



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 03:43 PM UTC

PDB ID : 3NB0 / pdb_00003nb0
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 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

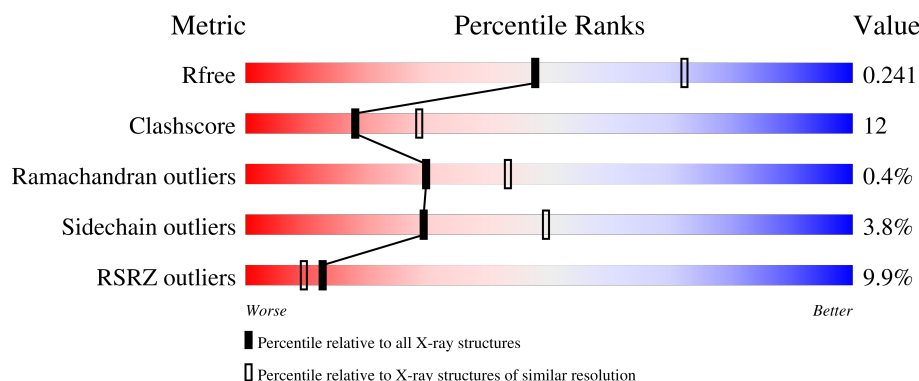
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (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 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	725	
1	B	725	
1	C	725	
1	D	725	

2 Entry composition [i](#)

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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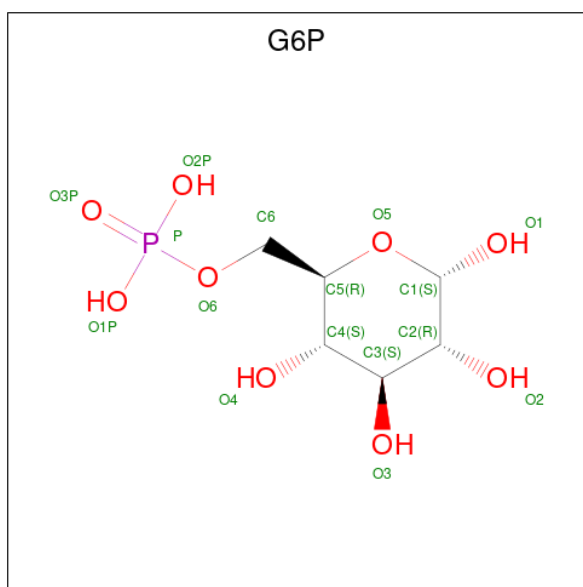
Chain	Residue	Modelled	Actual	Comment	Reference
A	589	ALA	ARG	engineered mutation	UNP P27472
A	592	ALA	ARG	engineered mutation	UNP P27472
B	-19	MET	-	expression tag	UNP P27472
B	-18	GLY	-	expression tag	UNP P27472
B	-17	SER	-	expression tag	UNP P27472
B	-16	SER	-	expression tag	UNP P27472
B	-15	HIS	-	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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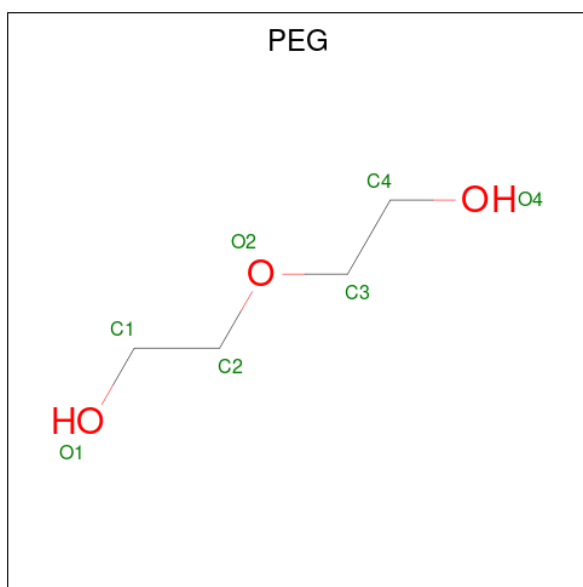
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP P27472
C	-1	SER	-	expression tag	UNP P27472
C	0	HIS	-	expression tag	UNP P27472
C	535	SER	ALA	SEE REMARK 999	UNP P27472
C	589	ALA	ARG	engineered mutation	UNP P27472
C	592	ALA	ARG	engineered mutation	UNP P27472
D	-19	MET	-	expression 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 (CCD ID: G6P) (formula: C₆H₁₃O₉P).



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 (CCD ID: 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		

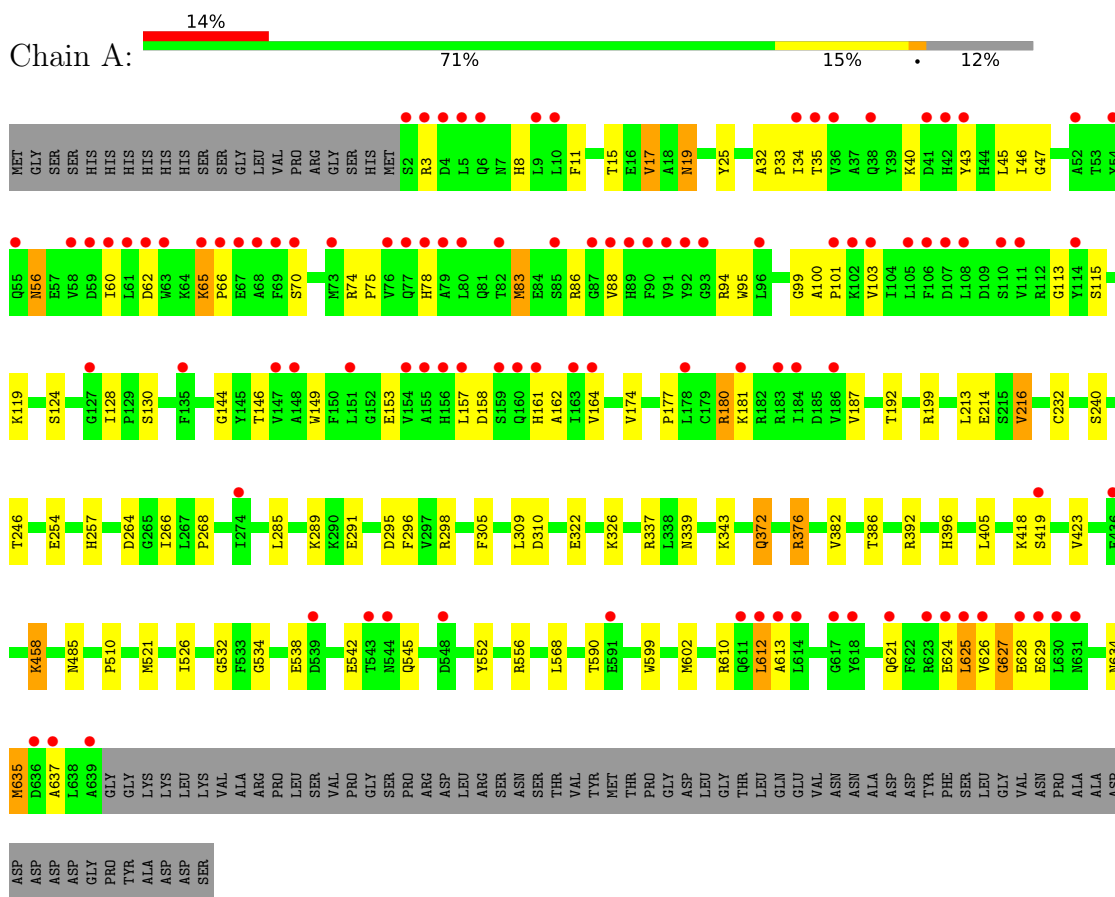
- Molecule 4 is water.

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		

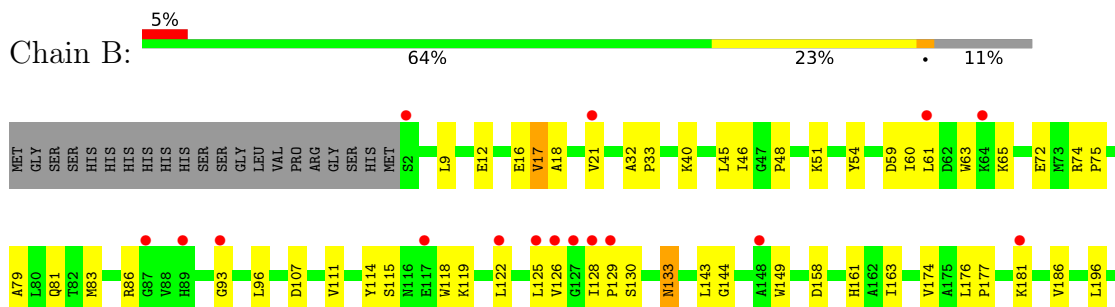
3 Residue-property plots

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: 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 0.211 , 0.241	Depositor DCC
R_{free} test set	7893 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)	0.33 , 39.0	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	0.95	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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, 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	0.32	0/5307	0.70	5/7191 (0.1%)
1	B	0.31	0/5349	0.70	5/7246 (0.1%)
1	C	0.30	0/5345	0.70	4/7239 (0.1%)
1	D	0.31	0/5262	0.70	2/7129 (0.0%)
All	All	0.31	0/21263	0.70	16/28805 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	485	ASN	CA-C-N	6.79	126.48	119.56
1	C	485	ASN	C-N-CA	6.79	126.48	119.56
1	B	40	LYS	CB-CA-C	-6.13	108.90	117.23
1	C	40	LYS	CB-CA-C	-5.86	109.83	116.63
1	B	65	LYS	CA-C-N	5.70	125.55	119.28
1	B	65	LYS	C-N-CA	5.70	125.55	119.28
1	D	485	ASN	CA-C-N	5.60	125.44	119.28
1	D	485	ASN	C-N-CA	5.60	125.44	119.28
1	B	618	TYR	CA-C-N	5.38	124.57	118.97
1	B	618	TYR	C-N-CA	5.38	124.57	118.97
1	A	612	LEU	N-CA-C	-5.18	105.81	111.82
1	A	40	LYS	CB-CA-C	-5.14	110.66	116.63
1	A	635	MET	N-CA-C	-5.10	106.56	112.89
1	A	65	LYS	CA-C-N	5.08	124.39	118.85
1	A	65	LYS	C-N-CA	5.08	124.39	118.85
1	C	213	LEU	N-CA-C	-5.07	107.04	113.18

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	5169	0	5079	77	0
1	B	5213	0	5136	130	0
1	C	5213	0	5139	130	0
1	D	5135	0	5046	169	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
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	504	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 (504) 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:493:ASP:HB2	1:C:521:MET:HE1	1.50	0.91
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:330:MET:HE2	1:C:568:LEU:HD12	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
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:HG3	1:C:180:ARG:NH1	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:A:83:MET:HE1	1:A:153:GLU:HG3	1.72	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:B:634:ASN:HD22	1:B:634:ASN:C	2.02	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:D:101:PRO:HD2	4:D:784:HOH:O	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:O	1:B:458:LYS:HD3	1.96	0.65
1:C:493:ASP:CB	1:C:521:MET:HE1	2.25	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:B:464:LEU:HD13	1:B:474:MET:HE3	1.79	0.65
1:B:126:VAL:HG22	1:B:181:LYS:HZ2	1.61	0.64
1:B:295:ASP:CG	1:B:376:ARG:HH22	2.06	0.64
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:D:183:ARG:O	1:D:183:ARG:HG3	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
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:C:485:ASN:CG	4:C:706:HOH:O	2.43	0.62
1:A:624:GLU:C	1:A:626:VAL:H	2.08	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:HE	1:D:180:ARG:HA	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:O	1:D:227:ILE:HG22	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:335:LEU:HD13	1:B:474:MET:HE1	1.84	0.59
1:A:32:ALA:HB3	1:A:33:PRO:HD3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295[A]:ASP:CG	1:A:376:ARG:HH22	2.09	0.59
1:C:209:PHE:O	1:C:213:LEU:HB3	2.02	0.59
1:C:458:LYS:HD3	1:C:458:LYS:O	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:97:ILE:HG12	4:D:784:HOH:O	2.03	0.58
1:D:109:ASP:HA	1:D:112:ARG:NH1	2.18	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
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: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:C:31:LYS:O	1:C:35:THR:HG23	2.05	0.57
1:A:32:ALA:HB1	1:A:101:PRO:HG3	1.87	0.57
1:B:299:GLY:HA2	1:B:375:VAL:HG21	1.86	0.56
1:B:331:PHE:CZ	1:B:335:LEU:HD11	2.41	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:458:LYS:O	1:A:458:LYS:HD3	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:B:391:LYS:HZ2	3:B:1001:PEG:H22	1.72	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
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:B:542:GLU:CD	1:B:544:ASN:H	2.15	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:HH11	1:C:180:ARG:CG	2.08	0.52
1:A:45:LEU:HB2	1:A:103:VAL:HG12	1.91	0.52
1:A:458:LYS:HD3	1:A:458:LYS:C	2.33	0.52
1:C:48:PRO:HG3	1:C:143:LEU:HD22	1.92	0.52
1:C:399:ARG:HD3	1:C:403:ASN:ND2	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:378:LEU:HD22	1:B:432:LEU:HD11	1.91	0.52
2:B:902:G6P:O3P	2:B:902:G6P:O4	2.12	0.52
1:C:507:TYR:HB2	1:C:556:ARG:NH2	2.23	0.52
1:D:31:LYS:HZ2	1:D:35:THR:HG21	1.73	0.52
1:C:578:LYS:HZ3	1:C:586:GLN:NE2	2.08	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:D:537:MET:HE2	1:D:593:LEU:HD13	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:C:458:LYS:HD3	1:C:458:LYS:C	2.36	0.51
1:B:458:LYS:HD3	1:B:458:LYS:C	2.34	0.51
1:C:449:ASP:OD2	1:C:452:ASN:HB2	2.11	0.51
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:A:99:GLY:O	1:A:100:ALA:C	2.54	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:A:291[A]:GLU:HA	1:A:291[A]:GLU:OE1	2.11	0.50
1:D:180:ARG:HA	1:D:180:ARG:NE	2.26	0.50
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:70:SER:O	1:A:74:ARG:HG2	2.12	0.50
1:A:266:ILE:HG22	1:A:268:PRO:HD3	1.94	0.50
1:D:214:GLU:HA	1:D:257:HIS:CD2	2.44	0.50
1:B:634:ASN:ND2	1:B:634:ASN:C	2.70	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:D:542:GLU:CD	1:D:544:ASN:H	2.20	0.49
1:A:213:LEU:HD12	1:A:213:LEU:C	2.38	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:A:113:GLY:C	1:A:115:SER:H	2.20	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:D:31:LYS:HZ3	1:D:35:THR:HG21	1.77	0.48
1:A:35:THR:HG21	1:A:43:TYR:CE1	2.48	0.48
1:A:78:HIS:CB	1:A:157:LEU:HD23	2.43	0.48
1:A:164:VAL:HG11	1:A:610:ARG:HG2	1.95	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
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:54:TYR:CE1	1:B:60:ILE:HD11	2.50	0.47
1:B:163:ILE:HB	1:B:186:VAL:HG12	1.96	0.47
1:B:208:ASP:CG	1:B:211:ASN:HB2	2.38	0.47
1:B:273:VAL:HG13	1:B:520:VAL:CG1	2.44	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:72:GLU:OE2	1:B:161:HIS:HE1	1.98	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:517:GLU:O	1:B:521:MET:HG2	2.13	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:ND2	1:B:133:ASN:N	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:C:542:GLU:OE1	1:C:542:GLU:C	2.58	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:OE1	1:D:436:GLU:HA	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:8:HIS:HB2	1:A:162:ALA:O	2.15	0.46
1:A:119:LYS:HE3	1:A:130:SER:OG	2.16	0.46
1:B:111:VAL:CG1	1:B:118:TRP:HH2	2.28	0.46
1:C:5:LEU:HD22	1:C:614:LEU:HD22	1.97	0.46
1:C:32:ALA:HB3	1:C:33:PRO:HD3	1.98	0.46
1:C:512:GLY:C	1:C:515:PRO:HD2	2.40	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
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:C:617:GLY:C	1:C:619:PRO:HD3	2.40	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:25:TYR:CE1	1:C:95:TRP:HZ2	2.34	0.45
1:C:207:PHE:CE2	1:C:209:PHE:HA	2.51	0.45
1:D:279:PHE:CE1	1:D:591:GLU:OE1	2.69	0.45
1:B:493:ASP:HB2	1:B:521:MET:HE3	1.99	0.45
1:A:285:LEU:O	1:A:289:LYS:HG2	2.17	0.45
1:B:111:VAL:O	1:B:111:VAL:HG12	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:337:ARG:HH21	1:C:566:GLU:HG2	1.81	0.45
1:D:54:TYR:HE1	1:D:60:ILE:HD11	1.82	0.45
1:D:450:ASP:OD1	1:D:460:ARG:NH2	2.49	0.45
1:A:74:ARG:N	1:A:75:PRO:CD	2.79	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:459:ILE:HG12	1:B:459:ILE:H	1.49	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: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:D:330:MET:HE3	1:D:330:MET:HB3	1.74	0.44
1:A:94:ARG:HD2	1:A:100:ALA:HB1	1.99	0.44
1:B:512:GLY:O	1:B:515:PRO:HD2	2.17	0.44
1:C:372[A]:GLN:HE22	1:C:376:ARG:HE	1.66	0.44
1:C:503:VAL:HG11	1:C:572:MET:HE3	1.98	0.44
1:C:542:GLU:C	1:C:542:GLU:CD	2.84	0.44
1:D:428:ARG:HD2	1:D:428:ARG:HA	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: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: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:B:597:LEU:HA	1:B:602:MET:HE2	2.00	0.44
1:C:305:PHE:HZ	1:C:309:LEU:HG	1.81	0.44
1:C:503:VAL:HG21	1:C:572:MET:HE2	2.00	0.44
1:B:74:ARG:N	1:B:75:PRO:CD	2.80	0.44
1:B:111:VAL:HG11	1:B:118:TRP:HH2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:N	1:B:16:GLU:OE2	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:222:ALA:HB1	1:C:228:TYR:HA	1.99	0.44
1:D:19:ASN:N	1:D:19:ASN:ND2	2.66	0.44
1:D:128:ILE:HA	1:D:129:PRO:HD2	1.91	0.44
1:D:274:ILE:HG23	1:D:274:ILE:O	2.17	0.44
1:C:92:TYR:OH	1:C:102:LYS:HD3	2.17	0.44
1:D:63:TRP:CZ3	1:D:80:LEU:HB3	2.53	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: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:B:79:ALA:O	1:B:83:MET:HB2	2.18	0.43
1:B:349:LYS:O	1:B:471:ARG:HG3	2.18	0.43
1:C:249:GLN:NE2	1:C:249:GLN:HA	2.32	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:320:ARG:HH21	1:B:322:GLU:CD	2.26	0.43
1:B:549:TYR:HB3	1:B:593:LEU:HD11	1.99	0.43
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:D:450:ASP:CG	1:D:460:ARG:HH22	2.26	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:HE	1:D:94:ARG:HB2	1.64	0.43
1:B:542:GLU:CD	1:B:542:GLU:C	2.86	0.43
1:C:9:LEU:HD13	1:C:161:HIS:CG	2.53	0.43
1:D:471:ARG:NE	1:D:471:ARG:HA	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:VAL:O	1:C:126:VAL:HG12	2.18	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:D:150:PHE:CD2	1:D:150:PHE:C	2.97	0.43
1:B:464:LEU:CD1	1:B:474:MET:HE3	2.48	0.43
1:C:199:ARG:HG3	1:C:508:TYR:HE2	1.80	0.43
1:D:54:TYR:CE1	1:D:60:ILE:HD11	2.54	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:A:86:ARG:HA	1:A:86:ARG:HH11	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:114:TYR:O	1:B:115:SER:C	2.62	0.42
1:B:517:GLU:HG2	1:B:517:GLU:H	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:16:GLU:CD	1:C:24:ILE:HB	2.44	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:A:187:VAL:HG11	1:A:613:ALA:C	2.44	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
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:D:330:MET:HE2	1:D:505:PRO:HB3	2.00	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:HD23	1:D:309:LEU:HA	1.90	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:LYS:HA	1:A:326:LYS:HD3	1.82	0.42
1:C:626:VAL:HG11	1:C:630:LEU:HD11	2.01	0.42
1:D:5:LEU:HD13	1:D:622:PHE:HD2	1.84	0.42
1:D:292:LYS:HD2	1:D:490:LEU:CD2	2.43	0.42
1:B:330:MET:HE1	1:B:564:SER:HB3	2.02	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: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:D:635:MET:HB3	1:D:635:MET:HE3	1.84	0.42
1:B:323:TYR:OH	1:B:458:LYS:HG2	2.20	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:HA	1:B:326:LYS:HD3	1.86	0.41
1:C:14:ALA:HB3	1:C:28:LEU:HD11	2.03	0.41
1:C:541:ILE:O	1:C:542:GLU:C	2.64	0.41
1:B:372:GLN:HE21	1:B:372:GLN:HB3	1.69	0.41
1:D:349:LYS:O	1:D:471:ARG:HD2	2.20	0.41
1:D:456:LEU:O	1:D:460:ARG:HG3	2.21	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:D:29:LYS:HG3	1:D:97:ILE:HD12	2.01	0.41
1:D:63:TRP:H	1:D:63:TRP:CD1	2.38	0.41
1:D:268:PRO:HB2	1:D:602:MET:CE	2.51	0.41
1:D:283:GLN:HG3	2:D:901:G6P:H1	2.03	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
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:D:11:PHE:CD1	1:D:44:HIS:HB2	2.56	0.41
1:C:542:GLU:CD	1:C:543:THR:N	2.79	0.41
1:D:283:GLN:HG3	2:D:901:G6P:O1	2.20	0.41
1:C:74:ARG:HB2	1:C:75:PRO:HD3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:177:PRO:HA	1:C:240:SER:OG	2.21	0.41
1:D:181:LYS:C	1:D:183:ARG:N	2.77	0.41
1:D:321:TYR:OH	1:D:455:ILE:HG12	2.21	0.41
1:A:621:GLN:O	1:A:625:LEU:HD13	2.21	0.41
1:B:326:LYS:NZ	1:B:509:GLU:HG3	2.35	0.41
1:D:283:GLN:NE2	1:D:587:ARG:NH2	2.65	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: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:C:537:MET:HE1	1:C:593:LEU:HD13	2.03	0.41
1:D:17:VAL:O	1:D:17:VAL:CG1	2.69	0.41
1:D:125:LEU:HD23	1:D:125:LEU:HA	1.83	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:OD1	1:A:56:ASN:N	2.53	0.40
1:B:61:LEU:HG	1:B:93:GLY:HA2	2.04	0.40
1:B:196:LEU:O	1:B:200:TYR:HD2	2.04	0.40
1:B:269:ASN:HD22	1:B:511:TRP:CD1	2.39	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:HD2	1:D:615:ARG:O	2.21	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:O	1:D:274:ILE:CG2	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
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%)	36	50
1	B	646/725 (89%)	622 (96%)	22 (3%)	2 (0%)	36	50
1	C	646/725 (89%)	618 (96%)	25 (4%)	3 (0%)	24	37
1	D	633/725 (87%)	597 (94%)	32 (5%)	4 (1%)	21	32
All	All	2565/2900 (88%)	2443 (95%)	111 (4%)	11 (0%)	30	43

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 [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	555/623 (89%)	536 (97%)	19 (3%)	32	54
1	B	558/623 (90%)	533 (96%)	25 (4%)	24	42
1	C	558/623 (90%)	539 (97%)	19 (3%)	32	54
1	D	550/623 (88%)	528 (96%)	22 (4%)	28	47
All	All	2221/2492 (89%)	2136 (96%)	85 (4%)	29	49

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	17	VAL
1	A	19	ASN
1	A	56	ASN
1	A	83	MET
1	A	124	SER
1	A	180	ARG
1	A	181	LYS
1	A	216	VAL
1	A	310	ASP
1	A	337	ARG
1	A	372[A]	GLN
1	A	372[B]	GLN
1	A	376	ARG
1	A	458	LYS
1	A	521	MET
1	A	556	ARG
1	A	568	LEU
1	A	590	THR
1	B	21	VAL
1	B	133	ASN
1	B	216	VAL
1	B	220	HIS
1	B	289	LYS
1	B	310	ASP
1	B	320	ARG
1	B	337	ARG
1	B	343	LYS
1	B	372	GLN
1	B	376	ARG
1	B	408	GLU
1	B	420	SER
1	B	458	LYS

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Mol	Chain	Res	Type
1	B	459	ILE
1	B	471	ARG
1	B	484	ASN
1	B	517	GLU
1	B	539	ASP
1	B	553	ILE
1	B	585	ASN
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	213	LEU
1	C	216	VAL
1	C	240	SER
1	C	254	GLU
1	C	310	ASP
1	C	320	ARG
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	531	SER
1	C	543	THR
1	C	568	LEU
1	C	590	THR
1	C	645	LYS
1	D	6	GLN
1	D	19	ASN
1	D	40	LYS
1	D	111	VAL
1	D	122	LEU
1	D	133	ASN
1	D	213	LEU
1	D	249	GLN
1	D	310	ASP
1	D	324	LYS
1	D	330	MET
1	D	337	ARG

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Mol	Chain	Res	Type
1	D	343	LYS
1	D	372	GLN
1	D	387	THR
1	D	417	LEU
1	D	471	ARG
1	D	475	ILE
1	D	482	ASN
1	D	540	LEU
1	D	543	THR
1	D	590	THR

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	249	GLN
1	A	283	GLN
1	A	311	ASN
1	A	403	ASN
1	A	484	ASN
1	A	544	ASN
1	A	545	GLN
1	A	586	GLN
1	A	621	GLN
1	B	133	ASN
1	B	161	HIS
1	B	396	HIS
1	B	402	HIS
1	B	452	ASN
1	B	484	ASN
1	B	582	GLN
1	B	634	ASN
1	C	6	GLN
1	C	19	ASN
1	C	81	GLN
1	C	133	ASN
1	C	249	GLN
1	C	257	HIS
1	C	284	ASN
1	C	362	ASN
1	C	403	ASN
1	C	477	HIS
1	C	545	GLN

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Mol	Chain	Res	Type
1	C	586	GLN
1	D	19	ASN
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
1	D	586	GLN
1	D	621	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	B	901	-	16,16,16	0.89	1 (6%)	23,24,24	1.34	2 (8%)
3	PEG	B	1001	-	6,6,6	0.56	0	5,5,5	0.43	0
2	G6P	A	902	-	16,16,16	0.90	1 (6%)	23,24,24	1.34	3 (13%)
3	PEG	D	1002	-	6,6,6	0.64	0	5,5,5	0.45	0
2	G6P	A	901	-	16,16,16	0.90	1 (6%)	23,24,24	1.39	2 (8%)
3	PEG	B	1002	-	6,6,6	0.58	0	5,5,5	0.44	0
3	PEG	B	706	-	6,6,6	0.62	0	5,5,5	0.44	0
2	G6P	D	901	-	16,16,16	0.87	0	23,24,24	1.34	3 (13%)
3	PEG	A	1001	-	6,6,6	0.62	0	5,5,5	0.39	0
3	PEG	C	1001	-	6,6,6	0.61	0	5,5,5	0.39	0
2	G6P	B	902	-	16,16,16	0.91	0	23,24,24	1.95	6 (26%)
2	G6P	C	901	-	16,16,16	0.86	1 (6%)	23,24,24	1.41	2 (8%)

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	G6P	P-O2P	-2.06	1.47	1.54
2	C	901	G6P	P-O2P	-2.05	1.47	1.54
2	A	901	G6P	P-O2P	-2.02	1.47	1.54
2	A	902	G6P	P-O2P	-2.00	1.47	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	902	G6P	C4-C3-C2	-5.18	101.74	110.83
2	B	902	G6P	O5-C5-C4	-4.17	102.19	109.70
2	A	901	G6P	C4-C3-C2	-3.65	104.42	110.83
2	C	901	G6P	C4-C3-C2	-3.36	104.94	110.83
2	B	901	G6P	C4-C3-C2	-3.04	105.50	110.83
2	D	901	G6P	C4-C3-C2	-2.94	105.68	110.83
2	A	902	G6P	C4-C3-C2	-2.49	106.47	110.83
2	B	902	G6P	O5-C1-C2	2.47	114.64	110.30
2	B	902	G6P	O6-C6-C5	2.34	116.97	108.99
2	A	902	G6P	C1-O5-C5	-2.34	109.12	113.65
2	B	901	G6P	O2P-P-O6	-2.25	100.81	106.67
2	B	902	G6P	O2P-P-O6	-2.24	100.84	106.67
2	A	901	G6P	O2-C2-C1	-2.23	104.11	109.25
2	C	901	G6P	O1P-P-O3P	2.19	119.36	110.83
2	A	902	G6P	O2-C2-C1	-2.14	104.31	109.25
2	D	901	G6P	O5-C5-C6	-2.11	102.51	106.69
2	D	901	G6P	O2P-P-O6	-2.07	101.26	106.67
2	B	902	G6P	C3-C4-C5	-2.02	106.57	110.23

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

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

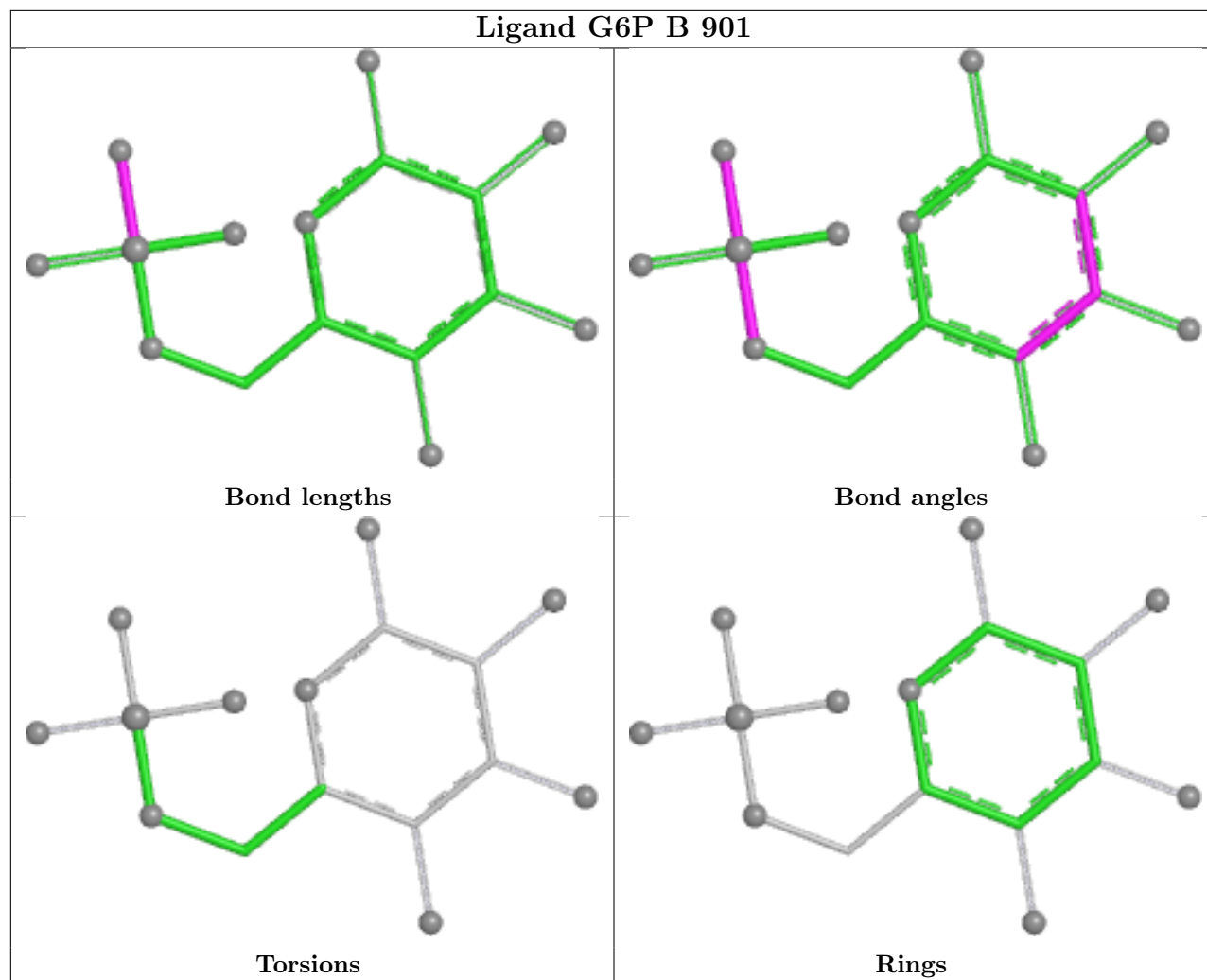
There are no ring outliers.

5 monomers are involved in 10 short contacts:

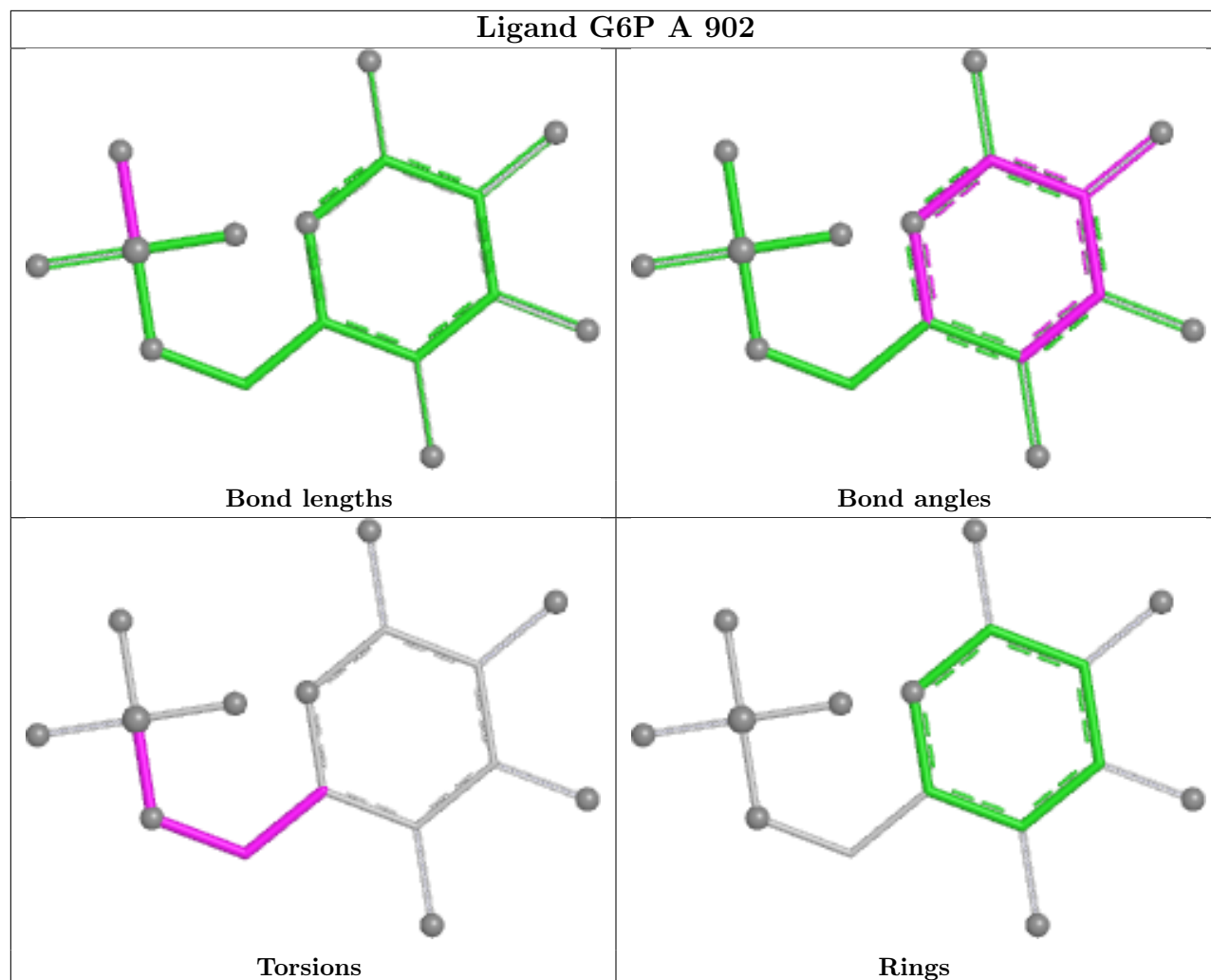
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1001	PEG	2	0
2	A	902	G6P	1	0
2	D	901	G6P	3	0
2	B	902	G6P	2	0
2	C	901	G6P	2	0

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.

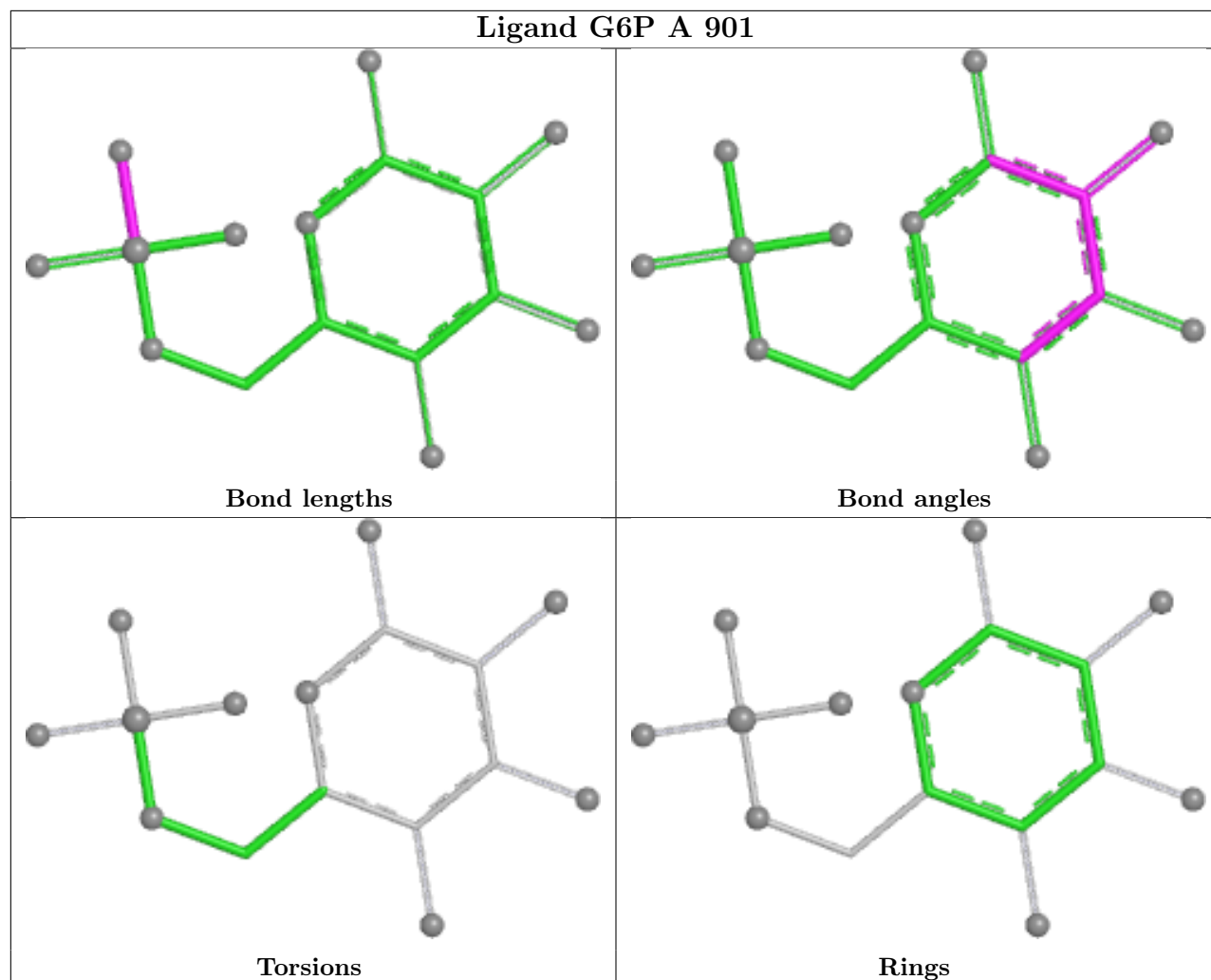
Ligand G6P B 901



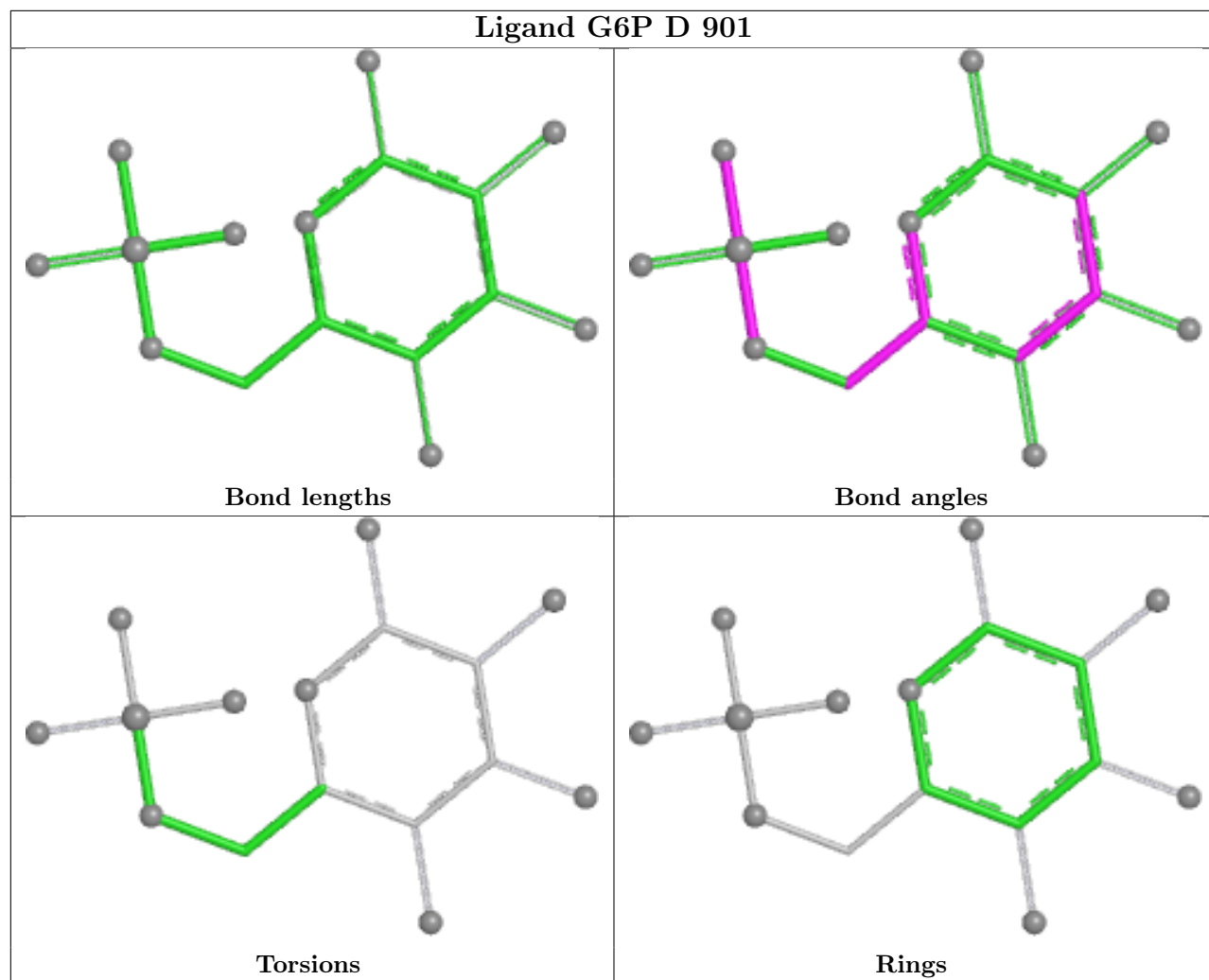
Ligand G6P A 902



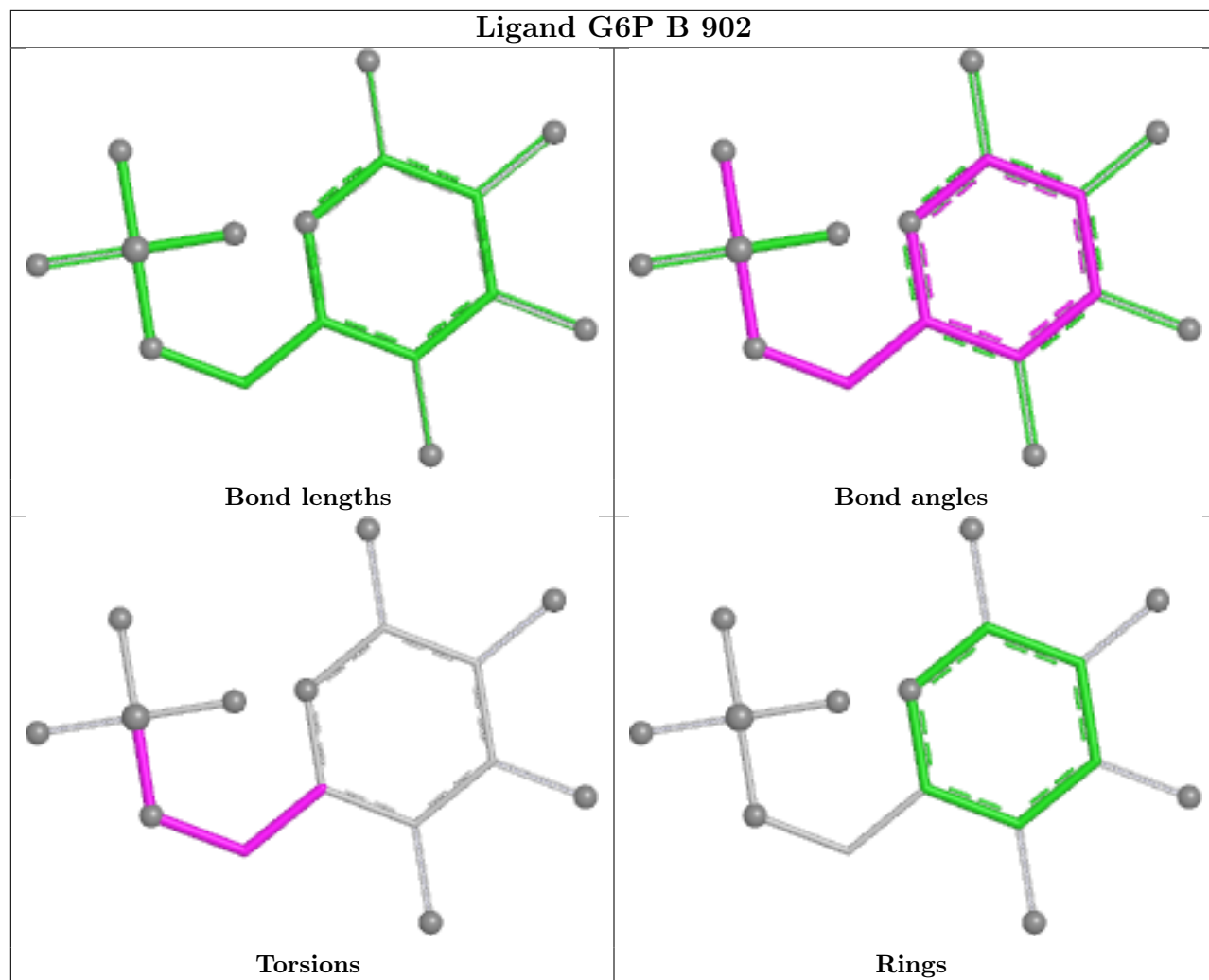
Ligand G6P A 901

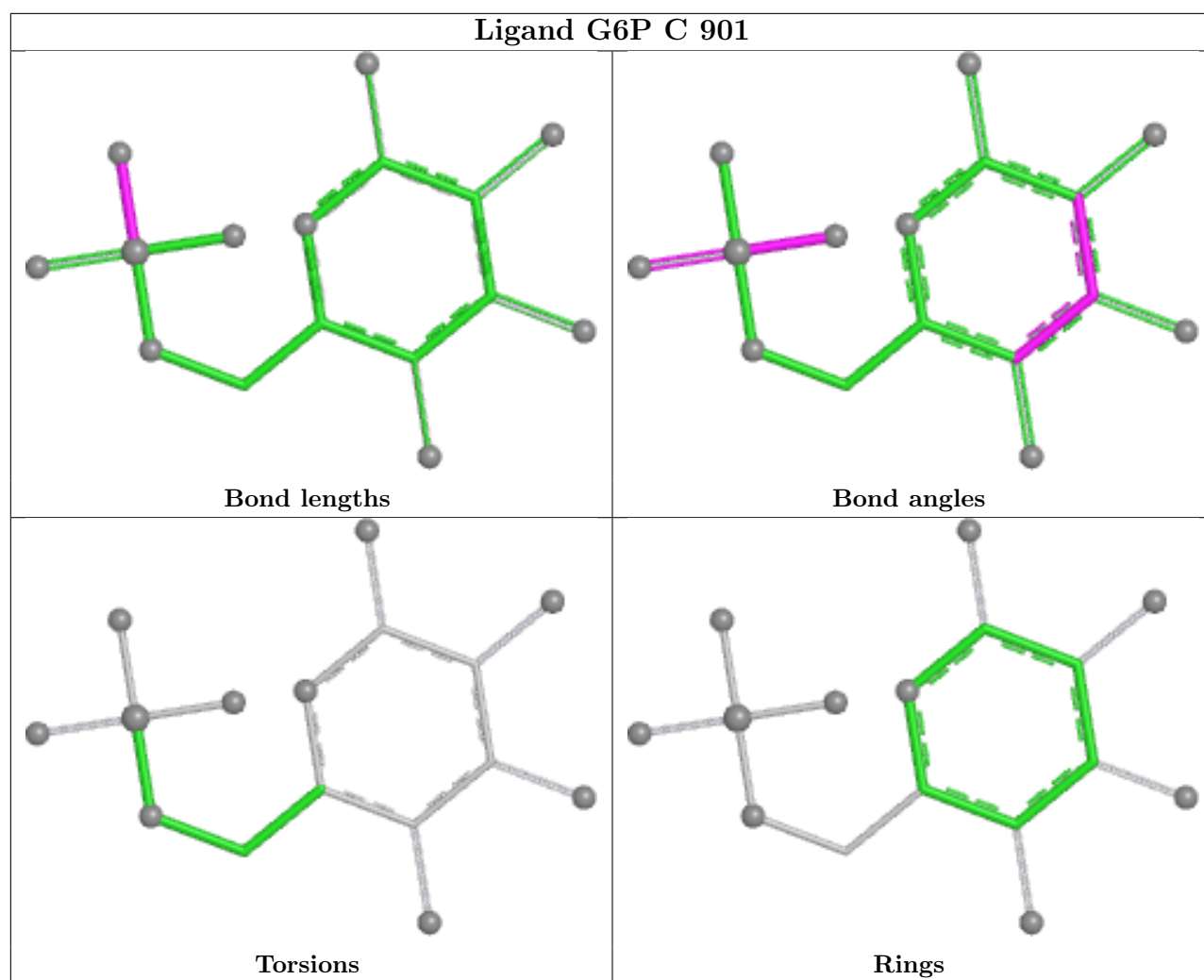


Ligand G6P D 901



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%)	0.60	100 (15%) 5 4	20, 49, 89, 101	4 (0%)
1	B	645/725 (88%)	0.43	34 (5%) 32 28	24, 50, 83, 96	3 (0%)
1	C	646/725 (89%)	0.53	36 (5%) 30 26	24, 53, 87, 101	2 (0%)
1	D	636/725 (87%)	0.78	84 (13%) 7 5	22, 61, 103, 114	1 (0%)
All	All	2565/2900 (88%)	0.59	254 (9%) 13 9	20, 52, 95, 114	10 (0%)

All (254) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	85	SER	5.7
1	D	630	LEU	5.5
1	D	205	GLY	5.3
1	D	125	LEU	5.3
1	D	622	PHE	4.7
1	C	135	PHE	4.5
1	A	623	ARG	4.5
1	A	90	PHE	4.4
1	A	155	ALA	4.2
1	A	91	VAL	4.2
1	D	212	CYS	4.0
1	A	63	TRP	4.0
1	A	2	SER	3.9
1	A	68	ALA	3.9
1	D	66	PRO	3.8
1	A	60	ILE	3.8
1	B	128	ILE	3.8
1	C	634	ASN	3.8
1	C	125	LEU	3.7
1	C	133	ASN	3.7
1	D	61	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	43	TYR	3.6
1	B	206	SER	3.6
1	C	645	LYS	3.6
1	A	624	GLU	3.6
1	A	625	LEU	3.5
1	A	637	ALA	3.5
1	A	160	GLN	3.5
1	B	542	GLU	3.4
1	A	148	ALA	3.4
1	D	123	TRP	3.4
1	D	122	LEU	3.3
1	C	82	THR	3.3
1	D	626	VAL	3.3
1	D	88	VAL	3.3
1	A	52	ALA	3.3
1	C	156	HIS	3.2
1	B	125	LEU	3.2
1	D	87	GLY	3.2
1	A	274	ILE	3.2
1	D	163	ILE	3.2
1	D	21	VAL	3.2
1	D	63	TRP	3.2
1	D	639	ALA	3.1
1	A	80	LEU	3.1
1	A	107	ASP	3.1
1	A	34	ILE	3.1
1	D	114	TYR	3.1
1	A	10	LEU	3.1
1	A	85	SER	3.1
1	A	76	VAL	3.1
1	A	630	LEU	3.0
1	A	151	LEU	3.0
1	D	625	LEU	3.0
1	D	179	CYS	3.0
1	B	2	SER	3.0
1	A	163	ILE	3.0
1	D	81	GLN	3.0
1	A	78	HIS	3.0
1	D	119	LYS	3.0
1	D	164	VAL	3.0
1	B	543	THR	3.0
1	C	83	MET	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	67	GLU	3.0
1	D	149	TRP	3.0
1	B	129	PRO	3.0
1	C	59	ASP	3.0
1	D	135	PHE	3.0
1	A	55	GLN	3.0
1	A	629	GLU	2.9
1	C	68	ALA	2.9
1	D	607	VAL	2.9
1	D	126	VAL	2.9
1	A	62	ASP	2.9
1	A	92	TYR	2.9
1	A	436	GLU	2.9
1	C	204	SER	2.9
1	A	103	VAL	2.9
1	A	154	VAL	2.9
1	D	121	ASP	2.9
1	A	544	ASN	2.8
1	B	181	LYS	2.8
1	A	111	VAL	2.8
1	A	114	TYR	2.8
1	A	93	GLY	2.8
1	A	88	VAL	2.8
1	A	61	LEU	2.8
1	C	132	GLU	2.8
1	D	62	ASP	2.8
1	D	60	ILE	2.8
1	A	618	TYR	2.7
1	C	66	PRO	2.7
1	D	129	PRO	2.7
1	B	257[A]	HIS	2.7
1	D	636	ASP	2.7
1	A	66	PRO	2.7
1	A	5	LEU	2.7
1	A	157	LEU	2.7
1	B	93	GLY	2.7
1	C	205	GLY	2.7
1	D	128	ILE	2.7
1	A	110	SER	2.7
1	C	51	LYS	2.7
1	D	204	SER	2.7
1	A	65	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	618	TYR	2.7
1	C	52	ALA	2.7
1	A	6	GLN	2.7
1	B	253	PHE	2.6
1	B	279	PHE	2.6
1	D	274	ILE	2.6
1	A	67	GLU	2.6
1	A	621	GLN	2.6
1	A	3	ARG	2.6
1	A	161	HIS	2.6
1	B	126	VAL	2.6
1	B	646	VAL	2.6
1	A	82	THR	2.6
1	C	18	ALA	2.6
1	A	89	HIS	2.6
1	D	86	ARG	2.6
1	A	108	LEU	2.6
1	A	628	GLU	2.6
1	D	637	ALA	2.6
1	B	219	ASP	2.5
1	A	159	SER	2.5
1	B	641	GLY	2.5
1	D	227	ILE	2.5
1	A	36	VAL	2.5
1	A	147	VAL	2.5
1	D	110	SER	2.5
1	A	617	GLY	2.5
1	A	105	LEU	2.5
1	A	611	GLN	2.5
1	A	42	HIS	2.5
1	C	279	PHE	2.5
1	C	86	ARG	2.5
1	A	73	MET	2.5
1	D	124	SER	2.5
1	A	59	ASP	2.4
1	A	636	ASP	2.4
1	D	178	LEU	2.4
1	A	135	PHE	2.4
1	B	544	ASN	2.4
1	C	21	VAL	2.4
1	D	174	VAL	2.4
1	D	72	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	87	GLY	2.4
1	D	93	GLY	2.4
1	D	113	GLY	2.4
1	B	487	ILE	2.4
1	D	261	ARG	2.4
1	D	116	ASN	2.4
1	D	246	THR	2.4
1	A	70	SER	2.4
1	B	127	GLY	2.4
1	D	619	PRO	2.4
1	A	156	HIS	2.4
1	A	69	PHE	2.4
1	A	591	GLU	2.4
1	A	543	THR	2.3
1	C	113	GLY	2.3
1	A	184	ILE	2.3
1	A	548	ASP	2.3
1	A	54	TYR	2.3
1	A	106	PHE	2.3
1	B	643	LYS	2.3
1	D	118	TRP	2.3
1	D	197	LEU	2.3
1	A	626	VAL	2.3
1	D	43	TYR	2.3
1	D	145	TYR	2.3
1	C	202	CYS	2.3
1	D	624	GLU	2.3
1	C	320	ARG	2.3
1	B	89	HIS	2.3
1	A	178	LEU	2.3
1	B	644	LEU	2.3
1	D	171	LEU	2.3
1	A	102	LYS	2.3
1	B	21	VAL	2.3
1	A	183	ARG	2.3
1	A	639	ALA	2.2
1	D	115	SER	2.2
1	D	130	SER	2.2
1	D	141	ILE	2.2
1	C	310	ASP	2.2
1	D	185	ASP	2.2
1	A	164	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	199	ARG	2.2
1	B	484	ASN	2.2
1	B	246	THR	2.2
1	A	9	LEU	2.2
1	A	96	LEU	2.2
1	D	638	LEU	2.2
1	A	419	SER	2.2
1	C	69	PHE	2.2
1	A	186	VAL	2.2
1	B	448	VAL	2.2
1	D	68	ALA	2.2
1	B	645	LYS	2.2
1	D	92	TYR	2.2
1	A	87	GLY	2.2
1	B	122	LEU	2.2
1	A	631	ASN	2.2
1	D	160	GLN	2.2
1	A	127	GLY	2.2
1	C	2	SER	2.2
1	A	539	ASP	2.2
1	C	517	GLU	2.2
1	D	131	PRO	2.2
1	D	133	ASN	2.1
1	D	142	LEU	2.1
1	C	180	ARG	2.1
1	A	38	GLN	2.1
1	D	621	GLN	2.1
1	A	613	ALA	2.1
1	C	203	ALA	2.1
1	D	609	ALA	2.1
1	D	192	THR	2.1
1	D	612	LEU	2.1
1	D	74	ARG	2.1
1	D	539	ASP	2.1
1	C	647	ALA	2.1
1	B	117	GLU	2.1
1	D	117	GLU	2.1
1	C	124	SER	2.1
1	C	206	SER	2.1
1	D	232	CYS	2.1
1	D	620	ASP	2.1
1	C	116	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	35	THR	2.1
1	D	219	ASP	2.1
1	A	101	PRO	2.1
1	D	161	HIS	2.1
1	A	77	GLN	2.0
1	A	614	LEU	2.0
1	B	61	LEU	2.0
1	A	181	LYS	2.0
1	B	64	LYS	2.0
1	D	228	TYR	2.0
1	D	73	MET	2.0
1	D	78	HIS	2.0
1	D	257	HIS	2.0
1	A	58	VAL	2.0
1	A	79	ALA	2.0
1	A	612	LEU	2.0
1	B	148	ALA	2.0
1	C	274	ILE	2.0
1	D	82	THR	2.0
1	A	4	ASP	2.0
1	A	41	ASP	2.0
1	B	424	MET	2.0
1	C	70	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

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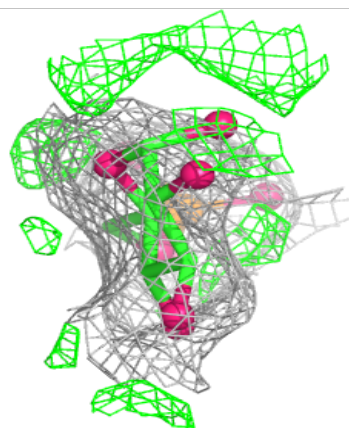
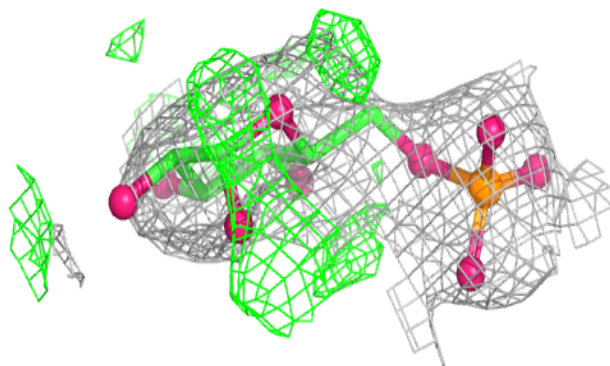
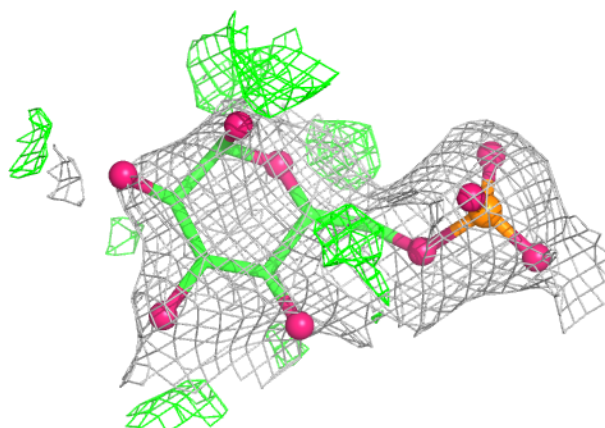
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PEG	B	1002	7/7	0.76	0.20	60,61,65,67	0
3	PEG	D	1002	7/7	0.81	0.15	47,50,55,58	0
3	PEG	A	1001	7/7	0.86	0.14	49,50,54,56	0
2	G6P	A	902	16/16	0.87	0.16	46,52,58,60	16
3	PEG	C	1001	7/7	0.88	0.11	43,45,49,50	0
2	G6P	B	902	16/16	0.89	0.13	51,54,56,60	0
3	PEG	B	1001	7/7	0.89	0.14	42,44,46,50	0
3	PEG	B	706	7/7	0.91	0.10	43,46,48,49	0
2	G6P	A	901	16/16	0.97	0.07	30,35,38,40	0
2	G6P	C	901	16/16	0.97	0.06	39,42,43,44	0
2	G6P	D	901	16/16	0.97	0.06	32,34,36,38	0
2	G6P	B	901	16/16	0.98	0.06	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.

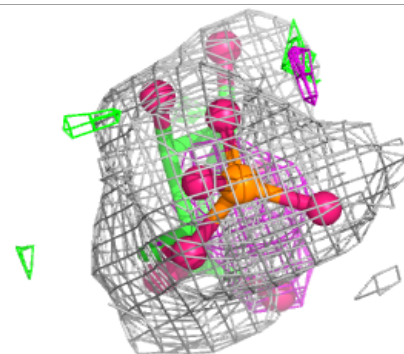
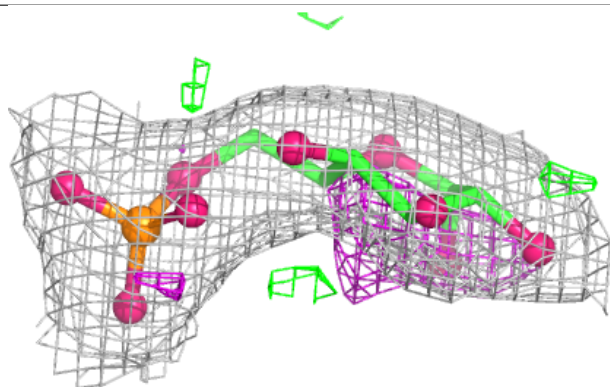
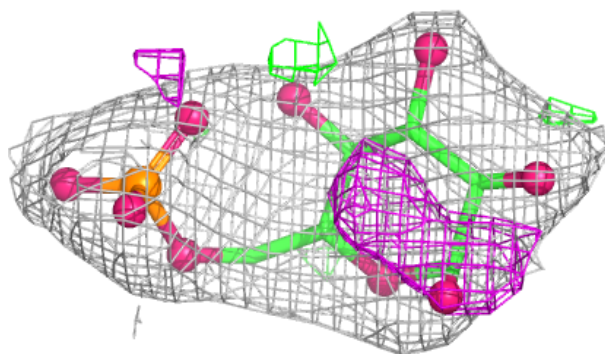
Electron density around G6P A 902:

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

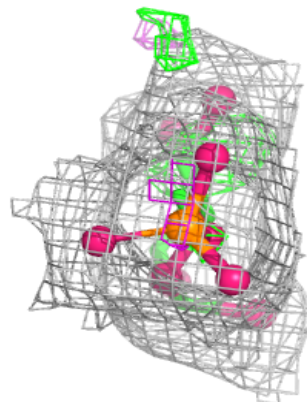
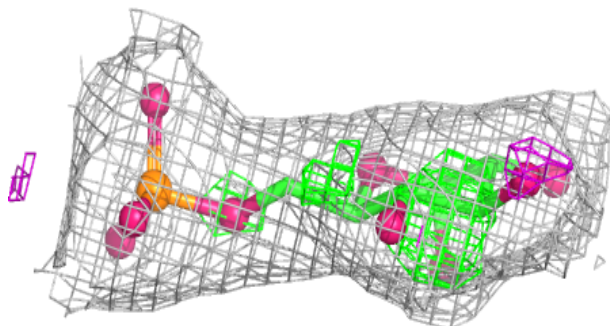
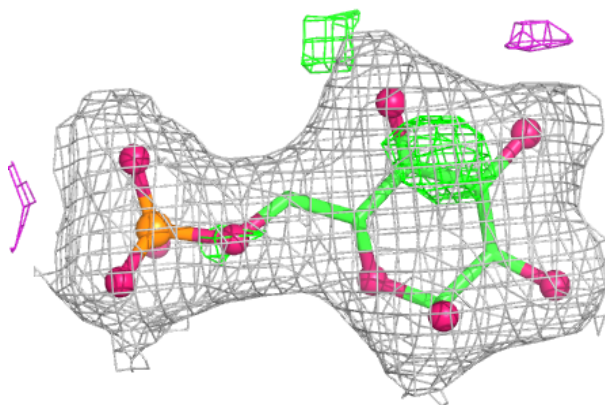


Electron density around G6P B 902:

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

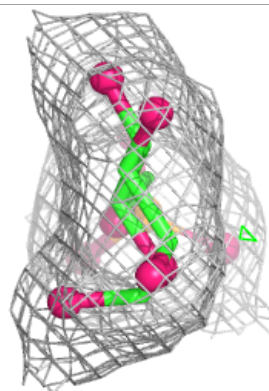
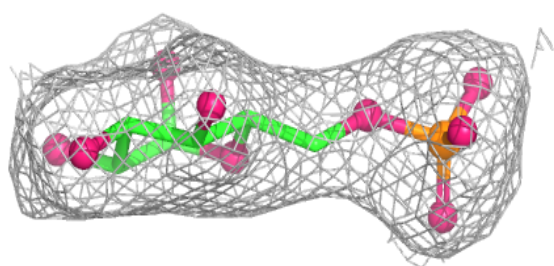
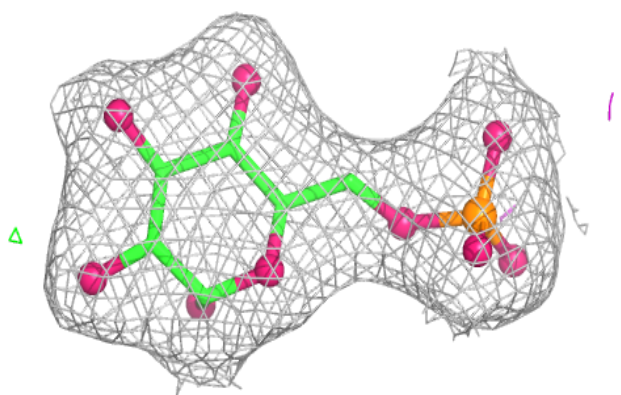
**Electron density around G6P A 901:**

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

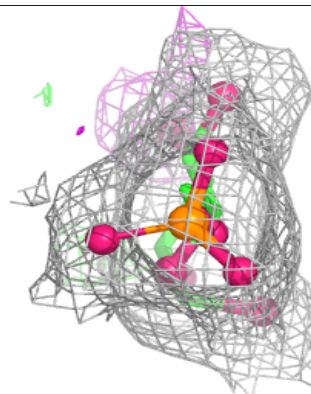
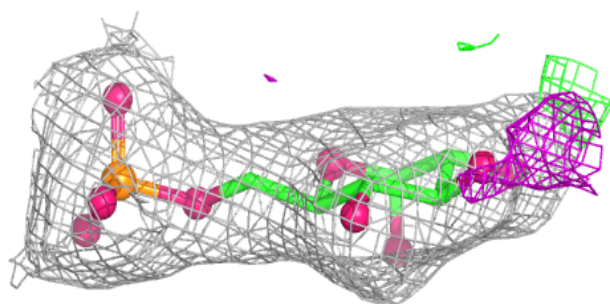
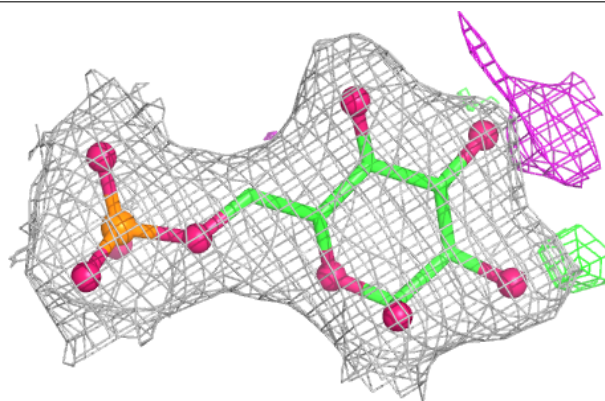


Electron density around G6P C 901:

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

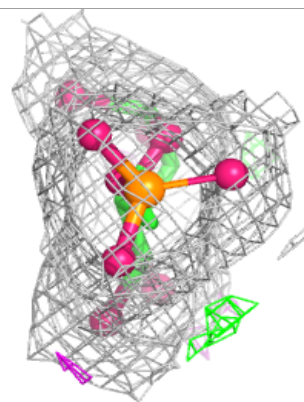
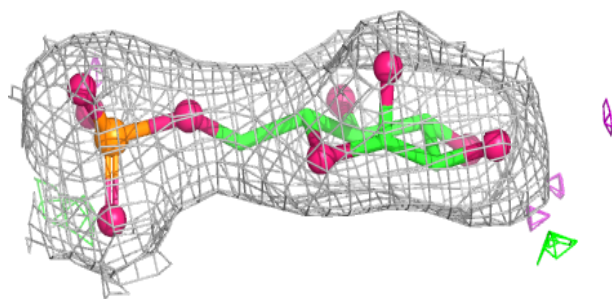
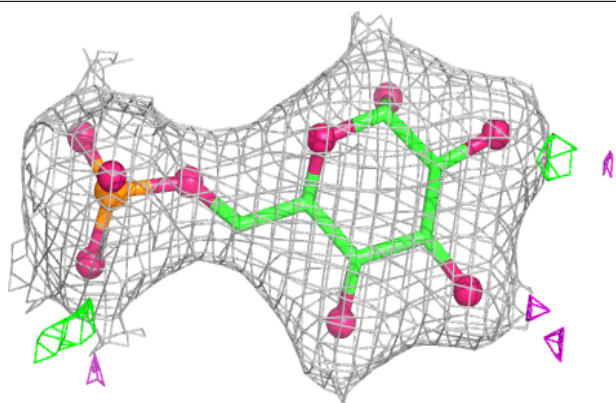
**Electron density around G6P D 901:**

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



Electron density around G6P 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.