



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 08:33 PM UTC

PDB ID : 1FJ9 / pdb_00001fj9
Title : FRUCTOSE-1,6-BISPHOSPHATASE (MUTANT Y57W) PROD-
UCTS/ZN/AMP COMPLEX (T-STATE)
Authors : Iancu, C.V.; Choe, J.Y.; Honzatko, R.B.
Deposited on : 2000-08-07
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

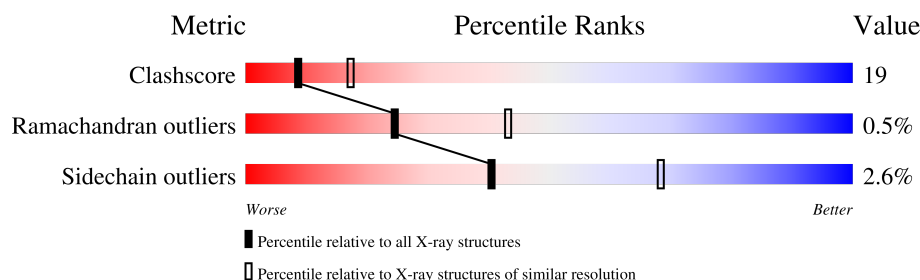
1 Overall quality at a glance

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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	337	
1	B	337	

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
4	PO4	B	412	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5190 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.

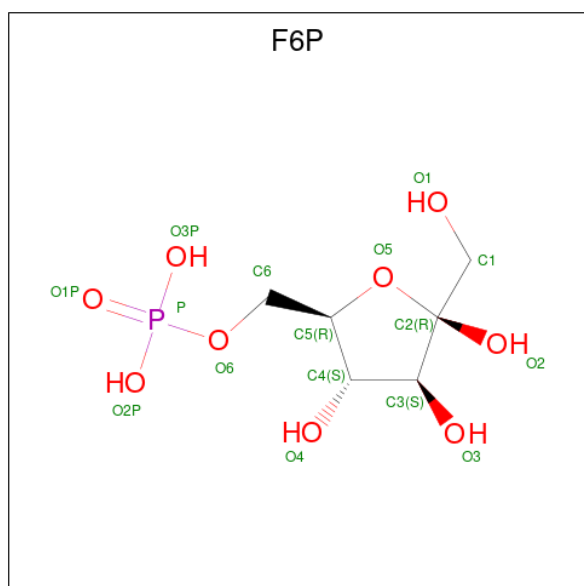
- Molecule 1 is a protein called FRUCTOSE-1,6-BISPHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2441	1556	411	459	15			
1	B	319	Total	C	N	O	S	0	0	0
			2441	1556	411	459	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	57	TRP	TYR	engineered mutation	UNP P00636
B	57	TRP	TYR	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		

Continued on next page...

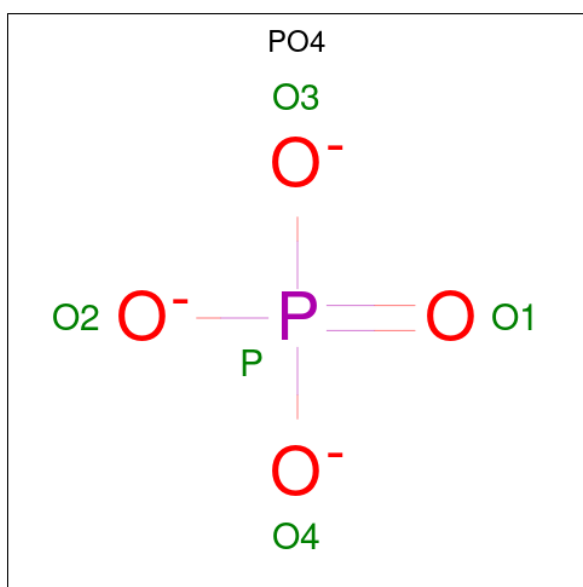
Continued from previous page...

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

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

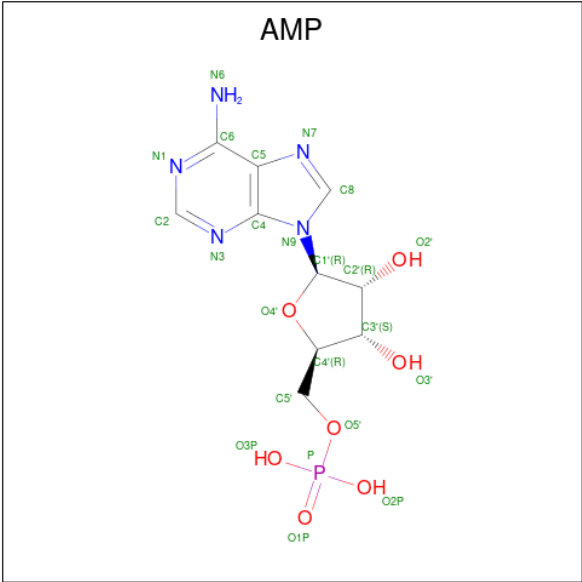
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 5 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: C₁₀H₁₄N₅O₇P).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	96	Total	O	0	0
			96	96		
6	B	120	Total	O	0	0
			120	120		

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	59.82Å 165.48Å 79.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 2.50	Depositor
% Data completeness (in resolution range)	87.9 (5.00-2.50)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.194 , 0.258	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5190	wwPDB-VP
Average B, all atoms (Å ²)	18.0	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: PO4, F6P, AMP, ZN

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	0.40	0/2483	0.96	9/3357 (0.3%)
1	B	0.43	0/2483	0.94	8/3357 (0.2%)
All	All	0.42	0/4966	0.95	17/6714 (0.3%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	278	LEU	N-CA-C	8.80	120.49	111.07
1	A	270	SER	CA-C-N	7.90	128.32	119.32
1	A	270	SER	C-N-CA	7.90	128.32	119.32
1	A	113	TYR	N-CA-C	7.03	121.50	110.32
1	B	298	GLY	N-CA-C	-6.96	106.19	114.48
1	B	113	TYR	N-CA-C	5.82	119.58	110.32
1	A	277	LEU	N-CA-C	5.79	117.28	110.97
1	A	263	MET	N-CA-C	5.67	118.72	108.69
1	B	278	LEU	N-CA-C	5.51	116.96	111.07
1	A	117	PHE	N-CA-C	5.48	117.41	109.07
1	B	277	LEU	N-CA-C	5.43	117.00	111.14
1	B	110	ARG	N-CA-C	5.29	118.11	110.28
1	A	111	GLY	N-CA-C	-5.26	103.33	112.06
1	B	297	THR	N-CA-C	-5.24	106.74	113.23
1	A	297	THR	N-CA-C	-5.04	106.97	113.23
1	B	270	SER	CA-C-N	5.01	124.62	119.56
1	B	270	SER	C-N-CA	5.01	124.62	119.56

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	2441	0	2499	86	0
1	B	2441	0	2499	105	0
2	A	16	0	11	0	0
2	B	16	0	11	3	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	2	0
5	A	23	0	12	0	0
5	B	23	0	12	0	0
6	A	96	0	0	3	0
6	B	120	0	0	3	0
All	All	5190	0	5044	188	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 19.

All (188) 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:198:ARG:HE	1:A:199:ASP:HB2	1.33	0.93
1:A:261:ILE:HD13	1:A:262:PHE:N	1.92	0.84
1:B:72:LYS:H	1:B:72:LYS:HD2	1.44	0.83
1:A:182:ASN:HD22	1:A:198:ARG:HA	1.41	0.83
1:A:182:ASN:HD21	1:A:198:ARG:HD2	1.42	0.82
1:A:308:THR:H	1:A:312:GLN:HE22	1.29	0.80
1:A:182:ASN:ND2	1:A:198:ARG:HD2	1.99	0.77
1:A:55:HIS:HA	1:A:59:ILE:HG22	1.66	0.77
1:A:223:ILE:O	1:A:227:ILE:HG12	1.85	0.76
1:B:288:MET:HE3	1:B:318:LEU:HD13	1.68	0.74
1:B:267:ASN:ND2	1:B:270:SER:H	1.86	0.74
1:A:182:ASN:ND2	1:A:199:ASP:H	1.85	0.74
1:A:308:THR:H	1:A:312:GLN:NE2	1.86	0.73
1:B:230:LYS:HA	1:B:230:LYS:HE2	1.70	0.73
1:B:276:ARG:HH11	1:B:313:ARG:HH12	1.33	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:LYS:HD3	1:B:269:LYS:C	2.16	0.70
1:B:276:ARG:HH11	1:B:313:ARG:NH1	1.90	0.69
2:B:410:F6P:H11	4:B:412:PO4:P	2.33	0.68
1:A:227:ILE:HG22	1:A:231:LYS:HE3	1.72	0.68
1:B:182:ASN:HD22	1:B:198:ARG:HA	1.58	0.68
1:A:148:SER:OG	1:A:150:LYS:HG3	1.94	0.68
1:A:182:ASN:ND2	1:A:198:ARG:HA	2.09	0.67
1:A:261:ILE:HD13	1:A:262:PHE:H	1.60	0.67
1:A:78:ASN:O	1:A:82:ILE:HG12	1.95	0.66
1:B:276:ARG:NH1	1:B:313:ARG:HH12	1.94	0.66
1:A:182:ASN:HD21	1:A:198:ARG:CD	2.07	0.66
2:B:410:F6P:H11	4:B:412:PO4:O4	1.97	0.65
1:A:59:ILE:HD12	1:B:10:ILE:HG21	1.79	0.64
1:B:182:ASN:ND2	1:B:199:ASP:H	1.95	0.64
1:B:218:GLU:OE2	1:B:268:LYS:HB2	1.98	0.64
1:A:95:VAL:HG22	1:A:310:ILE:HD11	1.81	0.63
1:B:317:ILE:HG21	1:B:327:LEU:HD23	1.81	0.63
1:A:276:ARG:HH21	1:A:311:HIS:HB3	1.64	0.62
1:A:102:ALA:HB2	1:A:149:GLU:HG3	1.80	0.62
1:B:216:ALA:HA	1:B:219:PHE:CD2	2.34	0.62
1:A:29:GLU:OE2	1:A:112:LYS:HG2	2.00	0.61
1:B:269:LYS:HD3	1:B:269:LYS:O	2.01	0.61
1:B:186:LEU:HB2	1:B:193:PHE:CE1	2.36	0.60
1:A:108:GLU:HG2	1:A:109:LYS:HG3	1.84	0.60
1:B:173:LEU:C	1:B:173:LEU:HD23	2.27	0.60
1:B:308:THR:H	1:B:312:GLN:NE2	1.98	0.60
1:A:182:ASN:HD21	1:A:199:ASP:H	1.50	0.59
1:A:10:ILE:HB	1:B:57:TRP:O	2.02	0.59
1:B:15:ARG:HG2	1:B:15:ARG:HH11	1.67	0.59
1:B:308:THR:H	1:B:312:GLN:HE22	1.51	0.58
1:A:45:SER:O	1:A:49:ARG:HD3	2.03	0.58
1:A:157:ARG:HD2	1:A:304:ASP:OD2	2.04	0.57
1:B:104:ILE:N	1:B:104:ILE:HD12	2.19	0.57
1:B:164:TYR:OH	1:B:253:HIS:HD2	1.87	0.57
1:B:184:PHE:HB3	1:B:193:PHE:HB3	1.87	0.57
1:A:186:LEU:HB2	1:A:193:PHE:CE1	2.39	0.56
1:B:106:GLU:H	1:B:106:GLU:CD	2.13	0.56
1:B:267:ASN:C	1:B:267:ASN:HD22	2.13	0.56
1:A:276:ARG:HG3	1:A:279:TYR:CE1	2.40	0.56
1:A:330:ILE:O	1:A:333:LYS:HB3	2.06	0.56
1:B:29:GLU:OE1	1:B:112:LYS:HG2	2.05	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:ARG:O	1:B:280:GLU:HB2	2.06	0.56
1:B:277:LEU:HA	1:B:281:CYS:HB2	1.88	0.56
1:B:142:ASN:H	1:B:142:ASN:HD22	1.52	0.56
1:A:216:ALA:HA	1:A:219:PHE:CD2	2.40	0.55
1:A:296:THR:HG21	1:A:328:LEU:HD21	1.88	0.55
1:B:102:ALA:HB3	1:B:149:GLU:HG2	1.89	0.55
1:B:182:ASN:HD21	1:B:199:ASP:H	1.54	0.54
1:A:102:ALA:CB	1:A:149:GLU:HG3	2.37	0.54
1:A:300:GLU:N	1:A:300:GLU:OE1	2.39	0.54
1:B:270:SER:HB3	1:B:273:GLY:C	2.32	0.54
1:B:270:SER:HB3	1:B:273:GLY:O	2.07	0.54
1:B:78:ASN:O	1:B:82:ILE:HG12	2.08	0.54
1:A:94:LEU:HB2	1:A:103:ILE:HB	1.90	0.53
1:B:274:LYS:HE2	2:B:410:F6P:O1	2.08	0.53
1:A:74:ASP:OD1	1:A:75:VAL:HG23	2.08	0.53
1:B:267:ASN:HD21	1:B:270:SER:H	1.56	0.53
1:A:269:LYS:N	1:A:269:LYS:HD2	2.24	0.53
1:B:267:ASN:HD21	1:B:270:SER:HB2	1.74	0.53
1:A:267:ASN:O	1:A:270:SER:N	2.41	0.53
1:B:117:PHE:CD2	1:B:119:PRO:HD3	2.44	0.53
1:B:218:GLU:HG2	1:B:268:LYS:HG3	1.91	0.52
1:B:279:TYR:OH	1:B:311:HIS:HD2	1.92	0.52
1:B:45:SER:O	1:B:49:ARG:HD3	2.10	0.51
1:B:318:LEU:C	1:B:318:LEU:HD12	2.35	0.51
1:A:299:LYS:HB2	1:A:300:GLU:OE1	2.10	0.51
1:A:36:SER:HB3	1:A:84:VAL:HG12	1.90	0.51
1:A:278:LEU:HD22	6:A:628:HOH:O	2.10	0.51
1:A:72:LYS:HG2	1:A:74:ASP:OD2	2.11	0.51
1:B:142:ASN:HD22	1:B:142:ASN:N	2.08	0.50
1:B:58:GLY:HA2	1:B:62:SER:C	2.36	0.50
1:A:223:ILE:HG12	1:A:265:PRO:HG3	1.93	0.49
1:B:72:LYS:HD2	1:B:72:LYS:N	2.21	0.49
1:B:142:ASN:H	1:B:142:ASN:ND2	2.10	0.49
1:B:205:LYS:HE2	1:B:322:GLU:OE1	2.13	0.49
1:B:126:ILE:HD13	1:B:132:ILE:HG21	1.94	0.49
1:B:229:ARG:NH1	1:B:330:ILE:HD11	2.26	0.49
1:A:251:ASP:OD1	1:A:254:ARG:NH2	2.46	0.49
1:B:194:ILE:HD12	1:B:194:ILE:N	2.28	0.49
1:A:37:LEU:HD23	1:A:85:LEU:HD21	1.95	0.48
1:A:150:LYS:C	1:A:150:LYS:HD2	2.39	0.48
1:A:318:LEU:C	1:A:318:LEU:HD12	2.38	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:ALA:C	1:B:111:GLY:HA3	2.38	0.48
1:B:112:LYS:HD3	1:B:140:ARG:NE	2.29	0.48
1:A:178:VAL:HA	1:A:290:LYS:NZ	2.29	0.48
1:B:72:LYS:H	1:B:72:LYS:CD	2.12	0.48
1:A:57:TRP:HB2	1:B:10:ILE:HD13	1.96	0.48
6:A:553:HOH:O	1:B:253:HIS:HE1	1.97	0.48
1:A:297:THR:HG21	1:A:305:ILE:HD11	1.95	0.47
1:B:332:GLN:HE21	1:B:332:GLN:HB3	1.46	0.47
1:A:150:LYS:HD2	1:A:151:ASP:N	2.29	0.47
1:B:277:LEU:HD12	1:B:307:PRO:HB3	1.96	0.47
1:A:157:ARG:NH2	1:A:289:GLU:OE2	2.47	0.47
1:A:242:ALA:HB3	1:B:212:ASN:OD1	2.14	0.47
1:B:198:ARG:O	1:B:199:ASP:C	2.56	0.47
1:B:307:PRO:HA	1:B:312:GLN:HE22	1.79	0.47
1:B:10:ILE:O	1:B:10:ILE:HG23	2.14	0.47
1:A:13:LEU:CD2	1:A:17:VAL:HG23	2.44	0.47
1:B:56:LEU:HD12	1:B:57:TRP:CE3	2.50	0.47
1:B:133:GLY:HA2	1:B:167:TYR:CD1	2.50	0.46
1:A:194:ILE:O	1:A:196:VAL:HG13	2.15	0.46
1:A:217:LYS:HG2	1:A:217:LYS:O	2.16	0.46
1:B:229:ARG:CZ	1:B:330:ILE:HD11	2.45	0.46
1:A:254:ARG:HD2	6:A:572:HOH:O	2.15	0.46
1:B:182:ASN:ND2	1:B:198:ARG:HA	2.27	0.46
1:A:44:ILE:O	1:A:48:VAL:HG23	2.16	0.46
1:A:310:ILE:HG23	1:A:311:HIS:ND1	2.31	0.46
1:A:121:ASP:OD1	1:A:249:VAL:HG22	2.16	0.45
1:A:305:ILE:O	1:A:307:PRO:HD3	2.16	0.45
1:B:221:PRO:HB2	1:B:334:HIS:HD2	1.81	0.45
1:B:316:ILE:HD11	1:B:318:LEU:HD23	1.98	0.45
1:A:133:GLY:HA2	1:A:167:TYR:CD1	2.51	0.45
1:B:157:ARG:NH2	1:B:289:GLU:OE1	2.50	0.45
1:B:267:ASN:ND2	1:B:267:ASN:C	2.74	0.45
1:A:90:ALA:C	1:A:111:GLY:HA3	2.41	0.45
1:A:223:ILE:HD11	1:A:331:TYR:OH	2.16	0.45
1:A:173:LEU:C	1:A:173:LEU:HD23	2.42	0.45
1:B:254:ARG:HG2	1:B:254:ARG:HH11	1.82	0.45
1:A:96:SER:HB2	1:A:117:PHE:CZ	2.52	0.44
1:B:128:CYS:SG	1:B:130:VAL:HG23	2.57	0.44
1:B:139:TYR:CE1	1:B:159:LEU:HG	2.53	0.44
1:B:221:PRO:HB2	1:B:334:HIS:CD2	2.53	0.44
1:A:97:GLU:HG2	1:A:97:GLU:O	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:SER:O	1:B:89:PHE:HB2	2.18	0.44
1:A:267:ASN:O	1:A:268:LYS:C	2.61	0.43
1:B:103:ILE:C	1:B:104:ILE:HD12	2.43	0.43
1:B:276:ARG:HG2	1:B:279:TYR:OH	2.19	0.43
1:B:298:GLY:C	1:B:299:LYS:HD2	2.43	0.43
1:B:36:SER:HB3	1:B:84:VAL:HG12	2.00	0.43
1:A:223:ILE:CD1	1:A:331:TYR:OH	2.66	0.43
1:B:276:ARG:HD2	6:B:624:HOH:O	2.17	0.43
1:B:326:GLU:O	1:B:330:ILE:HG12	2.19	0.43
1:B:119:PRO:HG2	6:B:669:HOH:O	2.18	0.43
1:A:261:ILE:HD12	1:A:263:MET:CG	2.48	0.43
1:B:164:TYR:OH	1:B:253:HIS:CD2	2.70	0.43
1:B:267:ASN:ND2	1:B:270:SER:HB2	2.33	0.43
1:B:316:ILE:CD1	1:B:318:LEU:HD23	2.49	0.43
1:A:10:ILE:O	1:A:15:ARG:NH2	2.52	0.42
1:B:56:LEU:HB3	1:B:57:TRP:CE3	2.54	0.42
1:A:276:ARG:HG3	1:A:279:TYR:HE1	1.83	0.42
1:A:225:GLU:OE1	1:A:334:HIS:HE1	2.03	0.42
1:B:263:MET:HG2	1:B:317:ILE:HG23	2.01	0.42
1:B:228:GLN:HA	1:B:228:GLN:OE1	2.20	0.42
1:B:212:ASN:HB2	1:B:244:TYR:CE2	2.55	0.42
1:B:278:LEU:HD13	6:B:555:HOH:O	2.19	0.42
1:A:276:ARG:HG2	1:A:280:GLU:OE1	2.20	0.42
1:B:223:ILE:O	1:B:227:ILE:HG13	2.20	0.42
1:A:178:VAL:HG13	1:A:179:ASN:N	2.35	0.41
1:B:112:LYS:HD3	1:B:140:ARG:HE	1.85	0.41
1:B:282:ASN:HB2	1:B:283:PRO:HD3	2.03	0.41
1:A:81:VAL:HG11	1:A:117:PHE:CE2	2.56	0.41
1:B:15:ARG:HG2	1:B:15:ARG:NH1	2.31	0.41
1:B:97:GLU:O	1:B:97:GLU:HG2	2.20	0.41
1:B:130:VAL:HG12	1:B:131:SER:N	2.34	0.41
1:B:80:LEU:O	1:B:84:VAL:HG23	2.20	0.41
1:A:220:ASP:OD2	1:A:223:ILE:HD13	2.20	0.41
1:B:13:LEU:HD23	1:B:184:PHE:CE1	2.56	0.41
1:A:276:ARG:NH1	1:A:313:ARG:CZ	2.84	0.41
1:A:53:ILE:HG22	1:B:187:ASP:HB2	2.02	0.41
1:B:230:LYS:HE2	1:B:230:LYS:CA	2.43	0.41
1:B:108:GLU:HG2	1:B:109:LYS:HG3	2.02	0.41
1:A:81:VAL:HG11	1:A:117:PHE:CD2	2.56	0.41
1:A:218:GLU:OE1	1:A:269:LYS:HD3	2.21	0.41
1:A:317:ILE:HG21	1:A:327:LEU:HD23	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:PRO:HA	1:B:234:PRO:HD2	1.88	0.41
1:B:317:ILE:HD13	1:B:327:LEU:CD2	2.51	0.41
1:A:218:GLU:OE1	1:A:269:LYS:HB2	2.21	0.40
1:A:114:VAL:HB	1:A:139:TYR:HB2	2.03	0.40
1:A:123:SER:HA	1:A:126:ILE:HG13	2.02	0.40
1:B:35:ASN:HD22	1:B:35:ASN:HA	1.66	0.40
1:A:282:ASN:HB2	1:A:283:PRO:HD3	2.04	0.40
1:B:274:LYS:O	1:B:313:ARG:NH1	2.51	0.40
1:A:266:ALA:O	1:A:267:ASN:HB3	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	315/337 (94%)	297 (94%)	16 (5%)	2 (1%)	21	38
1	B	315/337 (94%)	304 (96%)	10 (3%)	1 (0%)	36	55
All	All	630/674 (94%)	601 (95%)	26 (4%)	3 (0%)	24	43

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	LYS
1	B	268	LYS
1	A	270	SER

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	265/279 (95%)	259 (98%)	6 (2%)	44	72
1	B	265/279 (95%)	257 (97%)	8 (3%)	36	64
All	All	530/558 (95%)	516 (97%)	14 (3%)	40	68

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	37	LEU
1	A	74	ASP
1	A	249	VAL
1	A	261	ILE
1	A	277	LEU
1	B	37	LEU
1	B	72	LYS
1	B	142	ASN
1	B	157	ARG
1	B	267	ASN
1	B	277	LEU
1	B	278	LEU
1	B	332	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	182	ASN
1	A	253	HIS
1	A	311	HIS
1	A	312	GLN
1	A	332	GLN
1	B	35	ASN
1	B	125	ASN
1	B	142	ASN
1	B	154	GLN
1	B	158	ASN
1	B	182	ASN
1	B	253	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	267	ASN
1	B	311	HIS
1	B	312	GLN
1	B	332	GLN
1	B	334	HIS

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	$\# Z > 2$
5	AMP	B	411	-	25,25,25	0.34	0	37,38,38	0.38	0
2	F6P	B	410	-	15,16,16	0.57	0	16,25,25	0.59	0
4	PO4	A	402	3	4,4,4	1.93	3 (75%)	6,6,6	0.48	0
5	AMP	A	401	-	25,25,25	0.34	0	37,38,38	0.34	0
2	F6P	A	400	-	15,16,16	0.60	0	16,25,25	0.60	0
4	PO4	B	412	3	4,4,4	1.70	1 (25%)	6,6,6	0.47	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	400	-	-	8/9/28/28	0/1/1/1
2	F6P	B	410	-	-	3/9/28/28	0/1/1/1
5	AMP	B	411	-	-	0/10/26/26	0/3/3/3
5	AMP	A	401	-	-	0/10/26/26	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	402	PO4	P-O2	-2.28	1.48	1.54
4	A	402	PO4	P-O4	-2.18	1.48	1.54
4	B	412	PO4	P-O4	-2.13	1.48	1.54
4	A	402	PO4	P-O3	-2.12	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	F6P	O1-C1-C2-O2
2	A	400	F6P	O1-C1-C2-C3
2	A	400	F6P	C6-O6-P-O2P
2	A	400	F6P	C6-O6-P-O3P
2	B	410	F6P	O1-C1-C2-O2
2	B	410	F6P	O1-C1-C2-C3
2	A	400	F6P	C6-O6-P-O1P
2	A	400	F6P	C4-C5-C6-O6
2	A	400	F6P	O1-C1-C2-O5
2	B	410	F6P	O1-C1-C2-O5
2	A	400	F6P	O5-C5-C6-O6

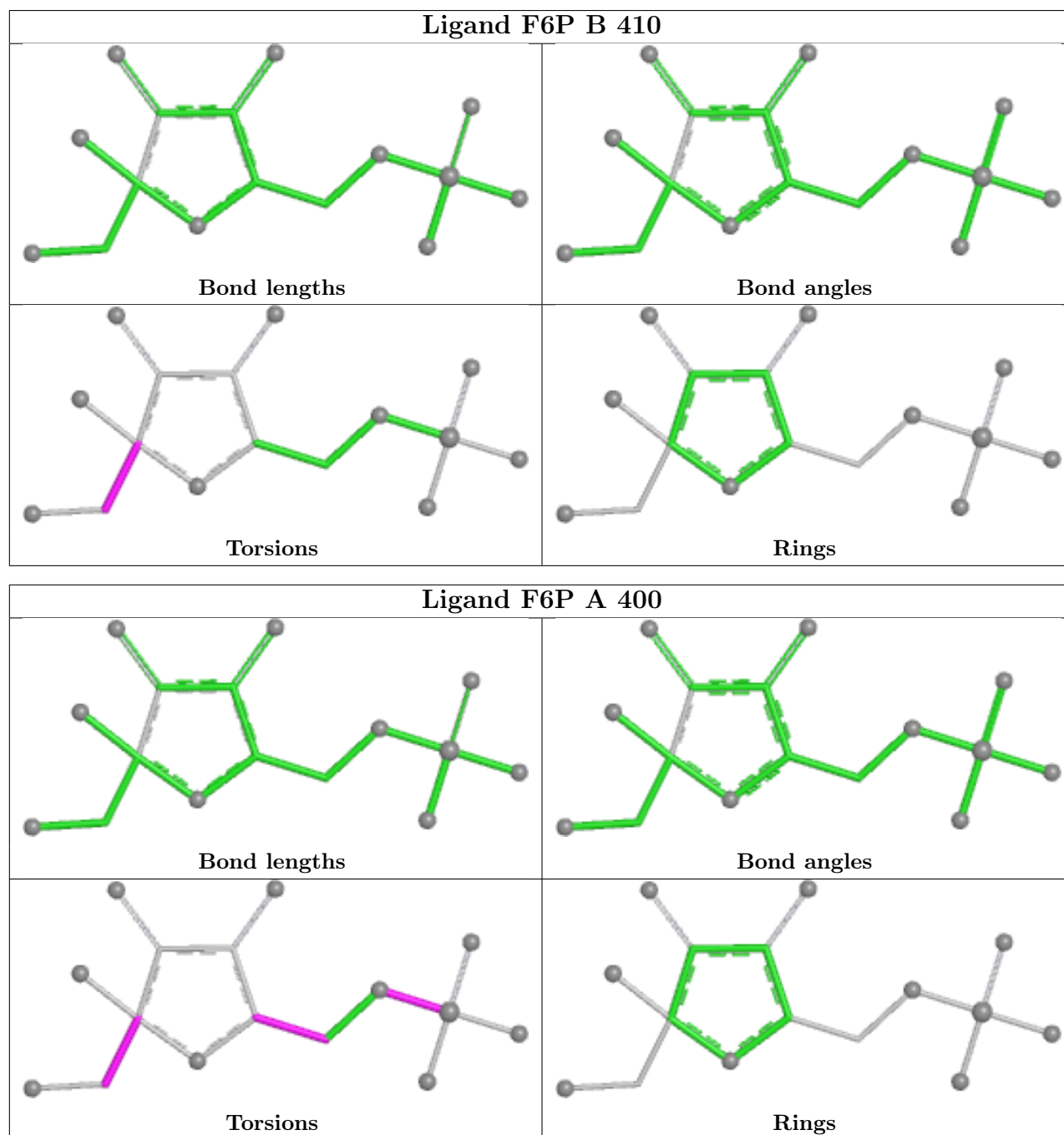
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	410	F6P	3	0
4	B	412	PO4	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

There are no such residues in this entry.

5.8 Polymer linkage issues

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.