



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:32 PM UTC

PDB ID : 3QDE / pdb_00003qde
Title : The structure of Cellobiose phosphorylase from *Clostridium thermocellum* in complex with phosphate
Authors : Bianchetti, C.M.; Elsen, N.L.; Horn, M.K.; Fox, B.G.; Phillips Jr., G.N.
Deposited on : 2011-01-18
Resolution : 2.40 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

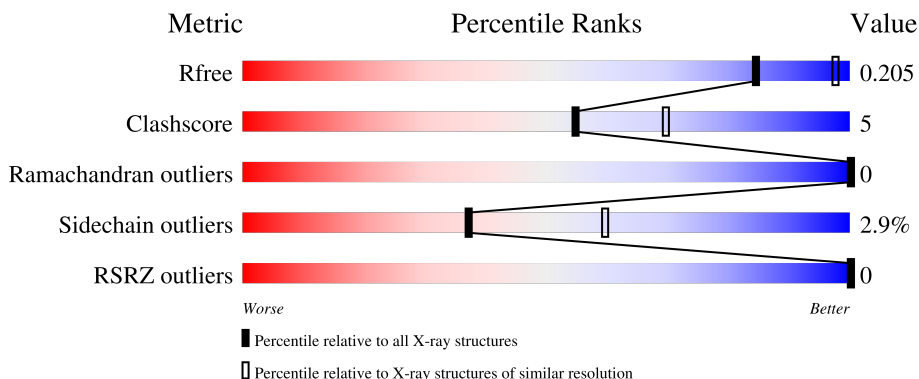
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	811	 84% 16%
1	B	811	 90% 10%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13881 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 Cellobiose phosphorylase.

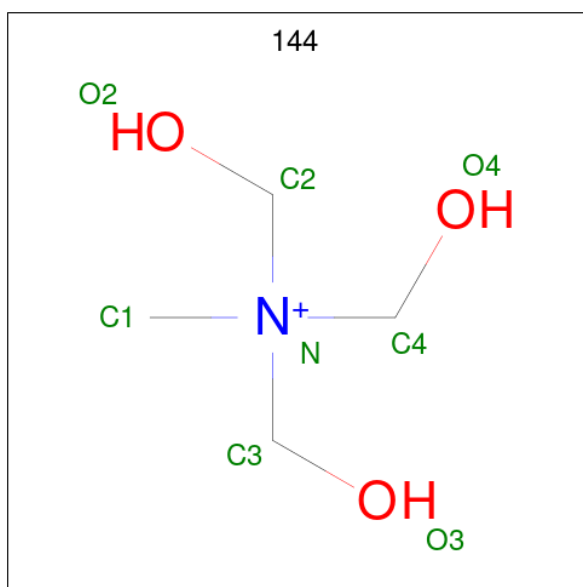
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	811	Total	C	N	O	S	0	1	0
			6571	4209	1098	1236	28			
1	B	811	Total	C	N	O	S	16	4	0
			6592	4222	1104	1237	29			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is TRIS-HYDROXYMETHYL-METHYL-AMMONIUM (CCD ID: 144) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	B	1	Total	C	N	O	0	0
			8	4	1	3		

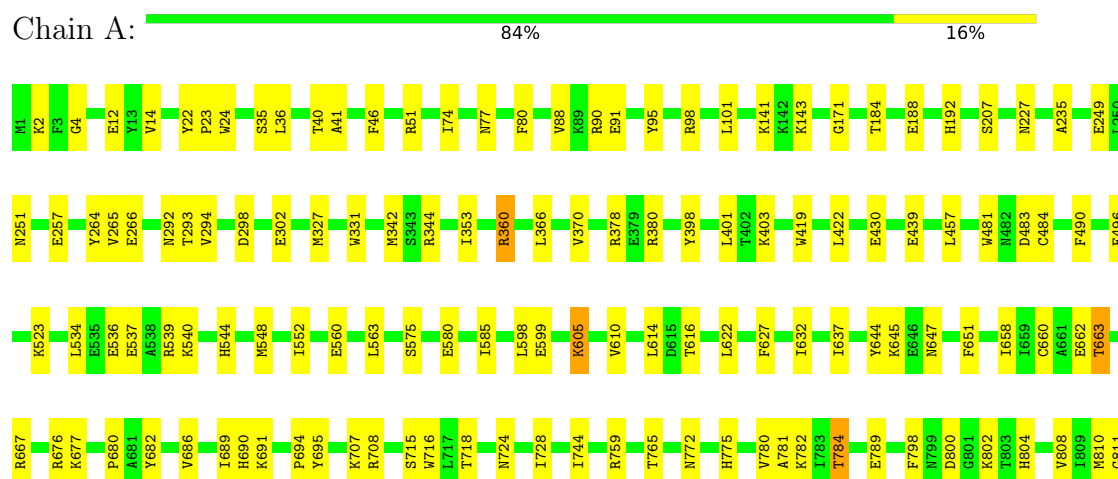
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	320	Total	O	0	0
			320	320		
4	B	372	Total	O	0	0
			372	372		

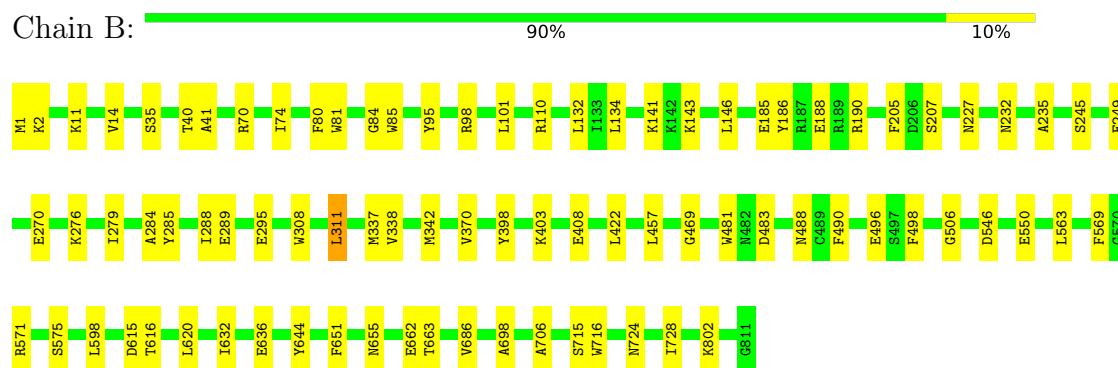
3 Residue-property plots

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: Cellobiose phosphorylase



• Molecule 1: Cellobiose phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.26Å 122.06Å 182.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.33 – 2.40 34.33 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.0 (34.33-2.40) 94.9 (34.33-2.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.94 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.151 , 0.208 0.151 , 0.205	Depositor DCC
R_{free} test set	3696 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13881	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 3.34% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 144, PO4

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	0.39	0/6747	0.72	1/9122 (0.0%)
1	B	0.39	0/6775	0.73	2/9159 (0.0%)
All	All	0.39	0/13522	0.72	3/18281 (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	B	469	GLY	CA-C-N	5.86	126.26	119.47
1	B	469	GLY	C-N-CA	5.86	126.26	119.47
1	A	331	TRP	N-CA-C	5.20	116.75	111.14

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	6571	0	6322	77	0
1	B	6592	0	6346	42	0
2	A	5	0	0	1	0
2	B	5	0	0	1	0
3	A	8	0	12	5	0
3	B	8	0	12	5	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	320	0	0	3	0
4	B	372	0	0	1	0
All	All	13881	0	12692	123	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 5.

All (123) 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:569:PHE:HD2	1:B:571:ARG:HH22	1.30	0.79
1:A:686:VAL:HG12	1:A:689:ILE:HB	1.67	0.76
1:A:481:TRP:O	3:A:813:144:H22	1.85	0.75
1:B:569:PHE:HD2	1:B:571:ARG:NH2	1.88	0.71
1:A:143:LYS:HG2	1:A:249:GLU:HG2	1.73	0.69
1:A:46:PHE:HZ	1:A:342:MET:HE1	1.60	0.66
1:A:483:ASP:H	3:A:813:144:C2	2.12	0.63
1:A:662:GLU:OE1	1:A:667:ARG:HD3	1.99	0.63
1:B:132:LEU:HD23	1:B:132:LEU:C	2.25	0.62
1:A:327:MET:HE2	1:A:680:PRO:HD2	1.84	0.60
1:A:544:HIS:HD2	4:A:1132:HOH:O	1.84	0.59
1:A:360:ARG:NH2	1:A:483:ASP:OD1	2.36	0.59
1:B:546:ASP:O	1:B:550:GLU:HG2	2.04	0.57
1:B:84:GLY:O	1:B:85:TRP:HB3	2.04	0.56
1:B:686:VAL:HG12	1:B:686:VAL:O	2.06	0.56
1:A:264:TYR:CE1	1:A:266:GLU:HG3	2.40	0.55
1:A:780:VAL:HG22	1:A:810:MET:HE1	1.89	0.55
1:B:143:LYS:HG2	1:B:249:GLU:HG2	1.89	0.54
3:A:813:144:H42	4:A:954:HOH:O	2.06	0.54
1:A:585:ILE:HD11	1:A:658:ILE:HG13	1.90	0.54
1:A:694:PRO:HD2	1:A:695:TYR:CD1	2.42	0.54
1:B:651:PHE:CE2	3:B:813:144:H12	2.43	0.54
1:A:360:ARG:HB2	1:A:419:TRP:CD1	2.43	0.53
1:A:12:GLU:OE1	1:A:98:ARG:HD3	2.08	0.52
1:B:132:LEU:HD21	1:B:134:LEU:HG	1.91	0.52
1:B:662:GLU:HA	1:B:662:GLU:OE1	2.09	0.52
1:A:490:PHE:HA	1:A:632:ILE:HG12	1.92	0.51
1:A:534:LEU:HB3	1:A:537:GLU:HG3	1.92	0.51
1:B:185:GLU:HB3	1:B:190:ARG:HG3	1.92	0.51
1:B:481:TRP:O	3:B:813:144:H41	2.11	0.51
1:A:563:LEU:CD2	1:A:575:SER:HB2	2.40	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:488:ASN:OD1	1:B:506:GLY:HA3	2.11	0.51
1:A:644:TYR:OH	1:A:708:ARG:HG3	2.10	0.50
1:A:782:LYS:HG3	1:A:789:GLU:OE2	2.09	0.50
1:A:560:GLU:HG2	1:A:605:LYS:HD2	1.93	0.50
1:A:800:ASP:O	1:A:802:LYS:HD2	2.11	0.49
1:B:615:ASP:HA	1:B:620:LEU:HD23	1.94	0.49
1:A:74:ILE:O	1:A:80:PHE:HA	2.12	0.49
1:A:24:TRP:CH2	1:A:694:PRO:HG3	2.48	0.49
1:A:660:CYS:O	1:A:663:THR:HB	2.13	0.49
1:A:40:THR:O	1:A:41:ALA:HB3	2.12	0.49
1:A:676:ARG:HG2	1:A:682:TYR:CD2	2.49	0.48
1:B:398:TYR:CD1	1:B:403:LYS:HA	2.49	0.48
1:A:51:ARG:HD2	1:A:353:ILE:O	2.13	0.48
1:A:36:LEU:N	1:A:36:LEU:HD12	2.28	0.48
1:A:143:LYS:HE2	1:A:249:GLU:OE1	2.12	0.48
1:B:483:ASP:OD2	3:B:813:144:H22	2.13	0.48
1:B:338:VAL:HG21	1:B:716:TRP:CZ3	2.49	0.48
1:B:563:LEU:HD21	1:B:575:SER:HB2	1.95	0.48
1:A:257:GLU:HB3	1:A:294:VAL:HG21	1.96	0.47
1:A:366:LEU:O	1:A:370:VAL:HG13	2.14	0.47
1:B:342:MET:HB3	1:B:342:MET:HE3	1.66	0.47
1:A:298:ASP:O	1:A:302:GLU:HG2	2.15	0.47
1:A:616:THR:O	1:A:677:LYS:HD3	2.14	0.47
1:A:88:VAL:HG12	1:A:90:ARG:HG3	1.96	0.47
1:A:536:GLU:O	1:A:540:LYS:HD2	2.15	0.46
1:B:207:SER:O	1:B:235:ALA:HB2	2.15	0.46
1:B:186:TYR:CE1	1:B:279:ILE:HG22	2.50	0.46
3:B:813:144:H32	4:B:913:HOH:O	2.14	0.46
1:A:686:VAL:HG13	1:A:689:ILE:HD12	1.97	0.46
1:A:496:GLU:HB3	4:A:1081:HOH:O	2.15	0.46
1:A:690:HIS:O	1:A:691:LYS:HB3	2.16	0.46
1:A:548:MET:O	1:A:552:ILE:HG13	2.16	0.46
1:A:780:VAL:HG22	1:A:810:MET:CE	2.45	0.46
1:A:772:ASN:HA	1:A:810:MET:HB2	1.98	0.46
2:B:812:PO4:O3	3:B:813:144:O3	2.34	0.46
1:B:644:TYR:CZ	1:B:706:ALA:HB1	2.51	0.45
1:A:398:TYR:CD1	1:A:403:LYS:HA	2.51	0.45
1:A:686:VAL:CG1	1:A:689:ILE:HD12	2.47	0.45
1:B:311:LEU:HD12	1:B:311:LEU:HA	1.83	0.45
1:B:146:LEU:O	1:B:245:SER:HA	2.17	0.45
1:A:715:SER:O	1:A:716:TRP:HB2	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:SER:HB3	1:B:70:ARG:NH1	2.32	0.45
1:A:171:GLY:HA2	1:A:184:THR:OG1	2.18	0.44
1:A:207:SER:O	1:A:235:ALA:HB2	2.17	0.44
1:A:327:MET:CE	1:A:680:PRO:HD2	2.47	0.44
1:B:14:VAL:HG22	1:B:98:ARG:HG2	2.00	0.44
1:A:484:CYS:HB3	1:A:637:ILE:N	2.33	0.44
1:A:580:GLU:OE1	1:A:627:PHE:HA	2.18	0.44
1:A:605:LYS:HE3	1:A:605:LYS:HA	2.00	0.44
1:A:366:LEU:HD23	1:A:366:LEU:HA	1.80	0.44
1:B:95:TYR:CD2	1:B:95:TYR:C	2.96	0.43
1:A:292:ASN:OD1	1:A:293:THR:HG23	2.18	0.43
1:A:483:ASP:CG	3:A:813:144:H21	2.42	0.43
1:A:637:ILE:O	1:A:647:ASN:HB2	2.18	0.43
1:A:686:VAL:HG12	1:A:686:VAL:O	2.18	0.43
1:A:724:ASN:O	1:A:728:ILE:HG22	2.19	0.43
1:A:798:PHE:CG	1:A:804:HIS:CE1	3.07	0.43
1:A:810:MET:HA	1:A:810:MET:HE2	1.99	0.43
1:A:676:ARG:HG2	1:A:682:TYR:CG	2.54	0.43
1:A:772:ASN:ND2	1:A:775:HIS:HA	2.33	0.43
2:A:812:PO4:O2	3:A:813:144:O3	2.37	0.43
1:B:284:ALA:O	1:B:288:ILE:HG13	2.19	0.43
1:A:2:LYS:HB3	1:A:2:LYS:HE3	1.79	0.42
1:A:539:ARG:NH2	1:A:540:LYS:HZ2	2.18	0.42
1:A:744:ILE:HG21	1:A:808:VAL:HG21	2.01	0.42
1:A:22:TYR:CG	1:A:23:PRO:HD2	2.55	0.42
1:A:257:GLU:CB	1:A:294:VAL:HG21	2.49	0.42
1:B:308:TRP:CH2	1:B:337:MET:HG2	2.54	0.42
1:B:40:THR:O	1:B:41:ALA:HB3	2.18	0.42
1:B:74:ILE:O	1:B:80:PHE:HA	2.20	0.42
1:B:81:TRP:CG	1:B:110:ARG:HD3	2.55	0.42
1:B:490:PHE:HA	1:B:632:ILE:HG12	2.01	0.42
1:A:781:ALA:HB2	1:A:811:GLY:HA3	2.02	0.41
1:A:378:ARG:NH2	1:A:439:GLU:OE2	2.54	0.41
1:B:186:TYR:CD1	1:B:279:ILE:HG22	2.55	0.41
1:B:285:TYR:O	1:B:289:GLU:HB2	2.20	0.41
1:A:192:HIS:HA	1:A:265:VAL:O	2.21	0.41
1:A:91:GLU:OE1	1:A:91:GLU:HA	2.20	0.41
1:B:205:PHE:O	1:B:232:ASN:HA	2.20	0.41
1:B:141:LYS:HE2	1:B:249:GLU:OE1	2.21	0.41
1:A:645:LYS:C	1:A:651:PHE:HE1	2.29	0.41
1:A:784:THR:HA	1:A:789:GLU:HA	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:498:PHE:CG	1:B:636:GLU:HB2	2.55	0.41
1:A:4:GLY:HA3	1:A:14:VAL:O	2.20	0.41
1:A:430:GLU:CD	1:A:759:ARG:HH22	2.28	0.41
1:A:610:VAL:HG13	1:A:614:LEU:HD12	2.03	0.41
1:B:569:PHE:CD2	1:B:571:ARG:NH2	2.79	0.41
1:B:698:ALA:HA	1:B:715:SER:HA	2.02	0.41
1:B:724:ASN:O	1:B:728:ILE:HG22	2.20	0.41
1:B:655:ASN:N	1:B:655:ASN:HD22	2.19	0.40
1:A:141:LYS:HE2	1:A:251:ASN:OD1	2.21	0.40
1:A:95:TYR:CD2	1:A:95:TYR:C	2.99	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	810/811 (100%)	776 (96%)	34 (4%)	0	100	100
1	B	813/811 (100%)	774 (95%)	39 (5%)	0	100	100
All	All	1623/1622 (100%)	1550 (96%)	73 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	694/693 (100%)	673 (97%)	21 (3%)	36	58
1	B	697/693 (101%)	678 (97%)	19 (3%)	39	62
All	All	1391/1386 (100%)	1351 (97%)	40 (3%)	37	60

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	SER
1	A	77	ASN
1	A	101	LEU
1	A	188	GLU
1	A	227	ASN
1	A	344	ARG
1	A	360	ARG
1	A	380	ARG
1	A	401	LEU
1	A	422	LEU
1	A	457	LEU
1	A	523	LYS
1	A	598	LEU
1	A	599	GLU
1	A	605	LYS
1	A	622	LEU
1	A	663	THR
1	A	707	LYS
1	A	718	THR
1	A	765	THR
1	A	784	THR
1	B	1	MET
1	B	2	LYS
1	B	11	LYS
1	B	101	LEU
1	B	188	GLU
1	B	227	ASN
1	B	270	GLU
1	B	276	LYS
1	B	295	GLU
1	B	311	LEU
1	B	370	VAL
1	B	408	GLU
1	B	422	LEU
1	B	457	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	496	GLU
1	B	598	LEU
1	B	616	THR
1	B	663	THR
1	B	802	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	332	ASN
1	A	407	ASN
1	A	542	GLN
1	A	544	HIS
1	B	267	ASN
1	B	332	ASN
1	B	542	GLN
1	B	624	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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	$\# Z > 2$
2	PO4	A	812	-	4,4,4	1.06	0	6,6,6	0.64	0
2	PO4	B	812	-	4,4,4	1.05	0	6,6,6	0.75	0
3	144	A	813	-	1,7,7	0.78	0	3,9,9	0.20	0
3	144	B	813	-	1,7,7	0.87	0	3,9,9	0.58	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	144	A	813	-	-	0/0/9/9	-
3	144	B	813	-	-	0/0/9/9	-

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.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	812	PO4	1	0
2	B	812	PO4	1	0
3	A	813	144	5	0
3	B	813	144	5	0

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	811/811 (100%)	-0.76	0 100 100	18, 33, 54, 73	1 (0%)
1	B	811/811 (100%)	-0.85	0 100 100	12, 31, 47, 64	4 (0%)
All	All	1622/1622 (100%)	-0.81	0 100 100	12, 32, 51, 73	5 (0%)

There are no RSRZ outliers to report.

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	144	B	813	8/8	0.86	0.15	38,42,49,57	0
3	144	A	813	8/8	0.93	0.11	36,40,43,44	0
2	PO4	A	812	5/5	0.98	0.05	32,33,34,36	0
2	PO4	B	812	5/5	0.99	0.04	25,33,34,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.