



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:35 PM UTC

PDB ID : 3PR3 / pdb_00003pr3
Title : Crystal structure of Plasmodium falciparum glucose-6-phosphate isomerase (PF14_0341) in complex with fructose-6-phosphate
Authors : Gileadi, T.; Wernimont, A.K.; Hutchinson, A.; Weadge, J.; Cossar, D.; Lew, J.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Weigelt, J.; Hui, R.; Hills, T.; Pizarro, J.C.; Structural Genomics Consortium (SGC)
Deposited on : 2010-11-29
Resolution : 2.45 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](#) ①) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (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

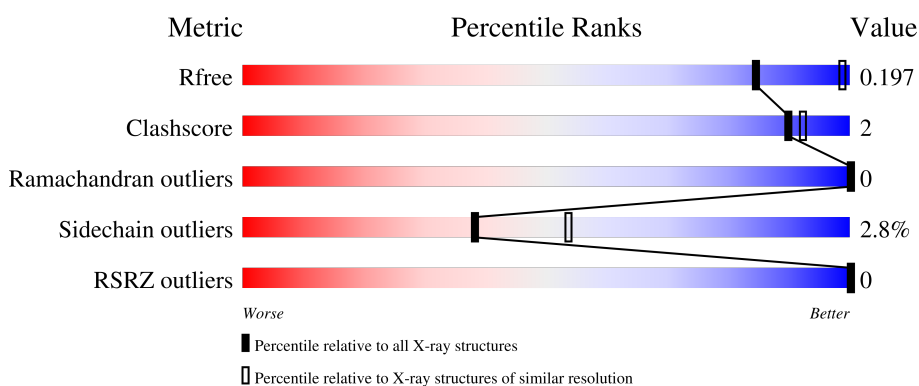
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	597	 87% 9% ..
1	B	597	 87% 9% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	F6P	A	580	X	-	-	-
2	F6P	B	580	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 19596 atoms, of which 9293 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	578	9369	3029	4640	793	889	18	0	13	0
1	B	577	9372	3030	4631	792	899	20	0	13	0

There are 36 discrepancies between the modelled and reference sequences:

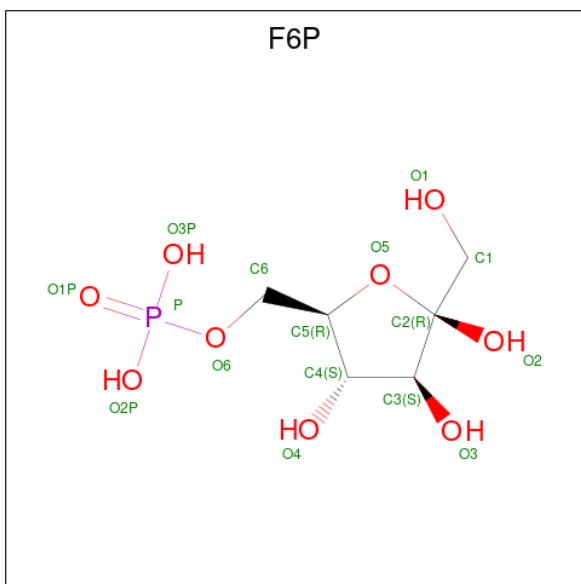
Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	expression tag	UNP Q8ILA4
A	-16	HIS	-	expression tag	UNP Q8ILA4
A	-15	HIS	-	expression tag	UNP Q8ILA4
A	-14	HIS	-	expression tag	UNP Q8ILA4
A	-13	HIS	-	expression tag	UNP Q8ILA4
A	-12	HIS	-	expression tag	UNP Q8ILA4
A	-11	HIS	-	expression tag	UNP Q8ILA4
A	-10	SER	-	expression tag	UNP Q8ILA4
A	-9	SER	-	expression tag	UNP Q8ILA4
A	-8	GLY	-	expression tag	UNP Q8ILA4
A	-7	ARG	-	expression tag	UNP Q8ILA4
A	-6	GLU	-	expression tag	UNP Q8ILA4
A	-5	ASN	-	expression tag	UNP Q8ILA4
A	-4	LEU	-	expression tag	UNP Q8ILA4
A	-3	TYR	-	expression tag	UNP Q8ILA4
A	-2	PHE	-	expression tag	UNP Q8ILA4
A	-1	GLN	-	expression tag	UNP Q8ILA4
A	0	GLY	-	expression tag	UNP Q8ILA4
B	-17	MET	-	expression tag	UNP Q8ILA4
B	-16	HIS	-	expression tag	UNP Q8ILA4
B	-15	HIS	-	expression tag	UNP Q8ILA4
B	-14	HIS	-	expression tag	UNP Q8ILA4
B	-13	HIS	-	expression tag	UNP Q8ILA4
B	-12	HIS	-	expression tag	UNP Q8ILA4
B	-11	HIS	-	expression tag	UNP Q8ILA4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	SER	-	expression tag	UNP Q8ILA4
B	-9	SER	-	expression tag	UNP Q8ILA4
B	-8	GLY	-	expression tag	UNP Q8ILA4
B	-7	ARG	-	expression tag	UNP Q8ILA4
B	-6	GLU	-	expression tag	UNP Q8ILA4
B	-5	ASN	-	expression tag	UNP Q8ILA4
B	-4	LEU	-	expression tag	UNP Q8ILA4
B	-3	TYR	-	expression tag	UNP Q8ILA4
B	-2	PHE	-	expression tag	UNP Q8ILA4
B	-1	GLN	-	expression tag	UNP Q8ILA4
B	0	GLY	-	expression tag	UNP Q8ILA4

- Molecule 2 is 6-O-phosphono-beta-D-fructofuranose (CCD ID: F6P) (formula: C₆H₁₃O₉P).



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

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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


- Molecule 4 is water.

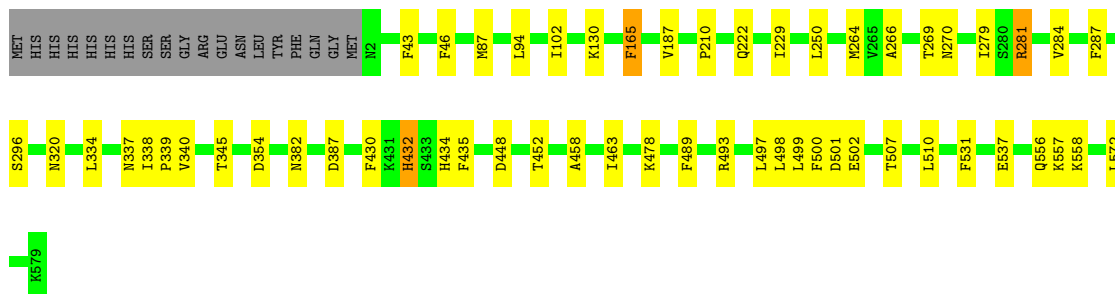
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	380	Total	O	0	0
			380	380		
4	B	411	Total	O	0	0
			411	411		

3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

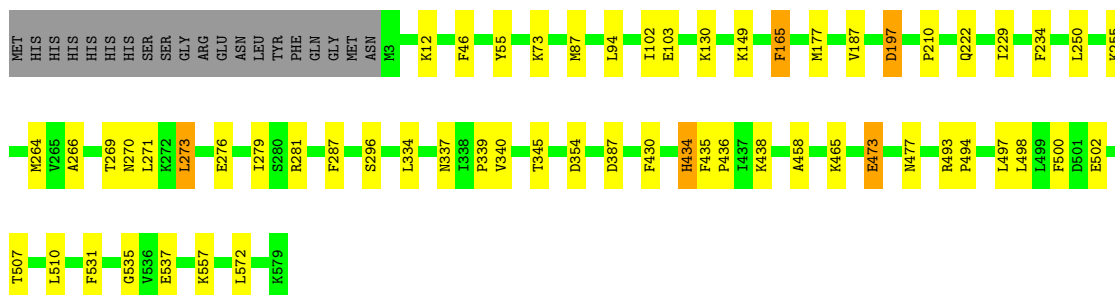
- Molecule 1: Glucose-6-phosphate isomerase

Chain A:  87% 9% ..



- Molecule 1: Glucose-6-phosphate isomerase

Chain B:  87% 9% ..



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.42Å 118.33Å 73.04Å 90.00° 117.30° 90.00°	Depositor
Resolution (Å)	32.18 – 2.45 32.18 – 2.45	Depositor EDS
% Data completeness (in resolution range)	(Not available) (32.18-2.45) 99.8 (32.18-2.45)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.20 (at 2.45Å)	Xtrriage
Refinement program	BUSTER-TNT BUSTER 2.8.0, BUSTER 2.8.0	Depositor
R, R_{free}	0.140 , 0.197 0.147 , 0.197	Depositor DCC
R_{free} test set	2014 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	24.7	Xtrriage
Anisotropy	0.169	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.003 for -h-l,k,h 0.003 for l,k,-h-l 0.022 for h,-k,-h-l 0.023 for -h-l,-k,l 0.017 for l,-k,h	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19596	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% 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: F6P, PO4

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	0.81	3/4865 (0.1%)	1.34	20/6574 (0.3%)
1	B	0.81	0/4884	1.32	16/6600 (0.2%)
All	All	0.81	3/9749 (0.0%)	1.33	36/13174 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	ILE	CA-CB	6.51	1.57	1.54
1	A	432	HIS	ND1-CE1	5.57	1.38	1.32
1	A	556	GLN	CD-OE1	5.16	1.33	1.23

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	432	HIS	CB-CG-CD2	-7.50	121.46	131.20
1	A	320	ASN	CA-CB-CG	7.08	119.68	112.60
1	A	269	THR	N-CA-C	-6.58	105.17	113.72
1	A	345	THR	CA-C-N	6.40	128.76	120.44
1	A	345	THR	C-N-CA	6.40	128.76	120.44
1	B	46[A]	PHE	N-CA-C	6.37	118.98	109.11
1	B	46[B]	PHE	N-CA-C	6.37	118.98	109.11
1	B	345	THR	CA-C-N	6.27	128.59	120.44
1	B	345	THR	C-N-CA	6.27	128.59	120.44
1	A	187	VAL	N-CA-C	6.12	116.72	110.05
1	A	537	GLU	N-CA-C	5.99	117.89	111.36
1	B	387	ASP	CA-CB-CG	5.98	118.58	112.60
1	B	296	SER	N-CA-C	5.97	120.56	113.28
1	B	537	GLU	N-CA-C	5.94	117.83	111.36
1	B	269	THR	N-CA-C	-5.92	106.10	113.38
1	A	387	ASP	CA-CB-CG	5.80	118.40	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	LEU	N-CA-C	5.76	118.25	110.88
1	A	43	PHE	N-CA-C	-5.61	101.22	109.18
1	A	296	SER	N-CA-C	5.45	120.08	113.38
1	B	197	ASP	N-CA-C	5.42	120.27	111.37
1	B	287	PHE	CA-CB-CG	5.41	119.21	113.80
1	B	94	LEU	N-CA-C	5.39	117.78	110.88
1	A	270	ASN	N-CA-C	5.35	117.96	107.57
1	A	463	ILE	N-CA-C	5.31	116.39	111.45
1	A	493	ARG	CB-CA-C	5.24	115.36	110.17
1	A	531	PHE	N-CA-C	5.22	119.80	113.38
1	A	434	HIS	N-CA-C	-5.19	106.80	113.23
1	B	494	PRO	N-CA-C	5.19	119.63	111.38
1	A	87	MET	N-CA-C	5.14	117.62	111.71
1	B	531	PHE	N-CA-C	5.11	119.67	113.38
1	A	287	PHE	CA-CB-CG	5.11	118.91	113.80
1	A	382	ASN	N-CA-C	5.09	119.58	113.17
1	B	255	LYS	N-CA-C	5.08	119.60	113.41
1	B	434	HIS	N-CA-C	-5.07	106.95	113.23
1	B	493	ARG	CB-CA-C	5.05	115.17	110.17
1	A	46	PHE	N-CA-C	5.02	115.93	108.60

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	4729	4640	4637	16	0
1	B	4741	4631	4618	21	0
2	A	16	11	11	1	0
2	B	16	11	11	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	380	0	0	2	0
4	B	411	0	0	1	0
All	All	10303	9293	9277	34	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 2.

All (34) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281[B]:ARG:NH1	4:A:609:HOH:O	2.36	0.56
1:B:270:ASN:ND2	1:B:273:LEU:HD22	2.24	0.52
1:B:271:LEU:HB3	1:B:281[B]:ARG:HH22	1.77	0.50
1:B:229:ILE:HD11	1:B:264:MET:HE2	1.95	0.49
1:B:334:LEU:HD22	1:B:340:VAL:HG11	1.94	0.49
1:A:435:PHE:O	1:B:210:PRO:HB2	2.13	0.49
1:B:273:LEU:HA	1:B:276:GLU:HG2	1.94	0.48
1:A:222:GLN:HG2	1:A:250:LEU:HA	1.95	0.47
1:B:465:LYS:HZ3	1:B:473[A]:GLU:HG2	1.79	0.47
1:B:222:GLN:HG2	1:B:250:LEU:HA	1.96	0.47
1:B:430:PHE:CD2	1:B:500:PHE:HB2	2.49	0.47
1:B:458:ALA:HB2	1:B:497:LEU:HD11	1.96	0.47
1:A:266:ALA:HB2	1:A:279:ILE:HD13	1.98	0.46
1:A:337:ASN:OD1	1:A:339:PRO:HD2	2.15	0.46
1:A:229:ILE:HD11	1:A:264:MET:HE2	1.97	0.45
1:A:489:PHE:CZ	1:B:535:GLY:HA2	2.51	0.45
1:B:266:ALA:HB2	1:B:279:ILE:HD13	1.99	0.45
1:A:430:PHE:CD2	1:A:500:PHE:HB2	2.51	0.45
1:A:210:PRO:HB2	1:B:435:PHE:O	2.17	0.44
1:B:12:LYS:HE3	4:B:717:HOH:O	2.17	0.43
1:A:165:PHE:CD1	1:A:165:PHE:C	2.96	0.43
1:A:502:GLU:O	1:A:507:THR:HG21	2.19	0.43
1:A:432:HIS:HD2	1:A:501:ASP:O	2.02	0.43
1:B:165:PHE:CD1	1:B:165:PHE:C	2.97	0.43
1:A:458:ALA:HB2	1:A:497:LEU:HD11	2.00	0.42
2:A:580:F6P:H4	4:A:743:HOH:O	2.20	0.42
1:B:177[A]:MET:HE1	1:B:434:HIS:HA	2.01	0.42
1:B:337:ASN:OD1	1:B:339:PRO:HD2	2.19	0.42
1:B:502:GLU:O	1:B:507:THR:HG21	2.19	0.42
1:B:87[B]:MET:H	1:B:87[B]:MET:HG3	1.66	0.42
1:A:334:LEU:HD22	1:A:340:VAL:HG11	2.01	0.41
1:A:499:LEU:HD23	1:B:572:LEU:HD22	2.02	0.41
1:B:187:VAL:HG12	1:B:436:PRO:HB2	2.01	0.41
1:A:448:ASP:O	1:A:452:THR:HG23	2.21	0.41

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	589/597 (99%)	573 (97%)	16 (3%)	0	100	100
1	B	588/597 (98%)	572 (97%)	16 (3%)	0	100	100
All	All	1177/1194 (99%)	1145 (97%)	32 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	523/550 (95%)	510 (98%)	13 (2%)	42	56
1	B	528/550 (96%)	510 (97%)	18 (3%)	32	47
All	All	1051/1100 (96%)	1020 (97%)	31 (3%)	38	52

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	ILE
1	A	130	LYS
1	A	165	PHE
1	A	281[A]	ARG
1	A	281[B]	ARG
1	A	284	VAL
1	A	354	ASP
1	A	478	LYS

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Mol	Chain	Res	Type
1	A	498	LEU
1	A	510	LEU
1	A	557	LYS
1	A	558	LYS
1	A	572	LEU
1	B	55	TYR
1	B	73	LYS
1	B	102	ILE
1	B	103	GLU
1	B	130	LYS
1	B	149	LYS
1	B	165	PHE
1	B	197	ASP
1	B	234	PHE
1	B	273	LEU
1	B	354	ASP
1	B	438	LYS
1	B	473[A]	GLU
1	B	473[B]	GLU
1	B	477	ASN
1	B	498	LEU
1	B	510	LEU
1	B	557	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	ASN
1	A	116	ASN
1	A	176	ASN
1	A	219	ASN
1	A	394	HIS
1	A	421	GLN
1	A	561	ASN
1	B	116	ASN
1	B	176	ASN
1	B	189	ASN
1	B	214	ASN
1	B	219	ASN
1	B	390	ASN
1	B	421	GLN
1	B	551	ASN

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Mol	Chain	Res	Type
1	B	555	ASN
1	B	561	ASN

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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	F6P	A	580	-	15,16,16	1.93	3 (20%)	16,25,25	2.90	6 (37%)
3	PO4	B	581	-	4,4,4	4.02	3 (75%)	6,6,6	0.94	0
2	F6P	B	580	-	15,16,16	1.11	1 (6%)	16,25,25	1.52	2 (12%)
3	PO4	A	581	-	4,4,4	3.19	3 (75%)	6,6,6	0.85	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	F6P	A	580	-	1/1/5/5	5/9/28/28	0/1/1/1
2	F6P	B	580	-	1/1/5/5	1/9/28/28	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	581	PO4	P-O1	6.44	1.65	1.50
2	A	580	F6P	O2-C2	5.46	1.50	1.40
3	A	581	PO4	P-O1	4.74	1.61	1.50
3	B	581	PO4	P-O2	3.70	1.65	1.54
2	B	580	F6P	O2-C2	3.42	1.46	1.40
2	A	580	F6P	O6-C6	-3.12	1.32	1.44
3	A	581	PO4	P-O2	3.08	1.63	1.54
3	B	581	PO4	P-O4	2.98	1.63	1.54
3	A	581	PO4	P-O3	2.64	1.62	1.54
2	A	580	F6P	O5-C2	2.33	1.47	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	580	F6P	O3P-P-O6	6.77	124.33	106.67
2	A	580	F6P	O2P-P-O6	-5.45	92.47	106.67
2	A	580	F6P	C6-C5-C4	-4.82	97.85	115.21
2	B	580	F6P	O1-C1-C2	-4.28	102.19	111.67
2	A	580	F6P	O1-C1-C2	-3.38	104.20	111.67
2	A	580	F6P	O6-C6-C5	-2.66	99.93	108.99
2	A	580	F6P	O6-P-O1P	2.24	112.49	106.44
2	B	580	F6P	O3-C3-C4	-2.18	105.57	113.25

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	580	F6P	C2
2	B	580	F6P	C2

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	580	F6P	C6-O6-P-O1P
2	A	580	F6P	C4-C5-C6-O6
2	A	580	F6P	O5-C5-C6-O6
2	A	580	F6P	C6-O6-P-O2P

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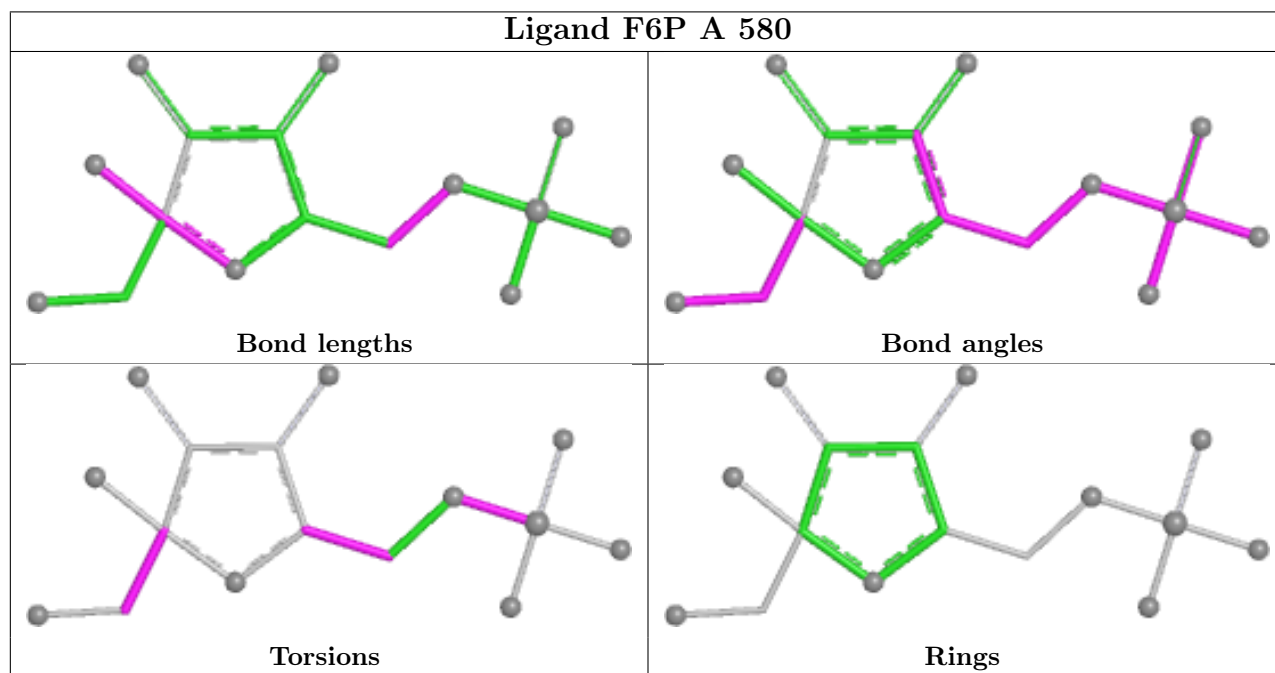
Mol	Chain	Res	Type	Atoms
2	A	580	F6P	O1-C1-C2-C3
2	B	580	F6P	O1-C1-C2-C3

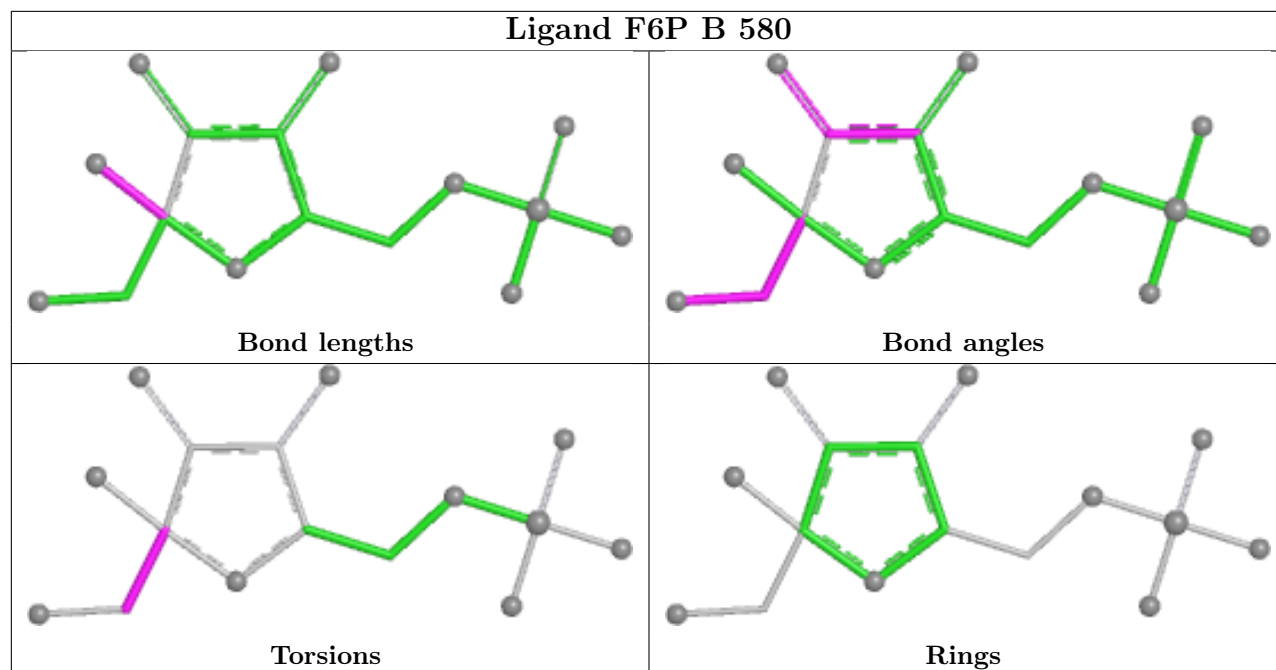
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	580	F6P	1	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.





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	578/597 (96%)	-0.63	0 100 100	12, 25, 49, 67	13 (2%)
1	B	577/597 (96%)	-0.69	0 100 100	13, 23, 48, 61	14 (2%)
All	All	1155/1194 (96%)	-0.66	0 100 100	12, 24, 48, 67	27 (2%)

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

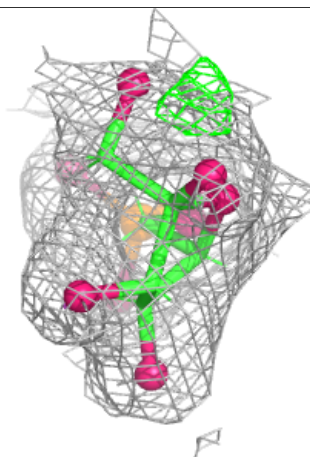
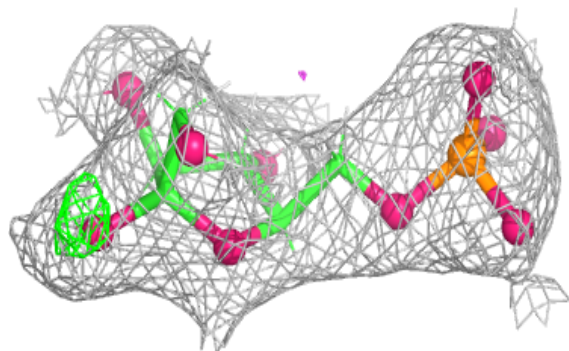
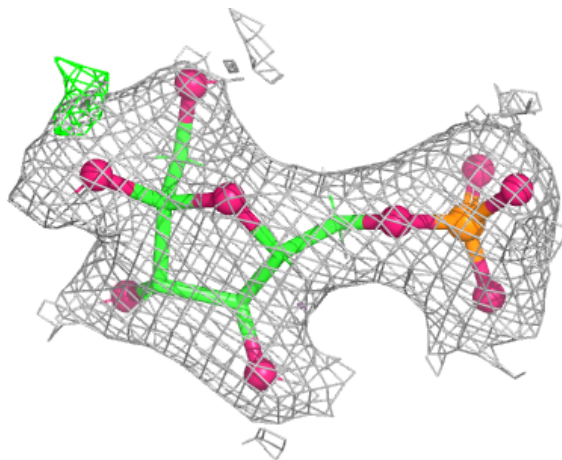
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PO4	A	581	5/5	0.87	0.12	50,52,57,58	0
3	PO4	B	581	5/5	0.94	0.10	33,35,38,42	0
2	F6P	A	580	16/16	0.96	0.06	24,27,33,44	0
2	F6P	B	580	16/16	0.98	0.05	18,21,27,35	0

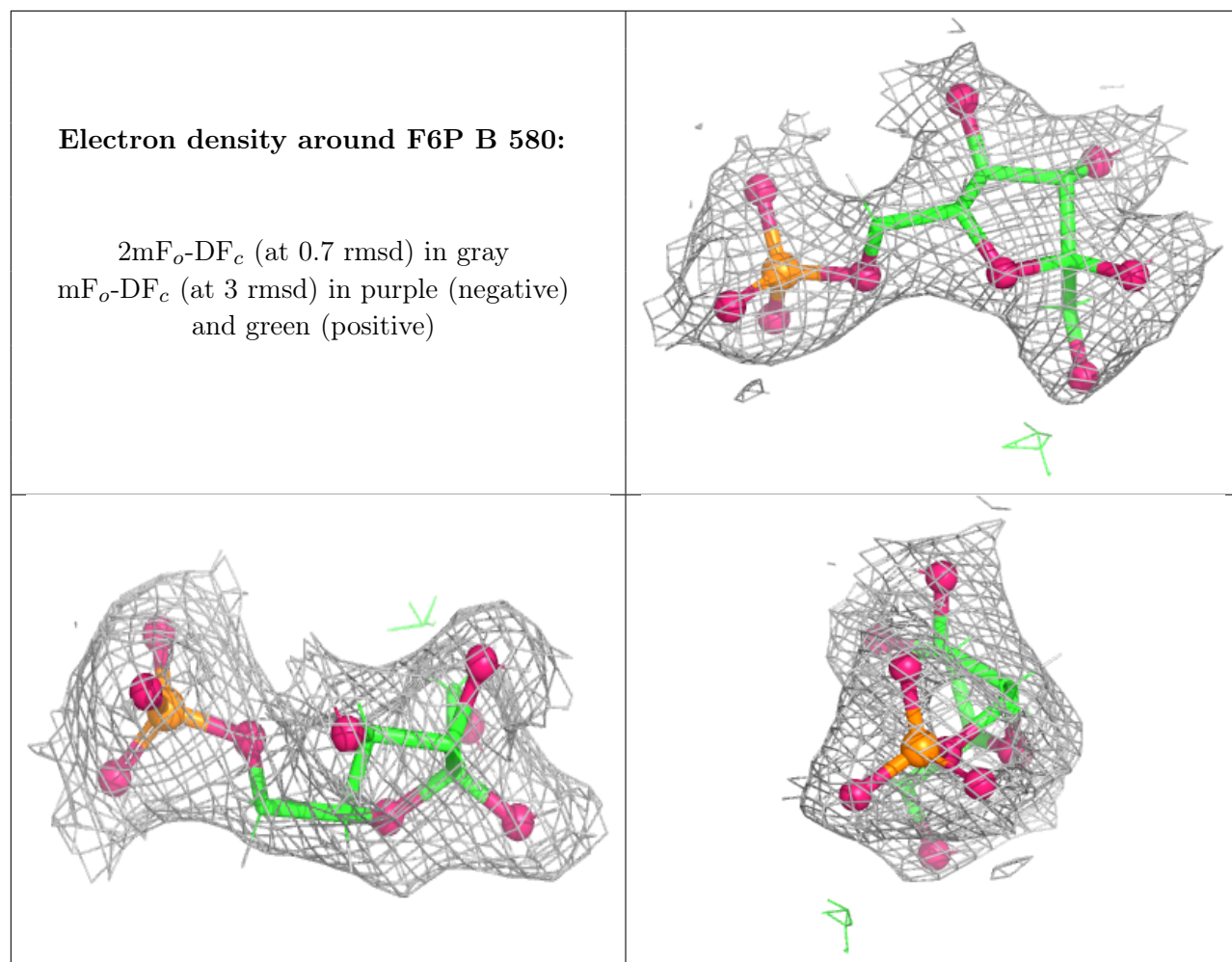
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

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 F6P A 580:

$2mF_o-DF_c$ (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.