Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 02:55 PM EDT

PDB ID	:	2I2W
Title	:	Crystal Structure of Escherichia Coli Phosphoheptose Isomerase
Authors	:	DeLeon, G.; Blakely, K.; Zhang, K.; Wright, G.; Junop, M.
Deposited on	:	2006-08-17
Resolution	:	1.95 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

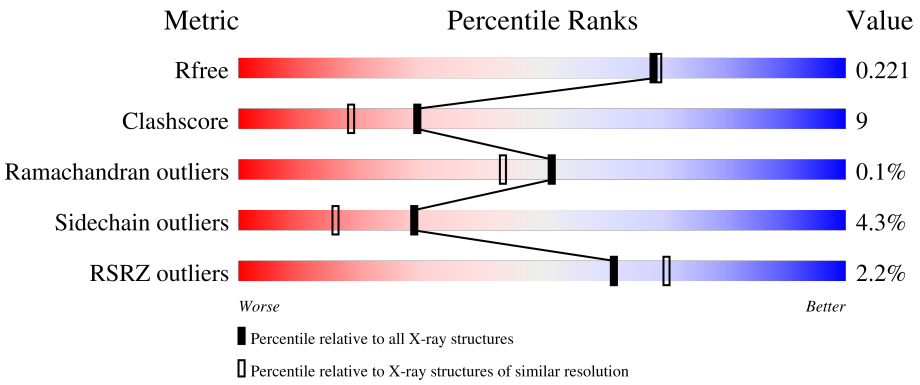
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	212	2% 62% 25% • • 9%
1	B	212	2% 58% 21% • • 17%
1	C	212	2% 65% 21% • 9%
1	D	212	% 62% 19% • 16%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6254 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 Phosphoheptose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	192	1461	914	261	278	8	0	0	0
			Total	C	N	O	S			
1	B	177	1351	846	243	255	7	0	0	0
			Total	C	N	O	S			
1	C	192	1461	914	261	278	8	0	0	0
			Total	C	N	O	S			
1	D	178	1357	849	244	257	7	0	0	0

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP P63224
A	-18	GLY	-	EXPRESSION TAG	UNP P63224
A	-17	SER	-	EXPRESSION TAG	UNP P63224
A	-16	SER	-	EXPRESSION TAG	UNP P63224
A	-15	HIS	-	EXPRESSION TAG	UNP P63224
A	-14	HIS	-	EXPRESSION TAG	UNP P63224
A	-13	HIS	-	EXPRESSION TAG	UNP P63224
A	-12	HIS	-	EXPRESSION TAG	UNP P63224
A	-11	HIS	-	EXPRESSION TAG	UNP P63224
A	-10	HIS	-	EXPRESSION TAG	UNP P63224
A	-9	SER	-	EXPRESSION TAG	UNP P63224
A	-8	SER	-	EXPRESSION TAG	UNP P63224
A	-7	GLY	-	EXPRESSION TAG	UNP P63224
A	-6	LEU	-	EXPRESSION TAG	UNP P63224
A	-5	VAL	-	EXPRESSION TAG	UNP P63224
A	-4	PRO	-	EXPRESSION TAG	UNP P63224
A	-3	ARG	-	EXPRESSION TAG	UNP P63224
A	-2	GLY	-	EXPRESSION TAG	UNP P63224
A	-1	SER	-	EXPRESSION TAG	UNP P63224
A	0	HIS	-	EXPRESSION TAG	UNP P63224
B	-19	MET	-	EXPRESSION TAG	UNP P63224

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	EXPRESSION TAG	UNP P63224
B	-17	SER	-	EXPRESSION TAG	UNP P63224
B	-16	SER	-	EXPRESSION TAG	UNP P63224
B	-15	HIS	-	EXPRESSION TAG	UNP P63224
B	-14	HIS	-	EXPRESSION TAG	UNP P63224
B	-13	HIS	-	EXPRESSION TAG	UNP P63224
B	-12	HIS	-	EXPRESSION TAG	UNP P63224
B	-11	HIS	-	EXPRESSION TAG	UNP P63224
B	-10	HIS	-	EXPRESSION TAG	UNP P63224
B	-9	SER	-	EXPRESSION TAG	UNP P63224
B	-8	SER	-	EXPRESSION TAG	UNP P63224
B	-7	GLY	-	EXPRESSION TAG	UNP P63224
B	-6	LEU	-	EXPRESSION TAG	UNP P63224
B	-5	VAL	-	EXPRESSION TAG	UNP P63224
B	-4	PRO	-	EXPRESSION TAG	UNP P63224
B	-3	ARG	-	EXPRESSION TAG	UNP P63224
B	-2	GLY	-	EXPRESSION TAG	UNP P63224
B	-1	SER	-	EXPRESSION TAG	UNP P63224
B	0	HIS	-	EXPRESSION TAG	UNP P63224
C	-19	MET	-	EXPRESSION TAG	UNP P63224
C	-18	GLY	-	EXPRESSION TAG	UNP P63224
C	-17	SER	-	EXPRESSION TAG	UNP P63224
C	-16	SER	-	EXPRESSION TAG	UNP P63224
C	-15	HIS	-	EXPRESSION TAG	UNP P63224
C	-14	HIS	-	EXPRESSION TAG	UNP P63224
C	-13	HIS	-	EXPRESSION TAG	UNP P63224
C	-12	HIS	-	EXPRESSION TAG	UNP P63224
C	-11	HIS	-	EXPRESSION TAG	UNP P63224
C	-10	HIS	-	EXPRESSION TAG	UNP P63224
C	-9	SER	-	EXPRESSION TAG	UNP P63224
C	-8	SER	-	EXPRESSION TAG	UNP P63224
C	-7	GLY	-	EXPRESSION TAG	UNP P63224
C	-6	LEU	-	EXPRESSION TAG	UNP P63224
C	-5	VAL	-	EXPRESSION TAG	UNP P63224
C	-4	PRO	-	EXPRESSION TAG	UNP P63224
C	-3	ARG	-	EXPRESSION TAG	UNP P63224
C	-2	GLY	-	EXPRESSION TAG	UNP P63224
C	-1	SER	-	EXPRESSION TAG	UNP P63224
C	0	HIS	-	EXPRESSION TAG	UNP P63224
D	-19	MET	-	EXPRESSION TAG	UNP P63224
D	-18	GLY	-	EXPRESSION TAG	UNP P63224
D	-17	SER	-	EXPRESSION TAG	UNP P63224

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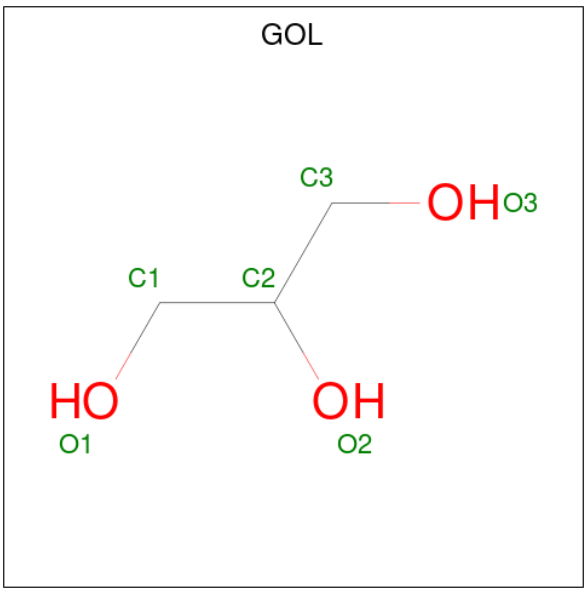
Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	EXPRESSION TAG	UNP P63224
D	-15	HIS	-	EXPRESSION TAG	UNP P63224
D	-14	HIS	-	EXPRESSION TAG	UNP P63224
D	-13	HIS	-	EXPRESSION TAG	UNP P63224
D	-12	HIS	-	EXPRESSION TAG	UNP P63224
D	-11	HIS	-	EXPRESSION TAG	UNP P63224
D	-10	HIS	-	EXPRESSION TAG	UNP P63224
D	-9	SER	-	EXPRESSION TAG	UNP P63224
D	-8	SER	-	EXPRESSION TAG	UNP P63224
D	-7	GLY	-	EXPRESSION TAG	UNP P63224
D	-6	LEU	-	EXPRESSION TAG	UNP P63224
D	-5	VAL	-	EXPRESSION TAG	UNP P63224
D	-4	PRO	-	EXPRESSION TAG	UNP P63224
D	-3	ARG	-	EXPRESSION TAG	UNP P63224
D	-2	GLY	-	EXPRESSION TAG	UNP P63224
D	-1	SER	-	EXPRESSION TAG	UNP P63224
D	0	HIS	-	EXPRESSION TAG	UNP P63224

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	D	1	6	3	3	0	0

- Molecule 3 is water.

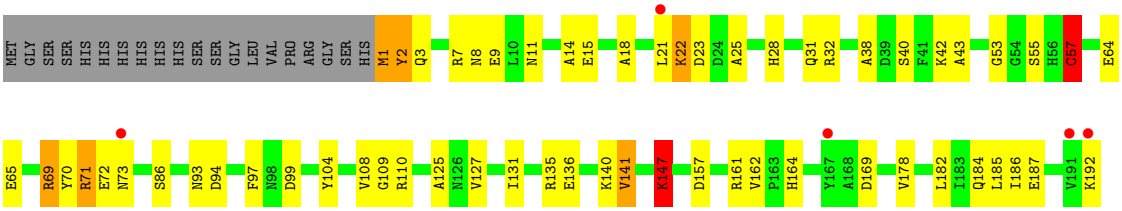
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	167	167	167	0	0
			Total	O		
3	B	142	142	142	0	0
			Total	O		
3	C	150	150	150	0	0
			Total	O		
3	D	159	159	159	0	0

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.

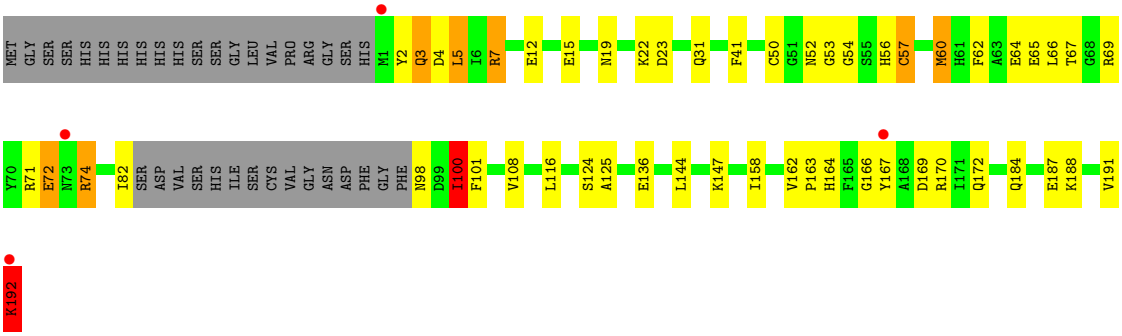
• Molecule 1: Phosphoheptose isomerase

Chain A:



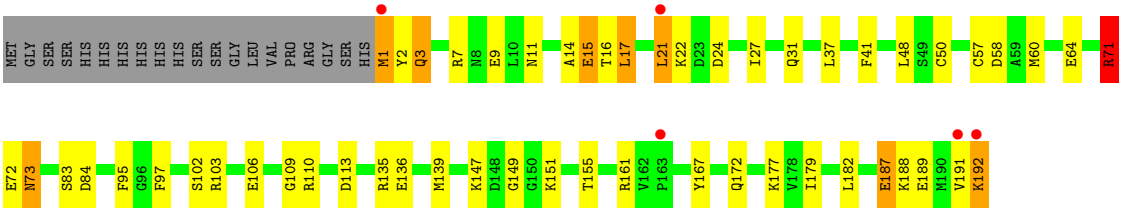
• Molecule 1: Phosphoheptose isomerase

Chain B:



• Molecule 1: Phosphoheptose isomerase

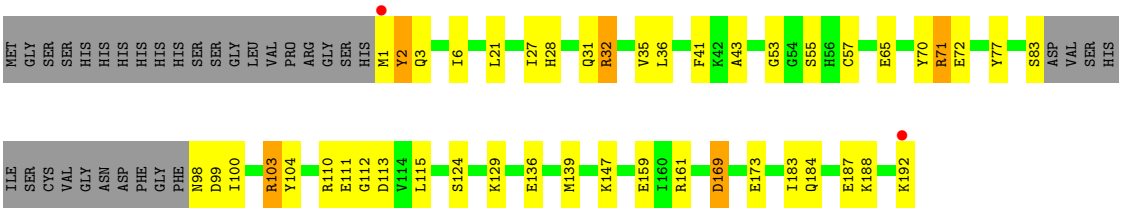
Chain C:



• Molecule 1: Phosphoheptose isomerase

Chain D:

%
62%
19%
16%


MET GLY SER SER HIS HIS HIS HIS HIS HIS SER SER GLY LEU VAL PRO ARG GLY SER HIS M1 Y2 Q3 I6 L21 I27 H28 Q31 R32 V35 L36 F41 K42 A43 G53 G54 S55 H56 C57 E65 Y70 R71 E72 Y77 S83 ASP VAL SER HIS
ILE SER CYS VAL GLY ASN ASP PHE GLY PHE N98 D99 I100 R103 Y104 R110 E111 G112 D113 V114 L115 S124 K129 E136 M139 K147 E159 I160 R161 D169 E173 I183 Q184 E187 K188 K192

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	83.93Å 89.61Å 106.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.90 – 1.95 45.91 – 1.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.90-1.95) 99.6 (45.91-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.65 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.170 , 0.219 0.174 , 0.221	Depositor DCC
R_{free} test set	2996 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	28.5	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 51.5	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	0.97	EDS
Total number of atoms	6254	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.42% 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 (i)

5.1 Standard geometry (i)

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

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	1.86	23/1482 (1.6%)	1.47	17/1993 (0.9%)
1	B	1.82	23/1368 (1.7%)	1.48	16/1837 (0.9%)
1	C	1.83	25/1482 (1.7%)	1.50	17/1993 (0.9%)
1	D	1.82	22/1374 (1.6%)	1.49	11/1845 (0.6%)
All	All	1.84	93/5706 (1.6%)	1.49	61/7668 (0.8%)

All (93) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	25	ALA	CA-CB	9.46	1.72	1.52
1	B	72	GLU	CG-CD	9.45	1.66	1.51
1	D	136	GLU	CD-OE1	9.21	1.35	1.25
1	C	9	GLU	CD-OE2	9.06	1.35	1.25
1	C	149	GLY	N-CA	-8.84	1.32	1.46
1	B	64	GLU	CG-CD	8.59	1.64	1.51
1	C	187	GLU	CG-CD	8.14	1.64	1.51
1	B	125	ALA	CA-CB	8.10	1.69	1.52
1	C	3	GLN	CG-CD	7.87	1.69	1.51
1	D	71	ARG	CD-NE	-7.80	1.33	1.46
1	D	55	SER	CB-OG	7.66	1.52	1.42
1	A	18	ALA	CA-CB	7.45	1.68	1.52
1	A	192	LYS	CB-CG	-7.29	1.32	1.52
1	D	124	SER	CB-OG	7.24	1.51	1.42
1	B	54	GLY	C-O	7.21	1.35	1.23
1	C	189	GLU	CB-CG	-7.01	1.38	1.52
1	D	65	GLU	CD-OE1	6.99	1.33	1.25
1	C	136	GLU	CB-CG	-6.85	1.39	1.52
1	D	70	TYR	CG-CD2	6.73	1.48	1.39
1	C	64	GLU	CB-CG	6.71	1.64	1.52
1	B	167	TYR	CE1-CZ	6.68	1.47	1.38
1	C	97	PHE	CE2-CZ	6.59	1.49	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	86	SER	CB-OG	6.48	1.50	1.42
1	C	151	LYS	CE-NZ	6.47	1.65	1.49
1	B	184	GLN	CD-NE2	6.46	1.49	1.32
1	A	161	ARG	CZ-NH1	-6.39	1.24	1.33
1	D	65	GLU	CG-CD	6.37	1.61	1.51
1	A	15	GLU	CD-OE1	6.35	1.32	1.25
1	A	72	GLU	CD-OE2	6.34	1.32	1.25
1	A	104	TYR	CD1-CE1	6.34	1.48	1.39
1	D	43	ALA	CA-CB	6.33	1.65	1.52
1	C	136	GLU	CD-OE2	6.33	1.32	1.25
1	B	167	TYR	CD1-CE1	6.27	1.48	1.39
1	B	15	GLU	CG-CD	6.16	1.61	1.51
1	D	83	SER	C-O	6.15	1.35	1.23
1	B	65	GLU	CG-CD	6.15	1.61	1.51
1	C	136	GLU	CD-OE1	6.14	1.32	1.25
1	C	41	PHE	CE1-CZ	6.12	1.49	1.37
1	C	83	SER	CA-CB	6.08	1.62	1.52
1	C	167	TYR	CE1-CZ	6.06	1.46	1.38
1	B	124	SER	CB-OG	6.03	1.50	1.42
1	A	109	GLY	N-CA	6.00	1.55	1.46
1	A	15	GLU	CG-CD	5.95	1.60	1.51
1	B	172	GLN	CD-NE2	5.94	1.47	1.32
1	B	101	PHE	CD2-CE2	5.92	1.51	1.39
1	D	188	LYS	CD-CE	5.84	1.65	1.51
1	A	40	SER	CB-OG	5.83	1.49	1.42
1	A	127	VAL	CA-CB	5.77	1.66	1.54
1	A	72	GLU	CG-CD	5.74	1.60	1.51
1	A	14	ALA	CA-CB	5.72	1.64	1.52
1	D	104	TYR	CE2-CZ	-5.68	1.31	1.38
1	A	43	ALA	CA-CB	5.67	1.64	1.52
1	A	104	TYR	CD2-CE2	5.64	1.47	1.39
1	D	77	TYR	CE2-CZ	5.62	1.45	1.38
1	C	109	GLY	CA-C	5.60	1.60	1.51
1	B	158	ILE	CA-CB	5.58	1.67	1.54
1	D	35	VAL	CB-CG1	5.56	1.64	1.52
1	C	2	TYR	CE1-CZ	5.56	1.45	1.38
1	D	53	GLY	C-O	5.53	1.32	1.23
1	B	12	GLU	CG-CD	5.52	1.60	1.51
1	D	159	GLU	CB-CG	5.52	1.62	1.52
1	D	77	TYR	CD1-CE1	5.50	1.47	1.39
1	C	187	GLU	CD-OE1	5.45	1.31	1.25
1	D	147	LYS	CE-NZ	5.43	1.62	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	141	VAL	CB-CG2	5.40	1.64	1.52
1	B	187	GLU	CG-CD	5.40	1.60	1.51
1	A	53	GLY	CA-C	5.39	1.60	1.51
1	D	41	PHE	CE2-CZ	5.38	1.47	1.37
1	C	14	ALA	CA-CB	5.37	1.63	1.52
1	C	95	PHE	CE2-CZ	5.34	1.47	1.37
1	C	95	PHE	CG-CD2	5.33	1.46	1.38
1	A	162	VAL	CB-CG2	5.31	1.64	1.52
1	A	55	SER	CB-OG	5.29	1.49	1.42
1	B	62	PHE	CE2-CZ	5.25	1.47	1.37
1	B	192	LYS	C-OXT	5.21	1.33	1.23
1	C	15	GLU	CG-CD	5.21	1.59	1.51
1	B	41	PHE	CG-CD2	5.21	1.46	1.38
1	D	72	GLU	CG-CD	5.19	1.59	1.51
1	B	167	TYR	CG-CD1	5.19	1.45	1.39
1	D	173	GLU	CD-OE1	5.19	1.31	1.25
1	A	64	GLU	CG-CD	5.18	1.59	1.51
1	B	53	GLY	N-CA	-5.17	1.38	1.46
1	B	187	GLU	CB-CG	-5.14	1.42	1.52
1	C	7	ARG	CG-CD	5.13	1.64	1.51
1	B	169	ASP	CG-OD1	5.11	1.37	1.25
1	A	140	LYS	CE-NZ	5.11	1.61	1.49
1	C	9	GLU	CG-CD	5.10	1.59	1.51
1	B	167	TYR	CD2-CE2	5.06	1.47	1.39
1	C	103	ARG	CZ-NH2	-5.06	1.26	1.33
1	D	111	GLU	CD-OE2	-5.04	1.20	1.25
1	A	127	VAL	C-O	5.02	1.32	1.23
1	D	112	GLY	CA-C	5.02	1.59	1.51
1	C	172	GLN	CG-CD	5.00	1.62	1.51

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	110	ARG	NE-CZ-NH1	-12.39	114.11	120.30
1	A	157	ASP	CB-CG-OD1	-12.18	107.34	118.30
1	D	71	ARG	NE-CZ-NH2	-11.66	114.47	120.30
1	A	157	ASP	CB-CG-OD2	11.47	128.63	118.30
1	D	71	ARG	NE-CZ-NH1	11.10	125.85	120.30
1	C	71	ARG	NE-CZ-NH1	-10.51	115.05	120.30
1	D	110	ARG	NE-CZ-NH1	-9.90	115.35	120.30
1	D	103	ARG	NE-CZ-NH2	-9.82	115.39	120.30
1	B	23	ASP	CB-CG-OD1	8.92	126.33	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	161	ARG	NE-CZ-NH1	8.26	124.43	120.30
1	C	103	ARG	NE-CZ-NH2	-7.91	116.34	120.30
1	C	110	ARG	NE-CZ-NH1	-7.83	116.38	120.30
1	D	99	ASP	CB-CG-OD1	7.82	125.33	118.30
1	C	7	ARG	NE-CZ-NH2	-7.68	116.46	120.30
1	C	37	LEU	CB-CG-CD2	-7.48	98.29	111.00
1	D	104	TYR	CB-CG-CD2	7.32	125.39	121.00
1	A	1	MET	CG-SD-CE	7.17	111.67	100.20
1	B	72	GLU	OE1-CD-OE2	-6.78	115.16	123.30
1	C	110	ARG	NE-CZ-NH2	6.75	123.68	120.30
1	A	69	ARG	NE-CZ-NH1	6.68	123.64	120.30
1	A	21	LEU	CA-CB-CG	6.57	130.41	115.30
1	A	178	VAL	CA-CB-CG2	-6.52	101.12	110.90
1	B	23	ASP	CB-CG-OD2	-6.41	112.53	118.30
1	C	182	LEU	CB-CG-CD1	-6.23	100.41	111.00
1	B	65	GLU	OE1-CD-OE2	-6.19	115.87	123.30
1	B	60	MET	CA-CB-CG	-6.12	102.90	113.30
1	B	144	LEU	CB-CG-CD2	6.10	121.38	111.00
1	B	4	ASP	CB-CG-OD2	-6.08	112.83	118.30
1	C	147	LYS	CD-CE-NZ	-5.93	98.06	111.70
1	C	48	LEU	CB-CG-CD1	-5.89	100.99	111.00
1	A	69	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	B	64	GLU	OE1-CD-OE2	-5.80	116.34	123.30
1	C	136	GLU	OE1-CD-OE2	5.78	130.23	123.30
1	B	12	GLU	OE1-CD-OE2	-5.74	116.42	123.30
1	C	155	THR	CA-CB-CG2	-5.66	104.48	112.40
1	A	140	LYS	CD-CE-NZ	-5.63	98.75	111.70
1	B	69	ARG	NE-CZ-NH2	5.57	123.08	120.30
1	A	104	TYR	CB-CG-CD1	-5.53	117.69	121.00
1	A	161	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	B	7	ARG	CG-CD-NE	-5.47	100.31	111.80
1	D	115	LEU	CB-CG-CD1	-5.46	101.73	111.00
1	B	170	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	C	135	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	C	17	LEU	CB-CG-CD1	-5.41	101.81	111.00
1	C	16	THR	CA-CB-CG2	-5.36	104.90	112.40
1	A	9	GLU	OE1-CD-OE2	-5.35	116.88	123.30
1	B	5	LEU	CB-CG-CD1	5.33	120.06	111.00
1	B	100	ILE	CB-CA-C	-5.30	101.01	111.60
1	C	161	ARG	NE-CZ-NH1	-5.26	117.67	120.30
1	A	57	CYS	CA-CB-SG	-5.17	104.69	114.00
1	B	74	ARG	NE-CZ-NH1	-5.15	117.72	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	36	LEU	CB-CG-CD2	-5.15	102.25	111.00
1	C	21	LEU	CB-CG-CD1	-5.12	102.30	111.00
1	D	32	ARG	NE-CZ-NH1	5.11	122.85	120.30
1	C	71	ARG	NH1-CZ-NH2	5.09	125.00	119.40
1	A	161	ARG	NE-CZ-NH2	5.08	122.84	120.30
1	A	147	LYS	CD-CE-NZ	-5.07	100.03	111.70
1	A	2	TYR	N-CA-CB	-5.07	101.47	110.60
1	A	94	ASP	CB-CG-OD1	5.06	122.86	118.30
1	B	116	LEU	CB-CG-CD1	-5.05	102.41	111.00
1	D	169	ASP	CB-CG-OD1	-5.01	113.79	118.30

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	1461	0	1460	38	0
1	B	1351	0	1366	30	0
1	C	1461	0	1460	28	0
1	D	1357	0	1371	30	0
2	D	6	0	8	0	0
3	A	167	0	0	16	0
3	B	142	0	0	10	0
3	C	150	0	0	1	0
3	D	159	0	0	9	0
All	All	6254	0	5665	106	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 (106) 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:82:ILE:CD1	1:B:82:ILE:CG1	1.79	1.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:249:HOH:O	1:D:184:GLN:HG3	1.43	1.15
1:D:98:ASN:HB2	3:D:301:HOH:O	1.46	1.13
1:B:98:ASN:HB2	3:B:324:HOH:O	1.46	1.12
1:B:100:ILE:HG12	3:B:322:HOH:O	1.60	1.01
1:C:73:ASN:HD22	1:C:73:ASN:N	1.66	0.92
1:D:169:ASP:HB2	3:D:260:HOH:O	1.79	0.82
1:B:3:GLN:HE21	1:B:3:GLN:HA	1.44	0.82
1:B:3:GLN:NE2	1:C:27:ILE:HG22	1.95	0.82
1:A:11:ASN:HB3	3:A:275:HOH:O	1.81	0.81
1:A:169:ASP:HB2	3:A:220:HOH:O	1.81	0.80
1:B:3:GLN:HA	1:B:3:GLN:NE2	1.95	0.80
1:D:183:ILE:HG21	3:D:338:HOH:O	1.82	0.78
1:B:56:HIS:ND1	3:B:326:HOH:O	2.22	0.73
1:A:184:GLN:HG2	3:A:290:HOH:O	1.91	0.70
1:C:187:GLU:OE2	1:C:187:GLU:O	2.09	0.70
1:C:73:ASN:HD22	1:C:73:ASN:H	1.39	0.70
1:C:73:ASN:N	1:C:73:ASN:ND2	2.37	0.69
1:D:1:MET:O	1:D:2:TYR:HB2	1.91	0.69
1:C:192:LYS:OXT	1:C:192:LYS:HG2	1.94	0.66
1:B:74:ARG:HD3	3:B:305:HOH:O	1.96	0.66
1:A:7:ARG:NH2	3:A:315:HOH:O	2.24	0.64
1:A:131:ILE:HG23	3:A:260:HOH:O	1.97	0.64
1:B:56:HIS:CE1	3:B:326:HOH:O	2.52	0.63
1:B:163:PRO:HG3	3:B:271:HOH:O	1.99	0.62
1:D:1:MET:O	1:D:2:TYR:CB	2.47	0.62
1:C:188:LYS:O	1:C:191:VAL:HG22	2.00	0.61
1:D:71:ARG:HD3	3:D:349:HOH:O	2.00	0.60
1:D:28:HIS:CE1	1:D:32:ARG:HD3	2.38	0.59
1:A:28:HIS:HB3	1:A:32:ARG:HH21	1.67	0.59
1:A:108:VAL:HG12	1:B:108:VAL:HB	1.85	0.59
1:B:3:GLN:HE21	1:C:27:ILE:HG22	1.66	0.59
1:A:141:VAL:HG21	3:A:260:HOH:O	2.03	0.58
1:D:192:LYS:HD2	3:D:341:HOH:O	2.03	0.58
1:C:71:ARG:HG2	1:D:103:ARG:CZ	2.34	0.57
1:A:97:PHE:HA	1:A:125:ALA:HB3	1.85	0.57
1:C:1:MET:SD	1:C:3:GLN:HB2	2.46	0.56
1:A:57:CYS:SG	1:D:57:CYS:SG	3.04	0.55
1:B:7:ARG:NH2	1:C:24:ASP:OD2	2.40	0.54
1:B:82:ILE:CD1	1:B:82:ILE:CB	2.82	0.54
1:B:3:GLN:NE2	1:B:3:GLN:CA	2.70	0.54
1:A:32:ARG:HD2	3:A:248:HOH:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:HIS:O	1:D:32:ARG:HG3	2.08	0.54
1:C:72:GLU:HG2	1:C:73:ASN:ND2	2.23	0.54
1:A:28:HIS:HB3	1:A:32:ARG:NH2	2.23	0.53
1:B:57:CYS:SG	1:C:60:MET:HE2	2.48	0.53
1:A:57:CYS:HG	1:D:57:CYS:CB	2.21	0.53
1:A:1:MET:HA	1:D:31:GLN:HE22	1.76	0.51
1:A:22:LYS:HG2	1:A:23:ASP:N	2.24	0.51
1:A:71:ARG:NH1	1:A:71:ARG:HG3	2.26	0.51
1:C:58:ASP:HB3	1:C:179:ILE:CD1	2.40	0.50
1:A:182:LEU:O	1:A:186:ILE:HG13	2.10	0.50
1:A:1:MET:SD	1:A:3:GLN:NE2	2.84	0.50
1:C:102:SER:O	1:C:106:GLU:HG3	2.13	0.49
1:B:19:ASN:HA	1:B:22:LYS:HD2	1.94	0.49
1:B:188:LYS:O	1:B:192:LYS:HB2	2.12	0.49
1:D:187:GLU:OE2	1:D:187:GLU:HA	2.12	0.49
1:A:2:TYR:OH	1:D:192:LYS:HE3	2.12	0.49
1:A:135:ARG:CZ	3:A:260:HOH:O	2.61	0.49
1:B:67:THR:O	1:B:71:ARG:HB2	2.12	0.49
1:A:1:MET:SD	1:A:3:GLN:HG2	2.54	0.48
1:C:187:GLU:OE2	1:C:191:VAL:HG13	2.14	0.48
1:B:57:CYS:CB	1:C:57:CYS:HG	2.24	0.48
1:A:2:TYR:HA	3:A:206:HOH:O	2.14	0.47
1:B:147:LYS:NZ	3:B:277:HOH:O	2.36	0.47
1:D:1:MET:HA	3:D:288:HOH:O	2.14	0.47
1:B:52:ASN:HB2	3:B:316:HOH:O	2.14	0.47
1:D:113:ASP:HB2	1:D:139:MET:HG2	1.97	0.47
1:C:11:ASN:O	1:C:15:GLU:HG3	2.15	0.46
1:A:38:ALA:O	1:A:42:LYS:HG3	2.15	0.46
1:B:98:ASN:N	3:B:322:HOH:O	2.48	0.46
1:A:3:GLN:OE1	1:D:27:ILE:HG22	2.15	0.46
1:A:135:ARG:NE	3:A:260:HOH:O	2.47	0.46
1:D:183:ILE:HD13	3:D:338:HOH:O	2.15	0.45
1:A:147:LYS:NZ	3:A:338:HOH:O	2.48	0.45
1:C:113:ASP:HB2	1:C:139:MET:HG2	1.99	0.45
1:A:2:TYR:OH	1:D:192:LYS:CE	2.65	0.45
1:A:164:HIS:HD2	3:A:340:HOH:O	1.98	0.45
1:A:3:GLN:HE22	1:D:28:HIS:HD2	1.65	0.45
1:A:65:GLU:O	1:A:69:ARG:HG3	2.17	0.45
1:A:31:GLN:NE2	1:D:3:GLN:HG3	2.32	0.44
1:C:192:LYS:OXT	1:C:192:LYS:CG	2.62	0.44
1:B:2:TYR:H	1:C:31:GLN:NE2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:CYS:HA	3:C:196:HOH:O	2.17	0.44
1:D:1:MET:O	1:D:1:MET:HG2	2.17	0.44
1:B:57:CYS:SG	1:C:60:MET:CE	3.06	0.43
1:B:66:LEU:HD23	1:B:66:LEU:HA	1.84	0.43
1:A:70:TYR:OH	1:A:187:GLU:HA	2.19	0.43
1:B:164:HIS:CD2	1:B:166:GLY:H	2.37	0.43
1:A:93:ASN:ND2	3:A:231:HOH:O	2.49	0.43
1:B:162:VAL:HA	1:B:163:PRO:HD3	1.87	0.42
1:D:1:MET:O	1:D:2:TYR:CG	2.72	0.42
1:A:8:ASN:HB3	3:A:262:HOH:O	2.20	0.42
1:C:1:MET:HE2	1:C:1:MET:HB3	1.73	0.42
1:D:1:MET:O	1:D:1:MET:CG	2.67	0.42
1:D:1:MET:HB3	3:D:321:HOH:O	2.18	0.42
1:C:1:MET:C	1:C:1:MET:HE3	2.40	0.41
1:B:31:GLN:HE22	1:C:1:MET:HG2	1.85	0.41
1:D:192:LYS:NZ	3:D:343:HOH:O	2.45	0.41
1:C:187:GLU:OE2	1:C:187:GLU:C	2.58	0.41
1:A:71:ARG:HG3	1:A:71:ARG:HH11	1.85	0.41
1:A:185:LEU:HD11	1:D:6:ILE:HD11	2.03	0.41
1:C:17:LEU:HD22	1:C:177:LYS:HE2	2.03	0.41
1:A:99:ASP:HB2	3:A:324:HOH:O	2.20	0.41
1:A:31:GLN:HE22	1:D:3:GLN:HG3	1.85	0.41
1:B:50:CYS:HA	3:B:326:HOH:O	2.22	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	190/212 (90%)	184 (97%)	6 (3%)	0	100 100
1	B	173/212 (82%)	171 (99%)	2 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	190/212 (90%)	189 (100%)	1 (0%)	0	100	100
1	D	174/212 (82%)	168 (97%)	5 (3%)	1 (1%)	25	14
All	All	727/848 (86%)	712 (98%)	14 (2%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	2	TYR

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	151/168 (90%)	145 (96%)	6 (4%)	31	19
1	B	138/168 (82%)	129 (94%)	9 (6%)	17	6
1	C	151/168 (90%)	144 (95%)	7 (5%)	27	14
1	D	139/168 (83%)	136 (98%)	3 (2%)	52	44
All	All	579/672 (86%)	554 (96%)	25 (4%)	29	16

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	57	CYS
1	A	71	ARG
1	A	73	ASN
1	A	136	GLU
1	A	147	LYS
1	B	3	GLN
1	B	5	LEU
1	B	57	CYS
1	B	60	MET
1	B	72	GLU

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Mol	Chain	Res	Type
1	B	100	ILE
1	B	136	GLU
1	B	191	VAL
1	B	192	LYS
1	C	1	MET
1	C	21	LEU
1	C	22	LYS
1	C	71	ARG
1	C	73	ASN
1	C	84	ASP
1	C	192	LYS
1	D	21	LEU
1	D	100	ILE
1	D	129	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	26	ASN
1	A	31	GLN
1	A	93	ASN
1	B	3	GLN
1	B	11	ASN
1	B	26	ASN
1	B	31	GLN
1	B	56	HIS
1	C	11	ASN
1	C	31	GLN
1	C	73	ASN
1	C	164	HIS
1	D	26	ASN
1	D	28	HIS
1	D	164	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is 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	GOL	D	196	-	5,5,5	0.96	0	5,5,5	2.66	3 (60%)

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	D	196	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	196	GOL	O1-C1-C2	-3.87	92.96	110.38
2	D	196	GOL	O2-C2-C1	3.55	123.86	109.18
2	D	196	GOL	O2-C2-C3	2.67	120.24	109.18

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	196	GOL	O2-C2-C3-O3
2	D	196	GOL	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	192/212 (90%)	-0.24	5 (2%)	56	65	26, 38, 62, 102	0
1	B	177/212 (83%)	-0.25	4 (2%)	60	69	26, 36, 62, 92	0
1	C	192/212 (90%)	-0.12	5 (2%)	56	65	24, 34, 63, 96	0
1	D	178/212 (83%)	-0.22	2 (1%)	80	85	25, 33, 55, 83	0
All	All	739/848 (87%)	-0.21	16 (2%)	62	70	24, 35, 62, 102	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1	MET	6.2
1	C	191	VAL	4.3
1	A	191	VAL	4.1
1	C	1	MET	3.4
1	B	167	TYR	3.2
1	D	192	LYS	2.7
1	A	73	ASN	2.7
1	A	21	LEU	2.6
1	C	192	LYS	2.6
1	A	192	LYS	2.6
1	C	21	LEU	2.5
1	A	167	TYR	2.5
1	C	163	PRO	2.3
1	B	73	ASN	2.2
1	B	1	MET	2.2
1	B	192	LYS	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, 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
2	GOL	D	196	6/6	0.91	0.14	44,51,53,63	0

6.5 Other polymers (i)

There are no such residues in this entry.