:ARG:HB2	2.11	0.49
1:A:113:THR:HA	1:A:117:HIS:O	2.13	0.48
1:A:137:VAL:HG13	1:A:194:CYS:SG	2.53	0.48
1:A:271:GLU:OE2	1:A:299:TYR:OH	2.32	0.48
1:B:137:VAL:HG13	1:B:194:CYS:SG	2.53	0.48
1:B:269:LYS:O	1:B:270:LEU:HD23	2.14	0.48
1:B:301:THR:OG1	1:B:315:ASN:HB2	2.14	0.48
1:A:35:ARG:HG2	1:A:131:TRP:CD1	2.49	0.48
1:A:262:LEU:HD13	1:A:386:ILE:HG21	1.96	0.47
1:A:228:TYR:CE2	1:A:264:ASP:O	2.68	0.47
1:B:94:PRO:O	1:B:95:GLU:C	2.56	0.47
1:A:103:LEU:HD12	1:A:107:SER:O	2.14	0.47
1:A:241:TRP:CE3	1:A:248:SER:HB3	2.50	0.47
1:A:269:LYS:HG3	1:A:291:ALA:HB2	1.95	0.47
1:B:61:GLY:N	1:B:64:GLN:HB3	2.29	0.47
1:A:79:LYS:HG2	1:A:181:PHE:CZ	2.50	0.47
1:A:393:TYR:C	1:A:393:TYR:CD1	2.92	0.47
1:B:282:TYR:C	1:B:284:ASP:H	2.22	0.47
1:A:101:LEU:HD11	1:A:108:LEU:HG	1.97	0.47
1:A:378:TYR:O	1:A:382:HIS:ND1	2.48	0.47
1:A:71:TRP:CH2	1:A:133:HIS:HB2	2.50	0.46
1:A:297:LEU:O	1:A:298:LYS:C	2.57	0.46
1:B:42:ILE:HG22	1:B:183:ILE:HG22	1.96	0.46
1:A:45:ASN:HD22	1:A:45:ASN:HA	1.62	0.46
1:A:139:ASN:OD1	1:A:139:ASN:C	2.58	0.46
1:B:75:PRO:HG3	1:B:182:PHE:CE1	2.51	0.46
1:B:91:SER:HA	1:B:166:THR:HG21	1.97	0.46
1:B:34:GLY:O	1:B:219:GLU:OE1	2.33	0.46
1:B:91:SER:CB	1:B:166:THR:HG21	2.46	0.46
1:B:313:ARG:NH2	1:B:314:LEU:O	2.49	0.46
1:A:91:SER:HB3	1:A:166:THR:HG21	1.98	0.46
1:B:291:ALA:C	1:B:292:ASP:OD1	2.58	0.46
1:A:228:TYR:CD1	1:A:228:TYR:C	2.94	0.46
1:B:163:ASN:HD21	1:B:165:ASN:HB2	1.81	0.46
1:B:20:SER:O	1:B:55:ASN:ND2	2.50	0.45
1:B:47:THR:O	1:B:48:ASP:C	2.60	0.45
1:B:252:LEU:HD13	1:B:361:LEU:HD21	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:VAL:O	1:B:144:MET:HA	2.15	0.45
1:A:106:GLU:OE1	1:A:127:ARG:HD3	2.16	0.45
1:A:121:LEU:HG	1:A:144:MET:HE3	1.99	0.45
1:A:223:TYR:C	1:A:225:THR:N	2.74	0.45
1:B:113:THR:HA	1:B:117:HIS:O	2.16	0.45
1:B:312:ILE:CD1	1:B:342:LEU:HD22	2.46	0.45
1:B:271:GLU:OE2	1:B:299:TYR:OH	2.36	0.44
1:A:45:ASN:O	1:A:46:THR:C	2.61	0.44
1:A:313:ARG:NH2	1:A:314:LEU:O	2.50	0.44
1:B:66:PHE:CZ	1:B:138:HIS:CD2	3.06	0.44
1:A:14:VAL:HB	1:A:24:ILE:HG21	1.99	0.44
1:A:65:ASP:CB	1:A:140:ARG:HG2	2.47	0.44
1:A:137:VAL:O	1:A:144:MET:HA	2.18	0.44
1:A:166:THR:C	1:A:168:GLU:H	2.24	0.44
1:A:42:ILE:HD11	1:A:172:MET:O	2.18	0.44
1:A:306:ASP:O	1:A:308:SER:N	2.51	0.43
1:A:346:LYS:O	1:A:347:LYS:C	2.61	0.43
1:A:282:TYR:C	1:A:284:ASP:H	2.25	0.43
1:A:34:GLY:O	1:A:219:GLU:OE1	2.37	0.43
1:A:228:TYR:CZ	1:A:264:ASP:N	2.84	0.43
1:B:91:SER:CA	1:B:166:THR:HG21	2.49	0.43
1:B:393:TYR:CD1	1:B:393:TYR:C	2.96	0.43
1:A:105:ASN:O	1:A:106:GLU:HB2	2.17	0.43
1:B:223:TYR:C	1:B:225:THR:N	2.77	0.43
1:B:205:ASP:C	1:B:207:TYR:H	2.26	0.43
1:B:317:ASN:O	1:B:320:ALA:HB2	2.18	0.43
1:A:252:LEU:HD13	1:A:361:LEU:HD21	2.01	0.42
1:A:269:LYS:O	1:A:270:LEU:HD23	2.18	0.42
1:B:320:ALA:O	1:B:321:SER:HB2	2.18	0.42
1:B:316:ASP:O	1:B:317:ASN:C	2.62	0.42
1:B:88:ILE:HD12	1:B:89:ALA:N	2.35	0.42
1:B:228:TYR:CZ	1:B:263:TYR:HA	2.54	0.42
1:B:71:TRP:CH2	1:B:133:HIS:HB2	2.55	0.42
1:B:45:ASN:HB2	1:B:178:GLU:O	2.19	0.42
1:B:228:TYR:HB3	1:B:386:ILE:O	2.20	0.42
1:A:45:ASN:OD1	1:A:179:THR:HG22	2.20	0.41
1:B:389:ASP:HB3	1:B:392:TRP:HB3	2.01	0.41
1:A:96:GLU:HA	1:A:113:THR:HG21	2.02	0.41
1:A:281:GLY:O	1:A:284:ASP:HB2	2.20	0.41
1:A:298:LYS:O	1:A:299:TYR:CD2	2.73	0.41
1:B:7:LEU:HD23	1:B:7:LEU:HA	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:GLN:O	1:B:64:GLN:HG3	2.20	0.41
1:A:60:PHE:HA	1:A:64:GLN:HG2	2.03	0.41
1:A:223:TYR:C	1:A:225:THR:H	2.27	0.41
1:B:389:ASP:OD1	1:B:389:ASP:N	2.53	0.41
1:A:322:HIS:HB2	1:A:345:ASP:OD2	2.20	0.41
1:B:66:PHE:CE2	1:B:138:HIS:CD2	3.09	0.41
1:A:299:TYR:O	1:A:314:LEU:HA	2.21	0.41
1:B:329:THR:C	1:B:331:MET:N	2.79	0.41
1:A:205:ASP:C	1:A:207:TYR:H	2.29	0.41
1:A:228:TYR:OH	1:A:264:ASP:CA	2.68	0.41
1:A:145:TYR:CD1	1:A:155:SER:HB3	2.56	0.40
1:B:105:ASN:O	1:B:106:GLU:HB2	2.21	0.40
1:A:166:THR:C	1:A:168:GLU:N	2.78	0.40
1:A:196:TYR:O	1:A:197:GLU:C	2.63	0.40
1:A:252:LEU:HB2	1:A:359:PHE:HB3	2.04	0.40
1:B:343:PRO:HD2	1:B:349:GLY:H	1.86	0.40
1:A:228:TYR:OH	1:A:264:ASP:C	2.62	0.40
1:B:271:GLU:CD	1:B:299:TYR:OH	2.65	0.40
1:A:162:ALA:O	1:A:163:ASN:C	2.64	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:HIS:HE1	2:A:501:NI:NI[2_565]	1.55	0.05

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	380/401 (95%)	311 (82%)	52 (14%)	17 (4%)	2 7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	380/401 (95%)	319 (84%)	46 (12%)	15 (4%)	2	8
All	All	760/802 (95%)	630 (83%)	98 (13%)	32 (4%)	2	7

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	SER
1	A	46	THR
1	A	321	SER
1	A	329	THR
1	B	33	SER
1	B	321	SER
1	B	329	THR
1	A	19	GLY
1	A	36	GLY
1	A	84	SER
1	A	97	LEU
1	A	389	ASP
1	B	19	GLY
1	B	37	GLU
1	B	56	SER
1	B	348	SER
1	B	375	ARG
1	A	296	THR
1	A	330	HIS
1	A	348	SER
1	B	84	SER
1	B	296	THR
1	B	299	TYR
1	A	375	ARG
1	B	55	ASN
1	B	330	HIS
1	A	18	THR
1	A	167	ASN
1	B	18	THR
1	B	283	ARG
1	A	115	GLY
1	A	278	GLY

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	288/366 (79%)	243 (84%)	45 (16%)	2	9
1	B	291/366 (80%)	251 (86%)	40 (14%)	3	13
All	All	579/732 (79%)	494 (85%)	85 (15%)	3	10

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	8	LYS
1	A	14	VAL
1	A	15	ILE
1	A	30	ILE
1	A	42	ILE
1	A	44	SER
1	A	45	ASN
1	A	49	SER
1	A	53	VAL
1	A	64	GLN
1	A	79	LYS
1	A	88	ILE
1	A	113	THR
1	A	130	ARG
1	A	137	VAL
1	A	153	LYS
1	A	156	LYS
1	A	160	GLN
1	A	179	THR
1	A	228	TYR
1	A	239	VAL
1	A	248	SER
1	A	252	LEU
1	A	254	THR
1	A	263	TYR
1	A	268	ILE

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Mol	Chain	Res	Type
1	A	286	ASN
1	A	290	LEU
1	A	292	ASP
1	A	303	GLU
1	A	308	SER
1	A	309	LYS
1	A	312	ILE
1	A	313	ARG
1	A	334	ASN
1	A	345	ASP
1	A	353	GLN
1	A	363	SER
1	A	370	LEU
1	A	372	LYS
1	A	386	ILE
1	A	389	ASP
1	A	393	TYR
1	A	394	GLU
1	B	4	ILE
1	B	18	THR
1	B	37	GLU
1	B	42	ILE
1	B	44	SER
1	B	45	ASN
1	B	46	THR
1	B	47	THR
1	B	48	ASP
1	B	54	ASN
1	B	56	SER
1	B	63	LEU
1	B	69	SER
1	B	79	LYS
1	B	88	ILE
1	B	113	THR
1	B	136	ILE
1	B	137	VAL
1	B	160	GLN
1	B	222	LYS
1	B	248	SER
1	B	251	LYS
1	B	252	LEU
1	B	254	THR

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Mol	Chain	Res	Type
1	B	263	TYR
1	B	286	ASN
1	B	290	LEU
1	B	309	LYS
1	B	312	ILE
1	B	313	ARG
1	B	324	THR
1	B	345	ASP
1	B	353	GLN
1	B	363	SER
1	B	372	LYS
1	B	386	ILE
1	B	388	LYS
1	B	389	ASP
1	B	390	GLU
1	B	394	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	64	GLN
1	A	117	HIS
1	A	138	HIS
1	A	237	GLN
1	A	330	HIS
1	B	55	ASN
1	B	64	GLN
1	B	138	HIS
1	B	163	ASN
1	B	237	GLN

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 [i](#)

There are no oligosaccharides 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 [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	390/401 (97%)	-0.96	0 100 100	28, 63, 104, 133	0
1	B	390/401 (97%)	-0.99	0 100 100	29, 64, 106, 136	0
All	All	780/802 (97%)	-0.97	0 100 100	28, 64, 106, 136	0

There are no RSRZ outliers to report.

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	NI	A	501	1/1	0.99	0.02	64,64,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.