



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 01:02 pm GMT

PDB ID : 2YF3
Title : Crystal structure of DR2231, the MazG-like protein from *Deinococcus radiodurans*, complex with manganese
Authors : Goncalves, A.M.D.; deSanctis, D.; McSweeney, S.M.
Deposited on : 2011-04-01
Resolution : 2.00 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk31020
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk31020

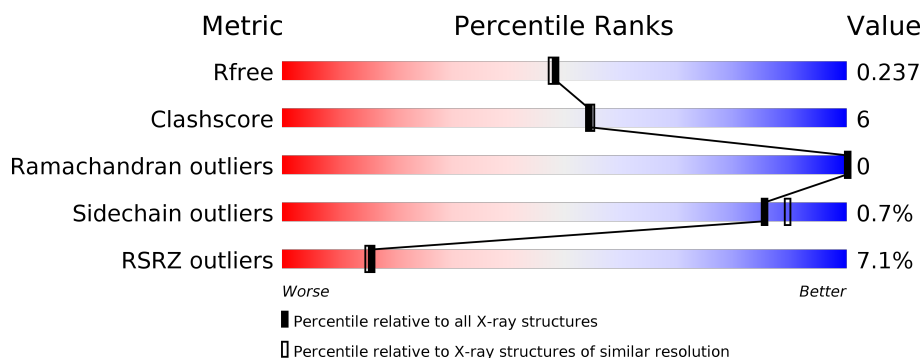
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	111664	7193 (2.00-2.00)
Clashscore	122126	8267 (2.00-2.00)
Ramachandran outliers	120053	8166 (2.00-2.00)
Sidechain outliers	120020	8165 (2.00-2.00)
RSRZ outliers	108989	7011 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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 $\leq 5\%$. 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	154	<div> <div>9%</div> <div> <div></div> <div>93%</div> <div>5%</div> <div>•</div> </div> </div>
1	B	154	<div> <div>7%</div> <div> <div></div> <div>87%</div> <div>8%</div> <div>• •</div> </div> </div>
1	C	154	<div> <div>6%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>• •</div> </div> </div>
1	D	154	<div> <div>6%</div> <div> <div></div> <div>94%</div> <div>5%</div> <div>• •</div> </div> </div>
1	E	154	<div> <div>7%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>• •</div> </div> </div>
1	F	154	<div> <div>6%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>• •</div> </div> </div>

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
4	GOL	C	1155	-	-	X	-
4	GOL	D	1155	-	-	X	-
4	GOL	F	1151	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15069 atoms, of which 7073 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 MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	150	Total	C	H	N	O	S	0	0	0
			2283	716	1135	212	218	2			
1	B	149	Total	C	H	N	O	S	0	0	0
			2270	713	1128	211	216	2			
1	C	153	Total	C	H	N	O	S	0	0	0
			2323	730	1151	215	225	2			
1	D	153	Total	C	H	N	O	S	0	0	0
			2318	728	1150	215	223	2			
1	E	153	Total	C	H	N	O	S	0	0	0
			2324	730	1152	215	225	2			
1	F	148	Total	C	H	N	O	S	0	0	0
			2251	707	1117	210	215	2			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP Q9RS96
A	-4	ILE	-	expression tag	UNP Q9RS96
A	-3	ASP	-	expression tag	UNP Q9RS96
A	-2	PRO	-	expression tag	UNP Q9RS96
A	-1	PHE	-	expression tag	UNP Q9RS96
A	0	THR	-	expression tag	UNP Q9RS96
B	-5	GLY	-	expression tag	UNP Q9RS96
B	-4	ILE	-	expression tag	UNP Q9RS96
B	-3	ASP	-	expression tag	UNP Q9RS96
B	-2	PRO	-	expression tag	UNP Q9RS96
B	-1	PHE	-	expression tag	UNP Q9RS96
B	0	THR	-	expression tag	UNP Q9RS96
C	-5	GLY	-	expression tag	UNP Q9RS96
C	-4	ILE	-	expression tag	UNP Q9RS96
C	-3	ASP	-	expression tag	UNP Q9RS96
C	-2	PRO	-	expression tag	UNP Q9RS96

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	PHE	-	expression tag	UNP Q9RS96
C	0	THR	-	expression tag	UNP Q9RS96
D	-5	GLY	-	expression tag	UNP Q9RS96
D	-4	ILE	-	expression tag	UNP Q9RS96
D	-3	ASP	-	expression tag	UNP Q9RS96
D	-2	PRO	-	expression tag	UNP Q9RS96
D	-1	PHE	-	expression tag	UNP Q9RS96
D	0	THR	-	expression tag	UNP Q9RS96
E	-5	GLY	-	expression tag	UNP Q9RS96
E	-4	ILE	-	expression tag	UNP Q9RS96
E	-3	ASP	-	expression tag	UNP Q9RS96
E	-2	PRO	-	expression tag	UNP Q9RS96
E	-1	PHE	-	expression tag	UNP Q9RS96
E	0	THR	-	expression tag	UNP Q9RS96
F	-5	GLY	-	expression tag	UNP Q9RS96
F	-4	ILE	-	expression tag	UNP Q9RS96
F	-3	ASP	-	expression tag	UNP Q9RS96
F	-2	PRO	-	expression tag	UNP Q9RS96
F	-1	PHE	-	expression tag	UNP Q9RS96
F	0	THR	-	expression tag	UNP Q9RS96

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	2	Total Mn 2 2	0	0
2	E	2	Total Mn 2 2	0	0
2	B	2	Total Mn 2 2	0	0
2	C	2	Total Mn 2 2	0	0
2	A	1	Total Mn 1 1	0	0
2	F	2	Total Mn 2 2	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	A	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	B	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	C	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		
4	D	1	Total	C	H	O	0	0
			14	3	8	3		
4	E	1	Total	C	H	O	0	0
			14	3	8	3		
4	E	1	Total	C	H	O	0	0
			14	3	8	3		
4	E	1	Total	C	H	O	0	0
			14	3	8	3		
4	F	1	Total	C	H	O	0	0
			14	3	8	3		
4	F	1	Total	C	H	O	0	0
			14	3	8	3		
4	F	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	133	Total	O	0	0
			133	133		
5	B	113	Total	O	0	0
			113	113		

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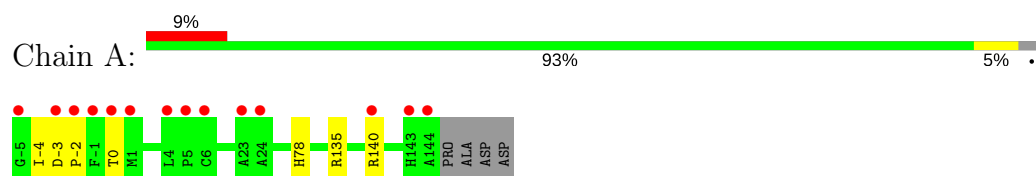
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	164	Total 164	O 164	0	0
5	D	155	Total 155	O 155	0	0
5	E	125	Total 125	O 125	0	0
5	F	109	Total 109	O 109	0	0

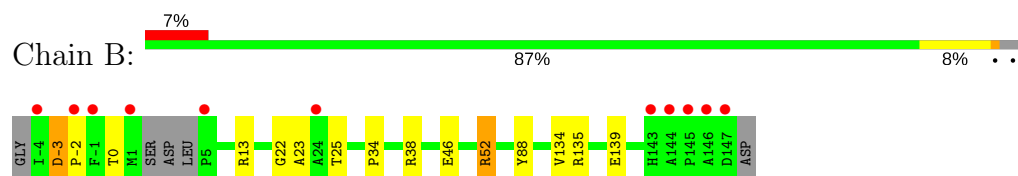
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

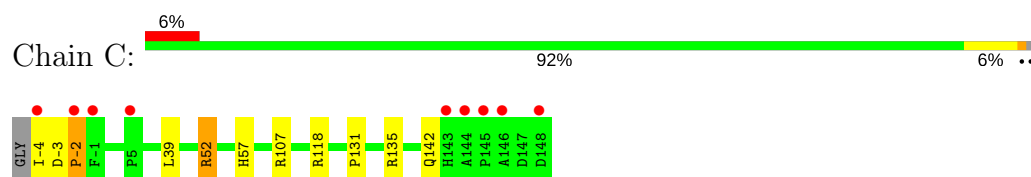
• Molecule 1: MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE



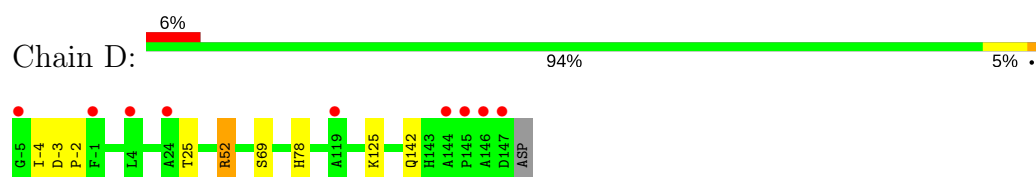
• Molecule 1: MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE



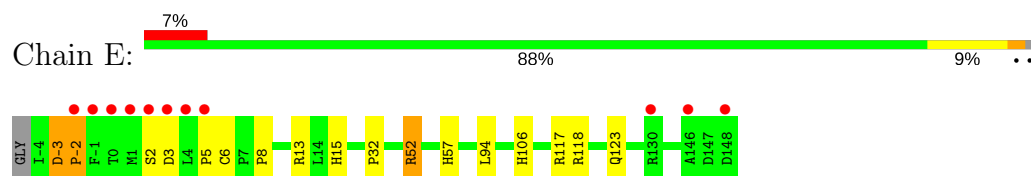
• Molecule 1: MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE



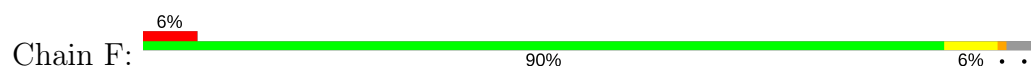
• Molecule 1: MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE

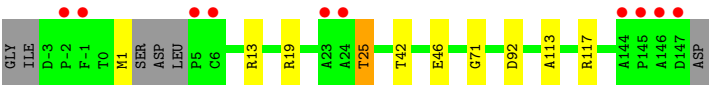


• Molecule 1: MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE



• Molecule 1: MAZG-LIKE NUCLEOSIDE TRIPHOSPHATE PYROPHOSPHOHYDROLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.11Å 111.40Å 166.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.12 – 2.00 43.12 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (43.12-2.00) 99.7 (43.12-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.10 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.191 , 0.238 0.192 , 0.237	Depositor DCC
R_{free} test set	4239 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.649	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15069	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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, MN, SO4

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	0.88	1/1174 (0.1%)	0.77	0/1600
1	B	1.01	3/1168 (0.3%)	0.80	3/1591 (0.2%)
1	C	0.88	1/1199 (0.1%)	0.88	3/1636 (0.2%)
1	D	0.87	0/1195	0.76	2/1630 (0.1%)
1	E	0.86	4/1199 (0.3%)	0.76	3/1636 (0.2%)
1	F	0.79	0/1160	0.72	0/1580
All	All	0.88	9/7095 (0.1%)	0.79	11/9673 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	-3	ASP	CG-OD2	11.64	1.52	1.25
1	E	-2	PRO	C-N	10.95	1.59	1.34
1	B	-3	ASP	CG-OD1	10.39	1.49	1.25
1	E	-3	ASP	CG-OD2	6.19	1.39	1.25
1	C	-2	PRO	C-O	-6.00	1.11	1.23
1	E	-3	ASP	CG-OD1	5.44	1.37	1.25
1	A	0	THR	C-N	5.41	1.46	1.34
1	E	-3	ASP	C-N	5.10	1.44	1.34
1	B	88	TYR	CD2-CE2	5.06	1.47	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	ARG	NE-CZ-NH2	-12.48	114.06	120.30
1	C	52	ARG	NE-CZ-NH1	8.86	124.73	120.30
1	B	-3	ASP	CB-CG-OD2	-7.95	111.15	118.30
1	E	52	ARG	NE-CZ-NH2	-7.35	116.63	120.30
1	D	52	ARG	NE-CZ-NH1	6.69	123.64	120.30
1	D	52	ARG	NE-CZ-NH2	-6.30	117.15	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	107	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	52	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	E	52	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	E	3	ASP	CB-CG-OD2	5.16	122.94	118.30
1	B	52	ARG	NE-CZ-NH1	5.12	122.86	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1148	1135	1134	5	0
1	B	1142	1128	1127	17	0
1	C	1172	1151	1150	15	0
1	D	1168	1150	1149	11	0
1	E	1172	1152	1151	15	1
1	F	1134	1117	1116	15	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	15	0	0	0	0
3	B	15	0	0	1	0
3	C	15	0	0	0	0
3	D	10	0	0	0	0
3	E	10	0	0	0	0
3	F	5	0	0	0	0
4	A	24	32	32	2	0
4	B	30	40	40	9	0
4	C	48	64	64	13	0
4	D	36	48	47	6	1
4	E	18	24	24	3	0
4	F	24	32	32	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	133	0	0	1	0
5	B	113	0	0	3	0
5	C	164	0	0	5	0
5	D	155	0	0	3	0
5	E	125	0	0	5	0
5	F	109	0	0	2	0
All	All	7996	7073	7066	81	1

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 6.

All (81) 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:52:ARG:HH22	4:C:1155:GOL:H12	0.98	1.09
1:F:25:THR:HG21	1:F:92:ASP:OD2	1.50	1.08
1:C:52:ARG:HH22	4:C:1155:GOL:C1	1.69	1.04
1:C:52:ARG:NH2	4:C:1155:GOL:H12	1.73	1.03
1:F:25:THR:CG2	1:F:92:ASP:OD2	2.09	1.00
1:F:113:ALA:O	4:F:1151:GOL:H2	1.62	0.99
1:C:118:ARG:HH12	4:C:1157:GOL:H32	1.28	0.96
1:B:23:ALA:O	5:B:2025:HOH:O	1.89	0.90
1:B:52:ARG:HH22	4:B:1153:GOL:H2	1.41	0.86
1:F:117:ARG:NE	4:F:1151:GOL:H31	1.91	0.84
1:F:117:ARG:HE	4:F:1151:GOL:H31	1.45	0.80
1:F:117:ARG:HE	4:F:1151:GOL:C3	1.96	0.79
1:F:71:GLY:H	4:F:1150:GOL:H31	1.49	0.76
4:C:1153:GOL:H2	1:E:-2:PRO:HG2	1.65	0.76
1:E:15:HIS:NE2	4:E:1152:GOL:H11	2.02	0.74
1:C:52:ARG:HD3	5:D:2060:HOH:O	1.87	0.73
1:C:118:ARG:NH1	4:C:1157:GOL:H32	2.04	0.69
1:A:140:ARG:NH1	5:A:2125:HOH:O	2.24	0.69
4:A:1149:GOL:H2	1:C:131:PRO:HG3	1.74	0.69
1:D:125:LYS:HD2	4:D:1155:GOL:H11	1.74	0.68
1:F:25:THR:HG22	1:F:92:ASP:OD2	1.94	0.67
1:B:134:VAL:HB	4:B:1154:GOL:H11	1.77	0.66
1:A:135:ARG:HH21	4:A:1149:GOL:H31	1.60	0.65
1:C:135:ARG:HE	4:C:1160:GOL:H12	1.62	0.65
4:C:1160:GOL:H11	5:C:2163:HOH:O	1.99	0.63
1:C:57:HIS:HB2	5:C:2093:HOH:O	2.00	0.62
4:E:1152:GOL:H12	5:E:2115:HOH:O	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:1155:GOL:H31	1:D:52:ARG:HH22	1.63	0.61
1:D:125:LYS:HE3	4:D:1155:GOL:H32	1.82	0.61
1:B:-3:ASP:HB2	1:B:-2:PRO:HD2	1.84	0.59
1:E:-3:ASP:HB2	1:E:-2:PRO:CD	2.34	0.58
1:D:69:SER:HB2	4:D:1152:GOL:H2	1.85	0.58
1:B:46:GLU:HG3	5:B:2037:HOH:O	2.04	0.58
4:D:1156:GOL:H32	5:D:2121:HOH:O	2.03	0.57
1:E:52:ARG:HD3	5:F:2021:HOH:O	2.03	0.57
1:F:71:GLY:HA3	4:F:1150:GOL:H11	1.86	0.56
1:E:118:ARG:HH12	4:E:1153:GOL:H11	1.70	0.56
1:C:135:ARG:HH21	4:C:1160:GOL:H32	1.72	0.55
1:D:25:THR:HG23	1:D:25:THR:O	2.07	0.54
1:B:34:PRO:CD	4:B:1157:GOL:H2	2.38	0.54
1:F:117:ARG:CZ	4:F:1151:GOL:H31	2.37	0.54
1:F:117:ARG:HE	4:F:1151:GOL:H32	1.73	0.53
4:C:1153:GOL:C2	1:E:-2:PRO:HG2	2.36	0.53
1:D:125:LYS:CE	4:D:1155:GOL:H32	2.39	0.52
1:E:8:PRO:O	1:E:13:ARG:NH1	2.44	0.51
1:B:135:ARG:O	1:B:139:GLU:HG3	2.10	0.51
1:E:106:HIS:CD2	1:F:13:ARG:HD3	2.45	0.51
1:C:142:GLN:HG2	5:C:2062:HOH:O	2.10	0.51
1:A:-4:ILE:C	1:A:-4:ILE:HD12	2.31	0.50
1:B:-3:ASP:H	1:B:0:THR:HG1	1.58	0.50
1:A:78:HIS:ND1	4:B:1154:GOL:H12	2.27	0.49
1:E:57:HIS:CD2	5:E:2067:HOH:O	2.65	0.49
1:B:25:THR:HG23	1:B:25:THR:O	2.11	0.49
1:B:34:PRO:HD2	4:B:1157:GOL:H2	1.95	0.49
1:D:142:GLN:HG2	5:D:2045:HOH:O	2.13	0.49
1:B:-3:ASP:HB2	1:B:-2:PRO:CD	2.44	0.47
1:D:-3:ASP:HB2	1:D:-2:PRO:CD	2.45	0.47
1:E:-3:ASP:HB2	1:E:-2:PRO:HD2	1.97	0.46
1:E:123:GLN:O	5:E:2098:HOH:O	2.21	0.46
1:F:117:ARG:NE	4:F:1151:GOL:C3	2.63	0.46
1:C:-4:ILE:N	5:C:2008:HOH:O	2.45	0.46
1:C:-3:ASP:HB2	1:C:-2:PRO:HD2	1.97	0.46
1:B:34:PRO:HD3	4:B:1157:GOL:H2	1.98	0.45
1:B:38:ARG:NH2	3:B:1150:SO4:O2	2.50	0.45
1:B:13:ARG:NE	4:B:1156:GOL:H11	2.31	0.45
1:B:52:ARG:HH22	4:B:1153:GOL:C2	2.22	0.44
1:B:22:GLY:HA2	5:B:2021:HOH:O	2.17	0.44
1:C:39:LEU:HD22	1:E:2:SER:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-3:ASP:HB2	1:A:-2:PRO:CD	2.48	0.43
1:F:113:ALA:O	4:F:1151:GOL:C2	2.50	0.43
1:F:42:THR:O	1:F:46:GLU:HG3	2.18	0.43
4:C:1153:GOL:H12	5:C:2158:HOH:O	2.18	0.43
1:B:13:ARG:HE	4:B:1156:GOL:H11	1.84	0.43
1:D:125:LYS:CD	4:D:1155:GOL:H32	2.49	0.42
1:E:5:PRO:HD3	5:E:2010:HOH:O	2.19	0.42
1:C:52:ARG:NH2	4:C:1155:GOL:C1	2.53	0.42
1:D:-3:ASP:HB2	1:D:-2:PRO:HD2	2.02	0.41
1:D:78:HIS:C	1:D:78:HIS:CD2	2.93	0.41
1:E:32:PRO:HG3	1:E:94:LEU:CD2	2.51	0.41
4:F:1151:GOL:C1	5:F:2092:HOH:O	2.69	0.41
1:E:5:PRO:CD	5:E:2010:HOH:O	2.69	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:117:ARG:HE	4:D:1156:GOL:H32[4_456]	0.79	0.81

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	148/154 (96%)	145 (98%)	3 (2%)	0	100	100
1	B	145/154 (94%)	142 (98%)	3 (2%)	0	100	100
1	C	151/154 (98%)	148 (98%)	3 (2%)	0	100	100
1	D	151/154 (98%)	147 (97%)	4 (3%)	0	100	100
1	E	151/154 (98%)	145 (96%)	6 (4%)	0	100	100
1	F	144/154 (94%)	141 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	890/924 (96%)	868 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	117/120 (98%)	117 (100%)	0	100	100
1	B	116/120 (97%)	116 (100%)	0	100	100
1	C	120/120 (100%)	120 (100%)	0	100	100
1	D	119/120 (99%)	118 (99%)	1 (1%)	83	87
1	E	120/120 (100%)	119 (99%)	1 (1%)	83	87
1	F	115/120 (96%)	112 (97%)	3 (3%)	49	50
All	All	707/720 (98%)	702 (99%)	5 (1%)	85	89

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	-4	ILE
1	E	6	CYS
1	F	1	MET
1	F	19	ARG
1	F	25	THR

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 55 ligands modelled in this entry, 11 are monoatomic - leaving 44 for Mogul analysis.

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	SO4	A	1146	-	4,4,4	0.30	0	6,6,6	0.44	0
3	SO4	A	1147	-	4,4,4	0.10	0	6,6,6	0.44	0
3	SO4	A	1148	-	4,4,4	0.21	0	6,6,6	0.25	0
4	GOL	A	1149	-	5,5,5	0.39	0	5,5,5	0.22	0
4	GOL	A	1150	-	5,5,5	0.59	0	5,5,5	0.53	0
4	GOL	A	1151	-	5,5,5	0.42	0	5,5,5	0.31	0
4	GOL	A	1152	-	5,5,5	0.92	0	5,5,5	0.67	0
3	SO4	B	1150	-	4,4,4	0.18	0	6,6,6	0.30	0
3	SO4	B	1151	-	4,4,4	0.08	0	6,6,6	0.44	0
3	SO4	B	1152	-	4,4,4	0.13	0	6,6,6	0.22	0
4	GOL	B	1153	-	5,5,5	1.65	1 (20%)	5,5,5	1.48	1 (20%)
4	GOL	B	1154	-	5,5,5	0.86	0	5,5,5	0.42	0
4	GOL	B	1155	-	5,5,5	0.65	0	5,5,5	0.48	0
4	GOL	B	1156	-	5,5,5	0.68	0	5,5,5	0.56	0
4	GOL	B	1157	-	5,5,5	0.35	0	5,5,5	0.46	0
3	SO4	C	1150	-	4,4,4	0.34	0	6,6,6	0.18	0
3	SO4	C	1151	-	4,4,4	0.12	0	6,6,6	0.35	0
3	SO4	C	1152	-	4,4,4	0.20	0	6,6,6	0.28	0
4	GOL	C	1153	-	5,5,5	0.47	0	5,5,5	0.66	0
4	GOL	C	1154	-	5,5,5	0.42	0	5,5,5	0.37	0
4	GOL	C	1155	-	5,5,5	0.62	0	5,5,5	0.55	0
4	GOL	C	1156	-	5,5,5	0.43	0	5,5,5	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	C	1157	-	5,5,5	0.58	0	5,5,5	0.22	0
4	GOL	C	1158	-	5,5,5	0.79	0	5,5,5	0.85	0
4	GOL	C	1159	-	5,5,5	0.24	0	5,5,5	0.44	0
4	GOL	C	1160	-	5,5,5	0.25	0	5,5,5	0.55	0
3	SO4	D	1150	-	4,4,4	0.16	0	6,6,6	0.30	0
3	SO4	D	1151	-	4,4,4	0.15	0	6,6,6	0.25	0
4	GOL	D	1152	-	5,5,5	0.67	0	5,5,5	0.93	0
4	GOL	D	1153	-	5,5,5	0.17	0	5,5,5	0.62	0
4	GOL	D	1154	-	5,5,5	0.64	0	5,5,5	0.44	0
4	GOL	D	1155	-	5,5,5	1.31	1 (20%)	5,5,5	0.90	0
4	GOL	D	1156	-	5,5,5	0.88	0	5,5,5	0.49	0
4	GOL	D	1157	-	5,5,5	0.51	0	5,5,5	0.55	0
3	SO4	E	1150	-	4,4,4	0.32	0	6,6,6	0.43	0
3	SO4	E	1151	-	4,4,4	0.13	0	6,6,6	0.50	0
4	GOL	E	1152	-	5,5,5	0.53	0	5,5,5	0.69	0
4	GOL	E	1153	-	5,5,5	0.47	0	5,5,5	0.29	0
4	GOL	E	1154	-	5,5,5	0.58	0	5,5,5	0.29	0
3	SO4	F	1149	-	4,4,4	0.13	0	6,6,6	0.29	0
4	GOL	F	1150	-	5,5,5	0.68	0	5,5,5	0.16	0
4	GOL	F	1151	-	5,5,5	0.60	0	5,5,5	0.49	0
4	GOL	F	1152	-	5,5,5	0.88	0	5,5,5	0.71	0
4	GOL	F	1153	-	5,5,5	0.19	0	5,5,5	0.36	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
3	SO4	A	1146	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1147	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1148	-	-	0/0/0/0	0/0/0/0
4	GOL	A	1149	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1150	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1151	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1152	-	-	0/4/4/4	0/0/0/0
3	SO4	B	1150	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1151	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1152	-	-	0/0/0/0	0/0/0/0
4	GOL	B	1153	-	-	0/4/4/4	0/0/0/0
4	GOL	B	1154	-	-	0/4/4/4	0/0/0/0
4	GOL	B	1155	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	1156	-	-	0/4/4/4	0/0/0/0
4	GOL	B	1157	-	-	0/4/4/4	0/0/0/0
3	SO4	C	1150	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1151	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1152	-	-	0/0/0/0	0/0/0/0
4	GOL	C	1153	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1154	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1155	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1156	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1157	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1158	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1159	-	-	0/4/4/4	0/0/0/0
4	GOL	C	1160	-	-	0/4/4/4	0/0/0/0
3	SO4	D	1150	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1151	-	-	0/0/0/0	0/0/0/0
4	GOL	D	1152	-	-	0/4/4/4	0/0/0/0
4	GOL	D	1153	-	-	0/4/4/4	0/0/0/0
4	GOL	D	1154	-	-	0/4/4/4	0/0/0/0
4	GOL	D	1155	-	-	0/4/4/4	0/0/0/0
4	GOL	D	1156	-	-	0/4/4/4	0/0/0/0
4	GOL	D	1157	-	-	0/4/4/4	0/0/0/0
3	SO4	E	1150	-	-	0/0/0/0	0/0/0/0
3	SO4	E	1151	-	-	0/0/0/0	0/0/0/0
4	GOL	E	1152	-	-	0/4/4/4	0/0/0/0
4	GOL	E	1153	-	-	0/4/4/4	0/0/0/0
4	GOL	E	1154	-	-	0/4/4/4	0/0/0/0
3	SO4	F	1149	-	-	0/0/0/0	0/0/0/0
4	GOL	F	1150	-	-	0/4/4/4	0/0/0/0
4	GOL	F	1151	-	-	0/4/4/4	0/0/0/0
4	GOL	F	1152	-	-	0/4/4/4	0/0/0/0
4	GOL	F	1153	-	-	0/4/4/4	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1155	GOL	O2-C2	-2.68	1.35	1.43
4	B	1153	GOL	O1-C1	2.15	1.51	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1153	GOL	O3-C3-C2	2.65	122.97	110.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

17 monomers are involved in 46 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1149	GOL	2	0
3	B	1150	SO4	1	0
4	B	1153	GOL	2	0
4	B	1154	GOL	2	0
4	B	1156	GOL	2	0
4	B	1157	GOL	3	0
4	C	1153	GOL	3	0
4	C	1155	GOL	5	0
4	C	1157	GOL	2	0
4	C	1160	GOL	3	0
4	D	1152	GOL	1	0
4	D	1155	GOL	4	0
4	D	1156	GOL	1	1
4	E	1152	GOL	2	0
4	E	1153	GOL	1	0
4	F	1150	GOL	2	0
4	F	1151	GOL	9	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, 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	150/154 (97%)	0.07	14 (9%) 8 8	9, 20, 54, 70	0
1	B	149/154 (96%)	0.07	11 (7%) 14 14	9, 22, 62, 93	0
1	C	153/154 (99%)	-0.15	9 (5%) 22 22	8, 17, 44, 55	0
1	D	153/154 (99%)	-0.10	9 (5%) 22 22	8, 17, 42, 52	0
1	E	153/154 (99%)	0.13	11 (7%) 15 15	12, 26, 57, 69	0
1	F	148/154 (96%)	0.08	10 (6%) 17 17	13, 25, 61, 91	0
All	All	906/924 (98%)	0.01	64 (7%) 16 15	8, 20, 53, 93	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	3	ASP	8.1
1	A	-1	PHE	7.1
1	F	145	PRO	6.7
1	B	-4	ILE	6.5
1	F	146	ALA	6.2
1	C	-1	PHE	6.1
1	E	4	LEU	5.8
1	B	-1	PHE	5.7
1	B	146	ALA	5.7
1	A	144	ALA	5.2
1	A	5	PRO	5.2
1	E	2	SER	5.2
1	B	145	PRO	4.8
1	B	1	MET	4.7
1	E	148	ASP	4.6
1	D	146	ALA	4.6
1	B	24	ALA	4.4
1	F	-1	PHE	4.4
1	E	1	MET	4.4

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Mol	Chain	Res	Type	RSRZ
1	F	144	ALA	4.1
1	F	147	ASP	4.0
1	A	4	LEU	4.0
1	E	-1	PHE	4.0
1	B	-2	PRO	3.9
1	D	145	PRO	3.9
1	C	5	PRO	3.8
1	A	143	HIS	3.8
1	D	-5	GLY	3.7
1	A	6	CYS	3.7
1	C	146	ALA	3.7
1	E	5	PRO	3.7
1	A	-5	GLY	3.6
1	C	-2	PRO	3.5
1	A	-2	PRO	3.5
1	D	144	ALA	3.4
1	B	144	ALA	3.3
1	F	24	ALA	3.1
1	D	-1	PHE	3.1
1	F	5	PRO	2.9
1	A	0	THR	2.8
1	C	143	HIS	2.8
1	F	6	CYS	2.8
1	A	23	ALA	2.8
1	C	145	PRO	2.7
1	D	24	ALA	2.7
1	B	143	HIS	2.7
1	A	140	ARG	2.7
1	E	130	ARG	2.6
1	E	0	THR	2.6
1	F	23	ALA	2.6
1	E	146	ALA	2.6
1	C	144	ALA	2.5
1	D	147	ASP	2.5
1	C	148	ASP	2.4
1	B	5	PRO	2.4
1	F	-2	PRO	2.4
1	A	1	MET	2.3
1	C	-4	ILE	2.2
1	D	4	LEU	2.2
1	A	-3	ASP	2.2
1	D	119	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	147	ASP	2.1
1	A	24	ALA	2.1
1	E	-2	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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(\AA^2)	Q<0.9
4	GOL	C	1157	6/6	0.68	0.20	48,49,50,50	0
4	GOL	B	1154	6/6	0.70	0.26	39,42,46,46	0
4	GOL	F	1152	6/6	0.72	0.36	30,38,42,42	0
4	GOL	E	1153	6/6	0.72	0.26	46,48,49,49	0
4	GOL	C	1158	6/6	0.73	0.28	29,40,45,45	0
4	GOL	A	1149	6/6	0.74	0.24	47,50,54,54	0
4	GOL	D	1157	6/6	0.75	0.17	45,48,49,49	0
4	GOL	D	1154	6/6	0.78	0.37	43,47,47,47	0
4	GOL	D	1156	6/6	0.78	0.29	30,37,40,40	0
4	GOL	E	1152	6/6	0.79	0.20	33,35,39,39	0
4	GOL	B	1153	6/6	0.81	0.20	26,30,33,33	0
4	GOL	B	1156	6/6	0.82	0.29	40,43,45,45	0
4	GOL	A	1150	6/6	0.82	0.21	36,41,45,45	0
4	GOL	C	1155	6/6	0.82	0.27	26,28,34,34	0
4	GOL	C	1160	6/6	0.82	0.25	38,39,43,43	0
4	GOL	D	1152	6/6	0.82	0.22	30,39,41,41	0
4	GOL	F	1150	6/6	0.83	0.19	35,37,40,40	0
4	GOL	A	1152	6/6	0.83	0.29	28,29,36,36	0
4	GOL	E	1154	6/6	0.84	0.55	19,36,41,41	0
3	SO4	A	1147	5/5	0.85	0.32	33,35,36,36	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	GOL	B	1157	6/6	0.86	0.34	37,40,43,43	0
4	GOL	C	1156	6/6	0.88	0.21	30,38,40,40	0
3	SO4	C	1151	5/5	0.88	0.20	25,26,28,28	5
4	GOL	B	1155	6/6	0.89	0.35	29,31,41,41	0
4	GOL	F	1153	6/6	0.91	0.17	22,27,32,32	0
4	GOL	F	1151	6/6	0.91	0.33	23,29,32,32	0
4	GOL	C	1159	6/6	0.92	0.20	32,39,45,45	0
4	GOL	D	1155	6/6	0.92	0.27	19,26,32,32	0
4	GOL	C	1154	6/6	0.92	0.11	16,21,22,22	0
3	SO4	F	1149	5/5	0.93	0.25	22,24,25,25	5
3	SO4	A	1148	5/5	0.93	0.17	33,34,34,34	5
4	GOL	C	1153	6/6	0.93	0.16	13,23,31,31	0
3	SO4	D	1150	5/5	0.93	0.18	26,28,30,30	5
3	SO4	B	1152	5/5	0.94	0.17	25,26,26,27	5
2	MN	F	1147	1/1	0.94	0.11	30,30,30,30	1
3	SO4	B	1150	5/5	0.94	0.14	26,28,29,29	5
4	GOL	D	1153	6/6	0.94	0.11	17,24,27,27	0
3	SO4	D	1151	5/5	0.94	0.14	25,26,28,31	5
3	SO4	C	1152	5/5	0.94	0.12	34,35,36,37	5
3	SO4	B	1151	5/5	0.95	0.17	14,17,20,21	5
4	GOL	A	1151	6/6	0.96	0.12	15,20,28,28	0
3	SO4	A	1146	5/5	0.98	0.14	35,36,37,39	0
3	SO4	E	1150	5/5	0.98	0.08	30,31,33,34	0
2	MN	B	1149	1/1	0.98	0.04	24,24,24,24	1
3	SO4	C	1150	5/5	0.98	0.06	27,27,28,29	0
3	SO4	E	1151	5/5	0.98	0.15	14,19,20,22	5
2	MN	F	1148	1/1	0.99	0.09	23,23,23,23	0
2	MN	E	1148	1/1	0.99	0.04	28,28,28,28	0
2	MN	D	1149	1/1	0.99	0.07	19,19,19,19	1
2	MN	E	1149	1/1	0.99	0.11	19,19,19,19	0
2	MN	B	1148	1/1	0.99	0.12	17,17,17,17	0
2	MN	C	1149	1/1	1.00	0.13	15,15,15,15	0
2	MN	A	1145	1/1	1.00	0.10	17,17,17,17	0
2	MN	D	1148	1/1	1.00	0.12	14,14,14,14	0
2	MN	C	1148	1/1	1.00	0.13	8,8,8,8	1

6.5 Other polymers

There are no such residues in this entry.