



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 12:55 AM GMT

PDB ID : 2F3X
Title : Crystal structure of FapR (in complex with effector)- a global regulator of fatty acid biosynthesis in B. subtilis
Authors : Buschiazzo, A.; Guerin, M.E.; Alzari, P.M.
Deposited on : 2005-11-22
Resolution : 3.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

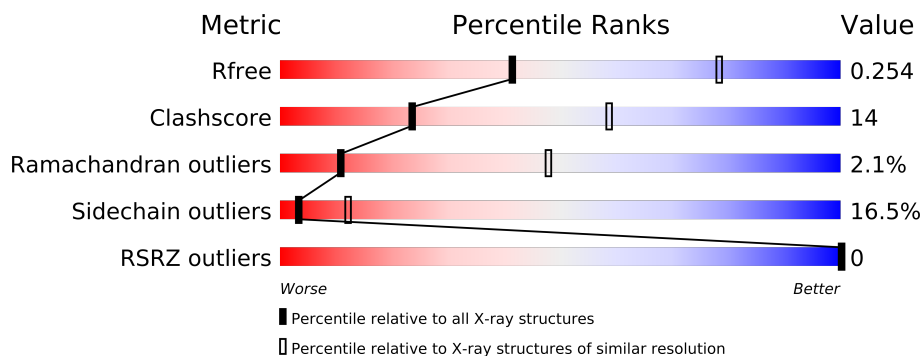
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.10 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}	66092	1007 (3.18-3.02)
Clashscore	79885	1078 (3.16-3.04)
Ramachandran outliers	78287	1044 (3.16-3.04)
Sidechain outliers	78261	1044 (3.16-3.04)
RSRZ outliers	66119	1008 (3.18-3.02)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	157	<div><div></div><div></div></div>
1	B	157	<div><div></div><div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2312 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

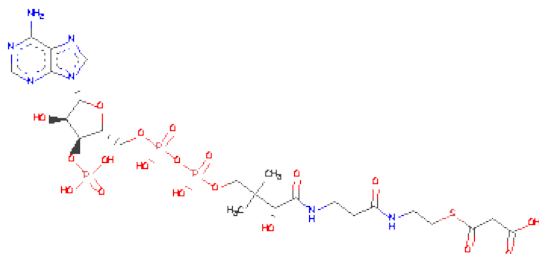
- Molecule 1 is a protein called Transcription factor fapR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	143	Total	C	N	O	S	0	0	0
			1110	694	193	222	1			
1	B	143	Total	C	N	O	S	0	0	0
			1128	705	198	224	1			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	MET	-	CLONING ARTIFACT	UNP O34835
A	33	ARG	-	CLONING ARTIFACT	UNP O34835
A	34	GLY	-	CLONING ARTIFACT	UNP O34835
A	35	SER	-	CLONING ARTIFACT	UNP O34835
A	36	HIS	-	EXPRESSION TAG	UNP O34835
A	37	HIS	-	EXPRESSION TAG	UNP O34835
A	38	HIS	-	EXPRESSION TAG	UNP O34835
A	39	HIS	-	EXPRESSION TAG	UNP O34835
A	40	HIS	-	EXPRESSION TAG	UNP O34835
A	41	HIS	-	EXPRESSION TAG	UNP O34835
A	42	GLY	-	CLONING ARTIFACT	UNP O34835
A	43	SER	-	CLONING ARTIFACT	UNP O34835
B	32	MET	-	CLONING ARTIFACT	UNP O34835
B	33	ARG	-	CLONING ARTIFACT	UNP O34835
B	34	GLY	-	CLONING ARTIFACT	UNP O34835
B	35	SER	-	CLONING ARTIFACT	UNP O34835
B	36	HIS	-	EXPRESSION TAG	UNP O34835
B	37	HIS	-	EXPRESSION TAG	UNP O34835
B	38	HIS	-	EXPRESSION TAG	UNP O34835
B	39	HIS	-	EXPRESSION TAG	UNP O34835
B	40	HIS	-	EXPRESSION TAG	UNP O34835
B	41	HIS	-	EXPRESSION TAG	UNP O34835
B	42	GLY	-	CLONING ARTIFACT	UNP O34835
B	43	SER	-	CLONING ARTIFACT	UNP O34835

- Molecule 2 is MALONYL-COENZYME A (three-letter code: MLC) (formula: $C_{24}H_{38}N_7O_{19}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	S	0	0
			32	14	2	13	2	1		
2	A	1	Total	C	N	O	P	S	0	0
			32	14	2	13	2	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total	O	0	0
			8	8		
3	B	2	Total	O	0	0
			2	2		

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	89.44Å 89.44Å 162.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.25 – 3.10 63.24 – 3.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (63.25-3.10) 100.0 (63.24-3.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.63 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.187 , 0.226 0.231 , 0.254	Depositor DCC
R_{free} test set	941 reflections (8.11%)	DCC
Wilson B-factor (Å ²)	69.9	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 12545 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2312	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1.01	3/1120 (0.3%)	0.99	4/1512 (0.3%)
1	B	0.91	0/1138	1.00	3/1532 (0.2%)
All	All	0.96	3/2258 (0.1%)	1.00	7/3044 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	77	GLU	CG-CD	7.74	1.63	1.51
1	A	172	GLU	CG-CD	6.93	1.62	1.51
1	A	172	GLU	CB-CG	6.41	1.64	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	ARG	NE-CZ-NH1	8.44	124.52	120.30
1	A	121	VAL	CB-CA-C	-6.93	98.23	111.40
1	B	121	VAL	CB-CA-C	-6.80	98.48	111.40
1	A	138	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	A	184	ARG	N-CA-C	5.56	126.02	111.00
1	B	72	ASP	CB-CA-C	-5.25	99.90	110.40
1	B	99	PHE	CB-CA-C	-5.00	100.39	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens

added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1110	0	1103	32	0
1	B	1128	0	1140	30	0
2	A	32	0	22	8	0
2	B	32	0	21	5	0
3	A	8	0	0	0	0
3	B	2	0	0	0	0
All	All	2312	0	2286	66	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (66) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:107:GLY:H	2:B:390:MLC:HM21	1.27	0.95
1:A:180:PHE:HB3	1:A:182:MET:CE	1.99	0.93
1:B:107:GLY:H	2:B:390:MLC:CM2	1.89	0.85
1:A:180:PHE:HB3	1:A:182:MET:HE3	1.65	0.78
1:B:107:GLY:N	2:B:390:MLC:HM21	2.05	0.70
1:B:129:THR:HA	1:B:182:MET:HE2	1.72	0.69
2:A:391:MLC:HP81	2:A:391:MLC:NP2	2.11	0.64
2:A:391:MLC:NP2	2:A:391:MLC:CP8	2.60	0.63
1:A:102:ASN:C	1:A:102:ASN:HD22	2.02	0.63
2:B:390:MLC:O5'	2:B:390:MLC:O7	2.17	0.62
1:A:98:VAL:CG1	1:A:103:GLN:HA	2.30	0.62
1:A:180:PHE:HB3	1:A:182:MET:HE2	1.79	0.62
1:B:102:ASN:C	1:B:102:ASN:HD22	2.03	0.62
1:A:56:ASN:HD22	1:A:56:ASN:N	1.96	0.61
1:A:98:VAL:HG13	1:A:103:GLN:HA	1.81	0.61
1:A:119:VAL:HG21	2:B:390:MLC:OM3	2.00	0.61
1:B:184:ARG:O	1:B:185:SER:HB2	2.00	0.61
1:B:169:TYR:CE2	1:B:174:ILE:HG12	2.37	0.59
1:B:122:ILE:HG21	1:B:127:ALA:HB2	1.85	0.58
2:A:391:MLC:HP81	2:A:391:MLC:HP2	1.69	0.57
1:A:67:LYS:O	1:B:101:ARG:NH2	2.38	0.57
1:B:47:ILE:N	1:B:48:PRO:HD2	2.19	0.57
1:B:161:ARG:C	1:B:161:ARG:HD2	2.28	0.55
1:B:98:VAL:HG13	1:B:103:GLN:HA	1.90	0.52
1:A:99:PHE:HZ	2:A:391:MLC:HM22	1.76	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:167:ASN:ND2	1:B:177:SER:HB2	2.24	0.51
1:A:98:VAL:HG21	1:A:142:GLN:HB2	1.93	0.51
1:A:82:GLU:O	1:A:85:ASP:HB3	2.10	0.51
1:A:124:ASP:OD1	1:A:185:SER:OG	2.28	0.51
1:B:129:THR:HG23	1:B:182:MET:HE1	1.93	0.50
1:A:85:ASP:C	1:A:85:ASP:OD1	2.50	0.50
1:A:180:PHE:CB	1:A:182:MET:HE3	2.40	0.50
1:B:98:VAL:HG21	1:B:142:GLN:HB2	1.95	0.48
1:B:47:ILE:HG22	1:B:51:ARG:HH21	1.78	0.48
2:A:391:MLC:OP1	2:A:391:MLC:HP12	2.14	0.47
1:A:169:TYR:CE2	1:A:174:ILE:HG12	2.51	0.46
1:A:56:ASN:ND2	1:A:56:ASN:N	2.61	0.46
1:B:83:LEU:HB3	1:B:123:ASP:HB2	1.98	0.45
1:A:99:PHE:CZ	2:A:391:MLC:HM22	2.52	0.45
1:A:167:ASN:ND2	1:A:177:SER:HB2	2.32	0.45
1:B:99:PHE:HB2	1:B:102:ASN:HD21	1.82	0.45
1:B:46:SER:HB2	1:B:48:PRO:HD2	1.99	0.45
1:A:113:GLN:HE22	1:A:146:VAL:CG1	2.30	0.44
1:B:180:PHE:HB3	1:B:182:MET:CE	2.47	0.44
1:B:102:ASN:HD22	1:B:103:GLN:N	2.16	0.44
1:B:102:ASN:ND2	1:B:104:ILE:H	2.16	0.43
1:A:183:TYR:C	1:A:183:TYR:CD2	2.92	0.43
1:A:98:VAL:HG21	1:A:142:GLN:CB	2.48	0.43
1:B:90:ILE:HG12	1:B:147:VAL:HG22	2.00	0.43
1:B:133:ASP:OD1	1:B:133:ASP:C	2.57	0.43
1:A:102:ASN:ND2	1:A:102:ASN:C	2.71	0.43
1:A:69:LEU:N	1:A:69:LEU:HD12	2.34	0.43
1:B:83:LEU:O	1:B:85:ASP:N	2.51	0.42
1:A:167:ASN:HD22	1:A:177:SER:CB	2.33	0.42
1:A:138:ARG:NE	1:A:173:GLU:OE2	2.52	0.42
1:A:136:PHE:O	2:A:391:MLC:HP81	2.19	0.42
1:A:127:ALA:HA	1:A:183:TYR:O	2.19	0.42
1:B:46:SER:OG	1:B:48:PRO:HG2	2.19	0.42
1:A:153:THR:O	1:A:154:ALA:HB2	2.20	0.42
1:B:98:VAL:HG22	1:B:103:GLN:C	2.39	0.41
1:B:176:PHE:CD2	1:B:176:PHE:C	2.93	0.41
1:B:117:LEU:O	1:B:121:VAL:HG22	2.21	0.41
1:A:106:ARG:HE	2:A:391:MLC:CM3	2.34	0.41
1:A:117:LEU:HD23	1:A:150:ALA:HB2	2.03	0.41
1:B:180:PHE:HB3	1:B:182:MET:HE3	2.03	0.40
1:A:103:GLN:N	1:A:103:GLN:CD	2.75	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	141/157 (90%)	132 (94%)	7 (5%)	2 (1%)	16	58
1	B	141/157 (90%)	125 (89%)	12 (8%)	4 (3%)	8	39
All	All	282/314 (90%)	257 (91%)	19 (7%)	6 (2%)	11	48

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	185	SER
1	B	84	ASP
1	B	158	GLU
1	B	185	SER
1	A	160	GLY
1	B	72	ASP

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	119/136 (88%)	104 (87%)	15 (13%)	7	24
1	B	123/136 (90%)	98 (80%)	25 (20%)	2	8
All	All	242/272 (89%)	202 (84%)	40 (16%)	3	12

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	LEU
1	A	98	VAL
1	A	102	ASN

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Mol	Chain	Res	Type
1	A	103	GLN
1	A	110	LEU
1	A	121	VAL
1	A	124	ASP
1	A	133	ASP
1	A	140	VAL
1	A	142	GLN
1	A	155	VAL
1	A	165	GLU
1	A	174	ILE
1	A	184	ARG
1	A	185	SER
1	B	45	LEU
1	B	51	ARG
1	B	52	GLU
1	B	55	LYS
1	B	66	VAL
1	B	69	LEU
1	B	70	SER
1	B	83	LEU
1	B	86	GLN
1	B	91	LEU
1	B	96	GLU
1	B	98	VAL
1	B	101	ARG
1	B	102	ASN
1	B	103	GLN
1	B	121	VAL
1	B	124	ASP
1	B	126	LEU
1	B	133	ASP
1	B	135	ARG
1	B	138	ARG
1	B	165	GLU
1	B	184	ARG
1	B	185	SER
1	B	186	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	ASN

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Mol	Chain	Res	Type
1	A	102	ASN
1	A	113	GLN
1	A	167	ASN
1	B	56	ASN
1	B	102	ASN
1	B	167	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MLC	A	391	-	31,31,56	1.22	2 (6%)	44,44,83	2.52	17 (38%)
2	MLC	B	390	-	31,31,56	1.97	4 (12%)	44,44,83	2.82	19 (43%)

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	MLC	A	391	-	-	0/40/40/71	0/0/0/3
2	MLC	B	390	-	-	2/40/40/71	0/0/0/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	390	MLC	CPA-CP7	-8.23	1.46	1.55
2	A	391	MLC	P1-O5'	4.06	1.64	1.50
2	A	391	MLC	CPA-CP7	-3.74	1.51	1.55
2	B	390	MLC	P1-O5'	3.68	1.63	1.50
2	B	390	MLC	CM2-CM1	-3.46	1.47	1.51
2	B	390	MLC	CM2-CM3	-2.46	1.47	1.51

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	390	MLC	CP4-CP5-NP2	-8.77	92.98	111.87
2	A	391	MLC	CP4-CP5-NP2	-7.63	95.43	111.87
2	B	390	MLC	CP5-NP2-CP6	7.25	137.47	122.57
2	A	391	MLC	P1-O6-P2	-5.69	108.53	132.95
2	A	391	MLC	CP5-NP2-CP6	5.39	133.64	122.57
2	B	390	MLC	CPA-CP7-CP6	-5.21	107.68	112.73
2	B	390	MLC	CPB-CPA-CP7	-5.03	101.34	108.70
2	B	390	MLC	CP7-CP6-NP2	-4.41	107.14	116.57
2	A	391	MLC	CP9-CPA-CPB	4.14	114.72	108.76
2	B	390	MLC	OP2-CP6-CP7	4.12	127.98	120.48
2	A	391	MLC	O5'-P1-O6	-3.85	94.76	106.65
2	B	390	MLC	OP3-CP7-CP6	3.81	120.59	110.76
2	A	391	MLC	OP2-CP6-CP7	3.61	127.07	120.48
2	A	391	MLC	CM2-CM1-S	-3.46	109.68	113.43
2	B	390	MLC	CP9-CPA-CPB	3.39	113.65	108.76
2	B	390	MLC	O6-P2-O7	-3.38	88.31	103.41
2	A	391	MLC	CP7-CP6-NP2	-3.36	109.37	116.57
2	A	391	MLC	CP8-CPA-CPB	-3.18	104.17	108.76
2	A	391	MLC	O7-CPB-CPA	3.16	115.86	110.57
2	B	390	MLC	P1-O6-P2	-3.11	119.59	132.95
2	A	391	MLC	CP2-NP1-CP3	3.03	129.01	122.84
2	B	390	MLC	OP3-CP7-CPA	-2.94	103.88	110.71
2	A	391	MLC	OM4-CM3-CM2	-2.77	112.59	121.95
2	A	391	MLC	CP5-CP4-CP3	2.74	116.92	112.25
2	A	391	MLC	OM2-CM1-CM2	2.71	126.98	122.75
2	B	390	MLC	O12-P1-O11	2.56	120.08	110.82
2	A	391	MLC	OP3-CP7-CP6	2.45	117.07	110.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	390	MLC	CP8-CPA-CP9	2.34	112.41	109.08
2	B	390	MLC	CM2-CM1-S	2.28	115.91	113.43
2	B	390	MLC	OM4-CM3-CM2	-2.27	114.29	121.95
2	B	390	MLC	O12-P1-O6	2.21	111.33	104.29
2	B	390	MLC	OM2-CM1-CM2	-2.14	119.40	122.75
2	B	390	MLC	OM3-CM3-OM4	2.09	128.61	123.30
2	B	390	MLC	O22-P2-O6	2.04	114.81	105.14
2	A	391	MLC	O22-P2-O21	2.03	123.54	112.21
2	A	391	MLC	CPA-CP7-CP6	-2.02	110.77	112.73

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	390	MLC	CM2-CM1-S-CP1
2	B	390	MLC	OM2-CM1-S-CP1

There are no ring outliers.

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 ⓘ

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	143/157 (91%)	-0.04	0 100 100	51, 59, 73, 87	0
1	B	143/157 (91%)	-0.02	0 100 100	49, 62, 80, 94	0
All	All	286/314 (91%)	-0.03	0 100 100	49, 60, 78, 94	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MLC	A	391	32/54	0.21	-0.94	59,66,102,104	0
2	MLC	B	390	32/54	0.19	-1.11	58,66,86,87	0

6.5 Other polymers ⓘ

There are no such residues in this entry.