



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:22 PM GMT

PDB ID : 1FBF
Title : CRYSTALLOGRAPHIC STUDIES OF THE CATALYTIC MECHANISM OF
THE NEUTRAL FORM OF FRUCTOSE-1,6-BISPHOSPHATASE
Authors : Zhang, Y.; Liang, J.-Y.; Huang, S.; Ke, H.; Lipscomb, W.N.
Deposited on : 1992-10-16
Resolution : 2.70 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

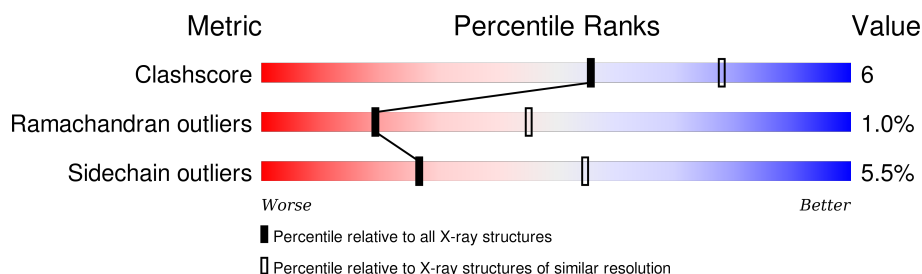
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

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	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6038 atoms, of which 1156 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	313	Total	C	H	N	O	S	0	0	1
			2920	1520	529	403	453	15			
1	B	315	Total	C	H	N	O	S	0	0	1
			2938	1532	531	405	455	15			

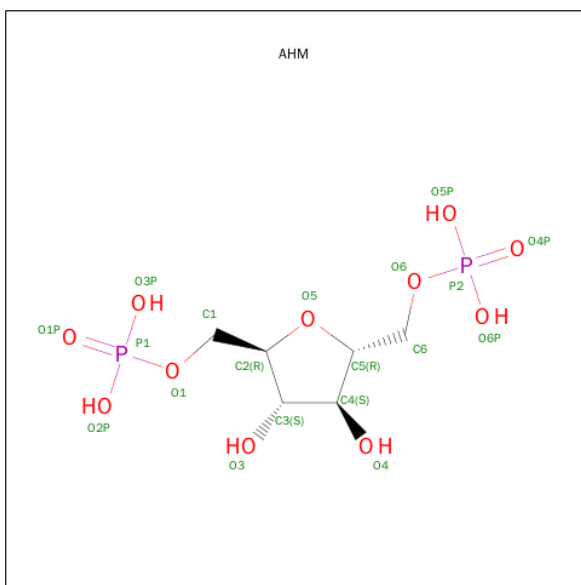
There are 6 discrepancies between the modelled and reference sequences:

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

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

- Molecule 3 is 2,5-ANHYDROMANNITOL-1,6-DIPHOSPHATE (three-letter code: AHM) (formula: C₆H₁₄O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	O	P	0	0
			23	6	4	11	2		
3	B	1	Total	C	H	O	P	0	0
			23	6	4	11	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	27	Total	H	O	0	0
			81	54	27		
4	B	17	Total	H	O	0	0
			51	34	17		

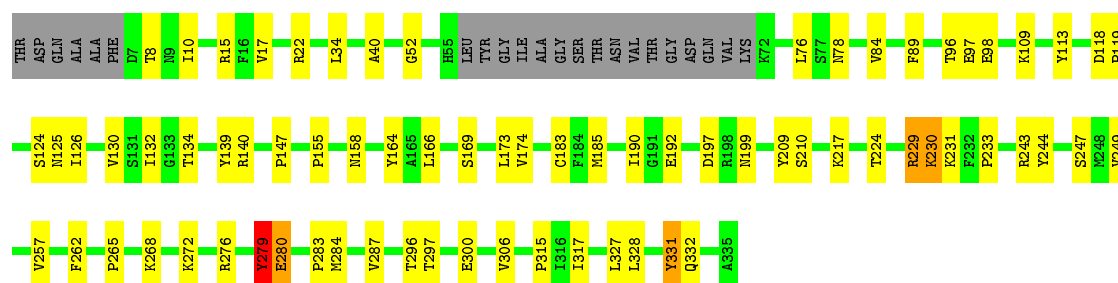
3 Residue-property plots

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.

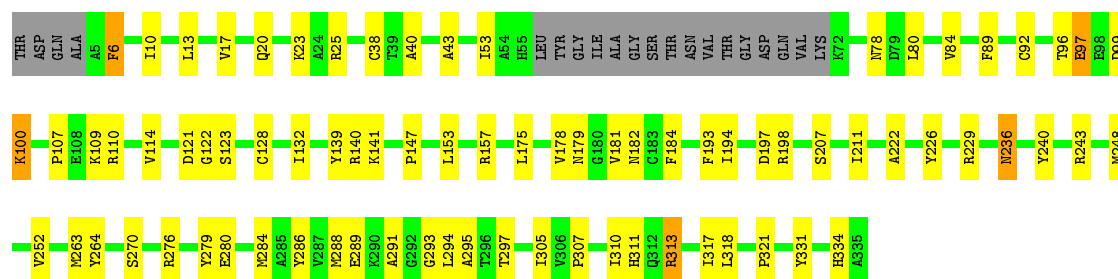
• Molecule 1: FRUCTOSE 1,6-BISPHOSPHATASE

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 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.40 Å 131.40 Å 69.20 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.189 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6038	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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: MG, AHM

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	0.75	0/2430	1.40	11/3286 (0.3%)
1	B	0.75	0/2447	1.44	18/3309 (0.5%)
All	All	0.75	0/4877	1.42	29/6595 (0.4%)

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	6
1	B	0	2
All	All	0	8

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	121	ASP	CB-CG-OD1	9.79	127.12	118.30
1	A	15	ARG	NE-CZ-NH1	8.88	124.74	120.30
1	B	313	ARG	NE-CZ-NH2	-8.07	116.26	120.30
1	A	243	ARG	NE-CZ-NH2	-7.86	116.37	120.30
1	B	121	ASP	CB-CG-OD2	-7.63	111.43	118.30
1	B	313	ARG	NE-CZ-NH1	7.27	123.93	120.30
1	B	240	TYR	CB-CG-CD2	-6.91	116.86	121.00
1	B	140	ARG	NE-CZ-NH2	-6.35	117.13	120.30
1	B	97	GLU	CA-CB-CG	6.33	127.34	113.40
1	A	22	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	B	243	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	B	157	ARG	NE-CZ-NH1	6.04	123.32	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	140	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	A	217	LYS	CA-CB-CG	5.96	126.52	113.40
1	B	279	TYR	CB-CG-CD2	-5.78	117.53	121.00
1	B	280	GLU	OE1-CD-OE2	-5.75	116.41	123.30
1	B	248	MET	CA-CB-CG	-5.69	103.63	113.30
1	A	140	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	A	284	MET	CG-SD-CE	-5.56	91.31	100.20
1	A	15	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	A	118	ASP	CB-CG-OD1	5.39	123.16	118.30
1	A	209	TYR	CB-CG-CD2	5.38	124.22	121.00
1	B	17	VAL	CA-CB-CG2	-5.37	102.85	110.90
1	B	229	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	B	38	CYS	CA-CB-SG	-5.31	104.45	114.00
1	A	209	TYR	CB-CG-CD1	-5.29	117.82	121.00
1	B	132	ILE	CA-C-N	5.19	126.59	116.20
1	B	25	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	A	332	GLN	CA-CB-CG	5.08	124.57	113.40

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	164	TYR	Sidechain
1	A	244	TYR	Sidechain
1	A	279	TYR	Sidechain
1	A	331	TYR	Sidechain
1	B	226	TYR	Sidechain
1	B	286	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	2391	529	2448	29	0
1	B	2407	531	2462	34	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	19	4	10	1	0
3	B	19	4	10	4	0
4	A	27	54	0	0	0
4	B	17	34	0	0	0
All	All	4882	1156	4930	61	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 6.

All (61) 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:B:182:ASN:HD22	1:B:198:ARG:HA	1.40	0.86
1:B:252:VAL:HG11	1:B:284:MET:SD	2.21	0.81
1:B:78:ASN:OD1	1:B:96:THR:HG21	1.80	0.80
1:A:125:ASN:HB3	1:A:130:VAL:HB	1.64	0.80
1:B:182:ASN:ND2	1:B:198:ARG:HA	1.99	0.77
3:B:336:AHM:O2P	3:B:336:AHM:H2	1.88	0.72
1:A:96:THR:HG22	1:A:98:GLU:H	1.56	0.70
1:B:310:ILE:HG13	1:B:311:HIS:CD2	2.36	0.61
1:B:100:LYS:O	1:B:310:ILE:HD11	2.02	0.59
1:B:114:VAL:HB	1:B:139:TYR:HB2	1.84	0.59
1:A:283:PRO:O	1:A:287:VAL:HG23	2.03	0.58
1:B:294:LEU:O	1:B:318:LEU:HA	2.03	0.58
1:A:297:THR:HG22	1:A:315:PRO:O	2.03	0.58
1:B:40:ALA:HB2	1:B:84:VAL:HG21	1.86	0.57
1:B:297:THR:HG21	1:B:305:ILE:HD11	1.86	0.56
1:A:174:VAL:HG22	1:A:183:CYS:SG	2.49	0.53
1:A:317:ILE:HG21	1:A:327:LEU:HD23	1.92	0.52
1:A:247:SER:HA	3:A:336:AHM:O3	2.11	0.51
1:B:288:MET:HG2	1:B:318:LEU:HB2	1.93	0.51
1:A:119:PRO:HA	1:A:134:THR:HG23	1.93	0.50
1:A:296:THR:HG21	1:A:328:LEU:HD21	1.93	0.50
1:A:185:MET:SD	1:B:53:ILE:HG13	2.51	0.50
1:A:210:SER:HB3	1:A:262:PHE:HA	1.94	0.50
1:B:89:PHE:HA	1:B:109:LYS:O	2.12	0.50
1:B:295:ALA:HA	1:B:317:ILE:O	2.11	0.50
1:A:17:VAL:HG11	1:A:34:LEU:HD12	1.94	0.50
1:A:40:ALA:HB2	1:A:84:VAL:HG21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:PRO:HG3	1:A:306:VAL:HG22	1.92	0.49
1:B:153:LEU:O	1:B:307:PRO:HD2	2.14	0.48
1:A:89:PHE:CE1	1:A:109:LYS:HG2	2.49	0.47
1:B:211:ILE:HD12	1:B:263:MET:HB2	1.97	0.47
1:B:264:TYR:CE2	3:B:336:AHM:H62	2.51	0.46
1:A:276:ARG:N	1:A:280:GLU:HG3	2.29	0.46
1:B:96:THR:HG22	1:B:97:GLU:N	2.30	0.46
1:B:10:ILE:HG22	1:B:194:ILE:HD12	1.97	0.46
1:B:107:PRO:HA	1:B:110:ARG:HD2	1.97	0.46
1:B:10:ILE:CG2	1:B:194:ILE:HD12	2.46	0.45
1:A:229:ARG:HG3	1:A:230:LYS:HD2	1.99	0.44
1:B:20:GLN:HA	1:B:23:LYS:HE2	1.99	0.44
1:A:166:LEU:HD13	1:A:249:VAL:HG12	2.00	0.44
1:B:276:ARG:NH1	1:B:313:ARG:HD3	2.32	0.44
1:A:257:VAL:HG21	1:B:128:CYS:HA	2.00	0.43
1:B:122:GLY:HA2	3:B:336:AHM:O2P	2.19	0.43
1:A:96:THR:HG22	1:A:97:GLU:N	2.34	0.43
1:A:126:ILE:HG12	1:A:132:ILE:HD13	2.00	0.42
1:A:185:MET:HG2	1:B:53:ILE:HD11	2.02	0.42
1:B:289:GLU:HA	1:B:293:GLY:O	2.20	0.42
1:B:43:ALA:HB3	1:B:80:LEU:HD21	2.01	0.42
1:A:231:LYS:O	1:A:233:PRO:HD3	2.20	0.41
1:B:181:VAL:HG21	1:B:291:ALA:HB2	2.03	0.41
1:B:141:LYS:HE3	1:B:147:PRO:HG3	2.02	0.41
1:B:293:GLY:HA2	1:B:321:PRO:HD3	2.03	0.41
1:B:264:TYR:HE2	3:B:336:AHM:H62	1.85	0.41
1:A:97:GLU:HG3	1:A:279:TYR:CE1	2.56	0.41
1:A:272:LYS:HA	1:A:272:LYS:HD3	1.76	0.40
1:B:222:ALA:HA	1:B:334:HIS:HD2	1.85	0.40
1:B:184:PHE:HB3	1:B:193:PHE:HB3	2.02	0.40
1:A:78:ASN:OD1	1:A:96:THR:HG21	2.21	0.40
1:A:97:GLU:HG3	1:A:279:TYR:HE1	1.87	0.40
1:A:155:PRO:HD2	1:A:158:ASN:ND2	2.37	0.40
1:A:190:ILE:HG13	1:A:192:GLU:HB2	2.04	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	309/335 (92%)	292 (94%)	16 (5%)	1 (0%)	46	75
1	B	311/335 (93%)	287 (92%)	19 (6%)	5 (2%)	12	30
All	All	620/670 (92%)	579 (93%)	35 (6%)	6 (1%)	19	45

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	236	ASN
1	B	178	VAL
1	A	52	GLY
1	B	6	PHE
1	B	123	SER
1	B	100	LYS

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	261/278 (94%)	243 (93%)	18 (7%)	19	43
1	B	262/278 (94%)	251 (96%)	11 (4%)	36	68
All	All	523/556 (94%)	494 (94%)	29 (6%)	27	55

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	A	10	ILE
1	A	76	LEU
1	A	124	SER
1	A	147	PRO
1	A	169	SER
1	A	173	LEU
1	A	197	ASP
1	A	199	ASN
1	A	224	THR
1	A	229	ARG
1	A	230	LYS
1	A	265	PRO
1	A	268	LYS
1	A	279	TYR
1	A	280	GLU
1	A	300	GLU
1	A	331	TYR
1	B	6	PHE
1	B	13	LEU
1	B	92	CYS
1	B	99	ASP
1	B	175	LEU
1	B	179	ASN
1	B	197	ASP
1	B	207	SER
1	B	236	ASN
1	B	270	SER
1	B	331	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	199	ASN
1	B	182	ASN
1	B	282	ASN
1	B	334	HIS

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
3	AHM	A	336	2	19,19,19	0.71	0	27,29,29	0.94	1 (3%)
3	AHM	B	336	2	19,19,19	0.80	0	27,29,29	0.95	1 (3%)

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	AHM	A	336	2	-	0/12/28/28	0/1/1/1
3	AHM	B	336	2	-	0/12/28/28	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	336	AHM	O2P-P1-O1P	2.64	119.09	110.58
3	A	336	AHM	O2P-P1-O1P	2.68	119.19	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	336	AHM	1	0
3	B	336	AHM	4	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.