Full wwPDB X-ray Structure Validation Report (i)

Jan 30, 2021 – 06:25 PM EST

PDB ID	:	3K85
Title	:	Crystal structure of a D-glycero-D-manno-heptose 1-phosphate kinase from Bacteriodes thetaiotaomicron
Authors	:	Palani, K.; Kumaran, D.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2009-10-13
Resolution	:	2.28 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 (i) symbol.

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

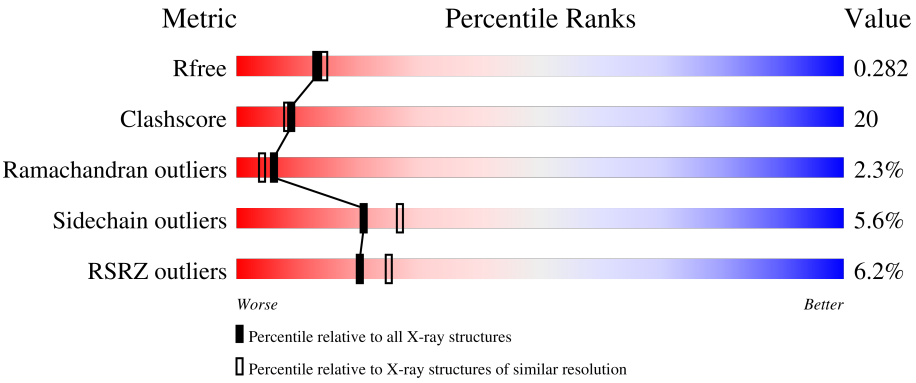
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

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 2.28 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	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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	357	7% 57% 25% . 14%
1	B	357	3% 55% 28% . 14%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4879 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 D-glycero-D-manno-heptose 1-phosphate kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	Se	0	0	0
			2371	1517	391	450	4	9			
1	B	306	Total	C	N	O	S	Se	0	0	0
			2371	1517	391	450	4	9			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MSE	-	expression tag	UNP Q8AAJ0
A	1	SER	-	expression tag	UNP Q8AAJ0
A	2	LEU	-	expression tag	UNP Q8AAJ0
A	349	GLU	-	expression tag	UNP Q8AAJ0
A	350	GLY	-	expression tag	UNP Q8AAJ0
A	351	HIS	-	expression tag	UNP Q8AAJ0
A	352	HIS	-	expression tag	UNP Q8AAJ0
A	353	HIS	-	expression tag	UNP Q8AAJ0
A	354	HIS	-	expression tag	UNP Q8AAJ0
A	355	HIS	-	expression tag	UNP Q8AAJ0
A	356	HIS	-	expression tag	UNP Q8AAJ0
B	0	MSE	-	expression tag	UNP Q8AAJ0
B	1	SER	-	expression tag	UNP Q8AAJ0
B	2	LEU	-	expression tag	UNP Q8AAJ0
B	349	GLU	-	expression tag	UNP Q8AAJ0
B	350	GLY	-	expression tag	UNP Q8AAJ0
B	351	HIS	-	expression tag	UNP Q8AAJ0
B	352	HIS	-	expression tag	UNP Q8AAJ0
B	353	HIS	-	expression tag	UNP Q8AAJ0
B	354	HIS	-	expression tag	UNP Q8AAJ0
B	355	HIS	-	expression tag	UNP Q8AAJ0
B	356	HIS	-	expression tag	UNP Q8AAJ0

- Molecule 2 is water.

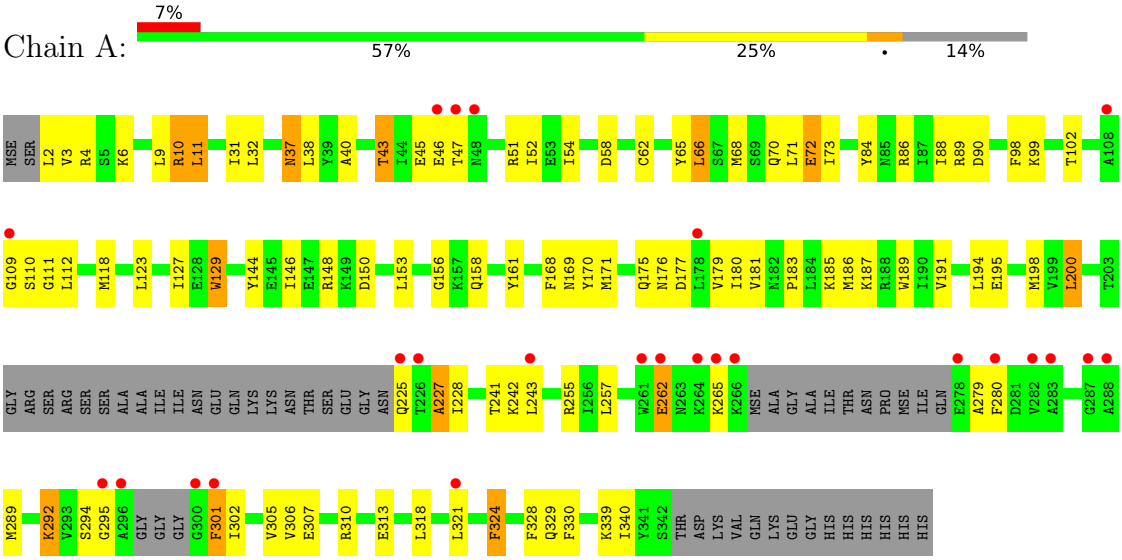
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	75	Total O 75 75	0	0
2	B	62	Total O 62 62	0	0

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.

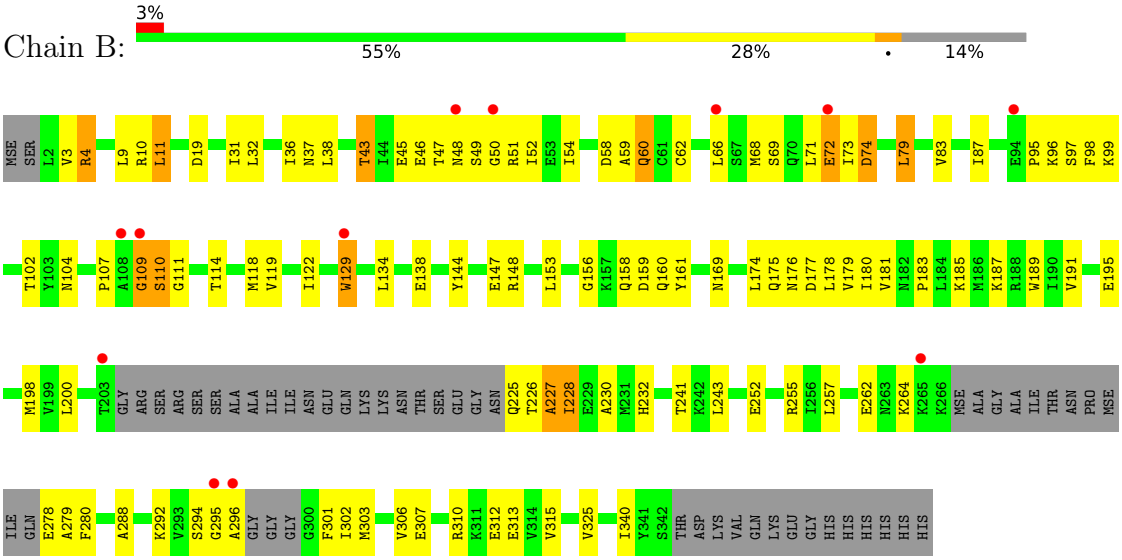
• Molecule 1: D-glycero-D-manno-heptose 1-phosphate kinase

Chain A:



• Molecule 1: D-glycero-D-manno-heptose 1-phosphate kinase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	71.89Å 149.47Å 141.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.95 – 2.28 40.95 – 2.04	Depositor EDS
% Data completeness (in resolution range)	94.1 (40.95-2.28) 88.4 (40.95-2.04)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.17 (at 2.05Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.254 , 0.282 0.254 , 0.282	Depositor DCC
R_{free} test set	1677 reflections (3.61%)	wwPDB-VP
Wilson B-factor (Å^2)	30.2	Xtriage
Anisotropy	0.682	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 45.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	4879	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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

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	0/2404	0.62	0/3224
1	B	0.39	0/2404	0.63	0/3224
All	All	0.39	0/4808	0.62	0/6448

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

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	2371	0	2364	90	0
1	B	2371	0	2364	107	0
2	A	75	0	0	1	0
2	B	62	0	0	4	0
All	All	4879	0	4728	191	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 (191) 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:257:LEU:HD22	1:B:292:LYS:HB2	1.51	0.91

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ILE:HD12	1:B:243:LEU:HD13	1.53	0.90
1:A:58:ASP:OD1	1:A:102:THR:HG22	1.71	0.89
1:A:185:LYS:HG2	1:B:185:LYS:HG2	1.60	0.83
1:A:2:LEU:HD11	1:A:4:ARG:NH1	1.93	0.83
1:A:3:VAL:HG12	1:A:339:LYS:HG2	1.58	0.82
1:B:174:LEU:HD12	1:B:178:LEU:HD23	1.61	0.82
1:A:70:GLN:HE22	1:A:89:ARG:HH11	1.27	0.82
1:B:225:GLN:HG2	1:B:226:THR:H	1.42	0.82
1:B:58:ASP:OD1	1:B:102:THR:HG22	1.79	0.82
1:B:187:LYS:HE2	1:B:189:TRP:HE1	1.44	0.81
1:A:4:ARG:NH1	1:A:340:ILE:HG13	1.97	0.80
1:B:11:LEU:HD13	1:B:36:ILE:HD13	1.62	0.78
1:B:52:ILE:HD13	1:B:71:LEU:HD11	1.66	0.77
1:A:52:ILE:HD13	1:A:71:LEU:HD11	1.71	0.73
1:B:181:VAL:O	1:B:183:PRO:HD3	1.90	0.72
1:A:294:SER:HB2	1:A:301:PHE:CD2	2.23	0.72
1:A:243:LEU:CD1	1:B:180:ILE:HD12	2.20	0.71
1:B:310:ARG:HG2	1:B:310:ARG:HH11	1.56	0.70
1:B:187:LYS:HE2	1:B:189:TRP:NE1	2.06	0.70
1:A:181:VAL:O	1:A:183:PRO:HD3	1.92	0.69
1:A:4:ARG:HH11	1:A:340:ILE:HG13	1.56	0.69
1:B:306:VAL:HG22	1:B:307:GLU:N	2.07	0.69
1:B:144:TYR:CE1	1:B:148:ARG:HG3	2.28	0.68
1:A:46:GLU:OE1	1:A:129:TRP:HH2	1.75	0.68
1:A:47:THR:HG22	1:A:98:PHE:HA	1.76	0.68
1:B:102:THR:HG21	2:B:380:HOH:O	1.94	0.68
1:A:243:LEU:HD13	1:B:180:ILE:HD12	1.76	0.68
1:B:228:ILE:HG22	1:B:232:HIS:NE2	2.10	0.67
1:B:148:ARG:NH1	1:B:156:GLY:O	2.28	0.66
1:B:294:SER:HB3	1:B:301:PHE:HD2	1.60	0.66
1:B:227:ALA:O	1:B:230:ALA:N	2.25	0.66
1:A:11:LEU:HD23	1:A:198:MSE:HE1	1.78	0.66
1:A:148:ARG:NH1	1:A:156:GLY:O	2.30	0.65
1:A:10:ARG:HD2	1:A:10:ARG:C	2.18	0.63
1:A:324:PHE:H	1:A:324:PHE:HD2	1.47	0.62
1:A:66:LEU:O	1:A:68:MSE:HG3	2.00	0.62
1:A:255:ARG:HH11	1:A:255:ARG:HG2	1.64	0.62
1:B:104:ASN:ND2	1:B:114:THR:HG23	2.15	0.61
1:B:54:ILE:O	1:B:62:CYS:HA	2.01	0.61
1:A:318:LEU:O	1:A:321:LEU:HB2	2.02	0.60
1:A:169:ASN:HD22	1:A:183:PRO:HA	1.68	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:TYR:CE1	1:A:148:ARG:HG3	2.37	0.59
1:B:107:PRO:HG2	1:B:110:SER:OG	2.03	0.59
1:A:310:ARG:HH11	1:A:310:ARG:HG2	1.68	0.58
1:A:31:ILE:CD1	1:A:158:GLN:HB3	2.32	0.58
1:A:4:ARG:HG2	1:A:43:THR:HB	1.86	0.58
1:B:161:TYR:CZ	1:B:179:VAL:HG11	2.38	0.58
1:B:294:SER:HB3	1:B:301:PHE:CD2	2.38	0.58
1:A:227:ALA:O	1:A:228:ILE:HB	2.04	0.57
1:B:10:ARG:C	1:B:10:ARG:HD2	2.25	0.57
1:A:169:ASN:ND2	1:A:183:PRO:HA	2.18	0.57
1:A:289:MSE:HB2	1:A:305:VAL:O	2.05	0.56
1:A:72:GLU:H	1:A:72:GLU:CD	2.09	0.56
1:B:47:THR:O	1:B:97:SER:HB3	2.06	0.56
1:B:148:ARG:HD3	1:B:153:LEU:O	2.05	0.56
1:A:243:LEU:HD11	1:B:180:ILE:HD12	1.89	0.55
1:A:68:MSE:HE3	1:A:71:LEU:CD2	2.35	0.55
1:B:144:TYR:CD1	1:B:148:ARG:HG3	2.42	0.55
1:B:47:THR:HG22	1:B:98:PHE:CA	2.37	0.55
1:B:9:LEU:HD12	1:B:38:LEU:HB3	1.88	0.55
1:A:180:ILE:CD1	1:B:243:LEU:HD13	2.33	0.54
1:B:79:LEU:O	1:B:83:VAL:HG23	2.07	0.54
1:B:46:GLU:HB2	1:B:129:TRP:CH2	2.43	0.54
1:A:9:LEU:HD12	1:A:38:LEU:HB3	1.89	0.54
1:A:179:VAL:O	1:A:180:ILE:HD13	2.08	0.54
1:A:86:ARG:HD3	1:A:150:ASP:OD2	2.07	0.54
1:A:191:VAL:O	1:A:195:GLU:HG3	2.08	0.53
1:B:43:THR:HG22	2:B:415:HOH:O	2.08	0.53
1:B:19:ASP:HB3	1:B:31:ILE:HD12	1.91	0.53
1:B:72:GLU:CD	1:B:72:GLU:N	2.63	0.53
1:A:58:ASP:OD1	1:A:102:THR:CG2	2.49	0.52
1:A:302:ILE:N	1:A:302:ILE:HD12	2.24	0.52
1:B:228:ILE:HG22	1:B:232:HIS:CE1	2.45	0.52
1:B:306:VAL:CG2	1:B:307:GLU:N	2.72	0.52
1:A:255:ARG:NH1	1:A:255:ARG:HG2	2.23	0.52
1:B:52:ILE:HG23	1:B:52:ILE:O	2.10	0.52
1:A:84:TYR:O	1:A:88:ILE:HG12	2.09	0.52
1:B:111:GLY:N	1:B:296:ALA:HB3	2.24	0.52
1:A:70:GLN:NE2	1:A:89:ARG:HH11	2.04	0.52
1:B:47:THR:HG22	1:B:98:PHE:HA	1.91	0.51
1:A:257:LEU:HD21	1:A:292:LYS:HE2	1.92	0.51
1:A:112:LEU:HD21	1:A:301:PHE:CE1	2.46	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:GLN:HE22	1:A:89:ARG:NH1	2.04	0.50
1:B:102:THR:HG23	1:B:118:MSE:CE	2.41	0.50
1:B:111:GLY:HA3	1:B:295:GLY:O	2.11	0.50
1:A:257:LEU:CD2	1:A:292:LYS:HE2	2.42	0.50
1:B:148:ARG:HA	1:B:153:LEU:O	2.10	0.50
1:B:68:MSE:O	1:B:95:PRO:HB3	2.12	0.50
1:B:138:GLU:H	1:B:138:GLU:CD	2.14	0.49
1:B:74:ASP:N	1:B:74:ASP:OD2	2.43	0.49
1:B:9:LEU:HD21	1:B:114:THR:HA	1.94	0.49
1:A:73:ILE:N	1:A:73:ILE:HD12	2.27	0.49
1:B:310:ARG:NH1	1:B:310:ARG:HG2	2.25	0.49
1:A:32:LEU:HD23	1:A:241:THR:HG22	1.95	0.49
1:B:47:THR:HG22	1:B:98:PHE:N	2.26	0.49
1:A:46:GLU:HG2	1:A:47:THR:N	2.26	0.49
1:B:306:VAL:HG22	1:B:307:GLU:H	1.77	0.49
1:B:73:ILE:N	1:B:73:ILE:HD12	2.29	0.48
1:A:6:LYS:HA	1:A:40:ALA:O	2.13	0.48
1:A:306:VAL:HG22	1:A:307:GLU:N	2.28	0.48
1:A:52:ILE:HG23	1:A:65:TYR:HB2	1.94	0.48
1:A:10:ARG:HD2	1:A:10:ARG:O	2.14	0.48
1:A:262:GLU:HA	1:A:265:LYS:HE2	1.95	0.48
1:A:328:PHE:O	1:A:329:GLN:NE2	2.47	0.48
2:A:391:HOH:O	1:B:185:LYS:HE2	2.13	0.48
1:B:315:VAL:HG13	1:B:325:VAL:HG21	1.96	0.48
1:A:2:LEU:C	1:A:2:LEU:HD12	2.34	0.48
1:A:31:ILE:HD13	1:A:158:GLN:HB3	1.95	0.48
1:B:175:GLN:O	1:B:176:ASN:HB2	2.14	0.47
1:B:161:TYR:OH	1:B:179:VAL:HG11	2.13	0.47
1:B:73:ILE:HG22	1:B:73:ILE:O	2.15	0.47
1:A:289:MSE:HE3	1:A:305:VAL:HG12	1.97	0.47
1:B:187:LYS:CE	1:B:189:TRP:HE1	2.21	0.47
1:B:302:ILE:N	1:B:302:ILE:HD12	2.29	0.47
1:B:3:VAL:HG21	1:B:129:TRP:HB2	1.97	0.47
1:B:83:VAL:O	1:B:87:ILE:HG12	2.15	0.47
1:B:47:THR:H	1:B:98:PHE:HA	1.80	0.47
1:B:47:THR:HG23	1:B:97:SER:OG	2.15	0.47
1:A:324:PHE:CD2	1:A:324:PHE:N	2.83	0.47
1:A:168:PHE:CG	1:A:186:MSE:HG3	2.50	0.46
1:B:148:ARG:NH1	1:B:160:GLN:HE22	2.14	0.46
1:A:294:SER:HB2	1:A:301:PHE:HD2	1.76	0.46
1:B:227:ALA:O	1:B:228:ILE:C	2.52	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:LYS:O	1:B:264:LYS:HG2	2.15	0.46
1:B:147:GLU:O	1:B:153:LEU:HB2	2.14	0.46
1:B:312:GLU:OE1	1:B:312:GLU:HA	2.15	0.46
1:A:257:LEU:HD22	1:A:292:LYS:HB2	1.98	0.46
1:A:187:LYS:HE2	1:A:189:TRP:HE1	1.81	0.46
1:B:46:GLU:OE1	1:B:129:TRP:CH2	2.68	0.46
1:A:102:THR:HG23	1:A:118:MSE:SE	2.66	0.46
1:A:112:LEU:HD21	1:A:301:PHE:CZ	2.51	0.46
1:B:109:GLY:O	1:B:111:GLY:N	2.45	0.46
1:A:31:ILE:HD12	1:A:158:GLN:HB3	1.98	0.45
1:A:146:ILE:O	1:A:150:ASP:HB2	2.16	0.45
1:B:95:PRO:O	1:B:96:LYS:HD2	2.17	0.45
1:A:11:LEU:HD11	1:A:200:LEU:HD11	1.97	0.45
1:A:200:LEU:HD22	1:A:200:LEU:HA	1.88	0.45
1:B:46:GLU:OE1	1:B:129:TRP:HH2	2.00	0.45
1:B:72:GLU:CD	1:B:72:GLU:H	2.20	0.45
1:A:170:TYR:CD1	1:A:242:LYS:HD3	2.52	0.45
1:A:225:GLN:C	1:A:227:ALA:N	2.71	0.44
1:B:129:TRP:O	1:B:129:TRP:CE3	2.70	0.44
1:A:68:MSE:HE3	1:A:71:LEU:HD23	1.97	0.44
1:B:200:LEU:HD22	1:B:301:PHE:CZ	2.52	0.44
1:B:110:SER:N	1:B:296:ALA:HB1	2.33	0.44
1:B:158:GLN:HG3	1:B:159:ASP:OD1	2.18	0.44
1:B:134:LEU:HD22	1:B:138:GLU:HB3	2.00	0.44
1:B:187:LYS:HE3	2:B:387:HOH:O	2.18	0.44
1:B:187:LYS:HE2	1:B:189:TRP:CD1	2.53	0.44
1:A:175:GLN:O	1:A:176:ASN:HB2	2.17	0.44
1:B:10:ARG:O	1:B:10:ARG:HD2	2.17	0.44
1:B:102:THR:HG23	1:B:118:MSE:SE	2.67	0.43
1:B:225:GLN:HG2	1:B:226:THR:N	2.21	0.43
1:A:45:GLU:HG3	1:A:99:LYS:HB3	2.01	0.43
1:B:32:LEU:HD23	1:B:241:THR:HG22	2.00	0.43
1:A:310:ARG:HG2	1:A:310:ARG:NH1	2.31	0.43
1:B:32:LEU:HD12	1:B:169:ASN:O	2.17	0.43
1:B:66:LEU:O	1:B:68:MSE:HG3	2.19	0.43
1:B:169:ASN:HD22	1:B:183:PRO:HA	1.84	0.43
1:A:109:GLY:O	1:A:111:GLY:N	2.51	0.43
1:A:37:ASN:ND2	1:A:329:GLN:HB2	2.33	0.43
1:B:110:SER:C	1:B:296:ALA:HB3	2.39	0.43
1:B:198:MSE:HE3	1:B:303:MSE:HG3	2.00	0.43
1:A:123:LEU:O	1:A:127:ILE:HG13	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ILE:HG22	1:A:73:ILE:O	2.18	0.43
1:A:148:ARG:HA	1:A:153:LEU:O	2.18	0.42
1:B:252:GLU:CD	1:B:255:ARG:HE	2.23	0.42
1:B:99:LYS:NZ	2:B:388:HOH:O	2.52	0.42
1:A:161:TYR:HB3	1:A:171:MSE:CE	2.49	0.42
1:A:46:GLU:HB2	1:A:129:TRP:CH2	2.53	0.42
1:B:306:VAL:CG2	1:B:307:GLU:H	2.32	0.42
1:A:52:ILE:CD1	1:A:71:LEU:HD11	2.46	0.42
1:B:288:ALA:CB	1:B:306:VAL:HB	2.50	0.42
1:A:86:ARG:NH1	1:A:90:ASP:OD2	2.52	0.41
1:B:49:SER:O	1:B:50:GLY:C	2.57	0.41
1:A:194:LEU:HD23	1:A:330:PHE:CE2	2.55	0.41
1:B:292:LYS:HD3	1:B:303:MSE:HE2	2.03	0.41
1:B:59:ALA:O	1:B:60:GLN:C	2.58	0.41
1:A:73:ILE:N	1:A:73:ILE:CD1	2.84	0.41
1:A:144:TYR:CD1	1:A:148:ARG:HG3	2.56	0.41
1:B:79:LEU:HG	1:B:119:VAL:CG2	2.51	0.41
1:A:47:THR:HG22	1:A:98:PHE:CA	2.50	0.40
1:B:19:ASP:HB3	1:B:31:ILE:CD1	2.50	0.40
1:B:252:GLU:HA	1:B:252:GLU:OE2	2.20	0.40
1:A:54:ILE:O	1:A:62:CYS:HA	2.21	0.40
1:B:191:VAL:O	1:B:195:GLU:HG3	2.21	0.40
1:B:118:MSE:O	1:B:122:ILE:HG13	2.21	0.40
1:B:4:ARG:HH11	1:B:340:ILE:HG13	1.84	0.40
1:B:228:ILE:HG13	1:B:228:ILE:H	1.65	0.40
1:B:45:GLU:HG3	1:B:99:LYS:HB3	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	298/357 (84%)	277 (93%)	14 (5%)	7 (2%)	6 4
1	B	298/357 (84%)	270 (91%)	21 (7%)	7 (2%)	6 4
All	All	596/714 (84%)	547 (92%)	35 (6%)	14 (2%)	6 4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	110	SER
1	A	279	ALA
1	A	295	GLY
1	B	177	ASP
1	B	228	ILE
1	A	177	ASP
1	B	279	ALA
1	A	37	ASN
1	A	227	ALA
1	A	301	PHE
1	B	37	ASN
1	B	109	GLY
1	B	110	SER
1	B	227	ALA

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	249/277 (90%)	236 (95%)	13 (5%)	23 30
1	B	249/277 (90%)	234 (94%)	15 (6%)	19 24
All	All	498/554 (90%)	470 (94%)	28 (6%)	21 27

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ARG
1	A	11	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	43	THR
1	A	51	ARG
1	A	66	LEU
1	A	72	GLU
1	A	129	TRP
1	A	200	LEU
1	A	262	GLU
1	A	280	PHE
1	A	292	LYS
1	A	313	GLU
1	A	324	PHE
1	B	4	ARG
1	B	11	LEU
1	B	43	THR
1	B	48	ASN
1	B	51	ARG
1	B	60	GLN
1	B	69	SER
1	B	72	GLU
1	B	74	ASP
1	B	79	LEU
1	B	129	TRP
1	B	262	GLU
1	B	278	GLU
1	B	280	PHE
1	B	313	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	70	GLN
1	A	169	ASN
1	B	60	GLN
1	B	70	GLN
1	B	169	ASN

5.3.3 RNA

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

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	297/357 (83%)	0.52	25 (8%)	11	14	18, 38, 77, 91	0
1	B	297/357 (83%)	0.44	12 (4%)	38	43	17, 40, 67, 82	0
All	All	594/714 (83%)	0.48	37 (6%)	20	25	17, 39, 73, 91	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	108	ALA	8.3
1	A	108	ALA	6.2
1	B	295	GLY	5.6
1	A	300	GLY	4.8
1	A	282	VAL	4.4
1	A	295	GLY	4.0
1	A	288	ALA	3.7
1	A	265	LYS	3.6
1	B	50	GLY	3.6
1	A	109	GLY	3.6
1	A	280	PHE	3.6
1	A	266	LYS	3.6
1	B	109	GLY	3.4
1	B	296	ALA	3.3
1	A	287	GLY	3.3
1	A	283	ALA	3.1
1	A	278	GLU	3.0
1	A	296	ALA	2.9
1	A	261	TRP	2.9
1	B	94	GLU	2.9
1	B	48	ASN	2.9
1	A	226	THR	2.8
1	A	301	PHE	2.7
1	A	178	LEU	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	321	LEU	2.5
1	B	72	GLU	2.5
1	A	225	GLN	2.4
1	B	203	THR	2.4
1	A	264	LYS	2.3
1	A	262	GLU	2.3
1	B	265	LYS	2.3
1	A	46	GLU	2.2
1	A	48	ASN	2.2
1	A	243	LEU	2.2
1	A	47	THR	2.1
1	B	129	TRP	2.0
1	B	66	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.