



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:13 AM GMT

PDB ID : 2G7K
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.; Crawford, A.; Janda, K.D.; Barbieri, J.T.; Kim, J.-J.P.
Deposited on : 2006-02-28
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

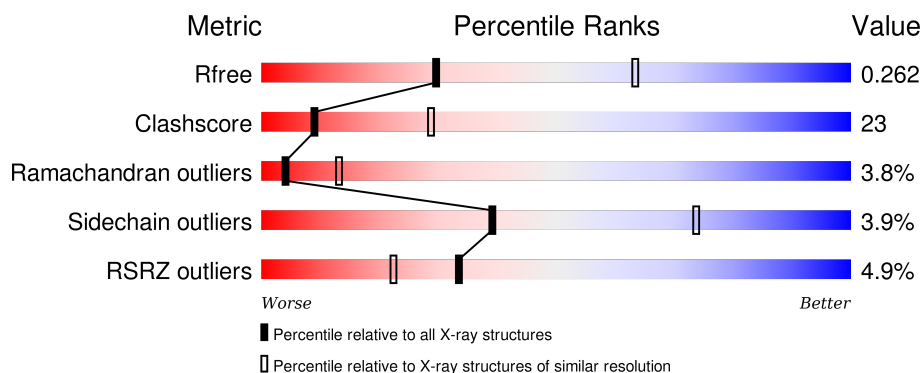
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	425	<div> <div>2%</div> <div>61%</div> <div>29%</div> <div>6%</div> </div>
1	B	425	<div> <div>7%</div> <div>53%</div> <div>36%</div> <div>6%</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6481 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
1	A	401	Total	C	N	O	S	0	0	0
			3247	2084	536	620	7			
1	B	399	Total	C	N	O	S	0	0	0
			3234	2076	534	617	7			

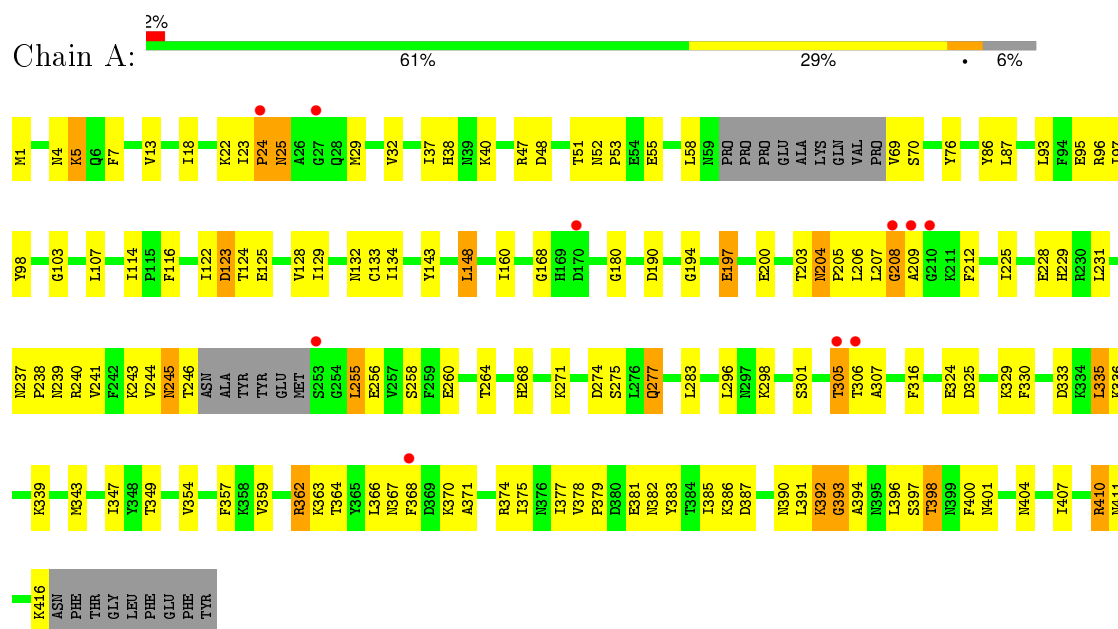
There are 2 discrepancies between the modelled and reference sequences:

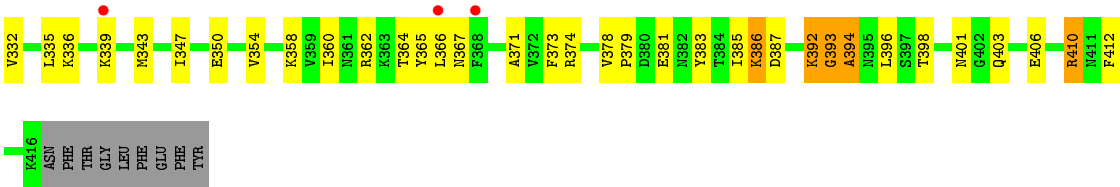
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP Q45894
B	1	MET	-	INITIATING METHIONINE	UNP Q45894

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.10Å 110.09Å 167.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.98 – 2.80 29.97 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.9 (29.98-2.80) 94.8 (29.97-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.80Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.223 , 0.265 0.223 , 0.262	Depositor DCC
R_{free} test set	3059 reflections (10.03%)	DCC
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30513 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6481	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.45	0/3317	0.67	0/4482
1	B	0.43	0/3304	0.61	0/4464
All	All	0.44	0/6621	0.64	0/8946

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

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	3247	0	3192	129	0
1	B	3234	0	3178	174	0
All	All	6481	0	6370	298	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 23.

All (298) 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:410:ARG:HB2	1:A:410:ARG:HH11	1.05	1.08
1:A:387:ASP:OD2	1:A:392:LYS:HG3	1.55	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:ARG:NH1	1:A:410:ARG:HB2	1.74	1.02
1:B:5:LYS:HD3	1:B:7:PHE:HE2	1.29	0.98
1:B:410:ARG:HH11	1:B:410:ARG:HB2	1.27	0.96
1:A:410:ARG:CB	1:A:410:ARG:HH11	1.80	0.95
1:A:343:MET:HA	1:A:347:ILE:HD12	1.44	0.95
1:A:225:ILE:O	1:A:228:GLU:HG2	1.66	0.94
1:A:209:ALA:H	1:A:404:ASN:HD21	1.18	0.89
1:B:265:PHE:HE2	1:B:362:ARG:HH12	1.20	0.88
1:B:122:ILE:HG22	1:B:124:THR:HG22	1.56	0.87
1:B:5:LYS:HD3	1:B:7:PHE:CE2	2.10	0.87
1:B:117:TRP:CE2	1:B:128:VAL:HB	2.11	0.85
1:B:362:ARG:HB2	1:B:362:ARG:HH11	1.42	0.84
1:B:374:ARG:HH11	1:B:374:ARG:HG3	1.44	0.82
1:B:5:LYS:CD	1:B:7:PHE:HE2	1.92	0.82
1:B:387:ASP:OD2	1:B:392:LYS:HG3	1.79	0.81
1:B:410:ARG:NH1	1:B:410:ARG:HB2	1.97	0.79
1:B:325:ASP:OD2	1:B:329:LYS:HB3	1.83	0.79
1:B:204:ASN:ND2	1:B:206:LEU:H	1.81	0.79
1:A:228:GLU:HG3	1:A:229:HIS:N	1.98	0.77
1:B:362:ARG:HH11	1:B:362:ARG:CB	1.96	0.77
1:B:228:GLU:HG3	1:B:229:HIS:H	1.51	0.76
1:A:391:LEU:O	1:A:397:SER:HB3	1.85	0.76
1:A:275:SER:HB2	1:B:243:LYS:HG2	1.66	0.76
1:B:362:ARG:NH1	1:B:362:ARG:HB2	2.01	0.75
1:B:5:LYS:CB	1:B:7:PHE:CE2	2.71	0.73
1:B:300:LYS:O	1:B:309:LEU:HD22	1.88	0.73
1:B:181:SER:O	1:B:231:LEU:HD23	1.89	0.73
1:A:268:HIS:CE1	1:B:268:HIS:HB3	2.24	0.73
1:B:255:LEU:HD12	1:B:255:LEU:O	1.89	0.72
1:B:5:LYS:HB2	1:B:7:PHE:CE2	2.24	0.71
1:A:70:SER:HA	1:A:160:ILE:HD11	1.72	0.71
1:B:5:LYS:HB2	1:B:7:PHE:CZ	2.26	0.70
1:B:122:ILE:CG2	1:B:124:THR:HG22	2.21	0.70
1:A:305:THR:C	1:A:307:ALA:H	1.95	0.70
1:A:397:SER:O	1:A:398:THR:HG23	1.92	0.70
1:A:209:ALA:N	1:A:404:ASN:HD21	1.90	0.70
1:B:172:LEU:H	1:B:172:LEU:HD23	1.56	0.70
1:A:239:ASN:HD21	1:A:240:ARG:HH12	1.39	0.69
1:B:78:SER:O	1:B:83:LYS:HE3	1.94	0.68
1:A:324:GLU:HB2	1:A:330:PHE:CE1	2.29	0.68
1:B:116:PHE:HA	1:B:316:PHE:CE1	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:ILE:HG23	1:B:148:LEU:HD21	1.76	0.67
1:B:266:GLY:HA2	1:B:270:ALA:HB2	1.75	0.67
1:B:204:ASN:C	1:B:204:ASN:HD22	1.97	0.67
1:A:245:ASN:HD22	1:A:246:THR:N	1.93	0.67
1:B:169:HIS:CD2	1:B:171:VAL:H	2.13	0.66
1:B:96:ARG:HA	1:B:385:ILE:HG23	1.78	0.66
1:A:382:ASN:HB3	1:A:411:ASN:OD1	1.95	0.66
1:A:225:ILE:O	1:A:228:GLU:CG	2.44	0.65
1:A:4:ASN:O	1:A:5:LYS:HB3	1.97	0.64
1:A:335:LEU:HD12	1:A:339:LYS:HE3	1.79	0.64
1:A:305:THR:O	1:A:307:ALA:N	2.31	0.64
1:A:354:VAL:HG12	1:B:201:VAL:HG21	1.80	0.64
1:B:312:MET:O	1:B:315:VAL:HG22	1.98	0.64
1:B:204:ASN:HD22	1:B:205:PRO:N	1.97	0.63
1:B:124:THR:HA	1:B:299:ALA:HA	1.78	0.63
1:B:228:GLU:HG3	1:B:229:HIS:N	2.13	0.62
1:A:335:LEU:O	1:A:339:LYS:HG3	1.99	0.62
1:A:239:ASN:HD21	1:A:240:ARG:NH1	1.98	0.62
1:B:130:ASP:OD2	1:B:130:ASP:N	2.30	0.62
1:B:396:LEU:HA	1:B:401:ASN:HB2	1.81	0.62
1:A:305:THR:C	1:A:307:ALA:N	2.51	0.61
1:A:204:ASN:HD22	1:A:205:PRO:N	1.98	0.61
1:B:125:GLU:HA	1:B:301:SER:O	2.02	0.60
1:A:225:ILE:CG2	1:A:264:THR:HG23	2.32	0.60
1:B:315:VAL:CG2	1:B:316:PHE:N	2.64	0.60
1:B:315:VAL:HG23	1:B:316:PHE:N	2.16	0.60
1:A:392:LYS:O	1:A:394:ALA:N	2.34	0.60
1:A:197:GLU:OE2	1:A:207:LEU:HB2	2.02	0.59
1:B:225:ILE:O	1:B:228:GLU:HG3	2.03	0.59
1:A:325:ASP:OD2	1:A:329:LYS:HB3	2.03	0.59
1:B:34:ALA:HB2	1:B:44:ILE:HG12	1.84	0.59
1:A:204:ASN:ND2	1:A:206:LEU:H	2.01	0.59
1:A:22:LYS:O	1:A:24:PRO:HD3	2.02	0.59
1:A:122:ILE:HG22	1:A:124:THR:HG22	1.83	0.59
1:A:47:ARG:CZ	1:A:58:LEU:HD13	2.33	0.58
1:B:50:PHE:O	1:B:52:ASN:N	2.37	0.58
1:B:224:LEU:O	1:B:227:ALA:HB3	2.03	0.58
1:A:228:GLU:HG3	1:A:229:HIS:H	1.69	0.58
1:B:127:LYS:HG2	1:B:128:VAL:N	2.18	0.58
1:B:106:LEU:O	1:B:110:ILE:HG13	2.04	0.57
1:A:204:ASN:HD22	1:A:205:PRO:CD	2.17	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:392:LYS:O	1:B:394:ALA:N	2.37	0.57
1:A:241:VAL:HG12	1:A:258:SER:HA	1.86	0.57
1:B:161:GLN:HG3	1:B:161:GLN:O	2.04	0.57
1:A:400:PHE:CZ	1:B:201:VAL:HG12	2.40	0.56
1:B:204:ASN:HB3	1:B:207:LEU:HD13	1.87	0.56
1:B:302:ILE:HB	1:B:309:LEU:HD12	1.87	0.56
1:B:358:LYS:NZ	1:B:358:LYS:HB2	2.20	0.56
1:B:22:LYS:HB2	1:B:135:ASN:HB2	1.87	0.56
1:A:13:VAL:HA	1:A:18:ILE:HG22	1.86	0.56
1:B:225:ILE:O	1:B:228:GLU:CG	2.54	0.56
1:B:101:ASP:HA	1:B:104:ARG:NH1	2.21	0.56
1:B:23:ILE:HG22	1:B:26:ALA:HB2	1.87	0.56
1:B:5:LYS:CD	1:B:7:PHE:CE2	2.78	0.56
1:B:5:LYS:HB3	1:B:7:PHE:CE2	2.40	0.56
1:A:324:GLU:HB2	1:A:330:PHE:CD1	2.41	0.56
1:A:134:ILE:HD13	1:A:148:LEU:HD13	1.88	0.55
1:B:169:HIS:CG	1:B:170:ASP:N	2.74	0.55
1:A:40:LYS:HE3	1:A:114:ILE:HD11	1.89	0.55
1:B:374:ARG:HH11	1:B:374:ARG:CG	2.18	0.55
1:A:204:ASN:C	1:A:204:ASN:HD22	2.10	0.55
1:A:69:VAL:HG23	1:A:416:LYS:HD3	1.87	0.55
1:B:134:ILE:HD13	1:B:148:LEU:HD13	1.89	0.55
1:B:266:GLY:CA	1:B:270:ALA:HB2	2.36	0.54
1:A:69:VAL:CG2	1:A:416:LYS:HD3	2.38	0.54
1:A:204:ASN:HD22	1:A:205:PRO:HD2	1.72	0.54
1:A:245:ASN:HD22	1:A:245:ASN:C	2.08	0.54
1:A:194:GLY:HA2	1:A:212:PHE:O	2.07	0.54
1:A:200:GLU:HB2	1:A:203:THR:HG23	1.90	0.54
1:A:93:LEU:HD21	1:A:377:ILE:HD13	1.89	0.53
1:B:343:MET:HA	1:B:347:ILE:HD12	1.89	0.53
1:B:7:PHE:HE1	1:B:37:ILE:HA	1.74	0.53
1:A:396:LEU:HA	1:A:401:ASN:HB2	1.91	0.53
1:B:406:GLU:OE1	1:B:406:GLU:HA	2.09	0.53
1:B:373:PHE:CD2	1:B:412:PHE:HB3	2.44	0.53
1:A:255:LEU:HD23	1:A:256:GLU:N	2.24	0.53
1:B:120:SER:OG	1:B:125:GLU:HB2	2.08	0.53
1:B:371:ALA:HB1	1:B:373:PHE:CE1	2.44	0.53
1:B:386:LYS:HB2	1:B:386:LYS:NZ	2.24	0.53
1:B:225:ILE:HA	1:B:228:GLU:HG2	1.90	0.53
1:B:190:ASP:O	1:B:374:ARG:HD2	2.08	0.52
1:B:169:HIS:CD2	1:B:172:LEU:HG	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:374:ARG:NH1	1:B:374:ARG:HG3	2.19	0.52
1:A:229:HIS:ND1	1:A:260:GLU:OE2	2.41	0.52
1:B:197:GLU:OE2	1:B:197:GLU:HA	2.10	0.52
1:A:359:VAL:O	1:B:201:VAL:HG11	2.09	0.52
1:A:343:MET:CA	1:A:347:ILE:HD12	2.31	0.52
1:A:357:PHE:HB3	1:A:359:VAL:HG13	1.92	0.52
1:B:50:PHE:C	1:B:52:ASN:H	2.13	0.52
1:B:17:ASP:OD1	1:B:18:ILE:HG13	2.10	0.52
1:B:290:LYS:HG2	1:B:332:VAL:HB	1.92	0.52
1:B:124:THR:CA	1:B:299:ALA:HA	2.40	0.51
1:B:101:ASP:HA	1:B:104:ARG:HH11	1.76	0.51
1:B:204:ASN:HD22	1:B:206:LEU:H	1.57	0.50
1:A:208:GLY:O	1:A:209:ALA:HB2	2.11	0.50
1:B:392:LYS:O	1:B:393:GLY:C	2.49	0.50
1:B:108:THR:O	1:B:112:ARG:HG3	2.10	0.50
1:B:387:ASP:OD2	1:B:392:LYS:CG	2.54	0.50
1:A:240:ARG:HG2	1:A:240:ARG:HH11	1.77	0.50
1:A:245:ASN:ND2	1:A:245:ASN:C	2.65	0.50
1:B:360:ILE:HG22	1:B:403:GLN:CD	2.32	0.49
1:B:50:PHE:C	1:B:52:ASN:N	2.66	0.49
1:A:129:ILE:HB	1:A:132:ASN:ND2	2.28	0.49
1:B:125:GLU:OE2	1:B:303:ILE:HD11	2.13	0.49
1:B:104:ARG:HB3	1:B:104:ARG:NH1	2.27	0.49
1:B:298:LYS:O	1:B:299:ALA:O	2.31	0.49
1:B:169:HIS:CG	1:B:170:ASP:H	2.30	0.49
1:B:169:HIS:HD2	1:B:171:VAL:H	1.57	0.49
1:B:204:ASN:HD22	1:B:205:PRO:CD	2.26	0.48
1:B:410:ARG:HH11	1:B:410:ARG:CB	2.12	0.48
1:A:23:ILE:O	1:A:24:PRO:C	2.50	0.48
1:B:378:VAL:HB	1:B:379:PRO:HD3	1.95	0.48
1:B:374:ARG:NH1	1:B:374:ARG:CG	2.76	0.48
1:B:24:PRO:CG	1:B:133:CYS:O	2.61	0.48
1:A:404:ASN:HB3	1:A:407:ILE:HB	1.95	0.48
1:B:34:ALA:CB	1:B:44:ILE:HG12	2.43	0.48
1:A:392:LYS:O	1:A:393:GLY:C	2.51	0.48
1:B:148:LEU:H	1:B:148:LEU:HD23	1.79	0.48
1:B:24:PRO:HG2	1:B:133:CYS:O	2.13	0.48
1:B:36:LYS:HB2	1:B:42:TRP:CE2	2.48	0.48
1:A:52:ASN:HD22	1:A:55:GLU:HG3	1.79	0.48
1:A:387:ASP:OD2	1:A:392:LYS:CG	2.44	0.48
1:B:47:ARG:CZ	1:B:58:LEU:HD13	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:ILE:O	1:A:103:GLY:HA3	2.14	0.48
1:A:381:GLU:N	1:A:381:GLU:CD	2.67	0.47
1:B:255:LEU:HD11	1:B:366:LEU:CD1	2.44	0.47
1:B:398:THR:O	1:B:401:ASN:ND2	2.47	0.47
1:A:37:ILE:O	1:A:38:HIS:HB2	2.15	0.47
1:B:101:ASP:N	1:B:101:ASP:OD1	2.41	0.47
1:B:302:ILE:HD11	1:B:312:MET:HG3	1.96	0.47
1:B:315:VAL:CG2	1:B:316:PHE:H	2.27	0.47
1:B:16:VAL:HG12	1:B:17:ASP:N	2.28	0.47
1:A:190:ASP:O	1:A:374:ARG:HD3	2.15	0.47
1:A:116:PHE:HE1	1:A:296:LEU:HD21	1.79	0.47
1:A:324:GLU:HB2	1:A:330:PHE:HE1	1.78	0.47
1:B:134:ILE:HG23	1:B:148:LEU:CD2	2.44	0.47
1:B:148:LEU:N	1:B:148:LEU:HD23	2.30	0.47
1:A:375:ILE:HD12	1:A:377:ILE:HG23	1.95	0.47
1:A:225:ILE:HG21	1:A:264:THR:HG23	1.96	0.47
1:B:13:VAL:HG11	1:B:20:TYR:HD1	1.78	0.47
1:A:96:ARG:HA	1:A:385:ILE:HG23	1.97	0.47
1:B:96:ARG:CA	1:B:385:ILE:HG23	2.44	0.47
1:B:265:PHE:CE2	1:B:365:TYR:HA	2.50	0.46
1:B:124:THR:O	1:B:299:ALA:HA	2.14	0.46
1:B:36:LYS:HE2	1:B:39:ASN:O	2.15	0.46
1:B:350:GLU:O	1:B:354:VAL:HG23	2.15	0.46
1:B:180:GLY:HA3	1:B:231:LEU:O	2.15	0.46
1:A:24:PRO:HD2	1:A:133:CYS:O	2.16	0.46
1:B:304:GLY:O	1:B:306:THR:N	2.48	0.46
1:A:7:PHE:HB3	1:A:87:LEU:HD21	1.98	0.46
1:B:288:LYS:O	1:B:292:VAL:HG23	2.15	0.46
1:A:4:ASN:O	1:A:5:LYS:CB	2.63	0.46
1:A:134:ILE:CD1	1:A:148:LEU:HD13	2.44	0.46
1:A:255:LEU:HD23	1:A:256:GLU:H	1.81	0.46
1:B:98:TYR:HD1	1:B:107:LEU:HD12	1.81	0.46
1:A:116:PHE:HA	1:A:316:PHE:CE1	2.50	0.46
1:B:3:VAL:HG13	1:B:37:ILE:HB	1.96	0.46
1:B:293:ALA:O	1:B:297:ASN:ND2	2.49	0.46
1:A:48:ASP:OD1	1:A:51:THR:HG23	2.15	0.46
1:B:279:ASN:O	1:B:280:GLU:C	2.53	0.46
1:A:225:ILE:C	1:A:228:GLU:HG2	2.34	0.46
1:B:204:ASN:C	1:B:204:ASN:ND2	2.65	0.46
1:A:52:ASN:ND2	1:A:55:GLU:HG3	2.31	0.46
1:B:176:ARG:HD2	1:B:237:ASN:HA	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:GLU:CG	1:A:229:HIS:N	2.75	0.45
1:B:265:PHE:CE2	1:B:362:ARG:NH1	2.79	0.45
1:A:1:MET:HA	1:A:4:ASN:OD1	2.16	0.45
1:B:371:ALA:HB1	1:B:373:PHE:HE1	1.80	0.45
1:A:378:VAL:HB	1:A:379:PRO:HD3	1.97	0.45
1:A:24:PRO:O	1:A:25:ASN:HB2	2.16	0.45
1:A:381:GLU:CD	1:A:381:GLU:H	2.19	0.45
1:B:100:THR:O	1:B:104:ARG:N	2.42	0.45
1:A:208:GLY:HA3	1:A:404:ASN:ND2	2.32	0.45
1:B:324:GLU:HA	1:B:329:LYS:O	2.16	0.45
1:A:1:MET:O	1:A:4:ASN:OD1	2.34	0.45
1:B:302:ILE:HB	1:B:309:LEU:CD1	2.46	0.45
1:A:47:ARG:NH1	1:A:76:TYR:HB3	2.32	0.45
1:A:239:ASN:ND2	1:A:240:ARG:NH1	2.65	0.44
1:A:370:LYS:NZ	1:A:370:LYS:HB2	2.32	0.44
1:A:237:ASN:OD1	1:A:238:PRO:HD2	2.18	0.44
1:B:124:THR:C	1:B:299:ALA:HA	2.38	0.44
1:A:52:ASN:HA	1:A:53:PRO:HD3	1.85	0.44
1:A:277:GLN:HB2	1:A:277:GLN:HE21	1.65	0.44
1:A:392:LYS:C	1:A:394:ALA:N	2.71	0.44
1:B:204:ASN:HD21	1:B:206:LEU:HB2	1.82	0.44
1:A:122:ILE:CG2	1:A:124:THR:HG22	2.47	0.44
1:A:363:LYS:HA	1:A:363:LYS:HE2	2.00	0.44
1:B:255:LEU:HD12	1:B:255:LEU:C	2.37	0.44
1:B:79:THR:O	1:B:83:LYS:HG3	2.18	0.44
1:B:225:ILE:CG2	1:B:264:THR:HG23	2.48	0.43
1:B:2:PHE:HB2	1:B:98:TYR:CG	2.53	0.43
1:A:335:LEU:CD1	1:A:339:LYS:HE3	2.47	0.43
1:B:29:MET:HG3	1:B:29:MET:O	2.18	0.43
1:A:225:ILE:HD13	1:A:349:THR:HA	2.00	0.43
1:B:128:VAL:HG22	1:B:129:ILE:N	2.33	0.43
1:A:362:ARG:HG3	1:A:362:ARG:HH11	1.83	0.43
1:B:321:LEU:O	1:B:336:LYS:HE3	2.19	0.43
1:B:362:ARG:HE	1:B:367:ASN:HB2	1.84	0.43
1:B:36:LYS:HB2	1:B:42:TRP:CZ2	2.54	0.43
1:B:54:GLU:H	1:B:54:GLU:CD	2.21	0.43
1:A:274:ASP:OD1	1:A:274:ASP:C	2.57	0.43
1:B:131:THR:HB	1:B:167:PHE:HB2	2.00	0.43
1:B:197:GLU:OE2	1:B:360:ILE:HD11	2.18	0.42
1:A:244:VAL:HG22	1:A:255:LEU:O	2.19	0.42
1:B:360:ILE:HG22	1:B:403:GLN:OE1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:SER:HB3	1:B:125:GLU:O	2.19	0.42
1:B:232:TYR:O	1:B:234:ILE:HG23	2.19	0.42
1:A:180:GLY:HA3	1:A:231:LEU:O	2.19	0.42
1:B:48:ASP:HA	1:B:153:ILE:HD11	2.02	0.42
1:B:138:GLN:HB3	1:B:139:PRO:HD2	2.02	0.42
1:B:96:ARG:HG3	1:B:385:ILE:O	2.19	0.42
1:A:243:LYS:HE3	1:A:256:GLU:OE2	2.20	0.42
1:A:143:TYR:C	1:A:143:TYR:CD1	2.92	0.42
1:A:29:MET:HB2	1:A:29:MET:HE3	1.92	0.42
1:A:125:GLU:HG3	1:A:301:SER:OG	2.19	0.42
1:A:225:ILE:HG23	1:A:228:GLU:OE2	2.19	0.42
1:A:271:LYS:HD3	1:A:271:LYS:HA	1.70	0.42
1:A:386:LYS:HE2	1:A:386:LYS:HB2	1.75	0.42
1:B:71:TYR:CE2	1:B:159:ILE:HD12	2.55	0.42
1:B:92:LYS:NZ	1:B:383:TYR:O	2.48	0.42
1:B:132:ASN:HA	1:B:182:THR:OG1	2.19	0.42
1:A:123:ASP:OD2	1:A:298:LYS:HE2	2.20	0.42
1:A:392:LYS:O	1:A:394:ALA:O	2.38	0.42
1:A:204:ASN:ND2	1:A:206:LEU:N	2.65	0.42
1:A:387:ASP:HB3	1:A:390:ASN:O	2.20	0.41
1:B:302:ILE:HG21	1:B:309:LEU:CA	2.50	0.41
1:A:204:ASN:HD21	1:A:206:LEU:HB3	1.85	0.41
1:A:95:GLU:HA	1:A:95:GLU:OE1	2.19	0.41
1:A:333:ASP:OD2	1:A:336:LYS:HG3	2.19	0.41
1:B:274:ASP:H	1:B:277:GLN:HE21	1.67	0.41
1:B:22:LYS:HB3	1:B:22:LYS:HE3	1.89	0.41
1:B:15:GLY:HA3	1:B:139:PRO:HD3	2.02	0.41
1:B:3:VAL:HG21	1:B:91:THR:HG23	2.02	0.41
1:B:302:ILE:HG22	1:B:309:LEU:HB2	2.01	0.41
1:B:96:ARG:HD3	1:B:383:TYR:OH	2.20	0.41
1:A:204:ASN:ND2	1:A:205:PRO:HD2	2.35	0.41
1:A:362:ARG:HG3	1:A:362:ARG:NH1	2.35	0.41
1:B:123:ASP:O	1:B:123:ASP:OD2	2.38	0.41
1:A:377:ILE:HG22	1:A:383:TYR:CD2	2.56	0.41
1:A:98:TYR:HD1	1:A:107:LEU:HD12	1.86	0.41
1:B:143:TYR:CD1	1:B:143:TYR:O	2.74	0.41
1:B:302:ILE:HG12	1:B:302:ILE:O	2.19	0.41
1:B:211:LYS:HD3	1:B:371:ALA:HB2	2.03	0.41
1:B:114:ILE:O	1:B:115:PRO:C	2.60	0.41
1:A:204:ASN:HD21	1:A:206:LEU:CB	2.34	0.40
1:A:364:THR:HG23	1:A:366:LEU:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:ILE:O	1:B:228:GLU:HG2	2.21	0.40
1:B:197:GLU:HG2	1:B:207:LEU:HD23	2.02	0.40
1:B:358:LYS:NZ	1:B:358:LYS:CB	2.83	0.40
1:B:47:ARG:NH1	1:B:58:LEU:HD13	2.36	0.40
1:B:335:LEU:HG	1:B:339:LYS:HE3	2.03	0.40
1:A:24:PRO:O	1:A:25:ASN:CB	2.70	0.40
1:A:86:TYR:C	1:A:86:TYR:CD1	2.93	0.40
1:B:35:PHE:CD1	1:B:35:PHE:N	2.89	0.40
1:B:242:PHE:HB2	1:B:259:PHE:HE1	1.87	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	395/425 (93%)	347 (88%)	36 (9%)	12 (3%)	5	18
1	B	393/425 (92%)	339 (86%)	36 (9%)	18 (5%)	3	9
All	All	788/850 (93%)	686 (87%)	72 (9%)	30 (4%)	4	13

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	305	THR
1	A	392	LYS
1	B	299	ALA
1	B	305	THR
1	B	392	LYS
1	A	25	ASN
1	A	168	GLY
1	A	197	GLU

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Mol	Chain	Res	Type
1	A	208	GLY
1	A	393	GLY
1	B	51	THR
1	B	199	LEU
1	B	393	GLY
1	A	123	ASP
1	A	306	THR
1	A	368	PHE
1	A	371	ALA
1	B	25	ASN
1	B	131	THR
1	B	156	SER
1	B	73	ASP
1	B	130	ASP
1	B	159	ILE
1	B	394	ALA
1	B	28	GLN
1	B	45	PRO
1	B	207	LEU
1	B	304	GLY
1	B	171	VAL

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	360/381 (94%)	346 (96%)	14 (4%)	39	74
1	B	358/381 (94%)	344 (96%)	14 (4%)	39	74
All	All	718/762 (94%)	690 (96%)	28 (4%)	39	74

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	PRO
1	A	32	VAL

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Mol	Chain	Res	Type
1	A	128	VAL
1	A	148	LEU
1	A	204	ASN
1	A	245	ASN
1	A	255	LEU
1	A	277	GLN
1	A	283	LEU
1	A	335	LEU
1	A	362	ARG
1	A	367	ASN
1	A	398	THR
1	A	410	ARG
1	B	6	GLN
1	B	28	GLN
1	B	50	PHE
1	B	101	ASP
1	B	130	ASP
1	B	148	LEU
1	B	172	LEU
1	B	204	ASN
1	B	243	LYS
1	B	310	GLN
1	B	364	THR
1	B	381	GLU
1	B	386	LYS
1	B	410	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	59	ASN
1	A	135	ASN
1	A	204	ASN
1	A	239	ASN
1	A	245	ASN
1	A	268	HIS
1	A	277	GLN
1	A	352	ASN
1	A	404	ASN
1	B	6	GLN
1	B	28	GLN

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Mol	Chain	Res	Type
1	B	59	ASN
1	B	138	GLN
1	B	169	HIS
1	B	204	ASN
1	B	277	GLN
1	B	287	ASN
1	B	355	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 ⓘ

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	401/425 (94%)	-0.31	10 (2%) 61 48	9, 25, 70, 89	0
1	B	399/425 (93%)	0.15	29 (7%) 18 10	16, 44, 85, 95	0
All	All	800/850 (94%)	-0.08	39 (4%) 33 22	9, 36, 81, 95	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	368	PHE	5.3
1	B	143	TYR	5.2
1	B	305	THR	5.0
1	B	142	SER	4.3
1	A	305	THR	4.0
1	B	123	ASP	3.7
1	B	28	GLN	3.7
1	B	25	ASN	3.5
1	B	26	ALA	3.5
1	B	311	TYR	3.5
1	A	208	GLY	3.3
1	A	170	ASP	3.3
1	B	141	GLY	3.3
1	A	27	GLY	3.2
1	B	368	PHE	3.1
1	A	306	THR	2.9
1	B	171	VAL	2.9
1	A	253	SER	2.9
1	B	121	THR	2.8
1	A	209	ALA	2.8
1	B	137	ILE	2.8
1	A	24	PRO	2.8
1	B	54	GLU	2.6
1	B	139	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	210	GLY	2.4
1	B	307	ALA	2.4
1	B	170	ASP	2.4
1	B	326	THR	2.3
1	B	304	GLY	2.3
1	B	56	GLY	2.3
1	B	57	ASP	2.2
1	B	6	GLN	2.2
1	B	327	SER	2.2
1	B	339	LYS	2.2
1	B	59	ASN	2.1
1	B	366	LEU	2.1
1	B	301	SER	2.1
1	B	308	SER	2.0
1	B	138	GLN	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.