



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:14 AM UTC

PDB ID : 5F9C / pdb_00005f9c
Title : Crystal structure of the G121R mutant of human phosphoglucomutase 1
Authors : Beamer, L.J.
Deposited on : 2015-12-09
Resolution : 2.50 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

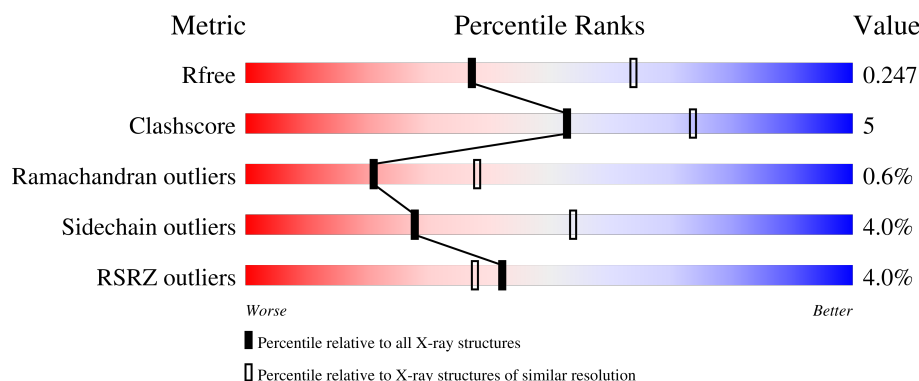
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	A	562	2% 84% 13% ..
1	B	562	6% 80% 14% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8196 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 Phosphoglucomutase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	550	Total	C	N	O	S	0	2	0
			4135	2633	695	791	16			
1	B	536	Total	C	N	O	S	0	0	0
			3922	2494	665	747	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	121	ARG	GLY	engineered mutation	UNP P36871
B	121	ARG	GLY	engineered mutation	UNP P36871

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: 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	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

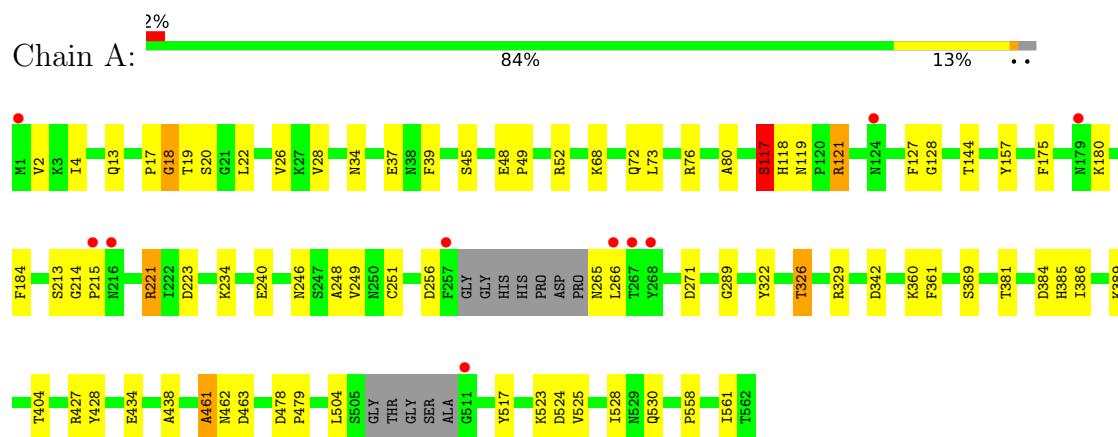
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	57	Total	O	0	0
			57	57		
5	B	23	Total	O	0	0
			23	23		

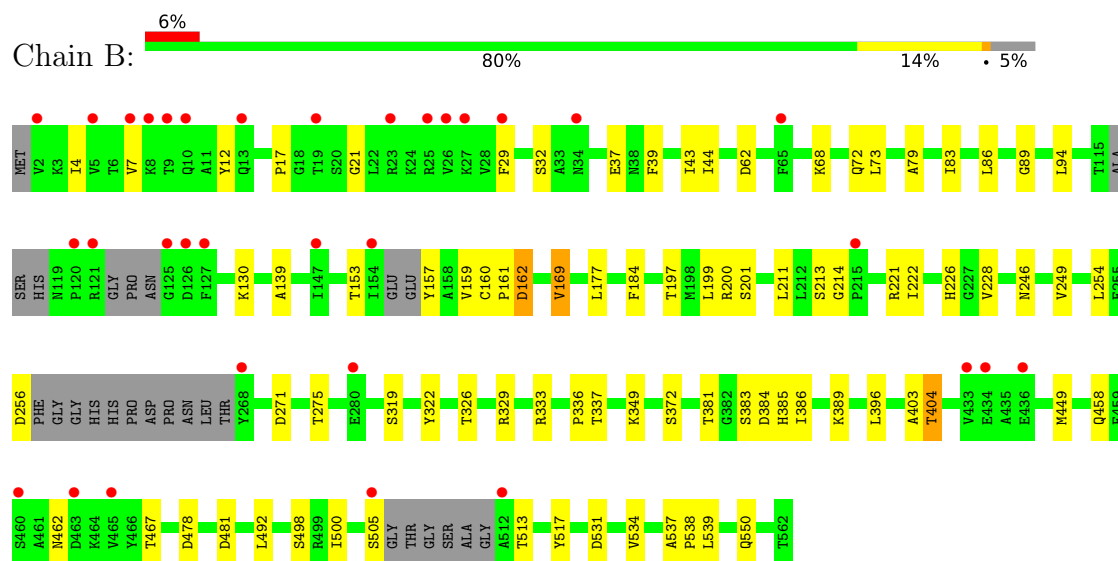
3 Residue-property plots [i](#)

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: Phosphoglucumutase-1



• Molecule 1: Phosphoglucumutase-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	171.98Å 171.98Å 99.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.94 – 2.50 60.94 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (60.94-2.50) 99.9 (60.94-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.192 , 0.243 0.197 , 0.247	Depositor DCC
R_{free} test set	3612 reflections (2.39%)	wwPDB-VP
Wilson B-factor (Å ²)	45.5	Xtriage
Anisotropy	0.439	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 78.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8196	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: MG, 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	A	0.57	0/4217	0.91	7/5719 (0.1%)
1	B	0.49	0/3990	0.87	6/5420 (0.1%)
All	All	0.53	0/8207	0.89	13/11139 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	214	GLY	CA-C-N	7.64	129.39	119.84
1	A	214	GLY	C-N-CA	7.64	129.39	119.84
1	A	461	ALA	N-CA-C	6.11	119.20	111.69
1	A	28	VAL	N-CA-C	-5.67	106.93	111.81
1	A	119	ASN	CA-C-N	5.61	125.55	119.78
1	A	119	ASN	C-N-CA	5.61	125.55	119.78
1	B	139	ALA	CA-C-N	5.58	125.57	120.21
1	B	139	ALA	C-N-CA	5.58	125.57	120.21
1	B	214	GLY	CA-C-N	5.55	126.78	119.84
1	B	214	GLY	C-N-CA	5.55	126.78	119.84
1	B	160	CYS	CA-C-N	5.22	125.27	119.32
1	B	160	CYS	C-N-CA	5.22	125.27	119.32
1	A	117	SER	CB-CA-C	-5.04	110.38	117.23

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	4135	0	4025	39	0
1	B	3922	0	3720	40	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	15	0	0	2	0
3	B	30	0	0	0	0
4	A	12	0	16	1	0
5	A	57	0	0	6	0
5	B	23	0	0	1	0
All	All	8196	0	7761	80	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 5.

All (80) 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:329:ARG:NH1	1:B:384:ASP:OD2	2.10	0.84
1:A:329:ARG:NH1	1:A:384:ASP:OD2	2.17	0.78
1:A:18:GLY:HA3	5:A:715:HOH:O	1.92	0.70
1:B:319:SER:HB2	1:B:404:THR:HG21	1.73	0.69
1:A:461:ALA:N	1:A:462:ASN:HA	2.07	0.69
3:A:603:SO4:O1	5:A:702:HOH:O	2.12	0.66
1:A:265:ASN:OD1	5:A:703:HOH:O	2.14	0.66
1:A:17:PRO:HG2	1:A:144:THR:HB	1.79	0.65
1:A:121:ARG:HB2	1:A:256:ASP:HA	1.79	0.64
1:A:234:LYS:HE3	1:A:251:CYS:HB2	1.82	0.60
1:B:458:GLN:HG2	1:B:467:THR:HG22	1.83	0.60
1:B:79:ALA:HB1	1:B:177:LEU:HD11	1.84	0.60
1:B:221:ARG:HH11	1:B:275:THR:HG21	1.69	0.58
1:A:34:ASN:OD1	1:A:157:TYR:OH	2.08	0.57
1:A:26:VAL:HG22	1:A:127:PHE:HB2	1.86	0.56
1:A:256:ASP:N	1:A:256:ASP:OD1	2.38	0.56
1:B:385:HIS:CD2	1:B:386:ILE:HG12	2.39	0.56
1:A:68:LYS:O	1:A:72:GLN:HG2	2.08	0.54
1:A:461:ALA:HB3	1:A:462:ASN:O	2.08	0.54
1:B:505:SER:HB2	1:B:513:THR:HB	1.90	0.53
1:A:322:TYR:O	1:A:326:THR:HG23	2.09	0.53
1:A:240:GLU:OE1	5:A:704:HOH:O	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:ILE:HG23	1:A:80:ALA:HB1	1.92	0.51
1:B:62:ASP:HA	1:B:228:VAL:HB	1.92	0.50
1:A:19:THR:N	5:A:715:HOH:O	2.45	0.50
1:B:44:ILE:HG23	1:B:83:ILE:HD11	1.94	0.49
1:A:524:ASP:O	1:A:528:ILE:HG23	2.12	0.49
1:B:322:TYR:O	1:B:326:THR:HG23	2.13	0.48
1:B:162:ASP:OD2	1:B:162:ASP:N	2.39	0.48
1:B:37:GLU:HG3	1:B:73:LEU:HD11	1.95	0.48
1:B:68:LYS:O	1:B:72:GLN:HG2	2.14	0.47
1:A:369:SER:HA	5:A:719:HOH:O	2.15	0.47
1:B:500:ILE:HD11	1:B:539:LEU:HD22	1.97	0.47
1:A:360:LYS:HG3	1:A:361:PHE:N	2.30	0.46
1:A:248:ALA:HB1	1:A:251:CYS:SG	2.57	0.45
1:B:211:LEU:HD22	1:B:403:ALA:HB2	1.97	0.45
1:B:385:HIS:NE2	1:B:386:ILE:HG12	2.30	0.45
1:B:7:VAL:O	1:B:157:TYR:N	2.49	0.45
1:B:159:VAL:O	1:B:161:PRO:HD3	2.17	0.44
1:B:86:LEU:HG	1:B:184:PHE:HZ	1.83	0.44
1:A:117:SER:HB3	1:A:118:HIS:H	1.56	0.44
1:B:12:TYR:OH	1:B:32:SER:HB3	2.18	0.44
1:A:17:PRO:HB3	1:A:39:PHE:HZ	1.82	0.43
1:A:428:TYR:CD2	1:A:558:PRO:HG3	2.53	0.43
1:A:117:SER:O	1:A:289:GLY:HA3	2.18	0.43
1:B:169:VAL:HA	5:B:713:HOH:O	2.19	0.43
1:A:561:ILE:CD1	4:A:606:GOL:H32	2.48	0.43
1:B:86:LEU:HG	1:B:184:PHE:CZ	2.54	0.43
1:B:79:ALA:HB2	1:B:184:PHE:CD2	2.54	0.43
1:B:226:HIS:HB3	1:B:254:LEU:HG	2.00	0.43
1:A:438:ALA:HB1	1:A:504:LEU:HD21	2.01	0.42
1:B:333:ARG:HD2	1:B:337:THR:HB	2.01	0.42
1:A:45:SER:HA	1:A:52:ARG:NH2	2.35	0.42
1:B:372:SER:O	1:B:383:SER:HA	2.20	0.42
1:A:73:LEU:HD23	1:A:76:ARG:NH1	2.35	0.42
1:A:523:LYS:HA	1:A:523:LYS:HD2	1.86	0.42
1:A:385:HIS:CD2	1:A:386:ILE:HG12	2.55	0.41
1:B:21:GLY:HA3	1:B:130:LYS:HA	2.01	0.41
1:B:498:SER:HB3	1:B:539:LEU:HD11	2.02	0.41
1:A:175:PHE:HB2	1:A:184:PHE:HB3	2.02	0.41
1:A:427:ARG:HD2	1:A:517:TYR:CE2	2.55	0.41
1:B:531:ASP:O	1:B:534:VAL:HG12	2.20	0.41
1:B:256:ASP:OD1	1:B:256:ASP:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:PHE:O	1:B:43:ILE:HG13	2.20	0.41
1:A:34:ASN:HA	1:A:37:GLU:OE1	2.21	0.41
1:A:434:GLU:O	1:A:434:GLU:HG3	2.20	0.41
1:B:89:GLY:HA3	1:B:94:LEU:HG	2.03	0.41
1:B:197:THR:HA	1:B:200:ARG:NH1	2.36	0.41
1:A:478:ASP:HA	1:A:479:PRO:HD3	1.83	0.41
1:B:449:MET:HG2	1:B:492:LEU:HD21	2.03	0.41
1:A:22:LEU:O	1:A:128:GLY:HA2	2.20	0.40
1:B:539:LEU:HA	1:B:539:LEU:HD23	1.88	0.40
1:A:48:GLU:HA	1:A:49:PRO:HD3	1.94	0.40
1:B:199:LEU:HD13	1:B:396:LEU:HD22	2.02	0.40
1:B:221:ARG:C	1:B:222:ILE:HD12	2.46	0.40
1:B:537:ALA:HB3	1:B:538:PRO:HD3	2.03	0.40
1:A:221:ARG:NH1	3:A:604:SO4:O3	2.55	0.40
1:A:223:ASP:HA	1:A:249:VAL:HG23	2.03	0.40
1:B:478:ASP:HB3	1:B:481:ASP:OD1	2.21	0.40
1:B:336:PRO:HB2	1:B:517:TYR:CE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	546/562 (97%)	524 (96%)	18 (3%)	4 (1%)	18	34
1	B	524/562 (93%)	501 (96%)	21 (4%)	2 (0%)	30	49
All	All	1070/1124 (95%)	1025 (96%)	39 (4%)	6 (1%)	21	38

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	GLY

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Mol	Chain	Res	Type
1	A	215	PRO
1	A	2	VAL
1	B	17	PRO
1	B	29	PHE
1	A	463	ASP

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	425/463 (92%)	408 (96%)	17 (4%)	28	54
1	B	385/463 (83%)	370 (96%)	15 (4%)	28	55
All	All	810/926 (88%)	778 (96%)	32 (4%)	28	54

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	20	SER
1	A	117	SER
1	A	121	ARG
1	A	180	LYS
1	A	213	SER
1	A	221	ARG
1	A	246	ASN
1	A	266	LEU
1	A	271	ASP
1	A	326	THR
1	A	342	ASP
1	A	381	THR
1	A	389	LYS
1	A	404	THR
1	A	525	VAL
1	A	530	GLN
1	B	4	ILE

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Mol	Chain	Res	Type
1	B	153	THR
1	B	162	ASP
1	B	169	VAL
1	B	201	SER
1	B	213	SER
1	B	246	ASN
1	B	249	VAL
1	B	271	ASP
1	B	349	LYS
1	B	381	THR
1	B	389	LYS
1	B	404	THR
1	B	462	ASN
1	B	550	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	226	HIS
1	A	246	ASN
1	A	294	ASN
1	A	300	HIS
1	A	423	ASN
1	B	294	ASN
1	B	423	ASN
1	B	439	ASN
1	B	488	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 13 ligands modelled in this entry, 2 are monoatomic - leaving 11 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	$\# Z > 2$
4	GOL	A	605	-	5,5,5	0.34	0	5,5,5	0.52	0
3	SO4	A	603	-	4,4,4	0.28	0	6,6,6	0.09	0
3	SO4	B	604	-	4,4,4	0.24	0	6,6,6	0.16	0
4	GOL	A	606	-	5,5,5	0.40	0	5,5,5	0.57	0
3	SO4	B	602	-	4,4,4	0.25	0	6,6,6	0.15	0
3	SO4	A	602	-	4,4,4	0.23	0	6,6,6	0.14	0
3	SO4	A	604	-	4,4,4	0.23	0	6,6,6	0.10	0
3	SO4	B	606	-	4,4,4	0.26	0	6,6,6	0.10	0
3	SO4	B	605	-	4,4,4	0.26	0	6,6,6	0.23	0
3	SO4	B	603	-	4,4,4	0.23	0	6,6,6	0.29	0
3	SO4	B	601	-	4,4,4	0.22	0	6,6,6	0.25	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
4	GOL	A	605	-	-	1/4/4/4	-
4	GOL	A	606	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	606	GOL	O1-C1-C2-C3
4	A	605	GOL	O1-C1-C2-O2
4	A	606	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	603	SO4	1	0
4	A	606	GOL	1	0
3	A	604	SO4	1	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	550/562 (97%)	-0.17	10 (1%)	67 64	31, 63, 100, 120	2 (0%)
1	B	536/562 (95%)	0.37	33 (6%)	26 23	44, 87, 139, 159	0
All	All	1086/1124 (96%)	0.09	43 (3%)	42 38	31, 73, 128, 159	2 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	127	PHE	4.8
1	B	9	THR	4.4
1	B	512	ALA	4.1
1	A	266	LEU	4.0
1	B	268	TYR	3.9
1	B	27	LYS	3.8
1	B	126	ASP	3.6
1	A	257	PHE	3.2
1	B	25	ARG	3.0
1	A	1	MET	3.0
1	A	179	ASN	2.9
1	B	147	ILE	2.9
1	B	436	GLU	2.8
1	B	154	ILE	2.8
1	B	34	ASN	2.8
1	B	8	LYS	2.8
1	B	13	GLN	2.8
1	B	121	ARG	2.8
1	B	26	VAL	2.7
1	B	463	ASP	2.7
1	A	268	TYR	2.7
1	B	434	GLU	2.7
1	B	125	GLY	2.6
1	B	215	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	215	PRO	2.5
1	B	120	PRO	2.5
1	B	280	GLU	2.5
1	B	2	VAL	2.5
1	B	10	GLN	2.4
1	A	511	GLY	2.4
1	A	216	ASN	2.4
1	A	124	ASN	2.3
1	B	505	SER	2.2
1	B	29	PHE	2.2
1	B	5	VAL	2.1
1	A	267	THR	2.1
1	B	465	VAL	2.1
1	B	7	VAL	2.1
1	B	433	VAL	2.1
1	B	23	ARG	2.1
1	B	65	PHE	2.0
1	B	19	THR	2.0
1	B	460	SER	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 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($\text{\AA}^2$)	Q<0.9
3	SO4	B	606	5/5	0.50	0.14	157,159,160,160	0
2	MG	B	607	1/1	0.57	0.19	105,105,105,105	0
3	SO4	A	603	5/5	0.61	0.18	151,151,153,155	0
3	SO4	B	604	5/5	0.70	0.14	107,108,110,111	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	A	606	6/6	0.81	0.22	57,70,72,73	6
3	SO4	B	603	5/5	0.82	0.14	93,94,96,97	5
3	SO4	A	604	5/5	0.87	0.14	115,115,117,118	5
4	GOL	A	605	6/6	0.89	0.15	68,80,84,87	0
3	SO4	B	602	5/5	0.89	0.16	94,95,98,99	5
3	SO4	B	601	5/5	0.90	0.09	89,96,97,99	5
3	SO4	B	605	5/5	0.91	0.10	85,85,89,93	5
3	SO4	A	602	5/5	0.94	0.09	69,69,72,75	0
2	MG	A	601	1/1	0.97	0.05	67,67,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.