



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:22 PM GMT

PDB ID : 1FBP
Title : CRYSTAL STRUCTURE OF FRUCTOSE-1,6-BISPHOSPHATASE COM-
PLEXED WITH FRUCTOSE 6-PHOSPHATE, AMP, AND MAGNESIUM
Authors : Ke, H.; Zhang, Y.; Lipscomb, W.N.
Deposited on : 1990-05-31
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

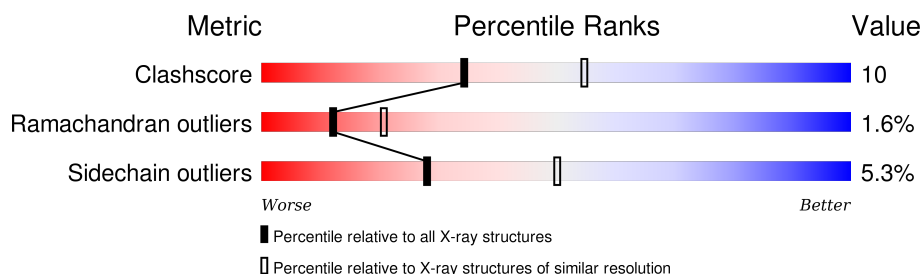
1 Overall quality at a glance

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))
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	335	
1	B	335	

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	336	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6016 atoms, of which 1076 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.

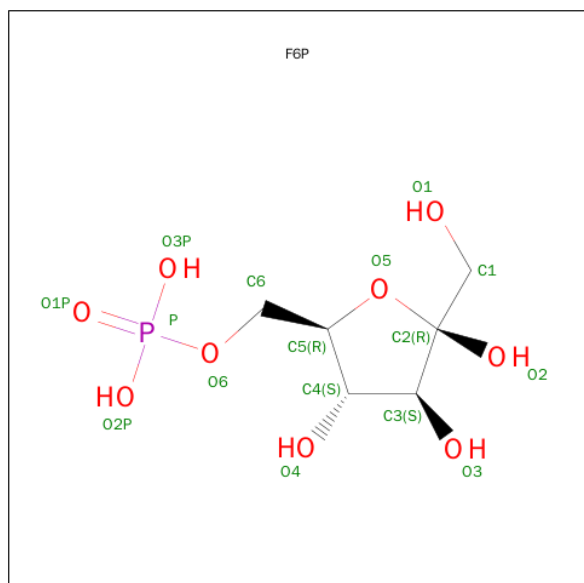
- Molecule 1 is a protein called FRUCTOSE 1,6-BISPHOSPHATASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	317	Total	C	H	N	O	S	0	0	1
			2968	1546	538	409	460	15			
1	B	317	Total	C	H	N	O	S	0	0	1
			2968	1546	538	409	460	15			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	GLN	GLU	CONFLICT	UNP P00636
A	96	THR	SER	CONFLICT	UNP P00636
A	199	ASN	ASP	CONFLICT	UNP P00636
B	20	GLN	GLU	CONFLICT	UNP P00636
B	96	THR	SER	CONFLICT	UNP P00636
B	199	ASN	ASP	CONFLICT	UNP P00636

- Molecule 2 is SUGAR (FRUCTOSE-6-PHOSPHATE) (three-letter code: F6P) (formula: $C_6H_{13}O_9P$).

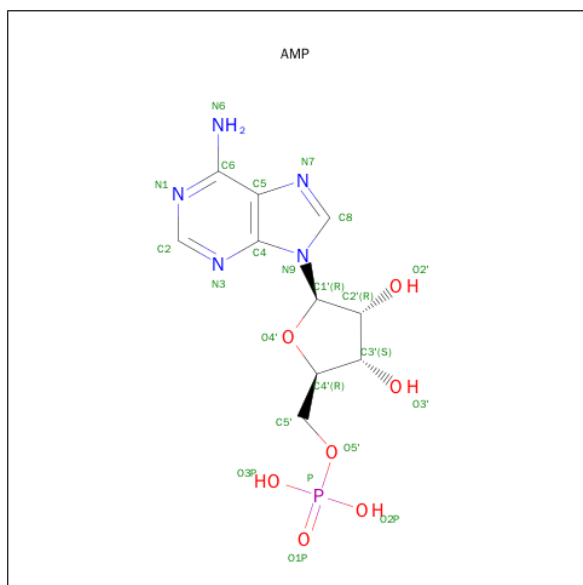


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			16	6	9	1		
2	B	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ADENOSINE MONOPHOSPHATE (three-letter code: 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		

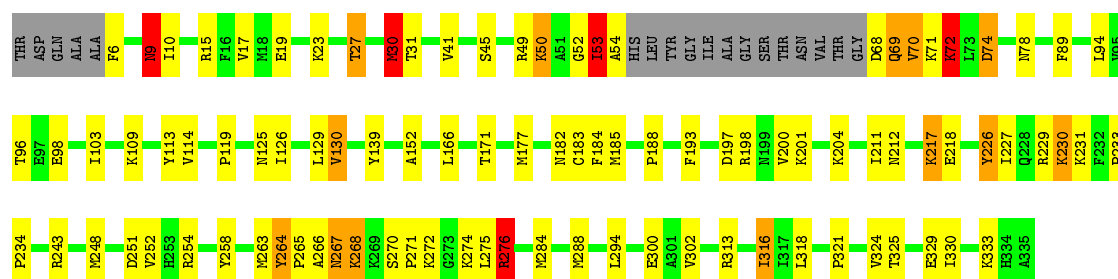
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

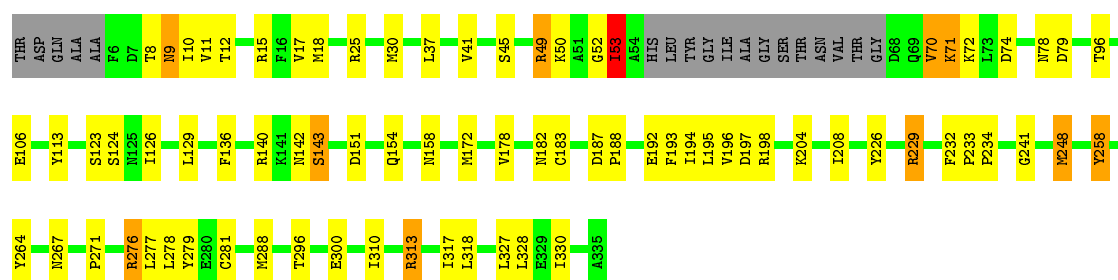
• Molecule 1: FRUCTOSE 1,6-BISPHOSPHATASE

Chain A: 



• Molecule 1: FRUCTOSE 1,6-BISPHOSPHATASE

Chain B: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	61.60Å 166.60Å 80.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.215 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6016	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AMP, MG, F6P

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	0.74	0/2470	1.51	22/3339 (0.7%)
1	B	0.75	0/2470	1.49	18/3339 (0.5%)
All	All	0.75	0/4940	1.50	40/6678 (0.6%)

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	5
1	B	0	3
All	All	0	8

There are no bond length outliers.

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	254	ARG	NE-CZ-NH2	-18.23	111.19	120.30
1	B	49	ARG	NE-CZ-NH2	-14.80	112.90	120.30
1	A	254	ARG	NE-CZ-NH1	10.73	125.66	120.30
1	B	248	MET	CG-SD-CE	-10.48	83.43	100.20
1	B	15	ARG	NE-CZ-NH2	-9.39	115.61	120.30
1	A	276	ARG	NE-CZ-NH1	9.03	124.81	120.30
1	B	276	ARG	NE-CZ-NH1	8.53	124.56	120.30
1	A	6	PHE	N-CA-C	-8.26	88.70	111.00
1	A	15	ARG	NE-CZ-NH2	-8.10	116.25	120.30
1	B	49	ARG	NE-CZ-NH1	7.94	124.27	120.30
1	A	50	LYS	N-CA-C	-7.76	90.05	111.00
1	B	229	ARG	NE-CZ-NH2	-7.37	116.62	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	313	ARG	NE-CZ-NH2	-7.07	116.76	120.30
1	A	251	ASP	CB-CG-OD2	6.81	124.43	118.30
1	B	18	MET	CG-SD-CE	-6.38	89.98	100.20
1	A	264	TYR	CB-CG-CD2	-6.32	117.21	121.00
1	A	53	ILE	CB-CA-C	-6.20	99.20	111.60
1	A	185	MET	CG-SD-CE	-6.14	90.38	100.20
1	A	30	MET	CA-CB-CG	-6.13	102.87	113.30
1	A	313	ARG	NE-CZ-NH2	-6.02	117.29	120.30
1	B	17	VAL	CA-CB-CG2	-5.77	102.25	110.90
1	A	70	VAL	N-CA-C	-5.75	95.46	111.00
1	A	27	THR	CA-CB-CG2	5.73	120.43	112.40
1	B	106	GLU	OE1-CD-OE2	-5.63	116.54	123.30
1	B	313	ARG	NE-CZ-NH1	5.63	123.11	120.30
1	B	187	ASP	CB-CG-OD1	5.55	123.30	118.30
1	A	9	ASN	CA-C-N	-5.46	105.20	117.20
1	B	53	ILE	CB-CA-C	-5.39	100.81	111.60
1	A	272	LYS	N-CA-C	-5.38	96.48	111.00
1	A	268	LYS	CB-CG-CD	-5.33	97.74	111.60
1	B	276	ARG	CA-CB-CG	-5.27	101.80	113.40
1	B	258	TYR	N-CA-CB	-5.24	101.17	110.60
1	A	177	MET	CG-SD-CE	-5.23	91.83	100.20
1	A	41	VAL	CA-CB-CG2	-5.15	103.18	110.90
1	A	30	MET	CG-SD-CE	-5.11	92.03	100.20
1	B	143	SER	N-CA-CB	-5.05	102.92	110.50
1	A	53	ILE	CA-CB-CG2	-5.04	100.82	110.90
1	B	151	ASP	CB-CG-OD2	5.04	122.83	118.30
1	A	130	VAL	CG1-CB-CG2	-5.03	102.85	110.90
1	B	30	MET	CG-SD-CE	-5.03	92.16	100.20

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	113	TYR	Sidechain
1	A	139	TYR	Sidechain
1	A	226	TYR	Sidechain
1	A	258	TYR	Sidechain
1	A	264	TYR	Sidechain
1	B	226	TYR	Sidechain
1	B	258	TYR	Sidechain
1	B	264	TYR	Sidechain

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	2430	538	2486	64	0
1	B	2430	538	2486	37	0
2	A	16	0	11	13	0
2	B	16	0	11	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	23	0	12	4	0
4	B	23	0	12	3	0
All	All	4940	1076	5018	98	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 10.

All (98) 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:27:THR:HB	4:A:337:AMP:O3P	1.34	1.26
1:A:274:LYS:NZ	2:A:336:F6P:H61	1.54	1.20
1:A:274:LYS:HZ3	2:A:336:F6P:H61	0.97	1.09
1:A:275:LEU:HD21	2:A:336:F6P:O2	1.53	1.09
1:A:274:LYS:NZ	2:A:336:F6P:H62	1.75	0.99
1:A:274:LYS:HZ3	2:A:336:F6P:C6	1.67	0.98
1:A:274:LYS:HZ2	2:A:336:F6P:C6	1.83	0.91
1:A:27:THR:CB	4:A:337:AMP:O3P	2.22	0.88
1:A:276:ARG:HH11	1:A:276:ARG:HG3	1.40	0.86
1:A:218:GLU:HB3	1:A:267:ASN:HB2	1.61	0.82
1:A:243:ARG:NH2	2:B:336:F6P:O1P	2.12	0.79
1:B:9:ASN:HB3	1:B:194:ILE:HG12	1.67	0.77
1:B:78:ASN:OD1	1:B:96:THR:HG21	1.85	0.77
1:A:125:ASN:HB3	1:A:130:VAL:HB	1.68	0.75
1:A:275:LEU:HD21	2:A:336:F6P:HO2	1.47	0.74
1:A:217:LYS:NZ	1:A:217:LYS:HB2	2.07	0.70
1:A:204:LYS:HA	1:A:321:PRO:HD2	1.74	0.70
1:A:248:MET:HB2	2:A:336:F6P:H12	1.73	0.70
1:A:78:ASN:OD1	1:A:96:THR:HG21	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:VAL:HA	1:B:126:ILE:HD12	1.75	0.68
1:A:274:LYS:HZ2	2:A:336:F6P:H62	1.48	0.67
1:B:229:ARG:NH2	1:B:330:ILE:HD11	2.10	0.67
1:B:195:LEU:HD21	1:B:198:ARG:HG2	1.78	0.65
1:B:277:LEU:HA	1:B:281:CYS:HB2	1.78	0.64
1:B:9:ASN:OD1	1:B:12:THR:HG23	1.98	0.64
1:A:316:ILE:HD11	1:A:318:LEU:HD23	1.82	0.61
1:A:30:MET:HB3	4:A:337:AMP:C2	2.35	0.60
1:A:212:ASN:ND2	2:A:336:F6P:O2P	2.33	0.60
1:A:218:GLU:HG2	1:A:268:LYS:HB2	1.83	0.59
1:B:182:ASN:HD22	1:B:198:ARG:HA	1.66	0.59
1:A:217:LYS:HZ2	1:A:217:LYS:HB2	1.70	0.57
1:A:302:VAL:HG21	1:A:316:ILE:HD12	1.85	0.57
1:A:96:THR:HG22	1:A:98:GLU:H	1.70	0.57
1:A:226:TYR:CZ	1:A:230:LYS:HD3	2.40	0.55
1:B:296:THR:HG21	1:B:328:LEU:HD21	1.88	0.55
1:A:50:LYS:HG2	1:B:188:PRO:HD2	1.87	0.55
1:A:248:MET:N	2:A:336:F6P:O1	2.40	0.54
1:A:17:VAL:HG12	1:A:31:THR:HG23	1.92	0.52
1:A:188:PRO:HD2	1:B:50:LYS:HG2	1.91	0.51
1:A:50:LYS:NZ	1:A:54:ALA:N	2.58	0.51
4:B:337:AMP:O4'	4:B:337:AMP:N3	2.43	0.51
4:A:337:AMP:N3	4:A:337:AMP:O4'	2.44	0.51
1:A:276:ARG:NH1	1:A:276:ARG:HG3	2.12	0.51
1:B:182:ASN:ND2	1:B:198:ARG:HA	2.26	0.50
1:A:227:ILE:HG22	1:A:231:LYS:HE2	1.93	0.49
1:A:252:VAL:HG11	1:A:284:MET:SD	2.53	0.49
1:A:288:MET:HG3	1:A:318:LEU:HB2	1.95	0.49
1:A:182:ASN:HD22	1:A:198:ARG:HA	1.79	0.47
1:A:183:CYS:HB2	1:A:197:ASP:HB3	1.96	0.47
1:A:218:GLU:CG	1:A:268:LYS:HB2	2.43	0.47
1:A:274:LYS:HZ1	2:A:336:F6P:H62	1.74	0.47
1:B:140:ARG:HH12	4:B:337:AMP:H3'	1.79	0.47
1:A:266:ALA:CB	1:A:271:PRO:HA	2.44	0.47
1:A:71:LYS:HA	1:A:74:ASP:OD1	2.15	0.47
1:A:229:ARG:HE	1:A:330:ILE:HD11	1.81	0.46
1:A:333:LYS:HG2	1:A:333:LYS:O	2.15	0.46
1:B:278:LEU:HD12	1:B:310:ILE:HA	1.98	0.46
1:A:217:LYS:HB3	1:B:232:PHE:CE2	2.51	0.46
1:A:325:THR:O	1:A:329:GLU:HG3	2.15	0.46
1:A:129:LEU:HD11	1:B:172:MET:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:LEU:HD12	1:A:324:VAL:HB	1.97	0.46
1:A:89:PHE:CD2	1:A:109:LYS:HA	2.51	0.46
1:B:194:ILE:O	1:B:196:VAL:HG13	2.16	0.46
1:A:266:ALA:HB1	1:A:271:PRO:HA	1.97	0.45
1:B:288:MET:HG3	1:B:318:LEU:HD13	1.97	0.45
1:B:126:ILE:O	1:B:129:LEU:HD23	2.16	0.45
1:B:71:LYS:HA	1:B:74:ASP:OD1	2.16	0.45
1:A:69:GLN:OE1	1:A:72:LYS:NZ	2.48	0.44
1:A:114:VAL:HG11	1:A:152:ALA:HA	1.99	0.44
1:B:113:TYR:OH	4:B:337:AMP:H5'1	2.17	0.44
1:B:267:ASN:O	1:B:271:PRO:HA	2.17	0.43
1:B:12:THR:HG22	1:B:192:GLU:HG3	2.00	0.43
1:B:229:ARG:HH22	1:B:330:ILE:HD11	1.82	0.43
1:A:96:THR:HG23	1:A:119:PRO:HD3	2.01	0.43
1:B:317:ILE:HG21	1:B:327:LEU:HD23	2.01	0.43
1:B:248:MET:N	2:B:336:F6P:O3	2.43	0.43
1:B:233:PRO:HA	1:B:234:PRO:HD3	1.89	0.43
1:A:45:SER:O	1:A:49:ARG:HD3	2.19	0.43
1:A:183:CYS:SG	1:A:200:VAL:HG21	2.59	0.42
1:A:211:ILE:HD12	1:A:263:MET:HB2	2.01	0.42
1:A:70:VAL:HA	1:A:126:ILE:HD12	2.01	0.42
1:B:37:LEU:HD21	1:B:136:PHE:CD2	2.54	0.42
1:A:19:GLU:HB3	1:A:23:LYS:NZ	2.35	0.42
1:A:71:LYS:O	1:A:72:LYS:HB2	2.19	0.42
1:B:52:GLY:O	1:B:53:ILE:HG13	2.20	0.42
1:B:154:GLN:HE21	1:B:158:ASN:HD22	1.67	0.41
1:A:166:LEU:O	1:A:171:THR:HA	2.20	0.41
1:A:50:LYS:CG	1:B:188:PRO:HD2	2.51	0.41
1:A:94:LEU:HB2	1:A:103:ILE:HB	2.01	0.41
1:B:208:ILE:HA	1:B:241:GLY:O	2.21	0.41
1:B:276:ARG:HD2	1:B:279:TYR:OH	2.21	0.41
1:B:204:LYS:HB2	1:B:204:LYS:HE3	1.91	0.41
1:A:184:PHE:HB3	1:A:193:PHE:HB3	2.03	0.41
1:A:233:PRO:HA	1:A:234:PRO:HD2	1.83	0.40
1:A:248:MET:CA	2:A:336:F6P:O1	2.70	0.40
1:B:183:CYS:HB2	1:B:197:ASP:HB3	2.02	0.40
1:B:10:ILE:HG23	1:B:11:VAL:HG23	2.02	0.40
1:B:41:VAL:HG12	1:B:193:PHE:HZ	1.87	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	313/335 (93%)	290 (93%)	16 (5%)	7 (2%)	8	13
1	B	313/335 (93%)	286 (91%)	24 (8%)	3 (1%)	19	34
All	All	626/670 (93%)	576 (92%)	40 (6%)	10 (2%)	12	21

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	10	ILE
1	A	53	ILE
1	B	70	VAL
1	A	72	LYS
1	B	9	ASN
1	B	178	VAL
1	A	52	GLY
1	A	267	ASN
1	A	270	SER

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	266/278 (96%)	252 (95%)	14 (5%)	28	50
1	B	266/278 (96%)	252 (95%)	14 (5%)	28	50
All	All	532/556 (96%)	504 (95%)	28 (5%)	28	50

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	30	MET
1	A	53	ILE
1	A	68	ASP
1	A	69	GLN
1	A	72	LYS
1	A	74	ASP
1	A	201	LYS
1	A	217	LYS
1	A	230	LYS
1	A	265	PRO
1	A	276	ARG
1	A	300	GLU
1	A	316	ILE
1	B	8	THR
1	B	25	ARG
1	B	45	SER
1	B	49	ARG
1	B	53	ILE
1	B	71	LYS
1	B	72	LYS
1	B	79	ASP
1	B	123	SER
1	B	124	SER
1	B	142	ASN
1	B	143	SER
1	B	300	GLU
1	B	313	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	142	ASN
1	A	154	GLN
1	A	182	ASN
1	B	154	GLN
1	B	182	ASN
1	B	282	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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$
2	F6P	A	336	1	15,16,16	1.03	0	16,25,25	0.84	0
4	AMP	A	337	-	20,25,25	2.51	6 (30%)	22,38,38	2.75	10 (45%)
2	F6P	B	336	-	15,16,16	1.03	0	16,25,25	0.84	0
4	AMP	B	337	1	20,25,25	2.51	6 (30%)	22,38,38	2.75	10 (45%)

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	336	1	-	0/9/28/28	0/1/1/1
4	AMP	A	337	-	-	0/6/26/26	0/3/3/3
2	F6P	B	336	-	-	0/9/28/28	0/1/1/1
4	AMP	B	337	1	-	0/6/26/26	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	337	AMP	C3'-C4'	-5.50	1.38	1.53
4	B	337	AMP	C3'-C4'	-5.49	1.38	1.53
4	A	337	AMP	P-O5'	-4.06	1.46	1.60
4	B	337	AMP	P-O5'	-4.06	1.46	1.60
4	A	337	AMP	O4'-C1'	3.41	1.45	1.41
4	B	337	AMP	O4'-C1'	3.41	1.45	1.41
4	B	337	AMP	O5'-C5'	3.58	1.59	1.44
4	A	337	AMP	O5'-C5'	3.59	1.59	1.44
4	A	337	AMP	O4'-C4'	4.27	1.54	1.45
4	B	337	AMP	O4'-C4'	4.27	1.54	1.45
4	B	337	AMP	O3'-C3'	4.66	1.54	1.43
4	A	337	AMP	O3'-C3'	4.71	1.54	1.43

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	337	AMP	C2'-C1'-N9	-4.62	107.24	114.29
4	A	337	AMP	C2'-C1'-N9	-4.58	107.29	114.29
4	A	337	AMP	N3-C2-N1	-4.15	125.72	128.89
4	B	337	AMP	N3-C2-N1	-4.14	125.72	128.89
4	B	337	AMP	O4'-C1'-N9	-3.83	100.08	108.10
4	A	337	AMP	O4'-C1'-N9	-3.81	100.11	108.10
4	A	337	AMP	O4'-C4'-C3'	-3.10	98.91	105.15
4	B	337	AMP	O3P-P-O1P	-3.08	100.66	110.58
4	A	337	AMP	O3P-P-O1P	-3.08	100.66	110.58
4	B	337	AMP	O4'-C4'-C3'	-3.08	98.95	105.15
4	B	337	AMP	C2-N1-C6	2.28	122.85	118.77
4	A	337	AMP	C2-N1-C6	2.29	122.85	118.77
4	A	337	AMP	C4-C5-N7	3.00	112.24	109.48
4	B	337	AMP	C4-C5-N7	3.03	112.27	109.48
4	B	337	AMP	O3P-P-O2P	3.05	118.99	107.38
4	A	337	AMP	O3P-P-O2P	3.06	119.03	107.38
4	A	337	AMP	O3P-P-O5'	3.35	116.21	106.56
4	B	337	AMP	O3P-P-O5'	3.36	116.24	106.56
4	B	337	AMP	C2'-C3'-C4'	7.04	117.08	102.61
4	A	337	AMP	C2'-C3'-C4'	7.06	117.12	102.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	336	F6P	13	0
4	A	337	AMP	4	0
2	B	336	F6P	2	0
4	B	337	AMP	3	0

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 ⓘ

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.