



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2020 – 09:31 AM BST

PDB ID : 3NB0
Title : Glucose-6-Phosphate activated form of Yeast Glycogen Synthase
Authors : Baskaran, S.; Hurley, T.D.
Deposited on : 2010-06-02
Resolution : 2.41 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

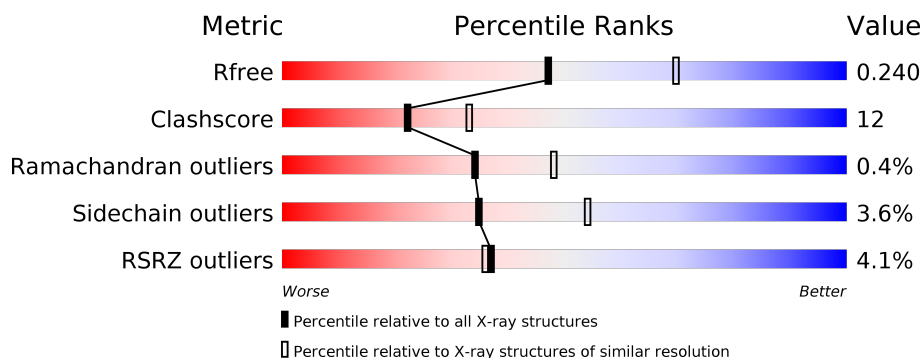
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	725	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>15%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	725	<div> <div>3%</div> <div> <div></div> <div>66%</div> <div>22%</div> <div>•</div> <div>11%</div> </div> </div>
1	C	725	<div> <div>3%</div> <div> <div></div> <div>66%</div> <div>21%</div> <div>•</div> <div>11%</div> </div> </div>
1	D	725	<div> <div>6%</div> <div> <div></div> <div>60%</div> <div>26%</div> <div>•</div> <div>12%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21392 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 Glycogen [starch] synthase isoform 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	638	Total	C	N	O	S	0	4	0
			5169	3302	899	949	19			
1	B	645	Total	C	N	O	S	0	3	0
			5213	3332	911	951	19			
1	C	646	Total	C	N	O	S	0	2	0
			5213	3331	910	953	19			
1	D	636	Total	C	N	O	S	0	1	0
			5135	3279	896	941	19			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P27472
A	-18	GLY	-	expression tag	UNP P27472
A	-17	SER	-	expression tag	UNP P27472
A	-16	SER	-	expression tag	UNP P27472
A	-15	HIS	-	expression tag	UNP P27472
A	-14	HIS	-	expression tag	UNP P27472
A	-13	HIS	-	expression tag	UNP P27472
A	-12	HIS	-	expression tag	UNP P27472
A	-11	HIS	-	expression tag	UNP P27472
A	-10	HIS	-	expression tag	UNP P27472
A	-9	SER	-	expression tag	UNP P27472
A	-8	SER	-	expression tag	UNP P27472
A	-7	GLY	-	expression tag	UNP P27472
A	-6	LEU	-	expression tag	UNP P27472
A	-5	VAL	-	expression tag	UNP P27472
A	-4	PRO	-	expression tag	UNP P27472
A	-3	ARG	-	expression tag	UNP P27472
A	-2	GLY	-	expression tag	UNP P27472
A	-1	SER	-	expression tag	UNP P27472
A	0	HIS	-	expression tag	UNP P27472
A	535	SER	ALA	SEE REMARK 999	UNP P27472

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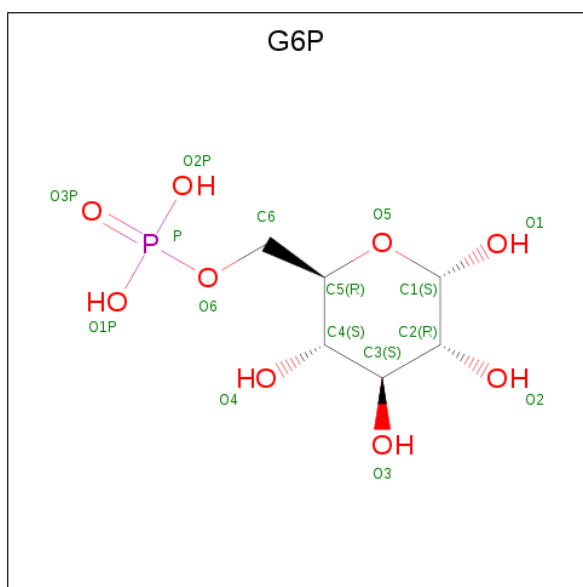
Chain	Residue	Modelled	Actual	Comment	Reference
A	589	ALA	ARG	engineered mutation	UNP P27472
A	592	ALA	ARG	engineered mutation	UNP P27472
B	-19	MET	-	expression tag	UNP P27472
B	-18	GLY	-	expression tag	UNP P27472
B	-17	SER	-	expression tag	UNP P27472
B	-16	SER	-	expression tag	UNP P27472
B	-15	HIS	-	expression tag	UNP P27472
B	-14	HIS	-	expression tag	UNP P27472
B	-13	HIS	-	expression tag	UNP P27472
B	-12	HIS	-	expression tag	UNP P27472
B	-11	HIS	-	expression tag	UNP P27472
B	-10	HIS	-	expression tag	UNP P27472
B	-9	SER	-	expression tag	UNP P27472
B	-8	SER	-	expression tag	UNP P27472
B	-7	GLY	-	expression tag	UNP P27472
B	-6	LEU	-	expression tag	UNP P27472
B	-5	VAL	-	expression tag	UNP P27472
B	-4	PRO	-	expression tag	UNP P27472
B	-3	ARG	-	expression tag	UNP P27472
B	-2	GLY	-	expression tag	UNP P27472
B	-1	SER	-	expression tag	UNP P27472
B	0	HIS	-	expression tag	UNP P27472
B	535	SER	ALA	SEE REMARK 999	UNP P27472
B	589	ALA	ARG	engineered mutation	UNP P27472
B	592	ALA	ARG	engineered mutation	UNP P27472
C	-19	MET	-	expression tag	UNP P27472
C	-18	GLY	-	expression tag	UNP P27472
C	-17	SER	-	expression tag	UNP P27472
C	-16	SER	-	expression tag	UNP P27472
C	-15	HIS	-	expression tag	UNP P27472
C	-14	HIS	-	expression tag	UNP P27472
C	-13	HIS	-	expression tag	UNP P27472
C	-12	HIS	-	expression tag	UNP P27472
C	-11	HIS	-	expression tag	UNP P27472
C	-10	HIS	-	expression tag	UNP P27472
C	-9	SER	-	expression tag	UNP P27472
C	-8	SER	-	expression tag	UNP P27472
C	-7	GLY	-	expression tag	UNP P27472
C	-6	LEU	-	expression tag	UNP P27472
C	-5	VAL	-	expression tag	UNP P27472
C	-4	PRO	-	expression tag	UNP P27472
C	-3	ARG	-	expression tag	UNP P27472

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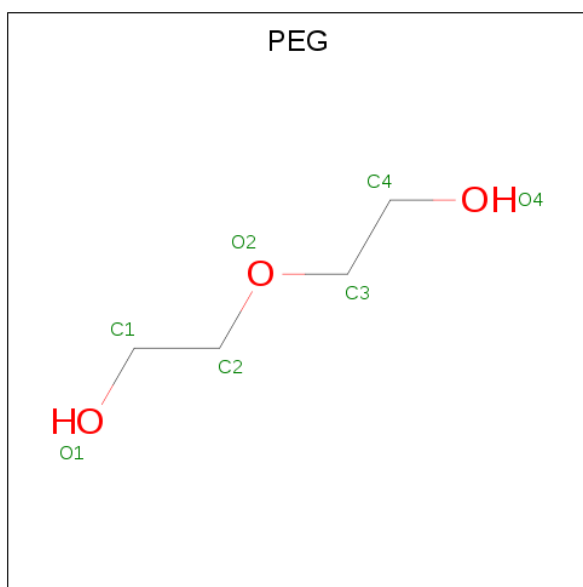
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP P27472
C	-1	SER	-	expression tag	UNP P27472
C	0	HIS	-	expression tag	UNP P27472
C	535	SER	ALA	SEE REMARK 999	UNP P27472
C	589	ALA	ARG	engineered mutation	UNP P27472
C	592	ALA	ARG	engineered mutation	UNP P27472
D	-19	MET	-	expression tag	UNP P27472
D	-18	GLY	-	expression tag	UNP P27472
D	-17	SER	-	expression tag	UNP P27472
D	-16	SER	-	expression tag	UNP P27472
D	-15	HIS	-	expression tag	UNP P27472
D	-14	HIS	-	expression tag	UNP P27472
D	-13	HIS	-	expression tag	UNP P27472
D	-12	HIS	-	expression tag	UNP P27472
D	-11	HIS	-	expression tag	UNP P27472
D	-10	HIS	-	expression tag	UNP P27472
D	-9	SER	-	expression tag	UNP P27472
D	-8	SER	-	expression tag	UNP P27472
D	-7	GLY	-	expression tag	UNP P27472
D	-6	LEU	-	expression tag	UNP P27472
D	-5	VAL	-	expression tag	UNP P27472
D	-4	PRO	-	expression tag	UNP P27472
D	-3	ARG	-	expression tag	UNP P27472
D	-2	GLY	-	expression tag	UNP P27472
D	-1	SER	-	expression tag	UNP P27472
D	0	HIS	-	expression tag	UNP P27472
D	535	SER	ALA	SEE REMARK 999	UNP P27472
D	589	ALA	ARG	engineered mutation	UNP P27472
D	592	ALA	ARG	engineered mutation	UNP P27472

- Molecule 2 is 6-O-phosphono-alpha-D-glucopyranose (three-letter code: G6P) (formula: $C_6H_{13}O_9P$).



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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).

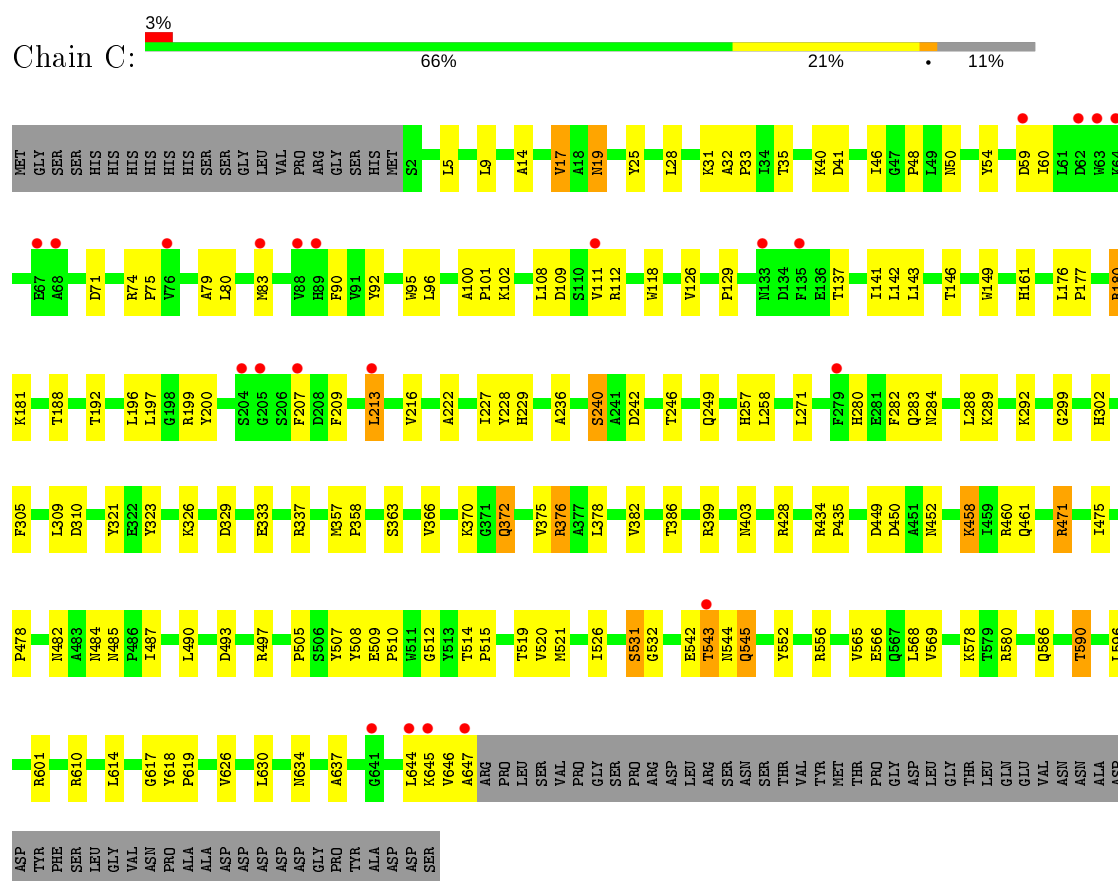


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	C	1	Total	C	O	0	0
			7	4	3		
3	D	1	Total	C	O	0	0
			7	4	3		

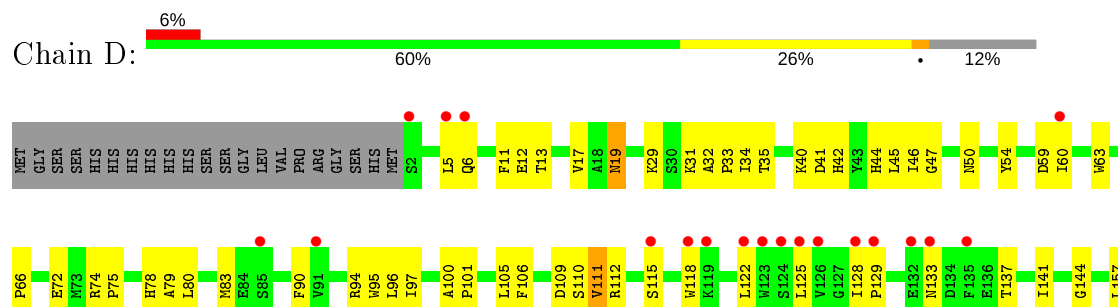
- Molecule 4 is water.

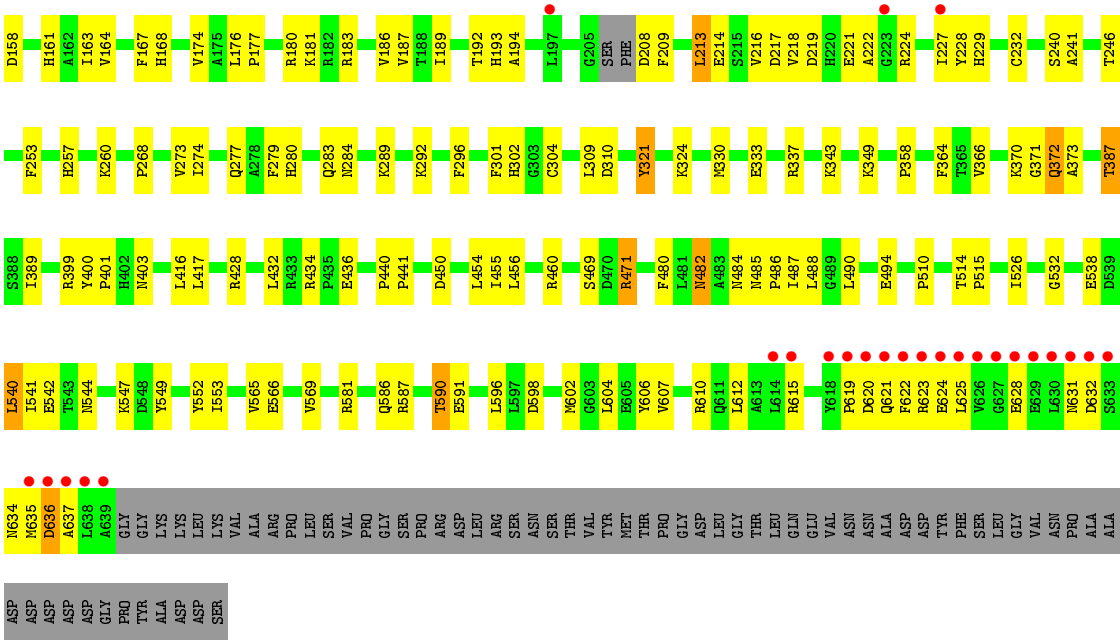
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	176	Total	O	0	0
			176	176		
4	B	127	Total	O	0	0
			127	127		
4	C	102	Total	O	0	0
			102	102		
4	D	119	Total	O	0	0
			119	119		

- Molecule 1: Glycogen [starch] synthase isoform 2



- Molecule 1: Glycogen [starch] synthase isoform 2





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	192.74Å 206.98Å 205.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.23 – 2.41 46.23 – 2.41	Depositor EDS
% Data completeness (in resolution range)	93.6 (46.23-2.41) 99.1 (46.23-2.41)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 2.39Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.211 , 0.242 0.210 , 0.240	Depositor DCC
R_{free} test set	7890 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.480	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21392	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.20% 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: G6P, PEG

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	0.26	0/5307	0.43	0/7191
1	B	0.26	0/5349	0.42	0/7246
1	C	0.25	0/5345	0.41	0/7239
1	D	0.25	0/5262	0.41	0/7129
All	All	0.25	0/21263	0.42	0/28805

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 [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	5169	0	5079	78	0
1	B	5213	0	5136	123	0
1	C	5213	0	5139	122	0
1	D	5135	0	5046	161	0
2	A	32	0	20	1	0
2	B	32	0	20	2	0
2	C	16	0	10	2	0
2	D	16	0	10	3	0
3	A	7	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	21	0	30	2	0
3	C	7	0	10	0	0
3	D	7	0	10	0	0
4	A	176	0	0	4	0
4	B	127	0	0	5	0
4	C	102	0	0	7	0
4	D	119	0	0	9	0
All	All	21392	0	20520	482	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 12.

All (482) 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:D:482:ASN:HD22	1:D:484:ASN:H	1.01	0.97
1:C:180:ARG:HG3	1:C:180:ARG:HH11	1.27	0.96
1:C:283:GLN:HG2	1:D:280:HIS:CE1	2.07	0.90
1:C:399:ARG:HD3	1:C:403:ASN:HD22	1.40	0.86
1:C:280:HIS:CE1	1:D:283:GLN:HG2	2.12	0.85
1:B:312:THR:HG22	1:B:350:THR:HB	1.59	0.83
1:C:59:ASP:HB2	1:C:96:LEU:HD21	1.61	0.83
1:C:542:GLU:HB3	1:C:545:GLN:HG3	1.61	0.82
1:D:540:LEU:HB3	1:D:541:ILE:HD12	1.62	0.81
1:D:482:ASN:ND2	1:D:484:ASN:H	1.76	0.81
1:C:372[A]:GLN:NE2	1:C:376:ARG:HH21	1.79	0.80
1:A:60:ILE:HD12	1:A:60:ILE:H	1.48	0.79
1:B:391:LYS:HD3	3:B:1001:PEG:H41	1.65	0.79
1:D:343:LYS:HD3	1:D:469:SER:O	1.83	0.78
1:D:440:PRO:HG3	4:D:763:HOH:O	1.83	0.78
1:C:519:THR:HG23	4:C:805:HOH:O	1.84	0.77
1:B:634:ASN:ND2	1:B:637:ALA:H	1.81	0.76
1:D:110:SER:O	1:D:111:VAL:HG23	1.86	0.76
1:B:626:VAL:HG11	1:B:630:LEU:HD11	1.68	0.75
1:D:292:LYS:HD2	1:D:490:LEU:HD21	1.68	0.75
1:A:187:VAL:CG1	1:A:613:ALA:HB1	2.17	0.74
1:B:86:ARG:HG2	1:B:149:TRP:HZ2	1.54	0.72
1:D:213:LEU:HD21	1:D:253:PHE:CE2	2.25	0.72
1:D:213:LEU:HD21	1:D:253:PHE:HE2	1.53	0.72
1:B:579:THR:H	1:B:582:GLN:NE2	1.87	0.71
1:B:459:ILE:HG21	1:B:474:MET:HG2	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:302:HIS:HD2	1:D:432:LEU:O	1.72	0.71
1:B:307:PHE:HD1	1:B:312:THR:HG21	1.55	0.71
1:C:180:ARG:NH1	1:C:180:ARG:HG3	2.05	0.71
1:D:283:GLN:NE2	1:D:587:ARG:HH21	1.88	0.70
1:D:192:THR:HG22	1:D:246:THR:HG22	1.71	0.70
1:D:615:ARG:HD3	1:D:622:PHE:CD1	2.26	0.70
1:D:214:GLU:HA	1:D:257:HIS:HD2	1.57	0.70
1:B:86:ARG:HG2	1:B:149:TRP:CZ2	2.28	0.69
1:C:236:ALA:O	1:C:240:SER:HB2	1.93	0.69
1:B:63:TRP:HZ3	1:B:81:GLN:HG3	1.58	0.68
1:B:493:ASP:O	1:B:497:ARG:HG3	1.93	0.68
1:C:450:ASP:OD1	1:C:460:ARG:NH2	2.26	0.68
1:B:579:THR:H	1:B:582:GLN:HE21	1.39	0.68
1:D:32:ALA:HB3	1:D:33:PRO:HD3	1.77	0.67
1:A:296:PHE:HA	1:A:372[B]:GLN:HE22	1.59	0.67
1:B:364:PHE:CD1	1:B:487:ILE:HD12	2.29	0.67
1:A:542:GLU:OE1	1:A:545:GLN:HB2	1.95	0.67
1:D:31:LYS:NZ	1:D:35:THR:HG21	2.10	0.66
1:B:304:CYS:SG	1:B:434:ARG:HD3	2.35	0.66
1:A:187:VAL:HG11	1:A:613:ALA:HB1	1.76	0.66
2:B:902:G6P:H1	4:B:823:HOH:O	1.96	0.66
1:B:126:VAL:HG22	1:B:181:LYS:HZ2	1.59	0.66
1:D:101:PRO:HD2	4:D:784:HOH:O	1.96	0.65
1:A:213:LEU:HA	1:A:216:VAL:HG13	1.78	0.65
1:D:100:ALA:HA	4:D:784:HOH:O	1.96	0.65
1:B:458:LYS:HD3	1:B:458:LYS:O	1.96	0.65
1:D:209:PHE:O	1:D:213:LEU:HB3	1.96	0.65
1:A:339:ASN:O	1:A:343:LYS:HG3	1.97	0.65
1:B:307:PHE:CD1	1:B:312:THR:HG21	2.32	0.65
1:B:458:LYS:HE2	1:B:461:GLN:OE1	1.96	0.65
1:A:8:HIS:HA	1:A:161:HIS:HB3	1.76	0.65
1:D:514:THR:HB	1:D:515:PRO:HD3	1.79	0.64
1:C:634:ASN:HB2	1:C:637:ALA:H	1.62	0.64
1:B:634:ASN:C	1:B:634:ASN:HD22	2.02	0.63
1:D:183:ARG:HG3	1:D:183:ARG:O	1.97	0.63
1:B:119:LYS:HE3	1:B:130:SER:O	1.98	0.63
1:A:254:GLU:HB3	4:A:878:HOH:O	1.98	0.63
1:B:126:VAL:HG22	1:B:181:LYS:NZ	2.13	0.63
1:D:612:LEU:HD22	1:D:635:MET:HG2	1.81	0.63
1:B:133:ASN:H	1:B:133:ASN:ND2	1.96	0.63
1:D:399:ARG:HD3	1:D:403:ASN:HD22	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:ASP:O	1:B:312:THR:HG23	1.98	0.62
1:D:128:ILE:HG12	1:D:232:CYS:HB3	1.80	0.62
1:B:513:TYR:O	1:B:517:GLU:HG2	2.00	0.61
1:B:273:VAL:HG13	1:B:520:VAL:HG13	1.83	0.61
1:D:180:ARG:HA	1:D:180:ARG:HE	1.65	0.61
1:B:246:THR:HB	4:B:793:HOH:O	1.99	0.61
1:B:213:LEU:O	1:B:216:VAL:HG13	2.00	0.60
1:A:128:ILE:HG12	1:A:232:CYS:HB3	1.83	0.60
1:C:199:ARG:HG3	1:C:508:TYR:CE2	2.35	0.60
1:C:284:ASN:HD21	1:D:284:ASN:HD21	1.47	0.60
1:D:606:TYR:O	1:D:610:ARG:HG3	2.01	0.60
1:C:458:LYS:HE2	1:C:461:GLN:OE1	2.00	0.60
1:D:227:ILE:HG22	1:D:227:ILE:O	2.00	0.60
1:D:482:ASN:HD22	1:D:484:ASN:N	1.86	0.60
1:D:289:LYS:HE3	1:D:494:GLU:HG2	1.83	0.60
1:D:296:PHE:HE1	1:D:487:ILE:HD12	1.67	0.60
1:B:295:ASP:CG	1:B:376:ARG:HH22	2.06	0.60
1:A:32:ALA:HB3	1:A:33:PRO:HD3	1.84	0.59
1:C:209:PHE:O	1:C:213:LEU:HB3	2.02	0.59
1:C:458:LYS:O	1:C:458:LYS:HD3	2.02	0.59
1:D:66:PRO:HA	1:D:74:ARG:HH12	1.67	0.59
1:B:174:VAL:O	1:B:177:PRO:HD2	2.03	0.59
1:B:450:ASP:OD1	1:B:460:ARG:NH2	2.36	0.59
1:C:493:ASP:O	1:C:497:ARG:HG3	2.02	0.59
1:B:295:ASP:OD1	1:B:376:ARG:NH2	2.35	0.59
1:C:299:GLY:HA2	1:C:375:VAL:HG21	1.85	0.59
1:A:78:HIS:HB3	1:A:157:LEU:HD23	1.85	0.59
1:C:372[A]:GLN:HE21	1:C:376:ARG:HH21	1.46	0.59
1:D:129:PRO:HG2	1:D:229:HIS:HB3	1.84	0.59
1:B:323:TYR:CZ	1:B:329:ASP:HB3	2.38	0.59
1:D:137:THR:HG21	1:D:229:HIS:HD2	1.67	0.59
1:A:83:MET:HG2	1:A:88:VAL:HG11	1.85	0.58
1:A:187:VAL:HG12	1:A:613:ALA:HB1	1.85	0.58
1:C:197:LEU:HD23	1:C:227:ILE:HD11	1.85	0.58
1:D:109:ASP:HA	1:D:112:ARG:NH1	2.18	0.58
1:D:97:ILE:HG12	4:D:784:HOH:O	2.03	0.58
1:B:396:HIS:HD2	1:B:415:GLU:OE2	1.87	0.58
1:B:430:LEU:HD12	1:B:433:ARG:HD3	1.84	0.58
1:B:111:VAL:CG1	1:B:118:TRP:CH2	2.87	0.57
1:D:586:GLN:O	1:D:590:THR:HG22	2.04	0.57
1:D:144:GLY:HA3	1:D:174:VAL:HB	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:111:VAL:HG12	1:C:111:VAL:O	2.04	0.57
1:A:146:THR:O	1:A:149:TRP:HB3	2.05	0.57
1:C:485:ASN:CG	4:C:706:HOH:O	2.43	0.57
1:C:213:LEU:HA	1:C:216:VAL:HG13	1.85	0.57
1:D:79:ALA:O	1:D:83:MET:HG2	2.05	0.57
1:A:624:GLU:C	1:A:626:VAL:H	2.08	0.57
1:C:31:LYS:O	1:C:35:THR:HG23	2.05	0.57
1:D:31:LYS:HZ2	1:D:35:THR:HG21	1.69	0.57
1:A:32:ALA:HB1	1:A:101:PRO:HG3	1.87	0.57
1:B:331:PHE:CZ	1:B:335:LEU:HD11	2.41	0.56
1:B:299:GLY:HA2	1:B:375:VAL:HG21	1.86	0.56
1:C:399:ARG:CD	1:C:403:ASN:HD22	2.17	0.56
1:B:323:TYR:CE1	1:B:329:ASP:HB3	2.41	0.56
1:B:59:ASP:HB2	1:B:96:LEU:HD21	1.86	0.56
1:D:213:LEU:O	1:D:216:VAL:HG22	2.06	0.56
1:C:109:ASP:HA	1:C:112:ARG:NH1	2.21	0.56
1:B:634:ASN:HD21	1:B:637:ALA:H	1.49	0.55
1:B:455:ILE:O	1:B:459:ILE:HG12	2.06	0.55
1:D:540:LEU:HD21	1:D:596:LEU:HD13	1.88	0.55
1:C:586:GLN:O	1:C:590:THR:HG23	2.06	0.55
1:D:301:PHE:HA	4:D:763:HOH:O	2.06	0.55
1:D:221:GLU:HA	1:D:224:ARG:HB3	1.87	0.55
1:D:31:LYS:O	1:D:34:ILE:HG22	2.07	0.55
1:A:485:ASN:HB3	4:A:722:HOH:O	2.07	0.55
1:B:144:GLY:HA3	1:B:174:VAL:HB	1.89	0.55
1:A:295[A]:ASP:CG	1:A:376:ARG:HH22	2.09	0.55
1:A:458:LYS:HD3	1:A:458:LYS:O	2.07	0.55
1:A:626:VAL:HG12	1:A:627:GLY:N	2.22	0.55
1:B:514:THR:HB	1:B:515:PRO:HD3	1.89	0.54
1:D:283:GLN:HE21	1:D:587:ARG:HH21	1.52	0.54
1:D:208:ASP:OD1	1:D:209:PHE:N	2.40	0.54
1:A:144:GLY:HA3	1:A:174:VAL:HB	1.89	0.54
1:A:3:ARG:NH2	1:A:158:ASP:O	2.41	0.54
1:C:192:THR:HG22	1:C:246:THR:HG22	1.89	0.54
1:C:323:TYR:CZ	1:C:329:ASP:HB3	2.42	0.54
1:D:623:ARG:HA	1:D:628:GLU:O	2.08	0.54
1:D:547:LYS:HG2	1:D:552:TYR:CD1	2.43	0.54
1:C:137:THR:O	1:C:141:ILE:HG13	2.08	0.53
1:D:31:LYS:O	1:D:35:THR:HG23	2.07	0.53
1:B:199:ARG:HB3	1:B:508:TYR:CE2	2.43	0.53
1:C:578:LYS:NZ	1:C:586:GLN:NE2	2.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:50:ASN:O	1:C:54:TYR:HB3	2.08	0.53
1:D:59:ASP:HB2	1:D:96:LEU:HD21	1.91	0.53
1:C:487:ILE:HG12	4:C:706:HOH:O	2.08	0.53
1:C:510:PRO:O	1:C:532:GLY:HA3	2.08	0.53
1:D:510:PRO:O	1:D:532:GLY:HA3	2.09	0.53
1:D:222:ALA:HB1	1:D:228:TYR:HA	1.91	0.53
1:D:615:ARG:HG3	1:D:632:ASP:OD2	2.08	0.53
1:C:382:VAL:O	1:C:386:THR:HG23	2.09	0.53
1:C:485:ASN:ND2	4:C:706:HOH:O	2.41	0.53
1:B:536:TYR:O	1:B:540:LEU:HD23	2.09	0.53
1:C:372[A]:GLN:NE2	1:C:376:ARG:NH2	2.55	0.53
1:A:199:ARG:HH22	2:A:902:G6P:P	2.33	0.52
1:C:146:THR:O	1:C:149:TRP:HB3	2.09	0.52
1:B:176:LEU:HB2	1:B:177:PRO:HD3	1.91	0.52
1:C:180:ARG:CG	1:C:180:ARG:HH11	2.08	0.52
1:A:45:LEU:HB2	1:A:103:VAL:HG12	1.91	0.52
1:C:399:ARG:HD3	1:C:403:ASN:ND2	2.18	0.52
1:C:48:PRO:HG3	1:C:143:LEU:HD22	1.92	0.52
1:B:378:LEU:HD22	1:B:432:LEU:HD11	1.91	0.52
2:B:902:G6P:O4	2:B:902:G6P:O3P	2.12	0.52
1:C:507:TYR:HB2	1:C:556:ARG:NH2	2.23	0.52
1:B:396:HIS:HE1	1:B:407:THR:O	1.93	0.52
1:D:164:VAL:HA	1:D:187:VAL:HG23	1.92	0.52
1:D:217:ASP:O	1:D:221:GLU:HG2	2.09	0.52
1:B:565:VAL:O	1:B:569:VAL:HG23	2.09	0.51
1:D:137:THR:O	1:D:141:ILE:HG13	2.10	0.51
1:D:621:GLN:O	1:D:625:LEU:HB2	2.10	0.51
1:C:176:LEU:HB2	1:C:177:PRO:HD3	1.92	0.51
1:B:122:LEU:HD11	1:B:128:ILE:HD12	1.91	0.51
1:C:288:LEU:HD23	1:C:289:LYS:HD3	1.92	0.51
1:C:596:LEU:CD1	1:C:646:VAL:HG21	2.40	0.51
1:C:610:ARG:HD2	4:C:729:HOH:O	2.10	0.51
1:B:463:GLN:HA	1:B:465:PHE:CE2	2.46	0.51
1:A:291[A]:GLU:OE1	1:A:291[A]:GLU:HA	2.11	0.51
1:C:449:ASP:OD2	1:C:452:ASN:HB2	2.11	0.51
1:A:83:MET:HE1	1:A:153:GLU:HG3	1.91	0.50
1:D:482:ASN:ND2	1:D:484:ASN:HB2	2.26	0.50
1:D:549:TYR:O	1:D:590:THR:HB	2.11	0.50
1:B:493:ASP:HB2	1:B:521:MET:CE	2.41	0.50
1:C:357:MET:O	1:C:478:PRO:HA	2.10	0.50
1:D:180:ARG:NE	1:D:180:ARG:HA	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:177:PRO:O	1:D:181:LYS:HG2	2.11	0.50
1:C:366:VAL:O	1:C:370:LYS:HB2	2.12	0.50
1:D:485:ASN:HB3	4:D:710:HOH:O	2.11	0.50
1:A:392:ARG:CZ	1:A:418:LYS:HG2	2.42	0.50
1:C:333:GLU:OE2	1:C:337:ARG:HD2	2.11	0.50
1:A:266:ILE:HG22	1:A:268:PRO:HD3	1.94	0.50
1:A:70:SER:O	1:A:74:ARG:HG2	2.12	0.50
1:D:214:GLU:HA	1:D:257:HIS:CD2	2.44	0.50
1:C:283:GLN:HG3	2:C:901:G6P:O1	2.12	0.50
1:D:109:ASP:HA	1:D:112:ARG:HH12	1.74	0.50
1:D:450:ASP:OD1	1:D:456:LEU:HD13	2.12	0.50
1:D:487:ILE:HG13	1:D:488:LEU:N	2.26	0.50
1:C:19:ASN:N	1:C:19:ASN:HD22	2.10	0.49
1:C:79:ALA:O	1:C:83:MET:HG2	2.12	0.49
1:D:434:ARG:NH1	4:D:763:HOH:O	2.44	0.49
1:B:542:GLU:CD	1:B:544:ASN:H	2.15	0.49
1:A:153:GLU:O	1:A:157:LEU:HD13	2.11	0.49
1:B:158:ASP:OD2	1:B:161:HIS:HD2	1.96	0.49
1:B:417:LEU:HD23	1:B:421:ASP:HB2	1.94	0.49
1:D:72:GLU:O	1:D:161:HIS:HE1	1.96	0.49
1:C:578:LYS:HZ1	1:C:586:GLN:NE2	2.11	0.49
1:A:298:ARG:NH2	4:A:760:HOH:O	2.46	0.49
1:C:526:ILE:HG12	1:C:552:TYR:HB2	1.95	0.49
1:A:25:TYR:CE1	1:A:95:TRP:HZ2	2.31	0.49
1:C:302:HIS:O	1:C:434:ARG:HD2	2.12	0.49
1:D:587:ARG:HA	1:D:590:THR:HG23	1.94	0.49
1:A:264:ASP:O	1:A:635:MET:HG3	2.13	0.48
1:B:293:ILE:O	1:B:297:VAL:HG23	2.13	0.48
1:B:349:LYS:O	1:B:471:ARG:CG	2.61	0.48
1:C:323:TYR:CE1	1:C:329:ASP:HB3	2.48	0.48
1:C:514:THR:HB	1:C:515:PRO:HD3	1.96	0.48
1:A:35:THR:HG21	1:A:43:TYR:CE1	2.48	0.48
1:C:493:ASP:HB2	1:C:521:MET:CE	2.44	0.48
1:A:458:LYS:HD3	1:A:458:LYS:C	2.33	0.48
1:A:164:VAL:HG11	1:A:610:ARG:HG2	1.95	0.48
1:A:78:HIS:CB	1:A:157:LEU:HD23	2.43	0.48
1:D:296:PHE:CE1	1:D:487:ILE:HD12	2.49	0.48
1:B:419:SER:O	1:B:423:VAL:HG23	2.14	0.48
1:A:305:PHE:HZ	1:A:309:LEU:HG	1.79	0.48
1:B:447:MET:HB2	1:B:450:ASP:HB2	1.95	0.48
1:C:80:LEU:HD22	1:C:90:PHE:CE1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:464:LEU:HD13	1:B:474:MET:HE3	1.96	0.48
1:C:17:VAL:HG21	1:C:46:ILE:O	2.14	0.47
1:D:17:VAL:HG21	1:D:46:ILE:O	2.14	0.47
1:B:163:ILE:HB	1:B:186:VAL:HG12	1.96	0.47
1:B:273:VAL:HG13	1:B:520:VAL:CG1	2.44	0.47
1:B:54:TYR:CE1	1:B:60:ILE:HD11	2.50	0.47
1:D:526:ILE:HG12	1:D:552:TYR:HB2	1.96	0.47
1:B:302:HIS:O	1:B:434:ARG:HD2	2.13	0.47
1:C:399:ARG:HH11	1:C:403:ASN:ND2	2.12	0.47
1:D:74:ARG:N	1:D:75:PRO:HD2	2.29	0.47
1:B:54:TYR:HE1	1:B:60:ILE:HD11	1.78	0.47
1:B:517:GLU:O	1:B:521:MET:HG2	2.13	0.47
1:B:72:GLU:OE2	1:B:161:HIS:HE1	1.98	0.47
1:D:163:ILE:HB	1:D:186:VAL:HG12	1.97	0.47
1:A:628:GLU:HG2	1:A:629:GLU:N	2.30	0.47
1:C:126:VAL:HG21	1:C:177:PRO:HB3	1.97	0.47
1:A:322:GLU:HB2	1:A:326:LYS:HG2	1.97	0.46
1:A:634:ASN:HB2	1:A:637:ALA:H	1.80	0.46
1:B:133:ASN:N	1:B:133:ASN:ND2	2.62	0.46
1:B:507:TYR:HB2	1:B:556:ARG:NH2	2.30	0.46
1:B:17:VAL:HG12	1:B:18:ALA:N	2.29	0.46
1:B:493:ASP:HB2	1:B:521:MET:HE1	1.97	0.46
1:D:19:ASN:HD22	1:D:19:ASN:N	2.11	0.46
1:B:129:PRO:HB2	1:B:229:HIS:HB3	1.98	0.46
1:B:408:GLU:HG2	4:B:799:HOH:O	2.15	0.46
1:D:12:GLU:OE2	1:D:168:HIS:HE1	1.98	0.46
1:D:436:GLU:HA	1:D:436:GLU:OE1	2.15	0.46
1:D:615:ARG:NH1	1:D:619:PRO:HB3	2.30	0.46
1:D:631:ASN:HB3	1:D:637:ALA:HB1	1.97	0.46
1:B:125:LEU:O	1:B:181:LYS:NZ	2.49	0.46
1:C:213:LEU:HD11	1:C:257:HIS:HB2	1.98	0.46
1:D:168:HIS:HD2	1:D:193:HIS:NE2	2.14	0.46
1:A:119:LYS:HE3	1:A:130:SER:OG	2.16	0.46
1:A:8:HIS:HB2	1:A:162:ALA:O	2.15	0.46
1:B:111:VAL:CG1	1:B:118:TRP:HH2	2.28	0.46
1:B:458:LYS:HD3	1:B:458:LYS:C	2.34	0.46
1:C:32:ALA:HB3	1:C:33:PRO:HD3	1.98	0.46
1:C:5:LEU:HD22	1:C:614:LEU:HD22	1.97	0.46
1:A:99:GLY:O	1:A:100:ALA:C	2.54	0.46
1:B:349:LYS:HG2	1:B:576:VAL:HG13	1.98	0.46
1:C:196:LEU:O	1:C:200:TYR:HD2	1.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:283:GLN:HG3	2:C:901:G6P:C1	2.46	0.46
1:D:74:ARG:N	1:D:75:PRO:CD	2.79	0.46
1:A:199:ARG:HD2	4:A:871:HOH:O	2.16	0.46
1:B:269:ASN:HB2	1:B:511:TRP:CD1	2.50	0.46
1:D:304:CYS:SG	1:D:434:ARG:HD3	2.56	0.46
1:D:615:ARG:HD3	1:D:622:PHE:HD1	1.77	0.46
1:C:207:PHE:CE2	1:C:209:PHE:HA	2.51	0.45
1:C:25:TYR:CE1	1:C:95:TRP:HZ2	2.34	0.45
1:D:279:PHE:CE1	1:D:591:GLU:OE1	2.69	0.45
1:A:285:LEU:O	1:A:289:LYS:HG2	2.17	0.45
1:B:111:VAL:HG12	1:B:111:VAL:O	2.16	0.45
1:C:337:ARG:HH21	1:C:566:GLU:HG2	1.81	0.45
1:D:450:ASP:OD1	1:D:460:ARG:NH2	2.49	0.45
1:D:54:TYR:HE1	1:D:60:ILE:HD11	1.82	0.45
1:A:74:ARG:N	1:A:75:PRO:CD	2.79	0.45
1:B:391:LYS:HZ2	3:B:1001:PEG:H22	1.80	0.45
1:D:218:VAL:HG23	1:D:219:ASP:N	2.32	0.45
1:D:490:LEU:HD22	1:D:494:GLU:HB3	1.97	0.45
1:B:206:SER:O	1:B:207:PHE:HB3	2.16	0.45
1:B:400:TYR:CD1	1:B:408:GLU:HA	2.51	0.45
1:B:634:ASN:ND2	1:B:634:ASN:C	2.70	0.45
1:B:459:ILE:H	1:B:459:ILE:HG12	1.49	0.45
1:C:458:LYS:C	1:C:458:LYS:HD3	2.36	0.45
1:C:71:ASP:O	1:C:74:ARG:HG2	2.17	0.45
1:D:137:THR:HG21	1:D:229:HIS:CD2	2.49	0.45
1:A:174:VAL:O	1:A:177:PRO:HG2	2.17	0.45
1:B:51:LYS:HE3	1:B:107:ASP:OD1	2.16	0.45
1:C:108:LEU:HB3	1:C:142:LEU:HD13	1.99	0.45
1:C:565:VAL:O	1:C:569:VAL:HG23	2.16	0.45
1:D:177:PRO:HA	1:D:240:SER:OG	2.16	0.45
1:D:366:VAL:O	1:D:370:LYS:HB2	2.17	0.45
1:D:485:ASN:HA	1:D:486:PRO:HD3	1.76	0.45
1:B:254:GLU:HG3	1:B:258:LEU:HD12	1.98	0.45
1:D:634:ASN:HB3	1:D:636:ASP:H	1.80	0.45
1:D:542:GLU:CD	1:D:544:ASN:H	2.20	0.45
1:A:268:PRO:HB2	1:A:602:MET:CE	2.47	0.44
1:A:382:VAL:O	1:A:386:THR:HG23	2.17	0.44
1:D:273:VAL:HB	1:D:598:ASP:OD1	2.17	0.44
1:A:94:ARG:HD2	1:A:100:ALA:HB1	1.99	0.44
1:A:213:LEU:C	1:A:213:LEU:HD12	2.38	0.44
1:B:512:GLY:O	1:B:515:PRO:HD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:372[A]:GLN:HE22	1:C:376:ARG:HE	1.66	0.44
1:D:31:LYS:HZ3	1:D:35:THR:HG21	1.82	0.44
1:D:428:ARG:HA	1:D:428:ARG:HD2	1.58	0.44
1:A:534:GLY:O	1:A:538:GLU:HB2	2.17	0.44
1:A:612:LEU:CD2	1:A:635:MET:HG2	2.48	0.44
1:C:249:GLN:NE2	1:C:249:GLN:HA	2.32	0.44
1:C:542:GLU:OE1	1:C:545:GLN:CG	2.65	0.44
1:D:581:ARG:HG3	4:D:724:HOH:O	2.15	0.44
1:A:113:GLY:C	1:A:115:SER:H	2.20	0.44
1:A:510:PRO:O	1:A:532:GLY:HA3	2.18	0.44
1:B:547:LYS:HG2	1:B:552:TYR:CE1	2.53	0.44
1:C:305:PHE:HZ	1:C:309:LEU:HG	1.81	0.44
1:B:111:VAL:HG11	1:B:118:TRP:HH2	1.82	0.44
1:B:74:ARG:N	1:B:75:PRO:CD	2.80	0.44
1:C:129:PRO:O	1:C:229:HIS:HB2	2.18	0.44
1:C:271:LEU:HD13	1:C:520:VAL:HG21	1.99	0.44
1:B:16:GLU:OE2	1:B:16:GLU:N	2.51	0.44
1:B:32:ALA:HB3	1:B:33:PRO:HD3	1.99	0.44
1:B:368:ALA:HB1	1:B:487:ILE:HD13	1.98	0.44
1:C:100:ALA:N	1:C:101:PRO:HD3	2.33	0.44
1:B:323:TYR:OH	1:B:458:LYS:CG	2.66	0.44
1:C:126:VAL:HG12	1:C:126:VAL:O	2.18	0.44
1:C:222:ALA:HB1	1:C:228:TYR:HA	1.99	0.44
1:D:128:ILE:HA	1:D:129:PRO:HD2	1.91	0.44
1:D:19:ASN:ND2	1:D:19:ASN:N	2.66	0.44
1:D:274:ILE:O	1:D:274:ILE:HG23	2.17	0.44
1:B:208:ASP:CG	1:B:211:ASN:HB2	2.38	0.44
1:C:92:TYR:OH	1:C:102:LYS:HD3	2.17	0.44
1:D:90:PHE:HB3	1:D:106:PHE:HD1	1.82	0.44
1:D:192:THR:CG2	1:D:246:THR:HG22	2.45	0.44
1:D:63:TRP:CZ3	1:D:80:LEU:HB3	2.53	0.44
1:B:119:LYS:HE3	1:B:130:SER:HB2	1.99	0.44
1:D:75:PRO:HG2	1:D:158:ASP:OD2	2.18	0.44
1:A:213:LEU:HA	1:A:216:VAL:CG1	2.47	0.43
1:A:83:MET:CE	1:A:153:GLU:HG3	2.48	0.43
1:B:349:LYS:O	1:B:471:ARG:HG3	2.18	0.43
1:B:79:ALA:O	1:B:83:MET:HB2	2.18	0.43
1:A:95:TRP:HB3	1:A:101:PRO:HD2	1.99	0.43
1:B:12:GLU:HB3	1:B:45:LEU:HD23	2.00	0.43
1:D:565:VAL:O	1:D:569:VAL:HG23	2.18	0.43
1:B:549:TYR:HB3	1:B:593:LEU:HD11	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:618:TYR:N	1:C:619:PRO:HD3	2.32	0.43
1:D:349:LYS:O	1:D:471:ARG:CD	2.66	0.43
1:A:180:ARG:HD2	1:A:240:SER:O	2.18	0.43
1:C:378:LEU:O	1:C:382:VAL:HG23	2.18	0.43
1:D:19:ASN:ND2	1:D:19:ASN:H	2.16	0.43
1:D:94:ARG:HB2	1:D:94:ARG:HE	1.64	0.43
1:C:9:LEU:HD13	1:C:161:HIS:CG	2.53	0.43
1:D:471:ARG:HA	1:D:471:ARG:NE	2.33	0.43
1:D:620:ASP:O	1:D:624:GLU:HG2	2.18	0.43
1:C:40:LYS:HB3	1:C:41:ASP:H	1.50	0.43
1:C:509:GLU:OE2	1:C:531:SER:HB2	2.19	0.43
1:A:65:LYS:HA	1:A:66:PRO:HD3	1.83	0.43
1:B:440:PRO:HA	1:B:441:PRO:HD3	1.85	0.43
1:A:526:ILE:HG12	1:A:552:TYR:HB2	2.01	0.43
1:C:428:ARG:HD3	1:C:428:ARG:HA	1.75	0.43
1:C:580:ARG:HB3	1:D:277:GLN:NE2	2.34	0.43
1:C:199:ARG:HG3	1:C:508:TYR:HE2	1.80	0.43
1:D:302:HIS:CD2	1:D:432:LEU:O	2.63	0.43
1:D:337:ARG:HH21	1:D:566:GLU:HG2	1.83	0.43
1:D:54:TYR:CE1	1:D:60:ILE:HD11	2.54	0.43
1:A:86:ARG:HH11	1:A:86:ARG:HA	1.84	0.42
1:C:601:ARG:HH21	1:C:644:LEU:HD23	1.83	0.42
1:D:12:GLU:HB3	1:D:45:LEU:HD23	2.01	0.42
1:D:78:HIS:HB2	1:D:157:LEU:HD13	2.00	0.42
1:D:604:LEU:O	1:D:607:VAL:HB	2.19	0.42
1:A:214:GLU:HG2	1:A:257:HIS:NE2	2.34	0.42
1:B:517:GLU:H	1:B:517:GLU:HG2	1.72	0.42
1:B:534:GLY:O	1:B:538:GLU:HB2	2.19	0.42
1:B:542:GLU:OE1	1:B:544:ASN:HB2	2.19	0.42
1:C:59:ASP:HB2	1:C:96:LEU:CD2	2.40	0.42
1:D:19:ASN:HB3	1:D:50:ASN:HD22	1.83	0.42
1:D:400:TYR:CD1	1:D:401:PRO:HA	2.54	0.42
1:C:14:ALA:O	1:C:17:VAL:HG23	2.19	0.42
1:C:326:LYS:HE3	4:C:804:HOH:O	2.20	0.42
1:C:634:ASN:HB2	1:C:637:ALA:CB	2.49	0.42
1:D:47:GLY:O	1:D:105:LEU:HA	2.19	0.42
1:D:111:VAL:HG13	1:D:118:TRP:CH2	2.54	0.42
1:D:176:LEU:HD22	1:D:241:ALA:HB2	2.01	0.42
1:D:358:PRO:HG2	1:D:480:PHE:CZ	2.54	0.42
1:A:192:THR:HG22	1:A:246:THR:HG22	2.01	0.42
1:B:17:VAL:HG21	1:B:46:ILE:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:292:LYS:HD3	1:C:490:LEU:HD21	2.00	0.42
1:D:302:HIS:O	1:D:434:ARG:HD2	2.20	0.42
1:D:321:TYR:CZ	1:D:455:ILE:HG12	2.55	0.42
1:A:34:ILE:HD13	1:A:599:TRP:HB3	2.02	0.42
1:A:62:ASP:OD2	1:A:65:LYS:HG3	2.18	0.42
1:D:309:LEU:HA	1:D:309:LEU:HD23	1.90	0.42
1:A:326:LYS:HA	1:A:326:LYS:HD3	1.82	0.42
1:C:493:ASP:HB2	1:C:521:MET:HE1	2.02	0.42
1:C:626:VAL:HG11	1:C:630:LEU:HD11	2.01	0.42
1:D:292:LYS:HD2	1:D:490:LEU:CD2	2.43	0.42
1:D:5:LEU:HD13	1:D:622:PHE:HD2	1.84	0.42
1:C:482:ASN:OD1	1:C:484:ASN:HB2	2.18	0.42
1:D:283:GLN:HG3	2:D:901:G6P:C1	2.49	0.42
1:A:419:SER:O	1:A:423:VAL:HG23	2.19	0.42
1:B:474:MET:HB2	1:B:474:MET:HE3	1.78	0.42
1:C:545:GLN:HE21	1:C:647:ALA:HA	1.85	0.42
1:D:13:THR:HB	1:D:167:PHE:CD1	2.55	0.42
1:B:125:LEU:HD11	4:B:738:HOH:O	2.19	0.42
1:D:364:PHE:CE2	1:D:486:PRO:HD2	2.55	0.42
1:B:323:TYR:OH	1:B:458:LYS:HG2	2.20	0.42
1:C:542:GLU:OE1	1:C:542:GLU:C	2.58	0.42
1:D:90:PHE:HB3	1:D:106:PHE:CD1	2.54	0.42
1:C:634:ASN:CB	1:C:637:ALA:H	2.30	0.41
1:A:396:HIS:CE1	1:A:405:LEU:HD22	2.55	0.41
1:C:95:TRP:HB3	1:C:101:PRO:HD2	2.02	0.41
1:C:357:MET:HA	1:C:358:PRO:HD3	1.92	0.41
1:C:590:THR:HB	4:C:805:HOH:O	2.20	0.41
1:B:326:LYS:HD3	1:B:326:LYS:HA	1.86	0.41
1:C:14:ALA:HB3	1:C:28:LEU:HD11	2.03	0.41
1:B:372:GLN:HE21	1:B:372:GLN:HB3	1.69	0.41
1:D:456:LEU:O	1:D:460:ARG:HG3	2.21	0.41
1:D:349:LYS:O	1:D:471:ARG:HD2	2.20	0.41
1:B:48:PRO:HG3	1:B:143:LEU:HD22	2.03	0.41
1:B:292:LYS:HD2	1:B:490:LEU:HD21	2.02	0.41
1:B:561:PRO:O	1:B:565:VAL:HG23	2.20	0.41
1:C:617:GLY:C	1:C:619:PRO:HD3	2.40	0.41
1:D:283:GLN:HG3	2:D:901:G6P:H1	2.03	0.41
1:D:29:LYS:HG3	1:D:97:ILE:HD12	2.01	0.41
1:D:268:PRO:HB2	1:D:602:MET:CE	2.51	0.41
1:A:11:PHE:HD1	1:A:46:ILE:HD11	1.85	0.41
1:B:9:LEU:HD13	1:B:161:HIS:CG	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:434:ARG:HB2	1:C:435:PRO:HD2	2.03	0.41
1:C:471:ARG:CZ	1:C:471:ARG:HA	2.50	0.41
1:C:512:GLY:C	1:C:515:PRO:HD2	2.40	0.41
1:D:11:PHE:CD1	1:D:44:HIS:HB2	2.56	0.41
1:D:283:GLN:HG3	2:D:901:G6P:O1	2.20	0.41
1:C:177:PRO:HA	1:C:240:SER:OG	2.21	0.41
1:C:74:ARG:HB2	1:C:75:PRO:HD3	2.02	0.41
1:D:321:TYR:OH	1:D:455:ILE:HG12	2.21	0.41
1:B:326:LYS:NZ	1:B:509:GLU:HG3	2.35	0.41
1:D:371:GLY:HA3	4:D:739:HOH:O	2.21	0.41
1:D:538:GLU:HB3	1:D:553:ILE:HD12	2.02	0.41
1:A:621:GLN:O	1:A:625:LEU:HD13	2.21	0.41
1:D:283:GLN:NE2	1:D:587:ARG:NH2	2.65	0.41
1:A:19:ASN:ND2	1:A:19:ASN:O	2.54	0.41
1:B:497:ARG:NH2	4:B:777:HOH:O	2.54	0.41
1:C:282:PHE:HB2	1:C:497:ARG:HH21	1.84	0.41
1:D:125:LEU:HA	1:D:125:LEU:HD23	1.83	0.41
1:D:17:VAL:O	1:D:17:VAL:CG1	2.69	0.41
1:D:389:ILE:HG23	1:D:416:LEU:HB3	2.03	0.41
1:A:17:VAL:HG22	1:A:47:GLY:HA3	2.03	0.40
1:A:56:ASN:N	1:A:56:ASN:OD1	2.53	0.40
1:B:196:LEU:O	1:B:200:TYR:HD2	2.04	0.40
1:B:61:LEU:HG	1:B:93:GLY:HA2	2.04	0.40
1:D:214:GLU:HG3	1:D:257:HIS:CD2	2.56	0.40
1:D:218:VAL:CG1	1:D:260:LYS:HE2	2.50	0.40
1:D:333:GLU:OE2	1:D:337:ARG:HD2	2.21	0.40
1:C:111:VAL:CG1	1:C:118:TRP:CH2	3.04	0.40
1:C:542:GLU:OE1	1:C:544:ASN:N	2.54	0.40
1:D:227:ILE:O	1:D:227:ILE:CG2	2.68	0.40
1:D:615:ARG:O	1:D:615:ARG:HD2	2.21	0.40
1:A:624:GLU:C	1:A:626:VAL:N	2.73	0.40
1:B:214:GLU:HA	1:B:257[B]:HIS:CE1	2.56	0.40
1:B:358:PRO:HG2	1:B:480:PHE:CZ	2.57	0.40
1:C:188:THR:O	1:C:242:ASP:HB2	2.21	0.40
1:C:197:LEU:HB2	1:C:258:LEU:HD13	2.02	0.40
1:D:189:ILE:HD11	1:D:610:ARG:HA	2.03	0.40
1:D:274:ILE:CG2	1:D:274:ILE:O	2.70	0.40
1:A:214:GLU:HG2	1:A:257:HIS:CD2	2.56	0.40
1:B:455:ILE:HG22	1:B:459:ILE:HD11	2.02	0.40
1:C:60:ILE:H	1:C:60:ILE:HG13	1.73	0.40
1:D:454:LEU:HA	1:D:454:LEU:HD23	1.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:LEU:HD21	1:B:254:GLU:HA	2.03	0.40
1:D:372:GLN:HG3	1:D:373:ALA:N	2.35	0.40
1:D:440:PRO:HA	1:D:441:PRO:HD3	1.96	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	640/725 (88%)	606 (95%)	32 (5%)	2 (0%)	41	55
1	B	646/725 (89%)	622 (96%)	22 (3%)	2 (0%)	41	55
1	C	646/725 (89%)	618 (96%)	25 (4%)	3 (0%)	29	41
1	D	633/725 (87%)	597 (94%)	32 (5%)	4 (1%)	25	36
All	All	2565/2900 (88%)	2443 (95%)	111 (4%)	11 (0%)	34	48

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	625	LEU
1	B	17	VAL
1	B	643	LYS
1	C	543	THR
1	D	111	VAL
1	C	363	SER
1	D	40	LYS
1	D	194	ALA
1	C	17	VAL
1	D	115	SER
1	A	627	GLY

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	555/623 (89%)	538 (97%)	17 (3%)	40	60
1	B	558/623 (90%)	533 (96%)	25 (4%)	27	44
1	C	558/623 (90%)	538 (96%)	20 (4%)	35	54
1	D	550/623 (88%)	530 (96%)	20 (4%)	35	54
All	All	2221/2492 (89%)	2139 (96%)	82 (4%)	35	53

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	19	ASN
1	A	56	ASN
1	A	83	MET
1	A	110	SER
1	A	124	SER
1	A	180	ARG
1	A	283	GLN
1	A	310	ASP
1	A	321	TYR
1	A	372[A]	GLN
1	A	372[B]	GLN
1	A	376	ARG
1	A	458	LYS
1	A	556	ARG
1	A	568	LEU
1	A	590	THR
1	B	19	ASN
1	B	40	LYS
1	B	133	ASN
1	B	199	ARG
1	B	216	VAL
1	B	220	HIS
1	B	289	LYS

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Mol	Chain	Res	Type
1	B	310	ASP
1	B	320	ARG
1	B	321	TYR
1	B	348	LYS
1	B	372	GLN
1	B	376	ARG
1	B	408	GLU
1	B	420	SER
1	B	458	LYS
1	B	459	ILE
1	B	471	ARG
1	B	484	ASN
1	B	518	CYS
1	B	539	ASP
1	B	590	THR
1	B	614	LEU
1	B	634	ASN
1	B	645	LYS
1	C	19	ASN
1	C	180	ARG
1	C	181	LYS
1	C	213	LEU
1	C	240	SER
1	C	310	ASP
1	C	321	TYR
1	C	372[A]	GLN
1	C	372[B]	GLN
1	C	376	ARG
1	C	458	LYS
1	C	471	ARG
1	C	475	ILE
1	C	505	PRO
1	C	531	SER
1	C	543	THR
1	C	545	GLN
1	C	568	LEU
1	C	590	THR
1	C	645	LYS
1	D	6	GLN
1	D	19	ASN
1	D	41	ASP
1	D	42	HIS

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Mol	Chain	Res	Type
1	D	95	TRP
1	D	122	LEU
1	D	133	ASN
1	D	213	LEU
1	D	310	ASP
1	D	321	TYR
1	D	324	LYS
1	D	330	MET
1	D	372	GLN
1	D	387	THR
1	D	417	LEU
1	D	471	ARG
1	D	482	ASN
1	D	540	LEU
1	D	590	THR
1	D	636	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	283	GLN
1	A	403	ASN
1	A	484	ASN
1	B	7	ASN
1	B	133	ASN
1	B	161	HIS
1	B	396	HIS
1	B	484	ASN
1	B	582	GLN
1	B	621	GLN
1	B	634	ASN
1	C	19	ASN
1	C	81	GLN
1	C	249	GLN
1	C	257	HIS
1	C	284	ASN
1	C	403	ASN
1	C	477	HIS
1	C	545	GLN
1	C	586	GLN
1	D	19	ASN

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Mol	Chain	Res	Type
1	D	50	ASN
1	D	161	HIS
1	D	168	HIS
1	D	211	ASN
1	D	249	GLN
1	D	257	HIS
1	D	283	GLN
1	D	302	HIS
1	D	403	ASN
1	D	482	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

12 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	G6P	A	901	-	16,16,16	0.90	0	24,24,24	1.43	3 (12%)
3	PEG	B	706	-	6,6,6	0.62	0	5,5,5	0.44	0
2	G6P	A	902	-	16,16,16	0.90	0	24,24,24	1.39	4 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	G6P	C	901	-	16,16,16	0.85	0	24,24,24	1.47	3 (12%)
3	PEG	A	1001	-	6,6,6	0.61	0	5,5,5	0.39	0
3	PEG	B	1001	-	6,6,6	0.56	0	5,5,5	0.44	0
3	PEG	C	1001	-	6,6,6	0.61	0	5,5,5	0.39	0
3	PEG	B	1002	-	6,6,6	0.57	0	5,5,5	0.45	0
2	G6P	D	901	-	16,16,16	0.86	0	24,24,24	1.39	4 (16%)
2	G6P	B	902	-	16,16,16	0.91	0	24,24,24	2.13	7 (29%)
3	PEG	D	1002	-	6,6,6	0.63	0	5,5,5	0.46	0
2	G6P	B	901	-	16,16,16	0.89	0	24,24,24	1.42	3 (12%)

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	G6P	A	901	-	-	0/6/26/26	0/1/1/1
3	PEG	B	706	-	-	4/4/4/4	-
2	G6P	A	902	-	-	6/6/26/26	0/1/1/1
2	G6P	C	901	-	-	0/6/26/26	0/1/1/1
3	PEG	A	1001	-	-	4/4/4/4	-
3	PEG	B	1001	-	-	3/4/4/4	-
3	PEG	C	1001	-	-	1/4/4/4	-
3	PEG	B	1002	-	-	3/4/4/4	-
2	G6P	D	901	-	-	0/6/26/26	0/1/1/1
2	G6P	B	902	-	-	6/6/26/26	0/1/1/1
3	PEG	D	1002	-	-	2/4/4/4	-
2	G6P	B	901	-	-	0/6/26/26	0/1/1/1

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	902	G6P	C4-C3-C2	-5.20	101.74	110.82
2	B	902	G6P	P-O6-C6	4.61	130.98	118.30
2	B	902	G6P	O5-C5-C4	-4.13	102.19	109.69
2	A	901	G6P	C4-C3-C2	-3.67	104.42	110.82
2	C	901	G6P	C4-C3-C2	-3.37	104.94	110.82
2	B	901	G6P	C4-C3-C2	-3.05	105.50	110.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	901	G6P	C4-C3-C2	-2.95	105.68	110.82
2	B	901	G6P	P-O6-C6	2.71	125.77	118.30
2	C	901	G6P	P-O6-C6	2.51	125.21	118.30
2	A	902	G6P	C4-C3-C2	-2.50	106.47	110.82
2	B	902	G6P	O5-C1-C2	2.44	114.64	110.28
2	A	902	G6P	C1-O5-C5	-2.41	109.12	113.66
2	B	902	G6P	O6-C6-C5	2.32	116.97	108.99
2	D	901	G6P	P-O6-C6	2.27	124.54	118.30
2	B	901	G6P	O2P-P-O6	-2.23	100.81	106.73
2	C	901	G6P	O1P-P-O3P	2.22	119.36	110.68
2	B	902	G6P	O2P-P-O6	-2.22	100.84	106.73
2	A	902	G6P	P-O6-C6	2.19	124.34	118.30
2	A	901	G6P	O2-C2-C1	-2.18	104.11	109.16
2	A	901	G6P	P-O6-C6	2.09	124.06	118.30
2	A	902	G6P	O2-C2-C1	-2.09	104.31	109.16
2	D	901	G6P	O2P-P-O6	-2.06	101.26	106.73
2	D	901	G6P	O5-C5-C6	-2.06	102.51	106.67
2	B	902	G6P	C3-C4-C5	-2.05	106.57	110.24

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	902	G6P	C4-C5-C6-O6
2	A	902	G6P	O5-C5-C6-O6
2	A	902	G6P	C6-O6-P-O1P
2	A	902	G6P	C6-O6-P-O2P
2	B	902	G6P	C4-C5-C6-O6
2	B	902	G6P	O5-C5-C6-O6
2	B	902	G6P	C5-C6-O6-P
2	B	902	G6P	C6-O6-P-O1P
2	B	902	G6P	C6-O6-P-O2P
2	B	902	G6P	C6-O6-P-O3P
3	C	1001	PEG	O1-C1-C2-O2
3	A	1001	PEG	O2-C3-C4-O4
3	B	1001	PEG	O1-C1-C2-O2
2	A	902	G6P	C6-O6-P-O3P
3	B	706	PEG	O2-C3-C4-O4
3	D	1002	PEG	O1-C1-C2-O2
3	B	1002	PEG	O1-C1-C2-O2
3	B	1002	PEG	O2-C3-C4-O4
3	B	1001	PEG	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
3	B	1001	PEG	O2-C3-C4-O4
3	B	706	PEG	C1-C2-O2-C3
3	B	706	PEG	C4-C3-O2-C2
3	B	706	PEG	O1-C1-C2-O2
2	A	902	G6P	C5-C6-O6-P
3	A	1001	PEG	O1-C1-C2-O2
3	D	1002	PEG	C4-C3-O2-C2
3	A	1001	PEG	C1-C2-O2-C3
3	B	1002	PEG	C4-C3-O2-C2
3	A	1001	PEG	C4-C3-O2-C2

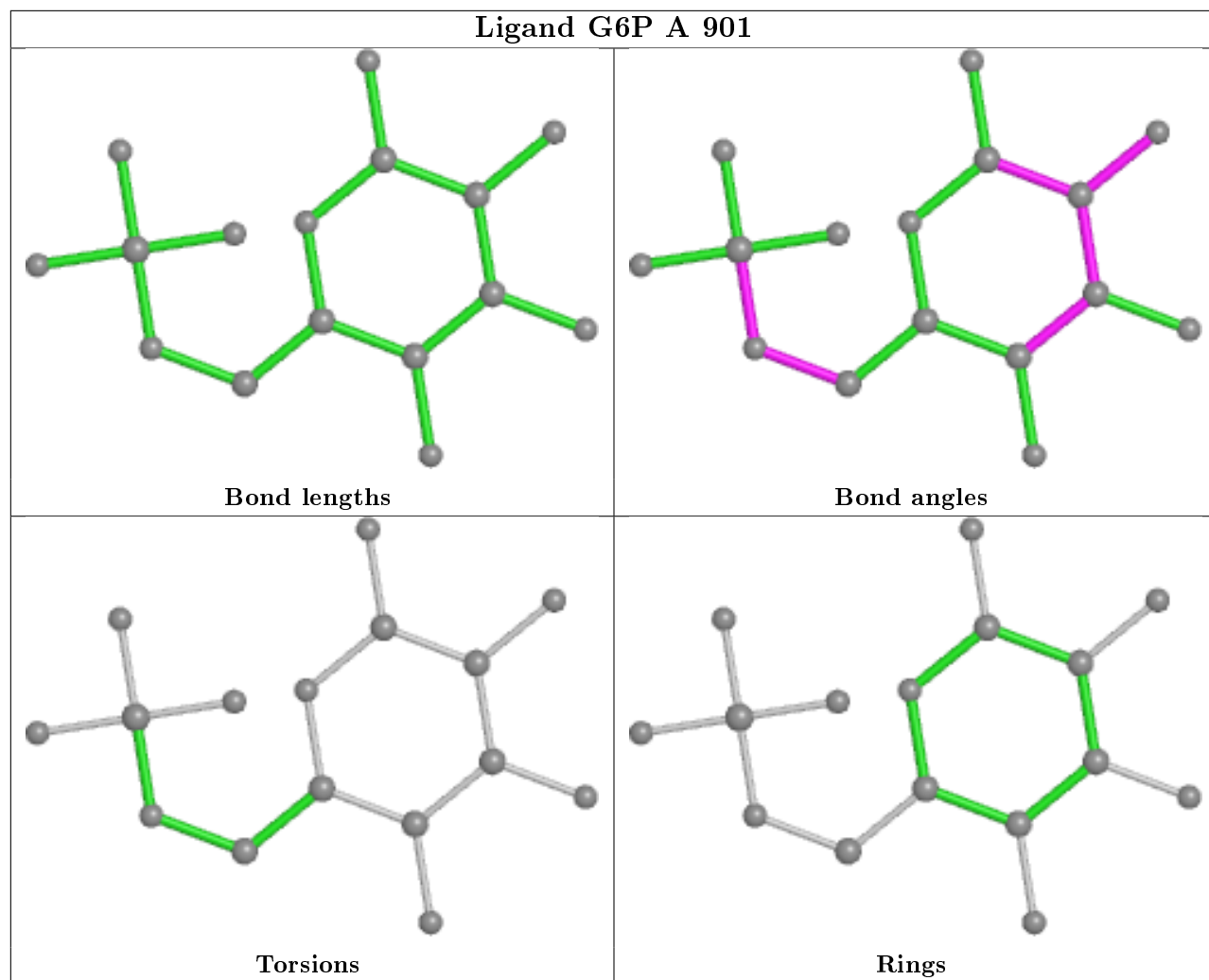
There are no ring outliers.

5 monomers are involved in 10 short contacts:

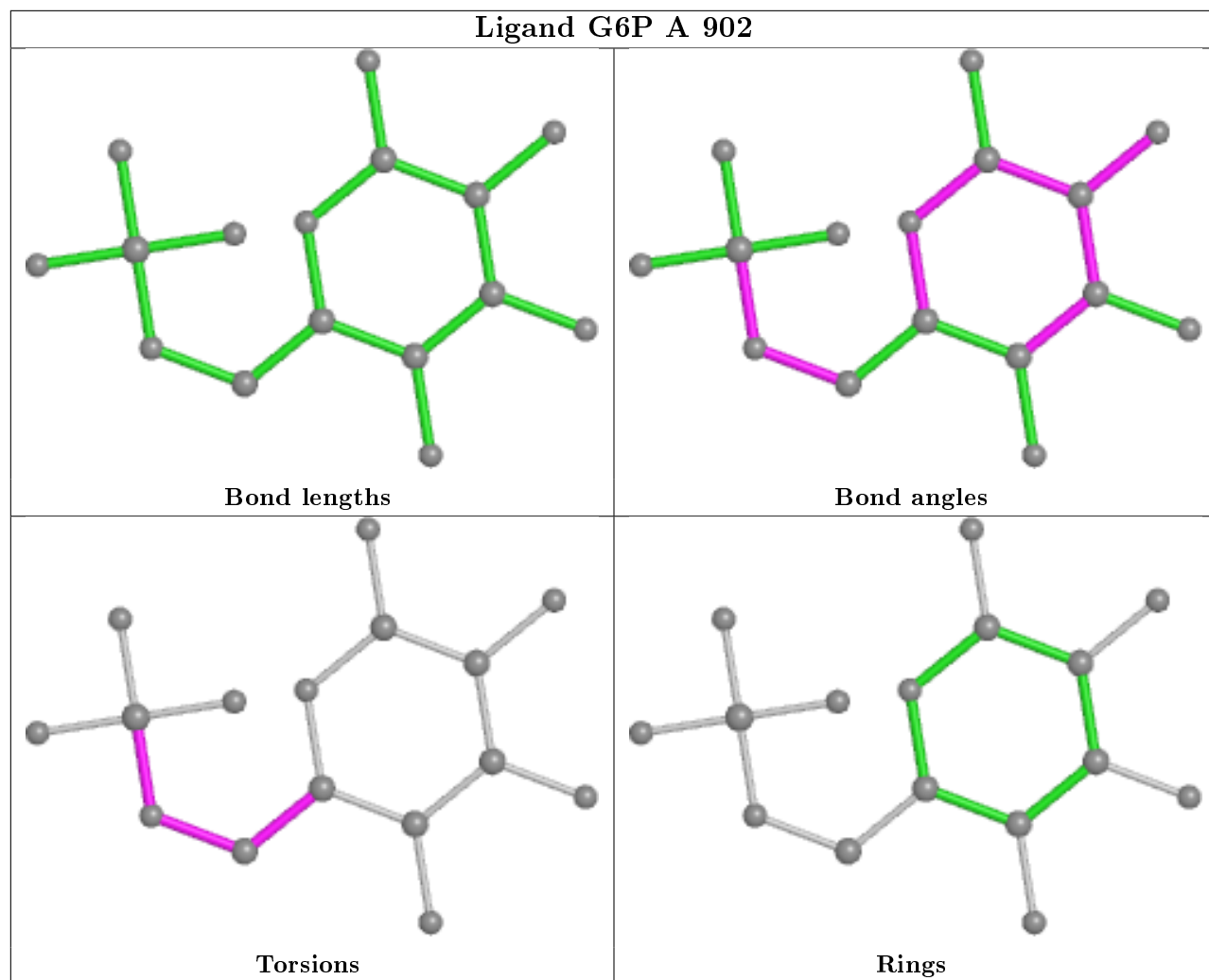
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	902	G6P	1	0
2	C	901	G6P	2	0
3	B	1001	PEG	2	0
2	D	901	G6P	3	0
2	B	902	G6P	2	0

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.

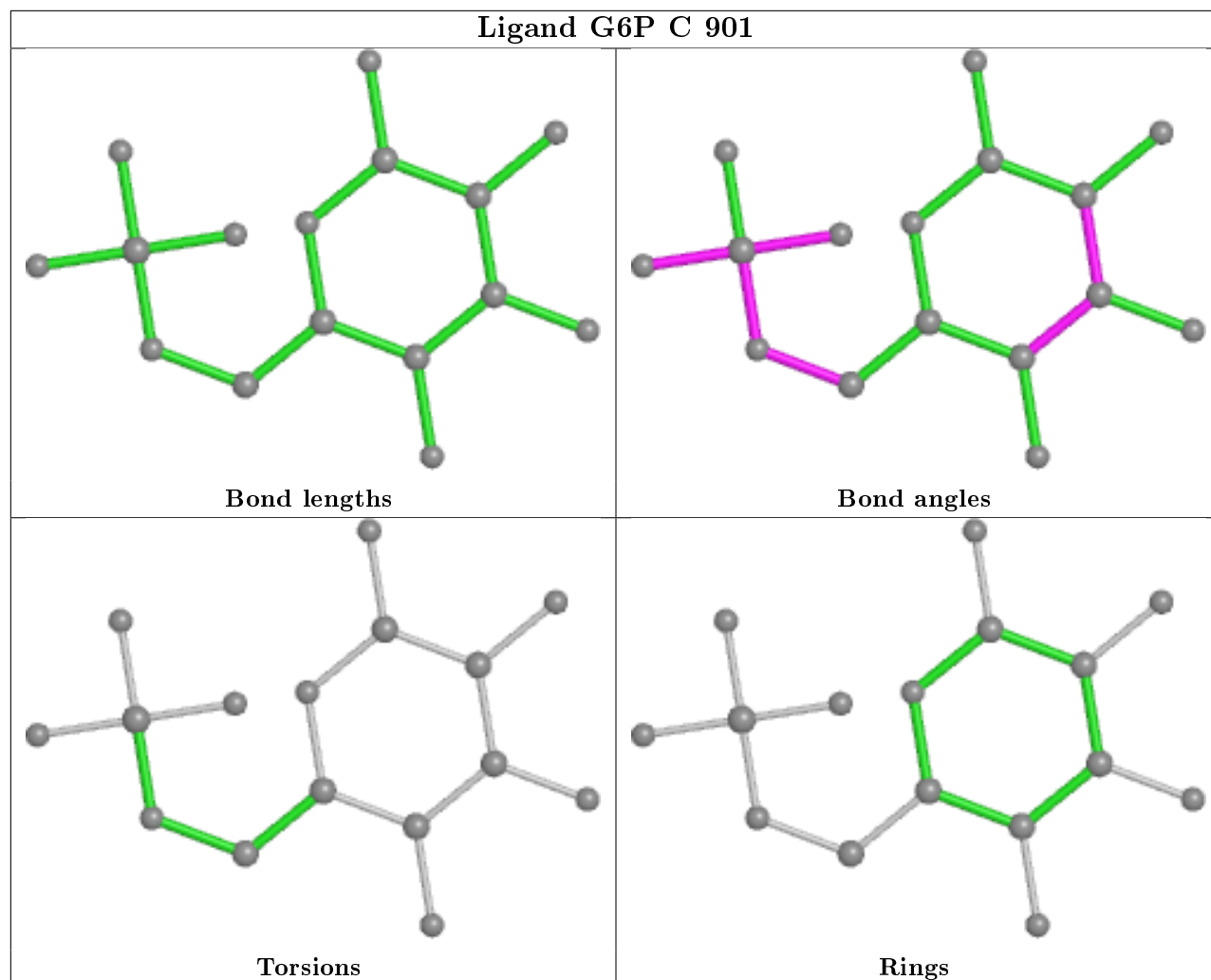
Ligand G6P A 901

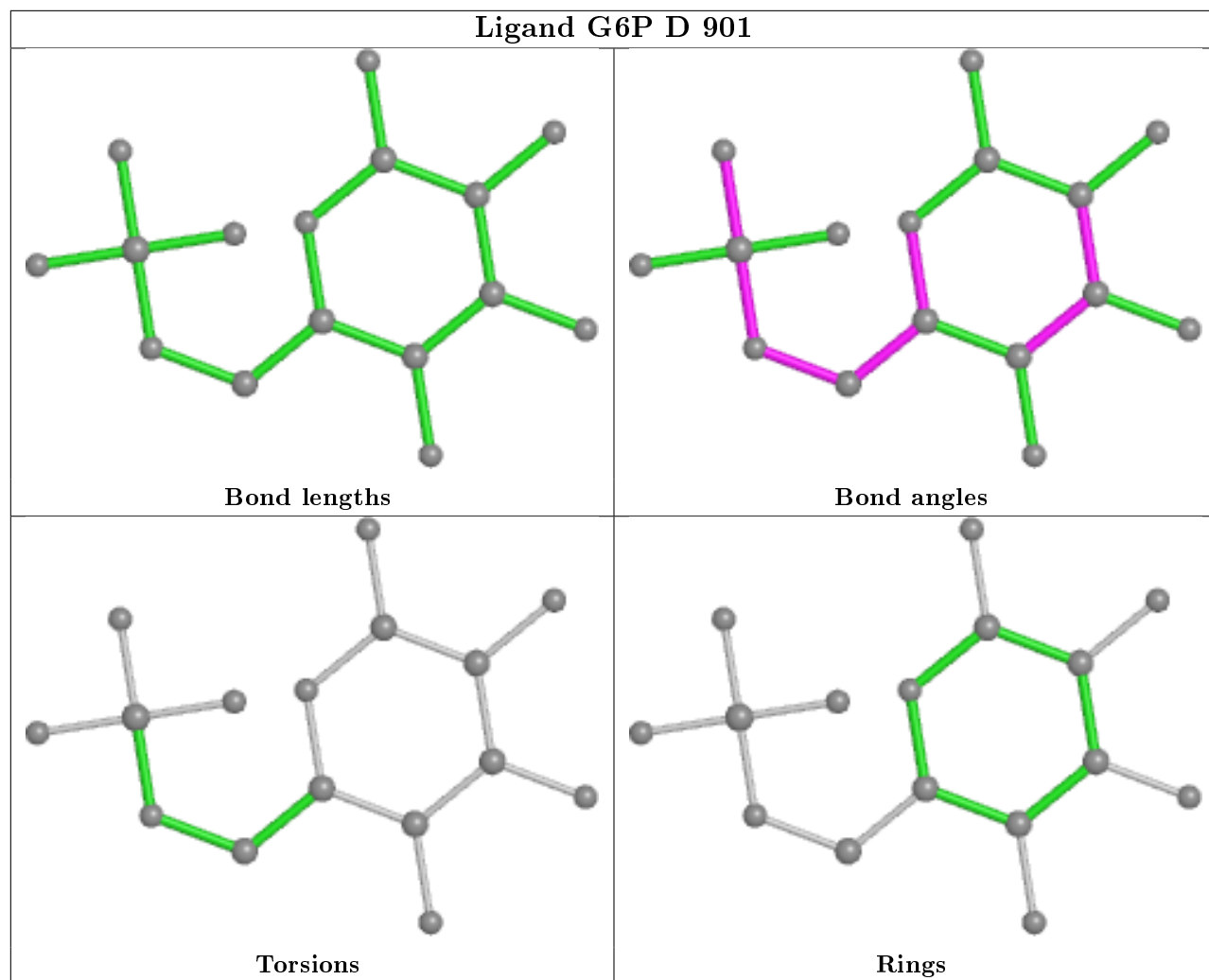


Ligand G6P A 902

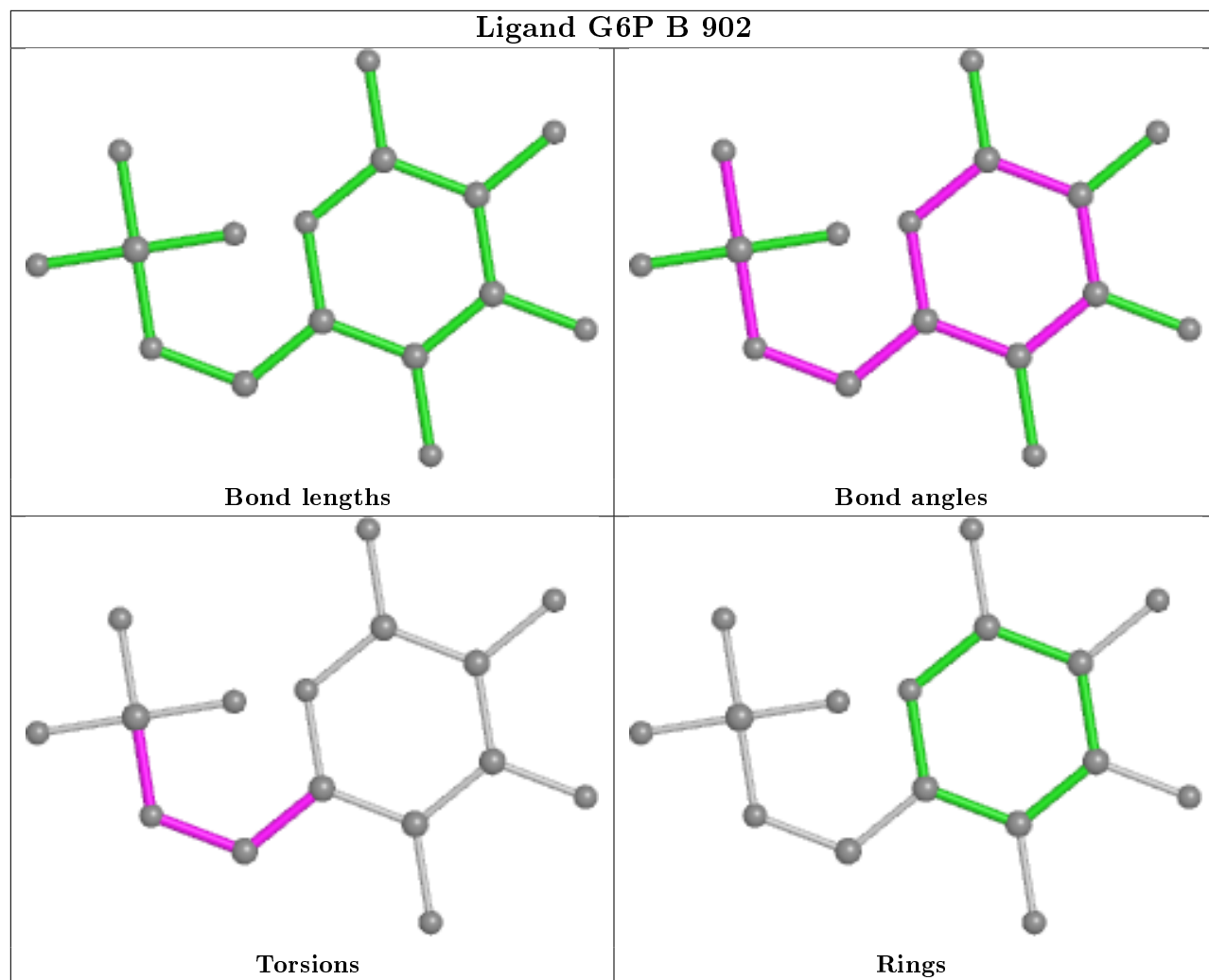


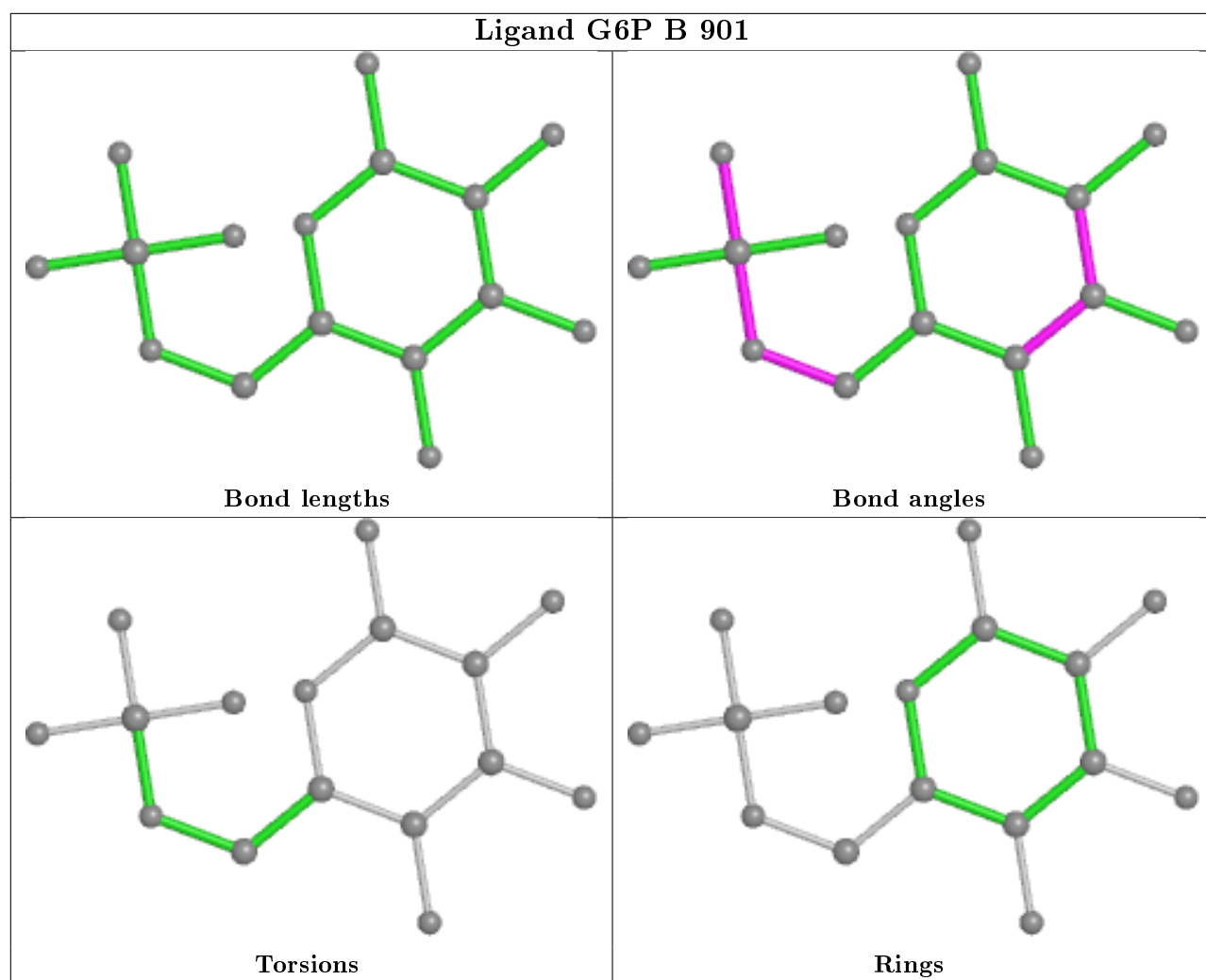
Ligand G6P C 901





Ligand G6P B 902





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	638/725 (88%)	0.17	14 (2%) 62 60	25, 49, 89, 101	0
1	B	645/725 (88%)	0.06	24 (3%) 41 41	33, 50, 83, 96	0
1	C	646/725 (89%)	0.15	23 (3%) 42 42	36, 53, 87, 101	0
1	D	636/725 (87%)	0.29	45 (7%) 16 14	28, 61, 103, 114	0
All	All	2565/2900 (88%)	0.17	106 (4%) 37 36	25, 52, 95, 114	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	630	LEU	8.0
1	D	627	GLY	7.1
1	D	622	PHE	6.5
1	D	126	VAL	6.0
1	D	623	ARG	6.0
1	B	644	LEU	5.9
1	B	646	VAL	5.6
1	D	626	VAL	5.6
1	D	638	LEU	5.5
1	D	125	LEU	5.3
1	B	645	LYS	5.3
1	D	629	GLU	5.2
1	B	128	ILE	5.1
1	D	625	LEU	5.0
1	D	122	LEU	5.0
1	D	133	ASN	4.8
1	B	125	LEU	4.7
1	A	623	ARG	4.2
1	C	645	LYS	4.1
1	D	128	ILE	4.0
1	C	205	GLY	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	628	GLU	3.8
1	D	637	ALA	3.8
1	D	639	ALA	3.7
1	D	624	GLU	3.6
1	D	129	PRO	3.6
1	B	543	THR	3.6
1	A	90	PHE	3.5
1	C	68	ALA	3.5
1	A	624	GLU	3.4
1	D	631	ASN	3.4
1	D	119	LYS	3.3
1	C	62	ASP	3.3
1	D	618	TYR	3.3
1	D	227	ILE	3.2
1	B	127	GLY	3.2
1	A	5	LEU	3.2
1	C	647	ALA	3.1
1	B	641	GLY	3.1
1	A	61	LEU	3.1
1	D	636	ASP	3.1
1	D	633	SER	3.1
1	C	59	ASP	3.0
1	D	632	ASP	3.0
1	B	642	LYS	3.0
1	B	544	ASN	3.0
1	C	133	ASN	2.9
1	A	630	LEU	2.9
1	B	130	SER	2.9
1	D	115	SER	2.9
1	D	2	SER	2.9
1	B	643	LYS	2.9
1	B	123	TRP	2.8
1	D	124	SER	2.8
1	D	132	GLU	2.8
1	A	94	ARG	2.8
1	B	541	ILE	2.8
1	B	126	VAL	2.7
1	A	63	TRP	2.7
1	A	2	SER	2.7
1	B	279	PHE	2.7
1	D	635	MET	2.7
1	C	279	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	132	GLU	2.6
1	D	60	ILE	2.6
1	C	76	VAL	2.6
1	C	207	PHE	2.6
1	D	6	GLN	2.6
1	A	67	GLU	2.5
1	D	197	LEU	2.5
1	D	615	ARG	2.5
1	D	91	VAL	2.5
1	A	132	GLU	2.5
1	D	135	PHE	2.5
1	C	213	LEU	2.5
1	D	85	SER	2.5
1	D	223	GLY	2.5
1	C	111	VAL	2.4
1	B	206	SER	2.4
1	C	641	GLY	2.4
1	C	204	SER	2.4
1	C	644	LEU	2.4
1	D	5	LEU	2.4
1	D	619	PRO	2.3
1	B	253	PHE	2.3
1	B	545	GLN	2.3
1	C	89	HIS	2.3
1	B	181	LYS	2.3
1	B	62	ASP	2.3
1	C	63	TRP	2.2
1	D	123	TRP	2.2
1	B	124	SER	2.2
1	C	83	MET	2.2
1	C	543	THR	2.2
1	D	620	ASP	2.1
1	D	614	LEU	2.1
1	A	637	ALA	2.1
1	C	64	LYS	2.1
1	C	135	PHE	2.1
1	B	113	GLY	2.1
1	A	631	ASN	2.1
1	C	67	GLU	2.1
1	A	636	ASP	2.0
1	C	88	VAL	2.0
1	D	118	TRP	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	621	GLN	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](#)

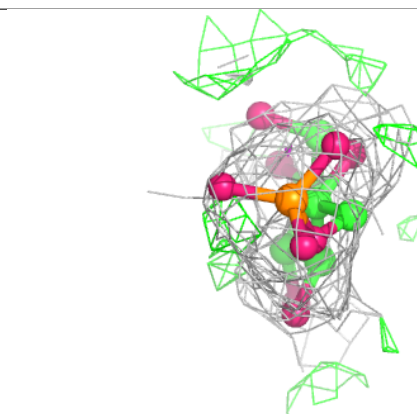
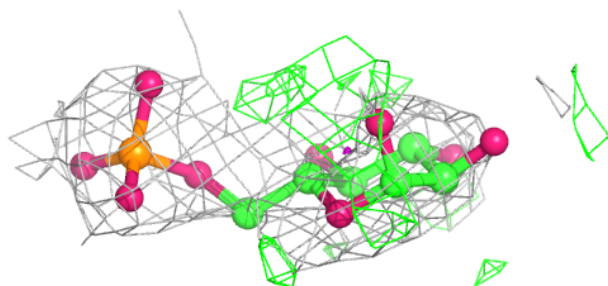
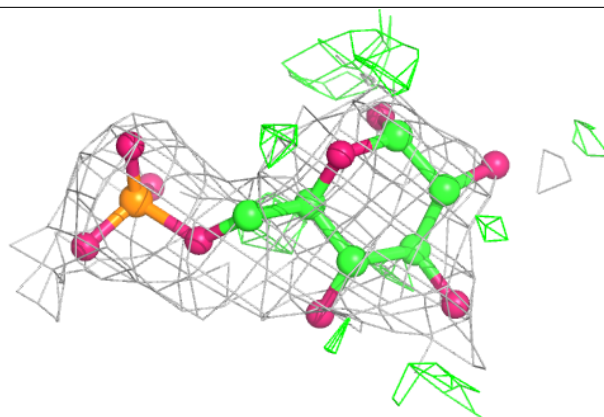
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	PEG	B	1002	7/7	0.80	0.29	60,61,65,67	0
3	PEG	A	1001	7/7	0.85	0.16	49,50,54,56	0
3	PEG	D	1002	7/7	0.86	0.16	47,50,55,58	0
3	PEG	B	1001	7/7	0.87	0.13	42,44,46,50	0
3	PEG	C	1001	7/7	0.90	0.12	43,45,49,50	0
2	G6P	A	902	16/16	0.92	0.21	46,52,58,60	16
2	G6P	B	902	16/16	0.93	0.15	51,54,56,60	0
3	PEG	B	706	7/7	0.93	0.11	43,46,48,49	0
2	G6P	D	901	16/16	0.99	0.14	32,34,36,38	0
2	G6P	A	901	16/16	0.99	0.18	30,35,38,40	0
2	G6P	C	901	16/16	0.99	0.16	39,42,43,44	0
2	G6P	B	901	16/16	0.99	0.13	35,38,42,42	0

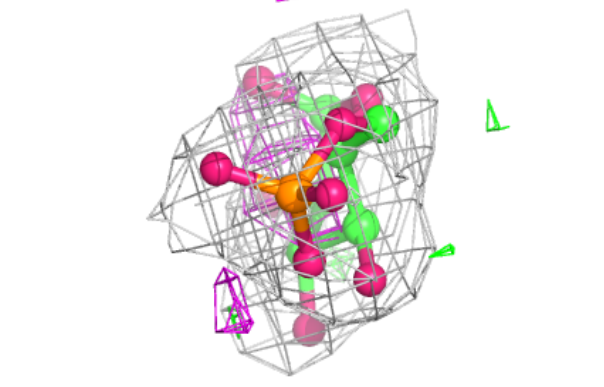
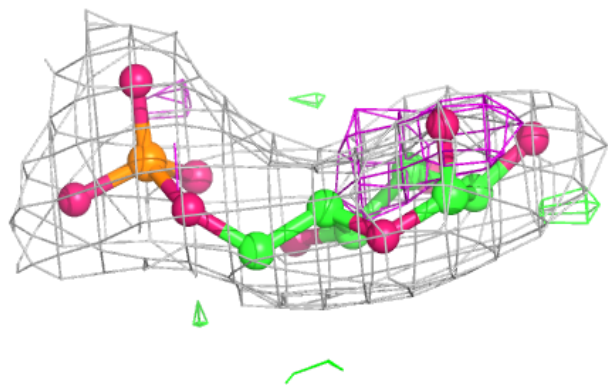
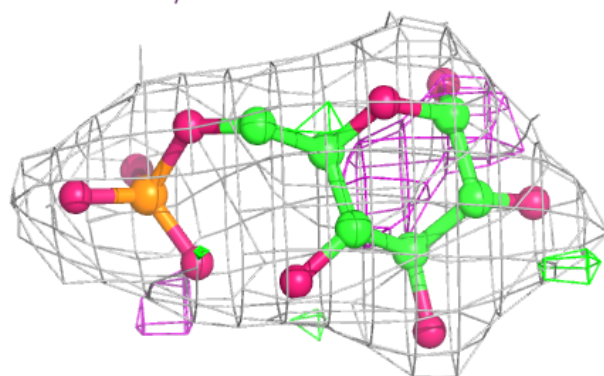
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around G6P A 902:

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

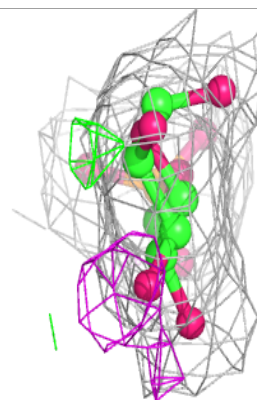
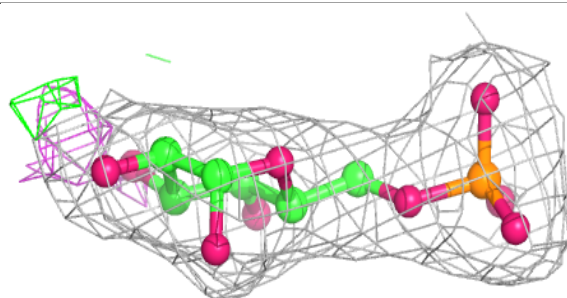
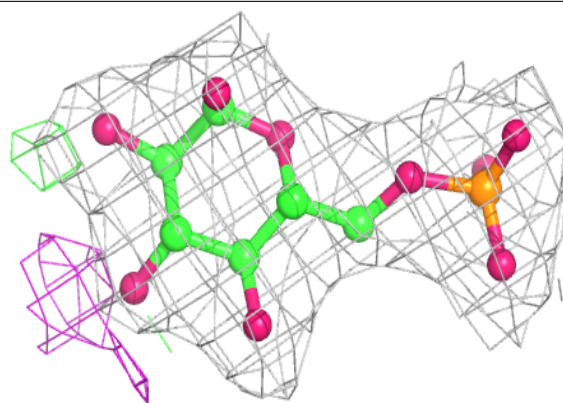
**Electron density around G6P B 902:**

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

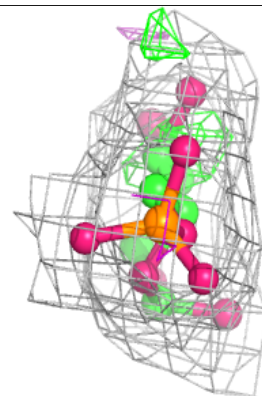
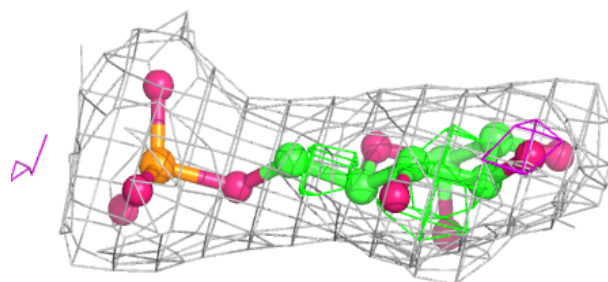
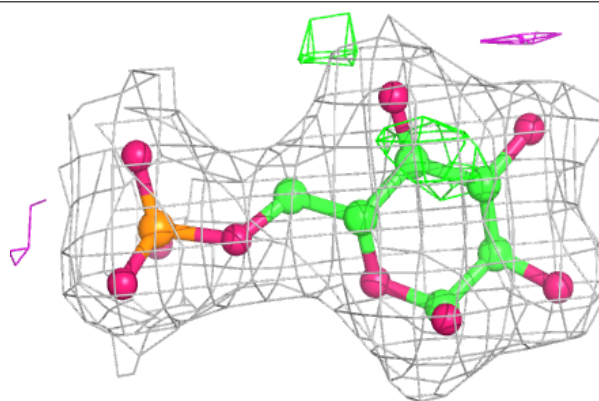


Electron density around G6P D 901:

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

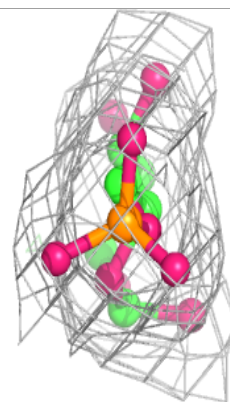
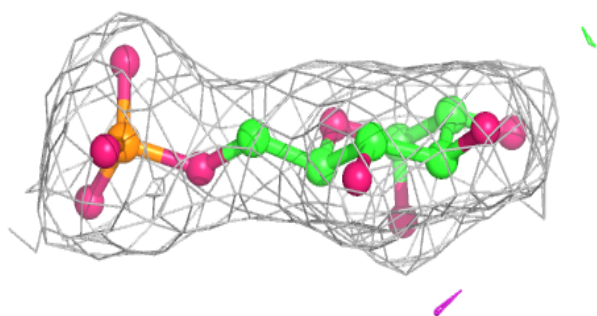
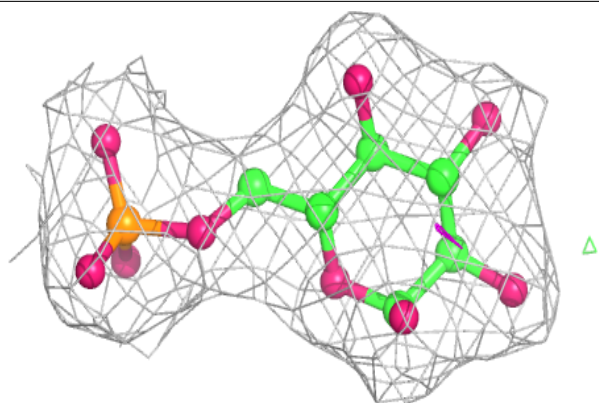
**Electron density around G6P A 901:**

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and green (positive)

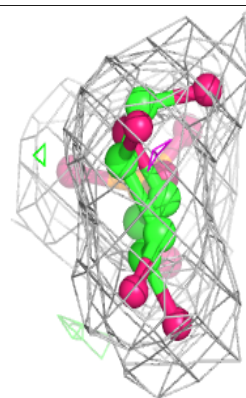
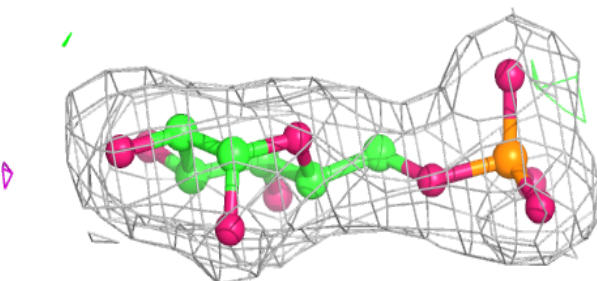
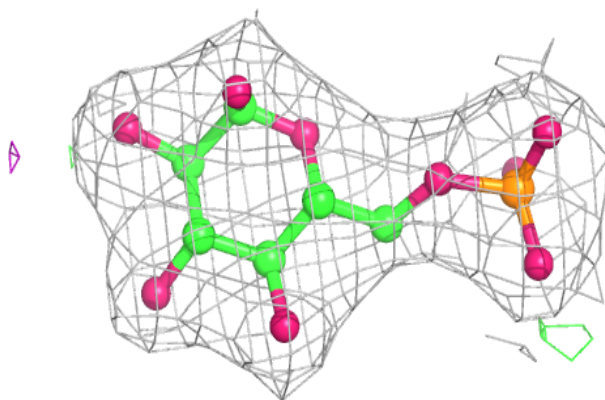


Electron density around G6P C 901:

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

**Electron density around G6P B 901:**

$2mF_o-DF_c$ (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.