



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2024 – 07:37 AM EDT

PDB ID : 1BTG
Title : CRYSTAL STRUCTURE OF BETA NERVE GROWTH FACTOR AT 2.5 Å
RESOLUTION IN C2 SPACE GROUP WITH ZN IONS BOUND
Authors : Holland, D.R.; Matthews, B.W.
Deposited on : 1995-08-29
Resolution : 2.50 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	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

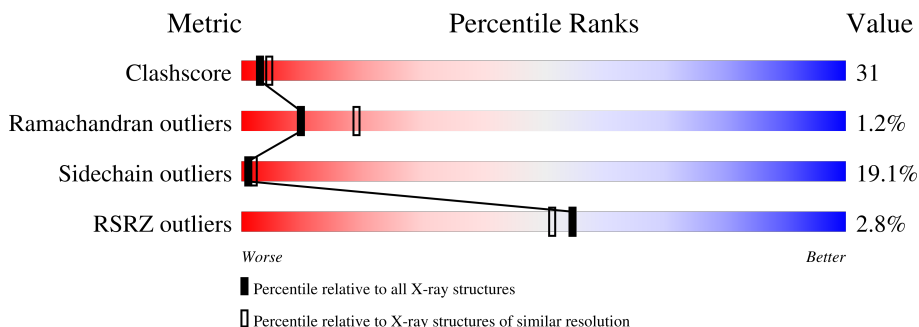
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (2.50-2.50)
RSRZ outliers	164620	5504 (2.50-2.50)

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	110	 46% 35% 16% 5%
1	B	110	 49% 40% 8% 5%
1	C	110	 36% 41% 20% 3%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2565 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 BETA NERVE GROWTH FACTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	108	Total	C	N	O	S	0	0	0
			842	526	147	163	6			
1	B	109	Total	C	N	O	S	0	0	0
			832	521	147	158	6			
1	C	109	Total	C	N	O	S	0	0	0
			834	520	147	161	6			

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

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

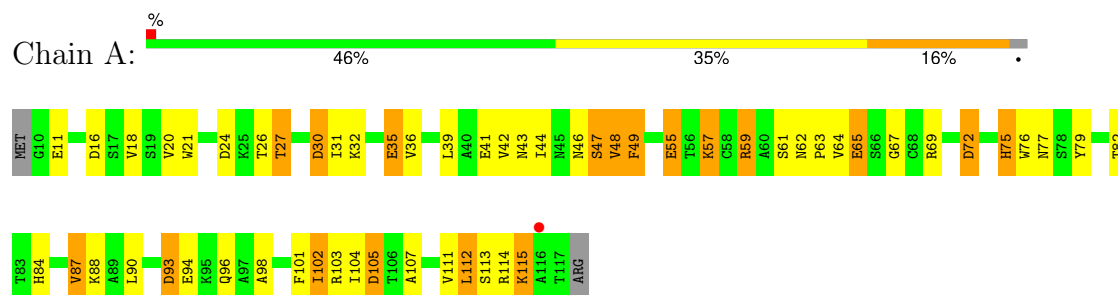
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	19	Total	O	0	0
			19	19		
3	B	20	Total	O	0	0
			20	20		
3	C	15	Total	O	0	0
			15	15		

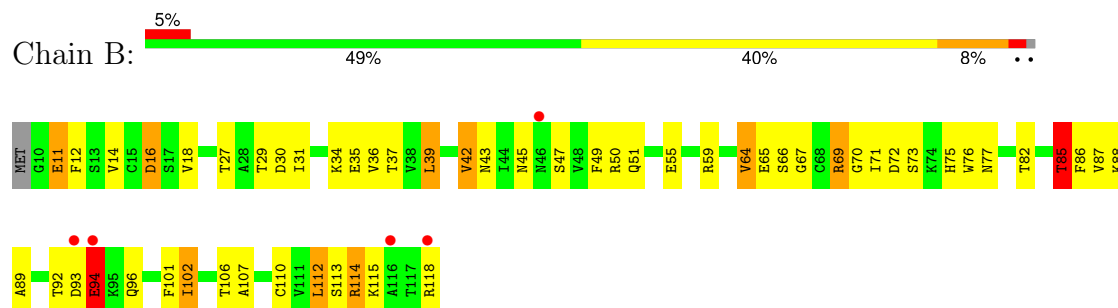
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

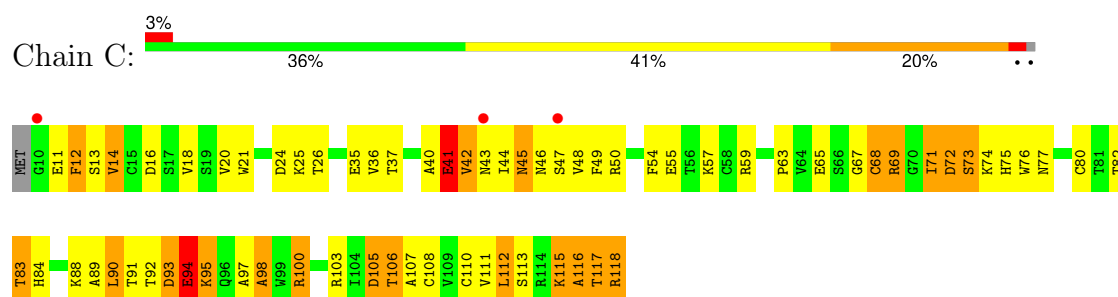
• Molecule 1: BETA NERVE GROWTH FACTOR



• Molecule 1: BETA NERVE GROWTH FACTOR



• Molecule 1: BETA NERVE GROWTH FACTOR



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.80Å 93.90Å 58.90Å 90.00° 123.10° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.50) 87.2 (20.00-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.85 (at 2.41Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.179 , (Not available) 0.173 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 107.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.008 for -k+l,-h-l,-l 0.011 for k+l,h+l,-l 0.036 for -h-2*l,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2565	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.30% 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	1.03	5/858 (0.6%)	1.36	11/1164 (0.9%)
1	B	1.09	5/848 (0.6%)	1.45	9/1152 (0.8%)
1	C	1.11	5/850 (0.6%)	1.44	11/1156 (1.0%)
All	All	1.08	15/2556 (0.6%)	1.42	31/3472 (0.9%)

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	1	0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	59	ARG	NE-CZ	11.41	1.47	1.33
1	B	94	GLU	CD-OE1	8.66	1.35	1.25
1	C	11	GLU	CD-OE2	7.56	1.33	1.25
1	A	11	GLU	CD-OE2	7.00	1.33	1.25
1	A	41	GLU	CD-OE1	6.70	1.33	1.25
1	C	41	GLU	CD-OE1	6.60	1.32	1.25
1	B	11	GLU	CD-OE2	6.45	1.32	1.25
1	C	35	GLU	CD-OE1	6.22	1.32	1.25
1	A	35	GLU	CD-OE1	6.15	1.32	1.25
1	C	94	GLU	CD-OE1	5.72	1.31	1.25
1	C	65	GLU	CD-OE1	5.67	1.31	1.25
1	A	55	GLU	CD-OE1	5.53	1.31	1.25
1	B	35	GLU	CD-OE1	5.47	1.31	1.25
1	A	65	GLU	CD-OE2	5.07	1.31	1.25
1	B	55	GLU	CD-OE1	5.03	1.31	1.25

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	118	ARG	NE-CZ-NH1	10.04	125.32	120.30
1	C	118	ARG	NE-CZ-NH2	-7.98	116.31	120.30
1	C	24	ASP	CB-CG-OD2	7.45	125.00	118.30
1	A	24	ASP	CB-CG-OD2	-7.05	111.95	118.30
1	B	59	ARG	CD-NE-CZ	-7.00	113.81	123.60
1	C	72	ASP	CB-CG-OD1	6.99	124.59	118.30
1	C	24	ASP	CB-CG-OD1	-6.91	112.08	118.30
1	A	30	ASP	CB-CG-OD1	6.69	124.32	118.30
1	C	72	ASP	CB-CG-OD2	-6.67	112.30	118.30
1	B	69	ARG	NE-CZ-NH1	6.57	123.58	120.30
1	C	69	ARG	NE-CZ-NH1	6.51	123.55	120.30
1	A	72	ASP	CB-CG-OD2	6.49	124.14	118.30
1	A	24	ASP	CB-CG-OD1	6.33	124.00	118.30
1	B	118	ARG	NE-CZ-NH2	-6.28	117.16	120.30
1	C	16	ASP	CB-CG-OD1	-6.20	112.72	118.30
1	C	93	ASP	N-CA-C	-6.15	94.39	111.00
1	C	69	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	114	ARG	NE-CZ-NH1	5.89	123.25	120.30
1	C	100	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	30	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	C	105	ASP	CB-CG-OD1	-5.67	113.19	118.30
1	A	105	ASP	CB-CG-OD1	-5.66	113.21	118.30
1	A	69	ARG	NE-CZ-NH1	5.63	123.11	120.30
1	A	93	ASP	CB-CG-OD1	-5.63	113.24	118.30
1	A	105	ASP	CB-CG-OD2	5.48	123.23	118.30
1	B	93	ASP	CB-CG-OD1	-5.48	113.36	118.30
1	B	16	ASP	CB-CG-OD1	-5.21	113.61	118.30
1	B	85	THR	N-CA-CB	5.12	120.03	110.30
1	A	93	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	16	ASP	CB-CG-OD2	5.03	122.82	118.30
1	B	93	ASP	CB-CG-OD2	5.01	122.81	118.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	82	THR	CB

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	842	0	810	52	0
1	B	832	0	784	48	0
1	C	834	0	775	60	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
3	A	19	0	0	1	1
3	B	20	0	0	3	0
3	C	15	0	0	0	0
All	All	2565	0	2369	151	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 31.

All (151) 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:45:ASN:ND2	1:C:45:ASN:H	1.52	1.08
1:B:45:ASN:HD21	1:C:45:ASN:N	1.62	0.96
1:A:47:SER:HB3	1:A:49:PHE:CZ	2.08	0.88
1:C:12:PHE:HD1	1:C:13:SER:H	1.19	0.87
1:B:45:ASN:HD21	1:C:45:ASN:H	0.81	0.81
1:C:92:THR:HG22	1:C:97:ALA:HA	1.62	0.79
1:B:89:ALA:HB3	1:B:102:ILE:HD13	1.67	0.77
1:C:90:LEU:HD12	1:C:91:THR:N	2.01	0.75
1:C:42:VAL:HG12	1:C:49:PHE:HB2	1.68	0.74
1:A:43:ASN:C	1:A:44:ILE:HD13	2.06	0.74
1:C:75:HIS:O	1:C:115:LYS:HG2	1.89	0.72
1:A:43:ASN:HD21	1:A:47:SER:N	1.89	0.70
1:A:43:ASN:O	1:A:44:ILE:HD13	1.90	0.70
1:B:94:GLU:HA	1:B:94:GLU:OE1	1.90	0.70
1:A:43:ASN:ND2	1:A:47:SER:N	2.41	0.69
1:C:83:THR:HG22	1:C:106:THR:HG22	1.76	0.68
1:B:39:LEU:O	1:B:51:GLN:NE2	2.28	0.67
1:C:73:SER:O	1:C:115:LYS:NZ	2.27	0.67
1:A:44:ILE:N	1:A:47:SER:O	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:ASP:OD2	1:C:75:HIS:ND1	2.29	0.66
1:C:98:ALA:HB3	1:C:100:ARG:NH1	2.11	0.66
1:B:42:VAL:N	1:B:49:PHE:O	2.29	0.65
1:A:42:VAL:HG12	1:A:49:PHE:O	1.97	0.65
1:A:72:ASP:OD2	1:A:75:HIS:NE2	2.30	0.65
1:B:75:HIS:NE2	3:B:904:HOH:O	2.29	0.65
1:B:77:ASN:HB2	1:B:113:SER:OG	1.97	0.65
1:C:42:VAL:N	1:C:49:PHE:O	2.29	0.65
1:C:94:GLU:OE1	1:C:95:LYS:N	2.27	0.65
3:B:954:HOH:O	1:C:113:SER:HB2	1.97	0.64
1:C:91:THR:OG1	1:C:100:ARG:NH1	2.30	0.64
1:C:57:LYS:HE2	1:C:105:ASP:OD2	1.98	0.63
1:B:64:VAL:HG12	1:B:67:GLY:H	1.66	0.61
1:A:43:ASN:HD21	1:A:46:ASN:CA	2.15	0.59
1:C:14:VAL:HG13	1:C:68:CYS:HB3	1.85	0.59
1:B:43:ASN:HA	1:B:47:SER:O	2.03	0.59
1:B:64:VAL:HG13	1:B:65:GLU:N	2.18	0.59
1:B:76:TRP:CH2	1:B:114:ARG:HD2	2.38	0.58
1:C:41:GLU:HB3	1:C:49:PHE:O	2.03	0.58
1:A:88:LYS:HG3	1:A:101:PHE:CZ	2.39	0.58
1:B:45:ASN:ND2	1:C:45:ASN:N	2.33	0.58
1:B:71:ILE:HG22	1:B:72:ASP:N	2.18	0.58
1:B:114:ARG:HH11	1:B:114:ARG:HG3	1.68	0.58
1:A:75:HIS:NE2	3:A:903:HOH:O	2.26	0.57
1:C:26:THR:O	1:C:37:THR:HB	2.05	0.57
1:C:90:LEU:HD12	1:C:91:THR:H	1.70	0.57
1:B:114:ARG:HG3	1:B:115:LYS:N	2.20	0.56
1:B:37:THR:O	1:B:92:THR:N	2.39	0.56
1:C:92:THR:HG22	1:C:97:ALA:CA	2.34	0.55
1:B:85:THR:HG22	1:C:54:PHE:HD2	1.71	0.55
1:A:43:ASN:HD22	1:A:47:SER:C	2.11	0.54
1:B:82:THR:HA	1:B:107:ALA:O	2.08	0.53
1:B:92:THR:HA	1:B:96:GLN:O	2.09	0.53
1:C:74:LYS:O	1:C:115:LYS:HG3	2.09	0.53
1:C:14:VAL:HG11	1:C:110:CYS:HB2	1.91	0.53
1:A:76:TRP:CZ3	1:A:114:ARG:HG2	2.44	0.52
1:C:77:ASN:HB2	1:C:113:SER:HB3	1.90	0.52
1:A:47:SER:HB3	1:A:49:PHE:HZ	1.71	0.52
1:C:18:VAL:CG1	1:C:59:ARG:HG3	2.39	0.52
1:C:41:GLU:OE1	1:C:48:VAL:HG11	2.10	0.52
1:C:42:VAL:HG23	1:C:90:LEU:HD23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:ASN:HA	1:C:47:SER:O	2.10	0.51
1:A:84:HIS:NE2	1:A:105:ASP:OD1	2.44	0.51
1:C:76:TRP:C	1:C:115:LYS:HD3	2.31	0.51
1:C:90:LEU:HD12	1:C:90:LEU:C	2.29	0.51
1:A:26:THR:OG1	1:A:27:THR:HG22	2.11	0.51
1:A:43:ASN:ND2	1:A:47:SER:H	2.08	0.51
1:B:30:ASP:OD1	1:B:34:LYS:HG2	2.10	0.51
1:A:82:THR:HA	1:A:107:ALA:O	2.11	0.50
1:C:116:ALA:O	1:C:117:THR:HG23	2.11	0.50
1:C:25:LYS:NZ	1:C:55:GLU:OE2	2.31	0.50
1:A:87:VAL:HG22	1:A:104:ILE:HD13	1.93	0.50
1:C:83:THR:CG2	1:C:106:THR:HG22	2.42	0.50
1:A:102:ILE:HG12	1:A:104:ILE:HD12	1.94	0.49
1:B:73:SER:O	1:B:115:LYS:NZ	2.38	0.49
1:A:18:VAL:O	1:A:18:VAL:HG13	2.13	0.49
1:C:14:VAL:CG1	1:C:110:CYS:HB2	2.43	0.49
1:A:102:ILE:HD11	1:A:104:ILE:HD11	1.95	0.49
1:C:12:PHE:HD1	1:C:13:SER:N	1.97	0.49
1:A:49:PHE:CD1	1:A:49:PHE:N	2.79	0.48
1:C:18:VAL:HG13	1:C:59:ARG:HG3	1.95	0.48
1:C:71:ILE:N	1:C:71:ILE:HD13	2.28	0.48
1:A:39:LEU:HD12	1:A:39:LEU:N	2.29	0.48
1:B:16:ASP:CG	1:B:69:ARG:HH22	2.16	0.48
1:B:85:THR:HG22	1:B:86:PHE:H	1.78	0.48
1:A:20:VAL:HG22	1:A:21:TRP:N	2.29	0.47
1:A:42:VAL:O	1:A:49:PHE:N	2.38	0.47
1:A:93:ASP:OD1	1:A:93:ASP:N	2.41	0.47
1:C:14:VAL:HG22	1:C:69:ARG:O	2.15	0.47
1:B:49:PHE:CD1	1:B:49:PHE:N	2.81	0.47
1:C:98:ALA:HB3	1:C:100:ARG:CZ	2.44	0.47
1:B:71:ILE:CG2	1:B:72:ASP:N	2.77	0.47
1:C:97:ALA:O	1:C:98:ALA:HB2	2.15	0.47
1:C:72:ASP:OD1	1:C:74:LYS:CG	2.63	0.47
1:C:82:THR:HA	1:C:107:ALA:O	2.15	0.46
1:B:70:GLY:HA3	1:C:112:LEU:HD21	1.98	0.46
1:B:112:LEU:HD12	1:B:112:LEU:HA	1.47	0.46
1:B:114:ARG:HG3	1:B:115:LYS:H	1.79	0.46
1:B:64:VAL:CG1	1:B:65:GLU:N	2.79	0.46
1:B:76:TRP:CE3	1:B:114:ARG:HA	2.51	0.46
1:C:36:VAL:HG13	1:C:91:THR:HB	1.97	0.45
1:A:31:ILE:HG23	1:A:32:LYS:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:SER:HB3	1:A:49:PHE:CE1	2.50	0.45
1:A:42:VAL:O	1:A:48:VAL:HA	2.16	0.45
1:A:43:ASN:HD21	1:A:46:ASN:C	2.20	0.45
1:B:87:VAL:O	1:B:101:PHE:HA	2.16	0.45
1:A:35:GLU:HG3	1:A:36:VAL:N	2.32	0.45
1:A:77:ASN:OD1	1:A:115:LYS:HE2	2.17	0.45
1:B:16:ASP:OD1	1:B:69:ARG:NH2	2.48	0.45
1:B:76:TRP:CZ3	1:B:114:ARG:HA	2.51	0.45
1:A:43:ASN:HA	1:A:48:VAL:HA	1.98	0.45
1:B:31:ILE:HD11	1:C:21:TRP:HZ2	1.81	0.45
1:A:43:ASN:HB2	1:A:48:VAL:HG23	1.99	0.44
1:A:55:GLU:HG2	1:A:104:ILE:HG23	1.99	0.44
1:A:93:ASP:O	1:A:94:GLU:HB2	2.17	0.44
1:C:89:ALA:O	1:C:91:THR:HG23	2.18	0.44
1:B:47:SER:CB	1:B:49:PHE:CZ	3.00	0.44
1:A:44:ILE:HD13	1:A:44:ILE:N	2.27	0.44
1:A:90:LEU:HD12	1:A:98:ALA:O	2.17	0.44
1:B:106:THR:HA	3:B:936:HOH:O	2.18	0.44
1:A:76:TRP:CE3	1:A:114:ARG:HG2	2.53	0.44
1:C:46:ASN:O	1:C:46:ASN:ND2	2.51	0.44
1:C:93:ASP:N	1:C:93:ASP:OD1	2.46	0.44
1:A:18:VAL:HG12	1:A:57:LYS:O	2.18	0.44
1:A:79:TYR:HD1	1:B:12:PHE:CZ	2.35	0.44
1:C:84:HIS:HB3	1:C:103:ARG:HG3	2.00	0.44
1:B:89:ALA:HB3	1:B:102:ILE:CD1	2.44	0.43
1:C:80:CYS:HB3	1:C:108:CYS:SG	2.59	0.43
1:A:30:ASP:HA	1:A:102:ILE:HA	2.01	0.43
1:C:55:GLU:OE1	1:C:105:ASP:HB2	2.19	0.43
1:A:62:ASN:OD1	1:A:67:GLY:HA2	2.19	0.43
1:B:86:PHE:C	1:B:87:VAL:HG23	2.38	0.43
1:B:14:VAL:HG12	1:B:110:CYS:HB2	2.00	0.43
1:A:75:HIS:HB3	1:A:114:ARG:NH2	2.34	0.42
1:C:40:ALA:O	1:C:50:ARG:HG3	2.19	0.42
1:B:73:SER:HA	1:B:76:TRP:O	2.20	0.42
1:B:114:ARG:HH11	1:B:115:LYS:H	1.66	0.42
1:A:59:ARG:CZ	1:A:59:ARG:HB3	2.46	0.42
1:A:82:THR:CG2	1:A:105:ASP:HB3	2.50	0.42
1:C:18:VAL:CG1	1:C:59:ARG:CG	2.98	0.42
1:C:42:VAL:CG2	1:C:90:LEU:HD23	2.50	0.41
1:A:43:ASN:HA	1:A:47:SER:O	2.20	0.41
1:C:63:PRO:HD2	1:C:67:GLY:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:SER:OG	1:C:117:THR:N	2.53	0.41
1:A:112:LEU:HA	1:A:112:LEU:HD12	1.86	0.41
1:B:106:THR:O	1:B:107:ALA:HB2	2.20	0.41
1:B:114:ARG:HG3	1:B:114:ARG:NH1	2.35	0.41
1:B:76:TRP:CZ3	1:B:114:ARG:HD2	2.56	0.41
1:B:89:ALA:CB	1:B:102:ILE:CD1	3.00	0.40
1:C:76:TRP:O	1:C:115:LYS:HD3	2.21	0.40
1:A:31:ILE:CG2	1:A:32:LYS:N	2.83	0.40
1:A:102:ILE:CG1	1:A:104:ILE:CD1	2.98	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)
3:A:922:HOH:O	3:A:922:HOH:O[2_654]	2.02	0.18

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	106/110 (96%)	102 (96%)	4 (4%)	0	100	100
1	B	107/110 (97%)	99 (92%)	8 (8%)	0	100	100
1	C	107/110 (97%)	100 (94%)	3 (3%)	4 (4%)	2	3
All	All	320/330 (97%)	301 (94%)	15 (5%)	4 (1%)	10	19

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	44	ILE
1	C	95	LYS
1	C	116	ALA

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Mol	Chain	Res	Type
1	C	98	ALA

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	94/97 (97%)	76 (81%)	18 (19%)	1	2
1	B	89/97 (92%)	74 (83%)	15 (17%)	1	3
1	C	89/97 (92%)	70 (79%)	19 (21%)	1	1
All	All	272/291 (94%)	220 (81%)	52 (19%)	1	2

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	THR
1	A	47	SER
1	A	48	VAL
1	A	49	PHE
1	A	57	LYS
1	A	59	ARG
1	A	63	PRO
1	A	64	VAL
1	A	65	GLU
1	A	75	HIS
1	A	87	VAL
1	A	96	GLN
1	A	102	ILE
1	A	103	ARG
1	A	111	VAL
1	A	112	LEU
1	A	113	SER
1	A	115	LYS
1	B	11	GLU
1	B	18	VAL
1	B	27	THR

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Mol	Chain	Res	Type
1	B	29	THR
1	B	36	VAL
1	B	39	LEU
1	B	42	VAL
1	B	50	ARG
1	B	64	VAL
1	B	66	SER
1	B	85	THR
1	B	88	LYS
1	B	94	GLU
1	B	102	ILE
1	B	112	LEU
1	C	12	PHE
1	C	14	VAL
1	C	20	VAL
1	C	41	GLU
1	C	42	VAL
1	C	45	ASN
1	C	68	CYS
1	C	71	ILE
1	C	73	SER
1	C	83	THR
1	C	88	LYS
1	C	90	LEU
1	C	94	GLU
1	C	106	THR
1	C	111	VAL
1	C	112	LEU
1	C	115	LYS
1	C	117	THR
1	C	118	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	96	GLN
1	B	45	ASN
1	B	77	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 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	108/110 (98%)	-0.64	1 (0%) 81 78	10, 25, 67, 80	0
1	B	109/110 (99%)	-0.35	5 (4%) 38 35	10, 32, 79, 80	0
1	C	109/110 (99%)	-0.41	3 (2%) 55 51	10, 28, 77, 80	0
All	All	326/330 (98%)	-0.47	9 (2%) 55 51	10, 28, 78, 80	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	93	ASP	3.2
1	B	94	GLU	3.0
1	B	116	ALA	2.8
1	B	118	ARG	2.7
1	C	43	ASN	2.7
1	B	46	ASN	2.2
1	C	10	GLY	2.1
1	A	116	ALA	2.0
1	C	47	SER	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(Å ²)	Q<0.9
2	ZN	C	900	1/1	0.99	0.02	36,36,36,36	0
2	ZN	B	902	1/1	1.00	0.03	42,42,42,42	0
2	ZN	C	901	1/1	1.00	0.02	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.