



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 29, 2020 – 04:02 PM BST

PDB ID : 6SXH  
Title : Crystal structure of the accessory translocation ATPase, SecA2, from *Clostridium difficile*  
Authors : Lindic, N.; Loboda, J.; Usenik, A.; Turk, D.  
Deposited on : 2019-09-26  
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](mailto: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 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	:	2.14.6
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)	:	2.14.6

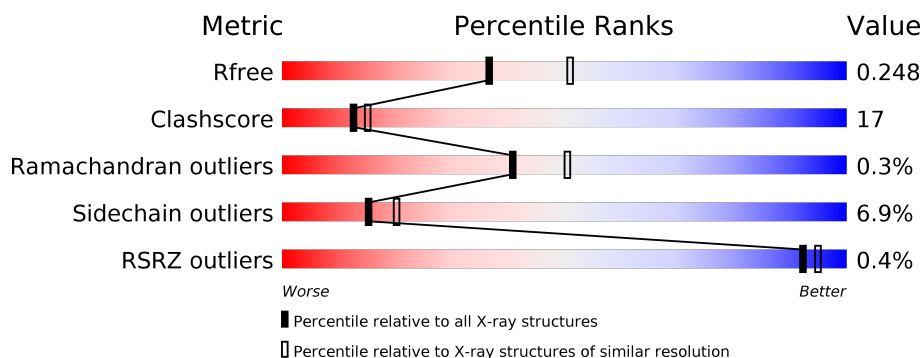
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


The reported resolution of this entry is 2.30 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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	786	 64% 32% . .

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8566 atoms, of which 2077 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 Protein translocase subunit SecA 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	769	Total	C	H	N	O	S	1604	0	0
			7621	3894	1447	1053	1205	22			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	SER	-	expression tag	UNP Q183M9
A	-15	ASN	-	expression tag	UNP Q183M9
A	-14	ALA	-	expression tag	UNP Q183M9
A	-13	ALA	-	expression tag	UNP Q183M9
A	-12	ALA	-	expression tag	UNP Q183M9

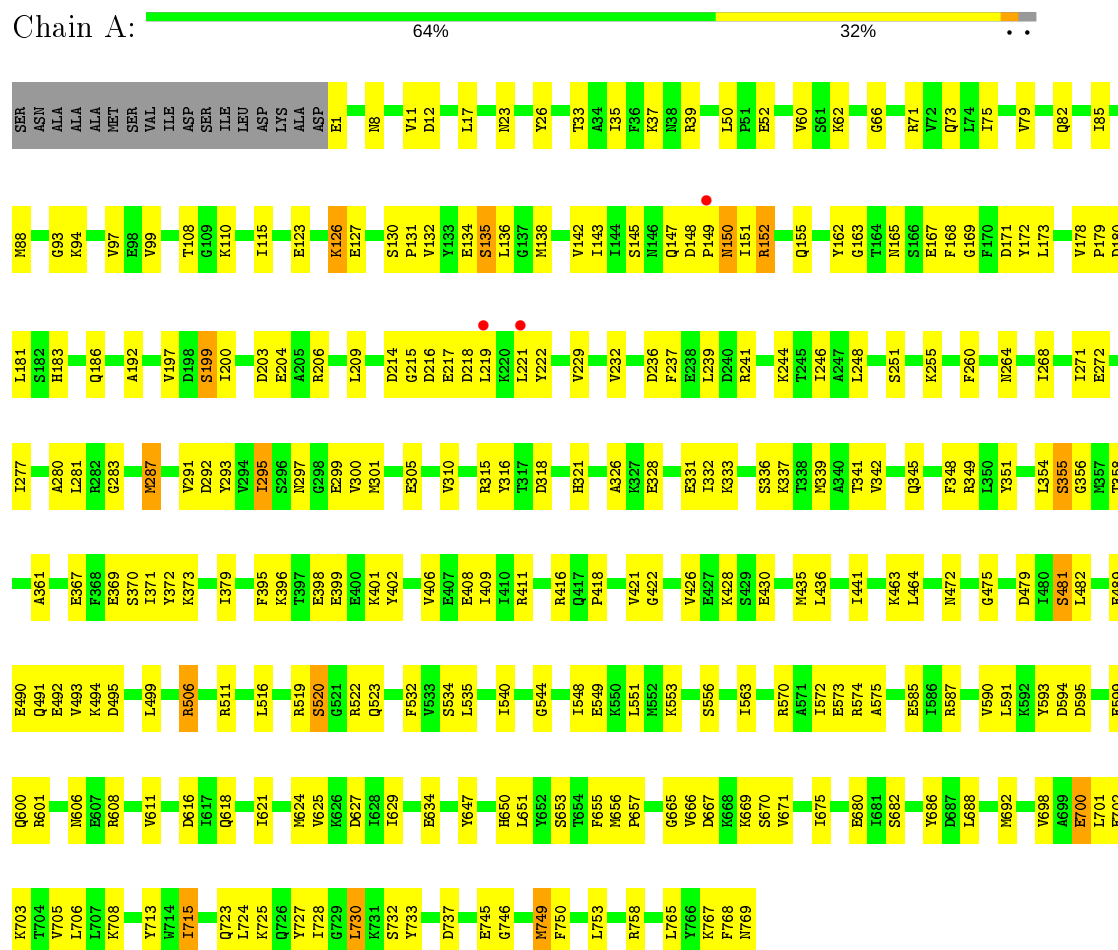
- Molecule 2 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	315	Total	H	O	630	0
			945	630	315		

### 3 Residue-property plots

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: Protein translocase subunit SecA 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.99 Å 96.85 Å 114.74 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.55 – 2.30 48.42 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.55-2.30) 99.8 (48.42-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.25 (at 2.29 Å)	Xtriage
Refinement program	MAIN	Depositor
R, $R_{free}$	0.243 , 0.274 0.241 , 0.248	Depositor DCC
$R_{free}$ test set	2100 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.6	Xtriage
Anisotropy	0.727	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 26.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8566	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.74	0/6258	0.91	0/8405

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	6174	1447	6259	202	0
2	A	315	630	0	22	0
All	All	6489	2077	6259	202	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 17.

All (202) 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:629:ILE:HG22	1:A:675:ILE:HG23	1.35	1.07
1:A:218:ASP:O	1:A:222:TYR:CD1	2.25	0.89
1:A:611:VAL:HG21	1:A:706:LEU:HD22	1.56	0.88
1:A:426:VAL:O	1:A:430:GLU:HG2	1.75	0.86

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:VAL:HG21	1:A:183:HIS:HB3	1.62	0.82
1:A:399:GLU:HG3	2:A:967:HOH:O	1.82	0.79
1:A:535:LEU:HA	1:A:540:ILE:HG21	1.65	0.78
1:A:248:LEU:HD13	1:A:281:LEU:HD11	1.65	0.77
1:A:506:ARG:HH22	1:A:573:GLU:CD	1.88	0.75
1:A:246:ILE:HD13	1:A:281:LEU:HB3	1.71	0.73
1:A:482:LEU:O	1:A:490:GLU:HB2	1.89	0.73
1:A:301:MET:HB3	1:A:310:VAL:HG21	1.70	0.72
1:A:11:VAL:HG11	1:A:71:ARG:HG2	1.72	0.72
1:A:305:GLU:HG3	2:A:988:HOH:O	1.89	0.71
1:A:315:ARG:HB2	1:A:723:GLN:HE22	1.53	0.71
1:A:315:ARG:NH2	1:A:333:LYS:HD3	2.05	0.71
1:A:656:MET:HE1	1:A:682:SER:CB	2.20	0.71
1:A:358:THR:HG21	1:A:361:ALA:HB2	1.73	0.71
1:A:301:MET:HB3	1:A:310:VAL:CG2	2.21	0.70
1:A:653:SER:HB2	2:A:1046:HOH:O	1.90	0.70
1:A:178:VAL:CG2	1:A:183:HIS:HB3	2.21	0.70
1:A:745:GLU:O	1:A:749:MET:HE2	1.92	0.69
1:A:173:LEU:HD21	1:A:342:VAL:HG12	1.74	0.69
1:A:316:TYR:H	1:A:321:HIS:HD2	1.41	0.68
1:A:115:ILE:HD13	1:A:168:PHE:CD1	2.28	0.68
1:A:316:TYR:H	1:A:321:HIS:CD2	2.12	0.68
1:A:60:VAL:HG21	1:A:136:LEU:HD21	1.75	0.68
1:A:199:SER:O	1:A:204:GLU:HG3	1.94	0.67
1:A:540:ILE:HG13	1:A:572:ILE:HD12	1.76	0.66
1:A:315:ARG:CZ	1:A:333:LYS:HD3	2.26	0.66
1:A:79:VAL:HG21	1:A:379:ILE:HD11	1.79	0.65
1:A:341:THR:HG22	2:A:1057:HOH:O	1.97	0.65
1:A:371:ILE:HB	1:A:372:TYR:CD1	2.32	0.64
1:A:178:VAL:HG21	1:A:183:HIS:CB	2.28	0.63
1:A:549:GLU:O	1:A:553:LYS:HG3	1.99	0.63
1:A:621:ILE:CD1	1:A:702:GLU:HB3	2.29	0.62
1:A:688:LEU:O	1:A:692:MET:HG2	1.99	0.62
1:A:724:LEU:HD22	1:A:749:MET:CE	2.30	0.62
1:A:354:LEU:HD12	1:A:355:SER:H	1.65	0.61
1:A:62:LYS:O	1:A:66:GLY:HA2	2.00	0.61
1:A:401:LYS:HD3	1:A:532:PHE:O	1.99	0.61
1:A:656:MET:HE1	1:A:682:SER:CA	2.30	0.60
1:A:629:ILE:CG2	1:A:675:ILE:HG23	2.20	0.60
1:A:618:GLN:HG2	1:A:702:GLU:OE1	2.01	0.60
1:A:93:GLY:O	1:A:97:VAL:HG23	2.01	0.60

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:GLY:HA2	1:A:287:MET:HG3	1.84	0.60
1:A:396:LYS:HG3	1:A:563:ILE:HD13	1.83	0.59
1:A:606:ASN:HB3	2:A:1065:HOH:O	2.01	0.59
1:A:180:ASP:O	1:A:183:HIS:HB2	2.03	0.59
1:A:475:GLY:O	1:A:519:ARG:NH2	2.36	0.58
1:A:621:ILE:HD13	1:A:702:GLU:HB3	1.86	0.58
1:A:656:MET:HE1	1:A:682:SER:HA	1.86	0.58
1:A:131:PRO:O	1:A:135:SER:HB2	2.03	0.58
1:A:150:ASN:HD22	1:A:151:ILE:HD12	1.68	0.58
1:A:369:GLU:O	1:A:373:LYS:HD3	2.04	0.58
1:A:108:THR:HB	1:A:110:LYS:HE2	1.85	0.57
1:A:271:ILE:HG21	1:A:708:LYS:HD2	1.86	0.57
1:A:724:LEU:HD22	1:A:749:MET:HE1	1.85	0.57
1:A:237:PHE:HE1	1:A:239:LEU:HD23	1.70	0.57
1:A:206:ARG:NH2	1:A:725:LYS:NZ	2.52	0.57
1:A:523:GLN:HA	2:A:882:HOH:O	2.05	0.57
1:A:326:ALA:HB2	1:A:332:ILE:HD11	1.87	0.56
1:A:152:ARG:HH12	1:A:167:GLU:HB3	1.71	0.56
1:A:179:PRO:HG3	2:A:1059:HOH:O	2.05	0.56
1:A:149:PRO:HA	1:A:152:ARG:HB2	1.87	0.56
1:A:767:LYS:HE2	2:A:866:HOH:O	2.05	0.55
1:A:401:LYS:HD3	1:A:532:PHE:C	2.26	0.55
1:A:608:ARG:HH11	1:A:608:ARG:HG2	1.71	0.55
1:A:337:LYS:CE	1:A:715:ILE:HD11	2.36	0.55
1:A:209:LEU:O	1:A:341:THR:HA	2.08	0.54
1:A:655:PHE:CE1	1:A:656:MET:HE3	2.42	0.54
1:A:178:VAL:CG2	1:A:183:HIS:CB	2.85	0.54
1:A:489:GLU:O	1:A:493:VAL:HG23	2.08	0.54
1:A:600:GLN:HG2	1:A:750:PHE:CD1	2.42	0.54
1:A:341:THR:HG21	2:A:1071:HOH:O	2.08	0.54
1:A:85:ILE:HA	1:A:356:GLY:O	2.08	0.53
1:A:540:ILE:HG13	1:A:572:ILE:CD1	2.38	0.53
1:A:85:ILE:HG12	1:A:197:VAL:HG11	1.90	0.53
1:A:650:HIS:HD2	2:A:1047:HOH:O	1.91	0.53
1:A:315:ARG:HB2	1:A:723:GLN:NE2	2.23	0.53
1:A:519:ARG:HD2	2:A:1038:HOH:O	2.08	0.52
1:A:535:LEU:HA	1:A:540:ILE:CG2	2.38	0.52
1:A:192:ALA:HB2	1:A:351:TYR:CD2	2.45	0.52
1:A:671:VAL:O	1:A:675:ILE:HG13	2.09	0.52
1:A:75:ILE:HG22	1:A:379:ILE:HD12	1.91	0.51
1:A:395:PHE:O	1:A:534:SER:HA	2.11	0.51

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:VAL:HA	1:A:82:GLN:NE2	2.25	0.51
1:A:402:TYR:HB3	1:A:435:MET:HE1	1.92	0.51
1:A:241:ARG:O	1:A:244:LYS:HE2	2.10	0.50
1:A:203:ASP:CG	1:A:511:ARG:HH22	2.15	0.50
1:A:88:MET:O	1:A:94:LYS:HE3	2.11	0.50
1:A:535:LEU:HD22	1:A:548:ILE:HD12	1.93	0.50
1:A:745:GLU:O	1:A:749:MET:CE	2.58	0.50
1:A:206:ARG:HH21	1:A:725:LYS:HZ3	1.58	0.50
1:A:544:GLY:HA2	1:A:575:ALA:HB2	1.95	0.49
1:A:172:TYR:HA	1:A:186:GLN:OE1	2.11	0.49
1:A:713:TYR:HB3	1:A:753:LEU:CD1	2.41	0.49
1:A:624:MET:HE3	1:A:758:ARG:HA	1.94	0.49
1:A:436:LEU:HB3	1:A:441:ILE:HB	1.94	0.49
1:A:192:ALA:HB2	1:A:351:TYR:CE2	2.48	0.49
1:A:608:ARG:HG2	1:A:608:ARG:NH1	2.28	0.49
1:A:345:GLN:O	1:A:349:ARG:HG3	2.13	0.48
1:A:706:LEU:C	1:A:706:LEU:HD23	2.33	0.48
1:A:358:THR:CG2	1:A:361:ALA:HB2	2.42	0.48
1:A:348:PHE:O	1:A:354:LEU:HD22	2.13	0.48
1:A:206:ARG:NH2	1:A:725:LYS:HZ3	2.11	0.48
1:A:746:GLY:HA2	1:A:749:MET:CE	2.44	0.48
1:A:656:MET:CE	1:A:682:SER:CB	2.90	0.48
1:A:371:ILE:HB	1:A:372:TYR:CE1	2.49	0.48
1:A:169:GLY:HA3	1:A:209:LEU:HD13	1.95	0.47
1:A:229:VAL:HG21	1:A:280:ALA:HB3	1.96	0.47
1:A:152:ARG:NH2	1:A:171:ASP:OD1	2.44	0.47
1:A:666:VAL:O	1:A:667:ASP:HB2	2.14	0.47
1:A:300:VAL:HB	1:A:321:HIS:CE1	2.48	0.47
1:A:260:PHE:CE1	1:A:277:ILE:HD11	2.49	0.47
1:A:337:LYS:NZ	1:A:715:ILE:HD11	2.29	0.47
1:A:209:LEU:HB2	1:A:342:VAL:O	2.14	0.47
1:A:79:VAL:HA	1:A:82:GLN:HE21	1.79	0.47
1:A:621:ILE:HD13	1:A:702:GLU:CD	2.35	0.47
1:A:593:TYR:HE2	1:A:728:ILE:HD11	1.80	0.46
1:A:232:VAL:HG13	1:A:236:ASP:HB2	1.95	0.46
1:A:408:GLU:CD	1:A:411:ARG:HH12	2.18	0.46
1:A:665:GLY:O	1:A:669:LYS:HG3	2.14	0.46
1:A:237:PHE:CE1	1:A:239:LEU:HD23	2.51	0.46
1:A:702:GLU:HA	1:A:765:LEU:HD21	1.98	0.46
1:A:595:ASP:HB3	2:A:1019:HOH:O	2.15	0.46
1:A:251:SER:O	1:A:255:LYS:HG3	2.16	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:499:LEU:HB3	1:A:520:SER:HB2	1.98	0.45
1:A:625:VAL:O	1:A:629:ILE:HG13	2.17	0.45
1:A:492:GLU:O	1:A:495:ASP:HB2	2.17	0.45
1:A:337:LYS:HE2	1:A:715:ILE:HD11	1.99	0.45
1:A:143:ILE:HD11	1:A:168:PHE:CE1	2.52	0.45
1:A:418:PRO:HB3	1:A:463:LYS:O	2.16	0.45
1:A:656:MET:CE	1:A:682:SER:HB2	2.47	0.45
1:A:271:ILE:HG23	1:A:272:GLU:N	2.31	0.45
1:A:295:ILE:HA	1:A:299:GLU:O	2.17	0.45
1:A:409:ILE:HD11	1:A:421:VAL:HG21	1.99	0.45
1:A:621:ILE:HD13	1:A:702:GLU:CG	2.47	0.45
1:A:17:LEU:CD1	1:A:52:GLU:HA	2.47	0.44
1:A:200:ILE:HA	1:A:200:ILE:HD13	1.70	0.44
1:A:551:LEU:HA	1:A:551:LEU:HD23	1.77	0.44
1:A:724:LEU:HD22	1:A:749:MET:HE3	1.99	0.44
1:A:416:ARG:O	1:A:464:LEU:O	2.35	0.44
1:A:591:LEU:HA	2:A:920:HOH:O	2.17	0.44
1:A:108:THR:CB	1:A:110:LYS:HE2	2.47	0.44
1:A:26:TYR:HE2	1:A:135:SER:HG	1.64	0.44
1:A:271:ILE:CG2	1:A:708:LYS:HD2	2.47	0.44
1:A:703:LYS:HE3	2:A:1030:HOH:O	2.18	0.44
1:A:143:ILE:HD11	1:A:168:PHE:CD1	2.53	0.44
1:A:126:LYS:HG3	1:A:127:GLU:N	2.32	0.44
1:A:181:LEU:HD23	1:A:181:LEU:O	2.17	0.44
1:A:219:LEU:HD12	1:A:219:LEU:HA	1.89	0.44
1:A:332:ILE:HG22	1:A:333:LYS:N	2.31	0.44
1:A:12:ASP:OD2	1:A:71:ARG:NH1	2.46	0.43
1:A:422:GLY:HA3	1:A:516:LEU:CD2	2.47	0.43
1:A:650:HIS:CD2	2:A:1047:HOH:O	2.68	0.43
1:A:209:LEU:HD23	1:A:209:LEU:HA	1.70	0.43
1:A:490:GLU:O	1:A:494:LYS:HB2	2.17	0.43
1:A:293:TYR:HE1	1:A:328:GLU:OE1	2.02	0.43
1:A:656:MET:HE1	1:A:682:SER:HB3	1.97	0.43
1:A:656:MET:CE	1:A:682:SER:HA	2.47	0.43
1:A:647:TYR:CZ	1:A:651:LEU:HD11	2.52	0.43
1:A:708:LYS:HD3	2:A:930:HOH:O	2.18	0.43
1:A:481:SER:HB2	2:A:842:HOH:O	2.18	0.43
1:A:570:ARG:NH1	2:A:816:HOH:O	2.51	0.43
1:A:402:TYR:O	1:A:406:VAL:HG23	2.19	0.43
1:A:656:MET:HA	1:A:657:PRO:HD3	1.94	0.42
1:A:725:LYS:O	1:A:728:ILE:HG13	2.19	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:688:LEU:HD21	2:A:1100:HOH:O	2.19	0.42
1:A:138:MET:HE2	1:A:138:MET:HB2	1.70	0.42
1:A:94:LYS:HD2	2:A:1002:HOH:O	2.18	0.42
1:A:99:VAL:HG11	1:A:132:VAL:HG21	2.02	0.42
1:A:33:THR:O	1:A:37:LYS:HG3	2.19	0.42
1:A:142:VAL:HA	1:A:162:TYR:O	2.19	0.42
1:A:506:ARG:NH2	1:A:573:GLU:OE1	2.53	0.42
1:A:698:VAL:O	1:A:702:GLU:HG3	2.20	0.42
1:A:214:ASP:OD2	1:A:215:GLY:N	2.53	0.42
1:A:700:GLU:HG3	1:A:701:LEU:N	2.35	0.42
1:A:123:GLU:O	1:A:127:GLU:HB2	2.21	0.41
1:A:342:VAL:HG23	1:A:601:ARG:NH2	2.35	0.41
1:A:50:LEU:HD23	1:A:50:LEU:C	2.41	0.41
1:A:519:ARG:HG2	1:A:519:ARG:NH1	2.35	0.41
1:A:130:SER:HB3	1:A:134:GLU:OE2	2.21	0.41
1:A:218:ASP:O	1:A:222:TYR:HD1	1.92	0.41
1:A:367:GLU:OE2	1:A:587:ARG:HD2	2.19	0.41
1:A:35:ILE:O	1:A:39:ARG:HG3	2.21	0.41
1:A:624:MET:CE	1:A:758:ARG:HA	2.50	0.41
1:A:8:ASN:HD22	1:A:71:ARG:HH21	1.68	0.41
1:A:115:ILE:HA	1:A:163:GLY:O	2.19	0.41
1:A:769:ASN:ND2	2:A:820:HOH:O	2.53	0.41
1:A:152:ARG:NH2	1:A:171:ASP:CG	2.74	0.41
1:A:398:GLU:OE2	1:A:428:LYS:NZ	2.50	0.41
1:A:594:ASP:HA	2:A:886:HOH:O	2.20	0.41
1:A:217:GLU:O	1:A:221:LEU:HB2	2.20	0.41
1:A:73:GLN:HG2	1:A:97:VAL:HG22	2.03	0.41
1:A:151:ILE:O	1:A:155:GLN:HG2	2.21	0.41
1:A:540:ILE:CG1	1:A:572:ILE:HD12	2.49	0.41
1:A:268:ILE:HG12	1:A:705:VAL:HG22	2.03	0.40
1:A:152:ARG:NH2	1:A:171:ASP:OD2	2.54	0.40
1:A:206:ARG:HD2	1:A:590:VAL:HG21	2.04	0.40
1:A:291:VAL:HG23	1:A:292:ASP:N	2.35	0.40
1:A:11:VAL:CG1	1:A:71:ARG:HG2	2.47	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	767/786 (98%)	727 (95%)	38 (5%)	2 (0%)	41	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	616	ASP
1	A	730	LEU

### 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	677/690 (98%)	630 (93%)	47 (7%)	15	20

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLU
1	A	23	ASN
1	A	126	LYS
1	A	135	SER
1	A	145	SER
1	A	147	GLN
1	A	148	ASP
1	A	150	ASN
1	A	152	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	165	ASN
1	A	199	SER
1	A	216	ASP
1	A	264	ASN
1	A	287	MET
1	A	295	ILE
1	A	297	ASN
1	A	318	ASP
1	A	331	GLU
1	A	336	SER
1	A	339	MET
1	A	355	SER
1	A	370	SER
1	A	472	ASN
1	A	479	ASP
1	A	481	SER
1	A	491	GLN
1	A	506	ARG
1	A	520	SER
1	A	522	ARG
1	A	556	SER
1	A	574	ARG
1	A	585	GLU
1	A	599	GLU
1	A	627	ASP
1	A	634	GLU
1	A	670	SER
1	A	680	GLU
1	A	686	TYR
1	A	700	GLU
1	A	715	ILE
1	A	727	TYR
1	A	730	LEU
1	A	732	SER
1	A	733	TYR
1	A	737	ASP
1	A	749	MET
1	A	768	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	82	GLN
1	A	147	GLN
1	A	150	ASN
1	A	154	GLN
1	A	183	HIS
1	A	284	HIS
1	A	321	HIS
1	A	439	GLN
1	A	447	ASN
1	A	560	ASN
1	A	609	ASN
1	A	650	HIS
1	A	723	GLN
1	A	756	ASN
1	A	769	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](#)

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	759/786 (96%)	-0.38	3 (0%) 92 95	10, 23, 68, 139	27 (3%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	149	PRO	2.8
1	A	219	LEU	2.3
1	A	221	LEU	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](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.