Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2023 – 05:22 AM EDT

PDB ID	:	3LG2
Title	:	A Ykr043C/ fructose-1,6-bisphosphate product complex following ligand soaking
Authors	:	Singer, A.; Xu, X.; Cui, H.; Dong, A.; Edwards, A.M.; Joachimiak, A.; Yakunin, A.F.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2010-01-19
Resolution	:	2.60 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:

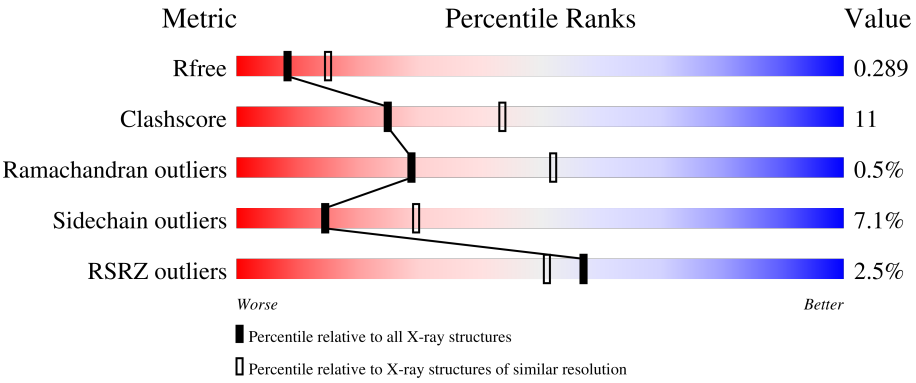
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	292	% 68% 22% • 8%
1	B	292	3% 68% 21% • 8%
1	C	292	5% 69% 23% • 8%
1	D	292	71% 19% • 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	B	300	-	-	X	-
2	PO4	B	302	-	-	X	-
2	PO4	C	272	-	-	X	-
2	PO4	C	302	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 8732 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 Uncharacterized protein YKR043C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	Se	0	0	0
			2172	1362	391	412	4	3			
1	B	269	Total	C	N	O	S	Se	0	0	0
			2172	1362	391	412	4	3			
1	C	269	Total	C	N	O	S	Se	0	0	0
			2172	1362	391	412	4	3			
1	D	269	Total	C	N	O	S	Se	0	0	0
			2172	1362	391	412	4	3			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MSE	-	expression tag	UNP P36136
A	-19	GLY	-	expression tag	UNP P36136
A	-18	SER	-	expression tag	UNP P36136
A	-17	SER	-	expression tag	UNP P36136
A	-16	HIS	-	expression tag	UNP P36136
A	-15	HIS	-	expression tag	UNP P36136
A	-14	HIS	-	expression tag	UNP P36136
A	-13	HIS	-	expression tag	UNP P36136
A	-12	HIS	-	expression tag	UNP P36136
A	-11	HIS	-	expression tag	UNP P36136
A	-10	SER	-	expression tag	UNP P36136
A	-9	SER	-	expression tag	UNP P36136
A	-8	GLY	-	expression tag	UNP P36136
A	-7	ARG	-	expression tag	UNP P36136
A	-6	GLU	-	expression tag	UNP P36136
A	-5	ASN	-	expression tag	UNP P36136
A	-4	LEU	-	expression tag	UNP P36136
A	-3	TYR	-	expression tag	UNP P36136
A	-2	PHE	-	expression tag	UNP P36136
A	-1	GLN	-	expression tag	UNP P36136
A	0	GLY	-	expression tag	UNP P36136

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MSE	-	expression tag	UNP P36136
B	-19	GLY	-	expression tag	UNP P36136
B	-18	SER	-	expression tag	UNP P36136
B	-17	SER	-	expression tag	UNP P36136
B	-16	HIS	-	expression tag	UNP P36136
B	-15	HIS	-	expression tag	UNP P36136
B	-14	HIS	-	expression tag	UNP P36136
B	-13	HIS	-	expression tag	UNP P36136
B	-12	HIS	-	expression tag	UNP P36136
B	-11	HIS	-	expression tag	UNP P36136
B	-10	SER	-	expression tag	UNP P36136
B	-9	SER	-	expression tag	UNP P36136
B	-8	GLY	-	expression tag	UNP P36136
B	-7	ARG	-	expression tag	UNP P36136
B	-6	GLU	-	expression tag	UNP P36136
B	-5	ASN	-	expression tag	UNP P36136
B	-4	LEU	-	expression tag	UNP P36136
B	-3	TYR	-	expression tag	UNP P36136
B	-2	PHE	-	expression tag	UNP P36136
B	-1	GLN	-	expression tag	UNP P36136
B	0	GLY	-	expression tag	UNP P36136
C	-20	MSE	-	expression tag	UNP P36136
C	-19	GLY	-	expression tag	UNP P36136
C	-18	SER	-	expression tag	UNP P36136
C	-17	SER	-	expression tag	UNP P36136
C	-16	HIS	-	expression tag	UNP P36136
C	-15	HIS	-	expression tag	UNP P36136
C	-14	HIS	-	expression tag	UNP P36136
C	-13	HIS	-	expression tag	UNP P36136
C	-12	HIS	-	expression tag	UNP P36136
C	-11	HIS	-	expression tag	UNP P36136
C	-10	SER	-	expression tag	UNP P36136
C	-9	SER	-	expression tag	UNP P36136
C	-8	GLY	-	expression tag	UNP P36136
C	-7	ARG	-	expression tag	UNP P36136
C	-6	GLU	-	expression tag	UNP P36136
C	-5	ASN	-	expression tag	UNP P36136
C	-4	LEU	-	expression tag	UNP P36136
C	-3	TYR	-	expression tag	UNP P36136
C	-2	PHE	-	expression tag	UNP P36136
C	-1	GLN	-	expression tag	UNP P36136
C	0	GLY	-	expression tag	UNP P36136

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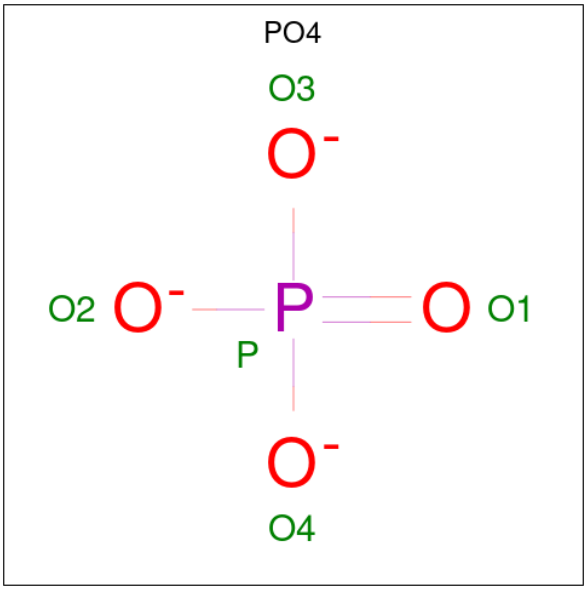
Chain	Residue	Modelled	Actual	Comment	Reference
D	-20	MSE	-	expression tag	UNP P36136
D	-19	GLY	-	expression tag	UNP P36136
D	-18	SER	-	expression tag	UNP P36136
D	-17	SER	-	expression tag	UNP P36136
D	-16	HIS	-	expression tag	UNP P36136
D	-15	HIS	-	expression tag	UNP P36136
D	-14	HIS	-	expression tag	UNP P36136
D	-13	HIS	-	expression tag	UNP P36136
D	-12	HIS	-	expression tag	UNP P36136
D	-11	HIS	-	expression tag	UNP P36136
D	-10	SER	-	expression tag	UNP P36136
D	-9	SER	-	expression tag	UNP P36136
D	-8	GLY	-	expression tag	UNP P36136
D	-7	ARG	-	expression tag	UNP P36136
D	-6	GLU	-	expression tag	UNP P36136
D	-5	ASN	-	expression tag	UNP P36136
D	-4	LEU	-	expression tag	UNP P36136
D	-3	TYR	-	expression tag	UNP P36136
D	-2	PHE	-	expression tag	UNP P36136
D	-1	GLN	-	expression tag	UNP P36136
D	0	GLY	-	expression tag	UNP P36136

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



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

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

- Molecule 3 is water.

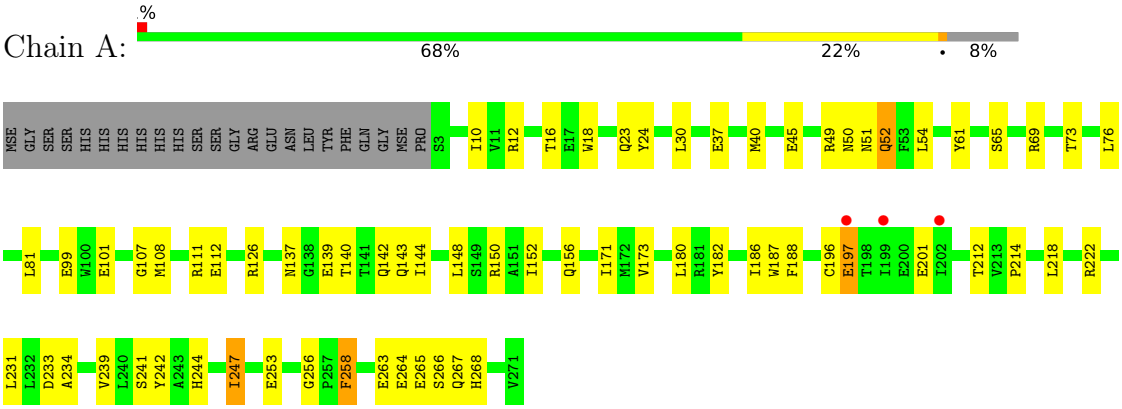
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	B	2	Total O 2 2	0	0
3	C	1	Total O 1 1	0	0

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.

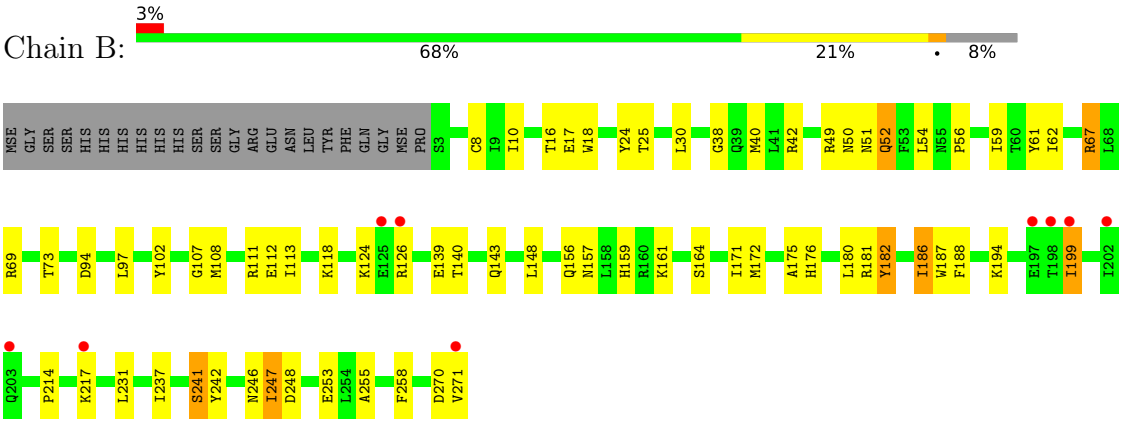
• Molecule 1: Uncharacterized protein YKR043C

Chain A:



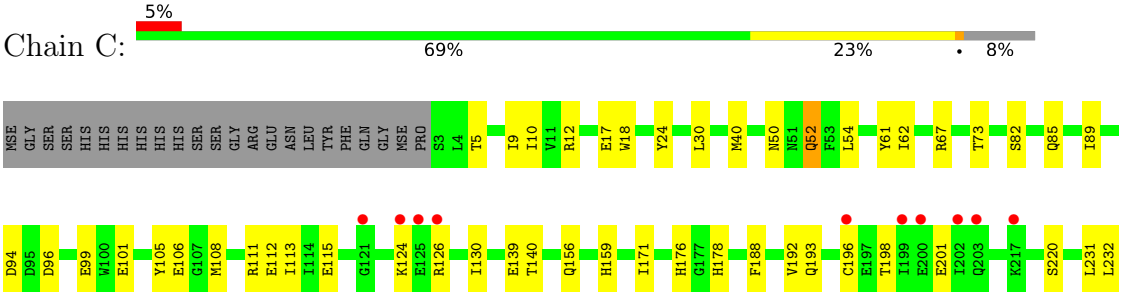
• Molecule 1: Uncharacterized protein YKR043C

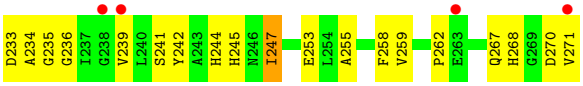
Chain B:



• Molecule 1: Uncharacterized protein YKR043C

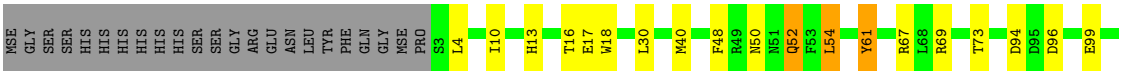
Chain C:

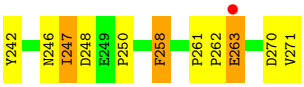




• Molecule 1: Uncharacterized protein YKR043C

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	83.30Å 83.39Å 104.69Å 90.00° 104.19° 90.00°	Depositor
Resolution (Å)	43.35 – 2.60 43.35 – 2.60	Depositor EDS
% Data completeness (in resolution range)	91.2 (43.35-2.60) 91.2 (43.35-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.45 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.220 , 0.270 0.242 , 0.289	Depositor DCC
R_{free} test set	2008 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å^2)	36.4	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 30.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	8732	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 62.41 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1303e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: 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	0.73	0/2218	0.78	1/2999 (0.0%)
1	B	0.71	0/2218	0.80	0/2999
1	C	0.81	1/2218 (0.0%)	0.84	0/2999
1	D	0.81	1/2218 (0.0%)	0.87	1/2999 (0.0%)
All	All	0.77	2/8872 (0.0%)	0.82	2/11996 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	61	TYR	CD1-CE1	5.56	1.47	1.39
1	C	12	ARG	CB-CG	5.32	1.66	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	147	ARG	NE-CZ-NH1	5.25	122.92	120.30
1	A	12	ARG	NE-CZ-NH1	5.18	122.89	120.30

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	2172	0	2114	45	0
1	B	2172	0	2114	55	0
1	C	2172	0	2114	48	0
1	D	2172	0	2114	44	0
2	A	5	0	0	0	0
2	B	15	0	0	7	0
2	C	15	0	0	4	0
2	D	5	0	0	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
All	All	8732	0	8456	186	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 11.

All (186) 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:C:247:ILE:O	1:C:247:ILE:HD12	1.47	1.15
1:C:50:ASN:HB2	1:C:52:GLN:NE2	1.63	1.13
1:D:108:MSE:HE2	1:D:113:ILE:HG13	1.28	1.10
1:D:108:MSE:HE1	1:D:113:ILE:HA	1.41	1.01
1:D:199:ILE:O	1:D:203:GLN:HG2	1.63	0.99
1:B:50:ASN:HB2	1:B:52:GLN:HE22	1.26	0.98
1:C:176:HIS:HB3	2:C:302:PO4:O2	1.64	0.96
1:B:108:MSE:HE3	1:B:112:GLU:HB3	1.48	0.95
1:C:50:ASN:HB2	1:C:52:GLN:HE22	1.19	0.93
1:C:82:SER:OG	1:C:85:GLN:HG3	1.74	0.88
1:D:108:MSE:HE2	1:D:113:ILE:CG1	2.06	0.84
1:C:50:ASN:CB	1:C:52:GLN:NE2	2.42	0.83
1:D:50:ASN:HB2	1:D:52:GLN:NE2	1.94	0.83
1:D:40:MSE:HG3	1:D:73:THR:HG23	1.64	0.80
1:B:181:ARG:NH2	2:B:300:PO4:O1	2.14	0.79
1:A:50:ASN:HB2	1:A:52:GLN:NE2	1.98	0.79
1:A:50:ASN:CB	1:A:52:GLN:HE22	1.97	0.78
1:A:50:ASN:HB2	1:A:52:GLN:HE22	1.47	0.78
1:C:61:TYR:HB2	1:C:171:ILE:HG12	1.67	0.76
1:C:50:ASN:CB	1:C:52:GLN:HE22	1.99	0.75
1:D:108:MSE:CE	1:D:113:ILE:HG13	2.14	0.74
1:B:108:MSE:HE1	1:B:113:ILE:HA	1.71	0.73
1:A:108:MSE:HE2	1:A:112:GLU:HB3	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:GLU:HG2	1:C:140:THR:CB	2.21	0.71
1:B:181:ARG:NH1	2:B:300:PO4:O1	2.23	0.70
1:C:268:HIS:ND1	2:C:272:PO4:O2	2.25	0.70
1:D:108:MSE:CE	1:D:113:ILE:HA	2.19	0.69
1:A:50:ASN:CB	1:A:52:GLN:NE2	2.55	0.69
1:B:69:ARG:NH2	2:B:302:PO4:O3	2.23	0.69
1:A:197:GLU:HG2	1:C:140:THR:HB	1.74	0.69
1:C:193:GLN:HE21	1:C:220:SER:HB2	1.58	0.69
1:A:101:GLU:HB2	1:A:139:GLU:HG3	1.76	0.68
1:B:157:ASN:OD1	1:B:161:LYS:HE2	1.94	0.67
1:D:50:ASN:CB	1:D:52:GLN:NE2	2.57	0.67
1:D:40:MSE:HE1	1:D:235:GLY:HA2	1.76	0.67
1:A:23:GLN:HG2	1:A:107:GLY:O	1.94	0.67
1:D:108:MSE:HE1	1:D:113:ILE:CA	2.21	0.66
1:C:101:GLU:HB2	1:C:139:GLU:HG3	1.78	0.66
1:B:16:THR:HG21	1:B:69:ARG:HD2	1.79	0.65
1:A:197:GLU:HG2	1:C:140:THR:CG2	2.27	0.64
1:A:197:GLU:HG2	1:C:140:THR:HG21	1.80	0.64
1:C:268:HIS:HB2	2:C:272:PO4:O2	1.97	0.64
1:B:62:ILE:HG12	1:B:172:MSE:HB2	1.80	0.64
1:A:268:HIS:HB2	2:B:300:PO4:O2	1.98	0.63
1:B:156:GLN:HG2	1:B:242:TYR:OH	1.99	0.62
1:D:61:TYR:HB2	1:D:171:ILE:HG12	1.82	0.62
1:C:40:MSE:HG3	1:C:73:THR:HG23	1.80	0.62
1:D:67:ARG:NE	1:D:67:ARG:HA	2.15	0.62
1:B:40:MSE:HG3	1:B:73:THR:HG23	1.82	0.61
1:C:242:TYR:CD1	1:C:247:ILE:HA	2.35	0.61
1:C:108:MSE:HE3	1:C:112:GLU:HB3	1.82	0.60
1:D:67:ARG:HA	1:D:67:ARG:HE	1.67	0.60
1:A:18:TRP:CG	1:A:30:LEU:HD11	2.37	0.60
1:D:108:MSE:CE	1:D:113:ILE:CA	2.79	0.59
1:B:50:ASN:HB2	1:B:52:GLN:NE2	2.08	0.59
1:B:40:MSE:HE2	1:B:73:THR:CG2	2.32	0.59
1:A:37:GLU:HG2	1:A:76:LEU:HD11	1.85	0.58
1:B:253:GLU:HG2	1:B:255:ALA:HB2	1.86	0.58
1:A:233:ASP:OD1	1:A:234:ALA:N	2.37	0.57
1:A:148:LEU:O	1:A:152:ILE:HG13	2.05	0.57
1:A:150:ARG:HG3	1:A:218:LEU:HD11	1.86	0.57
1:D:52:GLN:HB2	1:D:258:PHE:HB3	1.86	0.57
1:A:242:TYR:CD1	1:A:247:ILE:HA	2.39	0.57
1:D:108:MSE:CE	1:D:113:ILE:N	2.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:THR:HG21	1:A:69:ARG:HD2	1.86	0.56
1:A:40:MSE:HG3	1:A:73:THR:HG23	1.87	0.56
1:D:10:ILE:HB	1:D:180:LEU:HD13	1.87	0.56
1:D:242:TYR:CE1	1:D:247:ILE:HD12	2.41	0.56
1:B:181:ARG:CZ	2:B:300:PO4:O1	2.53	0.56
1:C:40:MSE:HE1	1:C:235:GLY:HA2	1.88	0.55
1:B:247:ILE:HG13	1:B:247:ILE:O	2.05	0.55
1:C:50:ASN:HB2	1:C:52:GLN:HE21	1.68	0.54
1:B:61:TYR:HB2	1:B:171:ILE:HG12	1.89	0.54
1:C:253:GLU:HG2	1:C:255:ALA:HB2	1.90	0.54
1:D:16:THR:HG21	1:D:69:ARG:HD2	1.89	0.53
1:C:247:ILE:HD12	1:C:247:ILE:C	2.24	0.53
1:A:24:TYR:N	1:A:24:TYR:CD1	2.76	0.53
1:C:108:MSE:HE2	1:C:113:ILE:HG13	1.91	0.53
1:D:4:LEU:HD23	1:D:262:PRO:HD3	1.90	0.53
1:B:242:TYR:CD1	1:B:247:ILE:HA	2.43	0.52
1:B:242:TYR:CZ	1:B:247:ILE:HD12	2.44	0.52
1:B:108:MSE:CE	1:B:113:ILE:N	2.72	0.52
1:D:48:PHE:CE1	1:D:54:LEU:HD13	2.45	0.52
1:B:56:PRO:HA	1:B:59:ILE:HG13	1.91	0.52
1:A:52:GLN:HB2	1:A:258:PHE:HB3	1.91	0.52
1:B:18:TRP:CZ3	1:B:107:GLY:HA2	2.45	0.51
1:B:157:ASN:HD21	1:B:214:PRO:HG2	1.75	0.51
1:C:108:MSE:HE1	1:C:112:GLU:C	2.31	0.51
1:C:232:LEU:HD12	1:C:236:GLY:HA3	1.92	0.51
1:C:159:HIS:CD2	1:C:171:ILE:HD12	2.46	0.51
1:A:50:ASN:HB3	1:A:52:GLN:HE22	1.72	0.50
1:D:198:THR:OG1	1:D:201:GLU:HG3	2.11	0.50
1:B:139:GLU:HA	1:B:143:GLN:OE1	2.11	0.50
1:D:132:ARG:HB2	1:D:225:VAL:O	2.12	0.50
1:D:94:ASP:OD1	1:D:96:ASP:HB2	2.11	0.50
1:A:186:ILE:O	1:A:187:TRP:C	2.50	0.50
1:A:256:GLY:HA2	1:B:237:ILE:HG13	1.92	0.49
1:B:157:ASN:ND2	1:B:214:PRO:HG2	2.27	0.49
1:A:10:ILE:HG22	1:A:173:VAL:HB	1.94	0.49
1:B:18:TRP:CG	1:B:30:LEU:HD11	2.48	0.49
1:C:108:MSE:HE3	1:C:112:GLU:CB	2.43	0.49
1:B:140:THR:OG1	1:B:143:GLN:HG3	2.12	0.49
1:C:18:TRP:CG	1:C:30:LEU:HD11	2.48	0.49
1:B:148:LEU:HD13	1:B:182:TYR:HB3	1.95	0.49
1:C:67:ARG:NE	1:C:67:ARG:HA	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:176:HIS:CB	2:C:302:PO4:O2	2.50	0.48
1:C:247:ILE:O	1:C:247:ILE:CD1	2.39	0.48
1:B:270:ASP:C	1:B:271:VAL:HG23	2.34	0.48
1:B:108:MSE:CE	1:B:113:ILE:HA	2.42	0.48
1:B:186:ILE:O	1:B:187:TRP:C	2.50	0.47
1:D:50:ASN:CB	1:D:52:GLN:HE21	2.27	0.47
1:D:99:GLU:OE1	1:D:178:HIS:HB2	2.14	0.47
1:A:239:VAL:HG12	1:A:253:GLU:HB3	1.97	0.47
1:D:18:TRP:CZ3	1:D:107:GLY:HA2	2.49	0.47
1:B:176:HIS:ND1	2:B:302:PO4:O3	2.48	0.47
1:D:61:TYR:O	1:D:171:ILE:HA	2.15	0.46
1:A:196:CYS:HB3	1:A:201:GLU:O	2.14	0.46
1:C:94:ASP:OD1	1:C:96:ASP:HB2	2.15	0.46
1:D:159:HIS:CD2	1:D:171:ILE:HD12	2.51	0.46
1:C:233:ASP:OD1	1:C:234:ALA:N	2.49	0.46
1:D:247:ILE:O	1:D:247:ILE:HG13	2.15	0.46
1:D:261:PRO:HB2	1:D:263:GLU:OE1	2.15	0.46
1:B:67:ARG:HG3	1:B:69:ARG:NH1	2.31	0.46
1:C:5:THR:OG1	1:C:245:HIS:ND1	2.49	0.45
1:B:52:GLN:HB2	1:B:258:PHE:HB3	1.98	0.45
1:C:50:ASN:CB	1:C:52:GLN:HE21	2.26	0.45
1:C:99:GLU:OE1	1:C:178:HIS:HB2	2.16	0.45
1:A:263:GLU:HG2	1:A:264:GLU:HG2	1.98	0.45
1:A:81:LEU:HD12	1:A:81:LEU:N	2.32	0.45
1:D:124:LYS:HD2	1:D:124:LYS:HA	1.87	0.45
1:C:156:GLN:HG2	1:C:242:TYR:OH	2.17	0.45
1:B:108:MSE:CE	1:B:112:GLU:C	2.86	0.44
1:A:108:MSE:HE2	1:A:112:GLU:CB	2.45	0.44
1:B:67:ARG:HH21	1:B:67:ARG:HA	1.83	0.44
1:C:52:GLN:HB2	1:C:258:PHE:HB3	1.99	0.44
1:A:50:ASN:HB3	1:A:52:GLN:NE2	2.32	0.44
1:C:239:VAL:HG12	1:C:253:GLU:HB3	1.99	0.44
1:D:246:ASN:OD1	1:D:248:ASP:HB2	2.17	0.44
1:A:212:THR:O	1:A:214:PRO:HD3	2.17	0.44
1:B:148:LEU:CD1	1:B:182:TYR:HB3	2.48	0.44
1:A:140:THR:HG23	1:A:143:GLN:OE1	2.18	0.43
1:B:108:MSE:HE3	1:B:112:GLU:CB	2.35	0.43
1:A:99:GLU:HA	1:A:99:GLU:OE1	2.18	0.43
1:C:105:TYR:O	1:C:106:GLU:C	2.57	0.43
1:B:159:HIS:CD2	1:B:171:ILE:HD12	2.54	0.43
1:B:108:MSE:HE1	1:B:113:ILE:CA	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:TYR:O	1:A:171:ILE:HA	2.18	0.43
1:B:38:GLY:O	1:B:42:ARG:HG3	2.18	0.43
1:D:228:PRO:HB2	1:D:230:PHE:CD2	2.54	0.43
1:C:270:ASP:O	1:C:271:VAL:HB	2.18	0.43
1:A:137:ASN:ND2	1:C:124:LYS:HE3	2.34	0.42
1:B:40:MSE:HE2	1:B:73:THR:HG21	2.00	0.42
1:C:24:TYR:CD1	1:C:24:TYR:N	2.87	0.42
1:A:244:HIS:H	1:A:267:GLN:NE2	2.17	0.42
1:C:196:CYS:HB3	1:C:201:GLU:O	2.18	0.42
1:B:242:TYR:CE1	1:B:247:ILE:HD12	2.54	0.42
1:B:8:CYS:HA	1:B:171:ILE:O	2.20	0.42
1:C:61:TYR:O	1:C:171:ILE:HA	2.19	0.42
1:A:24:TYR:N	1:A:24:TYR:HD1	2.17	0.42
1:A:156:GLN:HG2	1:A:242:TYR:OH	2.18	0.42
1:A:50:ASN:O	1:A:51:ASN:HB2	2.20	0.42
1:A:148:LEU:HD23	1:A:148:LEU:HA	1.89	0.42
1:D:108:MSE:HE3	1:D:113:ILE:N	2.34	0.42
1:B:241:SER:OG	1:B:242:TYR:N	2.52	0.42
1:A:268:HIS:ND1	2:B:300:PO4:O2	2.51	0.41
1:D:13:HIS:CG	1:D:69:ARG:HG3	2.55	0.41
1:B:50:ASN:O	1:B:51:ASN:HB2	2.20	0.41
1:D:40:MSE:CE	1:D:235:GLY:HA2	2.47	0.41
1:D:248:ASP:O	1:D:250:PRO:HD3	2.20	0.41
1:B:199:ILE:H	1:B:199:ILE:HG12	1.58	0.41
1:B:270:ASP:OD1	1:B:271:VAL:N	2.53	0.41
1:C:62:ILE:HD12	1:C:89:ILE:HG21	2.03	0.41
1:D:18:TRP:CG	1:D:30:LEU:HD11	2.55	0.41
1:D:152:ILE:CD1	1:D:186:ILE:HG13	2.51	0.41
1:D:222:ARG:HG3	1:D:222:ARG:NH2	2.36	0.41
1:A:180:LEU:HD23	1:A:180:LEU:HA	1.88	0.41
1:B:246:ASN:OD1	1:B:248:ASP:HB2	2.21	0.41
1:C:244:HIS:H	1:C:267:GLN:HE22	1.69	0.41
1:A:144:ILE:HG23	1:A:182:TYR:CE2	2.54	0.40
1:D:270:ASP:C	1:D:271:VAL:HG23	2.41	0.40
1:B:94:ASP:O	1:B:97:LEU:HB2	2.21	0.40
1:B:10:ILE:HB	1:B:180:LEU:HD13	2.03	0.40
1:D:105:TYR:O	1:D:106:GLU:C	2.58	0.40
1:B:24:TYR:HD2	1:B:102:TYR:CE1	2.39	0.40
1:B:180:LEU:HD23	1:B:180:LEU:HA	1.87	0.40
1:B:270:ASP:O	1:B:271:VAL:HG23	2.21	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	267/292 (91%)	257 (96%)	10 (4%)	0	100	100
1	B	267/292 (91%)	257 (96%)	8 (3%)	2 (1%)	22	43
1	C	267/292 (91%)	255 (96%)	11 (4%)	1 (0%)	34	57
1	D	267/292 (91%)	258 (97%)	7 (3%)	2 (1%)	22	43
All	All	1068/1168 (91%)	1027 (96%)	36 (3%)	5 (0%)	29	52

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	206	LYS
1	B	175	ALA
1	B	241	SER
1	D	175	ALA
1	C	241	SER

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	235/250 (94%)	218 (93%)	17 (7%)	14	29
1	B	235/250 (94%)	216 (92%)	19 (8%)	11	23
1	C	235/250 (94%)	219 (93%)	16 (7%)	16	32
1	D	235/250 (94%)	220 (94%)	15 (6%)	17	35
All	All	940/1000 (94%)	873 (93%)	67 (7%)	14	29

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	GLU
1	A	49	ARG
1	A	52	GLN
1	A	54	LEU
1	A	65	SER
1	A	111	ARG
1	A	126	ARG
1	A	142	GLN
1	A	188	PHE
1	A	197	GLU
1	A	222	ARG
1	A	231	LEU
1	A	241	SER
1	A	247	ILE
1	A	258	PHE
1	A	265	GLU
1	A	266	SER
1	B	17	GLU
1	B	25	THR
1	B	49	ARG
1	B	52	GLN
1	B	54	LEU
1	B	67	ARG
1	B	111	ARG
1	B	118	LYS
1	B	124	LYS
1	B	126	ARG
1	B	164	SER
1	B	182	TYR
1	B	186	ILE
1	B	188	PHE
1	B	194	LYS
1	B	199	ILE
1	B	217	LYS
1	B	231	LEU
1	B	247	ILE
1	C	9	ILE
1	C	10	ILE
1	C	17	GLU
1	C	52	GLN
1	C	54	LEU
1	C	111	ARG

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Mol	Chain	Res	Type
1	C	115	GLU
1	C	126	ARG
1	C	130	ILE
1	C	188	PHE
1	C	192	VAL
1	C	198	THR
1	C	231	LEU
1	C	247	ILE
1	C	259	VAL
1	C	262	PRO
1	D	17	GLU
1	D	52	GLN
1	D	54	LEU
1	D	111	ARG
1	D	115	GLU
1	D	118	LYS
1	D	125	GLU
1	D	126	ARG
1	D	142	GLN
1	D	164	SER
1	D	188	PHE
1	D	203	GLN
1	D	247	ILE
1	D	258	PHE
1	D	263	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	GLN
1	A	137	ASN
1	A	267	GLN
1	B	52	GLN
1	B	137	ASN
1	B	267	GLN
1	C	23	GLN
1	C	52	GLN
1	C	58	ASN
1	C	157	ASN
1	C	203	GLN
1	C	267	GLN
1	D	52	GLN

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Mol	Chain	Res	Type
1	D	137	ASN
1	D	142	GLN
1	D	193	GLN
1	D	267	GLN

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)

8 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	B	300	-	4,4,4	0.88	0	6,6,6	1.86	2 (33%)
2	PO4	A	302	-	4,4,4	0.85	0	6,6,6	1.30	1 (16%)
2	PO4	C	272	-	4,4,4	1.46	1 (25%)	6,6,6	2.01	2 (33%)
2	PO4	C	300	-	4,4,4	0.68	0	6,6,6	2.51	3 (50%)
2	PO4	B	302	-	4,4,4	0.64	0	6,6,6	1.49	2 (33%)
2	PO4	D	302	-	4,4,4	0.97	0	6,6,6	1.88	2 (33%)
2	PO4	B	272	-	4,4,4	0.84	0	6,6,6	1.36	2 (33%)
2	PO4	C	302	-	4,4,4	1.53	0	6,6,6	0.66	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	272	PO4	P-O2	-2.25	1.47	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	300	PO4	O4-P-O3	4.23	121.53	107.97
2	C	272	PO4	O2-P-O1	-3.87	96.74	110.89
2	C	300	PO4	O4-P-O1	-3.08	99.61	110.89
2	D	302	PO4	O3-P-O1	-2.97	100.04	110.89
2	B	300	PO4	O2-P-O1	-2.74	100.88	110.89
2	D	302	PO4	O2-P-O1	2.66	120.62	110.89
2	B	300	PO4	O4-P-O3	2.64	116.44	107.97
2	A	302	PO4	O4-P-O2	2.49	115.95	107.97
2	B	302	PO4	O4-P-O1	-2.48	101.81	110.89
2	C	272	PO4	O4-P-O1	2.43	119.80	110.89
2	C	300	PO4	O3-P-O2	-2.34	100.47	107.97
2	B	272	PO4	O3-P-O1	-2.26	102.61	110.89
2	B	272	PO4	O4-P-O3	2.07	114.62	107.97
2	B	302	PO4	O4-P-O2	2.00	114.40	107.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	300	PO4	5	0
2	C	272	PO4	2	0
2	B	302	PO4	2	0
2	C	302	PO4	2	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

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	266/292 (91%)	-0.11	3 (1%) 80 78	14, 24, 41, 56	0
1	B	266/292 (91%)	0.02	9 (3%) 45 38	14, 24, 41, 56	0
1	C	266/292 (91%)	0.02	14 (5%) 26 20	14, 24, 41, 56	0
1	D	266/292 (91%)	-0.07	1 (0%) 92 91	14, 24, 41, 55	0
All	All	1064/1168 (91%)	-0.03	27 (2%) 57 51	14, 24, 43, 56	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	125	GLU	3.9
1	A	197	GLU	3.3
1	C	200	GLU	3.0
1	B	198	THR	2.8
1	C	203	GLN	2.8
1	B	202	ILE	2.8
1	C	199	ILE	2.8
1	C	121	GLY	2.7
1	C	196	CYS	2.7
1	C	124	LYS	2.6
1	B	203	GLN	2.6
1	B	199	ILE	2.5
1	C	125	GLU	2.5
1	B	126	ARG	2.4
1	C	271	VAL	2.4
1	A	202	ILE	2.3
1	D	263	GLU	2.3
1	C	217	LYS	2.3
1	A	199	ILE	2.2
1	C	202	ILE	2.2
1	B	217	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	263	GLU	2.1
1	C	238	GLY	2.1
1	C	126	ARG	2.1
1	B	197	GLU	2.1
1	B	271	VAL	2.0
1	C	239	VAL	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)

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
2	PO4	B	300	5/5	0.89	0.19	32,37,42,43	0
2	PO4	B	302	5/5	0.96	0.20	16,24,25,27	0
2	PO4	D	302	5/5	0.96	0.16	4,9,15,18	0
2	PO4	A	302	5/5	0.97	0.18	15,22,23,24	0
2	PO4	C	300	5/5	0.97	0.15	14,15,19,21	0
2	PO4	C	272	5/5	0.97	0.12	22,22,22,22	0
2	PO4	B	272	5/5	0.97	0.15	27,27,29,29	0
2	PO4	C	302	5/5	0.99	0.13	2,5,9,9	0

6.5 Other polymers (i)

There are no such residues in this entry.