



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 02:20 PM EDT

PDB ID : 6LKG / pdb_00006lkg
Title : two-component system protein mediate signal transduction
Authors : Wang, M.; Tao, Y.
Deposited on : 2019-12-19
Resolution : 1.95 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

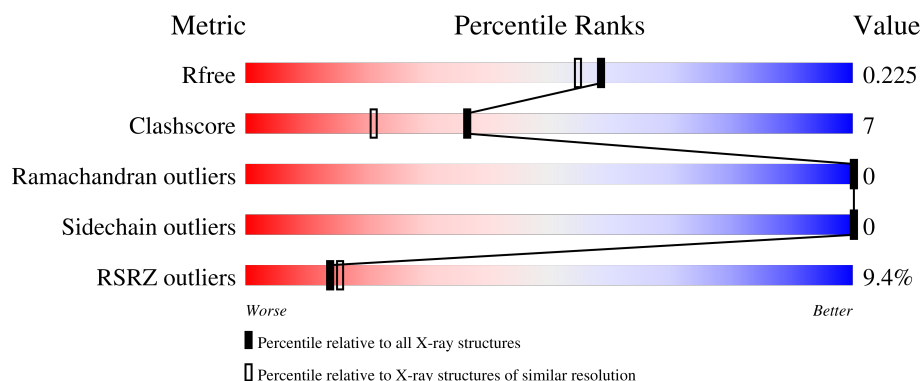
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	B	172	5% 85% 13% .
1	D	172	5% 92% 7% .
2	A	294	6% 85% 14%
2	C	294	17% 84% 13% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8224 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 Sensor protein kinase HptS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	170	Total	C	N	O	S	0	0	0
			1429	903	253	269	4			
1	D	171	Total	C	N	O	S	0	0	0
			1438	907	254	273	4			

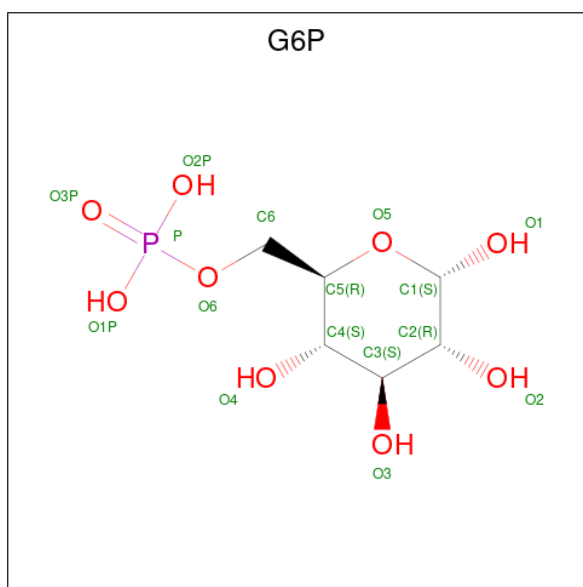
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	44	SER	-	expression tag	UNP Q2G1E0
B	68	SER	ILE	conflict	UNP Q2G1E0
D	44	SER	-	expression tag	UNP Q2G1E0
D	68	SER	ILE	conflict	UNP Q2G1E0

- Molecule 2 is a protein called ABC transporter, solute-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	294	Total	C	N	O	S	0	0	0
			2376	1506	413	450	7			
2	C	294	Total	C	N	O	S	0	0	0
			2376	1506	413	450	7			

- Molecule 3 is 6-O-phosphono-alpha-D-glucopyranose (CCD ID: G6P) (formula: C₆H₁₃O₉P) (labeled as "Ligand of Interest" by depositor).



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

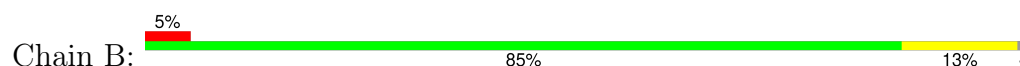
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	125	Total	O	0	0
			125	125		
4	A	165	Total	O	0	0
			165	165		
4	C	166	Total	O	0	0
			166	166		
4	D	117	Total	O	0	0
			117	117		

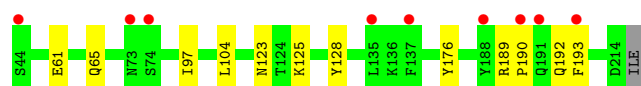
3 Residue-property plots [i](#)

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.

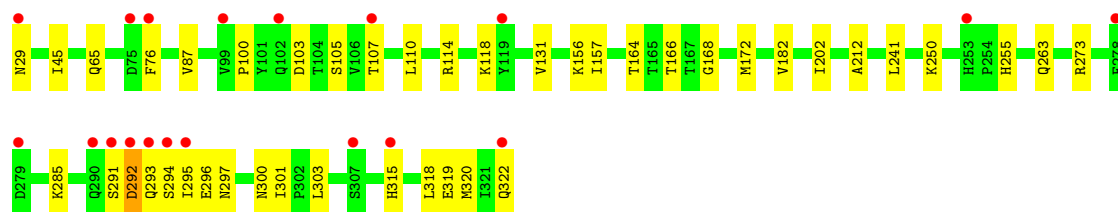
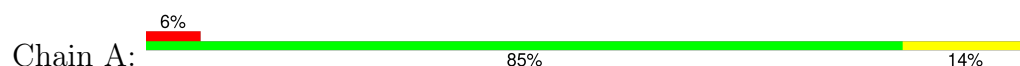
- Molecule 1: Sensor protein kinase HptS



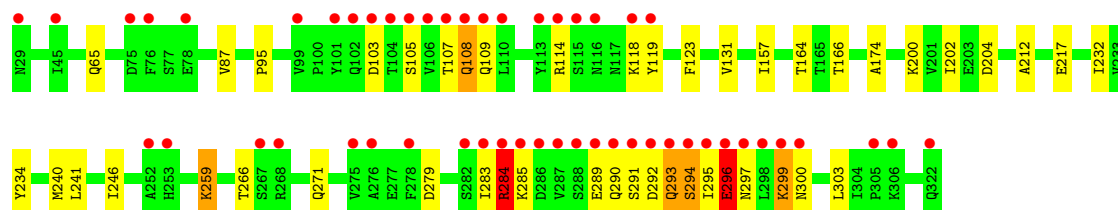
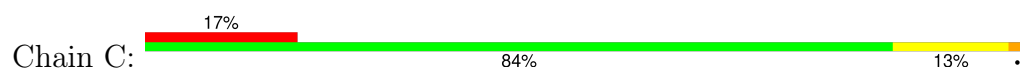
- Molecule 1: Sensor protein kinase HptS



- Molecule 2: ABC transporter, solute-binding protein



- Molecule 2: ABC transporter, solute-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.30Å 89.97Å 92.83Å 90.00° 100.87° 90.00°	Depositor
Resolution (Å)	38.16 – 1.95 38.16 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (38.16-1.95) 99.7 (38.16-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.204 , 0.225 0.204 , 0.225	Depositor DCC
R_{free} test set	4331 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.191	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8224	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: G6P

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	B	0.27	0/1463	0.39	0/1977
1	D	0.30	0/1472	0.40	1/1989 (0.1%)
2	A	0.15	0/2428	0.46	1/3288 (0.0%)
2	C	0.38	2/2428 (0.1%)	0.83	12/3288 (0.4%)
All	All	0.29	2/7791 (0.0%)	0.58	14/10542 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	283	ILE	C-N	11.76	1.50	1.33
2	C	299	LYS	C-O	-6.08	1.16	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	284	ARG	O-C-N	-19.14	100.95	123.16
2	C	283	ILE	CA-C-N	15.66	145.26	122.09
2	C	283	ILE	C-N-CA	15.66	145.26	122.09
2	C	285	LYS	N-CA-C	-11.08	99.28	111.36
2	C	296	GLU	N-CA-C	10.44	123.81	110.24
2	A	292	ASP	N-CA-C	-10.42	88.61	110.80
2	C	293	GLN	N-CA-C	-9.79	92.33	108.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	294	SER	N-CA-C	8.00	119.90	111.03
2	C	284	ARG	CB-CA-C	6.41	120.10	109.53
1	D	193	PHE	N-CA-C	5.75	118.59	109.50
2	C	108	GLN	CA-CB-CG	-5.75	102.60	114.10
2	C	259	LYS	CB-CG-CD	5.28	123.43	111.30
2	C	107	THR	CA-C-N	-5.22	114.46	122.60
2	C	107	THR	C-N-CA	-5.22	114.46	122.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	284	ARG	Mainchain
2	C	296	GLU	Peptide

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	B	1429	0	1377	15	0
1	D	1438	0	1384	9	0
2	A	2376	0	2373	39	0
2	C	2376	0	2373	46	0
3	A	16	0	11	0	0
3	C	16	0	11	0	0
4	A	165	0	0	1	0
4	B	125	0	0	3	0
4	C	166	0	0	5	0
4	D	117	0	0	0	0
All	All	8224	0	7529	109	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 7.

All (109) 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)
2:C:232:ILE:HB	2:C:295:ILE:HD13	1.21	1.21
2:C:232:ILE:HB	2:C:295:ILE:CD1	2.02	0.87
2:C:103:ASP:OD1	2:C:105:SER:HB3	1.75	0.87
2:C:232:ILE:HD12	2:C:295:ILE:HD12	1.56	0.85
2:A:294:SER:O	2:A:295:ILE:HD12	1.78	0.82
2:C:294:SER:C	2:C:295:ILE:HG13	2.04	0.82
2:A:293:GLN:HG2	2:A:294:SER:HA	1.64	0.79
1:B:189:ARG:NH1	1:B:192:GLN:OE1	2.17	0.78
2:C:290:GLN:NE2	2:C:292:ASP:OD2	2.19	0.76
2:A:295:ILE:HG23	2:A:296:GLU:OE2	1.86	0.76
2:A:294:SER:C	2:A:295:ILE:HD12	2.11	0.76
2:C:118:LYS:HG3	4:C:575:HOH:O	1.86	0.75
2:C:234:TYR:HE1	2:C:295:ILE:HG23	1.53	0.74
2:A:103:ASP:OD2	2:A:105:SER:OG	2.06	0.73
2:C:297:ASN:ND2	2:C:299:LYS:H	1.88	0.71
2:C:234:TYR:HE1	2:C:295:ILE:CG2	2.03	0.71
1:B:201:LYS:NZ	4:B:303:HOH:O	2.22	0.70
2:A:291:SER:O	2:A:292:ASP:C	2.31	0.70
2:C:294:SER:O	2:C:295:ILE:HG13	1.93	0.68
1:D:123:ASN:HD21	1:D:125:LYS:HE3	1.59	0.68
2:C:234:TYR:CE1	2:C:295:ILE:CG2	2.77	0.68
2:C:234:TYR:OH	2:C:295:ILE:HG22	1.94	0.67
2:A:293:GLN:HA	2:A:294:SER:O	1.94	0.67
2:C:232:ILE:CB	2:C:295:ILE:HD13	2.13	0.66
2:C:291:SER:CB	2:C:297:ASN:HB3	2.26	0.65
2:C:95:PRO:HB3	2:C:119:TYR:CE2	2.32	0.65
2:C:291:SER:HB3	2:C:297:ASN:HB3	1.79	0.65
2:A:293:GLN:CG	2:A:294:SER:HA	2.27	0.64
2:A:318:LEU:O	2:A:322:GLN:HG2	2.00	0.62
2:A:291:SER:O	2:A:293:GLN:N	2.33	0.61
2:C:279:ASP:OD2	4:C:501:HOH:O	2.16	0.61
2:C:296:GLU:CD	4:C:502:HOH:O	2.43	0.61
2:A:45:ILE:CD1	2:A:273:ARG:HD3	2.32	0.60
2:A:156:LYS:NZ	4:A:501:HOH:O	2.22	0.59
2:A:294:SER:O	2:A:295:ILE:CD1	2.48	0.59
2:C:95:PRO:HB3	2:C:119:TYR:HE2	1.68	0.59
2:A:315:HIS:O	2:A:319:GLU:HG3	2.04	0.57
2:C:109:GLN:HA	2:C:284:ARG:HA	1.86	0.56
2:A:87:VAL:HG22	2:A:166:THR:HG23	1.88	0.56
2:C:87:VAL:HG22	2:C:166:THR:HG23	1.87	0.55
1:B:47:HIS:HD2	1:B:198:GLU:OE2	1.89	0.55
2:C:289:GLU:OE2	2:C:297:ASN:HB2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:100:PRO:HB3	2:A:118:LYS:HD3	1.88	0.54
2:C:232:ILE:HD12	2:C:295:ILE:CD1	2.33	0.54
2:A:296:GLU:CG	2:A:300:ASN:HB2	2.37	0.54
2:C:234:TYR:OH	2:C:295:ILE:CG2	2.56	0.53
2:A:296:GLU:HG3	2:A:297:ASN:H	1.74	0.52
2:C:118:LYS:HG3	4:C:641:HOH:O	2.08	0.52
1:B:87:THR:OG1	4:B:301:HOH:O	2.18	0.52
2:A:293:GLN:HA	2:A:294:SER:C	2.34	0.52
2:C:109:GLN:O	2:C:114:ARG:NH1	2.37	0.52
1:D:123:ASN:ND2	1:D:125:LYS:HG2	2.24	0.52
2:C:241:LEU:HB2	2:C:303:LEU:HD22	1.92	0.51
2:A:296:GLU:HG3	2:A:297:ASN:N	2.24	0.51
1:D:189:ARG:HB3	1:D:192:GLN:HG3	1.93	0.51
1:B:128:TYR:HB2	1:B:176:TYR:HB2	1.93	0.50
2:A:295:ILE:HG23	2:A:296:GLU:HB2	1.94	0.50
2:A:157:ILE:HG22	2:A:212:ALA:HB3	1.94	0.49
2:C:259:LYS:HG3	4:C:625:HOH:O	2.13	0.49
1:D:128:TYR:HB2	1:D:176:TYR:HB2	1.94	0.49
1:B:82:ASN:OD1	1:B:84:HIS:HD2	1.96	0.49
2:A:107:THR:HA	2:A:114:ARG:NH1	2.29	0.48
2:C:297:ASN:HD22	2:C:299:LYS:H	1.60	0.48
1:B:48:GLN:OE1	4:B:302:HOH:O	2.20	0.47
1:B:47:HIS:CD2	1:B:198:GLU:OE2	2.67	0.47
1:B:164:PHE:HB2	1:B:212:TYR:CZ	2.49	0.47
2:C:103:ASP:OD2	2:C:266:THR:OG1	2.24	0.47
1:B:189:ARG:NH1	1:B:192:GLN:HB2	2.30	0.47
2:C:65:GLN:HB2	2:C:164:THR:HA	1.96	0.47
1:B:54:GLN:HB2	1:B:200:VAL:HG11	1.97	0.47
2:C:217:GLU:OE2	2:C:234:TYR:OH	2.26	0.47
2:A:65:GLN:HB2	2:A:164:THR:HA	1.96	0.46
2:C:108:GLN:HE21	2:C:284:ARG:HH22	1.63	0.46
2:A:29:ASN:O	2:A:255:HIS:NE2	2.46	0.46
2:C:294:SER:C	2:C:295:ILE:CG1	2.82	0.46
2:A:76:PHE:HE1	2:A:250:LYS:HG2	1.80	0.46
1:B:59:HIS:O	1:B:63:GLN:HG3	2.16	0.46
2:C:157:ILE:HG22	2:C:212:ALA:HB3	1.98	0.45
1:B:46:ILE:O	1:B:50:VAL:HG23	2.17	0.45
1:B:58:HIS:CE1	1:B:203:ASN:HD22	2.35	0.45
2:A:131:VAL:HG11	2:A:202:ILE:HD13	1.99	0.45
2:A:182:VAL:HG22	2:A:320:MET:HE3	1.97	0.45
2:A:45:ILE:HD11	2:A:273:ARG:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:294:SER:O	2:C:295:ILE:CG1	2.65	0.44
2:C:292:ASP:O	2:C:293:GLN:C	2.59	0.44
2:A:241:LEU:HB2	2:A:303:LEU:HD22	2.00	0.44
2:A:315:HIS:CD2	2:A:319:GLU:OE2	2.70	0.44
2:A:110:LEU:HD23	2:A:285:LYS:HG2	1.99	0.43
1:D:61:GLU:OE2	1:D:65:GLN:NE2	2.48	0.43
2:C:234:TYR:CE1	2:C:295:ILE:HG23	2.42	0.42
1:D:123:ASN:CG	1:D:125:LYS:HG2	2.45	0.42
2:C:290:GLN:HE22	2:C:292:ASP:CG	2.23	0.42
2:A:295:ILE:CG2	2:A:296:GLU:HB2	2.49	0.42
2:C:131:VAL:HG11	2:C:202:ILE:HD13	2.01	0.42
2:C:174:ALA:HB2	2:C:240:MET:HB2	2.02	0.42
2:A:103:ASP:OD1	2:A:263:GLN:NE2	2.53	0.42
2:A:297:ASN:O	2:A:301:ILE:HG13	2.20	0.41
2:C:123:PHE:HB2	2:C:246:ILE:HG13	2.02	0.41
2:A:168:GLY:O	2:A:172:MET:HG2	2.21	0.41
2:C:296:GLU:CD	2:C:300:ASN:O	2.64	0.41
2:C:200:LYS:HE3	2:C:204:ASP:OD2	2.20	0.41
1:D:97:ILE:HD12	1:D:104:LEU:HB2	2.03	0.41
1:B:72:ASN:O	1:B:76:GLN:HG3	2.21	0.41
1:D:189:ARG:HA	1:D:190:PRO:HD2	1.95	0.41
2:A:291:SER:OG	2:A:292:ASP:N	2.53	0.41
1:D:189:ARG:HA	1:D:189:ARG:HD3	1.88	0.40
2:C:271:GLN:HG3	2:C:284:ARG:HD3	2.02	0.40
2:A:292:ASP:HB2	2:A:293:GLN:H	1.76	0.40
2:A:296:GLU:HG2	2:A:300:ASN:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	B	168/172 (98%)	166 (99%)	2 (1%)	0	100	100
1	D	169/172 (98%)	168 (99%)	1 (1%)	0	100	100
2	A	292/294 (99%)	282 (97%)	10 (3%)	0	100	100
2	C	292/294 (99%)	279 (96%)	13 (4%)	0	100	100
All	All	921/932 (99%)	895 (97%)	26 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	B	160/163 (98%)	160 (100%)	0	100	100
1	D	162/163 (99%)	162 (100%)	0	100	100
2	A	271/271 (100%)	271 (100%)	0	100	100
2	C	271/271 (100%)	271 (100%)	0	100	100
All	All	864/868 (100%)	864 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	47	HIS
1	B	63	GLN
1	B	72	ASN
1	B	83	HIS
1	B	84	HIS
1	B	121	HIS
1	B	203	ASN
2	A	29	ASN
2	A	38	GLN
2	A	109	GLN

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Mol	Chain	Res	Type
2	A	163	ASN
2	A	194	GLN
2	A	263	GLN
2	A	290	GLN
2	A	322	GLN
2	C	38	GLN
2	C	108	GLN
2	C	116	ASN
2	C	137	GLN
2	C	163	ASN
2	C	297	ASN
2	C	315	HIS
1	D	123	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](#)

2 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	$\# Z > 2$
3	G6P	A	401	-	16,16,16	1.49	4 (25%)	23,24,24	1.49	3 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	G6P	C	401	-	16,16,16	1.50	3 (18%)	23,24,24	1.46	3 (13%)

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
3	G6P	A	401	-	-	0/6/26/26	0/1/1/1
3	G6P	C	401	-	-	0/6/26/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	G6P	P-O6	3.00	1.69	1.60
3	C	401	G6P	P-O6	2.91	1.69	1.60
3	A	401	G6P	O3-C3	2.85	1.50	1.43
3	C	401	G6P	O3-C3	2.83	1.50	1.43
3	A	401	G6P	C4-C3	-2.16	1.46	1.52
3	C	401	G6P	C4-C3	-2.08	1.46	1.52
3	A	401	G6P	O5-C1	2.02	1.47	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	401	G6P	O5-C1-C2	4.97	119.04	110.30
3	A	401	G6P	O5-C1-C2	4.96	119.03	110.30
3	C	401	G6P	C1-O5-C5	2.68	118.83	113.65
3	A	401	G6P	C1-C2-C3	2.67	115.80	110.36
3	A	401	G6P	C1-O5-C5	2.53	118.55	113.65
3	C	401	G6P	C1-C2-C3	2.45	115.36	110.36

There are no chirality outliers.

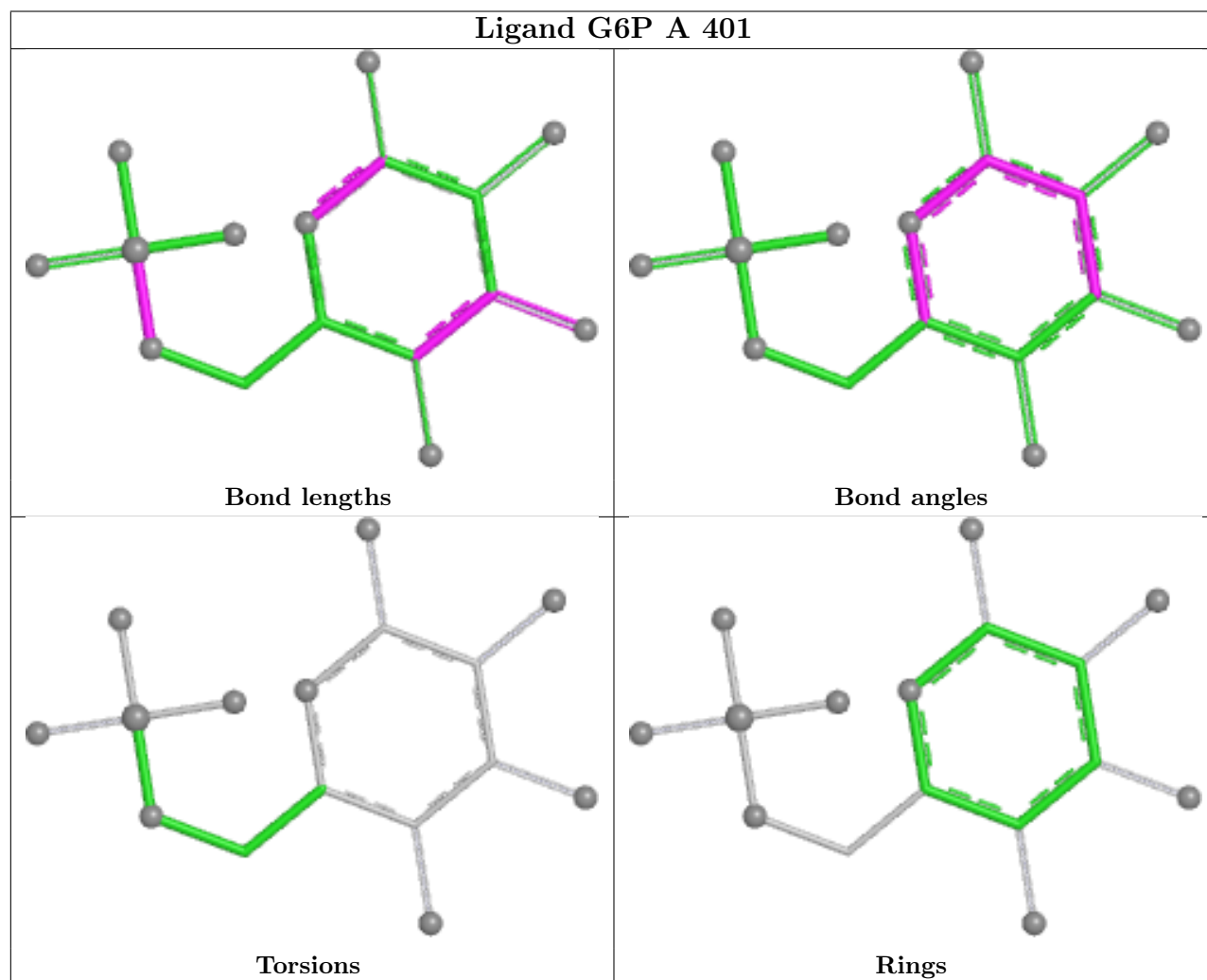
There are no torsion outliers.

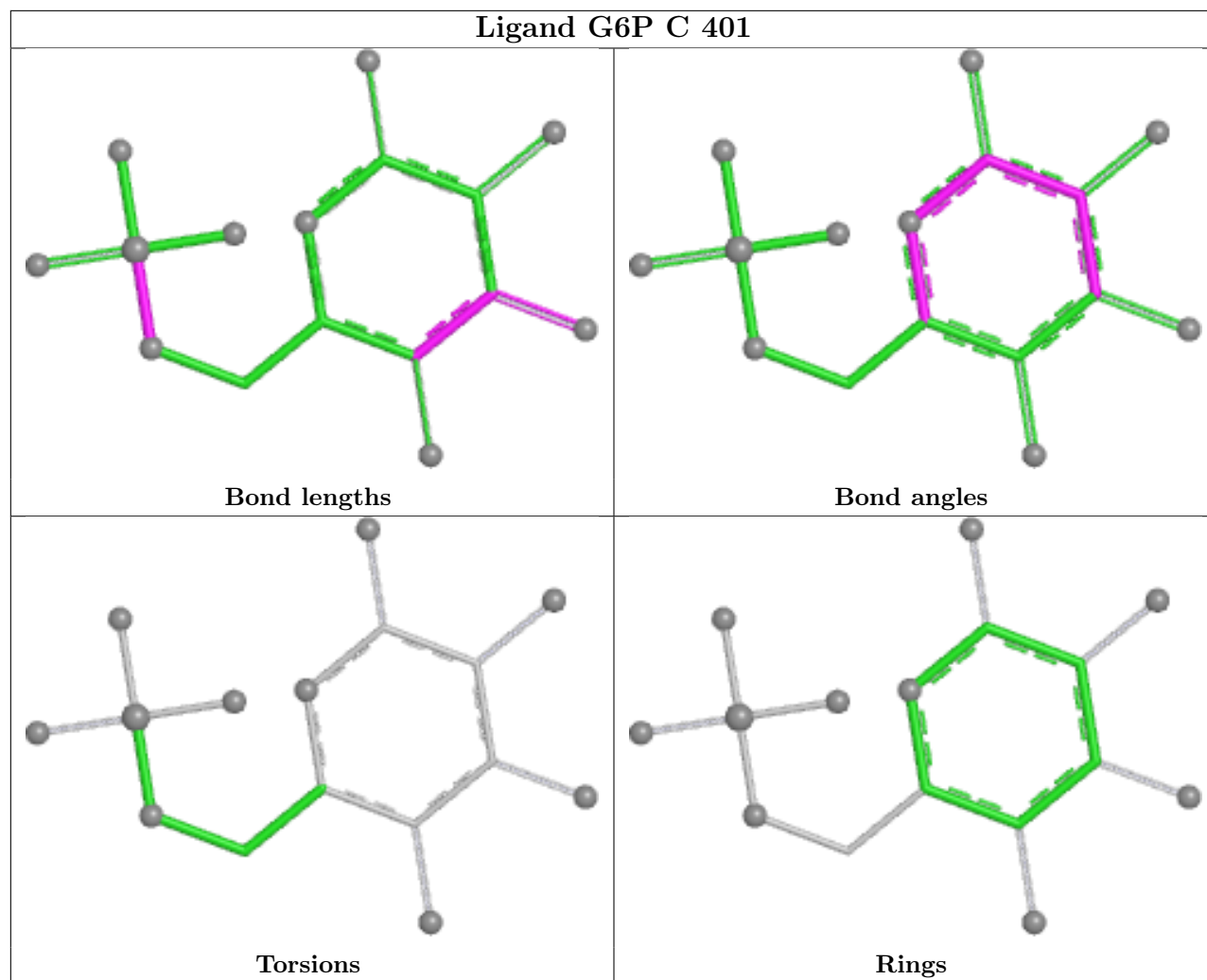
There are no ring outliers.

No monomer is involved in short contacts.

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

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	B	170/172 (98%)	0.57	8 (4%) 36 41	27, 39, 54, 71	0
1	D	171/172 (99%)	0.55	9 (5%) 32 36	26, 38, 62, 76	0
2	A	294/294 (100%)	0.54	19 (6%) 25 28	26, 38, 62, 109	0
2	C	294/294 (100%)	0.77	51 (17%) 4 4	23, 37, 76, 127	0
All	All	929/932 (99%)	0.62	87 (9%) 14 15	23, 38, 66, 127	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	295	ILE	7.3
2	A	295	ILE	7.2
2	C	296	GLU	6.8
2	C	110	LEU	5.6
2	C	105	SER	5.4
2	C	294	SER	5.1
1	B	111	PRO	4.9
2	C	106	VAL	4.8
2	C	109	GLN	4.5
2	C	283	ILE	4.4
2	C	289	GLU	4.4
1	B	112	ASN	4.4
2	A	294	SER	4.4
1	B	214	ASP	4.4
2	C	107	THR	4.4
1	B	83	HIS	4.2
2	A	29	ASN	4.1
2	C	287	VAL	4.1
2	C	291	SER	4.1
1	D	193	PHE	4.1
2	A	76	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
2	C	286	ASP	4.0
2	A	291	SER	4.0
2	C	284	ARG	3.9
2	C	118	LYS	3.8
2	C	288	SER	3.7
2	A	102	GLN	3.6
2	C	119	TYR	3.6
2	C	300	ASN	3.4
2	C	102	GLN	3.3
2	C	285	LYS	3.2
2	C	108	GLN	3.2
1	D	191	GLN	3.1
2	C	114	ARG	3.1
2	A	292	ASP	3.1
2	C	76	PHE	3.1
2	C	103	ASP	2.9
2	C	275	VAL	2.9
1	D	73	ASN	2.9
2	C	290	GLN	2.8
2	C	252	ALA	2.8
2	C	268	ARG	2.7
1	B	113	GLN	2.7
2	A	322	GLN	2.7
2	C	116	ASN	2.7
2	C	298	LEU	2.7
2	C	299	LYS	2.6
2	C	104	THR	2.6
2	C	297	ASN	2.5
2	A	307	SER	2.5
2	A	279	ASP	2.5
2	C	306	LYS	2.5
2	C	113	TYR	2.5
2	C	305	PRO	2.5
1	D	44	SER	2.4
2	A	107	THR	2.4
1	D	190	PRO	2.4
2	A	75	ASP	2.4
2	C	276	ALA	2.4
2	A	119	TYR	2.3
2	C	29	ASN	2.3
2	C	253	HIS	2.3
2	C	115	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	188	TYR	2.3
2	C	278	PHE	2.3
1	D	74	SER	2.2
2	C	282	SER	2.2
2	C	99	VAL	2.2
2	C	322	GLN	2.2
2	C	292	ASP	2.2
1	D	135	LEU	2.2
1	B	84	HIS	2.1
2	C	78	GLU	2.1
2	C	293	GLN	2.1
2	C	75	ASP	2.1
1	B	82	ASN	2.1
2	A	293	GLN	2.1
2	C	45	ILE	2.1
2	A	253	HIS	2.1
1	D	137	PHE	2.0
2	A	278	PHE	2.0
2	C	267	SER	2.0
1	B	196	MET	2.0
2	A	290	GLN	2.0
2	A	315	HIS	2.0
2	C	101	TYR	2.0
2	A	99	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

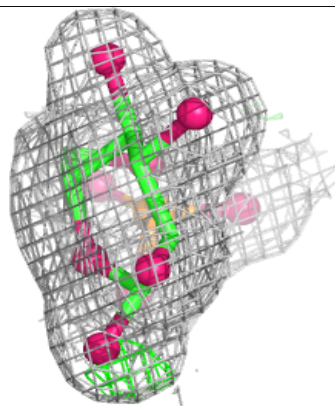
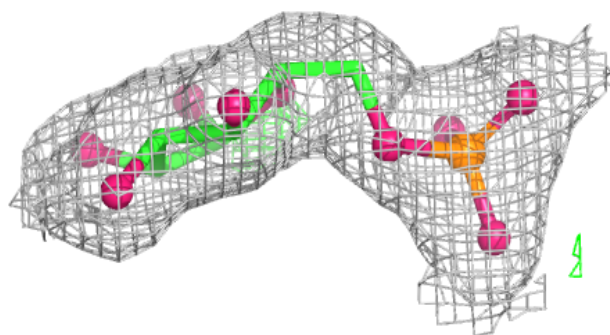
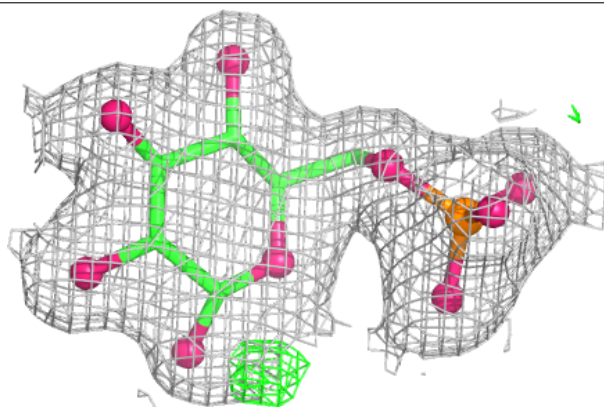
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	G6P	A	401	16/16	0.97	0.07	25,30,34,35	0
3	G6P	C	401	16/16	0.97	0.06	21,25,29,29	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.

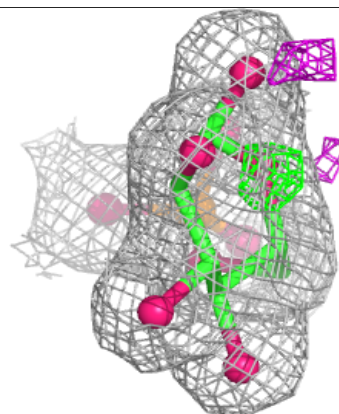
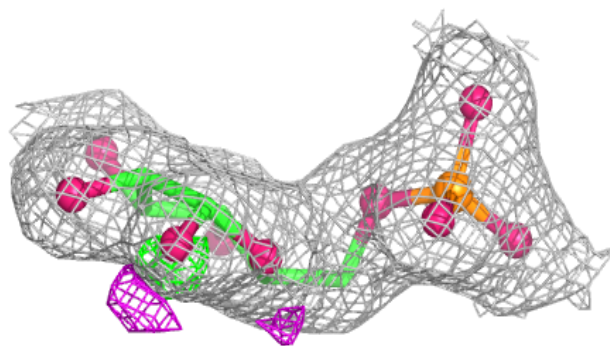
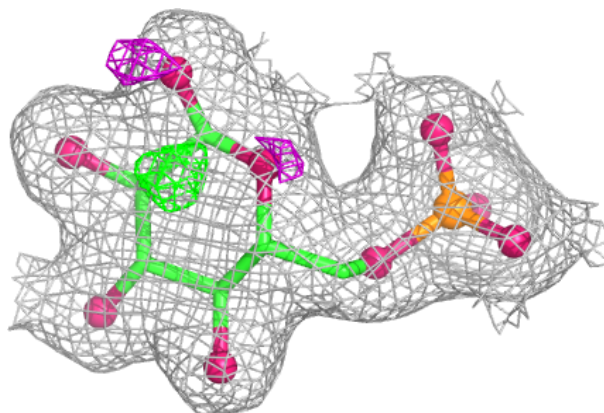
Electron density around G6P A 401:

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 401:

$2mF_o-DF_c$ (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.