Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2023 – 04:27 PM EDT

PDB ID	:	6XGZ
Title	:	Crystal structure of E. coli MlaFB ABC transport subunits in the monomeric state
Authors	:	Chang, Y.; Bhabha, G.; Ekiert, D.C.
Deposited on	:	2020-06-18
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:

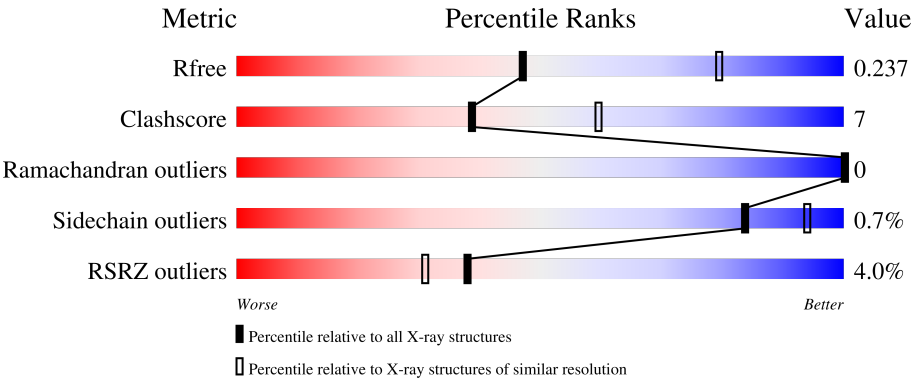
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 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	269	2% 85% 13% ·
1	C	269	2% 79% 17% ·
1	E	269	10% 75% 22% ·
1	G	269	5% 75% 23% ·
2	B	109	3% 82% 12% 6%

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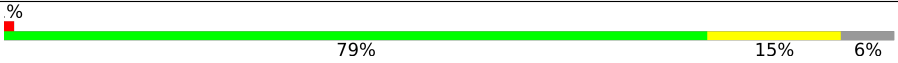
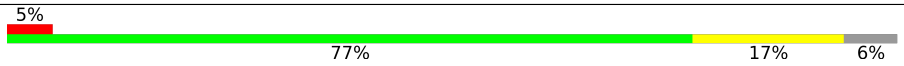
Mol	Chain	Length	Quality of chain
2	D	109	
2	F	109	
2	H	109	

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
3	PO4	E	301	-	-	X	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 11433 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 Organic solvent ABC transporter ATP-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	1	0
			2010	1276	356	365	13			
1	C	260	Total	C	N	O	S	0	3	0
			1999	1269	356	360	14			
1	E	259	Total	C	N	O	S	0	2	0
			1977	1252	354	358	13			
1	G	265	Total	C	N	O	S	0	3	0
			2024	1287	358	366	13			

- Molecule 2 is a protein called ABC transporter maintaining OM lipid asymmetry, cytoplasmic STAS component.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	102	Total	C	N	O	S	0	5	0
			850	540	144	163	3			
2	D	102	Total	C	N	O	S	0	4	0
			835	532	140	160	3			
2	F	102	Total	C	N	O	S	0	3	0
			827	528	138	158	3			
2	H	102	Total	C	N	O	S	0	4	0
			839	534	140	162	3			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	MET	-	expression tag	UNP W8T4U6
B	-10	HIS	-	expression tag	UNP W8T4U6
B	-9	HIS	-	expression tag	UNP W8T4U6
B	-8	HIS	-	expression tag	UNP W8T4U6
B	-7	HIS	-	expression tag	UNP W8T4U6
B	-6	HIS	-	expression tag	UNP W8T4U6
B	-5	HIS	-	expression tag	UNP W8T4U6
B	-4	GLU	-	expression tag	UNP W8T4U6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	ASN	-	expression tag	UNP W8T4U6
B	-2	LEU	-	expression tag	UNP W8T4U6
B	-1	TYR	-	expression tag	UNP W8T4U6
B	0	PHE	-	expression tag	UNP W8T4U6
B	1	GLN	-	expression tag	UNP W8T4U6
D	-11	MET	-	expression tag	UNP W8T4U6
D	-10	HIS	-	expression tag	UNP W8T4U6
D	-9	HIS	-	expression tag	UNP W8T4U6
D	-8	HIS	-	expression tag	UNP W8T4U6
D	-7	HIS	-	expression tag	UNP W8T4U6
D	-6	HIS	-	expression tag	UNP W8T4U6
D	-5	HIS	-	expression tag	UNP W8T4U6
D	-4	GLU	-	expression tag	UNP W8T4U6
D	-3	ASN	-	expression tag	UNP W8T4U6
D	-2	LEU	-	expression tag	UNP W8T4U6
D	-1	TYR	-	expression tag	UNP W8T4U6
D	0	PHE	-	expression tag	UNP W8T4U6
D	1	GLN	-	expression tag	UNP W8T4U6
F	-11	MET	-	expression tag	UNP W8T4U6
F	-10	HIS	-	expression tag	UNP W8T4U6
F	-9	HIS	-	expression tag	UNP W8T4U6
F	-8	HIS	-	expression tag	UNP W8T4U6
F	-7	HIS	-	expression tag	UNP W8T4U6
F	-6	HIS	-	expression tag	UNP W8T4U6
F	-5	HIS	-	expression tag	UNP W8T4U6
F	-4	GLU	-	expression tag	UNP W8T4U6
F	-3	ASN	-	expression tag	UNP W8T4U6
F	-2	LEU	-	expression tag	UNP W8T4U6
F	-1	TYR	-	expression tag	UNP W8T4U6
F	0	PHE	-	expression tag	UNP W8T4U6
F	1	GLN	-	expression tag	UNP W8T4U6
H	-11	MET	-	expression tag	UNP W8T4U6
H	-10	HIS	-	expression tag	UNP W8T4U6
H	-9	HIS	-	expression tag	UNP W8T4U6
H	-8	HIS	-	expression tag	UNP W8T4U6
H	-7	HIS	-	expression tag	UNP W8T4U6
H	-6	HIS	-	expression tag	UNP W8T4U6
H	-5	HIS	-	expression tag	UNP W8T4U6
H	-4	GLU	-	expression tag	UNP W8T4U6
H	-3	ASN	-	expression tag	UNP W8T4U6
H	-2	LEU	-	expression tag	UNP W8T4U6
H	-1	TYR	-	expression tag	UNP W8T4U6

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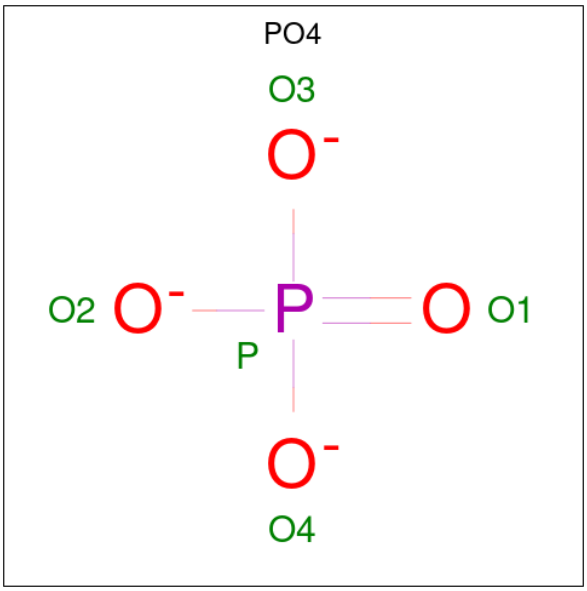
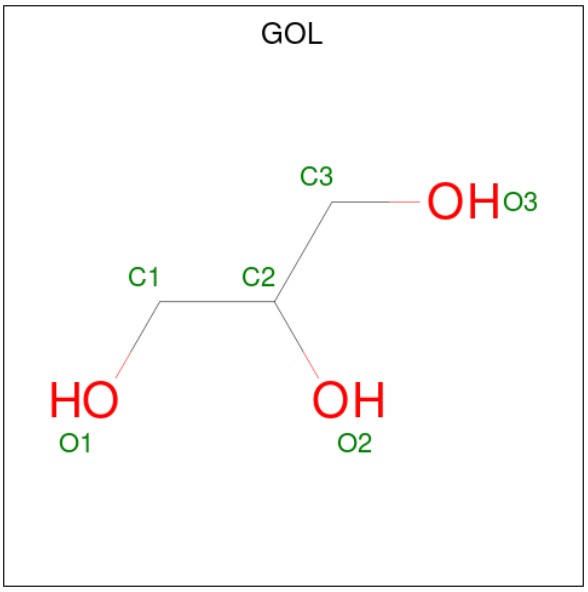
Chain	Residue	Modelled	Actual	Comment	Reference
H	0	PHE	-	expression tag	UNP W8T4U6
H	1	GLN	-	expression tag	UNP W8T4U6

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



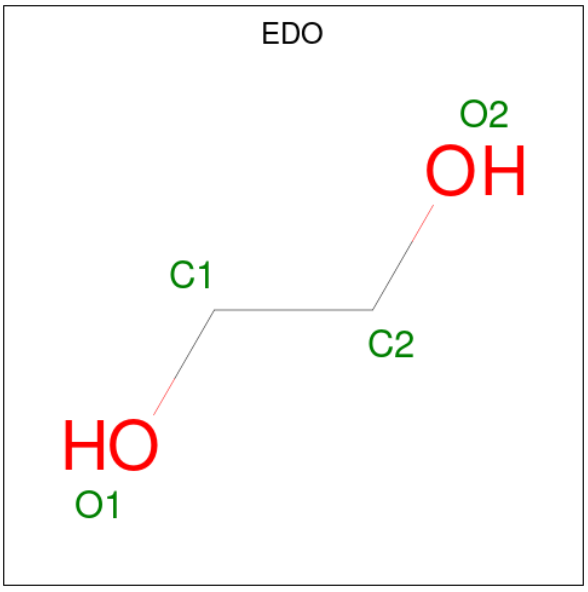
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	A	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0
3	E	1	Total O P 5 4 1	0	0
3	E	1	Total O P 5 4 1	0	0
3	G	1	Total O P 5 4 1	0	0
3	G	1	Total O P 5 4 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

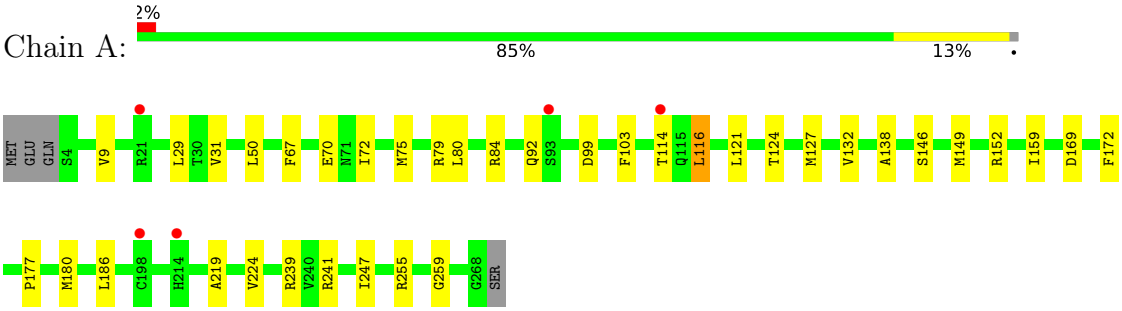
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total O 1 1	0	0
6	C	4	Total O 4 4	0	0
6	E	3	Total O 3 3	0	0
6	G	2	Total O 2 2	0	0

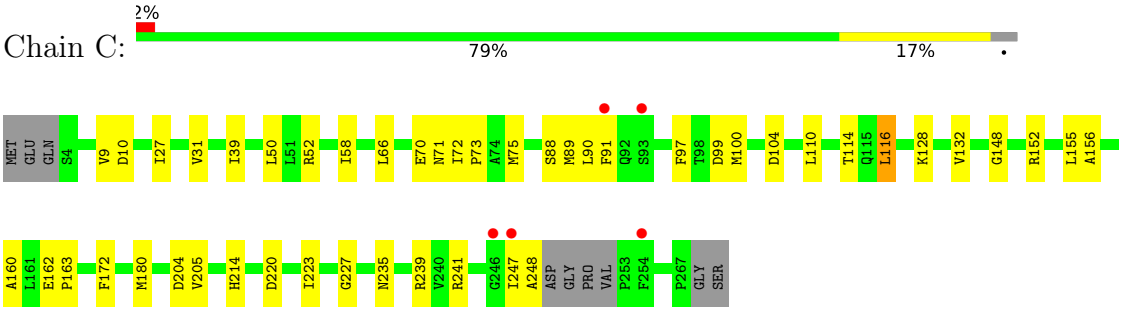
3 Residue-property plots (i)

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.

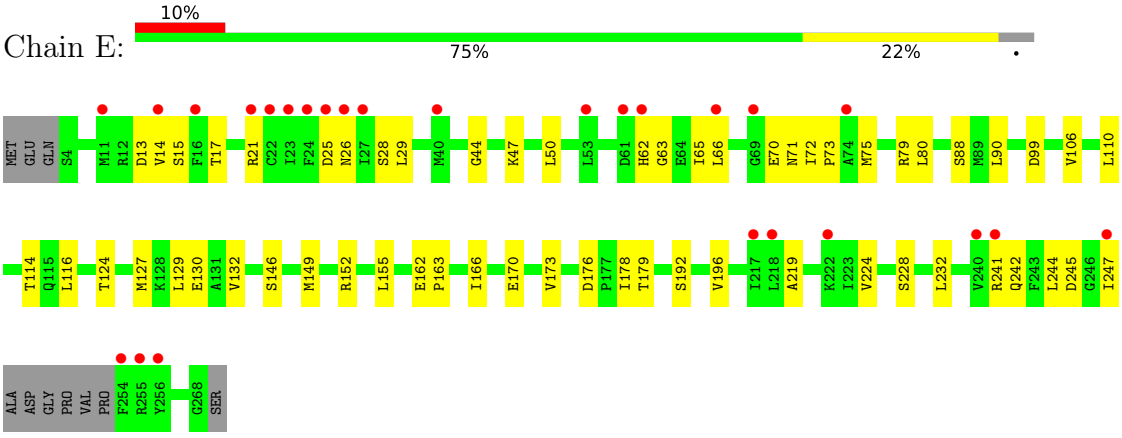
• Molecule 1: Organic solvent ABC transporter ATP-binding protein



• Molecule 1: Organic solvent ABC transporter ATP-binding protein

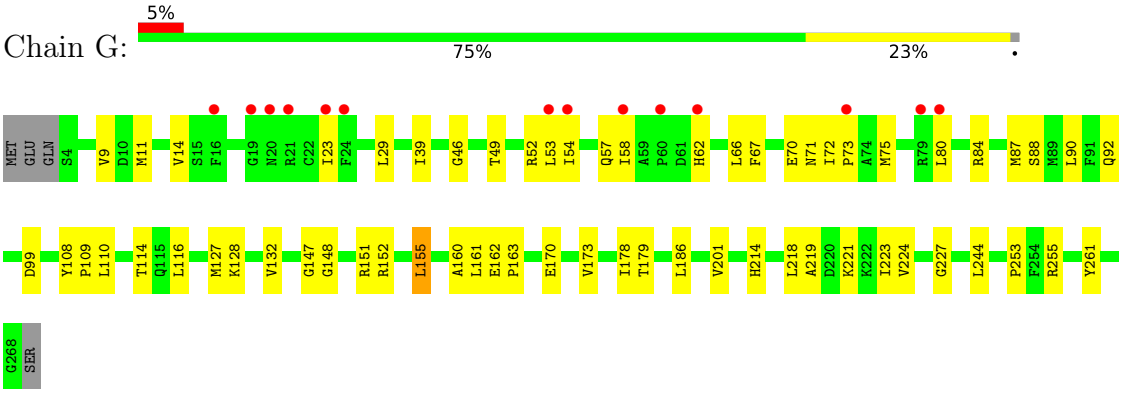


• Molecule 1: Organic solvent ABC transporter ATP-binding protein



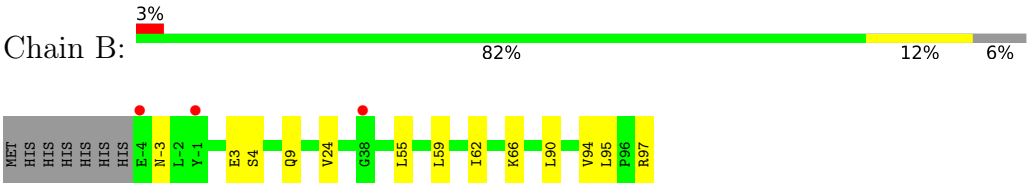
• Molecule 1: Organic solvent ABC transporter ATP-binding protein

Chain G:



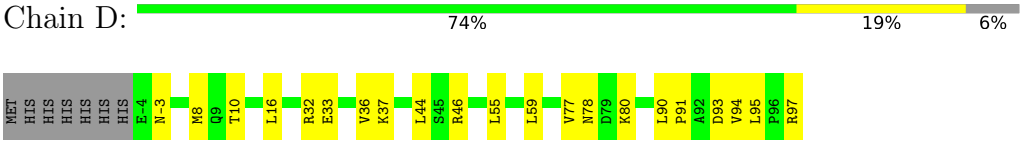
• Molecule 2: ABC transporter maintaining OM lipid asymmetry, cytoplasmic STAS component

Chain B:



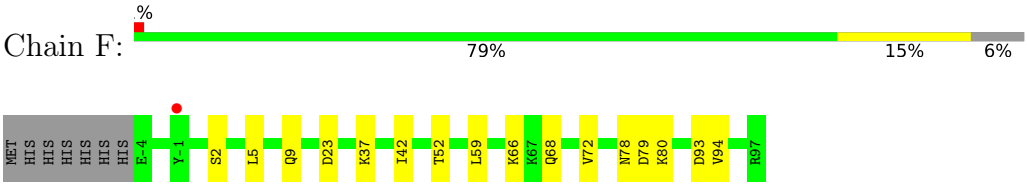
• Molecule 2: ABC transporter maintaining OM lipid asymmetry, cytoplasmic STAS component

Chain D:



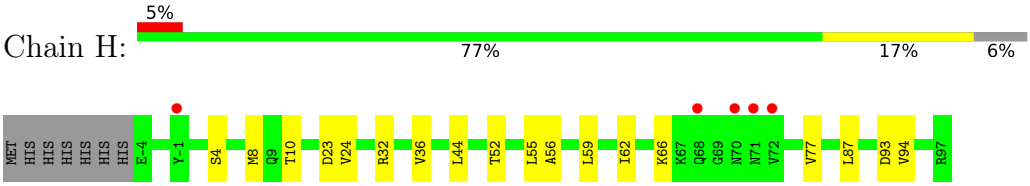
• Molecule 2: ABC transporter maintaining OM lipid asymmetry, cytoplasmic STAS component

Chain F:



• Molecule 2: ABC transporter maintaining OM lipid asymmetry, cytoplasmic STAS component

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	78.27Å 136.08Å 261.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.87 – 2.60 48.87 – 2.60	Depositor EDS
% Data completeness (in resolution range)	92.1 (48.87-2.60) 91.3 (48.87-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.27 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.203 , 0.237 0.203 , 0.237	Depositor DCC
R_{free} test set	1998 reflections (2.30%)	wwPDB-VP
Wilson B-factor (Å^2)	72.4	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 65.4	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	11433	wwPDB-VP
Average B, all atoms (Å^2)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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: PO4, GOL, EDO

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	0.28	0/2051	0.48	0/2785
1	C	0.29	0/2038	0.53	1/2762 (0.0%)
1	E	0.30	0/2014	0.55	1/2730 (0.0%)
1	G	0.31	0/2066	0.54	0/2807
2	B	0.29	0/864	0.50	0/1174
2	D	0.28	0/849	0.52	0/1155
2	F	0.27	0/841	0.50	0/1144
2	H	0.30	0/853	0.50	0/1160
All	All	0.29	0/11576	0.52	2/15717 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	116	LEU	CA-CB-CG	5.75	128.53	115.30
1	C	116	LEU	CA-CB-CG	5.69	128.40	115.30

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	2010	0	2042	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1999	0	2036	32	0
1	E	1977	0	2011	41	0
1	G	2024	0	2050	45	0
2	B	850	0	853	10	0
2	D	835	0	837	10	0
2	F	827	0	832	9	0
2	H	839	0	841	12	0
3	A	10	0	0	0	0
3	C	10	0	0	0	0
3	E	10	0	0	2	0
3	G	10	0	0	1	0
4	A	6	0	8	1	0
4	C	6	0	8	0	0
4	E	6	0	8	0	0
5	A	4	0	6	0	0
6	A	1	0	0	0	0
6	C	4	0	0	0	0
6	E	3	0	0	0	0
6	G	2	0	0	0	0
All	All	11433	0	11532	172	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 7.

All (172) 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)
2:B:95:LEU:O	2:B:97:ARG:NH1	2.10	0.84
1:E:146:SER:H	1:E:149:MET:HE3	1.42	0.82
1:E:75:MET:HE1	1:E:80:LEU:HD13	1.65	0.79
2:F:37:LYS:HA	2:F:68:GLN:HE22	1.50	0.75
1:G:52:ARG:NH2	1:G:57:GLN:OE1	2.20	0.74
1:E:47:LYS:N	3:E:301:PO4:O3	2.21	0.72
1:G:132:VAL:CG1	1:G:152:ARG:HB3	2.20	0.72
2:B:90:LEU:O	2:B:97:ARG:NH2	2.23	0.71
1:E:70:GLU:OE2	1:E:79:ARG:NE	2.22	0.71
1:C:235:ASN:O	1:C:241:ARG:NH1	2.24	0.70
1:A:146:SER:H	1:A:149:MET:HE3	1.56	0.70
1:C:172:PHE:HB3	1:C:180:MET:HG3	1.74	0.69
1:E:132:VAL:HG12	1:E:152:ARG:HB3	1.75	0.69
1:E:14:VAL:HA	1:E:63:GLY:HA3	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:67:PHE:HB2	1:G:72:ILE:HD11	1.75	0.66
2:B:59:LEU:HD22	2:B:94:VAL:HG12	1.77	0.66
1:E:178:ILE:HD12	1:G:244:LEU:HD11	1.78	0.65
2:B:3:GLU:N	2:B:3:GLU:OE1	2.27	0.65
2:H:62:ILE:HG22	2:H:66:LYS:HE2	1.79	0.65
1:E:65:ILE:HB	1:E:73:PRO:HD3	1.78	0.64
1:E:75:MET:HE3	1:E:80:LEU:HB2	1.80	0.64
1:G:219:ALA:HB3	1:G:224:VAL:HG21	1.79	0.63
1:A:132:VAL:CG1	1:A:152:ARG:HB3	2.29	0.63
2:B:4:SER:HB2	2:B:24:VAL:HG11	1.81	0.62
2:D:59:LEU:HD22	2:D:94:VAL:HG12	1.80	0.62
2:H:59:LEU:HD22	2:H:94:VAL:HG12	1.81	0.62
1:C:132:VAL:CG1	1:C:152:ARG:HB3	2.28	0.62
1:A:132:VAL:HG13	1:A:152:ARG:HB3	1.82	0.62
1:G:9:VAL:HG11	1:G:54:ILE:HD12	1.82	0.62
1:E:242:GLN:NE2	1:E:247:ILE:O	2.33	0.61
1:A:114:THR:HG22	1:A:116:LEU:HB3	1.82	0.60
1:A:219:ALA:HB3	1:A:224:VAL:HG21	1.83	0.60
1:E:75:MET:CE	1:E:80:LEU:HD13	2.31	0.60
1:E:132:VAL:CG1	1:E:152:ARG:HB3	2.32	0.59
1:E:44:GLY:N	3:E:301:PO4:O2	2.35	0.57
2:F:59:LEU:HD22	2:F:94:VAL:HG12	1.86	0.57
1:G:114:THR:HG23	1:G:116:LEU:H	1.68	0.57
2:B:-3[A]:ASN:HA	2:B:9:GLN:O	2.04	0.57
1:G:147:GLY:O	1:G:151:ARG:HG3	2.05	0.57
2:D:44:LEU:HD12	2:D:77:VAL:HG22	1.86	0.57
2:H:66:LYS:NZ	2:H:93:ASP:O	2.38	0.57
1:A:255:ARG:HG2	1:G:253:PRO:HB2	1.87	0.56
1:E:90:LEU:HD22	1:E:155:LEU:HD13	1.87	0.56
1:G:80:LEU:HD21	1:G:84[B]:ARG:HH11	1.68	0.56
1:A:172:PHE:HB3	1:A:180:MET:HG3	1.88	0.55
2:B:-3[B]:ASN:HA	2:B:9:GLN:O	2.05	0.55
1:E:219:ALA:HB3	1:E:224:VAL:HG21	1.89	0.55
1:C:114:THR:HG23	1:C:116:LEU:H	1.73	0.54
1:A:92:GLN:NE2	1:A:169:ASP:O	2.40	0.54
1:E:72:ILE:HA	1:E:75:MET:HG3	1.88	0.54
1:G:162:GLU:OE2	2:H:52:THR:OG1	2.21	0.53
1:G:132:VAL:HG13	1:G:152:ARG:HB3	1.91	0.53
1:A:224:VAL:HG13	1:A:239:ARG:HH21	1.73	0.53
1:G:148:GLY:O	1:G:152:ARG:HG3	2.09	0.53
2:B:95:LEU:HB3	2:B:97:ARG:HH12	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:15:SER:OG	1:E:26:ASN:N	2.42	0.52
1:G:46:GLY:N	3:G:302:PO4:O4	2.42	0.52
1:C:9:VAL:HB	1:C:31:VAL:HB	1.91	0.52
1:C:72:ILE:HA	1:C:75:MET:HE2	1.92	0.51
1:G:170:GLU:HB3	1:G:173:VAL:HG13	1.93	0.51
1:C:97:PHE:HB3	1:C:99:ASP:OD1	2.10	0.51
1:G:110:LEU:O	1:G:114:THR:HG22	2.09	0.51
1:A:9:VAL:HB	1:A:31:VAL:HB	1.93	0.51
1:A:70:GLU:OE2	1:A:79:ARG:NH2	2.34	0.51
2:H:4:SER:HB2	2:H:24:VAL:HG11	1.91	0.51
1:A:72:ILE:HA	1:A:75:MET:SD	2.51	0.50
1:C:110:LEU:O	1:C:114:THR:HG22	2.11	0.50
1:E:192:SER:O	2:F:80:LYS:HE3	2.12	0.50
2:F:2:SER:HB3	2:F:5:LEU:O	2.12	0.50
1:G:71:ASN:O	1:G:75:MET:HG3	2.12	0.50
1:G:84[A]:ARG:HA	1:G:87:MET:HE2	1.94	0.49
1:C:88:SER:HB2	1:C:163:PRO:HG2	1.94	0.49
1:C:132:VAL:HG21	1:C:156:ALA:HB2	1.94	0.49
1:C:52:ARG:HB3	1:C:58:ILE:HG12	1.94	0.49
1:G:84[B]:ARG:HA	1:G:87:MET:HE2	1.95	0.49
1:E:66:LEU:HA	1:E:70:GLU:O	2.13	0.49
1:C:99:ASP:OD1	1:C:99:ASP:N	2.40	0.49
2:F:66:LYS:NZ	2:F:93:ASP:O	2.46	0.49
1:G:66:LEU:HA	1:G:70:GLU:O	2.13	0.49
1:C:148:GLY:O	1:C:152:ARG:HG3	2.13	0.49
1:C:90:LEU:HD22	1:C:155:LEU:HD13	1.93	0.49
1:E:162:GLU:OE2	2:F:52:THR:OG1	2.20	0.49
1:A:67:PHE:HB2	1:A:72:ILE:HD11	1.95	0.48
1:C:10:ASP:HB3	1:C:66:LEU:HB2	1.96	0.48
1:C:132:VAL:HG12	1:C:152:ARG:HB3	1.94	0.48
1:C:220:ASP:OD2	1:C:239[B]:ARG:NH2	2.46	0.48
2:F:42:ILE:HG13	2:F:72:VAL:HG11	1.96	0.48
1:G:11:MET:HE2	1:G:29:LEU:HB3	1.95	0.48
2:H:8:MET:HG2	2:H:10:THR:HG23	1.96	0.48
1:G:99:ASP:N	1:G:99:ASP:OD1	2.46	0.48
1:E:110:LEU:O	1:E:114:THR:HG22	2.13	0.48
1:A:75:MET:HE3	1:A:79:ARG:HG3	1.97	0.47
1:A:177:PRO:HG3	1:C:248:ALA:HB1	1.96	0.47
1:C:89:MET:HB3	1:C:91:PHE:CE2	2.49	0.47
1:C:214[A]:HIS:CE1	1:C:227:GLY:HA2	2.49	0.47
2:H:55:LEU:O	2:H:59:LEU:HG	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:99:ASP:N	1:E:99:ASP:OD1	2.46	0.47
2:B:55:LEU:O	2:B:59:LEU:HG	2.15	0.47
1:G:152:ARG:HD3	1:G:179:THR:HG21	1.95	0.47
1:E:88:SER:HB2	1:E:163:PRO:HG2	1.97	0.47
1:G:88:SER:HB2	1:G:163:PRO:HG2	1.97	0.47
2:D:33:GLU:O	2:D:37:LYS:HG2	2.16	0.46
1:G:39:ILE:HB	1:G:201:VAL:HG22	1.98	0.46
2:D:91:PRO:HB2	2:D:93:ASP:OD1	2.14	0.46
2:D:78:ASN:HD21	2:D:80:LYS:NZ	2.15	0.45
1:E:17:THR:HG23	1:E:21:ARG:O	2.16	0.45
1:G:53:LEU:HD23	1:G:58:ILE:HB	1.98	0.45
1:A:124:THR:HA	1:A:127:MET:HE2	1.99	0.45
1:E:71:ASN:OD1	1:E:73:PRO:HD2	2.16	0.45
1:E:162:GLU:N	1:E:163:PRO:HD3	2.32	0.45
1:E:176:ASP:OD2	1:G:255:ARG:NH2	2.50	0.45
1:E:29:LEU:HD21	1:E:50:LEU:HD21	1.98	0.45
1:E:146:SER:H	1:E:149:MET:CE	2.19	0.45
2:H:32:ARG:O	2:H:36:VAL:HG22	2.17	0.45
1:G:128:LYS:HE3	1:G:160:ALA:O	2.17	0.44
1:E:228:SER:O	1:E:232:LEU:HD13	2.18	0.44
1:A:80:LEU:HD21	1:A:84:ARG:HH11	1.81	0.44
1:G:110:LEU:HD23	1:G:161:LEU:HD23	1.99	0.44
2:B:62:ILE:HG22	2:B:66:LYS:HE3	1.99	0.44
1:E:166:ILE:HD12	1:E:196:VAL:HG11	2.00	0.44
1:G:23:ILE:O	1:G:221:LYS:HD2	2.17	0.44
2:D:8:MET:HG2	2:D:10:THR:HG23	1.98	0.44
1:G:214[B]:HIS:NE2	1:G:227:GLY:HA2	2.33	0.44
1:A:99:ASP:N	1:A:99:ASP:OD1	2.47	0.44
1:E:130:GLU:HG2	1:G:261:TYR:CZ	2.52	0.44
1:E:152:ARG:HD3	1:E:179:THR:HG21	2.00	0.44
1:A:116:LEU:HD21	1:A:121:LEU:HB2	2.00	0.43
2:D:32:ARG:O	2:D:36:VAL:HG22	2.17	0.43
1:G:57:GLN:HE21	1:G:84[A]:ARG:HH22	1.66	0.43
1:G:127:MET:HE1	2:H:56:ALA:HB2	2.00	0.43
1:G:186:LEU:HG	2:H:87:LEU:HD11	2.00	0.43
1:C:27:ILE:HB	1:C:223:ILE:CD1	2.48	0.43
1:E:71:ASN:O	1:E:75:MET:HG3	2.18	0.43
1:G:14:VAL:HA	1:G:62:HIS:O	2.19	0.43
1:G:49:THR:O	1:G:53:LEU:HG	2.17	0.43
1:C:100:MET:HB2	1:C:104:ASP:HB2	2.01	0.43
2:D:95:LEU:HB3	2:D:97:ARG:HH12	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:23:ASP:OD1	2:H:23:ASP:N	2.50	0.43
1:C:71:ASN:OD1	1:C:73:PRO:HD2	2.19	0.43
1:E:15:SER:HB2	1:E:62:HIS:HB3	2.01	0.43
1:C:128:LYS:HE3	1:C:160:ALA:O	2.18	0.42
1:C:205:VAL:HB	1:C:247:ILE:HG21	2.00	0.42
1:C:39:ILE:HD13	1:C:50:LEU:HD23	2.00	0.42
1:A:103:PHE:HB2	1:A:138:ALA:HB1	2.01	0.42
1:C:162:GLU:N	1:C:163:PRO:HD3	2.33	0.42
1:A:259:GLY:H	4:A:303:GOL:H11	1.84	0.42
1:E:106:VAL:HG21	1:E:129:LEU:HD21	2.00	0.42
1:A:159:ILE:HD13	1:A:186:LEU:HD21	2.01	0.42
2:F:78:ASN:OD1	2:F:79:ASP:N	2.53	0.42
1:G:108:TYR:HB3	1:G:109:PRO:HD3	2.02	0.42
2:D:55:LEU:HD21	2:D:90:LEU:HD11	2.02	0.42
1:G:162:GLU:N	1:G:163:PRO:HD3	2.35	0.41
1:A:80:LEU:O	1:A:84:ARG:HG2	2.20	0.41
1:E:13:ASP:N	1:E:28:SER:OG	2.24	0.41
1:G:90:LEU:HD22	1:G:155:LEU:HD12	2.02	0.41
1:C:247:ILE:HD12	1:C:247:ILE:HA	1.85	0.41
1:A:29:LEU:HD21	1:A:50:LEU:HD21	2.02	0.41
2:D:16:LEU:O	2:D:46:ARG:HB2	2.19	0.41
1:G:72:ILE:N	1:G:73:PRO:HD2	2.36	0.41
1:C:172:PHE:HB2	1:C:204:ASP:OD2	2.20	0.41
1:E:241:ARG:O	1:E:245:ASP:CB	2.69	0.41
1:A:177:PRO:CG	1:C:248:ALA:HB1	2.50	0.41
1:E:244:LEU:HB2	1:G:178:ILE:HD11	2.02	0.41
1:G:57:GLN:HG2	1:G:84[B]:ARG:HH22	1.84	0.41
1:C:66:LEU:HA	1:C:70:GLU:O	2.21	0.41
1:E:124:THR:HA	1:E:127:MET:HE3	2.02	0.41
2:H:44:LEU:HD12	2:H:77:VAL:HG22	2.02	0.40
1:C:71:ASN:O	1:C:75:MET:HG3	2.21	0.40
1:E:170:GLU:HB3	1:E:173:VAL:CG2	2.52	0.40
1:G:57:GLN:CG	1:G:84[B]:ARG:HH22	2.34	0.40
1:G:218:LEU:HG	1:G:223:ILE:HD13	2.02	0.40
2:F:23:ASP:OD1	2:F:23:ASP:N	2.54	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	264/269 (98%)	264 (100%)	0	0	100	100
1	C	259/269 (96%)	258 (100%)	1 (0%)	0	100	100
1	E	257/269 (96%)	257 (100%)	0	0	100	100
1	G	266/269 (99%)	266 (100%)	0	0	100	100
2	B	105/109 (96%)	105 (100%)	0	0	100	100
2	D	104/109 (95%)	103 (99%)	1 (1%)	0	100	100
2	F	103/109 (94%)	103 (100%)	0	0	100	100
2	H	104/109 (95%)	104 (100%)	0	0	100	100
All	All	1462/1512 (97%)	1460 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	215/220 (98%)	212 (99%)	3 (1%)	67	85
1	C	214/220 (97%)	214 (100%)	0	100	100
1	E	211/220 (96%)	210 (100%)	1 (0%)	88	96
1	G	215/220 (98%)	213 (99%)	2 (1%)	78	91
2	B	94/96 (98%)	94 (100%)	0	100	100
2	D	92/96 (96%)	90 (98%)	2 (2%)	52	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	F	91/96 (95%)	90 (99%)	1 (1%)	73 88
2	H	93/96 (97%)	93 (100%)	0	100 100
All	All	1225/1264 (97%)	1216 (99%)	9 (1%)	84 94

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	LEU
1	A	241	ARG
1	A	247	ILE
2	D	-3[A]	ASN
2	D	-3[B]	ASN
1	E	25	ASP
2	F	9	GLN
1	G	92	GLN
1	G	155	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

12 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	PO4	A	302	-	4,4,4	0.86	0	6,6,6	0.53	0
3	PO4	G	301	-	4,4,4	0.93	0	6,6,6	0.51	0
4	GOL	C	303	-	5,5,5	0.89	0	5,5,5	1.03	0
3	PO4	C	302	-	4,4,4	1.07	0	6,6,6	0.27	0
3	PO4	A	301	-	4,4,4	0.93	0	6,6,6	0.45	0
4	GOL	A	303	-	5,5,5	0.95	0	5,5,5	0.98	0
3	PO4	E	302	-	4,4,4	0.92	0	6,6,6	0.46	0
3	PO4	C	301	-	4,4,4	0.94	0	6,6,6	0.38	0
3	PO4	G	302	-	4,4,4	0.92	0	6,6,6	0.62	0
5	EDO	A	304	-	3,3,3	0.48	0	2,2,2	0.29	0
3	PO4	E	301	-	4,4,4	0.98	0	6,6,6	0.34	0
4	GOL	E	303	-	5,5,5	0.97	0	5,5,5	1.01	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
5	EDO	A	304	-	-	0/1/1/1	-
4	GOL	C	303	-	-	0/4/4/4	-
4	GOL	E	303	-	-	0/4/4/4	-
4	GOL	A	303	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	GOL	1	0
3	G	302	PO4	1	0
3	E	301	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	265/269 (98%)	0.08	5 (1%) 66 62	55, 80, 122, 172	0
1	C	260/269 (96%)	0.08	5 (1%) 66 62	55, 77, 125, 174	0
1	E	259/269 (96%)	0.52	26 (10%) 7 4	55, 99, 167, 193	0
1	G	265/269 (98%)	0.33	14 (5%) 26 20	61, 96, 149, 188	0
2	B	102/109 (93%)	0.30	3 (2%) 51 45	63, 89, 129, 154	0
2	D	102/109 (93%)	0.21	0 100 100	56, 87, 127, 142	0
2	F	102/109 (93%)	0.13	1 (0%) 82 80	70, 94, 134, 141	0
2	H	102/109 (93%)	0.21	5 (4%) 29 23	69, 92, 128, 156	0
All	All	1457/1512 (96%)	0.24	59 (4%) 38 31	55, 88, 148, 193	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	24	PHE	7.7
1	C	246	GLY	5.9
1	E	254	PHE	5.2
1	G	80	LEU	4.8
1	G	54	ILE	4.8
2	B	-1[A]	TYR	4.5
1	E	218	LEU	4.5
1	E	21	ARG	4.1
1	G	53	LEU	4.0
1	E	62	HIS	3.8
1	E	23	ILE	3.8
1	E	27	ILE	3.7
1	G	21	ARG	3.6
1	E	16	PHE	3.6
1	E	66	LEU	3.4
1	E	240	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	256	TYR	3.2
1	G	19	GLY	3.1
1	G	73	PRO	3.1
1	A	21	ARG	3.0
1	E	26	ASN	3.0
2	H	68	GLN	2.9
1	E	22	CYS	2.9
1	C	247	ILE	2.9
2	F	-1[A]	TYR	2.9
1	G	16	PHE	2.8
1	G	62	HIS	2.8
1	C	254	PHE	2.8
1	G	23	ILE	2.7
1	E	247	ILE	2.7
1	E	25	ASP	2.7
1	E	255	ARG	2.6
1	A	93	SER	2.6
1	G	60	PRO	2.5
1	G	58	ILE	2.4
1	E	11	MET	2.4
1	E	241	ARG	2.4
2	B	38	GLY	2.3
2	H	70	ASN	2.3
1	E	61	ASP	2.3
1	E	222	LYS	2.3
2	H	72	VAL	2.3
1	A	198	CYS	2.3
2	H	-1[A]	TYR	2.2
1	E	14	VAL	2.2
1	G	20	ASN	2.2
1	E	53	LEU	2.2
1	E	74	ALA	2.2
2	B	-4	GLU	2.2
1	E	217	ILE	2.2
1	G	24	PHE	2.2
1	C	93	SER	2.1
1	E	40	MET	2.1
1	C	91	PHE	2.1
1	A	214[A]	HIS	2.1
1	E	69	GLY	2.1
2	H	71	ASN	2.1
1	G	79	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	114	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 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
5	EDO	A	304	4/4	0.67	0.22	133,139,154,169	0
4	GOL	E	303	6/6	0.78	0.29	113,124,130,135	0
4	GOL	A	303	6/6	0.84	0.24	90,103,123,124	0
4	GOL	C	303	6/6	0.87	0.22	77,92,97,103	0
3	PO4	E	302	5/5	0.91	0.13	105,110,118,129	0
3	PO4	C	302	5/5	0.95	0.09	113,113,117,121	0
3	PO4	G	302	5/5	0.95	0.13	79,106,110,113	0
3	PO4	E	301	5/5	0.95	0.08	98,110,126,128	0
3	PO4	G	301	5/5	0.96	0.14	82,98,112,114	0
3	PO4	A	302	5/5	0.97	0.12	92,95,98,100	0
3	PO4	C	301	5/5	0.98	0.10	90,93,98,100	0
3	PO4	A	301	5/5	0.98	0.11	70,81,94,100	0

6.5 Other polymers (i)

There are no such residues in this entry.