Full wwPDB X-ray Structure Validation Report (i)

Aug 19, 2023 – 06:20 PM EDT

PDB ID	:	2G7P
Title	:	Structure of the Light Chain of Botulinum Neurotoxin Serotype A Bound to Small Molecule Inhibitors
Authors	:	Fu, Z.; Baldwin, M.R.; Boldt, G.E.; Janda, K.D.; Barbieri, J.T.; Kim, J.-J.P.
Deposited on	:	2006-02-28
Resolution	:	2.30 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:

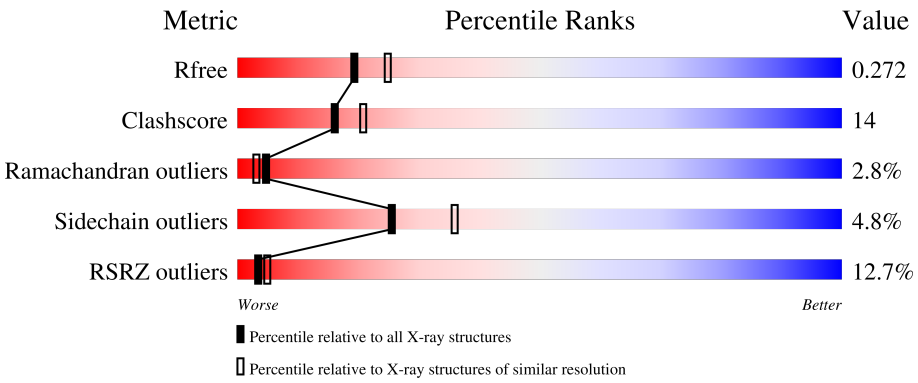
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	425	8% 73% 20% • •
1	B	425	16% 68% 24% • •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6864 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 A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	411	3323	2135	545	636	7	0	0	0
1	B	409	3300	2118	544	631	7	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q45894
A	362	ALA	ARG	engineered mutation	UNP Q45894
A	365	PHE	TYR	engineered mutation	UNP Q45894
B	1	MET	-	initiating methionine	UNP Q45894
B	362	ALA	ARG	engineered mutation	UNP Q45894
B	365	PHE	TYR	engineered mutation	UNP Q45894

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

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

- Molecule 3 is water.

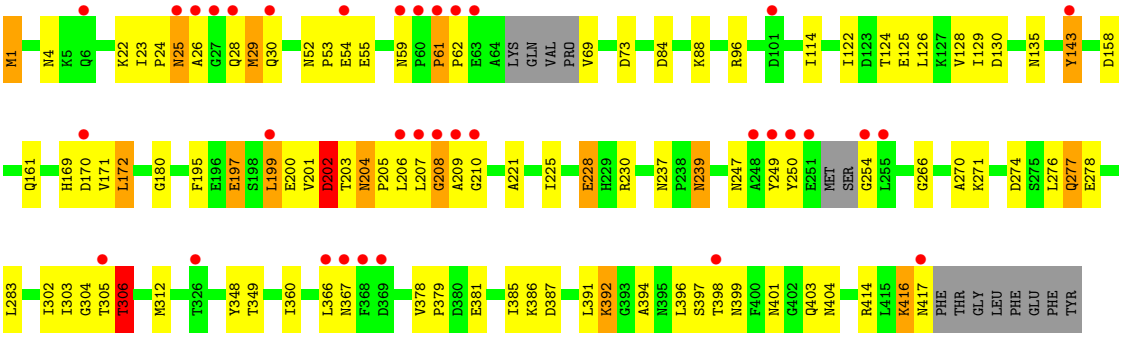
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	167	167	167	0	0
3	B	72	72	72	0	0

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.

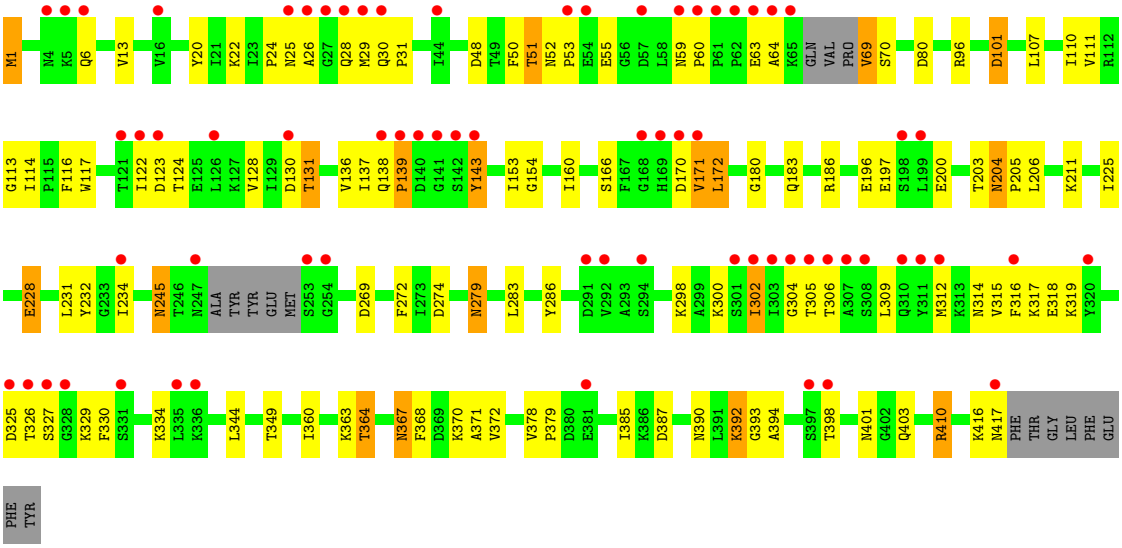
• Molecule 1: Botulinum neurotoxin type A

Chain A:



• Molecule 1: Botulinum neurotoxin type A

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	69.21Å 110.26Å 167.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.19 – 2.30 20.19 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.6 (20.19-2.30) 89.8 (20.19-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$< I/\sigma(I) >$ [1]	3.91 (at 2.30Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.275 0.230 , 0.272	Depositor DCC
R_{free} test set	5195 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å^2)	28.0	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 51.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6864	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 3.23% 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: 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	0.41	0/3398	0.64	0/4597
1	B	0.37	0/3373	0.58	0/4561
All	All	0.39	0/6771	0.61	0/9158

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3323	0	3252	92	0
1	B	3300	0	3241	99	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	167	0	0	5	0
3	B	72	0	0	2	0
All	All	6864	0	6493	190	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 14.

All (190) 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:416:LYS:HG2	1:B:417:ASN:H	1.14	1.08
1:A:225:ILE:O	1:A:228:GLU:HG3	1.69	0.93
1:A:171:VAL:HG23	1:A:172:LEU:HD13	1.55	0.89
1:A:61:PRO:HB2	1:A:62:PRO:HD2	1.55	0.88
1:B:410:ARG:HB2	1:B:410:ARG:HH11	1.38	0.85
1:B:416:LYS:HG2	1:B:417:ASN:N	1.91	0.85
1:A:199:LEU:HD23	1:A:199:LEU:H	1.42	0.84
1:B:387:ASP:OD2	1:B:392:LYS:HG3	1.78	0.84
1:B:204:ASN:ND2	1:B:206:LEU:H	1.77	0.83
1:A:204:ASN:HD22	1:A:205:PRO:N	1.82	0.77
1:A:417:ASN:ND2	3:A:698:HOH:O	2.21	0.74
1:A:204:ASN:HD22	1:A:204:ASN:C	1.92	0.73
1:B:22:LYS:HE2	1:B:31:PRO:HG3	1.73	0.70
1:B:410:ARG:HB2	1:B:410:ARG:NH1	2.05	0.70
1:B:143:TYR:H	1:B:143:TYR:HD2	1.40	0.69
1:A:387:ASP:OD2	1:A:392:LYS:HG3	1.92	0.69
1:B:138:GLN:HB3	1:B:139:PRO:HD2	1.74	0.68
1:A:391:LEU:O	1:A:397:SER:HB3	1.93	0.67
1:B:122:ILE:HG22	1:B:124:THR:HG22	1.75	0.67
1:B:325:ASP:OD2	1:B:329:LYS:HB3	1.95	0.66
1:A:52:ASN:ND2	1:A:55:GLU:HG3	2.11	0.66
1:A:22:LYS:HA	1:A:29:MET:HE1	1.77	0.66
1:A:52:ASN:HD22	1:A:55:GLU:HG3	1.62	0.65
1:B:300:LYS:O	1:B:309:LEU:HD22	1.97	0.65
1:B:160:ILE:HD13	1:B:372:VAL:HG21	1.79	0.64
1:B:371:ALA:HA	1:B:417:ASN:HB2	1.79	0.64
1:B:48:ASP:OD1	1:B:51:THR:HG22	1.97	0.64
1:A:1:MET:SD	1:A:1:MET:N	2.62	0.63
1:A:204:ASN:ND2	1:A:206:LEU:H	1.97	0.62
1:B:117:TRP:CE2	1:B:128:VAL:HB	2.34	0.62
1:B:52:ASN:HD22	1:B:55:GLU:HG3	1.65	0.62
1:A:23:ILE:HG23	1:A:24:PRO:HD2	1.80	0.62
1:A:225:ILE:HA	1:A:228:GLU:HG2	1.82	0.61
1:A:278:GLU:HG3	1:B:245:ASN:OD1	2.00	0.61
1:B:204:ASN:C	1:B:204:ASN:HD22	2.03	0.61
1:A:247:ASN:HB3	1:A:249:TYR:H	1.65	0.60
1:B:171:VAL:O	1:B:172:LEU:HB2	2.01	0.59
1:B:228:GLU:CG	1:B:344:LEU:HD22	2.33	0.59
1:A:61:PRO:HB2	1:A:62:PRO:CD	2.31	0.59
1:A:69:VAL:HG13	1:A:69:VAL:O	2.03	0.58
1:A:169:HIS:CE1	1:A:171:VAL:HG22	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:GLU:HG3	1:A:207:LEU:HD13	1.85	0.58
1:B:116:PHE:HA	1:B:316:PHE:CE1	2.39	0.58
1:A:225:ILE:O	1:A:228:GLU:CG	2.49	0.58
1:B:200:GLU:HB2	1:B:203:THR:HG23	1.86	0.58
1:B:371:ALA:HA	1:B:417:ASN:CB	2.34	0.58
1:B:13:VAL:HG11	1:B:20:TYR:CD1	2.39	0.58
1:B:378:VAL:HB	1:B:379:PRO:HD3	1.85	0.58
1:A:54:GLU:HA	3:A:667:HOH:O	2.04	0.57
1:B:228:GLU:OE1	1:B:344:LEU:O	2.23	0.57
1:A:199:LEU:H	1:A:199:LEU:CD2	2.16	0.57
1:A:237:ASN:CG	1:A:239:ASN:HD22	2.08	0.56
1:A:271:LYS:HE2	3:A:714:HOH:O	2.06	0.56
1:B:124:THR:HA	1:B:298:LYS:O	2.05	0.56
1:A:302:ILE:HD11	1:A:312:MET:HG3	1.87	0.56
1:B:136:VAL:HG12	1:B:137:ILE:N	2.21	0.56
1:A:398:THR:HG22	1:A:399:ASN:ND2	2.21	0.56
1:B:304:GLY:O	1:B:306:THR:HG22	2.06	0.56
1:A:228:GLU:OE2	1:A:348:TYR:HD2	1.89	0.56
1:B:138:GLN:HB3	1:B:139:PRO:CD	2.35	0.56
1:B:416:LYS:CG	1:B:417:ASN:H	1.99	0.56
1:A:381:GLU:H	1:A:381:GLU:CD	2.09	0.54
1:B:204:ASN:HD22	1:B:205:PRO:N	2.04	0.54
1:A:199:LEU:HD23	1:A:199:LEU:N	2.18	0.54
1:A:171:VAL:HG23	1:A:172:LEU:CD1	2.34	0.54
1:B:96:ARG:HA	1:B:385:ILE:HG23	1.88	0.54
1:A:158:ASP:OD2	1:A:161:GLN:HB2	2.08	0.53
1:A:416:LYS:HD2	1:A:417:ASN:OD1	2.09	0.53
1:A:200:GLU:HB2	1:A:203:THR:HG23	1.90	0.53
1:A:197:GLU:HG2	1:A:360:ILE:HD11	1.91	0.53
1:A:201:VAL:O	1:A:202:ASP:CB	2.57	0.53
1:A:52:ASN:HD22	1:A:55:GLU:CG	2.22	0.53
1:A:195:PHE:HZ	1:A:360:ILE:HG23	1.73	0.53
1:B:153:ILE:HG12	1:B:154:GLY:N	2.24	0.52
1:A:114:ILE:H	1:A:114:ILE:HD12	1.75	0.52
1:B:110:ILE:HG23	1:B:231:LEU:CD1	2.39	0.52
1:B:367:ASN:HD21	1:B:370:LYS:HZ1	1.58	0.52
1:B:113:GLY:HA2	1:B:319:LYS:HE3	1.92	0.52
1:B:1:MET:SD	1:B:1:MET:N	2.79	0.51
1:B:302:ILE:HD11	1:B:312:MET:HG3	1.92	0.51
1:A:303:ILE:HG22	1:A:304:GLY:N	2.26	0.51
1:A:161:GLN:HG3	3:A:604:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:MET:HE3	1:A:30:GLN:O	2.10	0.51
1:B:197:GLU:OE2	1:B:360:ILE:HD11	2.10	0.51
1:A:204:ASN:ND2	1:A:206:LEU:N	2.58	0.51
1:B:51:THR:CG2	1:B:186:ARG:HH21	2.23	0.51
1:A:201:VAL:O	1:A:202:ASP:HB3	2.10	0.50
1:B:325:ASP:OD1	1:B:327:SER:HB3	2.11	0.50
1:A:22:LYS:HA	1:A:29:MET:CE	2.40	0.50
1:B:130:ASP:OD1	1:B:131:THR:HG23	2.12	0.50
1:B:204:ASN:ND2	1:B:204:ASN:C	2.66	0.50
1:B:29:MET:O	1:B:29:MET:HG3	2.11	0.49
1:B:360:ILE:HG22	1:B:403:GLN:CD	2.33	0.49
1:A:250:TYR:HB3	1:A:254:GLY:N	2.27	0.49
1:B:63:GLU:O	1:B:64:ALA:HB3	2.12	0.49
1:B:228:GLU:HG2	1:B:344:LEU:HD22	1.93	0.49
1:A:26:ALA:CB	1:A:29:MET:HG2	2.43	0.49
1:B:410:ARG:HH11	1:B:410:ARG:CB	2.19	0.48
1:B:367:ASN:HD21	1:B:370:LYS:NZ	2.11	0.48
1:A:59:ASN:ND2	1:A:59:ASN:N	2.62	0.48
1:B:6:GLN:HE21	1:B:6:GLN:HA	1.77	0.48
1:A:171:VAL:CG2	1:A:172:LEU:HD13	2.36	0.48
1:A:274:ASP:OD1	1:A:277:GLN:NE2	2.47	0.48
1:B:166:SER:OG	1:B:183:GLN:NE2	2.46	0.48
1:A:24:PRO:O	1:A:26:ALA:N	2.47	0.48
1:B:392:LYS:O	1:B:394:ALA:N	2.47	0.47
1:A:24:PRO:O	1:A:25:ASN:C	2.52	0.47
1:A:122:ILE:HG22	1:A:124:THR:HG22	1.96	0.47
1:B:196:GLU:H	1:B:370:LYS:NZ	2.11	0.47
1:B:204:ASN:HD21	1:B:206:LEU:HB2	1.78	0.47
1:A:200:GLU:HB2	1:A:203:THR:CG2	2.43	0.47
1:A:237:ASN:CG	1:A:239:ASN:ND2	2.68	0.47
1:B:136:VAL:O	1:B:143:TYR:HA	2.15	0.47
1:B:279:ASN:O	1:B:283:LEU:HD23	2.14	0.47
1:B:325:ASP:O	1:B:327:SER:N	2.43	0.46
1:A:96:ARG:HA	1:A:385:ILE:HG23	1.97	0.46
1:B:24:PRO:O	1:B:26:ALA:N	2.48	0.46
1:B:136:VAL:HG12	1:B:137:ILE:H	1.81	0.46
1:B:225:ILE:O	1:B:228:GLU:OE2	2.32	0.46
1:A:360:ILE:HG22	1:A:403:GLN:CD	2.36	0.46
1:B:228:GLU:HG3	1:B:344:LEU:HD22	1.98	0.46
1:A:266:GLY:HA2	1:A:270:ALA:HB2	1.98	0.46
1:A:84:ASP:O	1:A:88:LYS:HD3	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:LEU:HD12	1:A:126:LEU:HA	1.76	0.45
1:A:26:ALA:HB1	1:A:29:MET:HG2	1.98	0.45
1:A:204:ASN:C	1:A:204:ASN:ND2	2.64	0.45
1:B:170:ASP:OD1	1:B:171:VAL:HG23	2.17	0.45
1:B:13:VAL:HG11	1:B:20:TYR:CE1	2.52	0.44
1:A:23:ILE:HG23	1:A:24:PRO:CD	2.46	0.44
1:B:363:LYS:HG3	3:B:580:HOH:O	2.16	0.44
1:A:221:ALA:O	1:A:225:ILE:HG13	2.17	0.44
1:A:228:GLU:OE2	1:A:348:TYR:HB2	2.18	0.44
1:B:107:LEU:O	1:B:111:VAL:HG23	2.17	0.44
1:A:125:GLU:OE2	1:A:303:ILE:HD11	2.17	0.44
1:A:305:THR:O	1:A:306:THR:OG1	2.32	0.44
1:A:180:GLY:HA2	1:A:230:ARG:O	2.18	0.43
1:B:392:LYS:C	1:B:394:ALA:H	2.22	0.43
1:B:13:VAL:O	1:B:13:VAL:HG23	2.17	0.43
1:B:130:ASP:OD1	1:B:131:THR:N	2.51	0.43
1:B:274:ASP:OD1	1:B:274:ASP:C	2.56	0.43
1:B:59:ASN:HA	1:B:60:PRO:HD3	1.85	0.43
1:B:70:SER:CA	1:B:160:ILE:HD11	2.49	0.43
1:B:101:ASP:OD1	1:B:101:ASP:N	2.49	0.43
1:B:314:ASN:O	1:B:318:GLU:HG3	2.19	0.43
1:B:360:ILE:HG22	1:B:403:GLN:OE1	2.19	0.43
1:A:197:GLU:CG	1:A:360:ILE:HD11	2.49	0.43
1:A:378:VAL:HB	1:A:379:PRO:HD3	2.00	0.43
1:B:269:ASP:O	1:B:272:PHE:HD2	2.01	0.43
1:B:317:LYS:HD2	1:B:330:PHE:HE1	1.84	0.43
1:A:128:VAL:HG22	1:A:129:ILE:N	2.34	0.42
1:B:114:ILE:HD12	1:B:315:VAL:HG13	2.01	0.42
1:A:52:ASN:HA	1:A:53:PRO:HD3	1.81	0.42
1:A:59:ASN:N	1:A:59:ASN:HD22	2.17	0.42
1:A:239:ASN:OD1	1:A:239:ASN:C	2.57	0.42
1:B:13:VAL:HG11	1:B:20:TYR:HD1	1.82	0.42
1:B:211:LYS:H	1:B:211:LYS:HG2	1.68	0.42
1:A:228:GLU:OE2	1:A:348:TYR:CD2	2.71	0.42
1:B:52:ASN:HA	1:B:53:PRO:HD3	1.90	0.42
1:B:368:PHE:O	1:B:370:LYS:NZ	2.51	0.42
1:A:84:ASP:OD1	1:A:88:LYS:HE2	2.19	0.42
1:A:396:LEU:HA	1:A:401:ASN:HB2	2.02	0.42
1:B:70:SER:HA	1:B:160:ILE:HD11	2.00	0.42
1:A:161:GLN:HA	1:A:161:GLN:HE21	1.85	0.42
1:B:143:TYR:CD2	1:B:143:TYR:N	2.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:TYR:CD1	1:A:143:TYR:N	2.87	0.42
1:A:204:ASN:HD21	1:A:206:LEU:H	1.68	0.42
1:A:199:LEU:HG	1:A:199:LEU:O	2.19	0.41
1:A:161:GLN:HA	1:A:161:GLN:NE2	2.35	0.41
1:A:52:ASN:HD22	1:A:55:GLU:CB	2.33	0.41
1:B:232:TYR:O	1:B:234:ILE:HG23	2.20	0.41
1:B:269:ASP:OD2	1:B:364:THR:HA	2.21	0.41
1:B:416:LYS:O	1:B:417:ASN:HB2	2.19	0.41
1:A:225:ILE:HD13	1:A:349:THR:HA	2.02	0.41
1:B:387:ASP:HB3	1:B:390:ASN:O	2.20	0.41
1:A:381:GLU:CD	1:A:381:GLU:N	2.74	0.41
1:B:22:LYS:HE3	1:B:137:ILE:HD11	2.02	0.41
1:B:410:ARG:HG2	3:B:627:HOH:O	2.20	0.41
1:A:208:GLY:HA3	1:A:404:ASN:ND2	2.36	0.41
1:B:225:ILE:HD13	1:B:349:THR:HA	2.02	0.41
1:B:286:TYR:CE1	1:B:334:LYS:HG2	2.56	0.41
1:A:1:MET:HB2	1:A:4:ASN:OD1	2.21	0.41
1:A:171:VAL:HG23	1:A:172:LEU:N	2.36	0.41
1:B:30:GLN:HA	1:B:30:GLN:OE1	2.20	0.41
1:A:208:GLY:HA3	3:A:576:HOH:O	2.21	0.40
1:A:135:ASN:HD22	1:A:135:ASN:HA	1.74	0.40
1:B:398:THR:O	1:B:401:ASN:ND2	2.54	0.40
1:B:69:VAL:HA	1:B:417:ASN:HD21	1.86	0.40
1:B:304:GLY:O	1:B:306:THR:N	2.41	0.40
1:B:180:GLY:HA3	1:B:231:LEU:O	2.22	0.40
1:B:367:ASN:ND2	1:B:370:LYS:NZ	2.69	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	405/425 (95%)	374 (92%)	19 (5%)	12 (3%)	4 2
1	B	403/425 (95%)	366 (91%)	26 (6%)	11 (3%)	5 3
All	All	808/850 (95%)	740 (92%)	45 (6%)	23 (3%)	5 3

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	29	MET
1	B	25	ASN
1	B	172	LEU
1	B	305	THR
1	B	392	LYS
1	A	25	ASN
1	A	202	ASP
1	A	208	GLY
1	A	209	ALA
1	A	306	THR
1	B	302	ILE
1	B	393	GLY
1	A	28	GLN
1	A	392	LYS
1	B	123	ASP
1	B	131	THR
1	A	73	ASP
1	A	210	GLY
1	A	394	ALA
1	B	51	THR
1	B	326	THR
1	B	171	VAL
1	A	61	PRO

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	367/380 (97%)	347 (95%)	20 (5%)	22 30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	366/380 (96%)	351 (96%)	15 (4%)	30	43
All	All	733/760 (96%)	698 (95%)	35 (5%)	25	36

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	130	ASP
1	A	143	TYR
1	A	170	ASP
1	A	172	LEU
1	A	197	GLU
1	A	199	LEU
1	A	202	ASP
1	A	204	ASN
1	A	228	GLU
1	A	239	ASN
1	A	276	LEU
1	A	277	GLN
1	A	283	LEU
1	A	306	THR
1	A	366	LEU
1	A	367	ASN
1	A	386	LYS
1	A	414	ARG
1	A	416	LYS
1	B	1	MET
1	B	28	GLN
1	B	50	PHE
1	B	69	VAL
1	B	80	ASP
1	B	101	ASP
1	B	139	PRO
1	B	143	TYR
1	B	204	ASN
1	B	228	GLU
1	B	245	ASN
1	B	279	ASN
1	B	364	THR
1	B	367	ASN
1	B	410	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	28	GLN
1	A	30	GLN
1	A	52	ASN
1	A	59	ASN
1	A	135	ASN
1	A	161	GLN
1	A	204	ASN
1	A	239	ASN
1	A	314	ASN
1	A	367	ASN
1	A	399	ASN
1	B	6	GLN
1	B	28	GLN
1	B	52	ASN
1	B	59	ASN
1	B	204	ASN
1	B	277	GLN
1	B	367	ASN
1	B	417	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 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	411/425 (96%)	0.41	35 (8%) 10 14	10, 26, 59, 74	0
1	B	409/425 (96%)	0.91	69 (16%) 1 2	18, 38, 68, 79	0
All	All	820/850 (96%)	0.66	104 (12%) 3 5	10, 32, 65, 79	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	25	ASN	10.3
1	A	305	THR	7.4
1	B	305	THR	7.4
1	B	253	SER	6.8
1	A	208	GLY	6.7
1	B	143	TYR	6.5
1	B	62	PRO	6.4
1	A	27	GLY	6.4
1	A	62	PRO	6.3
1	B	64	ALA	6.2
1	B	65	LYS	6.2
1	A	25	ASN	5.9
1	A	248	ALA	5.8
1	B	307	ALA	5.7
1	A	61	PRO	5.5
1	B	28	GLN	5.5
1	A	199	LEU	5.4
1	B	311	TYR	5.2
1	B	59	ASN	4.8
1	A	28	GLN	4.8
1	A	249	TYR	4.6
1	B	308	SER	4.5
1	B	139	PRO	4.5
1	B	417	ASN	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	207	LEU	4.4
1	B	121	THR	4.3
1	A	209	ALA	4.2
1	A	417	ASN	4.2
1	B	326	THR	4.2
1	B	53	PRO	4.2
1	B	247	ASN	4.1
1	B	168	GLY	4.1
1	B	122	ILE	4.1
1	B	130	ASP	4.0
1	A	143	TYR	3.9
1	B	142	SER	3.9
1	B	327	SER	3.8
1	A	326	THR	3.8
1	B	303	ILE	3.8
1	B	304	GLY	3.8
1	B	254	GLY	3.7
1	B	199	LEU	3.7
1	B	397	SER	3.7
1	B	27	GLY	3.7
1	B	325	ASP	3.6
1	B	141	GLY	3.6
1	B	138	GLN	3.6
1	A	63	GLU	3.5
1	A	398	THR	3.5
1	B	54	GLU	3.5
1	B	398	THR	3.5
1	B	4	ASN	3.5
1	A	60	PRO	3.3
1	B	126	LEU	3.3
1	B	306	THR	3.3
1	B	26	ALA	3.3
1	B	123	ASP	3.3
1	B	170	ASP	3.2
1	B	63	GLU	3.2
1	B	61	PRO	3.2
1	A	206	LEU	3.2
1	A	59	ASN	3.1
1	A	369	ASP	3.1
1	B	6	GLN	3.1
1	B	169	HIS	3.0
1	B	140	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	254	GLY	3.0
1	B	302	ILE	3.0
1	A	250	TYR	3.0
1	B	312	MET	2.9
1	A	251	GLU	2.9
1	A	210	GLY	2.9
1	B	5	LYS	2.8
1	B	198	SER	2.7
1	B	30	GLN	2.7
1	A	366	LEU	2.7
1	B	335	LEU	2.7
1	A	368	PHE	2.6
1	B	294	SER	2.6
1	A	54	GLU	2.6
1	A	6	GLN	2.6
1	B	331	SER	2.5
1	B	60	PRO	2.4
1	A	30	GLN	2.4
1	B	301	SER	2.4
1	A	170	ASP	2.4
1	B	310	GLN	2.3
1	A	101	ASP	2.3
1	A	367	ASN	2.3
1	B	44	ILE	2.3
1	B	291	ASP	2.3
1	B	320	TYR	2.3
1	B	171	VAL	2.2
1	B	234	ILE	2.2
1	A	255	LEU	2.2
1	B	292	VAL	2.2
1	A	26	ALA	2.2
1	B	16	VAL	2.2
1	B	336	LYS	2.2
1	B	316	PHE	2.1
1	B	328	GLY	2.1
1	B	29	MET	2.0
1	B	57	ASP	2.0
1	B	381	GLU	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	B	451	1/1	0.98	0.04	31,31,31,31	0
2	ZN	A	452	1/1	0.99	0.06	25,25,25,25	0

6.5 Other polymers (i)

There are no such residues in this entry.