



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:58 am BST

PDB ID : 4OW6
Title : Crystal structure of Diphtheria Toxin at acidic pH
Authors : Leka, O.; Vallese, F.; Pirazzini, M.; Berto, P.; Montecucco, C.; Zanutti, G.
Deposited on : 2014-01-31
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

<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 : 2.11
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) : 2.11

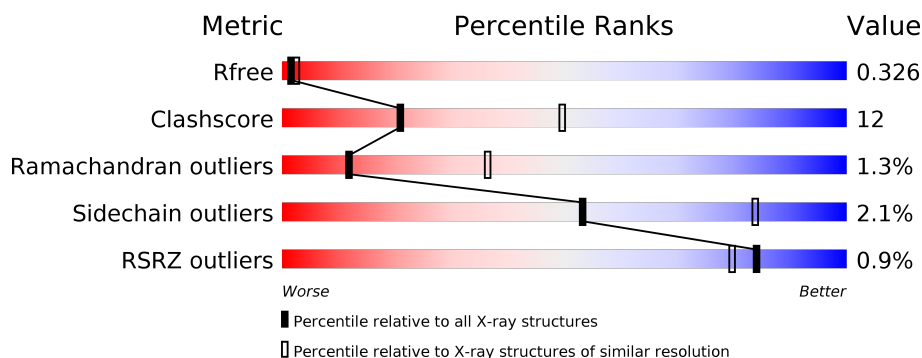
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 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	535	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 64%, yellow 24%, grey 11%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> % 64% 24% • 11% </div> </div>
1	B	535	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 65%, yellow 24%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> % 65% 24% • 9% </div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7374 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 Diphtheria toxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	477	Total	C	N	O	S	0	0	0
			3648	2298	620	719	11			
1	B	488	Total	C	N	O	S	0	0	0
			3726	2347	632	736	11			

R460	C461
I464	
F470	
K474	S475
P476	V477
G480	
H484	
S495	S496
E497	K498
I499	H500
S501	V502
E503	I504
S505	S506
Y514	Q515
K516	T517
V518	D519
H520	T521
K522	
L527	S528
L529	
I533	K534
	S535

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.44Å 141.28Å 176.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.13 – 2.80 45.13 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.2 (45.13-2.80) 94.2 (45.13-2.80)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.238 , 0.322 0.245 , 0.326	Depositor DCC
R_{free} test set	1452 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	7374	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% 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 [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	0.52	0/3720	0.73	2/5041 (0.0%)
1	B	0.49	0/3798	0.72	1/5147 (0.0%)
All	All	0.50	0/7518	0.72	3/10188 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	466	GLY	N-CA-C	-5.86	98.45	113.10
1	B	258	PRO	N-CA-CB	5.83	110.30	103.30
1	A	138	LEU	CA-CB-CG	5.80	128.65	115.30

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	3648	0	3581	90	0
1	B	3726	0	3654	88	0
All	All	7374	0	7235	178	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 12.

All (178) 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:106:LEU:O	1:B:126:ARG:NH2	1.98	0.97
1:A:45:ASN:O	1:A:51:LYS:NZ	2.13	0.81
1:B:515:GLN:OE1	1:B:522:LYS:NZ	2.14	0.81
1:B:271:PRO:HB3	1:B:274:ALA:HA	1.64	0.79
1:B:499:ILE:HD11	1:B:503:GLU:HB2	1.67	0.77
1:A:202:ILE:HD13	1:A:321:VAL:HG21	1.66	0.76
1:A:437:ILE:HG21	1:A:440:LYS:HD2	1.67	0.76
1:A:436:THR:O	1:A:467:ASP:HB3	1.86	0.76
1:B:400:THR:HG22	1:B:402:GLU:H	1.49	0.76
1:A:86:PRO:O	1:A:133:ARG:NH2	2.20	0.74
1:A:391:HIS:HB3	1:A:527:LEU:HG	1.70	0.73
1:B:349:GLU:O	1:B:351:VAL:N	2.22	0.73
1:B:445:LYS:HD2	1:B:461:CYS:HB2	1.71	0.72
1:B:411:GLN:HG2	1:B:496:SER:HB2	1.71	0.71
1:A:422:ALA:HB3	1:A:480:GLY:HA2	1.79	0.64
1:B:497:GLU:HG3	1:B:498:LYS:H	1.63	0.64
1:A:320:ALA:HB3	1:A:322:HIS:NE2	2.13	0.63
1:A:396:VAL:HG12	1:A:529:LEU:HD13	1.81	0.63
1:A:459:MET:HG2	1:A:471:CYS:HB3	1.80	0.63
1:A:337:SER:OG	1:A:356:ALA:O	2.14	0.62
1:B:402:GLU:HA	1:B:405:ILE:HD12	1.80	0.62
1:B:37:LYS:NZ	1:B:43:GLN:HB3	2.14	0.62
1:B:500:HIS:CD2	1:B:501:SER:N	2.69	0.61
1:A:408:THR:HG22	1:A:409:GLY:H	1.66	0.61
1:A:37:LYS:HE2	1:A:153:TRP:CD1	2.35	0.61
1:A:437:ILE:HD11	1:A:504:ILE:HG22	1.83	0.61
1:A:344:ILE:HG21	1:A:351:VAL:HB	1.83	0.60
1:B:396:VAL:HG22	1:B:420:ILE:HG12	1.82	0.60
1:A:173:ARG:HH21	1:A:207:ASP:HB2	1.66	0.60
1:B:349:GLU:C	1:B:351:VAL:H	2.05	0.60
1:B:337:SER:OG	1:B:356:ALA:O	2.19	0.59
1:A:425:THR:O	1:A:427:LEU:HD13	2.02	0.59
1:B:398:TRP:NE1	1:B:529:LEU:HG	2.18	0.58
1:A:314:MET:SD	1:A:371:VAL:HG21	2.43	0.58
1:A:323:HIS:HB3	1:A:329:VAL:HG22	1.85	0.58
1:B:267:THR:OG1	1:B:268:GLY:N	2.37	0.58
1:A:408:THR:HG21	1:A:499:ILE:HG22	1.86	0.58
1:B:42:THR:O	1:B:51:LYS:NZ	2.29	0.58
1:A:41:GLY:O	1:A:43:GLN:HG3	2.04	0.57
1:B:464:ILE:HD11	1:B:470:PHE:HB2	1.85	0.57
1:B:518:VAL:HG23	1:B:519:ASP:HB2	1.86	0.57
1:A:500:HIS:O	1:A:502:ASN:N	2.32	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:382:PRO:HG3	1:B:484:HIS:CE1	2.41	0.56
1:A:272:VAL:HG12	1:A:273:PHE:HB2	1.89	0.55
1:A:349:GLU:O	1:A:351:VAL:N	2.39	0.55
1:A:165:ILE:HD11	1:A:170:ARG:HH11	1.72	0.55
1:A:45:ASN:OD1	1:A:47:ASP:N	2.33	0.54
1:B:400:THR:HG22	1:B:402:GLU:N	2.21	0.54
1:B:408:THR:HG21	1:B:499:ILE:HG22	1.89	0.53
1:A:394:TYR:HB2	1:A:527:LEU:HD21	1.91	0.52
1:A:98:ASN:ND2	1:A:413:GLU:O	2.43	0.52
1:B:260:LEU:HD12	1:B:261:SER:HA	1.90	0.52
1:B:460:ARG:HH21	1:B:474:LYS:NZ	2.07	0.52
1:A:272:VAL:O	1:A:274:ALA:N	2.37	0.52
1:B:271:PRO:C	1:B:273:PHE:H	2.13	0.52
1:B:500:HIS:CD2	1:B:501:SER:H	2.28	0.52
1:A:361:VAL:O	1:A:365:ILE:HG12	2.09	0.52
1:A:408:THR:HG22	1:A:409:GLY:N	2.24	0.52
1:B:445:LYS:NZ	1:B:460:ARG:HD3	2.24	0.52
1:B:284:ASN:HD21	1:B:308:PRO:HD2	1.74	0.51
1:B:408:THR:HG22	1:B:409:GLY:H	1.75	0.51
1:B:445:LYS:HZ1	1:B:460:ARG:HD3	1.74	0.51
1:A:72:PRO:HG3	1:A:460:ARG:NH2	2.26	0.51
1:A:212:LYS:O	1:A:216:LYS:HG3	2.12	0.50
1:A:103:LYS:NZ	1:A:112:GLU:O	2.23	0.50
1:B:337:SER:HA	1:B:360:PHE:CE2	2.46	0.50
1:B:56:THR:HB	1:B:58:ASN:H	1.77	0.50
1:B:204:LEU:HD23	1:B:209:ILE:HD11	1.94	0.50
1:A:422:ALA:HB2	1:A:479:VAL:HG23	1.94	0.50
1:B:337:SER:HA	1:B:360:PHE:HE2	1.76	0.50
1:B:7:ASP:OD2	1:B:10:LYS:N	2.43	0.50
1:B:314:MET:HG2	1:B:368:PHE:CE1	2.47	0.50
1:B:445:LYS:NZ	1:B:461:CYS:H	2.10	0.49
1:A:352:ASP:CG	1:A:354:GLY:H	2.16	0.49
1:B:514:TYR:HE1	1:B:516:LYS:HE2	1.77	0.49
1:B:201:CYS:C	1:B:202:ILE:HG13	2.32	0.49
1:B:32:GLN:NE2	1:B:163:LEU:HB2	2.27	0.49
1:A:302:ALA:O	1:A:305:SER:OG	2.23	0.49
1:A:85:TYR:HB3	1:A:133:ARG:NH2	2.27	0.49
1:A:314:MET:HG2	1:A:368:PHE:CZ	2.47	0.49
1:A:437:ILE:HG23	1:A:438:PRO:HD2	1.94	0.49
1:A:449:HIS:CD2	1:A:456:LYS:HE3	2.48	0.49
1:B:94:LEU:HD11	1:B:136:LEU:HD13	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:VAL:O	1:B:287:GLN:HB2	2.13	0.48
1:A:380:TYR:CE2	1:A:419:LYS:HE2	2.48	0.48
1:B:460:ARG:HH21	1:B:474:LYS:HZ3	1.59	0.48
1:A:333:ILE:O	1:A:336:SER:HB3	2.13	0.48
1:A:170:ARG:HH21	1:A:172:LYS:NZ	2.11	0.48
1:A:90:LYS:HB2	1:A:131:ALA:HB3	1.96	0.48
1:A:49:ASP:OD1	1:A:113:PRO:HA	2.14	0.48
1:A:373:ASN:OD1	1:A:377:ARG:NE	2.39	0.47
1:B:349:GLU:HG3	1:B:350:LEU:N	2.28	0.47
1:A:395:ALA:HB3	1:A:421:THR:HB	1.96	0.47
1:B:32:GLN:HE22	1:B:163:LEU:HB2	1.79	0.47
1:B:315:GLY:HA3	1:B:323:HIS:CD2	2.49	0.47
1:B:500:HIS:HD2	1:B:501:SER:H	1.62	0.47
1:B:415:GLY:C	1:B:416:HIS:CD2	2.87	0.47
1:B:499:ILE:HG23	1:B:533:ILE:HG21	1.96	0.47
1:A:202:ILE:HD13	1:A:321:VAL:CG2	2.39	0.47
1:B:336:SER:O	1:B:340:VAL:HG23	2.14	0.47
1:A:131:ALA:O	1:A:133:ARG:N	2.44	0.46
1:A:170:ARG:HH21	1:A:172:LYS:HZ1	1.64	0.46
1:A:202:ILE:HG22	1:A:202:ILE:O	2.14	0.46
1:B:316:ILE:HD13	1:B:321:VAL:HA	1.96	0.46
1:A:106:LEU:O	1:A:126:ARG:NH2	2.49	0.46
1:B:457:ILE:HD12	1:B:477:VAL:HG22	1.97	0.46
1:B:519:ASP:HB3	1:B:521:THR:HB	1.96	0.46
1:A:48:ASP:HA	1:A:51:LYS:HE2	1.98	0.46
1:B:48:ASP:OD2	1:B:113:PRO:HG3	2.15	0.46
1:B:400:THR:N	1:B:403:ASP:OD2	2.48	0.46
1:B:391:HIS:HB3	1:B:527:LEU:HD23	1.98	0.46
1:B:389:PHE:HE2	1:B:527:LEU:HA	1.81	0.46
1:B:24:LYS:HE2	1:B:69:ASN:HD22	1.81	0.45
1:A:314:MET:HG2	1:A:368:PHE:CE1	2.52	0.45
1:A:353:ILE:H	1:A:353:ILE:HG13	1.57	0.45
1:A:500:HIS:CG	1:A:501:SER:N	2.84	0.45
1:A:32:GLN:HA	1:A:161:VAL:HB	1.97	0.45
1:A:287:GLN:HA	1:A:347:VAL:HG11	1.99	0.45
1:A:73:LEU:HD22	1:A:476:PRO:HD3	1.97	0.45
1:B:99:ALA:O	1:B:103:LYS:HG3	2.16	0.45
1:A:451:SER:O	1:A:485:ALA:HA	2.17	0.45
1:A:432:VAL:HG21	1:A:487:LEU:HD21	1.98	0.44
1:A:284:ASN:O	1:A:288:VAL:HG23	2.16	0.44
1:A:101:THR:O	1:A:105:GLU:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:HIS:CD2	1:A:384:HIS:H	2.34	0.44
1:A:421:THR:HG22	1:A:422:ALA:N	2.32	0.44
1:A:316:ILE:HD13	1:A:321:VAL:HA	2.00	0.44
1:B:86:PRO:O	1:B:133:ARG:NH1	2.40	0.44
1:B:106:LEU:HD22	1:B:127:PHE:CE1	2.53	0.43
1:A:204:LEU:HD11	1:A:329:VAL:HG21	1.99	0.43
1:B:504:ILE:HG23	1:B:506:SER:N	2.32	0.43
1:B:40:SER:HA	1:B:43:GLN:NE2	2.33	0.43
1:A:422:ALA:HB3	1:A:480:GLY:CA	2.48	0.43
1:B:141:ALA:O	1:B:144:SER:HB3	2.18	0.43
1:B:379:ALA:HA	1:B:453:ASN:ND2	2.33	0.43
1:A:172:LYS:NZ	1:A:180:GLU:OE1	2.37	0.43
1:A:2:ALA:HA	1:A:106:LEU:HD23	2.01	0.43
1:A:421:THR:HG23	1:A:484:HIS:ND1	2.33	0.43
1:A:173:ARG:HH21	1:A:207:ASP:CB	2.32	0.43
1:B:290:ASP:OD1	1:B:293:THR:OG1	2.23	0.43
1:A:184:GLN:C	1:A:186:CYS:H	2.23	0.42
1:B:422:ALA:HB3	1:B:480:GLY:HA2	2.01	0.42
1:B:181:TYR:O	1:B:184:GLN:HB2	2.20	0.42
1:B:284:ASN:O	1:B:288:VAL:HG23	2.20	0.42
1:A:76:LYS:HB2	1:A:76:LYS:HE3	1.78	0.42
1:B:457:ILE:HD13	1:B:475:SER:HB2	2.02	0.42
1:A:24:LYS:HA	1:A:66:SER:O	2.19	0.42
1:B:79:GLY:HA2	1:B:167:PHE:CD2	2.55	0.42
1:B:210:ARG:HD3	1:B:362:GLU:OE2	2.19	0.42
1:A:23:THR:HG21	1:A:28:VAL:HA	2.01	0.42
1:B:315:GLY:HA3	1:B:323:HIS:CG	2.55	0.42
1:B:529:LEU:HD12	1:B:529:LEU:HA	1.90	0.42
1:A:350:LEU:O	1:A:350:LEU:HG	2.20	0.42
1:A:352:ASP:O	1:A:355:PHE:HB3	2.20	0.42
1:A:381:SER:HA	1:A:382:PRO:HD3	1.78	0.42
1:A:402:GLU:OE2	1:A:407:ARG:NH1	2.53	0.42
1:B:73:LEU:HD22	1:B:476:PRO:HD3	2.01	0.42
1:A:408:THR:CG2	1:A:409:GLY:H	2.32	0.41
1:A:421:THR:HG22	1:A:422:ALA:H	1.85	0.41
1:B:344:ILE:HB	1:B:345:PRO:HD3	2.01	0.41
1:B:400:THR:O	1:B:403:ASP:N	2.50	0.41
1:A:42:THR:O	1:A:43:GLN:HB2	2.19	0.41
1:A:273:PHE:HE2	1:A:322:HIS:CE1	2.38	0.41
1:B:519:ASP:O	1:B:520:HIS:HB3	2.19	0.41
1:A:384:HIS:ND1	1:A:397:SER:OG	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:GLU:H	1:B:292:GLU:HG2	1.64	0.41
1:B:497:GLU:CG	1:B:498:LYS:H	2.31	0.41
1:A:529:LEU:HD12	1:A:529:LEU:HA	1.98	0.41
1:B:380:TYR:CE2	1:B:419:LYS:HE2	2.55	0.41
1:B:451:SER:HA	1:B:455:ARG:O	2.21	0.41
1:B:333:ILE:HG21	1:B:361:VAL:HG13	2.03	0.41
1:A:273:PHE:CE2	1:A:322:HIS:CE1	3.08	0.40
1:A:386:THR:HB	1:A:398:TRP:O	2.22	0.40
1:A:415:GLY:C	1:A:416:HIS:CD2	2.94	0.40
1:B:186:CYS:O	1:B:201:CYS:SG	2.79	0.40
1:B:407:ARG:HD3	1:B:534:LYS:HB2	2.03	0.40
1:B:445:LYS:HB3	1:B:445:LYS:HE3	1.24	0.40
1:B:166:ASN:HB3	1:B:169:THR:HB	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	471/535 (88%)	425 (90%)	39 (8%)	7 (2%)	10	33
1	B	482/535 (90%)	430 (89%)	47 (10%)	5 (1%)	15	44
All	All	953/1070 (89%)	855 (90%)	86 (9%)	12 (1%)	12	36

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	350	LEU
1	A	501	SER
1	A	519	ASP
1	B	350	LEU
1	A	467	ASP

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Mol	Chain	Res	Type
1	B	78	GLY
1	B	261	SER
1	B	263	LEU
1	B	496	SER
1	A	78	GLY
1	A	132	SER
1	A	382	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	399/452 (88%)	395 (99%)	4 (1%)	76	93
1	B	407/452 (90%)	394 (97%)	13 (3%)	39	73
All	All	806/904 (89%)	789 (98%)	17 (2%)	53	84

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ILE
1	A	51	LYS
1	A	138	LEU
1	A	186	CYS
1	B	56	THR
1	B	69	ASN
1	B	207	ASP
1	B	211	ASP
1	B	272	VAL
1	B	312	SER
1	B	338	LEU
1	B	408	THR
1	B	445	LYS
1	B	495	SER
1	B	500	HIS
1	B	504	ILE

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Mol	Chain	Res	Type
1	B	520	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	502	ASN
1	B	43	GLN
1	B	500	HIS

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	477/535 (89%)	-0.24	4 (0%) 86 81	11, 32, 71, 98	0
1	B	488/535 (91%)	-0.23	5 (1%) 82 77	13, 31, 68, 90	0
All	All	965/1070 (90%)	-0.24	9 (0%) 84 80	11, 32, 70, 98	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	335	LEU	3.4
1	B	218	GLU	2.7
1	B	342	GLN	2.7
1	A	505	SER	2.6
1	A	535	SER	2.6
1	B	275	GLY	2.5
1	B	268	GLY	2.4
1	A	502	ASN	2.4
1	A	39	LYS	2.2

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

There are no such residues in this entry.