Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 – 07:37 PM EDT

PDB ID	:	2NP0
Title	:	Crystal structure of the Botulinum neurotoxin type B complexed with synaptotagamin-II ectodomain
Authors	:	Chai, Q.; Arndt, J.W.; Stevens, R.C.
Deposited on	:	2006-10-26
Resolution	:	2.62 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 (i) 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 (1)) were used in the production of this report:

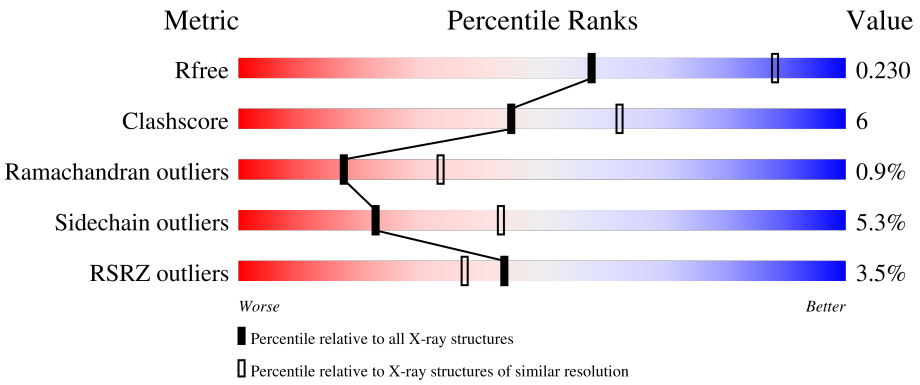
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3797 (2.64-2.60)
Clashscore	141614	4168 (2.64-2.60)
Ramachandran outliers	138981	4093 (2.64-2.60)
Sidechain outliers	138945	4093 (2.64-2.60)
RSRZ outliers	127900	3731 (2.64-2.60)

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 <=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	1290	3% 84% 13% •
2	B	21	33% 57% 14% 29%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 11182 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 B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1289	10640	6861	1713	2033	33	0	0	0

- Molecule 2 is a protein called Synaptotagmin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	15	131	87	20	23	1	0	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
3	A	1	1	1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
4	A	3	3	3	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
5	A	1	1	1	0	0

- Molecule 6 is water.

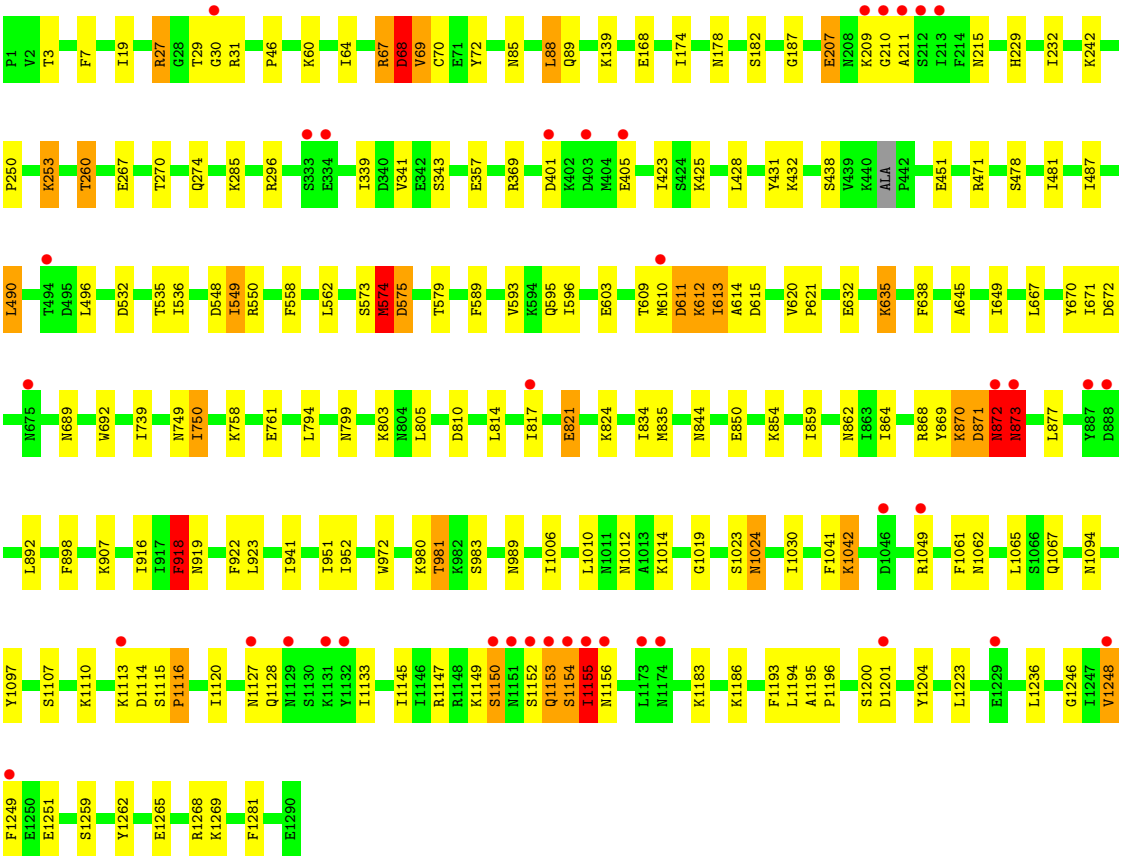
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	406	406	406	0	0

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.

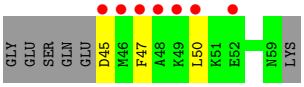
• Molecule 1: Botulinum neurotoxin type B

Chain A: 3% 84% 13%



• Molecule 2: Synaptotagmin-2

Chain B: 33% 57% 14% 29%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	170.41Å 351.97Å 101.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.62 39.31 – 2.62	Depositor EDS
% Data completeness (in resolution range)	90.0 (40.00-2.62) 90.0 (39.31-2.62)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.186 , 0.227 0.191 , 0.230	Depositor DCC
R_{free} test set	4177 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	44.8	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 42.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11182	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, CA, 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	0.70	2/10872 (0.0%)	0.68	0/14683
2	B	0.59	0/133	0.59	0/174
All	All	0.70	2/11005 (0.0%)	0.68	0/14857

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	A	0	13

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	168	GLU	CD-OE1	6.34	1.32	1.25
1	A	850	GLU	CD-OE1	5.21	1.31	1.25

There are no bond angle outliers.

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1150	SER	Peptide
1	A	1155	ILE	Peptide
1	A	209	LYS	Peptide
1	A	30	GLY	Peptide
1	A	549	ILE	Peptide
1	A	574	MET	Peptide

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Mol	Chain	Res	Type	Group
1	A	575	ASP	Peptide
1	A	68	ASP	Peptide
1	A	870	LYS	Peptide
1	A	872	ASN	Peptide
1	A	873	ASN	Peptide
1	A	918	PHE	Peptide
1	A	919	ASN	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	10640	0	10503	138	0
2	B	131	0	129	2	0
3	A	1	0	0	0	0
4	A	3	0	0	0	0
5	A	1	0	0	0	0
6	A	406	0	0	9	0
All	All	11182	0	10632	139	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 6.

All (139) 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:1155:ILE:HA	1:A:1156:ASN:CB	1.67	1.21
1:A:872:ASN:N	1:A:873:ASN:HB2	1.56	1.20
1:A:799:ASN:HB3	6:A:1540:HOH:O	1.48	1.13
1:A:1152:SER:HA	1:A:1153:GLN:C	1.69	1.10
1:A:1155:ILE:HA	1:A:1156:ASN:HB2	1.23	1.09
1:A:1153:GLN:H	1:A:1153:GLN:CD	1.62	1.01
1:A:1154:SER:O	1:A:1155:ILE:HG22	1.65	0.97
1:A:549:ILE:HG23	1:A:549:ILE:O	1.65	0.96
1:A:1155:ILE:HG23	1:A:1155:ILE:O	1.63	0.96
1:A:1155:ILE:HA	1:A:1156:ASN:HB3	1.53	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1155:ILE:CA	1:A:1156:ASN:HB2	2.01	0.90
1:A:873:ASN:H	1:A:892:LEU:CD1	1.85	0.89
1:A:1155:ILE:CA	1:A:1156:ASN:CB	2.50	0.88
1:A:869:TYR:O	1:A:870:LYS:HG2	1.74	0.88
1:A:1248:VAL:HG23	1:A:1249:PHE:H	1.38	0.87
1:A:871:ASP:C	1:A:873:ASN:HB2	1.96	0.85
1:A:1152:SER:HA	1:A:1153:GLN:O	1.79	0.80
1:A:579:THR:HG22	1:A:596:ILE:HD13	1.65	0.79
1:A:253:LYS:HG3	1:A:253:LYS:O	1.83	0.78
1:A:1153:GLN:CD	1:A:1153:GLN:N	2.37	0.78
1:A:1155:ILE:O	1:A:1155:ILE:CG2	2.32	0.77
1:A:432:LYS:HE3	6:A:1349:HOH:O	1.82	0.77
1:A:1154:SER:O	1:A:1155:ILE:CG2	2.32	0.77
1:A:873:ASN:H	1:A:892:LEU:HD13	1.52	0.75
1:A:253:LYS:CB	1:A:253:LYS:NZ	2.49	0.75
1:A:1010:LEU:HD12	1:A:1010:LEU:O	1.86	0.74
1:A:549:ILE:O	1:A:549:ILE:CG2	2.35	0.73
1:A:871:ASP:O	1:A:872:ASN:HB2	1.88	0.72
1:A:614:ALA:O	1:A:615:ASP:HB2	1.89	0.70
1:A:611:ASP:OD1	1:A:611:ASP:C	2.31	0.69
1:A:611:ASP:OD1	1:A:612:LYS:N	2.29	0.66
1:A:1248:VAL:HG23	1:A:1249:PHE:N	2.09	0.66
1:A:425:LYS:HD3	1:A:428:LEU:HD12	1.79	0.65
1:A:1154:SER:O	1:A:1155:ILE:CB	2.45	0.65
1:A:869:TYR:O	1:A:870:LYS:CG	2.47	0.62
1:A:916:ILE:HA	1:A:918:PHE:CE2	2.35	0.62
1:A:70:CYS:SG	1:A:428:LEU:HD13	2.41	0.60
1:A:253:LYS:NZ	1:A:253:LYS:HB3	2.16	0.60
1:A:267:GLU:HG3	6:A:1300:HOH:O	2.01	0.60
1:A:817:ILE:H	1:A:817:ILE:HD12	1.66	0.59
1:A:941:ILE:HD12	1:A:1133:ILE:HD13	1.84	0.59
1:A:952:ILE:HD13	1:A:1041:PHE:CE2	2.38	0.59
1:A:1152:SER:CA	1:A:1153:GLN:C	2.56	0.58
1:A:1183:LYS:HD3	1:A:1204:TYR:CE1	2.38	0.58
1:A:941:ILE:CD1	1:A:1133:ILE:HD13	2.34	0.58
1:A:980:LYS:HG2	1:A:1030:ILE:HG22	1.85	0.57
1:A:1265:GLU:OE1	1:A:1268:ARG:NH2	2.38	0.57
2:B:45:ASP:OD1	2:B:47:PHE:HB2	2.05	0.56
1:A:242:LYS:NZ	6:A:1347:HOH:O	2.34	0.55
1:A:229:HIS:ND1	1:A:357:GLU:OE1	2.38	0.54
1:A:870:LYS:O	1:A:871:ASP:CG	2.45	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:573:SER:O	1:A:575:ASP:N	2.42	0.53
1:A:573:SER:OG	1:A:603:GLU:OE2	2.26	0.53
1:A:870:LYS:C	1:A:871:ASP:OD1	2.46	0.53
1:A:1195:ALA:HB1	1:A:1196:PRO:HD2	1.92	0.52
1:A:868:ARG:HG3	1:A:877:LEU:HD12	1.90	0.52
1:A:19:ILE:HD12	1:A:139:LYS:HB3	1.90	0.52
1:A:1107:SER:HB3	1:A:1120:ILE:CG2	2.40	0.52
1:A:67:ARG:HB3	1:A:431:TYR:CD2	2.46	0.51
1:A:490:LEU:CD1	1:A:496:LEU:HD22	2.41	0.50
1:A:872:ASN:H	1:A:873:ASN:HB2	1.68	0.50
1:A:471:ARG:HD3	6:A:1573:HOH:O	2.12	0.50
1:A:253:LYS:NZ	1:A:253:LYS:HB2	2.24	0.49
1:A:72:TYR:CD1	1:A:423:ILE:HG21	2.48	0.49
1:A:871:ASP:O	1:A:873:ASN:ND2	2.45	0.49
1:A:64:ILE:HD12	1:A:64:ILE:O	2.11	0.49
1:A:211:ALA:HB3	1:A:215:ASN:HB3	1.95	0.49
1:A:296:ARG:HG2	1:A:339:ILE:HD12	1.94	0.49
1:A:1154:SER:O	1:A:1155:ILE:HB	2.13	0.49
1:A:834:ILE:HD12	1:A:835:MET:C	2.33	0.48
1:A:689:ASN:HB3	1:A:835:MET:HE1	1.95	0.48
1:A:253:LYS:CB	1:A:253:LYS:HZ1	2.23	0.48
1:A:1194:LEU:HD12	1:A:1194:LEU:N	2.29	0.48
1:A:922:PHE:HD2	1:A:1010:LEU:HD13	1.79	0.48
1:A:558:PHE:CZ	1:A:562:LEU:HD11	2.49	0.47
1:A:1193:PHE:C	1:A:1194:LEU:HD12	2.34	0.47
1:A:859:ILE:HA	1:A:862:ASN:HD22	1.80	0.47
1:A:739:ILE:O	1:A:739:ILE:CG2	2.63	0.47
1:A:907:LYS:HB3	1:A:1042:LYS:HG3	1.97	0.47
1:A:941:ILE:HD11	1:A:1133:ILE:HG23	1.96	0.47
1:A:1114:ASP:OD1	1:A:1114:ASP:N	2.47	0.46
1:A:667:LEU:CD1	1:A:805:LEU:HD23	2.45	0.46
1:A:870:LYS:O	1:A:871:ASP:OD1	2.32	0.46
1:A:1116:PRO:HB3	2:B:50:LEU:HD22	1.97	0.46
1:A:670:TYR:O	1:A:671:ILE:C	2.54	0.46
1:A:1094:ASN:HA	1:A:1145:ILE:HD11	1.97	0.46
1:A:750:ILE:HD13	1:A:750:ILE:HA	1.79	0.46
1:A:573:SER:C	1:A:575:ASP:N	2.69	0.46
1:A:981:THR:HG22	6:A:1521:HOH:O	2.15	0.46
1:A:285:LYS:HA	1:A:481:ILE:HD12	1.98	0.46
1:A:645:ALA:HB3	1:A:649:ILE:CG2	2.46	0.45
1:A:1183:LYS:CD	1:A:1204:TYR:CE1	2.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:972:TRP:CD2	1:A:1006:ILE:HG21	2.51	0.45
1:A:46:PRO:HB3	1:A:88:LEU:HD13	1.99	0.45
1:A:27:ARG:C	1:A:29:THR:H	2.20	0.45
1:A:250:PRO:HG3	1:A:260:THR:HG22	1.99	0.44
1:A:285:LYS:HA	1:A:481:ILE:CD1	2.48	0.44
1:A:589:PHE:O	1:A:593:VAL:HG23	2.17	0.44
1:A:174:ILE:HG13	6:A:1308:HOH:O	2.17	0.44
1:A:612:LYS:HE3	1:A:612:LYS:HB3	1.25	0.44
1:A:667:LEU:HD11	1:A:805:LEU:HD23	2.00	0.44
1:A:1012:ASN:OD1	1:A:1024:ASN:ND2	2.51	0.44
1:A:821:GLU:OE1	1:A:821:GLU:HA	2.18	0.43
1:A:182:SER:HA	1:A:187:GLY:HA2	2.00	0.43
1:A:573:SER:C	1:A:575:ASP:H	2.21	0.43
1:A:692:TRP:CE3	1:A:794:LEU:HD13	2.54	0.43
1:A:951:ILE:HG13	1:A:952:ILE:HG13	2.00	0.43
1:A:613:ILE:HG23	1:A:613:ILE:HD13	1.71	0.43
1:A:438:SER:O	1:A:532:ASP:HA	2.18	0.43
1:A:1259:SER:HB3	1:A:1262:TYR:CD2	2.53	0.43
1:A:548:ASP:OD1	1:A:548:ASP:N	2.41	0.43
1:A:490:LEU:HD12	1:A:490:LEU:O	2.19	0.43
1:A:573:SER:HB2	6:A:1487:HOH:O	2.18	0.43
1:A:232:ILE:CG2	1:A:270:THR:HG23	2.49	0.42
1:A:892:LEU:HD23	1:A:898:PHE:HB3	2.00	0.42
1:A:253:LYS:HB2	1:A:253:LYS:HE3	1.87	0.42
1:A:535:THR:OG1	1:A:536:ILE:N	2.53	0.42
1:A:68:ASP:OD1	1:A:69:VAL:HG22	2.19	0.42
1:A:1145:ILE:HD13	1:A:1147:ARG:NH2	2.35	0.42
1:A:1019:GLY:O	1:A:1067:GLN:NE2	2.48	0.42
1:A:85:ASN:O	1:A:89:GLN:HG2	2.19	0.42
1:A:207:GLU:OE1	1:A:210:GLY:HA2	2.20	0.42
1:A:253:LYS:HB3	1:A:253:LYS:HZ2	1.84	0.42
1:A:1061:PHE:CD2	1:A:1065:LEU:HD21	2.55	0.42
1:A:1248:VAL:CG2	1:A:1249:PHE:N	2.78	0.42
1:A:620:VAL:HA	1:A:621:PRO:HD3	1.85	0.41
1:A:1223:LEU:HA	1:A:1236:LEU:HD23	2.01	0.41
1:A:972:TRP:CG	1:A:1006:ILE:HG21	2.55	0.41
1:A:864:ILE:HD13	1:A:1062:ASN:HB3	2.01	0.41
1:A:692:TRP:CE2	1:A:794:LEU:HB3	2.56	0.41
1:A:1097:TYR:CG	1:A:1281:PHE:HB3	2.56	0.41
1:A:638:PHE:HA	6:A:1422:HOH:O	2.21	0.41
1:A:859:ILE:HD11	1:A:1065:LEU:HD23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:632:GLU:HB3	1:A:635:LYS:HG3	2.03	0.41
1:A:870:LYS:O	1:A:871:ASP:CB	2.69	0.41
1:A:490:LEU:HD12	1:A:496:LEU:HD22	2.02	0.41
1:A:1006:ILE:HA	1:A:1014:LYS:O	2.20	0.40
1:A:834:ILE:HD12	1:A:834:ILE:C	2.41	0.40
1:A:613:ILE:HG22	1:A:613:ILE:O	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	1285/1290 (100%)	1210 (94%)	63 (5%)	12 (1%)	17	33
2	B	13/21 (62%)	13 (100%)	0	0	100	100
All	All	1298/1311 (99%)	1223 (94%)	63 (5%)	12 (1%)	17	33

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	69	VAL
1	A	550	ARG
1	A	871	ASP
1	A	872	ASN
1	A	873	ASN
1	A	1155	ILE
1	A	1248	VAL
1	A	574	MET
1	A	1154	SER
1	A	1116	PRO
1	A	1201	ASP
1	A	1246	GLY

5.3.2 Protein sidechains (i)

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	1189/1189 (100%)	1125 (95%)	64 (5%)	22 42
2	B	14/19 (74%)	14 (100%)	0	100 100
All	All	1203/1208 (100%)	1139 (95%)	64 (5%)	22 43

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	THR
1	A	7	PHE
1	A	27	ARG
1	A	31	ARG
1	A	60	LYS
1	A	67	ARG
1	A	68	ASP
1	A	88	LEU
1	A	178	ASN
1	A	207	GLU
1	A	253	LYS
1	A	260	THR
1	A	274	GLN
1	A	341	VAL
1	A	343	SER
1	A	369	ARG
1	A	401	ASP
1	A	405	GLU
1	A	451	GLU
1	A	478	SER
1	A	487	ILE
1	A	490	LEU
1	A	574	MET
1	A	595	GLN
1	A	609	THR
1	A	610	MET
1	A	611	ASP

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Mol	Chain	Res	Type
1	A	612	LYS
1	A	613	ILE
1	A	635	LYS
1	A	672	ASP
1	A	749	ASN
1	A	750	ILE
1	A	758	LYS
1	A	761	GLU
1	A	803	LYS
1	A	810	ASP
1	A	814	LEU
1	A	821	GLU
1	A	824	LYS
1	A	844	ASN
1	A	854	LYS
1	A	918	PHE
1	A	923	LEU
1	A	981	THR
1	A	983	SER
1	A	989	ASN
1	A	1023	SER
1	A	1024	ASN
1	A	1042	LYS
1	A	1049	ARG
1	A	1110	LYS
1	A	1113	LYS
1	A	1115	SER
1	A	1127	ASN
1	A	1128	GLN
1	A	1149	LYS
1	A	1150	SER
1	A	1153	GLN
1	A	1155	ILE
1	A	1186	LYS
1	A	1200	SER
1	A	1251	GLU
1	A	1269	LYS

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	862	ASN

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Mol	Chain	Res	Type
1	A	912	GLN
1	A	1011	ASN
1	A	1069	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 5 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	1289/1290 (99%)	-0.14	39 (3%)	50	44	35, 45, 56, 68	0
2	B	15/21 (71%)	1.61	7 (46%)	0	0	50, 57, 61, 62	0
All	All	1304/1311 (99%)	-0.12	46 (3%)	44	37	35, 45, 57, 68	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	212	SER	5.2
2	B	46	MET	5.1
1	A	1152	SER	5.0
1	A	1153	GLN	4.9
1	A	1155	ILE	4.7
1	A	1151	ASN	4.5
1	A	1129	ASN	4.2
1	A	1249	PHE	4.1
1	A	1154	SER	3.9
2	B	47	PHE	3.6
1	A	209	LYS	3.4
1	A	1173	LEU	3.3
1	A	1156	ASN	3.2
1	A	401	ASP	3.2
1	A	1127	ASN	3.1
1	A	1248	VAL	3.1
1	A	873	ASN	3.0
1	A	213	ILE	3.0
1	A	333	SER	2.9
1	A	1131	LYS	2.9
1	A	211	ALA	2.9
1	A	888	ASP	2.9
1	A	334	GLU	2.9
1	A	887	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	210	GLY	2.8
2	B	49	LYS	2.7
1	A	403	ASP	2.6
1	A	405	GLU	2.6
1	A	1049	ARG	2.6
1	A	1174	ASN	2.5
1	A	1132	TYR	2.5
1	A	872	ASN	2.5
1	A	1201	ASP	2.4
1	A	494	THR	2.4
1	A	30	GLY	2.4
1	A	1150	SER	2.3
2	B	48	ALA	2.3
2	B	52	GLU	2.3
1	A	1046	ASP	2.3
2	B	50	LEU	2.2
1	A	610	MET	2.1
1	A	675	ASN	2.1
2	B	45	ASP	2.1
1	A	817	ILE	2.1
1	A	1229	GLU	2.1
1	A	1113	LYS	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
5	CL	A	1294	1/1	0.87	0.19	66,66,66,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	CA	A	1293	1/1	0.92	0.11	68,68,68,68	0
4	CA	A	1295	1/1	0.96	0.05	44,44,44,44	0
4	CA	A	1292	1/1	0.97	0.13	48,48,48,48	0
3	ZN	A	1291	1/1	0.98	0.04	60,60,60,60	0

6.5 Other polymers (i)

There are no such residues in this entry.