



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 05:36 PM BST

PDB ID : 5HUU
Title : Structure of Candida albicans trehalose-6-phosphate synthase in complex with
UDP and glucose-6-phosphate
Authors : Miao, Y.; Brennan, R.G.
Deposited on : 2016-01-27
Resolution : 2.37 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

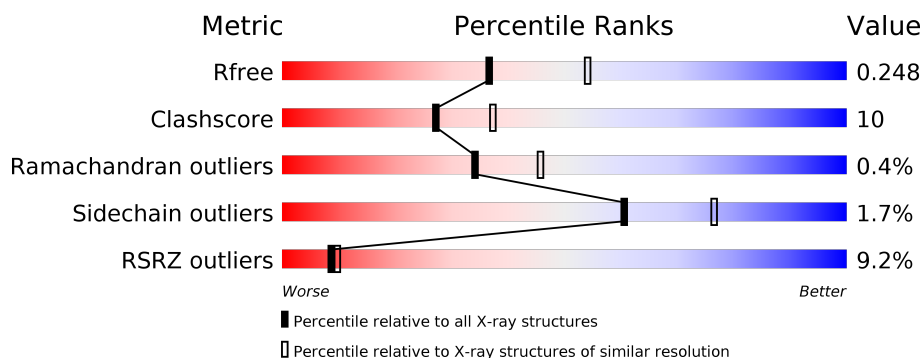
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



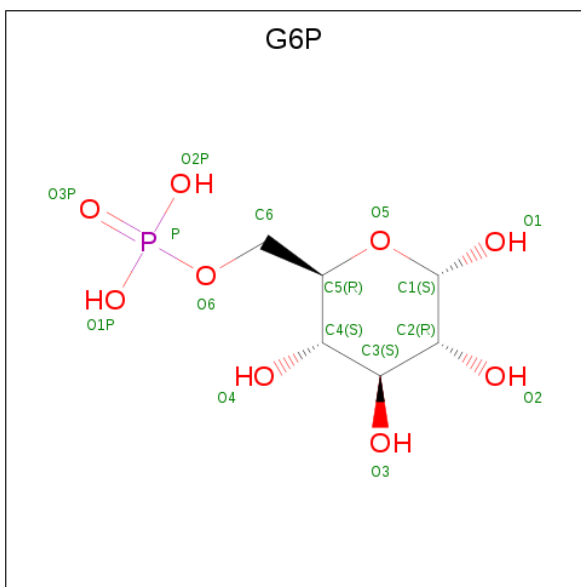
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 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	478	<div> <div>8%</div> <div>79%</div> <div>18%</div> <div>••</div> </div>
1	B	478	<div> <div>10%</div> <div>73%</div> <div>21%</div> <div>6%</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
3	G6P	B	502	-	-	X	-



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	O	P	0	0
			27	6	11	9	1		
3	B	1	Total	C	H	O	P	0	0
			27	6	11	9	1		

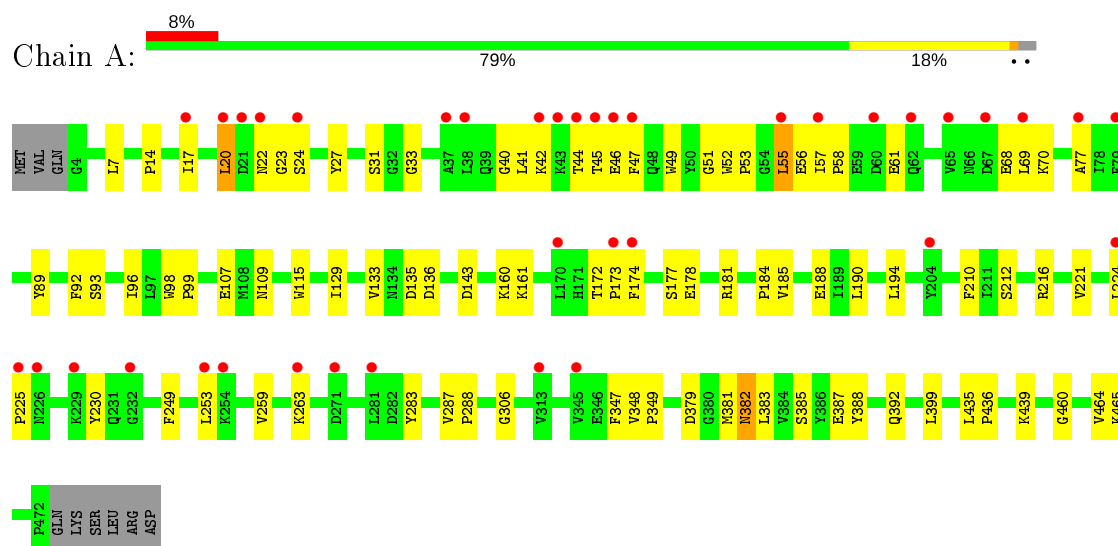
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	31	Total	O	0	0
			31	31		
4	B	25	Total	O	0	0
			25	25		

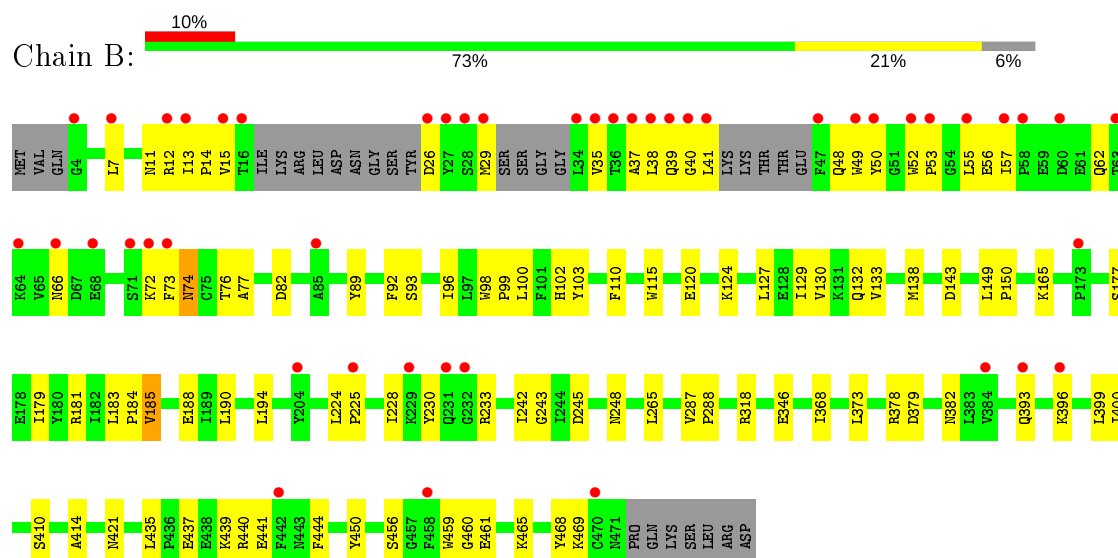
3 Residue-property plots

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: Alpha,alpha-trehalose-phosphate synthase [UDP-forming]



- Molecule 1: Alpha,alpha-trehalose-phosphate synthase [UDP-forming]



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	98.68Å 98.68Å 188.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.36 – 2.37 41.68 – 2.37	Depositor EDS
% Data completeness (in resolution range)	99.5 (35.36-2.37) 93.1 (41.68-2.37)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.37Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.200 , 0.247 0.204 , 0.248	Depositor DCC
R_{free} test set	1999 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	49.1	Xtriage
Anisotropy	0.480	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.057 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7584	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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: UDP, G6P

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	0.24	0/3864	0.40	0/5235
1	B	0.24	0/3717	0.40	0/5034
All	All	0.24	0/7581	0.40	0/10269

There are no bond length outliers.

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	3772	0	3754	67	0
1	B	3630	0	3607	84	0
2	A	25	11	10	0	0
2	B	25	11	10	2	0
3	A	16	11	11	0	0
3	B	16	11	11	6	0
4	A	31	0	0	0	0
4	B	25	0	0	0	0
All	All	7540	44	7403	151	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 10.

All (151) 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:38:LEU:HD21	1:B:460:GLY:HA3	1.27	1.10
1:A:41:LEU:HD11	1:A:460:GLY:HA3	1.41	1.02
1:B:53:PRO:HG3	1:B:77:ALA:HB1	1.43	0.99
1:A:115:TRP:CZ2	1:A:185:VAL:HG11	2.05	0.92
1:A:224:LEU:HB3	1:A:225:PRO:HD2	1.53	0.91
1:A:435:LEU:HD12	1:A:436:PRO:HD2	1.53	0.90
1:B:38:LEU:HD13	1:B:41:LEU:HD22	1.54	0.88
1:B:38:LEU:HD22	1:B:41:LEU:HB2	1.60	0.84
1:A:41:LEU:HD11	1:A:460:GLY:CA	2.08	0.83
1:B:7:LEU:HD22	1:B:50:TYR:HE1	1.43	0.82
1:B:39:GLN:N	1:B:40:GLY:HA3	1.92	0.82
1:A:22:ASN:HB2	1:A:24:SER:H	1.45	0.81
1:B:224:LEU:HB3	1:B:225:PRO:HD2	1.64	0.80
1:B:138:MET:HG2	1:B:165:LYS:HB2	1.68	0.76
1:B:265:LEU:HB3	1:B:368:ILE:HD13	1.71	0.71
1:A:185:VAL:HG12	1:A:188:GLU:HB2	1.71	0.71
1:A:22:ASN:N	1:A:23:GLY:HA2	2.05	0.70
1:A:185:VAL:CG1	1:A:188:GLU:HB2	2.22	0.70
1:B:177:SER:O	1:B:181:ARG:HG3	1.93	0.69
1:B:29:MET:CE	1:B:35:VAL:HG21	2.24	0.67
1:B:38:LEU:HD13	1:B:41:LEU:CD2	2.25	0.65
1:A:188:GLU:N	1:A:188:GLU:OE1	2.27	0.65
1:B:13:ILE:HG23	1:B:14:PRO:HD2	1.79	0.65
1:A:41:LEU:HA	1:A:44:THR:HB	1.77	0.64
1:A:212:SER:OG	1:A:216:ARG:NH1	2.31	0.64
1:A:115:TRP:CH2	1:A:185:VAL:HG11	2.33	0.64
1:B:245:ASP:HB3	1:B:248:ASN:OD1	1.98	0.63
1:B:53:PRO:CG	1:B:77:ALA:HB1	2.25	0.63
1:B:318:ARG:HH21	3:B:502:G6P:H61	1.63	0.63
1:B:102:HIS:O	1:B:378:ARG:HG3	1.98	0.63
1:B:230:TYR:O	1:B:233:ARG:HG2	1.98	0.62
1:A:20:LEU:HB2	1:A:24:SER:O	2.00	0.61
1:A:40:GLY:O	1:A:44:THR:OG1	2.18	0.61
1:B:224:LEU:CB	1:B:225:PRO:HD2	2.30	0.61
1:B:318:ARG:HH21	3:B:502:G6P:C6	2.13	0.61
1:A:55:LEU:HD12	1:A:56:GLU:N	2.15	0.61
1:A:27:TYR:CE2	1:A:68:GLU:HG2	2.35	0.61
1:B:52:TRP:HD1	1:B:53:PRO:HD2	1.66	0.60
1:B:37:ALA:O	1:B:456:SER:HB2	2.02	0.59
1:B:11:ASN:O	1:B:52:TRP:HB3	2.02	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:VAL:O	1:A:185:VAL:HG12	2.01	0.59
1:B:7:LEU:HD22	1:B:50:TYR:CE1	2.33	0.59
1:A:92:PHE:HA	1:A:96:ILE:HD12	1.85	0.58
1:B:243:GLY:O	1:B:456:SER:HB3	2.03	0.58
1:B:129:ILE:O	1:B:133:VAL:HG23	2.03	0.58
1:B:224:LEU:HB3	1:B:225:PRO:CD	2.33	0.58
1:B:410:SER:HB3	1:B:450:TYR:CZ	2.38	0.58
1:B:15:VAL:HG21	1:B:73:PHE:HD2	1.69	0.58
1:A:129:ILE:O	1:A:133:VAL:HG23	2.04	0.57
1:B:56:GLU:HG3	1:B:82:ASP:OD1	2.05	0.57
1:A:259:VAL:HG12	1:A:263:LYS:HE2	1.86	0.57
1:B:7:LEU:HD23	1:B:48:GLN:HB2	1.87	0.57
1:A:41:LEU:HA	1:A:44:THR:CB	2.34	0.56
1:B:120:GLU:O	1:B:124:LYS:HG2	2.05	0.56
1:B:38:LEU:HD21	1:B:460:GLY:CA	2.19	0.55
1:A:259:VAL:O	1:A:263:LYS:HG3	2.07	0.55
1:A:178:GLU:O	1:B:378:ARG:NH2	2.39	0.55
1:A:435:LEU:CD1	1:A:436:PRO:HD2	2.33	0.55
1:B:15:VAL:HG21	1:B:73:PHE:CD2	2.42	0.55
1:A:22:ASN:HB2	1:A:24:SER:N	2.19	0.55
1:B:48:GLN:NE2	1:B:74:ASN:OD1	2.39	0.55
1:A:46:GLU:CD	1:A:46:GLU:H	2.11	0.54
1:A:224:LEU:HB3	1:A:225:PRO:CD	2.33	0.53
1:B:461:GLU:O	1:B:465:LYS:HG3	2.09	0.53
1:A:190:LEU:O	1:A:194:LEU:HG	2.09	0.52
1:A:174:PHE:CE2	1:A:210:PHE:HA	2.45	0.52
1:B:55:LEU:HD11	1:B:57:ILE:HD11	1.92	0.52
1:B:265:LEU:CB	1:B:368:ILE:HD13	2.39	0.52
1:B:183:LEU:HD12	1:B:184:PRO:HD2	1.92	0.51
1:B:13:ILE:HG22	1:B:15:VAL:H	1.74	0.51
1:B:149:LEU:HB3	1:B:150:PRO:HD3	1.91	0.51
1:B:7:LEU:CD2	1:B:48:GLN:HB2	2.41	0.50
1:B:265:LEU:HB3	1:B:368:ILE:CD1	2.41	0.50
1:B:399:LEU:HD23	1:B:414:ALA:HB2	1.93	0.50
1:B:76:THR:HG21	1:B:132:GLN:HE22	1.77	0.50
1:B:26:ASP:OD1	1:B:72:LYS:NZ	2.44	0.50
1:B:465:LYS:O	1:B:469:LYS:HG3	2.12	0.50
1:B:12:ARG:NH1	3:B:502:G6P:O1P	2.45	0.49
1:A:58:PRO:HG2	1:A:61:GLU:OE1	2.12	0.49
1:B:39:GLN:N	1:B:40:GLY:CA	2.71	0.49
1:B:127:LEU:O	1:B:130:VAL:HG12	2.11	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:PRO:HG3	1:A:49:TRP:CE3	2.48	0.49
1:A:52:TRP:CD1	1:A:53:PRO:HD2	2.48	0.49
1:B:190:LEU:O	1:B:194:LEU:HG	2.12	0.49
1:B:110:PHE:CE1	1:B:184:PRO:HG2	2.48	0.48
1:B:29:MET:HE1	1:B:35:VAL:HG21	1.93	0.48
1:B:92:PHE:CD1	1:B:96:ILE:HD12	2.49	0.48
1:A:70:LYS:HB2	1:A:70:LYS:HE3	1.67	0.48
1:B:57:ILE:O	1:B:62:GLN:NE2	2.42	0.48
1:A:221:VAL:HG22	1:A:230:TYR:CD2	2.48	0.48
1:A:460:GLY:O	1:A:464:VAL:HG23	2.14	0.47
1:B:53:PRO:HG3	1:B:77:ALA:CB	2.29	0.47
1:A:42:LYS:HG2	1:A:47:PHE:CE2	2.50	0.47
1:A:55:LEU:HD11	1:A:57:ILE:HD13	1.96	0.47
1:B:396:LYS:HG2	1:B:444:PHE:CD2	2.50	0.47
1:A:249:PHE:O	1:A:253:LEU:HG	2.15	0.46
1:A:160:LYS:C	1:A:161:LYS:HG2	2.35	0.46
1:B:98:TRP:N	1:B:99:PRO:HD2	2.31	0.46
1:B:29:MET:HE3	1:B:35:VAL:HG21	1.96	0.46
2:B:501:UDP:O1B	3:B:502:G6P:H1	2.16	0.46
1:A:98:TRP:N	1:A:99:PRO:HD2	2.31	0.46
1:B:49:TRP:HD1	1:B:50:TYR:N	2.14	0.46
1:A:172:THR:HB	1:A:173:PRO:CD	2.45	0.45
1:A:7:LEU:N	1:A:7:LEU:HD12	2.31	0.45
1:A:172:THR:HB	1:A:173:PRO:HD2	1.98	0.45
1:A:107:GLU:N	1:A:107:GLU:OE1	2.49	0.45
1:A:177:SER:O	1:A:181:ARG:HG3	2.17	0.45
1:B:115:TRP:CE2	1:B:185:VAL:HG11	2.52	0.45
1:B:435:LEU:HD23	1:B:440:ARG:HG2	1.99	0.45
1:A:14:PRO:HG3	1:A:49:TRP:CZ3	2.52	0.44
1:A:89:TYR:O	1:A:93:SER:HB2	2.18	0.44
1:B:55:LEU:HD12	1:B:56:GLU:O	2.16	0.44
1:A:388:TYR:O	1:A:392:GLN:HG2	2.17	0.44
1:B:14:PRO:HB3	1:B:35:VAL:HB	1.98	0.44
1:A:184:PRO:HB3	1:B:421:ASN:ND2	2.32	0.44
1:B:439:LYS:HB2	1:B:439:LYS:HE2	1.79	0.44
1:B:13:ILE:CG2	1:B:14:PRO:HD2	2.48	0.44
1:B:89:TYR:O	1:B:93:SER:HB2	2.17	0.43
1:A:287:VAL:HB	1:A:288:PRO:HD3	2.00	0.43
1:A:55:LEU:HD12	1:A:56:GLU:C	2.38	0.43
1:B:242:ILE:HG13	1:B:459:TRP:CD2	2.53	0.43
1:B:62:GLN:O	1:B:66:ASN:HB2	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:LEU:O	1:A:387:GLU:HG3	2.19	0.43
1:B:96:ILE:HD11	1:B:110:PHE:HA	2.01	0.43
1:A:135:ASP:O	1:A:136:ASP:HB2	2.18	0.43
1:A:381:MET:O	1:A:382:ASN:HB3	2.19	0.42
1:B:100:LEU:HD11	1:B:179:ILE:O	2.18	0.42
1:B:38:LEU:HD23	1:B:38:LEU:HA	1.83	0.42
1:A:99:PRO:HB3	1:A:283:TYR:OH	2.19	0.42
1:A:17:ILE:HD11	1:A:69:LEU:HD21	2.02	0.42
1:A:51:GLY:O	1:A:77:ALA:HA	2.19	0.42
1:A:439:LYS:HD2	1:A:439:LYS:HA	1.74	0.42
1:B:102:HIS:O	1:B:103:TYR:HB2	2.20	0.42
1:A:20:LEU:HA	1:A:20:LEU:HD12	1.86	0.42
1:B:437:GLU:O	1:B:441:GLU:HG3	2.20	0.42
1:B:287:VAL:N	1:B:288:PRO:CD	2.83	0.41
1:A:348:VAL:HA	1:A:349:PRO:HD3	1.84	0.41
1:A:465:LYS:HA	1:A:465:LYS:HD3	1.78	0.41
1:B:224:LEU:H	1:B:228:ILE:HA	1.85	0.41
2:B:501:UDP:O2B	3:B:502:G6P:O1	2.27	0.41
1:B:55:LEU:HD11	1:B:57:ILE:CD1	2.50	0.41
1:B:89:TYR:OH	3:B:502:G6P:O2P	2.34	0.41
1:A:385:SER:HB2	1:A:399:LEU:HD21	2.02	0.41
1:B:373:LEU:HD13	1:B:400:ILE:HD12	2.02	0.41
1:B:14:PRO:HG3	1:B:49:TRP:CZ3	2.56	0.41
1:B:76:THR:HG21	1:B:132:GLN:NE2	2.35	0.41
1:A:306:GLY:HA2	1:A:347:PHE:CZ	2.55	0.40
1:A:178:GLU:OE1	1:A:181:ARG:NH1	2.54	0.40
1:A:46:GLU:HG2	1:A:46:GLU:O	2.22	0.40
1:A:57:ILE:HA	1:A:58:PRO:HD2	1.90	0.40
1:B:224:LEU:CB	1:B:225:PRO:CD	2.98	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	467/478 (98%)	449 (96%)	16 (3%)	2 (0%)	34	46
1	B	442/478 (92%)	423 (96%)	17 (4%)	2 (0%)	29	39
All	All	909/956 (95%)	872 (96%)	33 (4%)	4 (0%)	34	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	ASN
1	B	382	ASN
1	A	33	GLY
1	B	185	VAL

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	420/429 (98%)	413 (98%)	7 (2%)	60	76
1	B	404/429 (94%)	397 (98%)	7 (2%)	60	76
All	All	824/858 (96%)	810 (98%)	14 (2%)	60	76

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	31	SER
1	A	45	THR
1	A	55	LEU
1	A	109	ASN
1	A	143	ASP
1	A	379	ASP
1	B	74	ASN
1	B	143	ASP
1	B	188	GLU
1	B	346	GLU
1	B	379	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	393	GLN
1	B	468	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	48	GLN
1	B	74	ASN
1	B	132	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	UDP	B	501	-	20,26,26	5.48	10 (50%)	25,40,40	1.02	2 (8%)
2	UDP	A	501	-	20,26,26	5.50	10 (50%)	25,40,40	1.05	2 (8%)
3	G6P	B	502	-	16,16,16	1.02	1 (6%)	24,24,24	1.22	3 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	G6P	A	502	-	16,16,16	0.97	1 (6%)	24,24,24	0.71	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	UDP	B	501	-	-	4/14/32/32	0/2/2/2
2	UDP	A	501	-	-	4/14/32/32	0/2/2/2
3	G6P	B	502	-	-	5/6/26/26	0/1/1/1
3	G6P	A	502	-	-	2/6/26/26	0/1/1/1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	UDP	O4'-C1'	15.72	1.63	1.41
2	B	501	UDP	O4'-C1'	15.68	1.63	1.41
2	B	501	UDP	C2'-C1'	-13.81	1.32	1.53
2	A	501	UDP	C2'-C1'	-13.74	1.32	1.53
2	A	501	UDP	O4'-C4'	-6.47	1.30	1.45
2	B	501	UDP	O4'-C4'	-6.41	1.30	1.45
2	A	501	UDP	C6-N1	5.58	1.42	1.35
2	B	501	UDP	C6-N1	5.52	1.42	1.35
2	B	501	UDP	C4-N3	5.17	1.42	1.33
2	A	501	UDP	C4-N3	5.13	1.42	1.33
2	A	501	UDP	C6-C5	4.39	1.47	1.38
2	B	501	UDP	C6-C5	4.26	1.47	1.38
2	A	501	UDP	C2-N3	4.20	1.46	1.38
2	A	501	UDP	O3'-C3'	-4.16	1.33	1.43
2	B	501	UDP	C2-N3	4.15	1.46	1.38
2	B	501	UDP	O3'-C3'	-4.02	1.33	1.43
3	B	502	G6P	O5-C1	2.56	1.49	1.42
3	A	502	G6P	O5-C1	2.56	1.49	1.42
2	A	501	UDP	C3'-C4'	2.28	1.58	1.53
2	B	501	UDP	C3'-C4'	2.23	1.58	1.53
2	B	501	UDP	O2'-C2'	2.02	1.47	1.43
2	A	501	UDP	O2'-C2'	2.01	1.47	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	UDP	O4'-C1'-C2'	-3.22	102.22	106.93
3	B	502	G6P	O5-C5-C4	2.89	114.94	109.69
2	B	501	UDP	PA-O3A-PB	-2.62	123.82	132.83
3	B	502	G6P	C3-C4-C5	2.44	114.58	110.24
2	B	501	UDP	O4'-C1'-C2'	-2.37	103.47	106.93
3	B	502	G6P	O5-C1-C2	2.25	114.29	110.28
2	A	501	UDP	PA-O3A-PB	-2.07	125.72	132.83

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	UDP	O4'-C1'-N1-C6
3	B	502	G6P	C4-C5-C6-O6
3	B	502	G6P	O5-C5-C6-O6
3	B	502	G6P	C6-O6-P-O1P
3	B	502	G6P	C6-O6-P-O3P
3	A	502	G6P	C4-C5-C6-O6
3	A	502	G6P	O5-C5-C6-O6
2	B	501	UDP	C3'-C4'-C5'-O5'
2	B	501	UDP	O4'-C4'-C5'-O5'
2	A	501	UDP	O4'-C4'-C5'-O5'
2	A	501	UDP	C3'-C4'-C5'-O5'
2	B	501	UDP	PB-O3A-PA-O5'
2	A	501	UDP	PB-O3A-PA-O5'
2	B	501	UDP	PA-O3A-PB-O1B
3	B	502	G6P	C6-O6-P-O2P

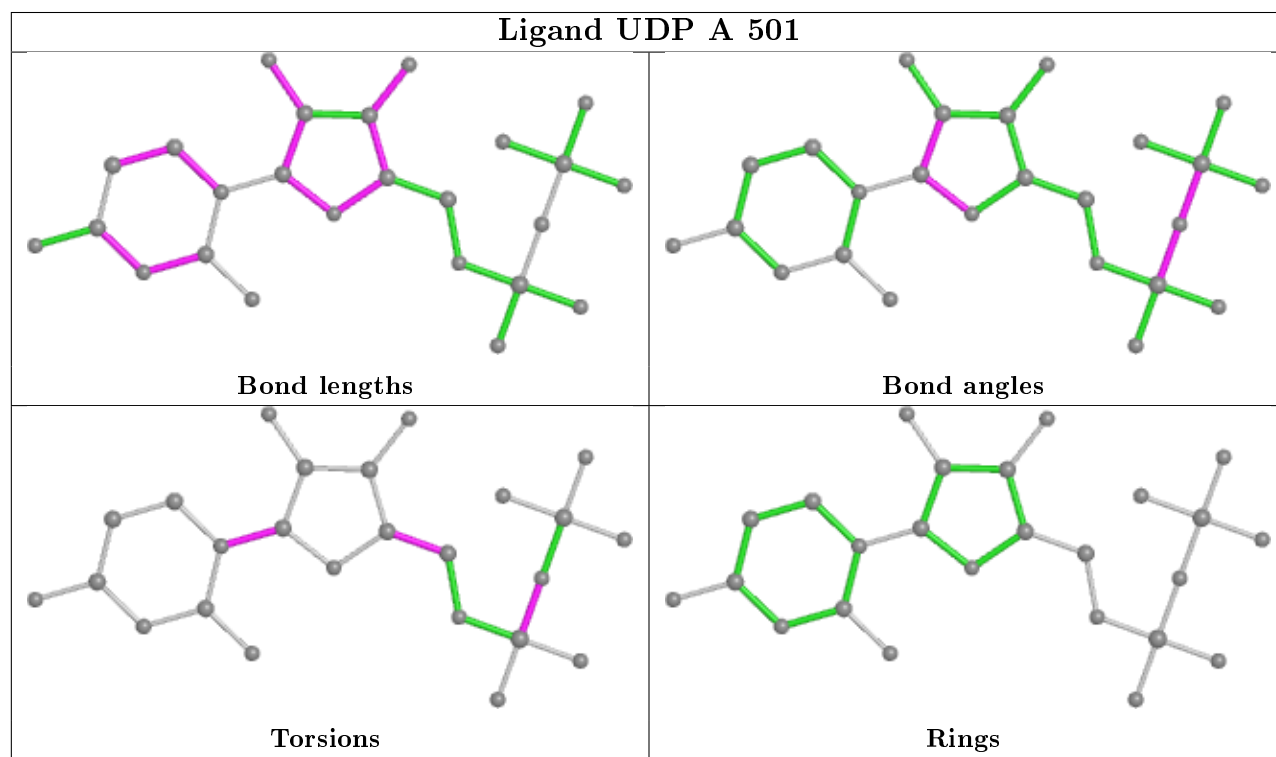
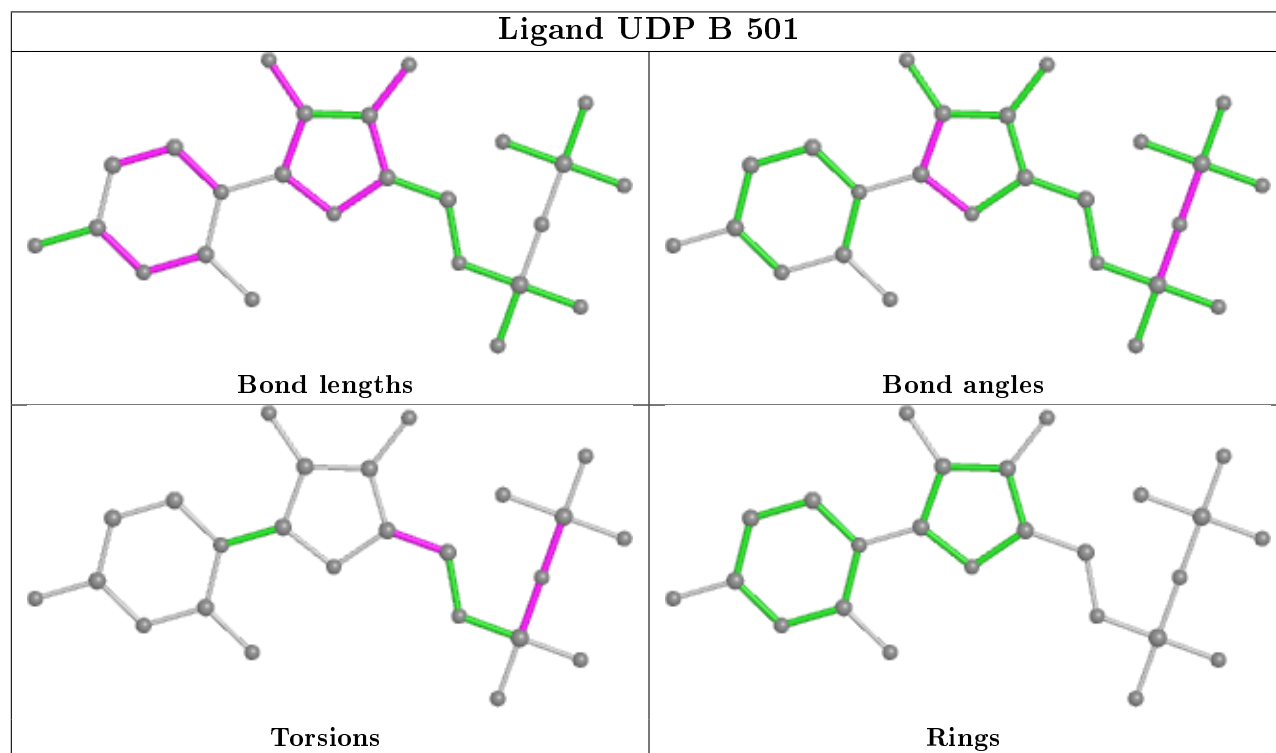
There are no ring outliers.

2 monomers are involved in 6 short contacts:

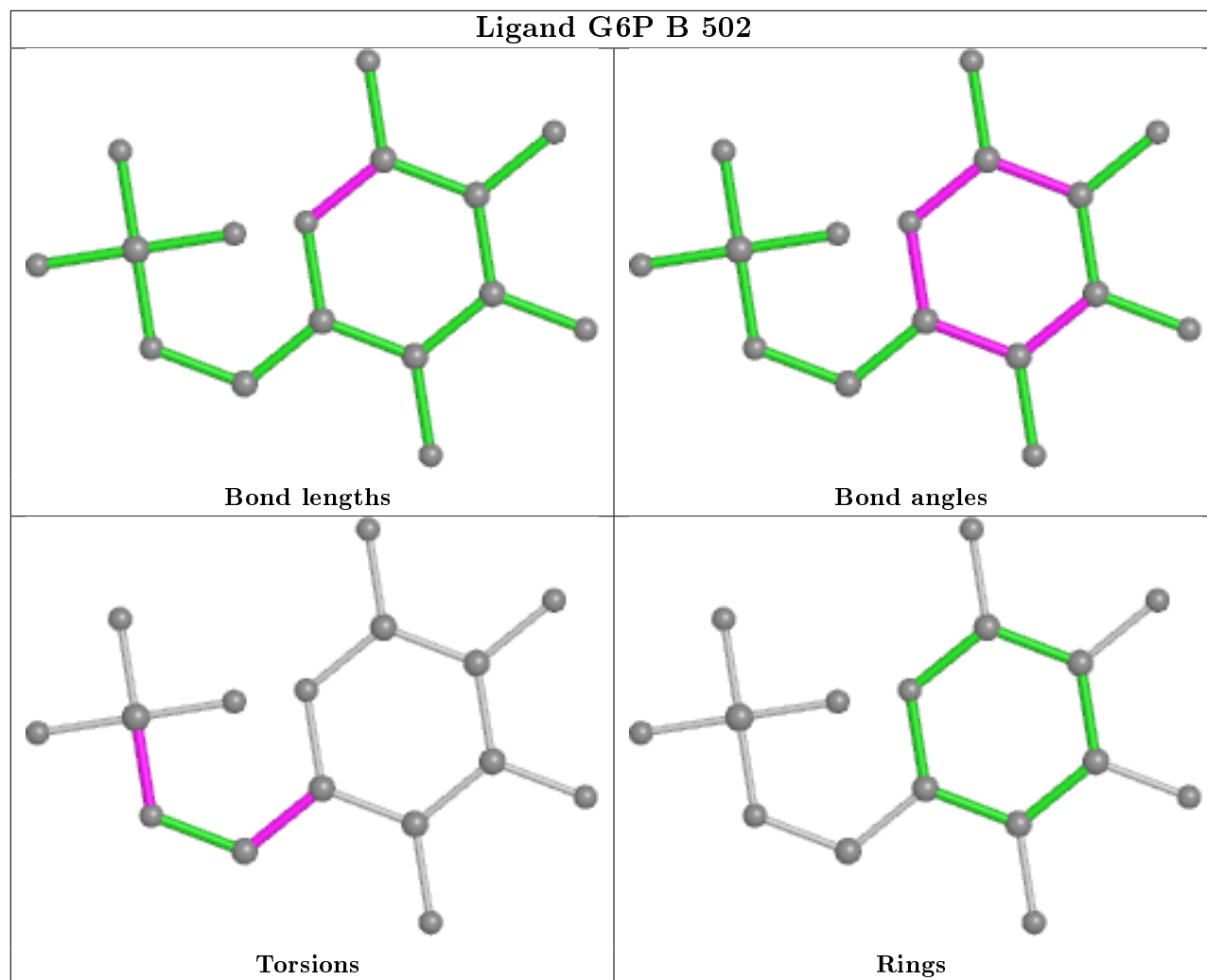
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	UDP	2	0
3	B	502	G6P	6	0

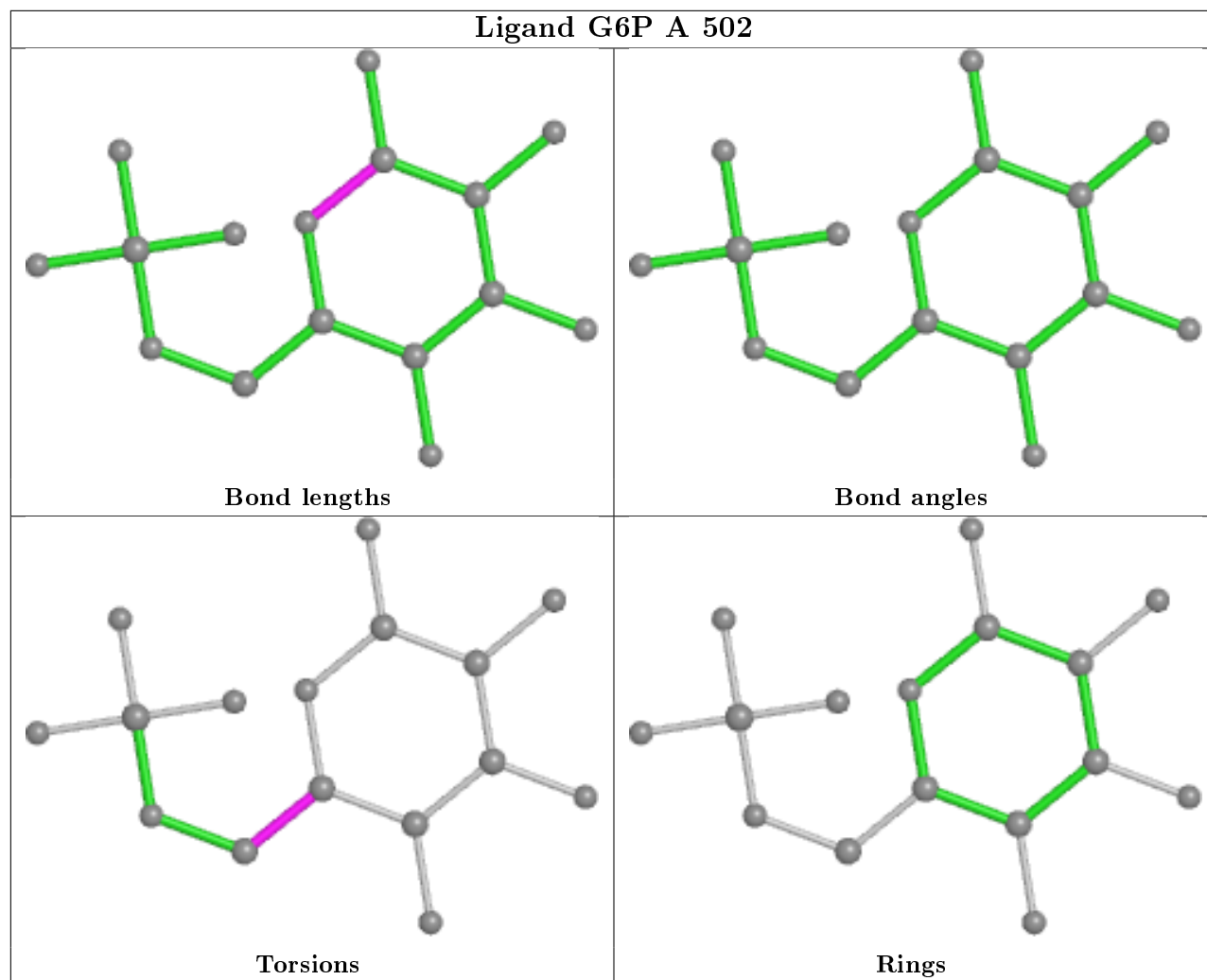
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 B 502





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	469/478 (98%)	0.49	38 (8%) 12 13	39, 57, 86, 111	0
1	B	450/478 (94%)	0.69	47 (10%) 6 7	41, 60, 107, 121	0
All	All	919/956 (96%)	0.58	85 (9%) 9 10	39, 59, 98, 121	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	37	ALA	10.6
1	B	35	VAL	7.2
1	B	73	PHE	7.1
1	B	29	MET	6.9
1	B	34	LEU	6.2
1	B	47	PHE	5.7
1	B	38	LEU	5.5
1	B	55	LEU	5.2
1	B	26	ASP	5.1
1	B	4	GLY	5.0
1	B	27	TYR	4.7
1	B	28	SER	4.6
1	B	64	LYS	4.4
1	B	58	PRO	4.4
1	A	45	THR	4.3
1	A	225	PRO	4.2
1	B	41	LEU	4.2
1	A	17	ILE	4.0
1	B	442	PHE	3.9
1	B	39	GLN	3.9
1	A	37	ALA	3.9
1	A	204	TYR	3.8
1	A	57	ILE	3.7
1	B	63	THR	3.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	47	PHE	3.7
1	B	72	LYS	3.7
1	A	232	GLY	3.6
1	B	57	ILE	3.5
1	B	71	SER	3.5
1	B	50	TYR	3.5
1	A	65	VAL	3.3
1	A	254	LYS	3.3
1	A	46	GLU	3.1
1	B	204	TYR	3.1
1	A	38	LEU	3.1
1	A	43	LYS	3.0
1	B	7	LEU	3.0
1	B	60	ASP	3.0
1	A	44	THR	2.9
1	B	229	LYS	2.9
1	A	22	ASN	2.9
1	A	77	ALA	2.8
1	A	226	ASN	2.8
1	A	69	LEU	2.8
1	A	24	SER	2.6
1	A	229	LYS	2.6
1	A	271	ASP	2.5
1	A	224	LEU	2.5
1	B	396	LYS	2.5
1	B	384	VAL	2.5
1	B	53	PRO	2.5
1	B	225	PRO	2.5
1	A	55	LEU	2.5
1	B	49	TRP	2.5
1	B	40	GLY	2.5
1	B	231	GLN	2.5
1	A	170	LEU	2.5
1	A	173	PRO	2.4
1	B	36	THR	2.4
1	A	20	LEU	2.3
1	A	42	LYS	2.3
1	B	15	VAL	2.3
1	B	393	GLN	2.3
1	B	66	ASN	2.3
1	B	68	GLU	2.3
1	B	85	ALA	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	52	TRP	2.3
1	B	458	PHE	2.3
1	A	67	ASP	2.3
1	A	62	GLN	2.3
1	A	263	LYS	2.2
1	A	313	VAL	2.2
1	A	79	PHE	2.2
1	B	13	ILE	2.2
1	B	470	CYS	2.2
1	B	173	PRO	2.2
1	A	21	ASP	2.2
1	A	60	ASP	2.2
1	A	174	PHE	2.1
1	B	16	THR	2.1
1	B	12	ARG	2.1
1	A	281	LEU	2.1
1	A	345	VAL	2.1
1	B	232	GLY	2.0
1	A	253	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 monosaccharides in this entry.

6.4 Ligands [i](#)

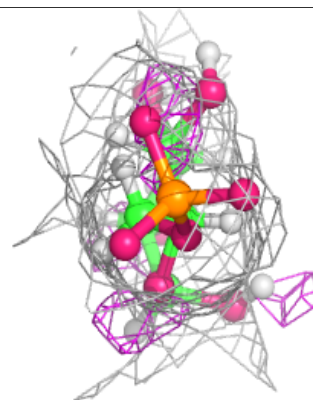
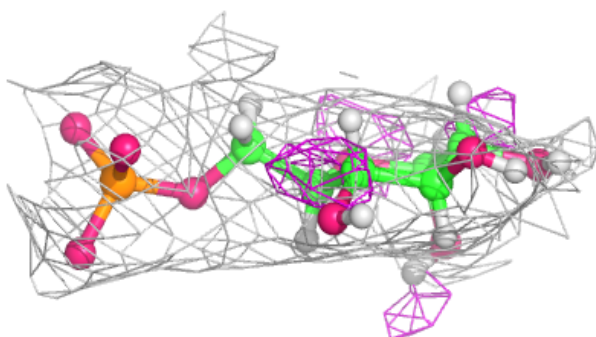
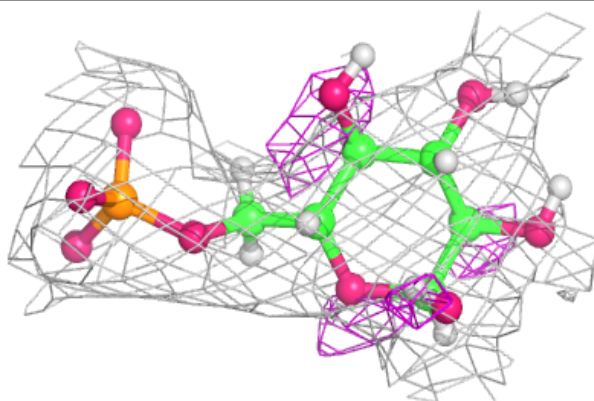
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	G6P	B	502	16/16	0.90	0.15	58,74,89,93	0
3	G6P	A	502	16/16	0.91	0.14	47,61,74,87	0
2	UDP	B	501	25/25	0.95	0.15	48,55,66,69	0
2	UDP	A	501	25/25	0.95	0.17	41,49,60,63	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.

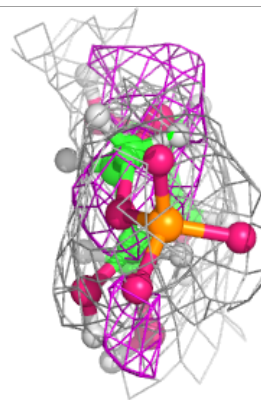
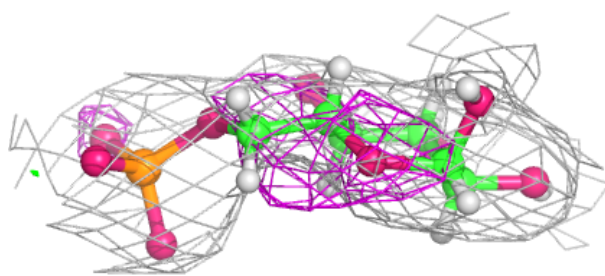
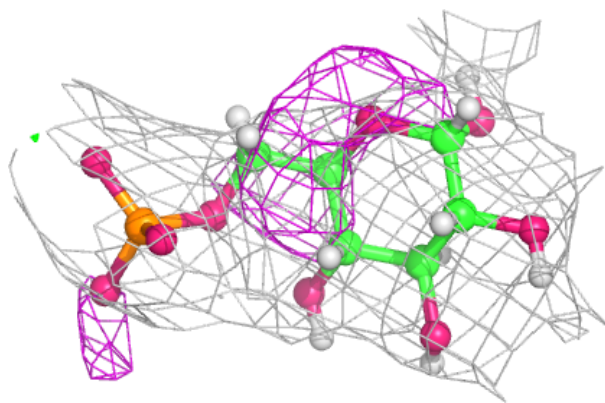
Electron density around G6P B 502:

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

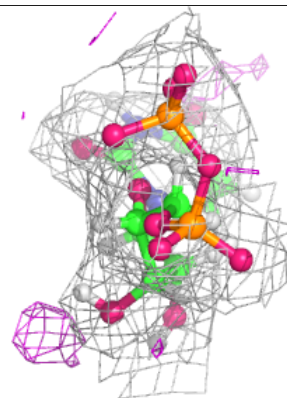
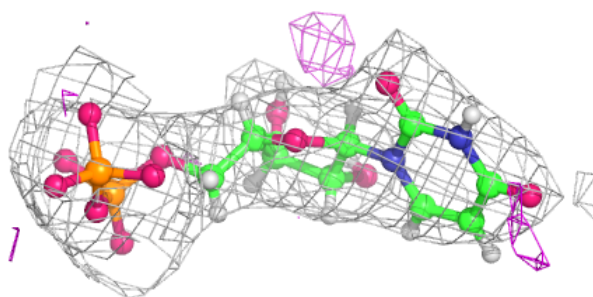
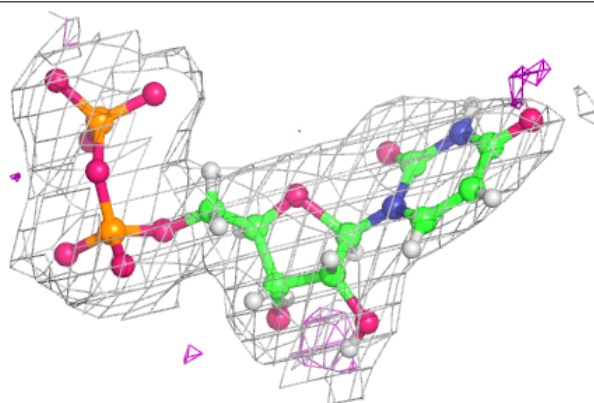


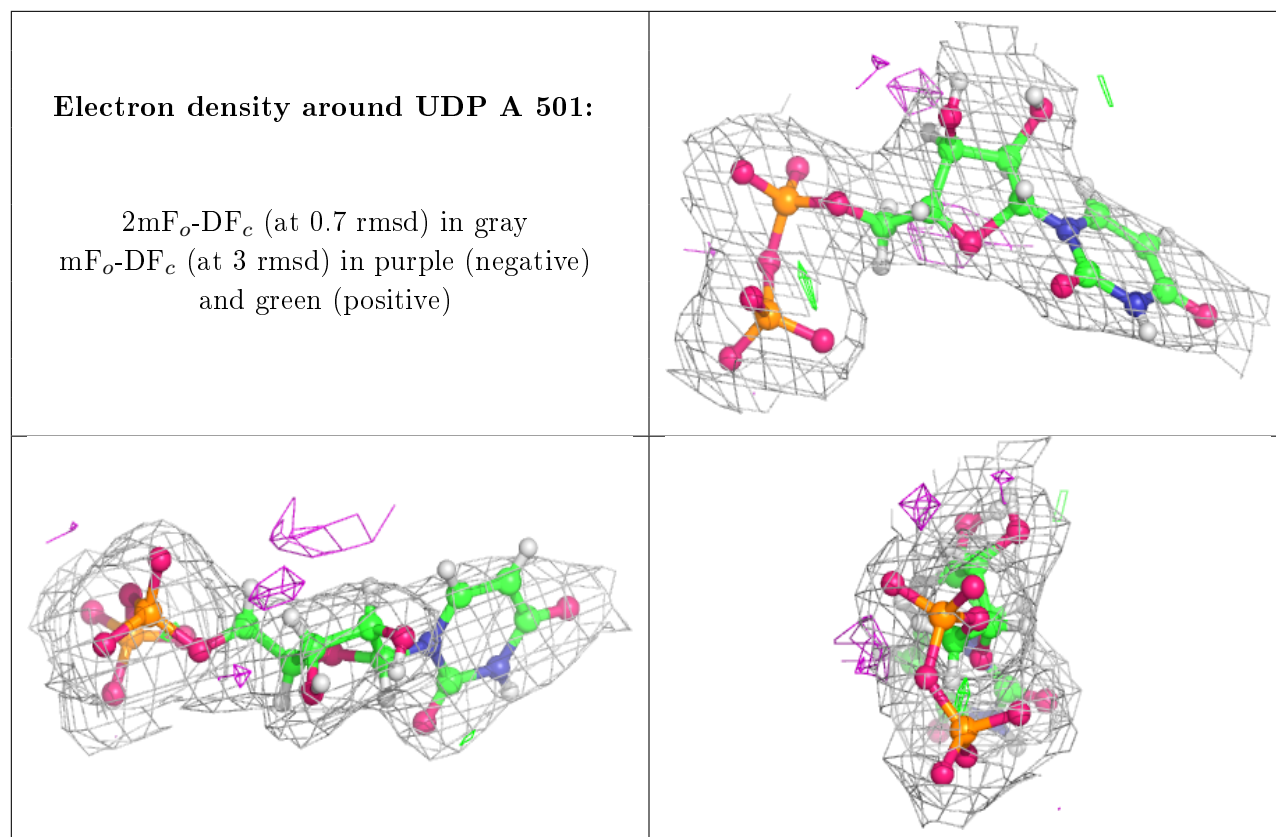
Electron density around G6P A 502:

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

**Electron density around UDP B 501:**

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





6.5 Other polymers ⓘ

There are no such residues in this entry.