



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:22 PM UTC

PDB ID : 1ZKW / pdb_00001zkw
Title : Crystal structure of Arg347Ala mutant of botulinum neurotoxin E catalytic domain
Authors : Agarwal, R.; Binz, T.; Swaminathan, S.
Deposited on : 2005-05-04
Resolution : 2.17 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

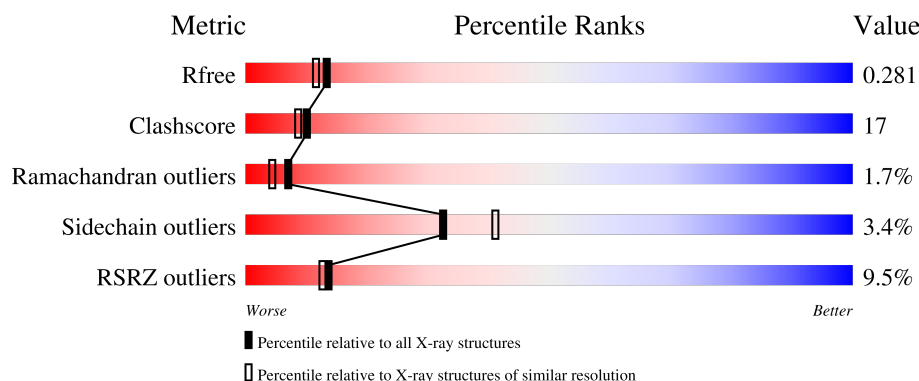
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	420	5% 70% 25% ..
1	B	420	13% 65% 29% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6841 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 botulinum neurotoxin type E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3293	2106	549	630	8			
1	B	409	Total	C	N	O	S	0	0	0
			3276	2094	547	628	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	347	ALA	ARG	engineered mutation	UNP Q00496
B	347	ALA	ARG	engineered mutation	UNP Q00496

- 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 CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	Cl	0	0
			6	6		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	165	Total	O	0	0
			165	165		

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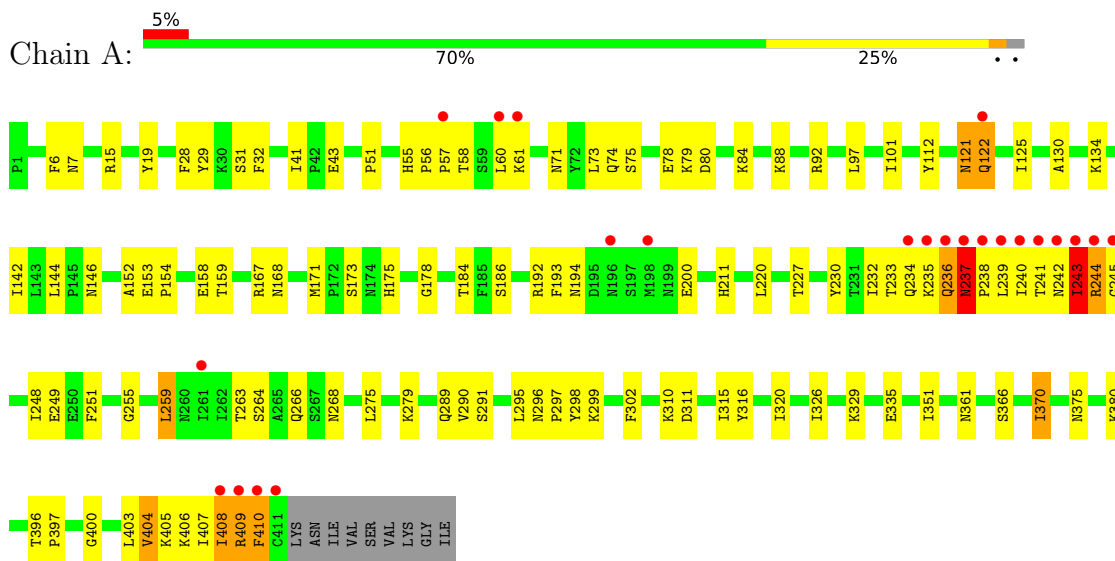
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	98	Total	O	0	0
			98	98		

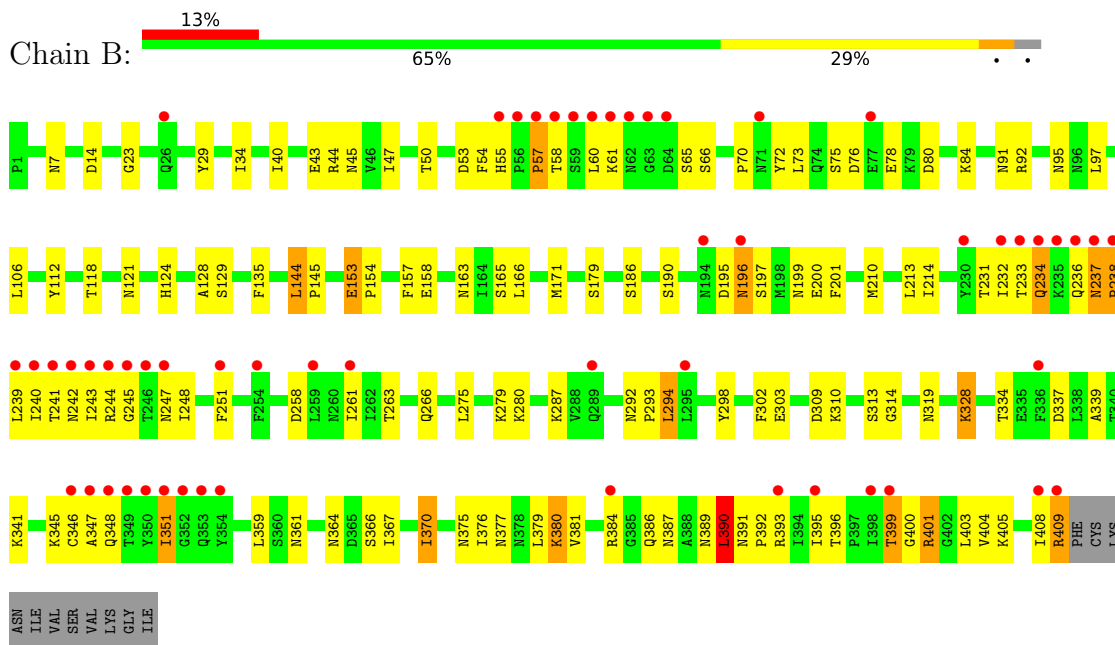
3 Residue-property plots

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: botulinum neurotoxin type E



• Molecule 1: botulinum neurotoxin type E



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	88.58Å 144.46Å 82.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.35 – 2.17 42.35 – 2.17	Depositor EDS
% Data completeness (in resolution range)	87.4 (42.35-2.17) 87.4 (42.35-2.17)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.77 (at 1.94Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.238 , 0.280 0.238 , 0.281	Depositor DCC
R_{free} test set	721 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	24.7	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6841	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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: CL, 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	0.43	0/3367	0.95	12/4562 (0.3%)
1	B	0.40	0/3349	0.93	11/4538 (0.2%)
All	All	0.41	0/6716	0.94	23/9100 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	197	SER	N-CA-C	8.04	120.55	109.54
1	B	399	THR	N-CA-C	8.02	122.73	113.19
1	A	320	ILE	N-CA-C	6.76	117.51	110.62
1	B	7	ASN	N-CA-C	-6.45	99.20	109.59
1	B	23	GLY	N-CA-C	-6.38	103.91	112.14
1	A	370	ILE	N-CA-C	6.30	116.98	110.36
1	A	237	ASN	CA-C-N	-6.19	112.53	118.97
1	A	237	ASN	C-N-CA	-6.19	112.53	118.97
1	A	366	SER	N-CA-C	-6.11	105.32	112.89
1	A	7	ASN	N-CA-C	-5.96	99.99	109.59
1	A	57	PRO	N-CA-C	-5.94	106.12	113.84
1	A	158	GLU	N-CA-C	5.91	119.02	110.28
1	B	401	ARG	N-CA-C	-5.86	104.97	111.36
1	B	303	GLU	N-CA-C	-5.75	104.93	111.14
1	A	122	GLN	N-CA-C	5.72	117.77	109.07
1	B	186	SER	CA-C-N	5.69	125.81	119.32
1	B	186	SER	C-N-CA	5.69	125.81	119.32
1	B	377	ASN	CB-CA-C	-5.59	110.11	116.54
1	B	370	ILE	N-CA-C	5.55	115.75	110.53
1	A	152	ALA	N-CA-C	5.52	117.85	110.35
1	B	158	GLU	N-CA-C	5.26	118.07	110.28
1	A	186	SER	CA-C-N	5.12	124.91	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	SER	C-N-CA	5.12	124.91	119.28

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	3293	0	3252	102	0
1	B	3276	0	3238	129	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	6	0	0	0	0
3	B	1	0	0	0	0
4	A	165	0	0	6	0
4	B	98	0	0	7	0
All	All	6841	0	6490	226	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 17.

All (226) 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:75:SER:HB3	1:B:78:GLU:HG3	1.54	0.89
1:B:351:ILE:HD13	1:B:351:ILE:H	1.41	0.85
1:B:196:ASN:HA	4:B:915:HOH:O	1.77	0.84
1:B:263:THR:H	1:B:266:GLN:HE21	1.21	0.83
1:B:345:LYS:H	1:B:386:GLN:HE22	1.27	0.83
1:A:238:PRO:C	1:A:240:ILE:H	1.86	0.82
1:B:263:THR:H	1:B:266:GLN:NE2	1.79	0.79
1:A:243:ILE:HA	4:A:686:HOH:O	1.82	0.78
1:A:234:GLN:NE2	1:A:243:ILE:H	1.83	0.75
1:B:234:GLN:HG3	1:B:243:ILE:HG12	1.68	0.74
1:B:381:VAL:O	1:B:384:ARG:HG3	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:GLN:HG3	1:A:291:SER:H	1.52	0.73
1:A:238:PRO:O	1:A:240:ILE:HG12	1.88	0.72
1:B:65:SER:HB2	4:B:904:HOH:O	1.88	0.71
1:B:54:PHE:O	1:B:70:PRO:HG3	1.91	0.71
1:A:234:GLN:HA	1:A:238:PRO:HG2	1.73	0.70
1:A:295:LEU:HD13	1:A:299:LYS:HE3	1.73	0.70
1:B:310:LYS:HE2	1:B:314:GLY:HA2	1.73	0.70
1:B:263:THR:OG1	1:B:266:GLN:HG3	1.91	0.70
1:A:409:ARG:HE	1:A:409:ARG:HA	1.57	0.69
1:B:233:THR:HA	1:B:245:GLY:HA2	1.75	0.69
1:B:334:THR:OG1	1:B:337:ASP:HB3	1.92	0.68
1:B:389:ASN:O	1:B:390:LEU:HB2	1.93	0.68
1:B:92:ARG:HA	1:B:370:ILE:HG23	1.76	0.67
1:B:231:THR:HG22	1:B:247:ASN:OD1	1.95	0.67
1:B:233:THR:HG22	1:B:234:GLN:N	2.08	0.67
1:B:55:HIS:ND1	1:B:70:PRO:HG2	2.10	0.65
1:A:238:PRO:HG2	1:A:239:LEU:H	1.60	0.65
1:B:351:ILE:HD13	1:B:351:ILE:N	2.10	0.64
1:B:351:ILE:H	1:B:351:ILE:CD1	2.11	0.64
1:A:405:LYS:O	1:A:408:ILE:HG12	1.97	0.64
1:A:130:ALA:HB2	1:A:144:LEU:HD23	1.81	0.63
1:B:118:THR:HG21	1:B:124:HIS:CD2	2.33	0.63
1:A:243:ILE:C	1:A:245:GLY:H	2.07	0.63
1:B:50:THR:HG22	1:B:53:ASP:CG	2.24	0.62
1:B:345:LYS:H	1:B:386:GLN:NE2	1.96	0.62
1:A:407:ILE:O	1:A:409:ARG:N	2.30	0.62
1:B:248:ILE:HA	1:B:251:PHE:HD1	1.65	0.62
1:A:125:ILE:HD11	1:B:294:LEU:CD1	2.29	0.62
1:A:232:ILE:HG23	1:A:236:GLN:HB2	1.82	0.62
1:A:130:ALA:HB2	1:A:144:LEU:CD2	2.30	0.62
1:A:406:LYS:O	1:A:410:PHE:HB2	1.99	0.62
1:B:57:PRO:HG2	1:B:58:THR:H	1.64	0.61
1:B:153:GLU:HB2	1:B:154:PRO:CD	2.30	0.61
1:B:50:THR:HG22	1:B:53:ASP:OD2	2.00	0.61
1:B:234:GLN:HE22	1:B:239:LEU:C	2.08	0.61
1:B:44:ARG:NH1	1:B:72:TYR:HB3	2.16	0.61
1:B:375:ASN:HB3	1:B:380:LYS:HA	1.83	0.60
1:A:153:GLU:HB2	1:A:154:PRO:CD	2.31	0.60
1:B:233:THR:HB	4:B:901:HOH:O	2.01	0.60
1:B:234:GLN:HA	1:B:241:THR:HB	1.82	0.60
1:A:71:ASN:O	1:A:74:GLN:HG2	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:ASP:OD1	1:B:341:LYS:HE2	2.03	0.59
1:A:410:PHE:N	4:A:689:HOH:O	2.35	0.59
1:B:80:ASP:OD2	1:B:84:LYS:HE2	2.02	0.59
1:A:121:ASN:C	1:A:121:ASN:HD22	2.10	0.59
1:A:238:PRO:C	1:A:240:ILE:N	2.56	0.59
1:B:58:THR:HG21	1:B:61:LYS:HD3	1.84	0.59
1:B:128:ALA:HB2	1:B:166:LEU:HD11	1.84	0.58
1:A:289:GLN:NE2	1:A:290:VAL:HG22	2.18	0.58
1:A:326:ILE:HA	1:A:329:LYS:HD3	1.85	0.58
1:A:233:THR:OG1	1:A:236:GLN:HG3	2.03	0.58
1:B:233:THR:HA	1:B:245:GLY:CA	2.32	0.58
1:A:255:GLY:HA2	1:A:259:LEU:HD13	1.85	0.58
1:A:236:GLN:C	1:A:238:PRO:CD	2.77	0.58
1:B:195:ASP:CG	1:B:199:ASN:HB2	2.30	0.57
1:A:88:LYS:HG3	1:A:370:ILE:HD11	1.85	0.57
1:A:61:LYS:HZ2	1:A:404:VAL:HG22	1.70	0.56
1:A:409:ARG:HE	1:A:409:ARG:CA	2.18	0.56
1:A:61:LYS:NZ	1:A:404:VAL:HG22	2.21	0.56
1:B:328:LYS:HB3	1:B:328:LYS:NZ	2.21	0.56
1:A:251:PHE:CZ	1:A:351:ILE:HD11	2.41	0.56
1:B:345:LYS:N	1:B:386:GLN:HE22	1.99	0.56
1:B:381:VAL:HB	1:B:384:ARG:NE	2.21	0.56
1:A:396:THR:HG22	1:A:397:PRO:HD2	1.88	0.54
1:A:295:LEU:CD1	1:A:299:LYS:HE3	2.38	0.54
1:B:129:SER:O	1:B:144:LEU:HD13	2.07	0.54
1:A:171:MET:O	1:A:175:HIS:HD2	1.91	0.54
1:A:125:ILE:HD11	1:B:294:LEU:HD11	1.89	0.54
1:A:238:PRO:O	1:A:240:ILE:N	2.41	0.53
1:B:387:ASN:HB3	1:B:390:LEU:HB3	1.89	0.53
1:A:121:ASN:HD22	1:A:122:GLN:N	2.06	0.53
1:B:292:ASN:ND2	1:B:293:PRO:HD2	2.23	0.53
1:B:45:ASN:C	1:B:45:ASN:HD22	2.17	0.53
1:B:201:PHE:CG	1:B:345:LYS:HG2	2.44	0.53
1:B:233:THR:CG2	1:B:234:GLN:N	2.72	0.53
1:B:112:TYR:HA	1:B:302:PHE:CE1	2.44	0.52
1:B:393:ARG:HB2	4:B:912:HOH:O	2.07	0.52
1:B:45:ASN:HD21	1:B:47:ILE:HB	1.74	0.52
1:B:153:GLU:HB2	1:B:154:PRO:HD2	1.90	0.52
1:B:190:SER:HB2	1:B:359:LEU:HD11	1.90	0.52
1:A:232:ILE:HG13	1:A:236:GLN:NE2	2.26	0.51
1:A:243:ILE:O	1:A:244:ARG:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:375:ASN:HD21	1:B:386:GLN:HG3	1.76	0.51
1:A:171:MET:O	1:A:175:HIS:CD2	2.63	0.51
1:B:234:GLN:HG2	1:B:243:ILE:H	1.75	0.51
1:B:237:ASN:HB3	1:B:238:PRO:HD3	1.93	0.51
1:A:153:GLU:HB2	1:A:154:PRO:HD2	1.94	0.50
1:B:237:ASN:CB	1:B:238:PRO:CD	2.89	0.50
1:B:381:VAL:HG21	1:B:384:ARG:NH2	2.25	0.50
1:B:234:GLN:HB2	1:B:243:ILE:HA	1.92	0.50
1:B:45:ASN:ND2	1:B:47:ILE:H	2.10	0.49
1:B:244:ARG:HH11	1:B:244:ARG:HB2	1.77	0.49
1:B:243:ILE:HG22	1:B:243:ILE:O	2.12	0.49
1:B:58:THR:C	1:B:60:LEU:H	2.21	0.49
1:A:243:ILE:O	1:A:245:GLY:N	2.44	0.49
1:B:251:PHE:CD1	1:B:261:ILE:HD11	2.47	0.49
1:A:125:ILE:HD11	1:B:294:LEU:HD13	1.95	0.49
1:B:66:SER:HA	1:B:157:PHE:CD2	2.47	0.49
1:B:408:ILE:O	1:B:409:ARG:HB2	2.13	0.49
1:B:233:THR:HG22	1:B:234:GLN:H	1.74	0.48
1:B:236:GLN:O	1:B:238:PRO:HD2	2.14	0.48
1:A:121:ASN:C	1:A:121:ASN:ND2	2.70	0.48
1:A:43:GLU:OE2	1:A:79:LYS:NZ	2.43	0.48
1:B:95:ASN:N	1:B:95:ASN:HD22	2.10	0.48
1:A:238:PRO:HG2	1:A:239:LEU:N	2.25	0.48
1:A:241:THR:HG22	1:A:242:ASN:N	2.29	0.48
1:B:248:ILE:HA	1:B:251:PHE:CD1	2.45	0.48
1:B:263:THR:N	1:B:266:GLN:HE21	2.00	0.48
1:B:376:ILE:O	1:B:379:LEU:HB2	2.14	0.48
1:A:234:GLN:O	1:A:239:LEU:HD23	2.13	0.47
1:A:234:GLN:NE2	1:A:243:ILE:N	2.58	0.47
1:B:91:ASN:HB3	1:B:370:ILE:HD13	1.95	0.47
1:B:292:ASN:HD22	1:B:293:PRO:CD	2.26	0.47
1:B:339:ALA:CB	1:B:346:CYS:HB2	2.44	0.47
1:A:92:ARG:HA	1:A:370:ILE:HG23	1.96	0.47
1:A:370:ILE:HG12	4:A:585:HOH:O	2.14	0.47
1:A:289:GLN:NE2	1:A:290:VAL:H	2.12	0.47
1:B:165:SER:OG	1:B:171:MET:HE3	2.15	0.47
1:B:75:SER:HB3	1:B:78:GLU:CG	2.36	0.47
1:A:329:LYS:CB	1:A:329:LYS:NZ	2.78	0.47
1:B:121:ASN:HA	1:B:287:LYS:O	2.14	0.47
1:A:263:THR:H	1:A:266:GLN:NE2	2.13	0.46
1:A:329:LYS:HB3	1:A:329:LYS:HZ3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:ASN:HB2	1:A:297:PRO:HD3	1.98	0.46
1:B:233:THR:CG2	1:B:234:GLN:H	2.28	0.46
1:B:234:GLN:N	1:B:241:THR:HG21	2.30	0.46
1:A:167:ARG:HG2	1:A:168:ASN:ND2	2.30	0.46
1:A:264:SER:O	1:A:268:ASN:HB2	2.15	0.46
1:B:92:ARG:CA	1:B:370:ILE:HG23	2.45	0.46
1:A:75:SER:OG	1:A:78:GLU:HG3	2.16	0.46
1:A:243:ILE:C	1:A:245:GLY:N	2.74	0.45
1:A:237:ASN:HD22	1:A:237:ASN:HA	1.57	0.45
1:B:14:ASP:OD1	1:B:135:PHE:HB3	2.16	0.45
1:B:391:ASN:N	1:B:392:PRO:HD3	2.32	0.45
1:A:19:TYR:HA	1:A:29:TYR:O	2.17	0.45
1:B:232:ILE:HD12	1:B:251:PHE:CE1	2.52	0.45
1:A:112:TYR:HA	1:A:302:PHE:CE1	2.52	0.45
1:B:237:ASN:CB	1:B:238:PRO:HD3	2.46	0.45
1:B:346:CYS:SG	1:B:347:ALA:N	2.90	0.45
1:B:34:ILE:HD12	1:B:40:ILE:HD11	1.97	0.45
1:B:392:PRO:HD2	4:B:912:HOH:O	2.15	0.45
1:B:400:GLY:O	1:B:404:VAL:HG23	2.16	0.45
1:A:142:ILE:HG13	1:B:293:PRO:HG3	1.99	0.45
1:A:173:SER:HA	1:A:178:GLY:HA2	1.99	0.45
1:A:192:ARG:HD2	1:A:200:GLU:HB3	1.98	0.45
1:A:230:TYR:N	1:A:230:TYR:CD1	2.84	0.45
1:B:309:ASP:OD2	1:B:319:ASN:HB2	2.17	0.45
1:A:235:LYS:C	1:A:237:ASN:H	2.24	0.44
1:B:393:ARG:HG3	1:B:393:ARG:HH11	1.83	0.44
1:B:337:ASP:O	1:B:341:LYS:HG2	2.18	0.44
1:B:401:ARG:O	1:B:405:LYS:HG2	2.18	0.44
1:A:6:PHE:CD2	1:A:32:PHE:HB3	2.53	0.44
1:B:364:ASN:HB3	1:B:367:ILE:HD12	2.00	0.44
1:B:395:ILE:HG22	1:B:396:THR:N	2.32	0.44
1:B:29:TYR:CG	1:B:43:GLU:HG3	2.53	0.44
1:A:409:ARG:C	4:A:689:HOH:O	2.60	0.43
1:B:409:ARG:HE	1:B:409:ARG:HA	1.83	0.43
1:A:146:ASN:O	1:A:220:LEU:HD21	2.18	0.43
1:A:236:GLN:C	1:A:238:PRO:HD2	2.42	0.43
1:B:54:PHE:C	1:B:70:PRO:HG3	2.42	0.43
1:B:58:THR:HG21	1:B:61:LYS:CD	2.48	0.43
1:B:242:ASN:O	1:B:244:ARG:HG3	2.17	0.43
1:A:375:ASN:HB3	1:A:380:LYS:HA	1.99	0.43
1:B:364:ASN:OD1	1:B:366:SER:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:LEU:HB2	4:A:590:HOH:O	2.17	0.43
1:A:211:HIS:ND1	1:A:335:GLU:OE1	2.52	0.43
1:A:31:SER:OG	1:A:41:ILE:HG12	2.18	0.43
1:A:97:LEU:O	1:A:101:ILE:HG12	2.19	0.43
1:B:275:LEU:HG	1:B:279:LYS:HE3	1.99	0.43
1:A:230:TYR:N	1:A:230:TYR:HD1	2.17	0.43
1:A:408:ILE:C	1:A:409:ARG:HG2	2.43	0.43
1:A:92:ARG:HG3	1:A:92:ARG:HH11	1.84	0.43
1:B:106:LEU:HD11	1:B:213:LEU:HB3	2.01	0.43
1:B:163:ASN:HB3	1:B:179:SER:OG	2.19	0.43
1:A:329:LYS:HG3	4:A:669:HOH:O	2.17	0.43
1:B:232:ILE:CD1	1:B:261:ILE:HD12	2.49	0.43
1:B:258:ASP:O	1:B:261:ILE:HG12	2.19	0.42
1:B:232:ILE:O	1:B:245:GLY:HA2	2.19	0.42
1:B:292:ASN:HD22	1:B:292:ASN:C	2.26	0.42
1:A:275:LEU:O	1:A:279:LYS:HG3	2.19	0.42
1:B:292:ASN:HD22	1:B:293:PRO:N	2.18	0.42
1:A:227:THR:HB	1:A:249:GLU:HB2	2.02	0.42
1:B:45:ASN:HD22	1:B:47:ILE:H	1.66	0.42
1:B:45:ASN:HD21	1:B:47:ILE:CB	2.32	0.42
1:A:58:THR:HG21	1:A:61:LYS:O	2.19	0.42
1:B:118:THR:HG21	1:B:124:HIS:CG	2.55	0.42
1:B:234:GLN:HE22	1:B:239:LEU:CA	2.33	0.42
1:B:45:ASN:C	1:B:45:ASN:ND2	2.78	0.42
1:B:292:ASN:HD22	1:B:293:PRO:HD2	1.82	0.42
1:B:240:ILE:HB	1:B:351:ILE:HG13	2.02	0.41
1:A:15:ARG:NH1	1:B:313:SER:O	2.53	0.41
1:A:19:TYR:HB3	1:A:28:PHE:HB3	2.03	0.41
1:A:55:HIS:HA	1:A:56:PRO:HD3	1.93	0.41
1:A:400:GLY:O	1:A:403:LEU:HB2	2.20	0.41
1:B:380:LYS:HG3	4:B:871:HOH:O	2.20	0.41
1:A:19:TYR:HB2	1:A:134:LYS:HB3	2.02	0.41
1:A:407:ILE:C	1:A:409:ARG:N	2.79	0.41
1:A:405:LYS:O	1:A:408:ILE:CG1	2.67	0.41
1:A:167:ARG:O	1:A:168:ASN:HB2	2.20	0.41
1:A:194:ASN:N	1:A:194:ASN:HD22	2.17	0.41
1:A:232:ILE:CD1	1:A:248:ILE:HG22	2.50	0.41
1:B:44:ARG:HD3	1:B:73:LEU:HB2	2.02	0.41
1:B:58:THR:HG21	1:B:61:LYS:CG	2.51	0.41
1:B:195:ASP:OD1	1:B:195:ASP:C	2.63	0.41
1:B:200:GLU:HG3	4:B:895:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:MET:O	1:B:214:ILE:HG13	2.21	0.41
1:B:80:ASP:CG	1:B:84:LYS:HE2	2.45	0.40
1:A:80:ASP:OD2	1:A:84:LYS:NZ	2.51	0.40
1:A:159:THR:HA	1:A:184:THR:O	2.21	0.40
1:A:237:ASN:N	1:A:238:PRO:HD2	2.36	0.40
1:B:234:GLN:C	1:B:236:GLN:H	2.29	0.40
1:A:153:GLU:CB	1:A:154:PRO:CD	2.97	0.40
1:A:193:PHE:C	1:A:194:ASN:HD22	2.30	0.40
1:A:310:LYS:HD2	1:A:316:TYR:CZ	2.56	0.40
1:A:311:ASP:OD2	1:A:315:ILE:HB	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	409/420 (97%)	384 (94%)	18 (4%)	7 (2%)	7	4
1	B	407/420 (97%)	367 (90%)	33 (8%)	7 (2%)	7	4
All	All	816/840 (97%)	751 (92%)	51 (6%)	14 (2%)	7	4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	237	ASN
1	A	243	ILE
1	B	237	ASN
1	B	390	LEU
1	A	236	GLN
1	B	238	PRO
1	B	348	GLN
1	A	244	ARG

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Mol	Chain	Res	Type
1	A	361	ASN
1	B	361	ASN
1	B	57	PRO
1	A	408	ILE
1	B	153	GLU
1	A	51	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	371/379 (98%)	362 (98%)	9 (2%)	43	55
1	B	369/379 (97%)	353 (96%)	16 (4%)	26	32
All	All	740/758 (98%)	715 (97%)	25 (3%)	32	41

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	LEU
1	A	121	ASN
1	A	237	ASN
1	A	243	ILE
1	A	259	LEU
1	A	298	TYR
1	A	404	VAL
1	A	409	ARG
1	A	410	PHE
1	B	76	ASP
1	B	97	LEU
1	B	144	LEU
1	B	145	PRO
1	B	196	ASN
1	B	234	GLN
1	B	280	LYS
1	B	294	LEU

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Mol	Chain	Res	Type
1	B	298	TYR
1	B	328	LYS
1	B	351	ILE
1	B	380	LYS
1	B	390	LEU
1	B	399	THR
1	B	403	LEU
1	B	409	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	121	ASN
1	A	124	HIS
1	A	168	ASN
1	A	174	ASN
1	A	175	HIS
1	A	194	ASN
1	A	234	GLN
1	A	236	GLN
1	A	237	ASN
1	A	260	ASN
1	A	266	GLN
1	A	289	GLN
1	A	296	ASN
1	A	348	GLN
1	A	375	ASN
1	A	377	ASN
1	A	378	ASN
1	B	45	ASN
1	B	71	ASN
1	B	95	ASN
1	B	121	ASN
1	B	122	GLN
1	B	124	HIS
1	B	174	ASN
1	B	199	ASN
1	B	234	GLN
1	B	236	GLN
1	B	242	ASN
1	B	266	GLN

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Mol	Chain	Res	Type
1	B	268	ASN
1	B	292	ASN
1	B	296	ASN
1	B	353	GLN
1	B	375	ASN
1	B	377	ASN
1	B	386	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 9 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 ⓘ

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	411/420 (97%)	0.26	23 (5%) 30 29	12, 28, 48, 64	0
1	B	409/420 (97%)	0.89	55 (13%) 7 6	20, 36, 63, 71	0
All	All	820/840 (97%)	0.57	78 (9%) 14 13	12, 32, 57, 71	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	60	LEU	6.6
1	A	239	LEU	6.4
1	A	243	ILE	6.1
1	B	240	ILE	6.0
1	B	243	ILE	6.0
1	B	239	LEU	5.6
1	B	351	ILE	5.4
1	A	237	ASN	5.4
1	B	59	SER	5.3
1	B	233	THR	5.1
1	B	238	PRO	5.0
1	A	61	LYS	4.8
1	B	241	THR	4.7
1	A	238	PRO	4.7
1	B	58	THR	4.5
1	B	242	ASN	4.3
1	A	235	LYS	4.2
1	B	409	ARG	4.1
1	B	354	TYR	4.0
1	B	245	GLY	3.9
1	B	237	ASN	3.9
1	A	408	ILE	3.8
1	B	235	LYS	3.8
1	B	63	GLY	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	240	ILE	3.7
1	B	346	CYS	3.7
1	B	234	GLN	3.7
1	B	399	THR	3.6
1	B	348	GLN	3.6
1	B	61	LYS	3.5
1	B	244	ARG	3.5
1	A	244	ARG	3.4
1	B	57	PRO	3.4
1	A	234	GLN	3.3
1	A	411	CYS	3.3
1	B	398	ILE	3.3
1	B	196	ASN	3.3
1	A	409	ARG	3.2
1	B	62	ASN	3.2
1	A	410	PHE	3.2
1	B	352	GLY	3.1
1	A	60	LEU	3.1
1	B	230	TYR	3.0
1	B	393	ARG	3.0
1	B	353	GLN	2.9
1	B	236	GLN	2.9
1	B	64	ASP	2.9
1	A	261	ILE	2.9
1	A	196	ASN	2.9
1	B	349	THR	2.8
1	B	251	PHE	2.8
1	B	347	ALA	2.8
1	A	122	GLN	2.7
1	B	350	TYR	2.7
1	A	57	PRO	2.7
1	B	56	PRO	2.7
1	B	55	HIS	2.7
1	A	236	GLN	2.6
1	A	242	ASN	2.6
1	B	194	ASN	2.6
1	B	232	ILE	2.4
1	B	246	THR	2.4
1	B	336	PHE	2.3
1	B	289	GLN	2.3
1	B	259	LEU	2.3
1	A	241	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	26	GLN	2.2
1	B	395	ILE	2.2
1	A	245	GLY	2.1
1	B	384	ARG	2.1
1	B	295	LEU	2.1
1	B	408	ILE	2.1
1	B	254	PHE	2.1
1	A	198	MET	2.0
1	B	71	ASN	2.0
1	B	261	ILE	2.0
1	B	77	GLU	2.0
1	B	247	ASN	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 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
3	CL	A	580	1/1	0.92	0.13	46,46,46,46	0
3	CL	A	578	1/1	0.93	0.11	52,52,52,52	0
3	CL	A	581	1/1	0.95	0.08	43,43,43,43	0
3	CL	A	576	1/1	0.96	0.15	46,46,46,46	0
3	CL	A	579	1/1	0.96	0.18	47,47,47,47	0
3	CL	B	577	1/1	0.96	0.13	53,53,53,53	0
2	ZN	B	822	1/1	0.97	0.04	49,49,49,49	0
3	CL	A	575	1/1	0.97	0.13	32,32,32,32	0
2	ZN	A	422	1/1	0.99	0.02	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.