



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:41 AM UTC

PDB ID : 8QAM / pdb_00008qam
Title : vaccinia virus Uracil DNA glycosidase mutant I197K-V200E-L204K
Authors : Tarbouriech, N.; Burmeister, W.P.
Deposited on : 2023-08-23
Resolution : 1.32 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

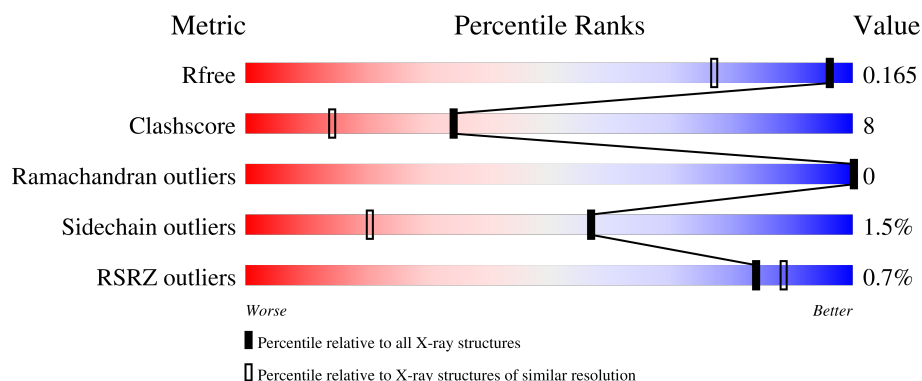
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.32 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}	180053	2531 (1.34-1.30)
Clashscore	190562	2585 (1.34-1.30)
Ramachandran outliers	187476	2528 (1.34-1.30)
Sidechain outliers	187428	2528 (1.34-1.30)
RSRZ outliers	180081	2528 (1.34-1.30)

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 $\leq 5\%$. 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	B	223	
1	C	223	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4648 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 Uracil-DNA glycosylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	218	Total	C	N	O	S	0	21	0
			1920	1245	316	353	6			
1	C	218	Total	C	N	O	S	0	26	0
			1934	1259	313	356	6			

There are 16 discrepancies between the modelled and reference sequences:

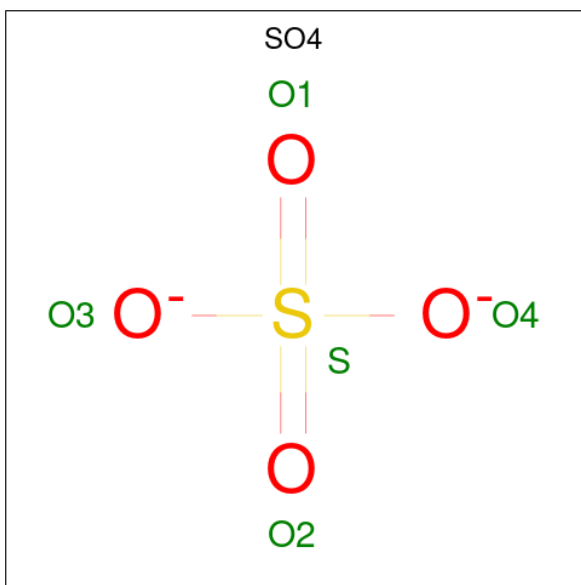
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	expression tag	UNP P20536
B	-3	ALA	-	expression tag	UNP P20536
B	-2	MET	-	expression tag	UNP P20536
B	-1	GLY	-	expression tag	UNP P20536
B	0	SER	-	expression tag	UNP P20536
B	197	LYS	ILE	engineered mutation	UNP P20536
B	200	GLU	VAL	engineered mutation	UNP P20536
B	204	LYS	LEU	engineered mutation	UNP P20536
C	-4	GLY	-	expression tag	UNP P20536
C	-3	ALA	-	expression tag	UNP P20536
C	-2	MET	-	expression tag	UNP P20536
C	-1	GLY	-	expression tag	UNP P20536
C	0	SER	-	expression tag	UNP P20536
C	197	LYS	ILE	engineered mutation	UNP P20536
C	200	GLU	VAL	engineered mutation	UNP P20536
C	204	LYS	LEU	engineered mutation	UNP P20536

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
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		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	353	Total	O	0	0
			353	353		
4	C	385	Total	O	0	0
			385	385		

4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	161.67Å 161.67Å 39.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.20 – 1.32 38.20 – 1.32	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.20-1.32) 99.8 (38.20-1.32)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 1.32Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.130 , 0.165 0.130 , 0.165	Depositor DCC
R_{free} test set	6094 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	18.8	Xtriage
Anisotropy	0.123	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 58.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	4648	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 3.78% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 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	B	1.04	1/2017 (0.0%)	1.15	6/2729 (0.2%)
1	C	1.13	2/2037 (0.1%)	1.21	11/2756 (0.4%)
All	All	1.09	3/4054 (0.1%)	1.18	17/5485 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	151	HIS	CE1-NE2	-6.61	1.25	1.32
1	C	147	HIS	CE1-NE2	6.19	1.38	1.32
1	B	8	HIS	CG-ND1	-5.06	1.32	1.38

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	58	ARG	NE-CZ-NH2	7.66	126.09	119.20
1	C	58	ARG	CD-NE-CZ	7.20	134.49	124.40
1	C	68	ASP	CA-CB-CG	6.56	119.16	112.60
1	C	58	ARG	NE-CZ-NH1	-6.45	115.05	121.50
1	C	17	ASP	CA-CB-CG	6.43	119.03	112.60
1	B	17	ASP	CA-CB-CG	6.41	119.01	112.60
1	B	68	ASP	CB-CA-C	5.99	118.07	110.76
1	C	146[A]	GLN	CB-CG-CD	-5.86	102.63	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	146[B]	GLN	CB-CG-CD	-5.86	102.63	112.60
1	B	120	ASN	CA-CB-CG	5.66	118.26	112.60
1	C	138	ASP	CA-CB-CG	-5.48	107.12	112.60
1	C	205	ASP	CA-CB-CG	5.42	118.02	112.60
1	B	189	PHE	CA-C-O	-5.30	114.93	120.55
1	C	160	LYS	CB-CA-C	-5.28	102.59	110.88
1	C	120	ASN	OD1-CG-ND2	5.27	127.87	122.60
1	B	68	ASP	CA-CB-CG	5.25	117.85	112.60
1	B	111	ASN	OD1-CG-ND2	5.04	127.64	122.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	165	ASN	Sidechain

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	B	1920	0	1954	36	0
1	C	1934	0	1991	22	0
2	B	18	0	24	4	0
2	C	18	0	24	3	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
4	B	353	0	0	17	1
4	C	385	0	0	16	1
All	All	4648	0	3993	60	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 8.

All (60) 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:167[C]:ARG:NH2	4:B:401:HOH:O	1.77	1.16
2:C:1001:GOL:H2	4:C:1105:HOH:O	1.50	1.12
1:B:52:GLN:NE2	4:B:402:HOH:O	1.88	0.98
2:C:1001:GOL:C2	4:C:1105:HOH:O	2.08	0.92
1:B:181:HIS:HD2	1:B:183:ALA:H	1.20	0.90
1:B:114[B]:ASP:OD2	4:B:403:HOH:O	1.95	0.81
1:C:16:HIS:HD2	1:C:18:ASP:H	1.27	0.80
1:B:138[B]:ASP:OD1	4:B:404:HOH:O	2.02	0.78
1:B:117:ILE:HD12	4:B:546:HOH:O	1.83	0.77
1:B:162:ASP:OD2	4:B:405:HOH:O	2.04	0.75
1:B:117:ILE:HD11	4:B:450:HOH:O	1.90	0.69
1:C:55:GLN:OE1	4:C:1101:HOH:O	2.11	0.68
1:B:167[C]:ARG:CZ	4:B:401:HOH:O	2.22	0.68
1:C:91[A]:GLU:OE1	4:C:1102:HOH:O	2.12	0.67
1:C:1:MET:N	4:C:1111:HOH:O	2.28	0.67
1:B:67:ILE:HD11	2:B:303:GOL:H32	1.77	0.67
1:C:16:HIS:CD2	1:C:18:ASP:H	2.11	0.66
1:C:138:ASP:HB2	4:C:1130:HOH:O	1.94	0.66
1:B:55[A]:GLN:HE21	1:B:55[A]:GLN:HA	1.61	0.65
1:B:55[A]:GLN:HA	1:B:55[A]:GLN:NE2	2.13	0.64
1:B:39[A]:ARG:HH12	1:C:207[A]:LYS:HE2	1.63	0.63
1:B:39[A]:ARG:NH1	1:C:207[A]:LYS:HE2	2.15	0.61
1:B:117:ILE:HG13	4:B:402:HOH:O	2.02	0.60
1:C:16:HIS:HE1	4:C:1407:HOH:O	1.86	0.58
1:C:5[A]:THR:HG23	4:C:1233:HOH:O	2.03	0.57
1:C:8:HIS:HD2	4:C:1193:HOH:O	1.87	0.56
1:B:117:ILE:HG13	4:B:671:HOH:O	2.07	0.55
1:C:14[B]:THR:OG1	1:C:54:LYS:HE3	2.08	0.54
1:B:33[A]:VAL:HG21	4:B:696:HOH:O	2.09	0.53
1:C:1:MET:CA	4:C:1111:HOH:O	2.57	0.53
1:C:1:MET:HA	4:C:1111:HOH:O	2.07	0.53
1:B:181:HIS:CD2	1:B:183:ALA:H	2.12	0.52
1:B:104:ASP:HB2	4:B:406:HOH:O	2.09	0.52
1:B:29:PHE:O	1:B:32[C]:GLU:HG2	2.09	0.52
1:B:117:ILE:CD1	4:B:450:HOH:O	2.51	0.51
2:B:302:GOL:H12	4:B:418:HOH:O	2.11	0.50
1:C:68:ASP:HB2	1:C:69:PRO:CD	2.42	0.49
1:C:164:SER:HA	4:C:1165:HOH:O	2.14	0.48
1:C:8:HIS:CD2	4:C:1193:HOH:O	2.65	0.47
1:C:142:LYS:HD2	1:C:169:LYS:HE2	1.95	0.47
1:B:33[A]:VAL:HG23	4:B:599:HOH:O	2.15	0.47
1:B:2:ASN:HB2	1:B:15:TYR:CZ	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:127:LEU:CD1	4:C:1404:HOH:O	2.63	0.46
1:B:32[A]:GLU:HG3	1:B:33[A]:VAL:HG23	1.96	0.46
1:B:48:LYS:HA	1:B:51[B]:ILE:HG12	1.98	0.46
1:B:167[C]:ARG:HD2	1:B:167[C]:ARG:HA	1.56	0.45
1:C:94[B]:SER:OG	1:C:98:ARG:NH2	2.50	0.45
1:B:79:PHE:CE1	2:B:301:GOL:H31	2.51	0.45
1:B:1:MET:N	4:B:415:HOH:O	2.50	0.45
1:B:197[A]:LYS:HG3	4:B:411:HOH:O	2.16	0.44
1:B:122:TYR:OH	2:B:303:GOL:H31	2.18	0.44
2:C:1002:GOL:H12	4:C:1199:HOH:O	2.18	0.44
1:B:185:ARG:H	1:B:185:ARG:NH1	2.15	0.43
1:B:142[B]:LYS:HE3	1:B:146:GLN:HE22	1.84	0.43
1:B:68:ASP:HB2	1:B:69:PRO:CD	2.50	0.42
1:C:1:MET:N	4:C:1131:HOH:O	2.52	0.42
1:B:32[B]:GLU:OE2	1:B:139[B]:LYS:NZ	2.41	0.41
1:C:35:SER:O	1:C:39:ARG:HG3	2.21	0.40
1:B:186:ASP:OD1	1:B:186:ASP:N	2.54	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)
4:B:628:HOH:O	4:C:1122:HOH:O[7_454]	2.09	0.11

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	B	237/223 (106%)	230 (97%)	7 (3%)	0	100	100
1	C	242/223 (108%)	239 (99%)	3 (1%)	0	100	100
All	All	479/446 (107%)	469 (98%)	10 (2%)	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	B	222/202 (110%)	218 (98%)	4 (2%)	51	15
1	C	226/202 (112%)	223 (99%)	3 (1%)	61	24
All	All	448/404 (111%)	441 (98%)	7 (2%)	57	18

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	55[A]	GLN
1	B	55[B]	GLN
1	B	185	ARG
1	B	186	ASP
1	C	127	LEU
1	C	169	LYS
1	C	185	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	146	GLN
1	B	181	HIS
1	B	188	GLN
1	C	8	HIS
1	C	16	HIS
1	C	151	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	B	302	-	5,5,5	0.20	0	5,5,5	0.32	0
2	GOL	B	301	-	5,5,5	0.25	0	5,5,5	1.02	0
2	GOL	C	1003	-	5,5,5	0.19	0	5,5,5	0.17	0
2	GOL	C	1002	-	5,5,5	0.53	0	5,5,5	0.71	0
2	GOL	B	303	-	5,5,5	0.37	0	5,5,5	0.74	0
3	SO4	B	305	-	4,4,4	0.26	0	6,6,6	0.33	0
2	GOL	C	1001	-	5,5,5	0.51	0	5,5,5	0.94	0
3	SO4	B	304	-	4,4,4	0.67	0	6,6,6	0.59	0
3	SO4	C	1005	-	4,4,4	1.75	1 (25%)	6,6,6	0.71	0
3	SO4	C	1004	-	4,4,4	0.51	0	6,6,6	0.26	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	302	-	-	0/4/4/4	-
2	GOL	B	301	-	-	4/4/4/4	-
2	GOL	C	1003	-	-	0/4/4/4	-
2	GOL	C	1002	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	303	-	-	2/4/4/4	-
2	GOL	C	1001	-	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1005	SO4	O1-S	2.84	1.61	1.44

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301	GOL	C1-C2-C3-O3
2	B	303	GOL	O1-C1-C2-C3
2	B	301	GOL	O1-C1-C2-C3
2	B	301	GOL	O1-C1-C2-O2
2	B	301	GOL	O2-C2-C3-O3
2	B	303	GOL	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	GOL	1	0
2	B	301	GOL	1	0
2	C	1002	GOL	1	0
2	B	303	GOL	2	0
2	C	1001	GOL	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 [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	B	218/223 (97%)	-0.16	3 (1%) 73 79	11, 18, 41, 113	21 (9%)
1	C	218/223 (97%)	-0.23	0 100 100	9, 17, 33, 72	26 (11%)
All	All	436/446 (97%)	-0.19	3 (0%) 84 88	9, 18, 36, 113	47 (10%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	ASP	2.6
1	B	184	ALA	2.3
1	B	187	ARG	2.1

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 oligosaccharides 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	B	303	6/6	0.86	0.14	42,50,58,62	0
2	GOL	C	1003	6/6	0.91	0.09	38,49,52,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	C	1001	6/6	0.93	0.10	30,34,43,49	0
3	SO4	C	1005	5/5	0.93	0.12	18,25,34,43	5
2	GOL	B	302	6/6	0.94	0.09	22,25,27,35	6
2	GOL	C	1002	6/6	0.94	0.10	18,24,29,41	6
3	SO4	B	305	5/5	0.96	0.08	26,33,39,62	5
2	GOL	B	301	6/6	0.96	0.08	15,38,47,50	0
3	SO4	B	304	5/5	0.97	0.08	29,32,46,46	5
3	SO4	C	1004	5/5	0.99	0.04	23,24,26,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.