



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:07 AM UTC

PDB ID : 4GX6 / pdb_00004gx6
Title : AMP Complexes of Porcine Liver Fructose-1,6-bisphosphatase with Mutation E192Q
Authors : Gao, Y.; Honzatko, R.B.
Deposited on : 2012-09-03
Resolution : 2.50 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

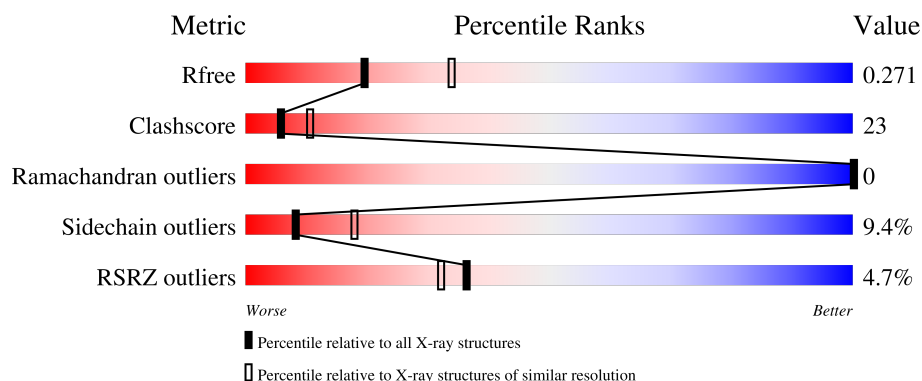
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	337	5% 62% 29% 6%
1	B	337	4% 55% 35% 6%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5304 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.

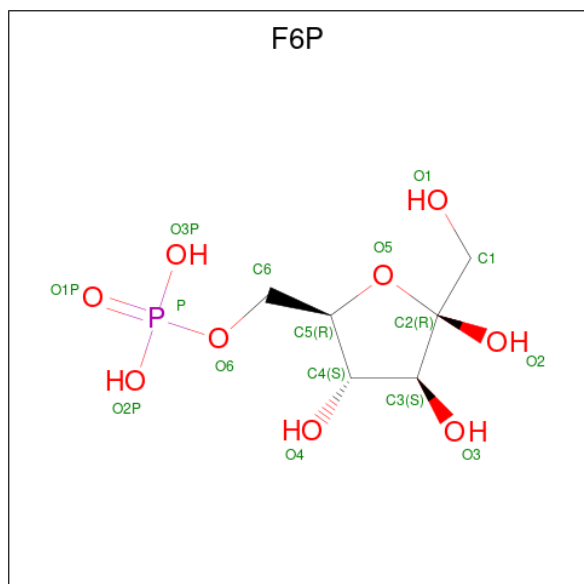
- Molecule 1 is a protein called Fructose-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	327	Total	C	N	O	S	0	0	0
			2496	1587	421	473	15			
1	B	327	Total	C	N	O	S	0	0	0
			2496	1587	421	473	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	192	GLN	GLU	engineered mutation	UNP P00636
B	192	GLN	GLU	engineered mutation	UNP P00636

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



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

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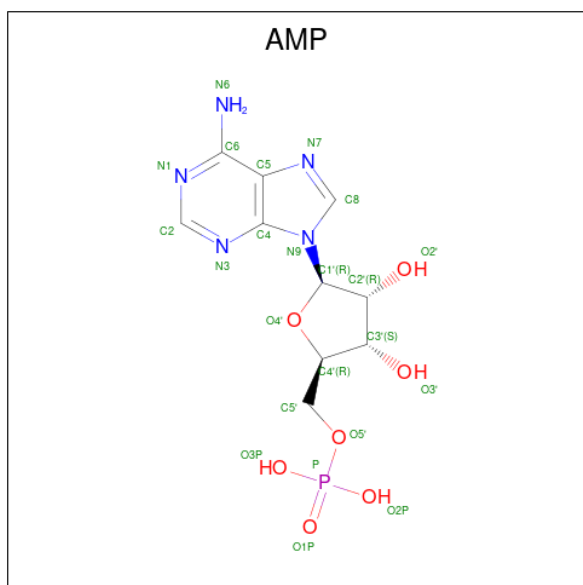
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

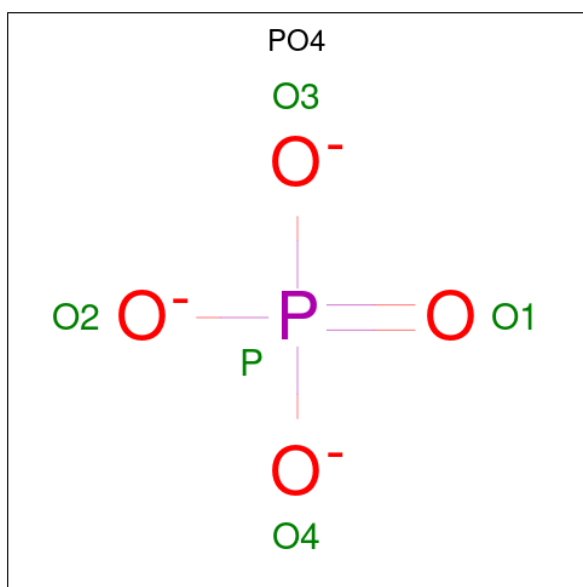
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	B	2	Total	Mg	0	0
			2	2		

- Molecule 4 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

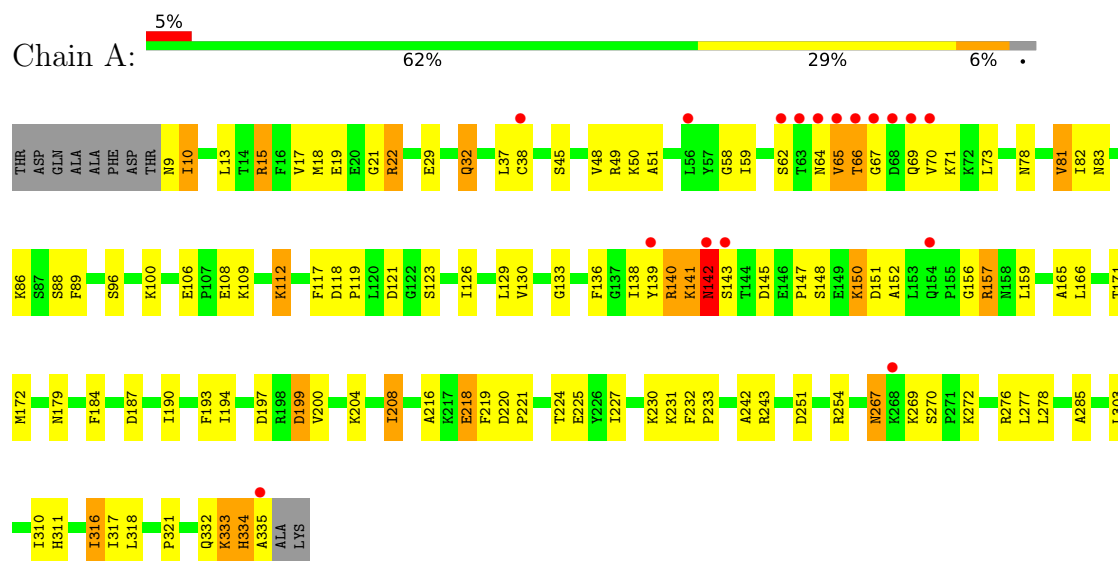
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	111	Total	O	0	0
			111	111		
6	B	109	Total	O	0	0
			109	109		

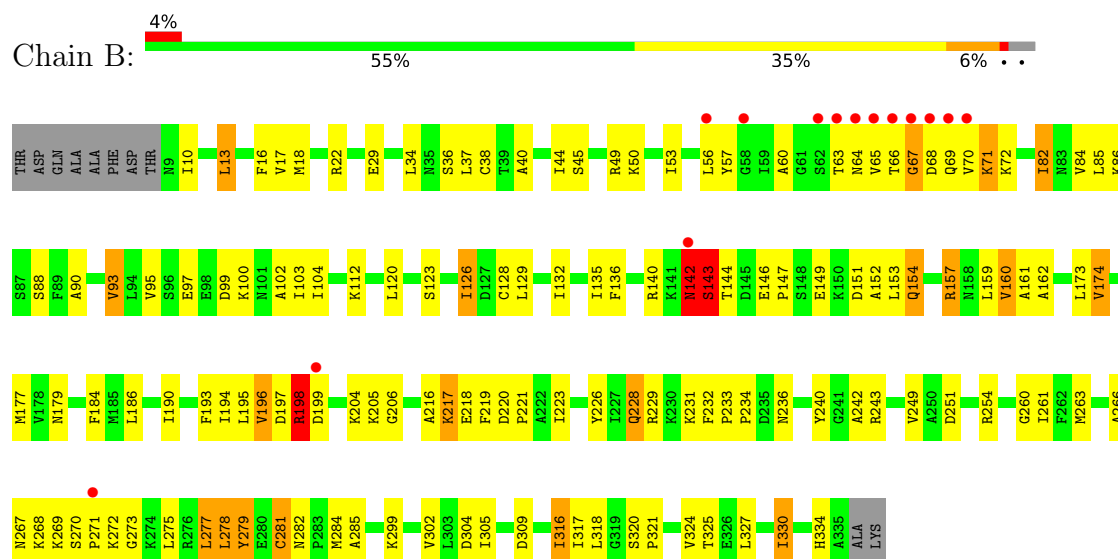
3 Residue-property plots [i](#)

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: Fructose-1,6-bisphosphatase 1



• Molecule 1: Fructose-1,6-bisphosphatase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	59.92Å 165.34Å 79.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.56 – 2.50 39.56 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.2 (39.56-2.50) 92.2 (39.56-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.211 , 0.273 0.211 , 0.271	Depositor DCC
R_{free} test set	1312 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5304	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.79% 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: F6P, PO4, AMP, MG

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	1.28	8/2538 (0.3%)	1.23	10/3434 (0.3%)
1	B	1.40	24/2538 (0.9%)	1.27	13/3434 (0.4%)
All	All	1.34	32/5076 (0.6%)	1.25	23/6868 (0.3%)

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
1	A	0	1
1	B	0	1
All	All	0	2

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	282	ASN	C-O	-8.11	1.16	1.24
1	B	197	ASP	C-O	-6.82	1.16	1.23
1	B	263	MET	C-O	-6.71	1.16	1.24
1	B	93	VAL	C-O	-6.58	1.17	1.24
1	B	279	TYR	C-O	-6.56	1.15	1.24
1	A	242	ALA	C-O	-6.34	1.16	1.24
1	B	95	VAL	C-O	-6.16	1.17	1.24
1	A	140	ARG	C-O	-6.10	1.16	1.23
1	B	160	VAL	C-O	-5.99	1.17	1.24
1	B	162	ALA	C-O	-5.91	1.16	1.23
1	B	90	ALA	C-O	-5.85	1.16	1.24
1	B	195	LEU	C-O	-5.84	1.17	1.24
1	B	196	VAL	C-O	-5.81	1.17	1.24
1	B	60	ALA	CA-CB	-5.79	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	281	CYS	C-O	-5.63	1.17	1.24
1	B	266	ALA	CA-CB	-5.44	1.44	1.53
1	A	243	ARG	C-O	-5.38	1.17	1.23
1	B	278	LEU	C-O	-5.36	1.17	1.24
1	A	200	VAL	C-O	-5.36	1.18	1.24
1	B	198	ARG	C-O	-5.34	1.17	1.24
1	B	159	LEU	C-O	-5.33	1.17	1.23
1	A	152	ALA	CA-CB	-5.33	1.44	1.53
1	A	270	SER	C-O	-5.30	1.20	1.24
1	A	51	ALA	CA-CB	-5.27	1.44	1.53
1	A	81	VAL	C-O	-5.26	1.18	1.24
1	B	153	LEU	C-O	-5.19	1.17	1.23
1	B	174	VAL	CA-C	-5.17	1.46	1.52
1	B	285	ALA	CA-CB	-5.17	1.44	1.53
1	B	161	ALA	C-O	-5.12	1.17	1.23
1	B	151	ASP	C-O	-5.08	1.17	1.24
1	B	16	PHE	C-O	-5.07	1.18	1.24
1	B	197	ASP	CA-C	-5.04	1.46	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	143	SER	N-CA-C	13.07	129.36	113.41
1	A	334	HIS	N-CA-C	-10.22	101.07	112.72
1	B	281	CYS	N-CA-C	9.85	121.78	111.14
1	B	142	ASN	N-CA-C	8.99	121.16	111.36
1	B	278	LEU	N-CA-C	8.85	120.70	111.14
1	A	142	ASN	N-CA-C	7.42	119.45	111.36
1	B	50	LYS	N-CA-C	7.16	121.31	112.58
1	B	160	VAL	N-CA-C	6.63	117.42	110.72
1	A	83	ASN	N-CA-C	6.58	118.54	111.36
1	B	67	GLY	N-CA-C	-6.17	99.54	111.02
1	A	199	ASP	N-CA-C	5.91	119.54	111.39
1	B	232	PHE	CA-C-N	-5.86	114.35	120.38
1	B	232	PHE	C-N-CA	-5.86	114.35	120.38
1	A	200	VAL	N-CA-C	5.83	117.24	109.37
1	B	126	ILE	N-CA-C	5.62	115.82	110.42
1	B	154	GLN	CA-C-N	-5.58	114.51	120.03
1	B	154	GLN	C-N-CA	-5.58	114.51	120.03
1	A	138	ILE	N-CA-C	5.30	115.54	108.11
1	B	152	ALA	N-CA-C	5.23	117.88	111.82
1	A	243	ARG	N-CA-C	5.12	116.76	108.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	THR	N-CA-C	5.04	116.78	111.28
1	A	285	ALA	N-CA-C	-5.04	105.87	111.36
1	A	129	LEU	N-CA-C	5.01	118.66	112.24

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	64	ASN	Peptide
1	B	142	ASN	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	A	2496	0	2554	104	0
1	B	2496	0	2554	134	0
2	A	16	0	11	0	0
2	B	16	0	11	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	23	0	12	1	0
4	B	23	0	12	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	111	0	0	16	0
6	B	109	0	0	24	0
All	All	5304	0	5154	236	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 23.

All (236) 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:13:LEU:HD23	1:A:38:CYS:SG	1.69	1.32

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ILE:HD11	1:B:249:VAL:CG2	1.68	1.23
1:A:13:LEU:CD2	1:A:38:CYS:SG	2.39	1.10
1:B:135:ILE:HD11	1:B:249:VAL:HG22	1.25	1.10
1:B:299:LYS:HE2	6:B:602:HOH:O	1.50	1.06
1:A:15:ARG:HD2	6:A:597:HOH:O	1.55	1.06
1:A:316:ILE:HD12	1:A:317:ILE:N	1.77	1.00
1:B:53:ILE:HG13	1:B:56:LEU:HD12	1.48	0.94
1:B:157:ARG:HG3	1:B:157:ARG:HH11	1.31	0.94
1:A:19:GLU:HG3	6:A:598:HOH:O	1.66	0.93
1:B:228:GLN:HA	1:B:228:GLN:OE1	1.71	0.90
1:B:135:ILE:CD1	1:B:249:VAL:HG22	2.02	0.86
1:B:65:VAL:HG13	1:B:69:GLN:HB2	1.58	0.85
1:B:53:ILE:CG1	1:B:56:LEU:HD12	2.06	0.85
1:A:70:VAL:HG22	6:A:578:HOH:O	1.76	0.83
1:B:302:VAL:HG21	1:B:316:ILE:HD13	1.61	0.82
1:B:157:ARG:HG3	1:B:157:ARG:NH1	1.94	0.80
1:A:18:MET:HE3	1:A:22:ARG:HD3	1.63	0.80
1:A:123:SER:HA	1:A:126:ILE:HG12	1.62	0.80
1:B:157:ARG:HH11	1:B:157:ARG:CG	1.95	0.79
1:A:204:LYS:O	1:A:321:PRO:HD2	1.83	0.79
1:B:120:LEU:HD11	1:B:132:ILE:HD12	1.63	0.79
1:A:139:TYR:CZ	1:A:159:LEU:HG	2.21	0.76
1:B:72:LYS:HB3	6:B:529:HOH:O	1.85	0.76
1:A:139:TYR:CZ	1:A:159:LEU:CD2	2.69	0.76
1:A:318:LEU:HD12	1:A:318:LEU:C	2.11	0.76
1:A:38:CYS:SG	1:A:193:PHE:CD2	2.79	0.76
1:B:135:ILE:CD1	1:B:249:VAL:CG2	2.60	0.75
1:B:330:ILE:HD13	1:B:330:ILE:N	2.03	0.73
1:B:218:GLU:HG2	1:B:268:LYS:HE3	1.71	0.72
1:B:70:VAL:HG12	1:B:71:LYS:N	2.05	0.71
1:B:18:MET:HE3	1:B:22:ARG:HD2	1.73	0.71
1:B:65:VAL:CG1	1:B:69:GLN:HB2	2.21	0.70
1:B:267:ASN:HA	6:B:596:HOH:O	1.91	0.69
1:A:65:VAL:O	1:A:65:VAL:CG1	2.39	0.69
1:B:330:ILE:CD1	6:B:518:HOH:O	2.40	0.69
1:A:221:PRO:HB2	1:A:334:HIS:CE1	2.28	0.69
1:B:18:MET:HE3	1:B:22:ARG:CD	2.22	0.69
1:B:63:THR:HG22	1:B:63:THR:O	1.92	0.69
1:B:321:PRO:O	1:B:325:THR:HG23	1.93	0.68
1:A:58:GLY:HA2	6:A:610:HOH:O	1.94	0.68
1:B:70:VAL:CG1	1:B:71:LYS:N	2.57	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:ASN:CA	6:B:596:HOH:O	2.44	0.66
1:B:68:ASP:O	1:B:69:GLN:NE2	2.29	0.65
1:A:316:ILE:HD12	1:A:317:ILE:H	1.58	0.65
1:A:254:ARG:HD2	6:A:515:HOH:O	1.96	0.65
1:B:86:LYS:HE2	6:B:505:HOH:O	1.96	0.65
1:B:204:LYS:O	1:B:320:SER:HB3	1.97	0.64
1:A:216:ALA:HA	1:A:219:PHE:CD2	2.32	0.64
1:B:65:VAL:HG13	1:B:69:GLN:CB	2.28	0.64
1:A:139:TYR:CE2	1:A:159:LEU:HD21	2.33	0.64
1:A:58:GLY:CA	6:A:610:HOH:O	2.47	0.63
1:B:173:LEU:HD23	1:B:173:LEU:C	2.24	0.63
1:B:146:GLU:HG2	1:B:147:PRO:HD2	1.79	0.62
1:A:139:TYR:CE1	1:A:159:LEU:HG	2.34	0.62
1:B:142:ASN:ND2	6:B:508:HOH:O	2.32	0.62
2:B:401:F6P:H11	6:B:551:HOH:O	1.99	0.62
1:B:135:ILE:HD12	1:B:284:MET:HE2	1.82	0.61
1:B:135:ILE:HD11	1:B:249:VAL:HG21	1.78	0.61
1:A:65:VAL:O	1:A:65:VAL:HG13	2.01	0.61
1:B:321:PRO:HB2	6:B:606:HOH:O	2.01	0.60
1:A:32:GLN:HA	1:A:32:GLN:OE1	2.00	0.60
1:B:268:LYS:HA	6:B:588:HOH:O	2.00	0.60
1:A:48:VAL:HA	1:A:73:LEU:HD21	1.84	0.60
1:A:9:ASN:N	6:A:609:HOH:O	2.35	0.60
1:A:38:CYS:SG	1:A:193:PHE:CE2	2.95	0.60
1:B:120:LEU:HD11	1:B:132:ILE:CD1	2.32	0.60
1:A:139:TYR:CE2	1:A:159:LEU:CD2	2.85	0.59
1:A:139:TYR:CZ	1:A:159:LEU:CG	2.85	0.59
1:A:139:TYR:OH	1:A:156:GLY:HA2	2.02	0.59
1:B:217:LYS:CG	1:B:217:LYS:O	2.50	0.59
1:B:157:ARG:NE	1:B:304:ASP:OD2	2.34	0.59
1:B:65:VAL:HG11	1:B:69:GLN:O	2.01	0.59
1:A:19:GLU:OE2	1:A:22:ARG:NH2	2.36	0.58
1:B:97:GLU:HB2	1:B:279:TYR:CE1	2.37	0.58
1:A:333:LYS:O	1:A:333:LYS:HG2	2.02	0.58
1:B:194:ILE:O	1:B:196:VAL:HG13	2.03	0.58
1:A:142:ASN:N	1:A:142:ASN:OD1	2.30	0.58
1:B:275:LEU:HD12	1:B:316:ILE:HG21	1.86	0.58
1:A:139:TYR:OH	1:A:159:LEU:HG	2.03	0.58
1:A:316:ILE:HD11	1:A:318:LEU:HB3	1.84	0.58
1:B:277:LEU:HA	1:B:281:CYS:HB2	1.85	0.58
1:A:29:GLU:OE1	1:A:112:LYS:HG2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ILE:HG12	1:B:57:TYR:O	2.04	0.58
1:B:66:THR:O	1:B:67:GLY:C	2.44	0.58
1:B:29:GLU:OE1	1:B:112:LYS:HG2	2.03	0.58
1:A:165:ALA:HA	1:A:172:MET:O	2.03	0.57
1:B:179:ASN:HB2	6:B:533:HOH:O	2.03	0.57
1:B:309:ASP:OD2	1:B:309:ASP:C	2.47	0.57
1:A:224:THR:HB	6:A:603:HOH:O	2.04	0.57
1:A:58:GLY:N	6:A:610:HOH:O	2.37	0.57
1:B:37:LEU:HD12	1:B:85:LEU:HD21	1.87	0.56
1:A:139:TYR:CZ	1:A:159:LEU:HD21	2.39	0.56
1:A:316:ILE:HD12	1:A:316:ILE:C	2.31	0.56
1:B:218:GLU:HG2	1:B:268:LYS:HG3	1.86	0.56
1:B:157:ARG:HE	1:B:304:ASP:CG	2.13	0.56
1:B:13:LEU:HD23	1:B:184:PHE:CZ	2.40	0.56
1:B:53:ILE:HG13	1:B:56:LEU:CD1	2.30	0.55
1:A:204:LYS:C	1:A:321:PRO:HD2	2.32	0.55
1:A:227:ILE:HG22	1:A:231:LYS:HE3	1.88	0.55
1:B:22:ARG:HG2	1:B:22:ARG:HH11	1.72	0.55
1:A:208:ILE:CD1	6:A:552:HOH:O	2.55	0.54
1:A:10:ILE:CD1	1:A:194:ILE:HG13	2.36	0.54
1:A:82:ILE:O	1:A:86:LYS:HG3	2.08	0.54
1:B:157:ARG:HD2	1:B:304:ASP:OD2	2.08	0.54
1:B:251:ASP:OD1	1:B:254:ARG:NH2	2.41	0.54
1:A:21:GLY:HA3	6:A:596:HOH:O	2.08	0.54
1:B:173:LEU:HD23	1:B:174:VAL:N	2.23	0.53
1:B:299:LYS:CE	6:B:602:HOH:O	2.26	0.53
1:A:96:SER:HB2	1:A:117:PHE:CZ	2.43	0.53
1:B:318:LEU:C	1:B:318:LEU:HD12	2.34	0.52
1:A:106:GLU:OE1	1:A:109:LYS:NZ	2.31	0.52
1:B:267:ASN:C	6:B:596:HOH:O	2.51	0.52
1:A:13:LEU:HD21	1:A:38:CYS:HB2	1.92	0.52
1:B:198:ARG:HH11	1:B:198:ARG:CG	2.23	0.52
1:B:218:GLU:CD	1:B:268:LYS:HG3	2.34	0.52
1:B:177:MET:HB3	6:B:521:HOH:O	2.10	0.51
1:B:271:PRO:HD2	6:B:597:HOH:O	2.10	0.51
1:A:332:GLN:O	1:A:335:ALA:N	2.44	0.51
1:B:218:GLU:CG	1:B:268:LYS:HG3	2.41	0.51
1:B:223:ILE:HG13	6:B:591:HOH:O	2.10	0.50
1:B:53:ILE:HG12	1:B:56:LEU:HD12	1.92	0.50
1:B:157:ARG:CD	1:B:304:ASP:OD2	2.59	0.50
1:B:330:ILE:N	1:B:330:ILE:CD1	2.74	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:GLU:HB2	1:A:267:ASN:HB2	1.91	0.50
2:B:401:F6P:C1	6:B:551:HOH:O	2.57	0.50
1:A:218:GLU:OE1	1:A:269:LYS:HB2	2.10	0.50
1:A:184:PHE:HB3	1:A:193:PHE:HB3	1.92	0.50
1:A:141:LYS:HE3	1:A:151:ASP:OD1	2.12	0.49
1:B:261:ILE:HD12	1:B:318:LEU:O	2.12	0.49
1:B:198:ARG:HH11	1:B:198:ARG:HG2	1.77	0.49
1:A:334:HIS:O	1:A:335:ALA:C	2.56	0.49
1:B:216:ALA:HA	1:B:219:PHE:CD2	2.48	0.49
1:B:22:ARG:HG2	1:B:22:ARG:NH1	2.28	0.49
1:B:102:ALA:HB3	1:B:149:GLU:HG2	1.95	0.49
1:A:18:MET:O	1:A:22:ARG:HG3	2.13	0.48
1:B:267:ASN:O	1:B:270:SER:C	2.57	0.48
1:A:37:LEU:HD21	1:A:136:PHE:CE2	2.49	0.48
1:B:82:ILE:HD12	1:B:103:ILE:HD13	1.94	0.48
1:B:221:PRO:HB2	1:B:334:HIS:ND1	2.29	0.48
1:B:63:THR:O	1:B:63:THR:CG2	2.60	0.48
1:B:226:TYR:HB2	1:B:327:LEU:HD13	1.95	0.48
1:A:208:ILE:HD12	6:A:552:HOH:O	2.13	0.48
1:B:270:SER:HB3	1:B:273:GLY:O	2.14	0.48
1:B:261:ILE:HD11	1:B:317:ILE:CG2	2.43	0.47
1:A:78:ASN:OD1	1:A:119:PRO:HG3	2.14	0.47
1:B:40:ALA:O	1:B:44:ILE:HG13	2.13	0.47
1:B:45:SER:O	1:B:49:ARG:HD3	2.14	0.47
1:A:38:CYS:SG	1:A:193:PHE:HD2	2.35	0.47
1:A:49:ARG:O	1:A:50:LYS:HB2	2.14	0.47
1:B:275:LEU:HD12	1:B:316:ILE:CG2	2.45	0.47
1:B:299:LYS:HG2	6:B:604:HOH:O	2.14	0.47
1:A:123:SER:HA	1:A:126:ILE:CG1	2.40	0.46
1:B:299:LYS:HB2	6:B:577:HOH:O	2.15	0.46
1:B:217:LYS:O	1:B:217:LYS:HG2	2.15	0.46
1:A:45:SER:O	1:A:49:ARG:HD3	2.15	0.46
1:B:36:SER:OG	1:B:88:SER:HB3	2.15	0.46
1:A:147:PRO:HA	1:A:151:ASP:OD1	2.16	0.46
6:A:527:HOH:O	1:B:217:LYS:HB3	2.15	0.46
1:B:330:ILE:HD13	1:B:330:ILE:H	1.76	0.46
1:A:276:ARG:NH1	1:A:311:HIS:O	2.45	0.46
1:A:218:GLU:CB	1:A:267:ASN:HB2	2.46	0.46
1:B:206:GLY:HA3	1:B:260:GLY:N	2.31	0.46
1:A:232:PHE:CE1	1:B:217:LYS:HB2	2.51	0.46
1:A:66:THR:HG23	1:A:67:GLY:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ASP:CG	1:A:199:ASP:O	2.58	0.45
1:B:72:LYS:CB	6:B:529:HOH:O	2.54	0.45
1:A:70:VAL:HA	6:A:578:HOH:O	2.16	0.45
1:A:10:ILE:HD11	1:A:194:ILE:HG13	1.99	0.45
1:A:148:SER:OG	1:A:150:LYS:HG2	2.16	0.45
1:B:217:LYS:HG2	1:B:218:GLU:HG3	1.99	0.45
1:A:156:GLY:HA3	1:A:303:LEU:HD22	1.97	0.45
1:B:37:LEU:HD21	1:B:136:PHE:CD2	2.52	0.45
1:B:13:LEU:HB2	1:B:193:PHE:HB2	1.98	0.45
1:B:220:ASP:HB2	1:B:221:PRO:HD2	1.99	0.45
1:A:70:VAL:CG1	1:A:71:LYS:N	2.80	0.44
1:B:10:ILE:HG23	1:B:10:ILE:O	2.15	0.44
1:B:93:VAL:HG22	1:B:104:ILE:CD1	2.47	0.44
1:A:199:ASP:O	1:A:199:ASP:OD1	2.35	0.44
1:A:227:ILE:CG2	1:A:231:LYS:HE3	2.47	0.44
1:A:70:VAL:HG12	1:A:71:LYS:N	2.32	0.44
1:A:230:LYS:HA	1:A:230:LYS:HD3	1.68	0.44
1:A:218:GLU:HB2	1:A:267:ASN:CB	2.47	0.44
1:B:64:ASN:OD1	1:B:64:ASN:O	2.36	0.43
1:B:302:VAL:HA	1:B:305:ILE:HD12	2.00	0.43
1:A:118:ASP:OD1	1:A:118:ASP:C	2.61	0.43
1:A:166:LEU:O	1:A:171:THR:HA	2.18	0.43
1:B:204:LYS:HA	1:B:321:PRO:HD2	2.00	0.43
1:B:267:ASN:C	1:B:267:ASN:OD1	2.60	0.43
1:B:330:ILE:HD11	6:B:518:HOH:O	2.14	0.43
1:A:73:LEU:HD12	1:A:73:LEU:HA	1.85	0.43
1:B:36:SER:HB3	1:B:84:VAL:HG12	2.00	0.43
1:B:123:SER:HA	1:B:126:ILE:CD1	2.49	0.43
1:B:233:PRO:HA	1:B:234:PRO:HD2	1.90	0.43
1:A:187:ASP:HB2	1:B:53:ILE:HG22	1.99	0.43
1:A:333:LYS:C	1:A:335:ALA:H	2.25	0.43
1:A:96:SER:HB2	1:A:117:PHE:CE2	2.54	0.43
1:A:179:ASN:HB2	6:A:576:HOH:O	2.19	0.43
1:A:231:LYS:C	1:A:233:PRO:HD3	2.44	0.43
1:B:330:ILE:HD13	6:B:518:HOH:O	2.14	0.43
1:A:10:ILE:CG1	1:B:57:TYR:O	2.67	0.43
1:B:261:ILE:CD1	1:B:324:VAL:HG22	2.49	0.43
1:B:270:SER:HA	6:B:597:HOH:O	2.19	0.43
1:A:121:ASP:N	1:A:133:GLY:O	2.46	0.42
1:A:157:ARG:HA	1:A:157:ARG:HD3	1.85	0.42
1:B:205:LYS:HE2	1:B:240:TYR:OH	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:ARG:CG	1:B:198:ARG:NH1	2.81	0.42
1:A:199:ASP:OD1	1:A:199:ASP:C	2.62	0.42
1:B:231:LYS:HD2	6:B:515:HOH:O	2.19	0.42
1:B:13:LEU:HD12	1:B:38:CYS:SG	2.60	0.42
1:B:143:SER:O	1:B:144:THR:OG1	2.30	0.42
1:B:13:LEU:HD23	1:B:184:PHE:CE2	2.55	0.42
1:B:22:ARG:NH1	1:B:22:ARG:CG	2.79	0.42
1:B:143:SER:C	1:B:144:THR:HG23	2.44	0.42
1:A:216:ALA:HA	1:A:219:PHE:CE2	2.54	0.41
1:A:88:SER:O	1:A:89:PHE:HB2	2.20	0.41
1:A:267:ASN:OD1	1:A:269:LYS:HB3	2.20	0.41
1:B:242:ALA:O	1:B:243:ARG:HG2	2.20	0.41
1:A:139:TYR:CE1	1:A:159:LEU:N	2.88	0.41
1:A:251:ASP:HA	1:A:254:ARG:NH2	2.35	0.41
1:B:65:VAL:CG1	1:B:66:THR:N	2.83	0.41
1:B:216:ALA:HA	1:B:219:PHE:CE2	2.55	0.41
1:A:220:ASP:OD1	1:A:220:ASP:C	2.63	0.41
1:B:204:LYS:HA	1:B:204:LYS:HD3	1.77	0.41
1:A:225:GLU:HA	6:A:591:HOH:O	2.20	0.41
1:B:18:MET:HE3	1:B:22:ARG:HD3	2.00	0.41
1:A:81:VAL:HG11	1:A:117:PHE:CD2	2.56	0.41
1:A:59:ILE:HB	1:B:10:ILE:HD13	2.01	0.41
1:A:106:GLU:H	1:A:106:GLU:CD	2.28	0.41
1:B:186:LEU:HB2	1:B:193:PHE:CE1	2.56	0.41
1:B:17:VAL:HG11	1:B:34:LEU:HB2	2.02	0.40
1:B:199:ASP:OD1	1:B:199:ASP:O	2.39	0.40
1:B:229:ARG:HG2	1:B:234:PRO:HD2	2.03	0.40
1:B:128:CYS:O	1:B:129:LEU:HB2	2.21	0.40
1:A:318:LEU:C	1:A:318:LEU:CD1	2.86	0.40
4:A:404:AMP:H8	4:A:404:AMP:O5'	2.05	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	A	325/337 (96%)	323 (99%)	2 (1%)	0	100	100
1	B	325/337 (96%)	318 (98%)	7 (2%)	0	100	100
All	All	650/674 (96%)	641 (99%)	9 (1%)	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	A	272/279 (98%)	242 (89%)	30 (11%)	6	13
1	B	272/279 (98%)	251 (92%)	21 (8%)	12	25
All	All	544/558 (98%)	493 (91%)	51 (9%)	8	18

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ILE
1	A	15	ARG
1	A	17	VAL
1	A	22	ARG
1	A	32	GLN
1	A	62	SER
1	A	65	VAL
1	A	69	GLN
1	A	100	LYS
1	A	108	GLU
1	A	112	LYS
1	A	130	VAL
1	A	140	ARG
1	A	141	LYS
1	A	142	ASN

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Mol	Chain	Res	Type
1	A	143	SER
1	A	145	ASP
1	A	150	LYS
1	A	157	ARG
1	A	190	ILE
1	A	197	ASP
1	A	208	ILE
1	A	218	GLU
1	A	267	ASN
1	A	272	LYS
1	A	277	LEU
1	A	278	LEU
1	A	310	ILE
1	A	316	ILE
1	A	333	LYS
1	B	13	LEU
1	B	71	LYS
1	B	82	ILE
1	B	99	ASP
1	B	100	LYS
1	B	140	ARG
1	B	143	SER
1	B	154	GLN
1	B	157	ARG
1	B	160	VAL
1	B	190	ILE
1	B	198	ARG
1	B	217	LYS
1	B	228	GLN
1	B	236	ASN
1	B	269	LYS
1	B	272	LYS
1	B	277	LEU
1	B	278	LEU
1	B	316	ILE
1	B	330	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	179	ASN
1	A	228	GLN

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Mol	Chain	Res	Type
1	A	334	HIS
1	B	142	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](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

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$
4	AMP	A	404	-	25,25,25	1.45	5 (20%)	37,38,38	1.99	12 (32%)
5	PO4	A	405	3	4,4,4	0.95	0	6,6,6	1.78	1 (16%)
2	F6P	A	401	3	15,16,16	1.08	1 (6%)	16,25,25	1.31	2 (12%)
2	F6P	B	401	3	15,16,16	0.94	1 (6%)	16,25,25	1.34	3 (18%)
4	AMP	B	404	-	25,25,25	1.56	4 (16%)	37,38,38	2.08	11 (29%)
5	PO4	B	405	3	4,4,4	0.90	0	6,6,6	1.64	1 (16%)

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	B	401	3	-	1/9/28/28	0/1/1/1
4	AMP	A	404	-	-	0/10/26/26	0/3/3/3
4	AMP	B	404	-	-	0/10/26/26	0/3/3/3
2	F6P	A	401	3	-	2/9/28/28	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	404	AMP	C5-C4	5.11	1.48	1.39
4	A	404	AMP	C5-C4	4.40	1.46	1.39
2	A	401	F6P	O2-C2	3.50	1.46	1.40
4	B	404	AMP	C8-N7	3.45	1.38	1.31
2	B	401	F6P	O2-C2	3.01	1.45	1.40
4	A	404	AMP	C8-N7	2.80	1.37	1.31
4	A	404	AMP	C5-C6	2.46	1.47	1.41
4	A	404	AMP	C5-N7	-2.43	1.34	1.39
4	B	404	AMP	C5-N7	-2.38	1.34	1.39
4	A	404	AMP	C4-N9	-2.34	1.32	1.37
4	B	404	AMP	C5-C6	2.28	1.47	1.41

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	404	AMP	C5-C4-N3	-4.93	119.92	126.72
4	A	404	AMP	C2'-C1'-N9	-4.58	101.92	113.30
4	B	404	AMP	C5-C4-N3	-4.46	120.58	126.72
4	B	404	AMP	C2'-C1'-N9	-4.03	103.28	113.30
4	B	404	AMP	O2P-P-O5'	-3.92	96.44	106.67
4	B	404	AMP	N3-C4-N9	3.90	133.81	127.17
4	A	404	AMP	N3-C2-N1	-3.63	123.09	128.58
4	B	404	AMP	N3-C2-N1	-3.59	123.14	128.58
4	A	404	AMP	N3-C4-N9	3.55	133.21	127.17
4	A	404	AMP	C4-C5-N7	-3.47	106.61	110.58
5	B	405	PO4	O4-P-O2	3.47	118.72	107.91
5	A	405	PO4	O3-P-O2	3.40	118.49	107.91
4	B	404	AMP	C4-N9-C8	3.38	109.29	105.74
4	A	404	AMP	C2-N3-C4	3.26	119.79	111.83
4	B	404	AMP	C2-N1-C6	3.16	123.92	118.73
4	B	404	AMP	C2-N3-C4	3.07	119.32	111.83
2	A	401	F6P	O1-C1-C2	-2.83	105.42	111.67
4	A	404	AMP	C2-N1-C6	2.81	123.34	118.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	404	AMP	C4-N9-C8	2.74	108.62	105.74
2	B	401	F6P	O1-C1-C2	-2.71	105.67	111.67
4	B	404	AMP	N9-C8-N7	-2.55	110.32	113.94
4	A	404	AMP	C5-N7-C8	2.48	107.34	103.45
2	A	401	F6P	O4-C4-C5	-2.39	104.23	111.08
4	A	404	AMP	N9-C8-N7	-2.36	110.59	113.94
4	B	404	AMP	C4-C5-N7	-2.30	107.95	110.58
2	B	401	F6P	O2P-P-O6	-2.28	100.72	106.67
4	B	404	AMP	C5-N7-C8	2.14	106.82	103.45
2	B	401	F6P	O3P-P-O2P	2.14	115.83	107.80
4	A	404	AMP	C6-C5-N7	2.03	136.00	132.09
4	A	404	AMP	O3P-P-O1P	2.01	118.65	110.83

There are no chirality outliers.

All (3) torsion outliers are listed below:

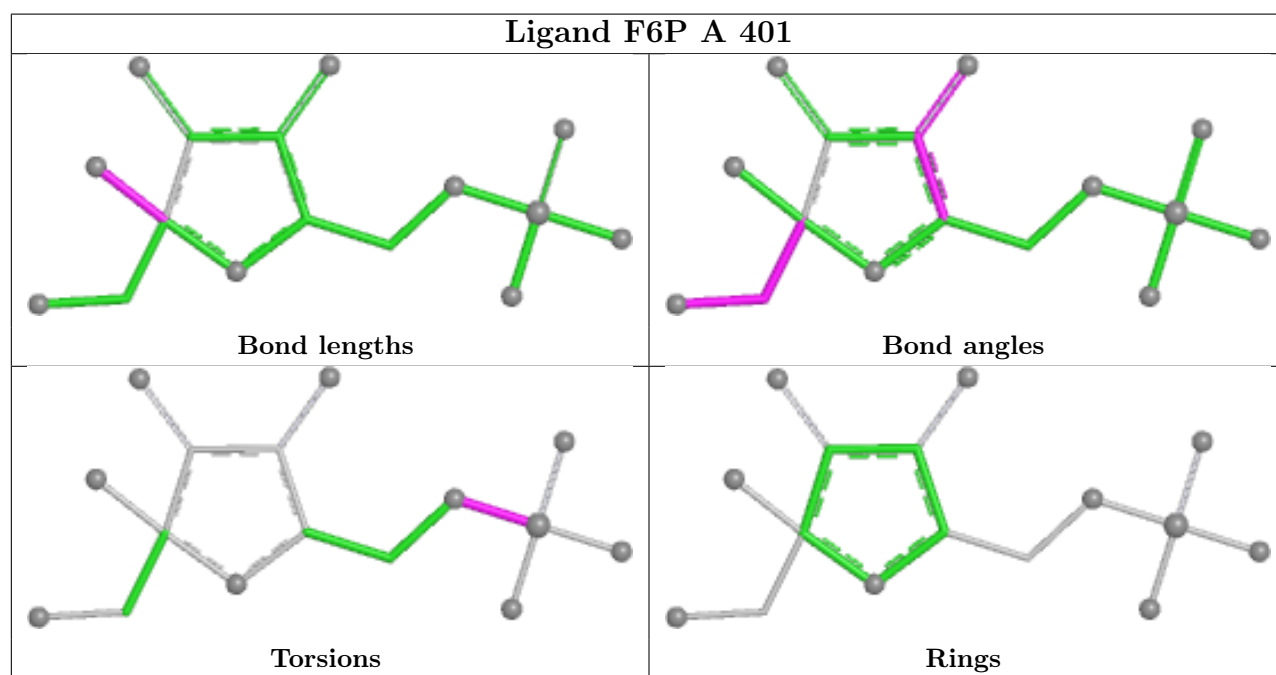
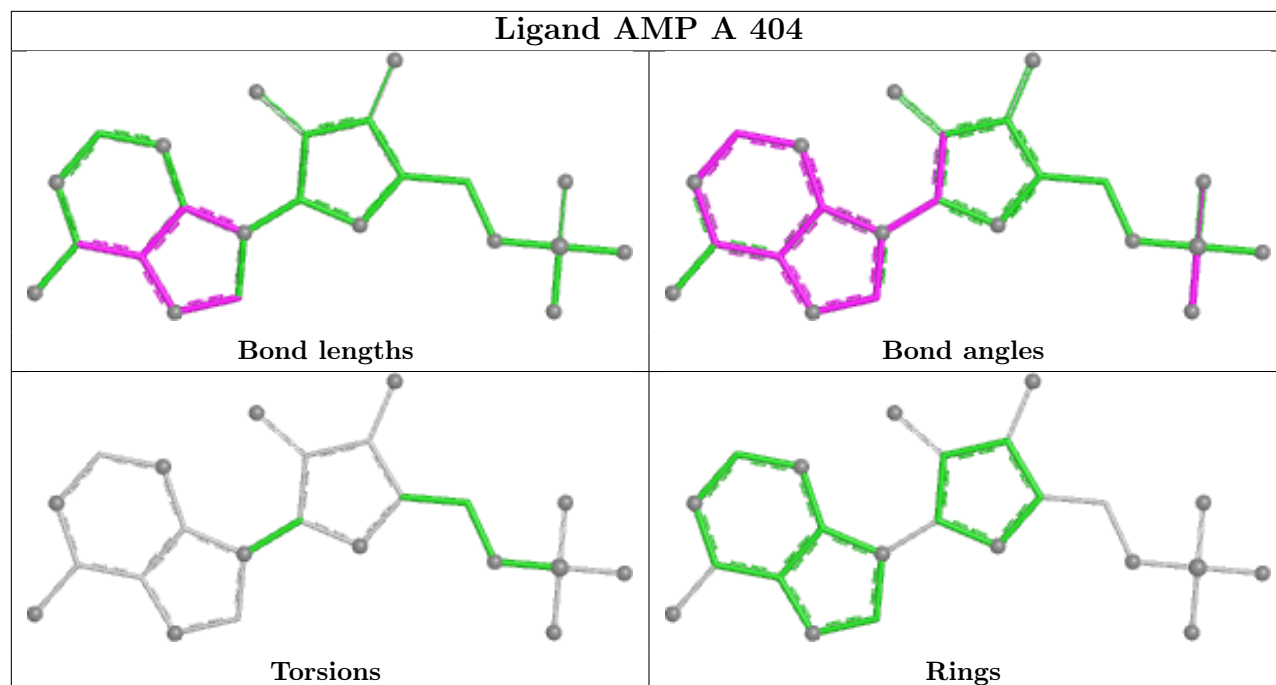
Mol	Chain	Res	Type	Atoms
2	A	401	F6P	C6-O6-P-O1P
2	A	401	F6P	C6-O6-P-O2P
2	B	401	F6P	C4-C5-C6-O6

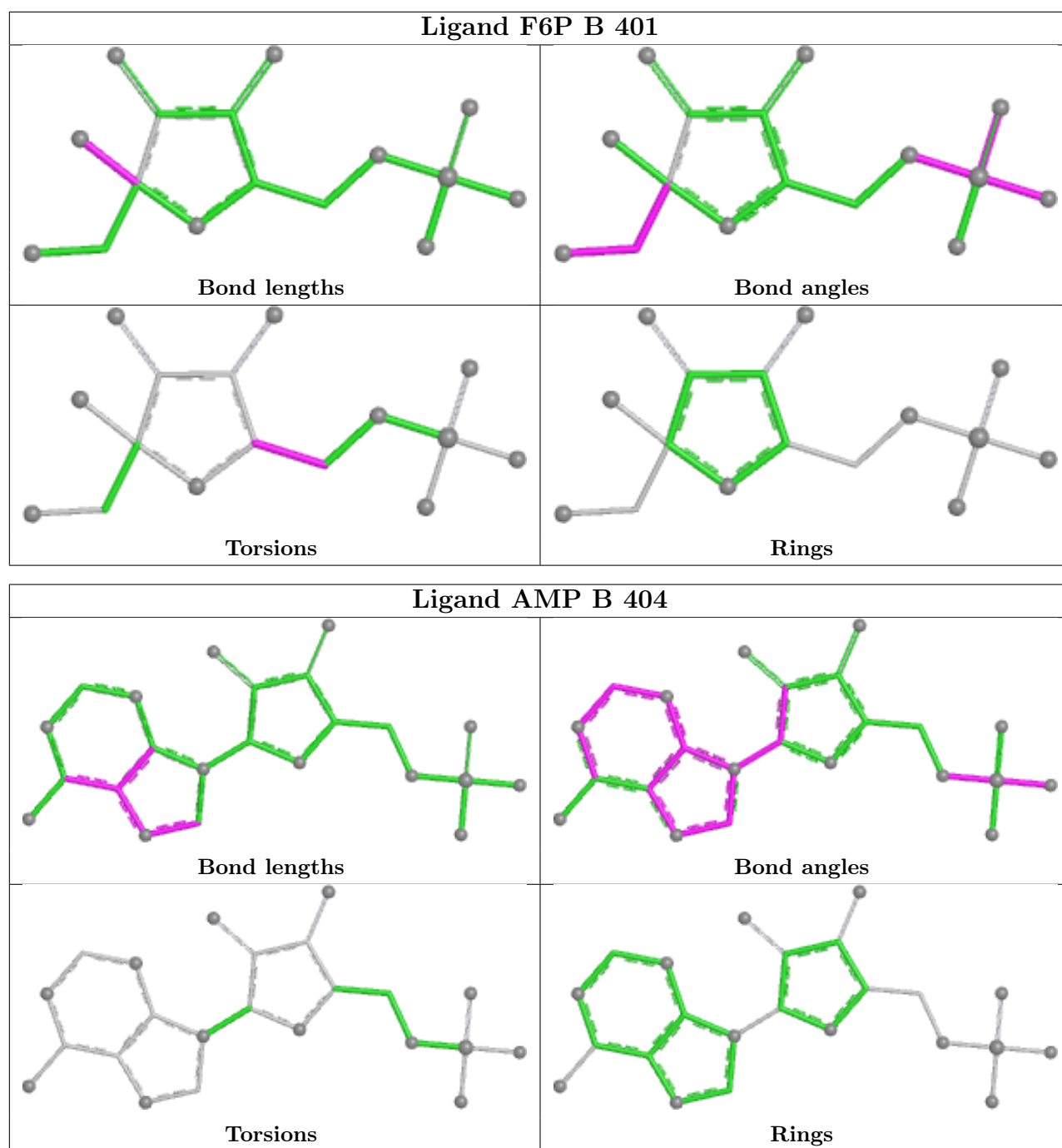
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	404	AMP	1	0
2	B	401	F6P	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.





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	327/337 (97%)	0.00	17 (5%) 33 29	13, 24, 51, 86	0
1	B	327/337 (97%)	-0.06	14 (4%) 40 35	13, 23, 45, 86	0
All	All	654/674 (97%)	-0.03	31 (4%) 36 32	13, 24, 51, 86	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	65	VAL	7.2
1	B	66	THR	7.0
1	A	63	THR	6.3
1	A	64	ASN	6.3
1	A	65	VAL	5.8
1	A	70	VAL	4.9
1	A	67	GLY	4.8
1	A	62	SER	4.6
1	B	142	ASN	4.4
1	B	67	GLY	4.4
1	B	70	VAL	4.3
1	A	66	THR	4.3
1	B	64	ASN	4.0
1	B	199	ASP	3.7
1	B	62	SER	3.6
1	B	69	GLN	3.5
1	B	58	GLY	3.4
1	A	69	GLN	3.2
1	B	68	ASP	3.2
1	A	142	ASN	3.2
1	B	63	THR	3.0
1	A	139	TYR	3.0
1	A	335	ALA	2.9
1	A	38	CYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	154	GLN	2.4
1	A	56	LEU	2.4
1	A	143	SER	2.4
1	A	68	ASP	2.4
1	A	268	LYS	2.4
1	B	56	LEU	2.3
1	B	271	PRO	2.3

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](#)

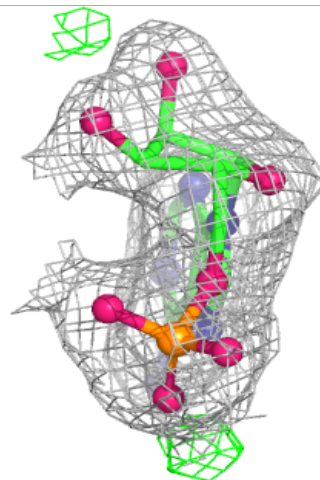
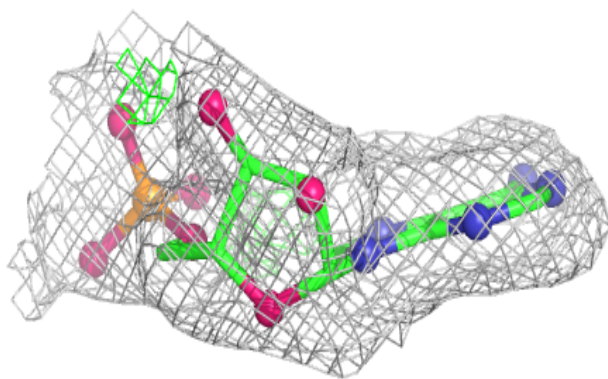
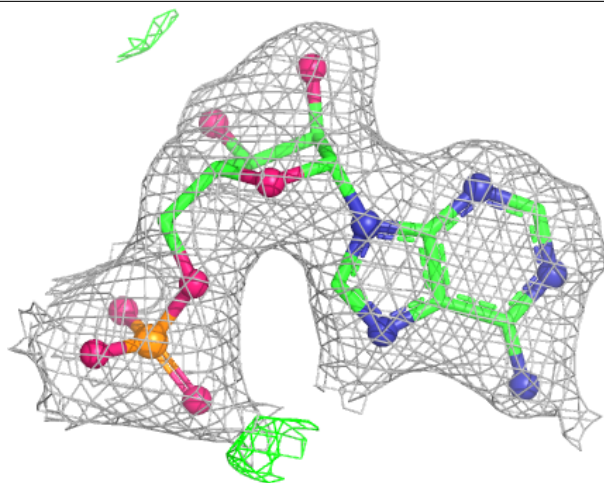
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	PO4	B	405	5/5	0.92	0.11	39,40,42,43	0
5	PO4	A	405	5/5	0.94	0.10	44,45,48,49	0
4	AMP	A	404	23/23	0.96	0.06	18,25,28,29	0
4	AMP	B	404	23/23	0.96	0.07	20,25,29,30	0
2	F6P	A	401	16/16	0.97	0.06	19,23,25,26	0
2	F6P	B	401	16/16	0.97	0.07	11,21,23,27	0
3	MG	A	403	1/1	0.97	0.11	19,19,19,19	0
3	MG	B	402	1/1	0.97	0.10	5,5,5,5	0
3	MG	A	402	1/1	0.98	0.09	10,10,10,10	0
3	MG	B	403	1/1	0.98	0.12	9,9,9,9	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.

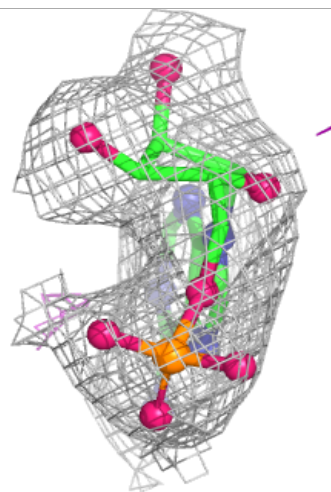
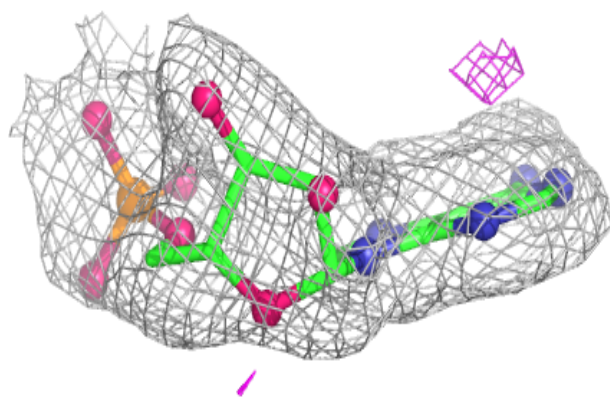
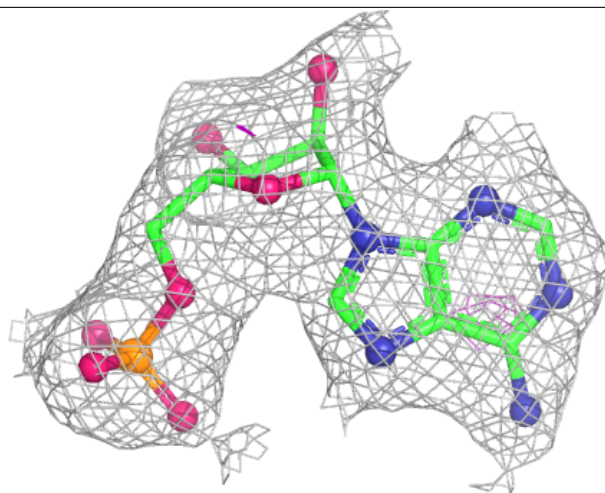
Electron density around AMP A 404:

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



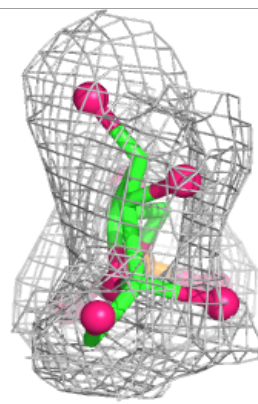
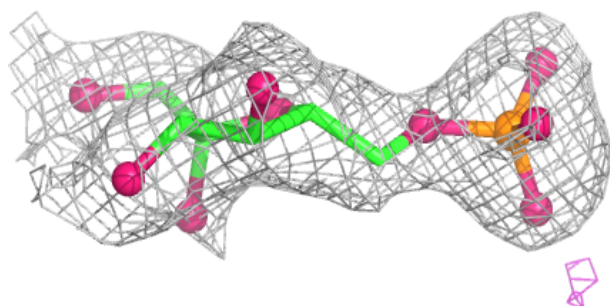
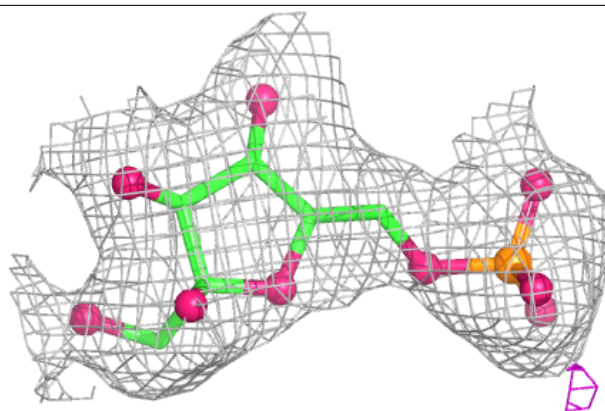
Electron density around AMP B 404:

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

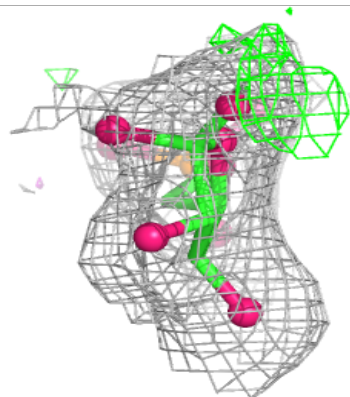
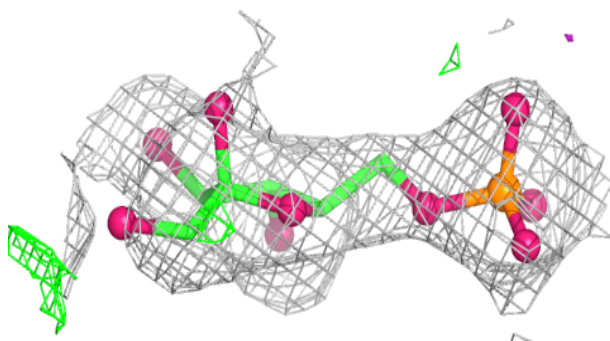
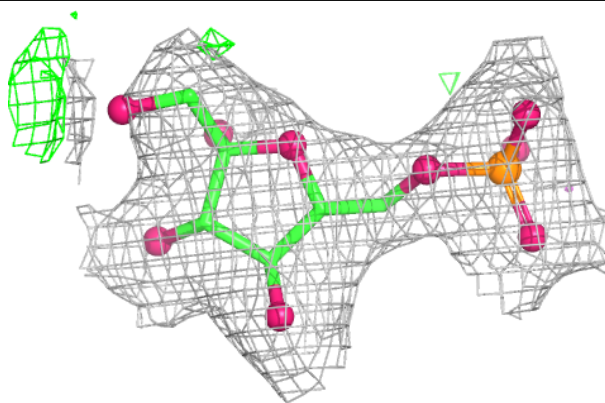


Electron density around F6P A 401:

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

**Electron density around F6P B 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.