Full wwPDB X-ray Structure Validation Report (i)

May 14, 2020 – 09:45 pm BST

PDB ID	:	3CZQ
Title	:	Crystal structure of putative polyphosphate kinase 2 from Sinorhizobium meliloti
Authors	:	Osipiuk, J.; Evdokimova, E.; Nocek, B.; Kudritska, M.; Savchenko, A.; Edwards, A.M.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2008-04-29
Resolution	:	2.23 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 (1)) were used in the production of this report:

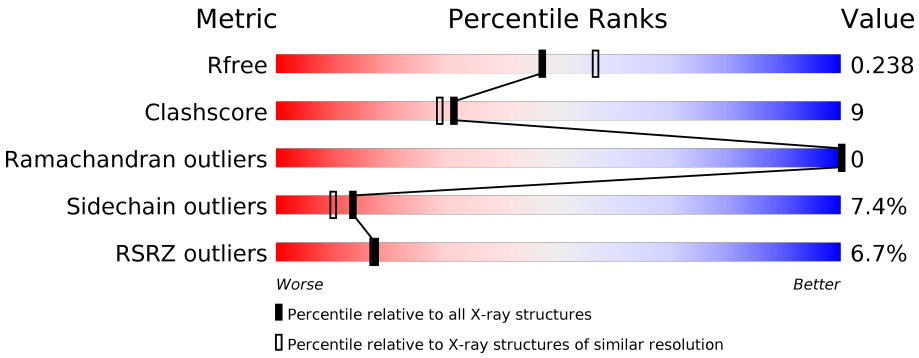
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.23 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	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (2.26-2.22)

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 <=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	304	4% 75% 17% • 5%
1	B	304	6% 78% 13% •• 6%
1	C	304	7% 68% 21% • 7%
1	D	304	8% 74% 18% • 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9706 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.

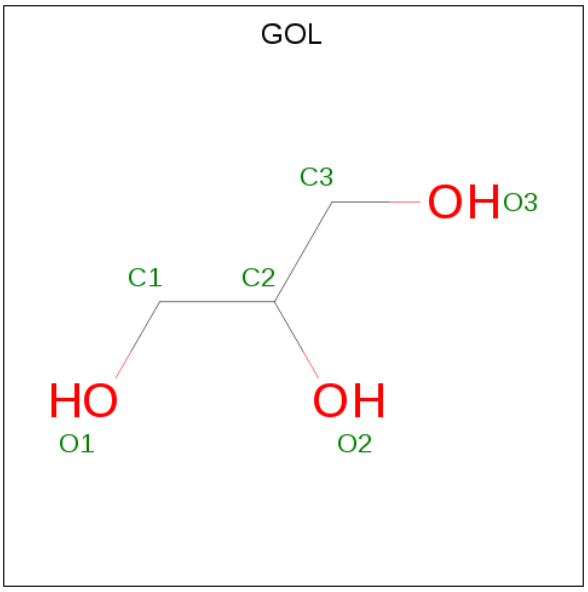
- Molecule 1 is a protein called Putative polyphosphate kinase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	Se	0	1	0
			2371	1507	421	432	1	10			
1	B	287	Total	C	N	O	S	Se	0	0	0
			2359	1498	420	431	1	9			
1	C	284	Total	C	N	O	S	Se	0	1	0
			2348	1493	417	428	1	9			
1	D	286	Total	C	N	O	S	Se	0	0	0
			2355	1496	419	430	1	9			

There are 16 discrepancies between the modelled and reference sequences:

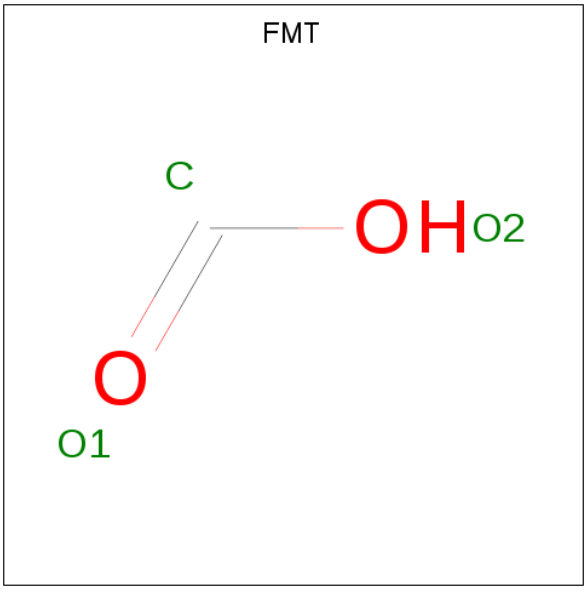
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q92SA6
A	0	HIS	-	EXPRESSION TAG	UNP Q92SA6
A	301	GLY	-	EXPRESSION TAG	UNP Q92SA6
A	302	SER	-	EXPRESSION TAG	UNP Q92SA6
B	-1	GLY	-	EXPRESSION TAG	UNP Q92SA6
B	0	HIS	-	EXPRESSION TAG	UNP Q92SA6
B	301	GLY	-	EXPRESSION TAG	UNP Q92SA6
B	302	SER	-	EXPRESSION TAG	UNP Q92SA6
C	-1	GLY	-	EXPRESSION TAG	UNP Q92SA6
C	0	HIS	-	EXPRESSION TAG	UNP Q92SA6
C	301	GLY	-	EXPRESSION TAG	UNP Q92SA6
C	302	SER	-	EXPRESSION TAG	UNP Q92SA6
D	-1	GLY	-	EXPRESSION TAG	UNP Q92SA6
D	0	HIS	-	EXPRESSION TAG	UNP Q92SA6
D	301	GLY	-	EXPRESSION TAG	UNP Q92SA6
D	302	SER	-	EXPRESSION TAG	UNP Q92SA6

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 12 6 6	0	1

- Molecule 3 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	1	Total C O 3 1 2	0	0
3	D	1	Total C O 3 1 2	0	0

- Molecule 4 is water.

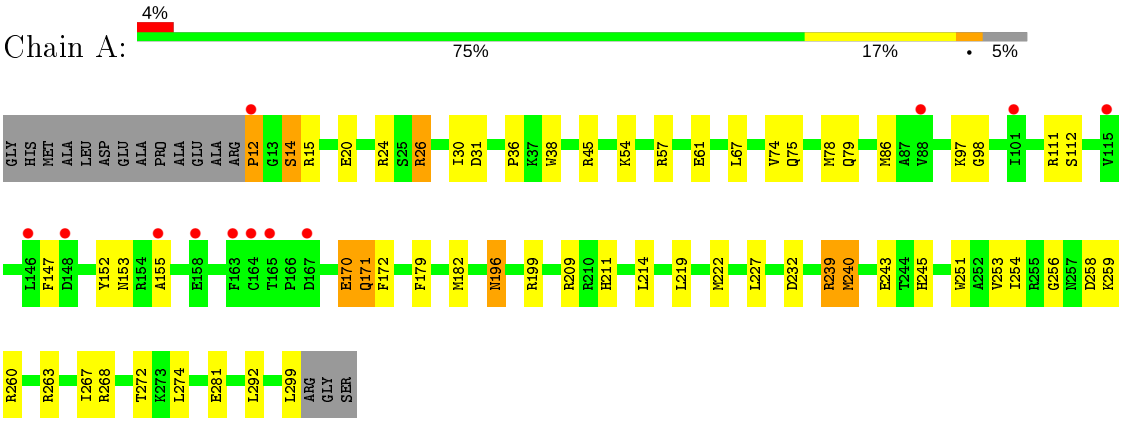
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	104	Total O 104 104	0	0
4	B	84	Total O 84 84	0	0
4	C	31	Total O 31 31	0	0
4	D	36	Total O 36 36	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

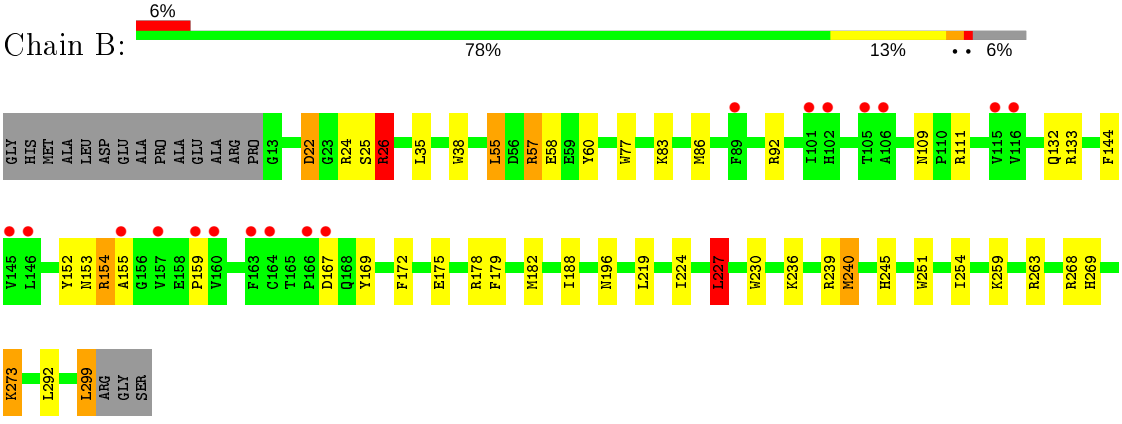
- Molecule 1: Putative polyphosphate kinase 2

Chain A:



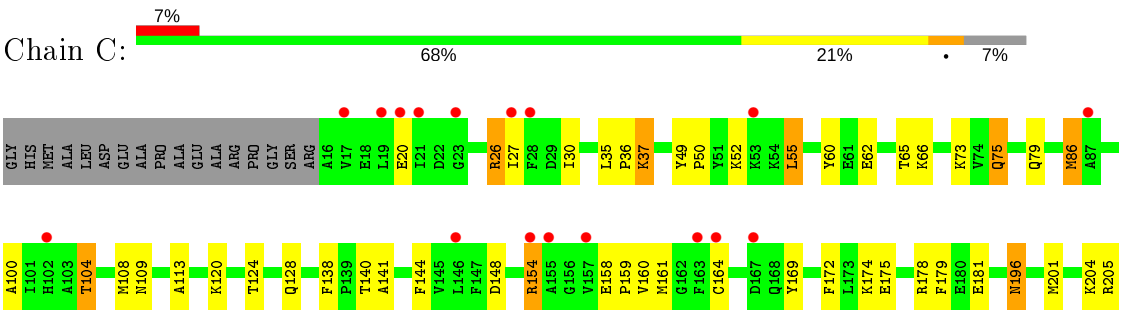
- Molecule 1: Putative polyphosphate kinase 2

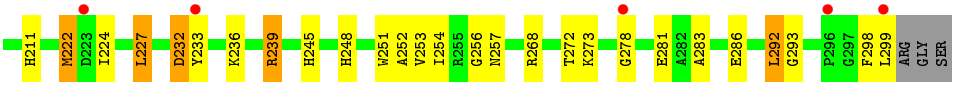
Chain B:



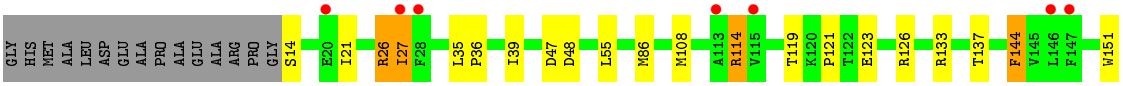
- Molecule 1: Putative polyphosphate kinase 2

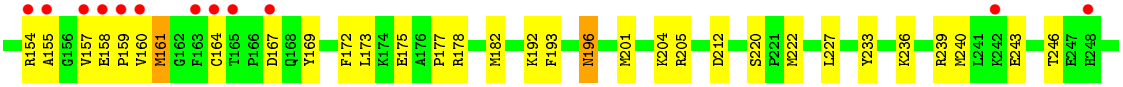
Chain C:

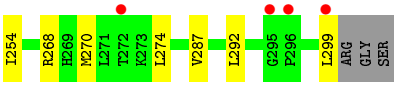




- Molecule 1: Putative polyphosphate kinase 2







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	59.61Å 71.67Å 89.47Å 75.86° 85.97° 65.39°	Depositor
Resolution (Å)	34.70 – 2.23 34.68 – 2.23	Depositor EDS
% Data completeness (in resolution range)	89.3 (34.70-2.23) 89.2 (34.68-2.23)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.07 (at 2.24Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.187 , 0.240 0.186 , 0.238	Depositor DCC
R_{free} test set	2840 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	47.8	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 52.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9706	wwPDB-VP
Average B, all atoms (Å^2)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% 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: GOL, FMT

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	0.76	0/2425	0.79	2/3253 (0.1%)
1	B	0.74	0/2409	0.81	7/3232 (0.2%)
1	C	0.80	3/2399 (0.1%)	0.71	2/3220 (0.1%)
1	D	0.64	0/2405	0.68	1/3227 (0.0%)
All	All	0.74	3/9638 (0.0%)	0.75	12/12932 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	26	ARG	NE-CZ	23.59	1.63	1.33
1	C	26	ARG	C-O	6.08	1.34	1.23
1	C	26	ARG	C-N	5.33	1.46	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	26	ARG	NE-CZ-NH2	-12.24	114.18	120.30
1	C	26	ARG	NE-CZ-NH1	7.83	124.21	120.30
1	A	260	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	B	133	ARG	NE-CZ-NH2	-6.50	117.05	120.30
1	B	133	ARG	NE-CZ-NH1	6.48	123.54	120.30
1	B	26	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	A	258	ASP	CB-CG-OD1	5.54	123.29	118.30
1	D	212	ASP	CB-CG-OD1	5.33	123.10	118.30
1	B	227	LEU	CA-CB-CG	5.32	127.53	115.30
1	B	299	LEU	CA-CB-CG	5.16	127.17	115.30
1	B	268	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	B	268	ARG	NE-CZ-NH2	-5.08	117.76	120.30

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	2371	0	2324	48	0
1	B	2359	0	2307	31	0
1	C	2348	0	2292	54	0
1	D	2355	0	2304	40	0
2	A	12	0	16	6	0
3	C	3	0	1	0	0
3	D	3	0	1	0	0
4	A	104	0	0	5	0
4	B	84	0	0	2	0
4	C	31	0	0	0	0
4	D	36	0	0	1	0
All	All	9706	0	9245	169	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 9.

All (169) 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:D:201:MSE:HE2	1:D:205:ARG:HG3	1.14	1.09
1:C:37:LYS:H	1:C:37:LYS:HD3	0.93	1.05
1:C:201:MSE:HE2	1:C:205:ARG:HG3	1.42	1.01
1:C:37:LYS:HD3	1:C:37:LYS:N	1.77	0.98
1:A:245:HIS:HD2	1:A:251:TRP:H	0.99	0.98
1:D:201:MSE:CE	1:D:205:ARG:HG3	1.94	0.96
1:C:37:LYS:H	1:C:37:LYS:CD	1.77	0.94
1:A:245:HIS:CD2	1:A:251:TRP:H	1.85	0.92
1:A:209:ARG:HH22	2:A:303[B]:GOL:H2	1.35	0.92
1:A:12:PRO:HD2	1:A:31:ASP:HB3	1.51	0.90
1:D:254:ILE:HD11	1:D:292:LEU:HD12	1.54	0.88
1:C:75:GLN:HE22	1:C:109:ASN:H	1.18	0.88
1:B:254:ILE:CD1	1:B:292:LEU:HD12	2.05	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:239:ARG:HG2	1:C:239:ARG:HH11	1.40	0.86
1:B:245:HIS:HD2	1:B:251:TRP:H	1.22	0.86
1:C:86:MSE:HE1	1:C:138:PHE:CE1	2.11	0.85
1:D:201:MSE:HE2	1:D:205:ARG:CG	2.04	0.83
1:C:245:HIS:HD2	1:C:251:TRP:H	1.25	0.81
1:B:254:ILE:HD11	1:B:292:LEU:HD12	1.63	0.78
1:D:254:ILE:CD1	1:D:292:LEU:HD12	2.12	0.77
1:C:245:HIS:CD2	1:C:251:TRP:H	2.04	0.76
1:D:201:MSE:HE3	1:D:201:MSE:O	1.85	0.76
1:A:209:ARG:NH2	2:A:303[B]:GOL:H2	2.03	0.74
1:C:201:MSE:HE3	1:C:204:LYS:HB3	1.72	0.71
1:D:26:ARG:CG	1:D:26:ARG:HH21	2.04	0.70
1:A:268:ARG:O	1:A:272:THR:HG23	1.90	0.70
1:D:108:MSE:HE1	1:D:144:PHE:CD2	2.27	0.70
1:D:108:MSE:HE1	1:D:144:PHE:HD2	1.56	0.69
1:C:174:LYS:HG2	1:C:178:ARG:HH12	1.57	0.69
1:C:159:PRO:HG3	1:C:169:TYR:CD1	2.29	0.67
1:D:154:ARG:HA	1:D:158:GLU:HB2	1.76	0.67
1:A:12:PRO:HB3	1:A:14:SER:HB2	1.75	0.67
1:C:268:ARG:O	1:C:272:THR:HG23	1.93	0.67
1:B:159:PRO:HB2	1:B:236:LYS:HE2	1.77	0.66
1:C:252:ALA:HB1	1:C:292:LEU:HD21	1.78	0.65
1:B:109:ASN:OD1	1:B:111:ARG:HG3	1.97	0.65
1:A:254:ILE:CD1	1:A:292:LEU:HD12	2.27	0.65
1:A:254:ILE:HD11	1:A:292:LEU:HD12	1.79	0.65
1:B:86:MSE:HE2	1:B:188:ILE:HG21	1.81	0.63
1:A:30:ILE:HD11	1:A:219:LEU:HD11	1.79	0.63
1:B:245:HIS:CD2	1:B:251:TRP:H	2.10	0.62
1:A:86:MSE:HE3	1:A:147:PHE:CE1	2.33	0.62
1:D:155:ALA:HB2	1:D:172:PHE:CE2	2.34	0.62
1:C:161:MSE:HE2	1:C:222:MSE:SE	2.50	0.61
1:C:86:MSE:HE1	1:C:138:PHE:CD1	2.36	0.61
1:A:245:HIS:HD2	1:A:251:TRP:N	1.83	0.61
1:B:153:ASN:H	1:B:153:ASN:HD22	1.48	0.61
1:B:154:ARG:HG2	1:B:172:PHE:HB2	1.82	0.61
1:D:86:MSE:HE2	1:D:151:TRP:HH2	1.67	0.60
1:C:175:GLU:OE2	1:C:178:ARG:NH1	2.35	0.60
1:D:119:THR:O	1:D:133:ARG:NH2	2.35	0.59
1:D:175:GLU:OE2	1:D:178:ARG:NH1	2.35	0.59
1:C:201:MSE:HE1	1:C:204:LYS:HD3	1.85	0.59
1:A:26:ARG:HG2	1:A:38:TRP:CE2	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:26:ARG:HG3	1:D:26:ARG:HH21	1.67	0.59
1:D:196:ASN:HB2	1:D:299:LEU:HD21	1.85	0.58
1:D:287:VAL:HG11	1:D:292:LEU:HD22	1.85	0.58
1:D:201:MSE:HE1	1:D:204:LYS:HD3	1.84	0.58
1:A:54:LYS:HE3	4:A:331:HOH:O	2.03	0.58
1:A:12:PRO:HD2	1:A:31:ASP:CB	2.29	0.58
1:C:154:ARG:HG2	1:C:172:PHE:HB2	1.86	0.58
1:A:97:LYS:HE3	2:A:303[A]:GOL:O1	2.04	0.57
1:C:201:MSE:HE3	1:C:201:MSE:HA	1.87	0.57
1:C:100:ALA:O	1:C:104:THR:HG22	2.05	0.57
1:B:152:TYR:HB3	1:B:240:MSE:HE2	1.87	0.57
1:D:268:ARG:HB3	1:D:287:VAL:HG22	1.86	0.57
1:A:111:ARG:NH1	1:C:148:ASP:OD2	2.38	0.56
1:D:177:PRO:HB3	1:D:246:THR:HG21	1.85	0.56
1:B:22:ASP:OD1	1:B:22:ASP:N	2.38	0.56
1:C:239:ARG:HG2	1:C:239:ARG:NH1	2.12	0.56
1:B:55:LEU:HD13	1:B:60:TYR:HB2	1.88	0.56
1:A:179:PHE:HB2	1:B:182:MSE:HE1	1.88	0.55
1:A:245:HIS:CD2	1:A:251:TRP:N	2.67	0.55
1:D:160:VAL:HG23	1:D:236:LYS:HG3	1.88	0.55
1:B:155:ALA:HB2	1:B:172:PHE:HE2	1.72	0.54
1:C:254:ILE:HD13	1:C:292:LEU:HD23	1.89	0.54
1:A:36:PRO:HB2	1:A:38:TRP:CD1	2.43	0.54
1:D:193:PHE:HZ	1:D:270:MSE:HG3	1.73	0.53
1:C:100:ALA:O	1:C:104:THR:CG2	2.57	0.53
1:D:173:LEU:HD13	1:D:243:GLU:HG3	1.91	0.53
1:A:74:VAL:O	1:A:78:MSE:HG3	2.09	0.53
1:A:20:GLU:O	1:A:199:ARG:NH1	2.42	0.52
1:B:153:ASN:N	1:B:153:ASN:HD22	2.05	0.52
2:A:303[A]:GOL:H12	4:A:373:HOH:O	2.07	0.52
1:D:201:MSE:HE3	1:D:204:LYS:HB3	1.91	0.52
1:D:114:ARG:NH2	1:D:137:THR:O	2.43	0.52
1:A:253:VAL:HG11	1:A:299:LEU:HD22	1.92	0.51
1:C:52:LYS:HB2	1:C:52:LYS:NZ	2.26	0.51
1:D:27:ILE:O	1:D:27:ILE:HG13	2.10	0.51
1:A:98:GLY:H	2:A:303[A]:GOL:H2	1.75	0.51
1:D:86:MSE:CE	1:D:151:TRP:HH2	2.24	0.50
1:D:155:ALA:HB2	1:D:172:PHE:HE2	1.76	0.50
1:B:26:ARG:HG2	1:B:38:TRP:CH2	2.46	0.50
1:A:86:MSE:HE3	1:A:147:PHE:CZ	2.47	0.50
1:C:124:THR:O	1:C:128:GLN:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:253:VAL:HG11	1:C:299:LEU:HG	1.94	0.50
1:D:159:PRO:HG3	1:D:169:TYR:CD1	2.47	0.50
1:C:49:TYR:CD2	1:C:257:ASN:HB2	2.47	0.49
1:D:36:PRO:HD2	1:D:39:ILE:HD12	1.93	0.49
1:A:30:ILE:CD1	1:A:219:LEU:HD11	2.42	0.49
1:A:30:ILE:O	1:A:211:HIS:CE1	2.66	0.49
1:C:75:GLN:O	1:C:79:GLN:HG2	2.13	0.49
1:A:209:ARG:HH22	2:A:303[B]:GOL:C2	2.14	0.49
1:A:155:ALA:HB2	1:A:172:PHE:CE2	2.48	0.49
1:A:30:ILE:O	1:A:211:HIS:HE1	1.96	0.49
1:B:154:ARG:HG2	1:B:172:PHE:CB	2.43	0.48
1:C:108:MSE:HE1	1:C:144:PHE:CD2	2.48	0.48
1:C:224:ILE:HA	1:C:227:LEU:HD22	1.95	0.48
1:C:232:ASP:O	1:C:236:LYS:HD3	2.13	0.48
1:A:86:MSE:CE	1:A:147:PHE:CE1	2.96	0.48
1:A:240:MSE:HE1	4:A:348:HOH:O	2.15	0.47
1:A:155:ALA:HB2	1:A:172:PHE:HE2	1.78	0.47
1:D:239:ARG:O	1:D:243:GLU:HG2	2.14	0.47
1:A:75:GLN:O	1:A:79:GLN:HG2	2.15	0.47
1:D:121:PRO:HB2	1:D:126:ARG:HG3	1.97	0.47
1:A:171:GLN:HE21	1:A:171:GLN:HA	1.80	0.46
1:B:153:ASN:ND2	1:B:153:ASN:N	2.63	0.46
1:C:154:ARG:HD2	1:C:164:CYS:SG	2.55	0.46
1:A:240:MSE:CE	4:A:348:HOH:O	2.64	0.46
1:B:155:ALA:HB2	1:B:172:PHE:CE2	2.50	0.46
1:C:179:PHE:HB2	1:D:182:MSE:HE1	1.98	0.46
1:A:239:ARG:O	1:A:243:GLU:HG2	2.16	0.46
1:D:158:GLU:HB3	1:D:164:CYS:SG	2.56	0.45
1:C:154:ARG:O	1:C:169:TYR:HD1	1.98	0.45
1:D:157:VAL:O	1:D:161:MSE:HB2	2.17	0.45
1:D:86:MSE:CE	1:D:151:TRP:CH2	2.99	0.45
1:B:154:ARG:O	1:B:169:TYR:HD1	2.00	0.45
1:C:201:MSE:CE	1:C:204:LYS:HB3	2.43	0.45
1:B:175:GLU:OE2	1:B:178:ARG:NH1	2.50	0.44
1:C:154:ARG:HG2	1:C:172:PHE:CB	2.47	0.44
1:A:182:MSE:HE1	1:B:179:PHE:HB2	1.98	0.44
1:C:174:LYS:HE3	1:C:178:ARG:HH12	1.82	0.44
1:C:49:TYR:CD2	1:C:50:PRO:HD2	2.53	0.44
1:D:160:VAL:CG2	1:D:236:LYS:HG3	2.48	0.44
1:A:196:ASN:O	1:A:256:GLY:N	2.48	0.44
1:A:222[A]:MSE:HE2	4:A:391:HOH:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:LYS:O	1:B:263:ARG:HG3	2.17	0.43
1:A:57:ARG:O	1:A:61:GLU:HG2	2.17	0.43
1:C:73:LYS:HD3	1:C:278:GLY:O	2.18	0.43
1:A:152:TYR:HB3	1:A:240:MSE:HE2	2.00	0.43
1:C:181:GLU:HG3	1:C:248:HIS:CE1	2.54	0.43
1:C:140:THR:O	1:C:141:ALA:C	2.56	0.43
1:C:35:LEU:HD12	1:C:36:PRO:HD2	2.00	0.43
1:A:15:ARG:HD3	1:A:31:ASP:OD2	2.19	0.43
1:A:214:LEU:HD11	1:C:65:THR:HG23	2.00	0.43
1:A:259:LYS:O	1:A:263:ARG:HG3	2.20	0.42
1:B:92:ARG:HD3	1:B:230:TRP:CZ3	2.55	0.42
1:C:55:LEU:HD13	1:C:60:TYR:HB2	2.00	0.42
1:A:170:GLU:CD	1:A:239:ARG:HH22	2.19	0.42
1:B:224:ILE:O	1:B:227:LEU:HB2	2.20	0.42
1:C:196:ASN:O	1:C:256:GLY:N	2.49	0.42
1:C:30:ILE:O	1:C:211:HIS:CE1	2.72	0.42
1:B:269:HIS:CE1	1:B:273:LYS:HE3	2.55	0.41
1:C:108:MSE:HE2	1:C:113:ALA:HB2	2.01	0.41
1:B:182:MSE:CB	4:B:320:HOH:O	2.68	0.41
1:D:133:ARG:HD3	4:D:323:HOH:O	2.20	0.41
1:D:192:LYS:HB2	1:D:192:LYS:HE3	1.92	0.41
1:A:75:GLN:NE2	1:A:112:SER:OG	2.53	0.41
1:C:154:ARG:HD3	1:C:158:GLU:OE1	2.20	0.41
1:D:196:ASN:HB3	1:D:254:ILE:O	2.20	0.41
1:B:132:GLN:NE2	4:B:311:HOH:O	2.54	0.41
1:A:67:LEU:HB3	1:A:267:ILE:HG21	2.03	0.41
1:B:57:ARG:HG3	1:B:58:GLU:N	2.36	0.41
1:B:77:TRP:CH2	1:B:83:LYS:HG2	2.56	0.40
1:C:293:GLY:HA3	1:C:298:PHE:CD2	2.57	0.40
1:B:26:ARG:HG2	1:B:38:TRP:CZ2	2.57	0.40
1:C:160:VAL:HG11	1:C:233:TYR:CE2	2.56	0.40
1:C:66:LYS:HE3	1:C:283:ALA:HA	2.02	0.40
1:C:75:GLN:HE22	1:C:109:ASN:N	2.01	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	287/304 (94%)	282 (98%)	5 (2%)	0	100	100
1	B	285/304 (94%)	279 (98%)	6 (2%)	0	100	100
1	C	283/304 (93%)	272 (96%)	11 (4%)	0	100	100
1	D	284/304 (93%)	274 (96%)	10 (4%)	0	100	100
All	All	1139/1216 (94%)	1107 (97%)	32 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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	247/247 (100%)	232 (94%)	15 (6%)	18	16
1	B	245/247 (99%)	228 (93%)	17 (7%)	15	12
1	C	244/247 (99%)	224 (92%)	20 (8%)	11	8
1	D	245/247 (99%)	225 (92%)	20 (8%)	11	8
All	All	981/988 (99%)	909 (93%)	72 (7%)	13	10

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	PRO
1	A	14	SER

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Mol	Chain	Res	Type
1	A	24	ARG
1	A	26	ARG
1	A	45	ARG
1	A	153	ASN
1	A	170	GLU
1	A	171	GLN
1	A	196	ASN
1	A	227	LEU
1	A	232	ASP
1	A	239	ARG
1	A	240	MSE
1	A	274	LEU
1	A	281	GLU
1	B	22	ASP
1	B	24	ARG
1	B	25	SER
1	B	26	ARG
1	B	35	LEU
1	B	55	LEU
1	B	57	ARG
1	B	144	PHE
1	B	154	ARG
1	B	167	ASP
1	B	196	ASN
1	B	219	LEU
1	B	227	LEU
1	B	239	ARG
1	B	240	MSE
1	B	273	LYS
1	B	299	LEU
1	C	20	GLU
1	C	26	ARG
1	C	27	ILE
1	C	37	LYS
1	C	55	LEU
1	C	62	GLU
1	C	75	GLN
1	C	86	MSE
1	C	104	THR
1	C	120	LYS
1	C	154	ARG
1	C	196	ASN

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Mol	Chain	Res	Type
1	C	222	MSE
1	C	227	LEU
1	C	232	ASP
1	C	239	ARG
1	C	273	LYS
1	C	281	GLU
1	C	286	GLU
1	C	292	LEU
1	D	14	SER
1	D	21	ILE
1	D	26	ARG
1	D	27	ILE
1	D	35	LEU
1	D	47	ASP
1	D	48	ASP
1	D	55	LEU
1	D	114	ARG
1	D	123	GLU
1	D	144	PHE
1	D	161	MSE
1	D	167	ASP
1	D	196	ASN
1	D	220	SER
1	D	222	MSE
1	D	227	LEU
1	D	233	TYR
1	D	240	MSE
1	D	274	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	GLN
1	A	132	GLN
1	A	153	ASN
1	A	171	GLN
1	A	185	ASN
1	A	211	HIS
1	A	245	HIS
1	B	132	GLN
1	B	153	ASN
1	B	211	HIS

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Mol	Chain	Res	Type
1	B	245	HIS
1	B	269	HIS
1	C	68	GLN
1	C	75	GLN
1	C	107	ASN
1	C	132	GLN
1	C	153	ASN
1	C	211	HIS
1	C	245	HIS
1	C	248	HIS
1	D	132	GLN
1	D	211	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry

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	$\#\|Z\| > 2$
3	FMT	C	303	-	0,2,2	0.00	-	0,1,1	0.00	-
2	GOL	A	303[B]	-	5,5,5	0.35	0	5,5,5	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	FMT	D	304	-	0,2,2	0.00	-	0,1,1	0.00	-
2	GOL	A	303[A]	-	5,5,5	0.41	0	5,5,5	0.45	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
2	GOL	A	303[A]	-	-	2/4/4/4	-
2	GOL	A	303[B]	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	303[A]	GOL	C1-C2-C3-O3
2	A	303[A]	GOL	O2-C2-C3-O3
2	A	303[B]	GOL	C1-C2-C3-O3
2	A	303[B]	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	303[B]	GOL	3	0
2	A	303[A]	GOL	3	0

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	279/304 (91%)	0.26	12 (4%)	35	34	35, 49, 66, 75	0
1	B	278/304 (91%)	0.28	17 (6%)	21	20	36, 50, 67, 73	0
1	C	275/304 (90%)	0.60	22 (8%)	12	12	50, 65, 82, 87	0
1	D	277/304 (91%)	0.55	23 (8%)	11	11	47, 60, 76, 83	0
All	All	1109/1216 (91%)	0.42	74 (6%)	17	17	35, 58, 76, 87	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	163	PHE	6.8
1	D	164	CYS	6.3
1	C	28	PHE	4.6
1	A	12	PRO	4.0
1	C	155	ALA	3.9
1	C	164	CYS	3.8
1	D	295	GLY	3.8
1	C	157	VAL	3.8
1	C	296	PRO	3.7
1	D	155	ALA	3.5
1	C	21	ILE	3.5
1	D	115	VAL	3.5
1	B	159	PRO	3.4
1	A	164	CYS	3.3
1	C	19	LEU	3.3
1	D	160	VAL	3.3
1	D	272	THR	3.3
1	D	296	PRO	3.2
1	C	163	PHE	3.2
1	A	115	VAL	3.2
1	D	159	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	146	LEU	3.1
1	B	167	ASP	3.1
1	D	165	THR	3.1
1	B	163	PHE	3.0
1	B	146	LEU	2.9
1	B	157	VAL	2.9
1	A	146	LEU	2.9
1	C	17	VAL	2.9
1	D	157	VAL	2.9
1	B	160	VAL	2.7
1	C	102[A]	HIS	2.7
1	D	248	HIS	2.6
1	B	115	VAL	2.6
1	D	299	LEU	2.6
1	D	154	ARG	2.6
1	D	167	ASP	2.6
1	D	147	PHE	2.6
1	C	154	ARG	2.6
1	D	28	PHE	2.5
1	C	223	ASP	2.5
1	C	27	ILE	2.5
1	C	167	ASP	2.4
1	A	155	ALA	2.4
1	A	101	ILE	2.4
1	D	27	ILE	2.4
1	A	148	ASP	2.4
1	D	158	GLU	2.3
1	D	242	LYS	2.3
1	D	113	ALA	2.3
1	C	23	GLY	2.3
1	B	166	PRO	2.3
1	B	101	ILE	2.3
1	A	163	PHE	2.2
1	B	105	THR	2.2
1	B	155	ALA	2.2
1	B	164	CYS	2.2
1	A	88	VAL	2.1
1	B	102	HIS	2.1
1	B	145	VAL	2.1
1	C	233	TYR	2.1
1	D	20	GLU	2.1
1	C	299	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	106	ALA	2.1
1	C	87	ALA	2.1
1	A	158	GLU	2.1
1	B	116	VAL	2.1
1	C	278	GLY	2.0
1	A	167	ASP	2.0
1	C	53	LYS	2.0
1	C	146	LEU	2.0
1	C	20	GLU	2.0
1	B	89	PHE	2.0
1	A	165	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 carbohydrates 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	FMT	D	304	3/3	0.83	0.21	62,62,63,63	0
3	FMT	C	303	3/3	0.89	0.20	70,70,70,70	0
2	GOL	A	303[B]	6/6	0.91	0.30	47,49,50,51	6
2	GOL	A	303[A]	6/6	0.91	0.30	26,36,37,38	6

6.5 Other polymers (i)

There are no such residues in this entry.