



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:21 PM UTC

PDB ID : 2G7Q / pdb_00002g7q
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 : 2.41 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

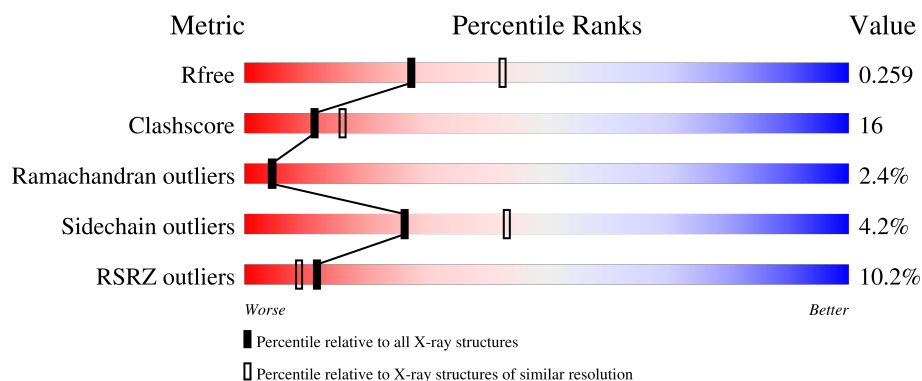
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	425	8% 70% 22% . .
1	B	425	12% 61% 30% 5% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6797 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	411	Total	C	N	O	S	0	0	0
			3323	2135	545	636	7			
1	B	409	Total	C	N	O	S	0	0	0
			3300	2118	544	631	7			

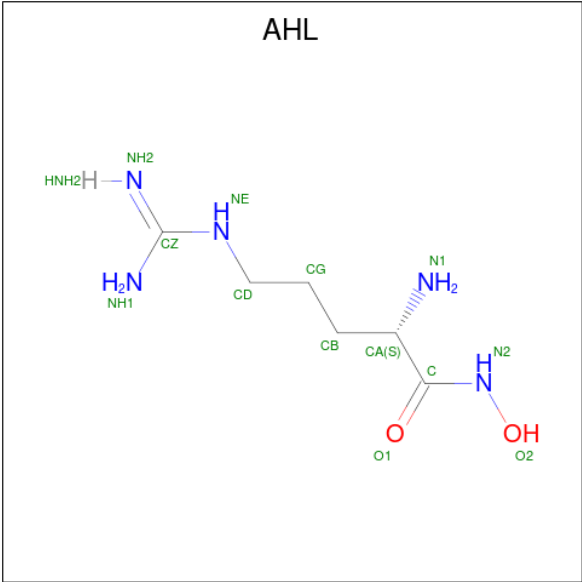
There are 6 discrepancies between the modelled and reference sequences:

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 (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is N-HYDROXY-L-ARGININAMIDE (CCD ID: AHL) (formula: C₆H₁₅N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	6	5	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	104	Total	O	0	0
			104	104		
4	B	55	Total	O	0	0
			55	55		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.32Å 109.94Å 166.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.55 – 2.41 21.55 – 2.41	Depositor EDS
% Data completeness (in resolution range)	98.5 (21.55-2.41) 98.7 (21.55-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.17 (at 2.41Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.261 0.223 , 0.259	Depositor DCC
R_{free} test set	4984 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6797	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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, AHL

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	0.47	0/3398	0.96	18/4597 (0.4%)
1	B	0.41	0/3373	0.90	13/4561 (0.3%)
All	All	0.44	0/6771	0.93	31/9158 (0.3%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	ALA	N-CA-C	-8.69	102.45	113.23
1	A	304	GLY	N-CA-C	-7.31	102.33	114.76
1	B	200	GLU	N-CA-C	-7.31	97.33	109.46
1	A	305	THR	N-CA-C	6.84	119.93	110.35
1	B	391	LEU	N-CA-C	6.65	120.00	110.10
1	B	199	LEU	N-CA-C	6.57	120.32	107.57
1	A	167	PHE	N-CA-C	6.54	118.20	111.14
1	A	174	LEU	N-CA-C	6.41	119.08	111.33
1	A	371	ALA	N-CA-C	6.19	117.97	109.18
1	B	11	ASP	CA-C-N	6.14	126.10	119.78
1	B	11	ASP	C-N-CA	6.14	126.10	119.78
1	B	368	PHE	N-CA-C	6.11	118.65	110.35
1	A	317	LYS	N-CA-C	-6.05	104.76	111.36
1	A	286	TYR	N-CA-C	-5.91	104.74	111.07
1	A	30	GLN	N-CA-C	-5.91	102.14	110.29
1	A	47	ARG	N-CA-C	-5.79	101.71	110.28
1	B	26	ALA	N-CA-C	-5.68	106.01	113.17
1	B	174	LEU	N-CA-C	5.59	118.91	111.75
1	A	395	ASN	N-CA-C	-5.55	106.17	113.17
1	B	52	ASN	CA-C-N	5.51	126.72	119.84
1	B	52	ASN	C-N-CA	5.51	126.72	119.84
1	A	391	LEU	N-CA-C	5.50	118.33	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	52	ASN	CA-C-N	5.36	124.54	118.97
1	A	52	ASN	C-N-CA	5.36	124.54	118.97
1	A	249	TYR	N-CA-C	-5.32	107.09	113.21
1	B	286	TYR	N-CA-C	-5.30	105.67	111.82
1	A	211	LYS	N-CA-C	5.30	117.97	110.50
1	A	128	VAL	CB-CA-C	-5.21	105.86	111.59
1	B	159	ILE	N-CA-C	5.17	117.51	111.05
1	B	371	ALA	N-CA-C	5.06	115.99	108.60
1	A	386	LYS	N-CA-C	5.04	117.89	111.69

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	3323	0	3252	94	0
1	B	3300	0	3241	120	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	13	0	13	0	0
4	A	104	0	0	0	0
4	B	55	0	0	3	0
All	All	6797	0	6506	210	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 16.

All (210) 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:387:ASP:OD2	1:B:392:LYS:HG3	1.63	0.95
1:A:204:ASN:HD22	1:A:205:PRO:N	1.72	0.87
1:B:228:GLU:HG3	1:B:344:LEU:HD22	1.55	0.86
1:A:171:VAL:HG23	1:A:172:LEU:HD13	1.58	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ASN:HD22	1:A:55:GLU:HG3	1.43	0.83
1:A:209:ALA:H	1:A:404:ASN:HD21	1.27	0.80
1:A:61:PRO:HB2	1:A:62:PRO:HD2	1.65	0.79
1:A:199:LEU:HD23	1:A:199:LEU:H	1.45	0.79
1:A:387:ASP:OD2	1:A:392:LYS:HG3	1.84	0.77
1:B:325:ASP:OD2	1:B:329:LYS:HB3	1.85	0.77
1:B:22:LYS:HE3	1:B:137:ILE:HD11	1.70	0.74
1:A:391:LEU:O	1:A:397:SER:HB3	1.88	0.73
1:B:410:ARG:HB2	1:B:410:ARG:HH11	1.52	0.73
1:A:225:ILE:O	1:A:228:GLU:HG2	1.88	0.73
1:B:325:ASP:OD1	1:B:327:SER:HB3	1.89	0.72
1:B:172:LEU:HD23	1:B:172:LEU:H	1.54	0.70
1:B:6:GLN:HE21	1:B:6:GLN:HA	1.57	0.69
1:B:64:ALA:HB2	4:B:535:HOH:O	1.91	0.69
1:A:29:MET:HE3	1:A:30:GLN:O	1.92	0.69
1:A:204:ASN:ND2	1:A:206:LEU:H	1.90	0.69
1:A:1:MET:SD	1:A:1:MET:N	2.63	0.69
1:A:204:ASN:HD22	1:A:204:ASN:C	1.96	0.69
1:B:299:ALA:HB3	1:B:313:LYS:HE2	1.76	0.68
1:A:161:GLN:HE21	1:A:161:GLN:HA	1.57	0.67
1:B:302:ILE:HD11	1:B:312:MET:HG3	1.77	0.67
1:B:52:ASN:HD22	1:B:55:GLU:HG3	1.61	0.66
1:B:304:GLY:O	1:B:306:THR:HG22	1.94	0.66
1:B:196:GLU:H	1:B:370:LYS:NZ	1.92	0.66
1:B:360:ILE:HG22	1:B:403:GLN:CD	2.22	0.65
1:A:52:ASN:ND2	1:A:55:GLU:HG3	2.10	0.64
1:A:197:GLU:OE1	1:A:360:ILE:HD11	1.97	0.64
1:A:114:ILE:H	1:A:114:ILE:HD12	1.61	0.64
1:B:116:PHE:HA	1:B:316:PHE:CE1	2.33	0.63
1:A:225:ILE:O	1:A:228:GLU:CG	2.46	0.63
1:B:196:GLU:H	1:B:370:LYS:HZ1	1.47	0.62
1:B:415:LEU:O	1:B:416:LYS:HB2	1.99	0.62
1:A:23:ILE:HG23	1:A:24:PRO:HD2	1.79	0.62
1:B:197:GLU:OE2	1:B:360:ILE:HD11	2.00	0.62
1:B:204:ASN:ND2	1:B:206:LEU:H	1.98	0.62
1:A:69:VAL:HG23	1:A:417:ASN:OD1	2.00	0.62
1:B:300:LYS:O	1:B:309:LEU:HD22	1.99	0.61
1:A:211:LYS:HE3	1:A:370:LYS:HB2	1.81	0.61
1:B:232:TYR:O	1:B:234:ILE:HG23	2.01	0.61
1:B:117:TRP:CE2	1:B:128:VAL:HB	2.36	0.61
1:B:196:GLU:HB2	1:B:370:LYS:HE2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:ALA:N	1:A:404:ASN:HD21	1.98	0.60
1:A:22:LYS:HA	1:A:29:MET:HE2	1.83	0.59
1:A:69:VAL:HG13	1:A:69:VAL:O	2.00	0.59
1:B:138:GLN:HB3	1:B:139:PRO:HD2	1.84	0.59
1:B:6:GLN:HA	1:B:6:GLN:NE2	2.17	0.59
1:A:228:GLU:HG3	1:A:229:HIS:N	2.16	0.59
1:B:304:GLY:C	1:B:306:THR:H	2.10	0.58
1:A:385:ILE:HG22	1:A:386:LYS:HD2	1.86	0.58
1:A:199:LEU:HD23	1:A:199:LEU:N	2.18	0.58
1:B:13:VAL:HA	1:B:18:ILE:HG22	1.85	0.57
1:B:228:GLU:CG	1:B:344:LEU:HD22	2.30	0.57
1:A:200:GLU:HB2	1:A:203:THR:HG23	1.87	0.57
1:A:204:ASN:HD22	1:A:205:PRO:CD	2.17	0.57
1:A:24:PRO:O	1:A:26:ALA:N	2.38	0.56
1:A:204:ASN:ND2	1:A:204:ASN:C	2.64	0.55
1:B:169:HIS:CE1	1:B:171:VAL:HG12	2.41	0.55
1:A:169:HIS:CE1	1:A:171:VAL:HG22	2.42	0.55
1:B:51:THR:HG22	1:B:186:ARG:HH21	1.70	0.55
1:B:228:GLU:HG3	1:B:344:LEU:CD2	2.32	0.55
1:B:124:THR:HA	1:B:298:LYS:O	2.05	0.55
1:B:377:ILE:HG12	4:B:520:HOH:O	2.05	0.54
1:B:378:VAL:HB	1:B:379:PRO:HD3	1.90	0.54
1:A:61:PRO:CB	1:A:62:PRO:HD2	2.36	0.54
1:B:21:ILE:HG22	1:B:136:VAL:HA	1.90	0.54
1:A:22:LYS:HA	1:A:29:MET:CE	2.37	0.54
1:B:200:GLU:HB2	1:B:203:THR:HG23	1.90	0.54
1:B:153:ILE:HG12	1:B:154:GLY:N	2.23	0.53
1:B:371:ALA:HA	1:B:417:ASN:HB3	1.91	0.53
1:A:321:LEU:HD12	1:A:340:LEU:HB2	1.89	0.53
1:B:166:SER:OG	1:B:183:GLN:NE2	2.42	0.53
1:B:225:ILE:O	1:B:228:GLU:OE2	2.27	0.53
1:A:237:ASN:CG	1:A:239:ASN:ND2	2.66	0.52
1:A:247:ASN:HB2	1:A:250:TYR:HD2	1.74	0.52
1:A:362:ALA:HB1	1:A:367:ASN:HB3	1.91	0.52
1:B:302:ILE:HG23	1:B:302:ILE:O	2.10	0.52
1:B:120:SER:HB3	1:B:125:GLU:O	2.10	0.52
1:B:228:GLU:OE1	1:B:344:LEU:O	2.27	0.52
1:B:387:ASP:HB3	1:B:390:ASN:O	2.09	0.52
1:B:107:LEU:O	1:B:111:VAL:HG23	2.10	0.52
1:B:136:VAL:HG12	1:B:137:ILE:N	2.26	0.51
1:B:268:HIS:O	1:B:271:LYS:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:416:LYS:CG	1:B:417:ASN:H	2.23	0.51
1:A:396:LEU:HA	1:A:401:ASN:HB2	1.92	0.50
1:A:204:ASN:HD21	1:A:206:LEU:H	1.60	0.50
1:A:228:GLU:OE1	1:A:344:LEU:O	2.29	0.50
1:B:130:ASP:OD1	1:B:131:THR:HG23	2.11	0.50
1:B:176:ARG:HD3	1:B:237:ASN:HA	1.94	0.50
1:A:279:ASN:O	1:A:283:LEU:HD23	2.12	0.50
1:A:114:ILE:HD12	1:A:114:ILE:N	2.26	0.50
1:B:180:GLY:HA3	1:B:231:LEU:O	2.12	0.50
1:A:209:ALA:H	1:A:404:ASN:ND2	2.03	0.49
1:B:181:SER:O	1:B:231:LEU:HD23	2.11	0.49
1:A:199:LEU:H	1:A:199:LEU:CD2	2.19	0.49
1:A:381:GLU:N	1:A:381:GLU:CD	2.71	0.49
1:B:416:LYS:HG2	1:B:417:ASN:H	1.77	0.49
1:A:237:ASN:CG	1:A:239:ASN:HD22	2.20	0.49
1:B:96:ARG:HA	1:B:385:ILE:HG23	1.95	0.49
1:B:29:MET:O	1:B:29:MET:HG3	2.13	0.48
1:A:274:ASP:OD2	1:A:276:LEU:HB2	2.13	0.48
1:B:225:ILE:HA	1:B:228:GLU:OE2	2.13	0.48
1:B:279:ASN:ND2	1:B:282:ARG:NH1	2.61	0.48
1:B:25:ASN:HD22	1:B:131:THR:HG22	1.78	0.48
1:A:335:LEU:HD12	1:A:336:LYS:N	2.28	0.48
1:A:259:PHE:HE2	1:A:277:GLN:HG2	1.78	0.48
1:B:304:GLY:O	1:B:306:THR:N	2.43	0.48
1:B:368:PHE:O	1:B:370:LYS:NZ	2.46	0.48
1:A:195:PHE:HZ	1:A:360:ILE:HG23	1.78	0.48
1:A:171:VAL:HG23	1:A:172:LEU:CD1	2.37	0.47
1:B:278:GLU:OE1	1:B:282:ARG:NH2	2.47	0.47
1:A:198:SER:O	1:A:200:GLU:HG3	2.15	0.47
1:B:117:TRP:CD2	1:B:312:MET:HE2	2.50	0.47
1:B:134:ILE:HG23	1:B:148:LEU:CD2	2.45	0.47
1:B:241:VAL:HG12	1:B:258:SER:HA	1.97	0.47
1:A:161:GLN:O	1:A:161:GLN:HG3	2.14	0.47
1:A:416:LYS:CD	1:A:417:ASN:H	2.28	0.47
1:A:225:ILE:HA	1:A:228:GLU:HG2	1.96	0.47
1:B:136:VAL:O	1:B:143:TYR:HA	2.14	0.47
1:B:143:TYR:H	1:B:143:TYR:HD2	1.61	0.46
1:A:274:ASP:OD1	1:A:277:GLN:NE2	2.47	0.46
1:B:311:TYR:O	1:B:315:VAL:HG23	2.15	0.46
1:A:381:GLU:CD	1:A:381:GLU:H	2.22	0.46
1:B:48:ASP:OD2	1:B:153:ILE:HD13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:ILE:C	1:B:124:THR:H	2.23	0.46
1:B:229:HIS:NE2	1:B:344:LEU:O	2.48	0.46
1:A:378:VAL:HB	1:A:379:PRO:HD3	1.96	0.46
1:B:274:ASP:OD1	1:B:276:LEU:N	2.49	0.46
1:B:61:PRO:HB2	1:B:63:GLU:OE2	2.15	0.46
1:B:125:GLU:HB2	1:B:301:SER:HB3	1.97	0.46
1:B:6:GLN:NE2	1:B:6:GLN:CA	2.79	0.46
1:B:279:ASN:O	1:B:283:LEU:HD23	2.16	0.46
1:A:360:ILE:HG22	1:A:403:GLN:CD	2.41	0.45
1:B:13:VAL:HG11	1:B:20:TYR:CD1	2.51	0.45
1:B:204:ASN:C	1:B:204:ASN:HD22	2.25	0.45
1:B:360:ILE:HG22	1:B:403:GLN:NE2	2.30	0.45
1:A:22:LYS:HG2	1:A:29:MET:HE2	1.97	0.45
1:B:2:PHE:HB3	1:B:95:GLU:OE2	2.17	0.45
1:B:15:GLY:O	1:B:138:GLN:NE2	2.49	0.45
1:A:24:PRO:O	1:A:25:ASN:C	2.59	0.45
1:A:171:VAL:CG2	1:A:172:LEU:HD13	2.39	0.45
1:A:23:ILE:HG23	1:A:24:PRO:CD	2.46	0.45
1:B:86:TYR:CE1	1:B:90:VAL:HG21	2.52	0.44
1:B:274:ASP:OD1	1:B:274:ASP:C	2.59	0.44
1:B:312:MET:O	1:B:315:VAL:HB	2.16	0.44
1:B:410:ARG:HB2	1:B:410:ARG:NH1	2.25	0.44
1:A:59:ASN:O	1:A:61:PRO:HD3	2.16	0.44
1:B:255:LEU:HD22	1:B:256:GLU:N	2.32	0.44
1:B:416:LYS:CE	1:B:417:ASN:H	2.29	0.44
1:A:225:ILE:C	1:A:228:GLU:HG2	2.42	0.44
1:A:52:ASN:HD22	1:A:55:GLU:CG	2.22	0.44
1:B:324:GLU:HA	1:B:329:LYS:O	2.17	0.44
1:B:171:VAL:HG13	1:B:172:LEU:N	2.33	0.44
1:B:169:HIS:ND1	1:B:170:ASP:N	2.65	0.44
1:B:30:GLN:OE1	1:B:30:GLN:HA	2.16	0.44
1:A:209:ALA:HB3	1:A:406:GLU:HG3	2.00	0.43
1:B:396:LEU:HA	1:B:401:ASN:HB2	2.00	0.43
1:B:85:ASN:OD1	1:B:378:VAL:HG21	2.18	0.43
1:A:275:SER:HA	1:B:245:ASN:HB2	2.01	0.43
1:A:365:PHE:C	1:A:367:ASN:N	2.75	0.43
1:B:224:LEU:O	1:B:227:ALA:HB3	2.19	0.43
1:A:239:ASN:OD1	1:A:239:ASN:C	2.61	0.43
1:A:52:ASN:HA	1:A:53:PRO:HD3	1.86	0.43
1:A:60:PRO:HA	1:A:61:PRO:HD2	1.95	0.43
1:B:156:SER:O	1:B:157:ALA:C	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:ASN:HD22	1:B:205:PRO:N	2.17	0.43
1:A:23:ILE:HG22	1:A:24:PRO:O	2.19	0.42
1:A:200:GLU:HB2	1:A:203:THR:CG2	2.49	0.42
1:B:98:TYR:HD1	1:B:107:LEU:HD12	1.84	0.42
1:B:138:GLN:HB3	1:B:139:PRO:CD	2.49	0.42
1:B:319:LYS:HD3	1:B:320:TYR:CE2	2.53	0.42
1:A:122:ILE:HG22	1:A:124:THR:HG22	2.00	0.42
1:B:258:SER:OG	1:B:261:GLU:HB2	2.19	0.42
1:B:304:GLY:C	1:B:306:THR:N	2.77	0.42
1:A:248:ALA:C	1:A:250:TYR:H	2.27	0.42
1:A:255:LEU:C	1:A:255:LEU:HD13	2.43	0.42
1:B:277:GLN:HE21	1:B:277:GLN:HB2	1.67	0.42
1:A:416:LYS:HD2	1:A:417:ASN:OD1	2.18	0.42
1:B:315:VAL:HA	4:B:655:HOH:O	2.18	0.42
1:B:1:MET:SD	1:B:1:MET:N	2.82	0.42
1:B:13:VAL:HG11	1:B:20:TYR:HD1	1.84	0.42
1:B:357:PHE:HB3	1:B:359:VAL:HG13	2.02	0.42
1:A:359:VAL:O	1:B:201:VAL:HG21	2.19	0.42
1:B:360:ILE:HG22	1:B:403:GLN:OE1	2.19	0.42
1:A:59:ASN:HA	1:A:60:PRO:HD3	1.83	0.42
1:A:271:LYS:HG3	1:B:366:LEU:HD11	2.01	0.42
1:A:35:PHE:CD1	1:A:35:PHE:N	2.88	0.42
1:B:392:LYS:C	1:B:394:ALA:H	2.28	0.42
1:B:286:TYR:O	1:B:290:LYS:HG3	2.20	0.42
1:A:161:GLN:HE21	1:A:161:GLN:CA	2.23	0.41
1:A:171:VAL:HG23	1:A:172:LEU:N	2.35	0.41
1:B:1:MET:N	1:B:1:MET:HE3	2.35	0.41
1:A:180:GLY:HA2	1:A:230:ARG:O	2.20	0.41
1:A:128:VAL:HG22	1:A:129:ILE:N	2.36	0.41
1:A:333:ASP:CG	1:A:336:LYS:HG3	2.46	0.41
1:A:365:PHE:CG	1:A:366:LEU:N	2.88	0.41
1:A:383:TYR:HA	1:A:388:GLY:O	2.21	0.41
1:B:51:THR:CG2	1:B:186:ARG:HH21	2.33	0.41
1:B:125:GLU:OE2	1:B:303:ILE:HG12	2.21	0.41
1:A:271:LYS:HG3	1:B:366:LEU:CD1	2.51	0.41
1:A:302:ILE:HD11	1:A:312:MET:HG3	2.02	0.41
1:B:28:GLN:HG2	1:B:28:GLN:O	2.21	0.41
1:B:134:ILE:HG23	1:B:148:LEU:HD22	2.03	0.40
1:B:391:LEU:O	1:B:397:SER:HB3	2.21	0.40
1:B:176:ARG:HG2	1:B:235:ALA:O	2.22	0.40
1:A:219:THR:O	1:A:222:HIS:HB3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:LYS:HD3	1:A:417:ASN:H	1.86	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%)	373 (92%)	24 (6%)	8 (2%)	6	7
1	B	403/425 (95%)	363 (90%)	29 (7%)	11 (3%)	4	3
All	All	808/850 (95%)	736 (91%)	53 (7%)	19 (2%)	4	4

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	A	208	GLY
1	A	392	LYS
1	B	305	THR
1	B	392	LYS
1	B	416	LYS
1	A	199	LEU
1	A	306	THR
1	B	51	THR
1	B	123	ASP
1	B	307	ALA
1	A	210	GLY
1	A	394	ALA
1	B	322	LEU
1	B	64	ALA
1	B	25	ASN
1	B	302	ILE

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Mol	Chain	Res	Type
1	B	139	PRO
1	A	168	GLY

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	367/380 (97%)	352 (96%)	15 (4%)	27	44
1	B	366/380 (96%)	350 (96%)	16 (4%)	25	41
All	All	733/760 (96%)	702 (96%)	31 (4%)	26	43

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	54	GLU
1	A	126	LEU
1	A	143	TYR
1	A	161	GLN
1	A	172	LEU
1	A	199	LEU
1	A	201	VAL
1	A	276	LEU
1	A	277	GLN
1	A	366	LEU
1	A	369	ASP
1	A	395	ASN
1	A	398	THR
1	A	416	LYS
1	B	1	MET
1	B	28	GLN
1	B	46	GLU
1	B	50	PHE
1	B	51	THR
1	B	63	GLU

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Mol	Chain	Res	Type
1	B	101	ASP
1	B	139	PRO
1	B	172	LEU
1	B	196	GLU
1	B	228	GLU
1	B	255	LEU
1	B	274	ASP
1	B	278	GLU
1	B	364	THR
1	B	410	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	6	GLN
1	A	28	GLN
1	A	30	GLN
1	A	39	ASN
1	A	52	ASN
1	A	59	ASN
1	A	135	ASN
1	A	161	GLN
1	A	204	ASN
1	A	367	ASN
1	A	399	ASN
1	B	6	GLN
1	B	25	ASN
1	B	28	GLN
1	B	52	ASN
1	B	59	ASN
1	B	204	ASN
1	B	226	HIS
1	B	245	ASN
1	B	277	GLN
1	B	279	ASN
1	B	367	ASN
1	B	417	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	AHL	A	500	2	11,12,12	0.99	1 (9%)	11,14,14	0.88	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	AHL	A	500	2	-	1/13/13/13	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	AHL	O2-N2	-2.57	1.33	1.40

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	500	AHL	NH1-CZ-NE-CD

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	411/425 (96%)	0.19	33 (8%) 18 15	10, 26, 57, 65	0
1	B	409/425 (96%)	0.71	51 (12%) 8 6	15, 39, 61, 69	0
All	All	820/850 (96%)	0.45	84 (10%) 12 9	10, 32, 59, 69	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	25	ASN	6.1
1	A	305	THR	5.5
1	A	208	GLY	5.1
1	B	168	GLY	5.1
1	B	417	ASN	4.8
1	B	304	GLY	4.6
1	A	199	LEU	4.4
1	A	209	ALA	4.4
1	A	228	GLU	4.1
1	A	398	THR	4.1
1	A	381	GLU	4.0
1	A	27	GLY	4.0
1	B	305	THR	4.0
1	B	126	LEU	4.0
1	A	207	LEU	3.7
1	A	368	PHE	3.6
1	B	65	LYS	3.6
1	B	139	PRO	3.6
1	A	62	PRO	3.6
1	A	397	SER	3.6
1	B	307	ALA	3.5
1	B	199	LEU	3.4
1	B	138	GLN	3.4
1	B	228	GLU	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	397	SER	3.3
1	A	248	ALA	3.3
1	B	198	SER	3.2
1	A	64	ALA	3.2
1	A	417	ASN	3.2
1	A	369	ASP	3.2
1	A	25	ASN	3.1
1	B	381	GLU	3.1
1	B	137	ILE	3.1
1	B	303	ILE	3.1
1	A	61	PRO	3.0
1	B	143	TYR	3.0
1	B	312	MET	3.0
1	A	60	PRO	3.0
1	B	62	PRO	3.0
1	A	69	VAL	3.0
1	B	140	ASP	3.0
1	B	311	TYR	2.9
1	B	323	SER	2.8
1	A	28	GLN	2.8
1	A	250	TYR	2.7
1	A	26	ALA	2.7
1	A	393	GLY	2.7
1	B	301	SER	2.7
1	B	253	SER	2.6
1	A	57	ASP	2.6
1	B	398	THR	2.6
1	B	180	GLY	2.6
1	A	365	PHE	2.6
1	B	59	ASN	2.5
1	B	64	ALA	2.5
1	B	142	SER	2.5
1	B	122	ILE	2.5
1	A	366	LEU	2.5
1	A	251	GLU	2.4
1	B	325	ASP	2.4
1	A	24	PRO	2.3
1	B	302	ILE	2.3
1	A	249	TYR	2.3
1	B	130	ASP	2.3
1	B	117	TRP	2.3
1	B	27	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	143	TYR	2.2
1	B	309	LEU	2.1
1	B	335	LEU	2.1
1	B	308	SER	2.1
1	B	254	GLY	2.1
1	B	20	TYR	2.1
1	B	116	PHE	2.1
1	A	168	GLY	2.1
1	B	121	THR	2.1
1	B	410	ARG	2.1
1	B	315	VAL	2.1
1	A	63	GLU	2.1
1	B	179	TYR	2.1
1	B	6	GLN	2.0
1	B	332	VAL	2.0
1	B	61	PRO	2.0
1	B	247	ASN	2.0
1	B	306	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	AHL	A	500	13/13	0.79	0.15	37,38,42,42	0
2	ZN	A	452	1/1	0.96	0.04	22,22,22,22	0
2	ZN	B	451	1/1	0.98	0.03	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.