



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:39 PM UTC

PDB ID : 1U0F / pdb_00001u0f
Title : Crystal structure of mouse phosphoglucose isomerase in complex with glucose 6-phosphate
Authors : Solomons, J.T.G.; Zimmerly, E.M.; Burns, S.; Krishnamurthy, N.; Swan, M.K.; Krings, S.; Muirhead, H.; Chirgwin, J.; Davies, C.
Deposited on : 2004-07-13
Resolution : 1.60 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

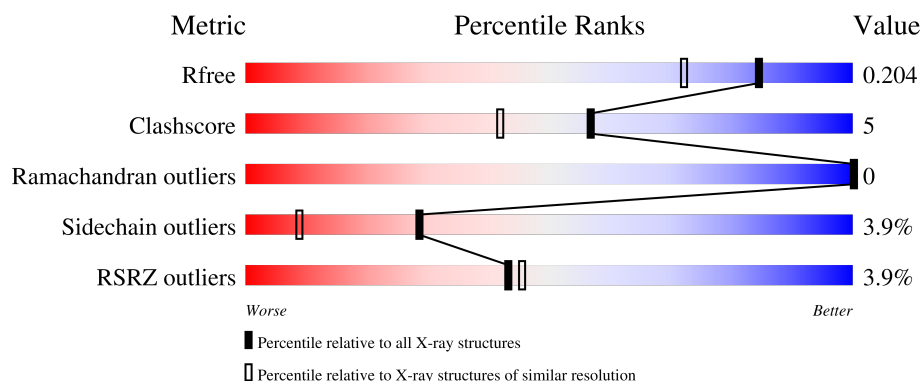
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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 $\leq 5\%$. 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	564	2% 88% 10% .
1	B	564	5% 86% 11% ..

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	SO4	B	1008	-	-	X	-
5	GOL	A	1026	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9903 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.

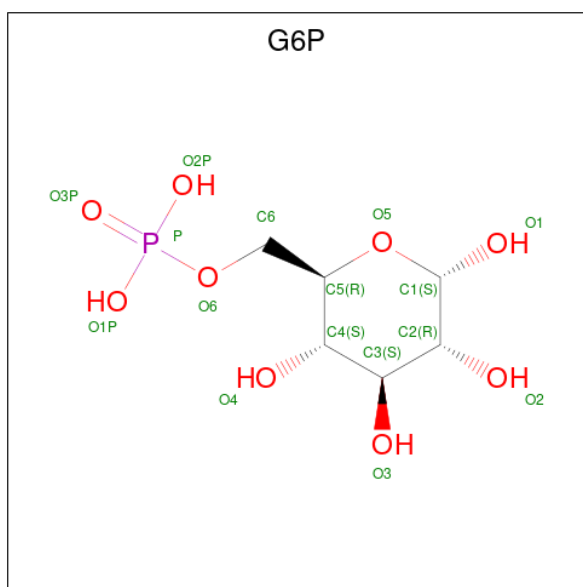
- Molecule 1 is a protein called Glucose-6-phosphate isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	556	Total	C	N	O	S	0	15	0
			4450	2843	769	819	19			
1	B	556	Total	C	N	O	S	0	13	0
			4446	2842	770	815	19			

There are 14 discrepancies between the modelled and reference sequences:

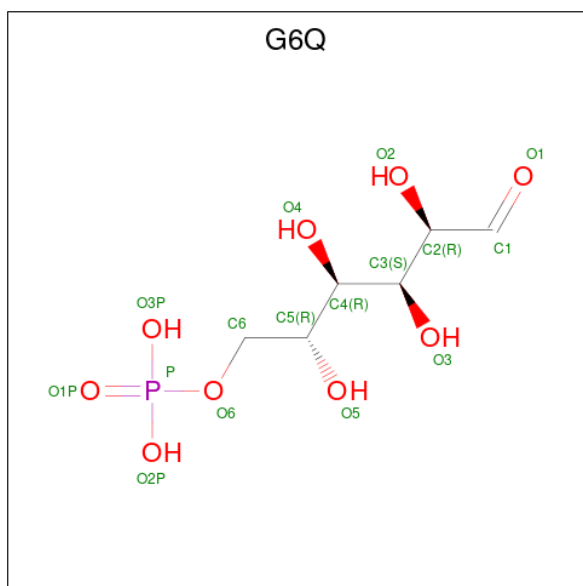
Chain	Residue	Modelled	Actual	Comment	Reference
A	263	LEU	PHE	SEE REMARK 999	UNP P06745
A	558	HIS	-	expression tag	UNP P06745
A	559	HIS	-	expression tag	UNP P06745
A	560	HIS	-	expression tag	UNP P06745
A	561	HIS	-	expression tag	UNP P06745
A	562	HIS	-	expression tag	UNP P06745
A	563	HIS	-	expression tag	UNP P06745
B	263	LEU	PHE	SEE REMARK 999	UNP P06745
B	558	HIS	-	expression tag	UNP P06745
B	559	HIS	-	expression tag	UNP P06745
B	560	HIS	-	expression tag	UNP P06745
B	561	HIS	-	expression tag	UNP P06745
B	562	HIS	-	expression tag	UNP P06745
B	563	HIS	-	expression tag	UNP P06745

- Molecule 2 is 6-O-phosphono-alpha-D-glucopyranose (CCD ID: G6P) (formula: C₆H₁₃O₉P).



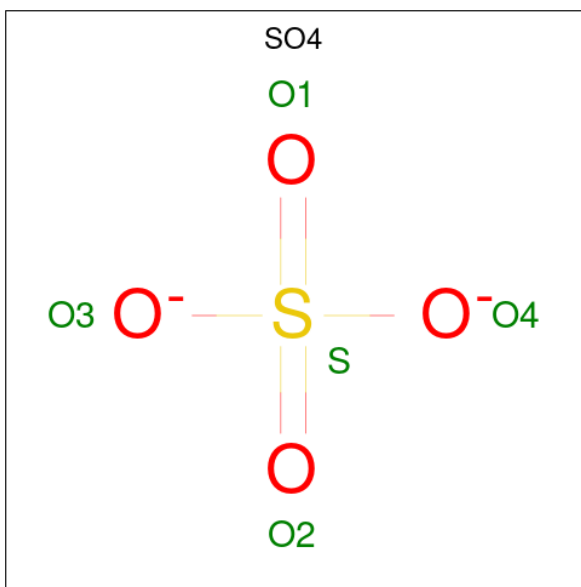
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	1
			17	6	10	1		

- Molecule 3 is GLUCOSE-6-PHOSPHATE (CCD ID: G6Q) (formula: $C_6H_{13}O_9P$).



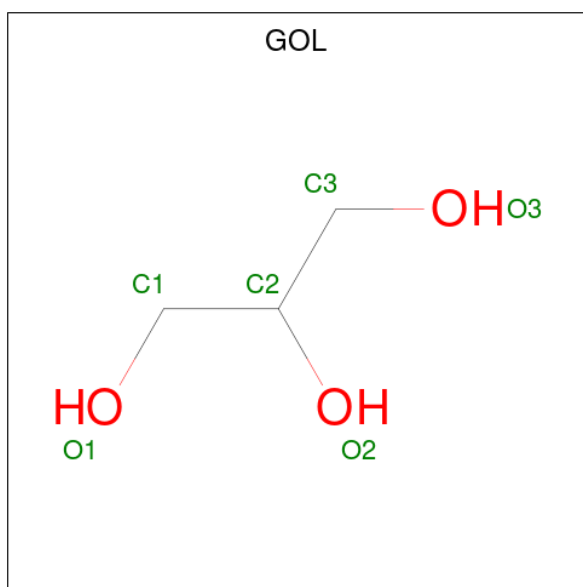
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	1
			16	6	9	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



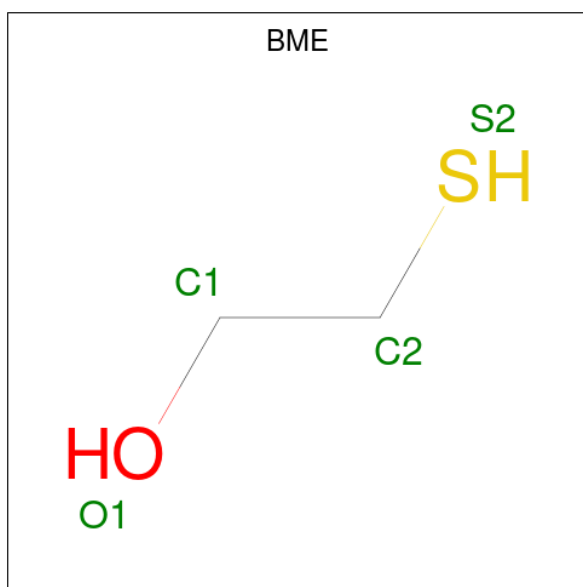
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 6 is BETA-MERCAPTOETHANOL (CCD ID: BME) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	O	S	0	0
			4	2	1	1		

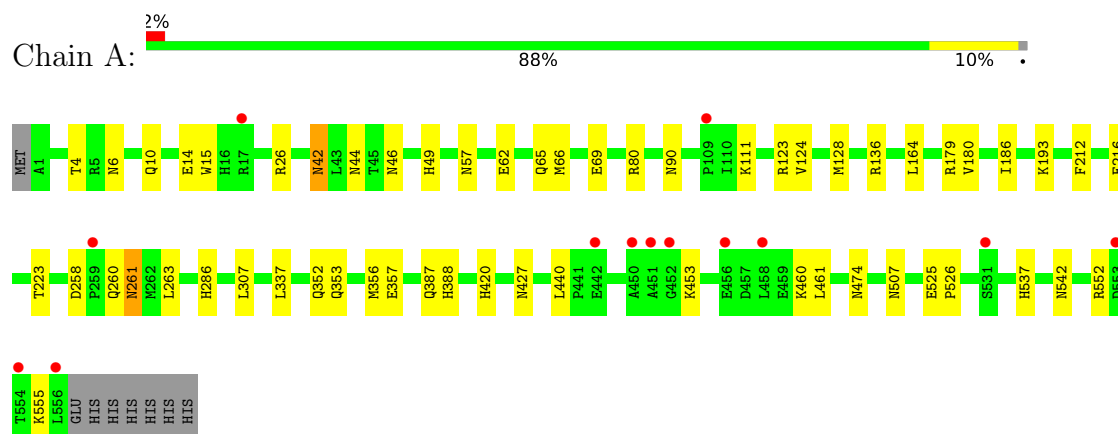
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	441	Total	O	0	1
			441	441		
7	B	417	Total	O	0	1
			417	417		

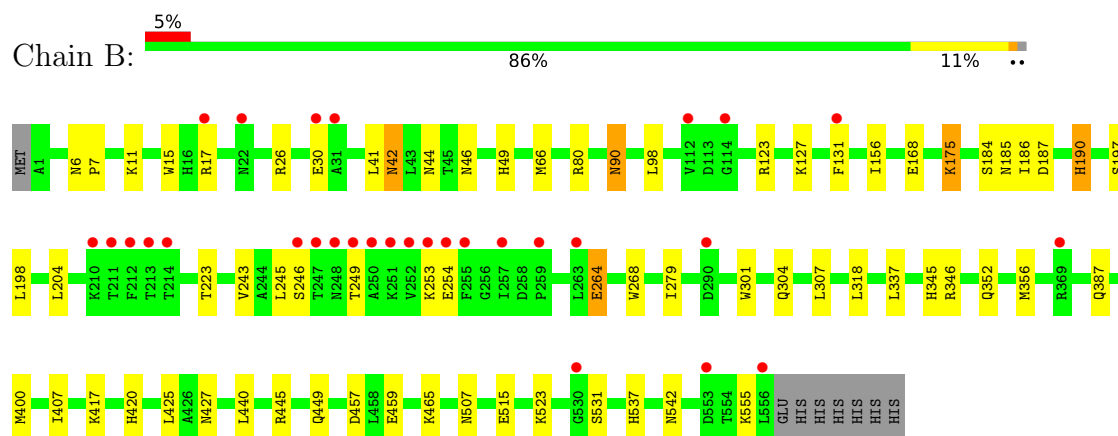
3 Residue-property plots [i](#)

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: Glucose-6-phosphate isomerase



- Molecule 1: Glucose-6-phosphate isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.40Å 116.80Å 73.90Å 90.00° 101.60° 90.00°	Depositor
Resolution (Å)	45.00 – 1.60 45.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	94.6 (45.00-1.60) 94.7 (45.00-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.21 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.178 , 0.207 (Not available) , 0.204	Depositor DCC
R_{free} test set	7210 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	18.1	Xtriage
Anisotropy	0.609	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9903	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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: G6P, SO4, GOL, BME, G6Q

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	0.62	0/4631	0.81	0/6269
1	B	0.58	1/4617 (0.0%)	0.80	0/6247
All	All	0.60	1/9248 (0.0%)	0.80	0/12516

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	400	MET	SD-CE	-6.11	1.64	1.79

There are no bond angle outliers.

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	4450	0	4434	40	0
1	B	4446	0	4438	41	0
2	A	17	0	12	0	0
3	A	16	0	11	2	0
4	A	10	0	0	0	0
4	B	30	0	0	2	0
5	A	54	0	72	14	0
5	B	18	0	24	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	4	0	6	0	0
7	A	441	0	0	5	0
7	B	417	0	0	10	0
All	All	9903	0	8997	84	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 (84) 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:A:136:ARG:HE	5:A:1026:GOL:H32	1.07	1.09
1:B:449[A]:GLN:NE2	7:B:1383:HOH:O	1.92	1.00
5:B:1029:GOL:H32	7:B:1264:HOH:O	1.77	0.85
1:B:449[B]:GLN:NE2	7:B:1337:HOH:O	2.11	0.83
1:A:136:ARG:NE	5:A:1026:GOL:H32	1.92	0.80
1:B:301:TRP:O	1:B:304[B]:GLN:HG3	1.81	0.79
1:B:531:SER:H	5:B:1029:GOL:H12	1.51	0.74
1:B:449[B]:GLN:NE2	7:B:1383:HOH:O	2.21	0.74
1:B:537:HIS:H	1:B:542:ASN:HD21	1.36	0.73
1:A:537:HIS:H	1:A:542:ASN:HD21	1.38	0.70
4:B:1008:SO4:O2	7:B:1266:HOH:O	2.08	0.69
1:A:42:ASN:HD21	1:A:49:HIS:HD2	1.40	0.67
4:B:1008:SO4:O4	7:B:1342:HOH:O	2.13	0.67
1:A:136:ARG:HH21	5:A:1026:GOL:H12	1.60	0.65
1:B:185:ASN:H	1:B:190:HIS:HD2	1.42	0.65
1:B:445:ARG:O	1:B:449[B]:GLN:HG3	1.98	0.64
1:A:136:ARG:HE	5:A:1026:GOL:C3	1.99	0.63
1:A:179:ARG:HG2	5:A:1030:GOL:H11	1.84	0.59
1:B:345:HIS:HD2	7:B:1139:HOH:O	1.85	0.59
1:A:42:ASN:ND2	1:A:49:HIS:HD2	2.00	0.59
1:A:387:GLN:HE22	1:A:427:ASN:HB3	1.68	0.58
1:B:168:GLU:OE2	1:B:345:HIS:HE1	1.87	0.58
1:A:57:ASN:HD21	1:A:474:ASN:ND2	2.02	0.57
5:A:1026:GOL:O2	7:A:1207:HOH:O	2.16	0.57
1:B:175:LYS:HB2	7:B:1150:HOH:O	2.05	0.56
1:A:42:ASN:HD21	1:A:49:HIS:CD2	2.22	0.56
1:A:42:ASN:HD22	1:A:42:ASN:C	2.13	0.56
1:A:261:ASN:ND2	5:A:1027:GOL:O1	2.39	0.55
1:A:388:HIS:CE1	1:B:156:ILE:HD12	2.43	0.54
1:A:223:THR:OG1	1:B:420:HIS:HE1	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:420:HIS:HE1	1:B:223:THR:OG1	1.91	0.52
1:A:6:ASN:O	1:A:10[B]:GLN:HG3	2.09	0.52
7:A:1323:HOH:O	1:B:457:ASP:HB3	2.09	0.52
1:B:184:SER:H	1:B:190:HIS:CD2	2.28	0.52
5:A:1032:GOL:H32	7:A:1323:HOH:O	2.10	0.51
1:B:131:PHE:CZ	1:B:243:VAL:HG11	2.45	0.51
1:B:42:ASN:HD22	1:B:42:ASN:C	2.20	0.50
1:A:90:ASN:ND2	1:A:507:ASN:HD21	2.10	0.49
1:A:186:ILE:HB	1:A:216:GLU:HG3	1.94	0.49
1:A:186:ILE:O	1:B:420:HIS:HD2	1.96	0.48
1:A:525:GLU:HB2	1:A:526:PRO:HD3	1.96	0.48
1:A:357:GLU:OE2	3:A:901[B]:G6Q:H2	2.13	0.48
1:A:164[B]:LEU:HD11	5:A:1025:GOL:H31	1.94	0.47
1:B:44:ASN:ND2	1:B:46:ASN:H	2.12	0.47
1:B:80:ARG:HD2	1:B:307:LEU:HA	1.95	0.47
1:B:245:LEU:HD13	1:B:279:ILE:HA	1.96	0.47
1:A:15:TRP:CG	1:A:66:MET:HE1	2.49	0.47
1:B:42:ASN:ND2	1:B:49:HIS:HD2	2.12	0.47
1:A:353:GLN:OE1	3:A:901[B]:G6Q:H1	2.15	0.47
1:B:352:GLN:O	1:B:356:MET:HB2	2.14	0.47
1:A:62[A]:GLU:HG2	1:A:66:MET:HE2	1.97	0.46
1:B:42:ASN:HD21	1:B:49:HIS:CD2	2.33	0.46
1:A:258:ASP:OD2	1:A:260:GLN:HG2	2.16	0.46
1:A:65:GLN:O	1:A:69:GLU:HG2	2.16	0.46
1:A:44:ASN:ND2	1:A:46:ASN:H	2.14	0.46
1:B:346:ARG:HD3	7:B:1273:HOH:O	2.16	0.46
1:B:387:GLN:HE22	1:B:427:ASN:HB3	1.79	0.46
1:A:164[B]:LEU:CD1	5:A:1025:GOL:H31	2.47	0.45
1:A:420:HIS:HD2	1:B:186:ILE:O	1.99	0.45
1:B:42:ASN:HD21	1:B:49:HIS:HD2	1.62	0.45
1:B:407:ILE:HD13	1:B:425:LEU:HD23	1.98	0.45
1:A:111:LYS:HE2	7:A:1423:HOH:O	2.17	0.45
1:A:128:MET:HG2	1:A:263:LEU:HD13	1.99	0.44
5:A:1032:GOL:C3	7:A:1323:HOH:O	2.65	0.44
1:B:90:ASN:HD22	1:B:90:ASN:C	2.25	0.44
1:B:98:LEU:HB2	1:B:268:TRP:CE3	2.53	0.44
1:A:90:ASN:HD22	1:A:507:ASN:HD21	1.66	0.43
1:A:180[B]:VAL:HG22	1:A:286:HIS:CE1	2.53	0.43
1:B:44:ASN:C	1:B:44:ASN:HD22	2.24	0.43
1:A:352:GLN:O	1:A:356:MET:HB2	2.18	0.43
1:A:90:ASN:HD22	1:A:90:ASN:C	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:ARG:HH22	1:B:264:GLU:CD	2.27	0.42
1:A:124:VAL:O	1:A:128:MET:HG3	2.19	0.42
1:B:41:LEU:HD11	1:B:318:LEU:HD12	2.01	0.42
1:B:6:ASN:HA	1:B:7:PRO:HD3	1.93	0.42
1:B:246:SER:C	7:B:1446:HOH:O	2.61	0.41
1:A:261:ASN:CG	5:A:1027:GOL:O1	2.63	0.41
1:B:15:TRP:CG	1:B:66:MET:HE1	2.56	0.41
1:B:90:ASN:HD22	1:B:507:ASN:HD21	1.69	0.41
1:A:80:ARG:HD2	1:A:307:LEU:HA	2.02	0.41
1:B:249:THR:HG22	1:B:253:LYS:HE2	2.03	0.41
1:B:186:ILE:O	1:B:187:ASP:C	2.64	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	569/564 (101%)	557 (98%)	12 (2%)	0	100	100
1	B	567/564 (100%)	550 (97%)	17 (3%)	0	100	100
All	All	1136/1128 (101%)	1107 (97%)	29 (3%)	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	489/482 (102%)	474 (97%)	15 (3%)	35	13
1	B	487/482 (101%)	463 (95%)	24 (5%)	22	5
All	All	976/964 (101%)	937 (96%)	39 (4%)	28	8

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	14	GLU
1	A	26	ARG
1	A	42	ASN
1	A	123	ARG
1	A	193	LYS
1	A	212	PHE
1	A	261	ASN
1	A	337	LEU
1	A	440	LEU
1	A	453	LYS
1	A	460	LYS
1	A	461	LEU
1	A	552	ARG
1	A	555	LYS
1	B	11	LYS
1	B	17	ARG
1	B	26	ARG
1	B	30	GLU
1	B	42	ASN
1	B	90	ASN
1	B	127	LYS
1	B	175	LYS
1	B	190	HIS
1	B	197	SER
1	B	198	LEU
1	B	204	LEU
1	B	254	GLU
1	B	264	GLU
1	B	337	LEU
1	B	417	LYS
1	B	440	LEU
1	B	459	GLU
1	B	465[A]	LYS
1	B	465[B]	LYS

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Mol	Chain	Res	Type
1	B	515[A]	GLU
1	B	515[B]	GLU
1	B	523	LYS
1	B	555	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	37	ASN
1	A	38	ASN
1	A	42	ASN
1	A	44	ASN
1	A	49	HIS
1	A	57	ASN
1	A	65	GLN
1	A	82	ASN
1	A	90	ASN
1	A	104	ASN
1	A	107	ASN
1	A	122	ASN
1	A	153	ASN
1	A	261	ASN
1	A	304	GLN
1	A	353	GLN
1	A	387	GLN
1	A	420	HIS
1	A	449	GLN
1	A	474	ASN
1	A	542	ASN
1	B	19	ASN
1	B	22	ASN
1	B	42	ASN
1	B	44	ASN
1	B	49	HIS
1	B	57	ASN
1	B	82	ASN
1	B	90	ASN
1	B	107	ASN
1	B	122	ASN
1	B	153	ASN
1	B	190	HIS

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Mol	Chain	Res	Type
1	B	215	GLN
1	B	335	HIS
1	B	345	HIS
1	B	353	GLN
1	B	373	GLN
1	B	385	ASN
1	B	420	HIS
1	B	542	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 24 ligands modelled in this entry, 1 is modelled with single atom - leaving 23 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	$\# Z > 2$
5	GOL	B	1028	-	5,5,5	0.34	0	5,5,5	0.34	0
4	SO4	B	1004	-	4,4,4	0.25	0	6,6,6	0.13	0
2	G6P	A	900[A]	-	16,16,16	1.01	1 (6%)	23,24,24	2.85	9 (39%)
5	GOL	A	1023	-	5,5,5	0.44	0	5,5,5	0.57	0
5	GOL	A	1024	-	5,5,5	0.33	0	5,5,5	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	B	1007	-	4,4,4	0.25	0	6,6,6	0.11	0
5	GOL	A	1021	-	5,5,5	0.34	0	5,5,5	0.48	0
4	SO4	B	1006	-	4,4,4	0.25	0	6,6,6	0.12	0
6	BME	B	1011	-	3,3,3	0.32	0	2,2,2	0.14	0
4	SO4	B	1005	-	4,4,4	0.24	0	6,6,6	0.10	0
5	GOL	B	1029	-	5,5,5	0.35	0	5,5,5	0.45	0
3	G6Q	A	901[B]	-	14,15,15	1.63	1 (7%)	19,21,21	1.05	0
5	GOL	A	1030	-	5,5,5	0.43	0	5,5,5	0.36	0
4	SO4	A	1002	-	4,4,4	0.24	0	6,6,6	0.07	0
5	GOL	A	1031	-	5,5,5	0.56	0	5,5,5	0.67	0
4	SO4	B	1003	-	4,4,4	0.22	0	6,6,6	0.06	0
5	GOL	A	1026	-	5,5,5	0.33	0	5,5,5	0.54	0
4	SO4	B	1008	-	4,4,4	0.23	0	6,6,6	0.15	0
5	GOL	A	1027	-	5,5,5	0.38	0	5,5,5	0.44	0
5	GOL	A	1025	-	5,5,5	0.37	0	5,5,5	0.34	0
4	SO4	A	1001	-	4,4,4	0.26	0	6,6,6	0.24	0
5	GOL	A	1032	-	5,5,5	0.40	0	5,5,5	0.36	0
5	GOL	B	1022	-	5,5,5	0.33	0	5,5,5	0.33	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	GOL	A	1021	-	-	0/4/4/4	-
5	GOL	A	1027	-	-	0/4/4/4	-
5	GOL	B	1028	-	-	2/4/4/4	-
3	G6Q	A	901[B]	-	-	1/19/20/20	-
5	GOL	A	1025	-	-	2/4/4/4	-
6	BME	B	1011	-	-	1/1/1/1	-
5	GOL	B	1022	-	-	0/4/4/4	-
2	G6P	A	900[A]	-	-	1/6/26/26	0/1/1/1
5	GOL	A	1023	-	-	4/4/4/4	-
5	GOL	A	1030	-	-	2/4/4/4	-
5	GOL	A	1031	-	-	2/4/4/4	-
5	GOL	B	1029	-	-	1/4/4/4	-
5	GOL	A	1032	-	-	2/4/4/4	-
5	GOL	A	1024	-	-	2/4/4/4	-
5	GOL	A	1026	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901[B]	G6Q	O1-C1	5.82	1.42	1.20
2	A	900[A]	G6P	P-O3P	2.79	1.59	1.50

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900[A]	G6P	O5-C5-C4	6.50	121.42	109.70
2	A	900[A]	G6P	O5-C1-C2	5.15	119.36	110.30
2	A	900[A]	G6P	C1-C2-C3	5.14	120.85	110.36
2	A	900[A]	G6P	C3-C4-C5	4.45	118.30	110.23
2	A	900[A]	G6P	C1-O5-C5	4.03	121.45	113.65
2	A	900[A]	G6P	O2-C2-C3	-3.90	101.17	110.38
2	A	900[A]	G6P	C4-C3-C2	3.45	116.89	110.83
2	A	900[A]	G6P	O5-C5-C6	2.99	112.62	106.69
2	A	900[A]	G6P	O2-C2-C1	-2.69	103.03	109.25

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	900[A]	G6P	C4-C5-C6-O6
3	A	901[B]	G6Q	O1-C1-C2-C3
5	A	1023	GOL	C1-C2-C3-O3
5	A	1025	GOL	O1-C1-C2-C3
5	A	1031	GOL	O1-C1-C2-C3
5	A	1032	GOL	C1-C2-C3-O3
5	B	1028	GOL	C1-C2-C3-O3
6	B	1011	BME	O1-C1-C2-S2
5	A	1032	GOL	O2-C2-C3-O3
5	B	1028	GOL	O2-C2-C3-O3
5	A	1023	GOL	O1-C1-C2-C3
5	A	1024	GOL	O1-C1-C2-C3
5	A	1030	GOL	O1-C1-C2-C3
5	A	1023	GOL	O2-C2-C3-O3
5	A	1024	GOL	O1-C1-C2-O2
5	A	1025	GOL	O1-C1-C2-O2
5	A	1031	GOL	O1-C1-C2-O2
5	A	1030	GOL	O1-C1-C2-O2
5	B	1029	GOL	O2-C2-C3-O3
5	A	1023	GOL	O1-C1-C2-O2

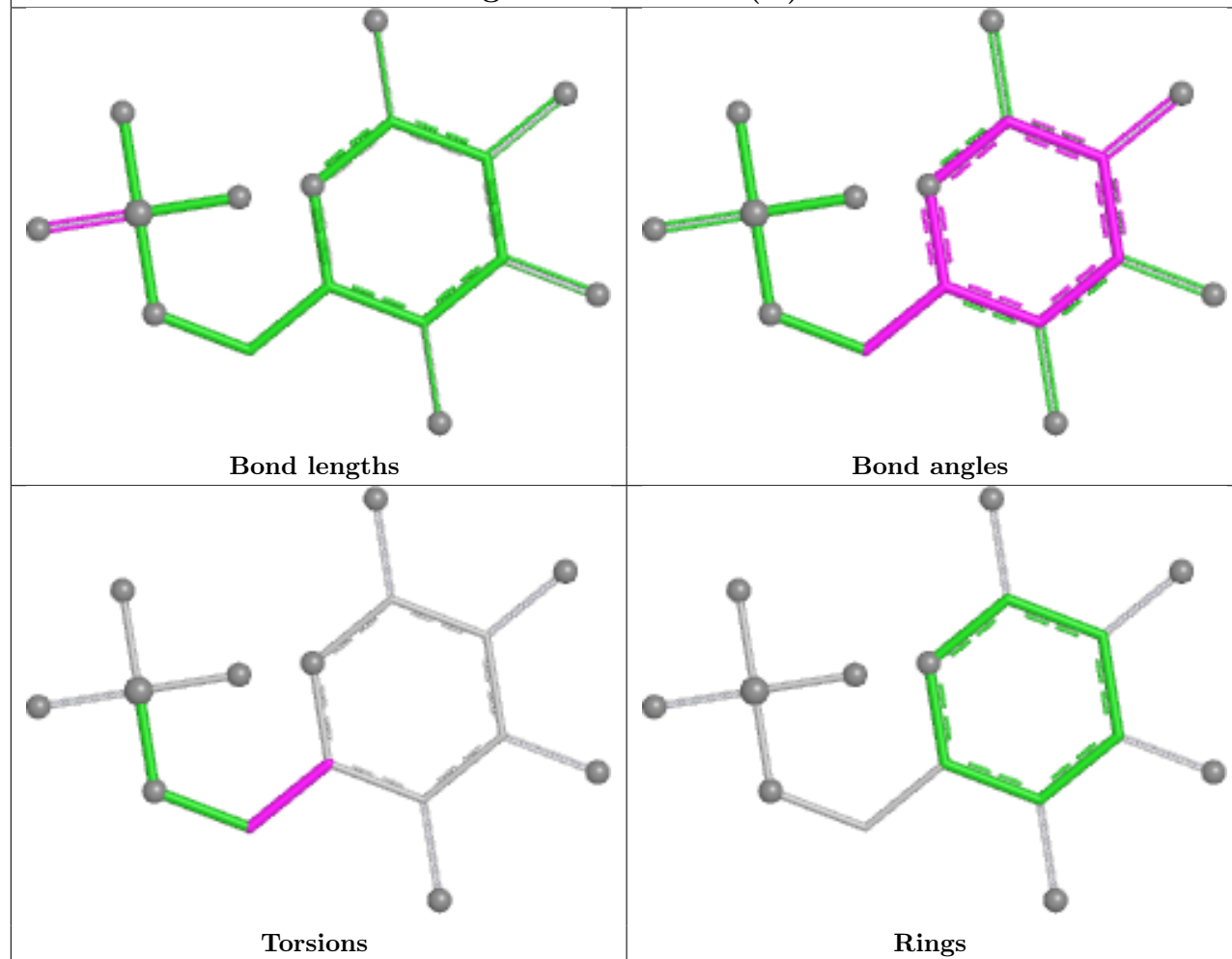
There are no ring outliers.

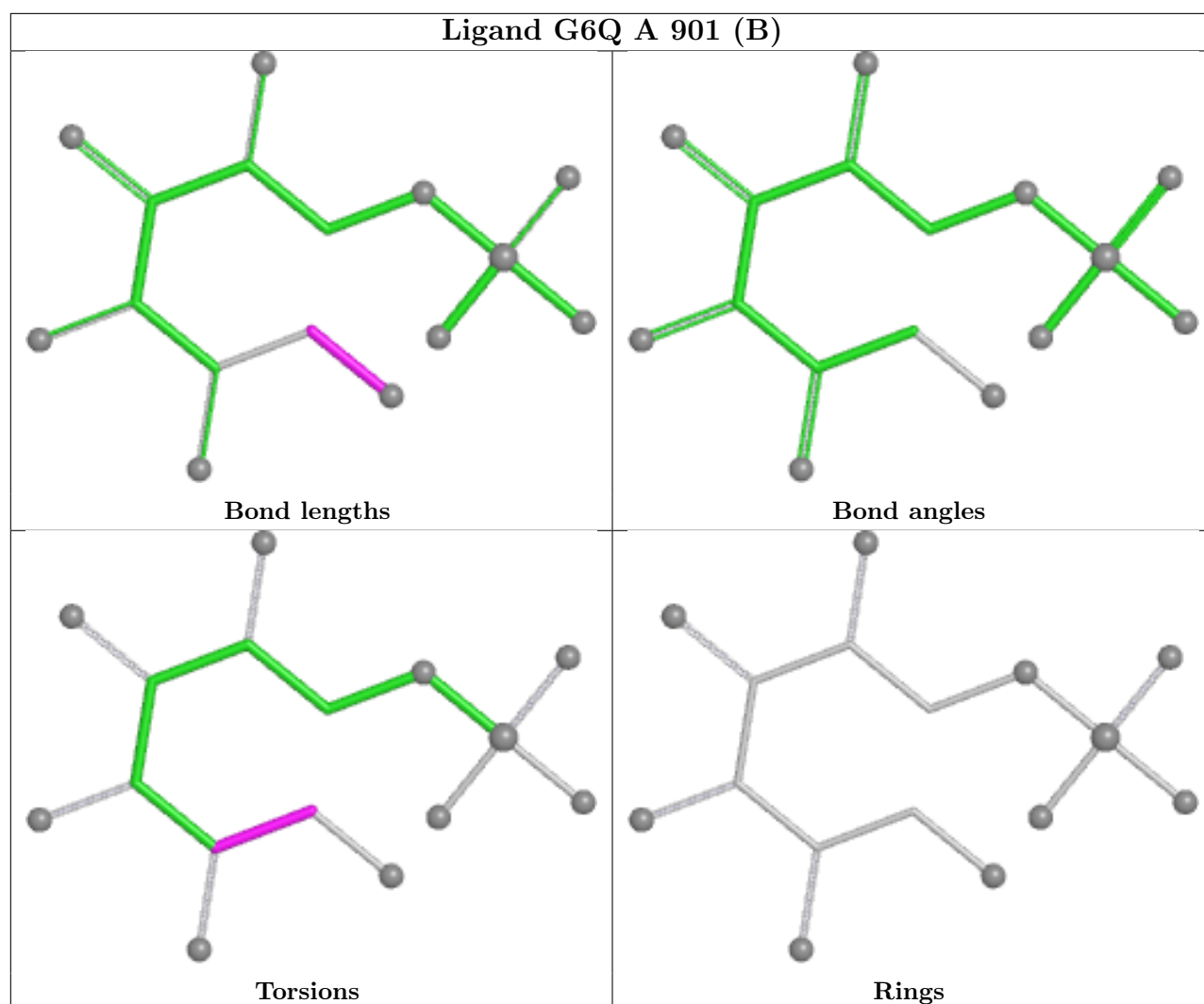
8 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1029	GOL	2	0
3	A	901[B]	G6Q	2	0
5	A	1030	GOL	3	0
5	A	1026	GOL	5	0
4	B	1008	SO4	2	0
5	A	1027	GOL	2	0
5	A	1025	GOL	2	0
5	A	1032	GOL	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand G6P A 900 (A)





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	556/564 (98%)	-0.03	13 (2%) 61 64	10, 19, 34, 50	15 (2%)
1	B	556/564 (98%)	0.32	30 (5%) 31 32	12, 23, 37, 45	13 (2%)
All	All	1112/1128 (98%)	0.14	43 (3%) 43 45	10, 21, 36, 50	28 (2%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	252	VAL	5.5
1	B	131	PHE	4.9
1	A	554	THR	4.2
1	B	255	PHE	3.9
1	B	250	ALA	3.8
1	B	211	THR	3.7
1	A	553	ASP	3.4
1	B	253	LYS	3.4
1	B	249	THR	3.2
1	B	369[A]	ARG	3.2
1	B	212	PHE	3.1
1	B	210	LYS	3.1
1	B	251	LYS	3.1
1	B	246	SER	3.0
1	B	290	ASP	2.9
1	B	556	LEU	2.9
1	B	114	GLY	2.8
1	B	31	ALA	2.7
1	B	553	ASP	2.7
1	B	259	PRO	2.7
1	B	254	GLU	2.6
1	B	214	THR	2.6
1	B	112	VAL	2.6
1	B	248	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	17	ARG	2.6
1	A	456	GLU	2.6
1	A	451	ALA	2.4
1	B	247	THR	2.4
1	B	22	ASN	2.4
1	A	458	LEU	2.4
1	A	259	PRO	2.3
1	A	556	LEU	2.3
1	B	17	ARG	2.3
1	B	213	THR	2.3
1	A	109	PRO	2.3
1	B	530	GLY	2.2
1	B	257	ILE	2.2
1	A	531	SER	2.2
1	A	442	GLU	2.2
1	A	452	GLY	2.1
1	A	450	ALA	2.1
1	B	30	GLU	2.1
1	B	263	LEU	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(Å ²)	Q<0.9
5	GOL	A	1032	6/6	0.48	0.18	41,43,45,47	0
5	GOL	A	1030	6/6	0.63	0.18	29,34,38,39	0
5	GOL	A	1026	6/6	0.67	0.17	38,39,41,41	0
4	SO4	B	1005	5/5	0.69	0.15	79,79,79,79	0

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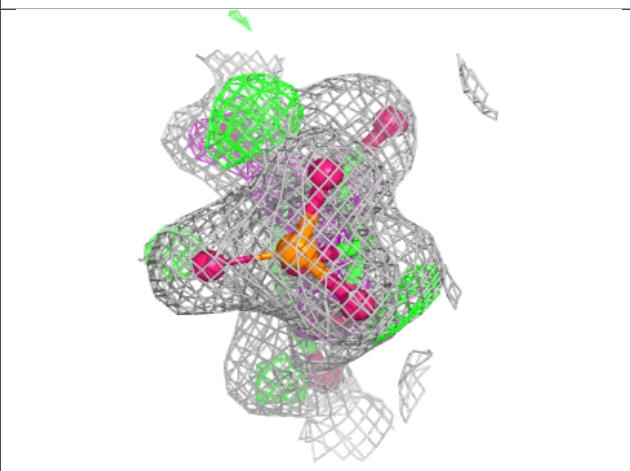
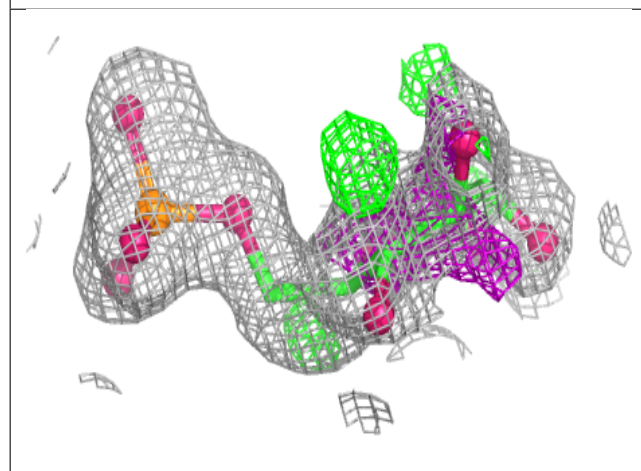
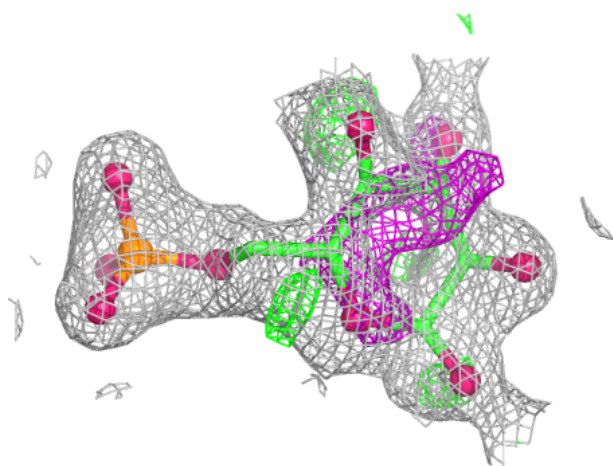
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	B	1029	6/6	0.69	0.14	39,40,41,43	0
5	GOL	A	1023	6/6	0.76	0.16	34,35,36,38	0
5	GOL	A	1031	6/6	0.77	0.15	27,35,37,38	0
5	GOL	A	1025	6/6	0.79	0.13	40,41,42,43	0
4	SO4	B	1008	5/5	0.79	0.15	58,58,58,58	0
4	SO4	B	1003	5/5	0.79	0.12	65,65,65,66	0
5	GOL	B	1028	6/6	0.80	0.12	40,41,42,43	0
6	BME	B	1011	4/4	0.80	0.18	40,41,42,43	0
5	GOL	A	1027	6/6	0.81	0.11	38,39,40,40	0
4	SO4	B	1006	5/5	0.82	0.12	55,55,56,56	0
4	SO4	B	1007	5/5	0.84	0.10	64,65,65,65	0
4	SO4	A	1002	5/5	0.85	0.11	75,75,76,76	0
5	GOL	A	1021	6/6	0.86	0.11	29,29,32,33	0
5	GOL	A	1024	6/6	0.87	0.10	33,35,36,37	0
2	G6P	A	900[A]	16/16	0.90	0.12	26,30,32,32	16
2	G6P	A	900[B]	1/16	0.90	0.12	29,29,29,29	1
3	G6Q	A	901[B]	16/16	0.90	0.12	8,17,20,22	16
5	GOL	B	1022	6/6	0.91	0.10	32,33,34,35	0
4	SO4	B	1004	5/5	0.91	0.10	36,37,38,39	0
4	SO4	A	1001	5/5	0.93	0.09	41,41,42,43	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

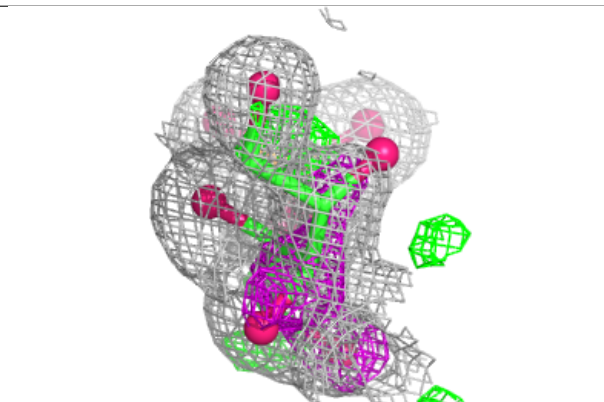
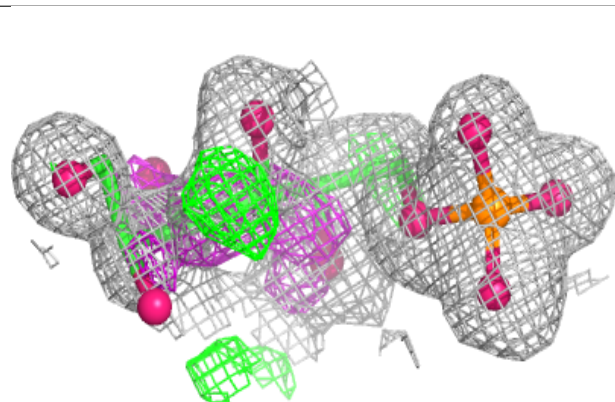
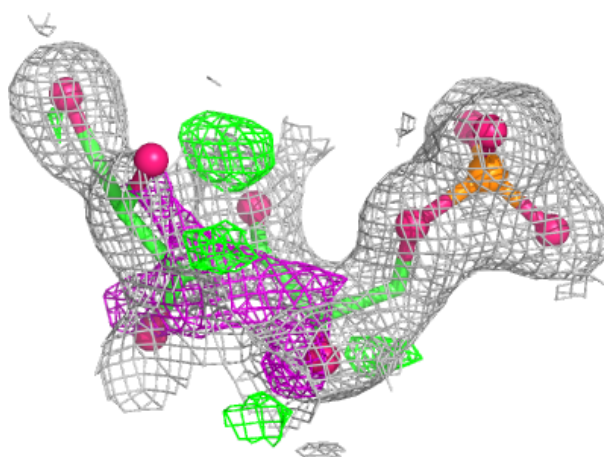
Electron density around G6P A 900 (A):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around G6Q A 901 (B):

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.