Full wwPDB X-ray Structure Validation Report ⓘ

Aug 16, 2023 – 01:39 AM EDT

PDB ID	:	1ZKX
Title	:	Crystal structure of Glu158Ala/Thr159Ala/Asn160Ala- a triple mutant of Clostridium botulinum neurotoxin E catalytic domain
Authors	:	Agarwal, R.; Binz, T.; Swaminathan, S.
Deposited on	:	2005-05-04
Resolution	:	2.52 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

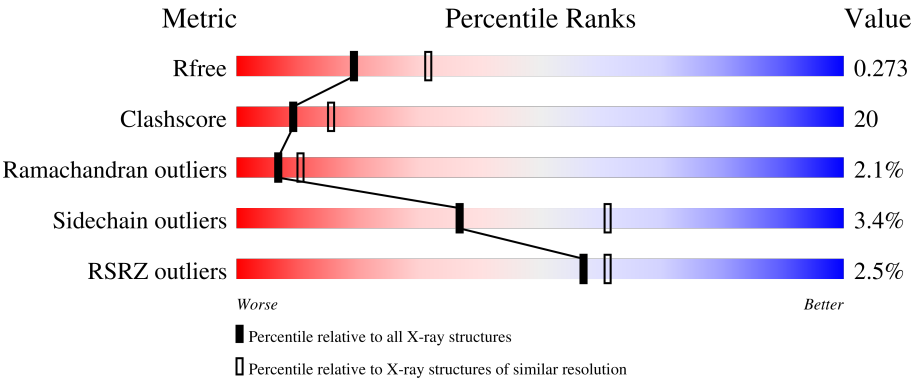
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation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

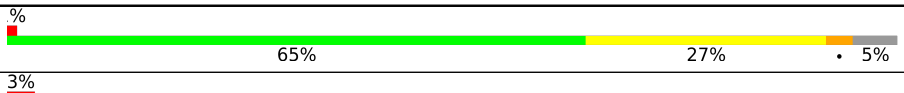
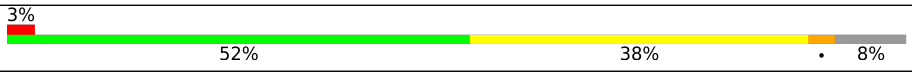
The reported resolution of this entry is 2.52 Å.

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	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (2.54-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	420	
1	B	420	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CL	B	427	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6548 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 E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	400	Total	C	N	O	S	0	0	0
			3198	2048	532	610	8			
1	B	387	Total	C	N	O	S	0	0	0
			3101	1988	515	591	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	ALA	GLU	engineered mutation	UNP Q00496
A	159	ALA	THR	engineered mutation	UNP Q00496
A	160	ALA	ASN	engineered mutation	UNP Q00496
B	158	ALA	GLU	engineered mutation	UNP Q00496
B	159	ALA	THR	engineered mutation	UNP Q00496
B	160	ALA	ASN	engineered mutation	UNP Q00496

- Molecule 2 is ZINC ION (three-letter code: 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 CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Cl	0	0
			3	3		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

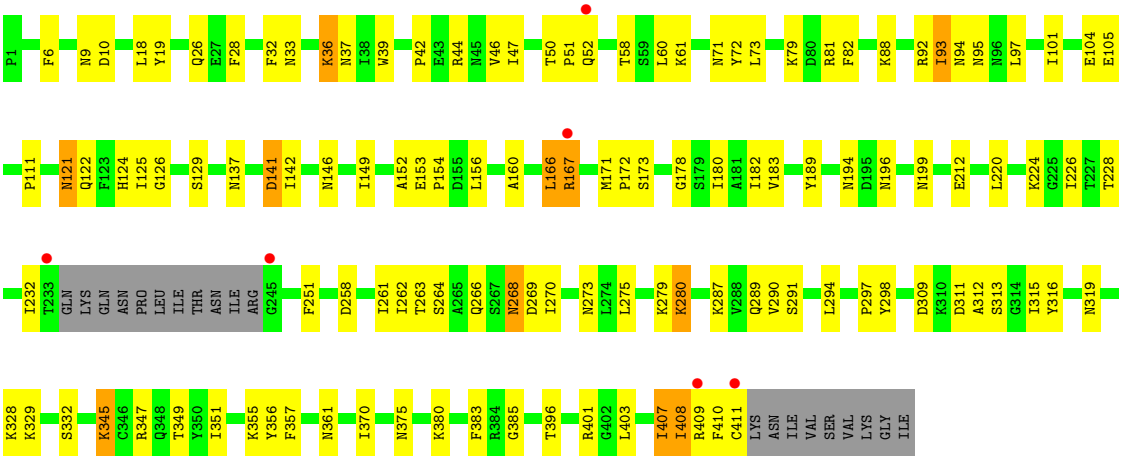
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	130	Total O 130 130	0	0
4	B	113	Total O 113 113	0	0

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.

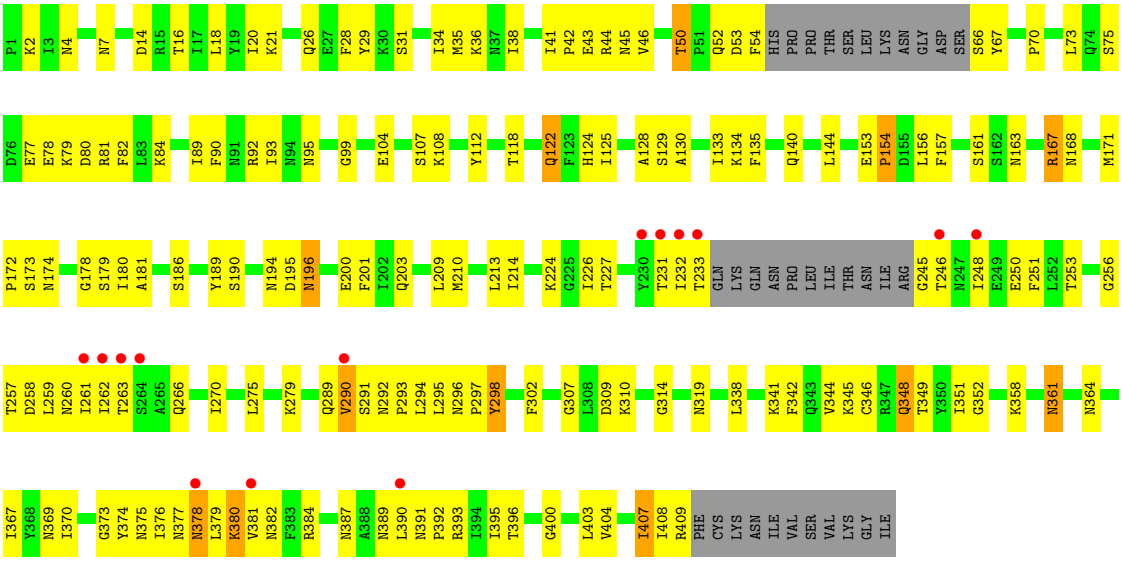
• Molecule 1: botulinum neurotoxin type E

Chain A:



• Molecule 1: botulinum neurotoxin type E

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	88.90Å 144.78Å 83.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.50 – 2.52 46.50 – 2.52	Depositor EDS
% Data completeness (in resolution range)	72.8 (46.50-2.52) 83.3 (46.50-2.52)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.23 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.208 , 0.259 0.224 , 0.273	Depositor DCC
R_{free} test set	928 reflections (2.81%)	wwPDB-VP
Wilson B-factor (Å^2)	33.4	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 43.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6548	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 3.94% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CL

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	0.43	1/3270 (0.0%)	0.68	3/4428 (0.1%)
1	B	0.39	0/3168	0.63	0/4287
All	All	0.41	1/6438 (0.0%)	0.66	3/8715 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	166	LEU	C-O	-9.05	1.06	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	166	LEU	O-C-N	-8.14	109.68	122.70
1	A	167	ARG	N-CA-CB	-7.05	97.91	110.60
1	A	93	ILE	N-CA-C	-5.67	95.70	111.00

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	3198	0	3154	103	0
1	B	3101	0	3064	151	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	3	0	0	0	0
3	B	1	0	0	2	0
4	A	130	0	0	19	0
4	B	113	0	0	19	0
All	All	6548	0	6218	248	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 20.

All (248) 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:379:LEU:HD23	1:B:384:ARG:HD2	1.55	0.89
1:B:263:THR:H	1:B:266:GLN:HE21	1.18	0.88
1:B:338:LEU:HA	4:B:882:HOH:O	1.73	0.88
1:B:232:ILE:HD13	1:B:261:ILE:HG22	1.55	0.87
1:B:231:THR:HG22	4:B:930:HOH:O	1.74	0.85
1:B:292:ASN:OD1	1:B:294:LEU:HB2	1.80	0.81
1:B:263:THR:H	1:B:266:GLN:NE2	1.79	0.78
1:A:105:GLU:CD	1:A:329:LYS:HD2	2.05	0.77
1:A:332:SER:HB2	4:A:554:HOH:O	1.87	0.74
1:A:125:ILE:HD11	1:B:294:LEU:CD2	2.18	0.73
1:B:130:ALA:HB2	1:B:144:LEU:CD2	2.20	0.72
1:B:92:ARG:HA	1:B:370:ILE:HG23	1.74	0.70
1:B:50:THR:HG23	1:B:52:GLN:HE21	1.57	0.69
1:A:409:ARG:NE	4:A:451:HOH:O	2.23	0.69
1:B:290:VAL:CG1	1:B:296:ASN:HD21	2.05	0.69
1:B:122:GLN:HE21	1:B:122:GLN:C	1.95	0.69
1:B:75:SER:OG	1:B:78:GLU:HG3	1.93	0.69
1:A:196:ASN:HB3	4:A:550:HOH:O	1.92	0.69
1:A:294:LEU:HD21	1:B:125:ILE:HD11	1.75	0.69
1:A:232:ILE:HD13	1:A:261:ILE:HG21	1.75	0.69
1:B:31:SER:OG	1:B:41:ILE:HG12	1.92	0.69
1:B:167:ARG:NH2	3:B:427:CL:CL	2.63	0.68
1:A:258:ASP:HA	1:A:261:ILE:HD13	1.74	0.67
1:B:409:ARG:HG3	1:B:409:ARG:HH11	1.59	0.67
1:A:97:LEU:O	1:A:101:ILE:HG12	1.94	0.67
1:B:378:ASN:ND2	1:B:384:ARG:HH12	1.91	0.67
1:B:66:SER:HA	1:B:157:PHE:CD2	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:545:HOH:O	1:B:36:LYS:HE2	1.95	0.67
1:A:409:ARG:HB2	4:A:517:HOH:O	1.92	0.67
1:B:309:ASP:OD2	1:B:319:ASN:HB2	1.95	0.66
1:A:125:ILE:HD11	1:B:294:LEU:HD23	1.77	0.66
1:A:275:LEU:HG	1:A:279:LYS:NZ	2.11	0.66
1:A:58:THR:HG21	1:A:61:LYS:HB2	1.77	0.66
1:B:275:LEU:HG	1:B:279:LYS:HE3	1.77	0.66
1:B:400:GLY:O	1:B:404:VAL:HG23	1.96	0.65
3:B:427:CL:CL	4:B:874:HOH:O	2.51	0.65
1:B:376:ILE:O	1:B:379:LEU:HB2	1.97	0.65
1:A:124:HIS:HA	4:A:483:HOH:O	1.97	0.65
1:B:196:ASN:HA	4:B:832:HOH:O	1.95	0.65
1:A:328:LYS:HE2	4:A:449:HOH:O	1.97	0.65
1:A:409:ARG:HG2	4:A:451:HOH:O	1.96	0.64
1:B:341:LYS:HB2	4:B:882:HOH:O	1.97	0.64
1:B:374:TYR:O	1:B:376:ILE:HG13	1.98	0.63
1:A:311:ASP:OD1	1:A:313:SER:N	2.32	0.62
1:B:66:SER:HA	1:B:157:PHE:HD2	1.63	0.62
1:B:128:ALA:O	1:B:180:ILE:HG13	1.99	0.62
1:A:275:LEU:HG	1:A:279:LYS:HZ1	1.65	0.62
1:B:54:PHE:O	1:B:70:PRO:HG3	2.00	0.61
1:B:248:ILE:HA	1:B:251:PHE:HD1	1.66	0.61
1:A:289:GLN:HG3	1:A:291:SER:H	1.64	0.61
1:A:153:GLU:HB2	1:A:154:PRO:CD	2.31	0.60
1:A:104:GLU:HG2	4:A:540:HOH:O	2.01	0.60
1:A:266:GLN:O	1:A:270:ILE:HG12	2.02	0.60
1:B:364:ASN:HB3	1:B:367:ILE:HD12	1.83	0.59
1:B:21:LYS:HE3	1:B:28:PHE:CZ	2.37	0.59
1:A:19:TYR:HB3	1:A:28:PHE:HB3	1.84	0.59
1:A:370:ILE:HG12	4:A:527:HOH:O	2.03	0.59
1:B:409:ARG:HG3	1:B:409:ARG:NH1	2.17	0.59
1:A:60:LEU:HB2	4:A:510:HOH:O	2.02	0.59
1:A:232:ILE:CD1	1:A:261:ILE:HG21	2.32	0.59
1:B:226:ILE:CG2	1:B:270:ILE:HD13	2.33	0.59
1:A:407:ILE:HG22	1:A:407:ILE:O	2.03	0.58
1:B:341:LYS:HD2	4:B:882:HOH:O	2.03	0.58
1:B:20:ILE:HG12	1:B:133:ILE:HG22	1.84	0.57
1:B:21:LYS:HE3	1:B:28:PHE:CE2	2.39	0.57
1:B:256:GLY:O	1:B:258:ASP:N	2.36	0.57
1:B:44:ARG:O	1:B:46:VAL:HG13	2.05	0.57
1:A:226:ILE:CG2	1:A:270:ILE:HD13	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:PRO:HB2	1:B:297:PRO:HB2	1.88	0.56
1:B:50:THR:HG22	1:B:52:GLN:H	1.70	0.56
1:A:396:THR:HG22	4:A:470:HOH:O	2.06	0.56
1:B:389:ASN:O	1:B:390:LEU:HD23	2.06	0.56
1:A:81:ARG:HG3	1:A:81:ARG:HH11	1.71	0.56
1:B:261:ILE:HG13	4:B:859:HOH:O	2.06	0.56
1:A:269:ASP:O	1:A:273:ASN:HB2	2.06	0.56
1:B:263:THR:OG1	1:B:266:GLN:HG3	2.05	0.56
1:A:73:LEU:HD21	1:A:82:PHE:HB2	1.88	0.55
1:B:130:ALA:HB2	1:B:144:LEU:HD23	1.87	0.55
1:B:190:SER:OG	1:B:203:GLN:HB3	2.06	0.55
1:A:275:LEU:CG	1:A:279:LYS:HZ1	2.20	0.55
1:B:153:GLU:HB2	1:B:154:PRO:HD2	1.89	0.54
1:B:163:ASN:HB3	1:B:179:SER:OG	2.07	0.54
1:B:369:ASN:OD1	1:B:373:GLY:N	2.40	0.54
1:A:289:GLN:NE2	1:A:290:VAL:H	2.05	0.54
1:B:167:ARG:HB2	1:B:167:ARG:HH11	1.73	0.54
1:A:280:LYS:NZ	1:A:280:LYS:HA	2.22	0.54
1:B:260:ASN:N	4:B:859:HOH:O	2.40	0.54
1:A:153:GLU:HB2	1:A:154:PRO:HD2	1.89	0.54
1:A:375:ASN:HB3	1:A:380:LYS:HA	1.89	0.54
1:B:232:ILE:HD13	1:B:261:ILE:CG2	2.34	0.54
1:B:232:ILE:O	1:B:233:THR:O	2.26	0.54
1:B:29:TYR:CG	1:B:43:GLU:HG3	2.43	0.53
1:B:168:ASN:N	1:B:168:ASN:HD22	2.07	0.53
1:A:262:ILE:HG23	1:A:266:GLN:HE21	1.74	0.53
1:B:256:GLY:C	1:B:258:ASP:H	2.13	0.53
1:B:93:ILE:O	1:B:99:GLY:HA3	2.09	0.52
1:B:310:LYS:HE2	1:B:314:GLY:HA2	1.90	0.52
1:B:378:ASN:HD22	1:B:384:ARG:HH12	1.56	0.52
1:B:80:ASP:O	1:B:84:LYS:HD3	2.10	0.52
1:B:292:ASN:ND2	1:B:293:PRO:HD2	2.24	0.52
1:A:311:ASP:OD2	1:A:315:ILE:HB	2.09	0.52
1:A:166:LEU:HD22	1:A:167:ARG:HH12	1.75	0.52
1:B:73:LEU:HD21	1:B:82:PHE:HB2	1.90	0.52
1:B:393:ARG:HG2	4:B:890:HOH:O	2.10	0.52
1:B:403:LEU:HD13	1:B:403:LEU:O	2.10	0.52
1:A:403:LEU:O	1:A:407:ILE:HG13	2.09	0.51
1:B:387:ASN:ND2	1:B:390:LEU:HD12	2.25	0.51
1:A:356:TYR:O	1:A:401:ARG:NH2	2.36	0.51
1:B:95:ASN:ND2	4:B:830:HOH:O	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:ILE:O	1:B:93:ILE:HG13	2.11	0.51
1:B:210:MET:O	1:B:214:ILE:HG13	2.11	0.50
1:B:358:LYS:O	1:B:396:THR:HG23	2.11	0.50
1:B:214:ILE:CG2	1:B:253:THR:HG23	2.41	0.50
1:B:50:THR:CG2	1:B:52:GLN:H	2.25	0.50
1:B:118:THR:HG21	1:B:124:HIS:CD2	2.46	0.50
1:A:121:ASN:HA	1:A:287:LYS:O	2.12	0.50
1:B:50:THR:CG2	1:B:52:GLN:HB2	2.42	0.50
1:B:77:GLU:HA	1:B:77:GLU:OE1	2.11	0.50
1:B:377:ASN:O	1:B:379:LEU:N	2.44	0.49
1:A:345:LYS:HD2	1:A:383:PHE:CD2	2.48	0.49
1:B:122:GLN:C	1:B:122:GLN:NE2	2.65	0.49
1:B:369:ASN:O	1:B:373:GLY:HA2	2.12	0.49
1:B:307:GLY:HA2	4:B:862:HOH:O	2.11	0.49
1:B:52:GLN:NE2	1:B:52:GLN:N	2.61	0.49
1:B:201:PHE:CD2	1:B:345:LYS:HG2	2.48	0.49
1:A:121:ASN:HD22	1:A:122:GLN:N	2.11	0.49
1:B:346:CYS:HB3	4:B:873:HOH:O	2.13	0.49
1:B:403:LEU:O	1:B:407:ILE:HG13	2.13	0.49
1:A:297:PRO:HG2	1:B:298:TYR:CD2	2.48	0.48
1:B:374:TYR:O	1:B:376:ILE:N	2.46	0.48
1:B:381:VAL:HG21	1:B:384:ARG:NH2	2.27	0.48
1:A:264:SER:O	1:A:268:ASN:HB2	2.13	0.48
1:B:34:ILE:HG23	1:B:35:MET:HG2	1.96	0.48
1:A:50:THR:O	1:A:52:GLN:N	2.46	0.48
1:A:126:GLY:HA3	4:A:505:HOH:O	2.12	0.48
1:A:146:ASN:O	1:A:220:LEU:HD21	2.14	0.48
1:B:348:GLN:HG3	1:B:349:THR:N	2.27	0.48
1:B:377:ASN:C	1:B:379:LEU:H	2.17	0.48
1:B:50:THR:N	1:B:53:ASP:OD2	2.43	0.48
1:A:280:LYS:O	1:A:280:LYS:HD3	2.13	0.48
1:B:34:ILE:HD13	1:B:90:PHE:CG	2.48	0.48
1:A:42:PRO:HB2	1:A:79:LYS:HB3	1.95	0.48
1:B:129:SER:C	1:B:180:ILE:HD11	2.34	0.47
1:B:134:LYS:HE2	1:B:140:GLN:HG3	1.96	0.47
1:B:28:PHE:CE2	1:B:134:LYS:HE3	2.50	0.47
1:B:35:MET:SD	1:B:38:ILE:HD12	2.55	0.47
1:A:129:SER:HA	1:A:180:ILE:HG13	1.97	0.47
1:B:194:ASN:CB	4:B:885:HOH:O	2.62	0.47
1:B:290:VAL:HG11	1:B:296:ASN:HD21	1.78	0.47
1:A:26:GLN:NE2	4:A:525:HOH:O	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:PHE:CZ	1:A:351:ILE:HD11	2.50	0.47
1:B:261:ILE:N	4:B:859:HOH:O	2.44	0.47
1:A:44:ARG:NE	1:A:152:ALA:O	2.39	0.47
1:A:156:LEU:HD23	1:A:189:TYR:CE2	2.50	0.47
1:A:329:LYS:HD3	4:A:477:HOH:O	2.13	0.47
1:B:256:GLY:C	1:B:258:ASP:N	2.68	0.47
1:B:50:THR:CG2	1:B:52:GLN:HE21	2.25	0.47
1:B:361:ASN:HD22	1:B:364:ASN:HB2	1.79	0.47
1:A:52:GLN:HG3	4:A:452:HOH:O	2.13	0.46
1:B:42:PRO:O	1:B:79:LYS:HG2	2.16	0.46
1:B:54:PHE:CZ	1:B:154:PRO:HD3	2.51	0.46
1:B:81:ARG:HH11	1:B:81:ARG:HG3	1.80	0.46
1:A:93:ILE:O	1:A:94:ASN:HB3	2.16	0.46
1:A:261:ILE:HD12	1:A:261:ILE:N	2.30	0.46
1:A:289:GLN:HE21	1:A:290:VAL:H	1.63	0.46
1:B:130:ALA:HB2	1:B:144:LEU:HD22	1.96	0.46
1:B:246:THR:CG2	1:B:250:GLU:HB2	2.45	0.46
1:B:290:VAL:HG12	1:B:296:ASN:HD21	1.81	0.46
1:A:409:ARG:O	4:A:517:HOH:O	2.20	0.46
1:B:351:ILE:CG2	1:B:352:GLY:N	2.79	0.46
1:A:121:ASN:HD22	1:A:121:ASN:C	2.17	0.45
1:A:312:ALA:O	1:B:16:THR:HG21	2.16	0.45
1:B:201:PHE:CE2	1:B:345:LYS:HE2	2.52	0.45
1:A:72:TYR:O	1:A:73:LEU:HB2	2.16	0.45
1:A:36:LYS:O	1:A:37:ASN:HB2	2.16	0.45
1:A:92:ARG:O	1:A:95:ASN:HB3	2.17	0.45
1:A:111:PRO:HG3	1:A:220:LEU:HD23	1.99	0.45
1:A:171:MET:HA	1:A:172:PRO:HD2	1.87	0.45
1:A:93:ILE:O	1:A:94:ASN:CB	2.62	0.44
1:A:309:ASP:OD2	1:A:319:ASN:HB2	2.17	0.44
1:B:50:THR:HG22	1:B:52:GLN:HB2	1.98	0.44
1:A:280:LYS:HA	1:A:280:LYS:HZ2	1.81	0.44
1:B:67:TYR:HE2	1:B:189:TYR:HH	1.64	0.44
1:B:391:ASN:N	1:B:392:PRO:HD3	2.33	0.44
1:A:93:ILE:C	1:A:95:ASN:H	2.21	0.44
1:A:312:ALA:HA	1:B:4:ASN:OD1	2.17	0.44
1:B:266:GLN:O	1:B:270:ILE:HG12	2.17	0.44
1:B:153:GLU:HB2	1:B:154:PRO:CD	2.47	0.44
1:B:275:LEU:CG	1:B:279:LYS:HE3	2.46	0.44
1:A:105:GLU:OE2	1:A:329:LYS:HD2	2.18	0.44
1:B:214:ILE:HG21	1:B:253:THR:HG23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:ASN:N	1:B:168:ASN:ND2	2.65	0.43
1:B:380:LYS:HD2	1:B:381:VAL:N	2.32	0.43
1:B:14:ASP:OD1	1:B:135:PHE:HB3	2.18	0.43
1:B:36:LYS:NZ	1:B:107:SER:O	2.51	0.43
1:A:46:VAL:O	1:A:47:ILE:C	2.57	0.43
1:A:88:LYS:HG3	1:A:370:ILE:HD11	2.00	0.43
1:B:112:TYR:HA	1:B:302:PHE:CE1	2.54	0.43
1:A:347:ARG:HG2	4:A:523:HOH:O	2.18	0.43
1:B:395:ILE:HG22	1:B:396:THR:N	2.34	0.43
1:A:71:ASN:HB2	4:A:516:HOH:O	2.18	0.43
1:A:355:LYS:HD3	1:A:357:PHE:CZ	2.54	0.43
1:B:122:GLN:NE2	1:B:122:GLN:O	2.51	0.42
1:B:201:PHE:CG	1:B:345:LYS:HG2	2.54	0.42
1:A:309:ASP:O	1:A:316:TYR:HA	2.19	0.42
1:A:408:ILE:O	1:A:408:ILE:HG22	2.19	0.42
1:A:32:PHE:N	1:A:32:PHE:CD2	2.87	0.42
1:B:259:LEU:O	1:B:262:ILE:HD12	2.19	0.42
1:B:7:ASN:HA	1:B:84:LYS:HE3	2.02	0.42
1:B:29:TYR:CD1	1:B:43:GLU:HG3	2.54	0.42
1:B:387:ASN:HB3	1:B:390:LEU:HB2	2.00	0.42
1:A:160:ALA:HA	1:A:212:GLU:OE1	2.19	0.42
1:A:263:THR:H	1:A:266:GLN:NE2	2.17	0.42
1:B:181:ALA:HB1	4:B:868:HOH:O	2.18	0.42
1:B:186:SER:HA	4:B:834:HOH:O	2.18	0.42
1:A:121:ASN:C	1:A:121:ASN:ND2	2.73	0.42
1:A:149:ILE:HD13	1:A:183:VAL:HB	2.02	0.42
1:B:153:GLU:O	1:B:154:PRO:C	2.56	0.42
1:A:182:ILE:N	1:A:182:ILE:HD12	2.35	0.42
1:A:194:ASN:HA	1:A:199:ASN:O	2.20	0.41
1:B:233:THR:HA	1:B:245:GLY:HA3	2.02	0.41
1:B:296:ASN:N	1:B:297:PRO:CD	2.83	0.41
1:A:141:ASP:OD1	1:A:142:ILE:N	2.53	0.41
1:B:174:ASN:ND2	1:B:224:LYS:HB2	2.35	0.41
1:A:226:ILE:HD11	1:A:273:ASN:ND2	2.35	0.41
1:A:261:ILE:HD12	1:A:261:ILE:H	1.86	0.41
1:B:227:THR:HG22	1:B:270:ILE:HD12	2.03	0.41
1:B:342:PHE:HB3	1:B:344:VAL:HG23	2.02	0.41
1:A:173:SER:HA	1:A:178:GLY:HA2	2.02	0.41
1:B:104:GLU:OE2	1:B:108:LYS:HD3	2.20	0.41
1:B:232:ILE:C	1:B:233:THR:O	2.59	0.41
1:B:403:LEU:HD13	1:B:403:LEU:C	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:GLU:CB	1:A:154:PRO:CD	2.98	0.41
1:A:9:ASN:O	1:A:10:ASP:C	2.58	0.41
1:A:81:ARG:HG3	1:A:81:ARG:NH1	2.34	0.41
1:A:6:PHE:CD2	1:A:32:PHE:HB3	2.56	0.41
1:A:33:ASN:HB2	1:A:39:TRP:CH2	2.55	0.41
1:B:194:ASN:HB2	4:B:885:HOH:O	2.21	0.41
1:B:209:LEU:O	1:B:213:LEU:HG	2.21	0.41
1:A:232:ILE:CD1	1:A:261:ILE:CG2	2.99	0.41
1:B:171:MET:HA	1:B:172:PRO:HD2	1.87	0.41
1:A:375:ASN:HD22	1:A:385:GLY:H	1.68	0.40
1:B:156:LEU:HD23	1:B:189:TYR:CE2	2.56	0.40
1:B:200:GLU:HG2	4:B:885:HOH:O	2.21	0.40
1:A:224:LYS:HB3	1:A:228:THR:HG23	2.03	0.40
1:B:75:SER:O	1:B:78:GLU:HB2	2.22	0.40
1:B:289:GLN:NE2	4:B:828:HOH:O	2.53	0.40
1:B:173:SER:HA	1:B:178:GLY:HA2	2.02	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	396/420 (94%)	368 (93%)	22 (6%)	6 (2%)	10 17
1	B	381/420 (91%)	346 (91%)	25 (7%)	10 (3%)	5 7
All	All	777/840 (92%)	714 (92%)	47 (6%)	16 (2%)	7 10

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	290	VAL
1	B	257	THR

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Mol	Chain	Res	Type
1	B	375	ASN
1	B	378	ASN
1	B	382	ASN
1	A	36	LYS
1	A	51	PRO
1	A	361	ASN
1	A	407	ILE
1	A	408	ILE
1	B	291	SER
1	B	361	ASN
1	B	408	ILE
1	B	45	ASN
1	A	141	ASP
1	B	407	ILE

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	358/377 (95%)	348 (97%)	10 (3%)	43	68
1	B	346/377 (92%)	332 (96%)	14 (4%)	31	54
All	All	704/754 (93%)	680 (97%)	24 (3%)	37	61

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	121	ASN
1	A	137	ASN
1	A	268	ASN
1	A	280	LYS
1	A	298	TYR
1	A	345	LYS
1	A	349	THR
1	A	410	PHE

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Mol	Chain	Res	Type
1	A	411	CYS
1	B	2	LYS
1	B	18	LEU
1	B	26	GLN
1	B	50	THR
1	B	122	GLN
1	B	154	PRO
1	B	161	SER
1	B	167	ARG
1	B	195	ASP
1	B	196	ASN
1	B	295	LEU
1	B	298	TYR
1	B	348	GLN
1	B	380	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	A	71	ASN
1	A	121	ASN
1	A	163	ASN
1	A	174	ASN
1	A	194	ASN
1	A	266	GLN
1	A	273	ASN
1	A	289	GLN
1	A	296	ASN
1	A	348	GLN
1	A	353	GLN
1	A	375	ASN
1	A	378	ASN
1	A	389	ASN
1	B	52	GLN
1	B	95	ASN
1	B	122	GLN
1	B	124	HIS
1	B	140	GLN
1	B	168	ASN
1	B	174	ASN
1	B	199	ASN

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Mol	Chain	Res	Type
1	B	266	GLN
1	B	289	GLN
1	B	296	ASN
1	B	375	ASN
1	B	378	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 monosaccharides in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 6 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	400/420 (95%)	-0.19	6 (1%) 73 76	15, 30, 49, 74	0
1	B	387/420 (92%)	0.05	14 (3%) 42 46	21, 38, 61, 87	0
All	All	787/840 (93%)	-0.07	20 (2%) 57 61	15, 34, 59, 87	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	245	GLY	4.0
1	B	261	ILE	3.2
1	B	246	THR	3.2
1	A	233	THR	3.1
1	B	378	ASN	3.0
1	B	262	ILE	2.9
1	B	248	ILE	2.8
1	A	409	ARG	2.8
1	B	381	VAL	2.7
1	B	290	VAL	2.7
1	B	232	ILE	2.6
1	B	390	LEU	2.4
1	B	230	TYR	2.4
1	A	411	CYS	2.3
1	B	231	THR	2.3
1	B	233	THR	2.3
1	A	167	ARG	2.2
1	B	264	SER	2.0
1	A	52	GLN	2.0
1	B	263	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 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, 95^{th} 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	CL	B	427	1/1	0.51	0.14	43,43,43,43	0
3	CL	A	424	1/1	0.70	0.33	66,66,66,66	0
3	CL	A	426	1/1	0.94	0.20	31,31,31,31	0
3	CL	A	425	1/1	0.95	0.09	36,36,36,36	0
2	ZN	B	822	1/1	0.96	0.07	46,46,46,46	0
2	ZN	A	423	1/1	0.99	0.07	32,32,32,32	0

6.5 Other polymers (i)

There are no such residues in this entry.