Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:16 PM GMT

PDB ID	:	4L4Z
Title	:	Crystal structures of the LsrR proteins complexed with phospho-AI-2 and its two different analogs reveal distinct mechanisms for ligand recognition
Authors	:	Ryu, K.S.; Ha, J.H.; Eo, Y.
Deposited on	:	2013-06-10
Resolution	:	2.30 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

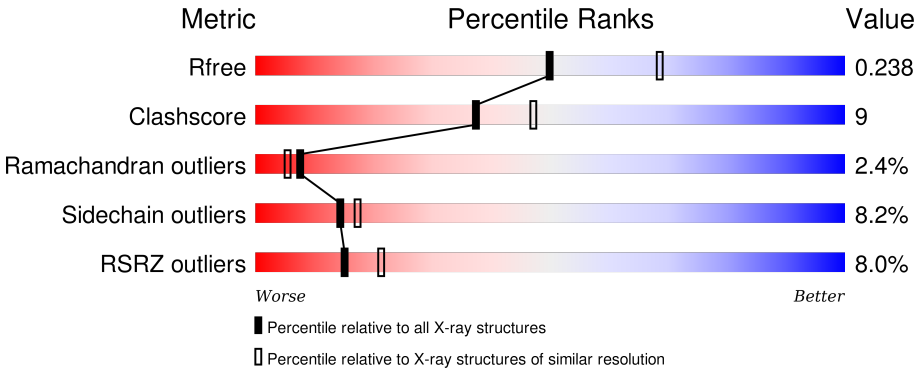
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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 <=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	266	8% 73% 17% • • 6%
1	B	266	8% 79% 13% • • 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3877 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.

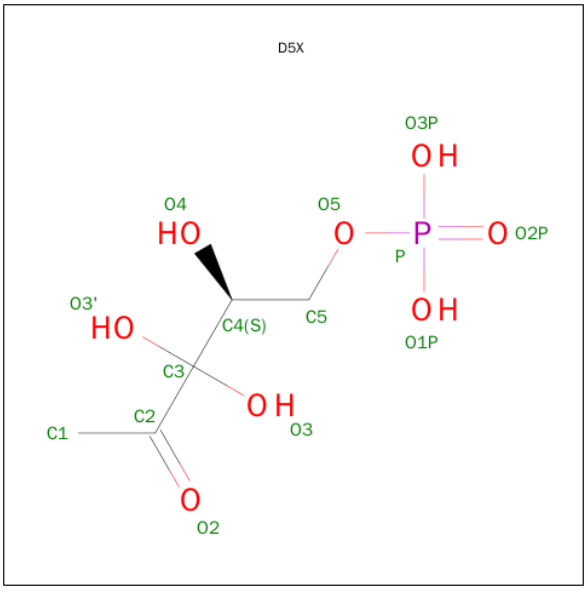
- Molecule 1 is a protein called Transcriptional regulator LsrR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			1820	1138	324	348	10			
1	B	250	Total	C	N	O	S	0	4	0
			1856	1164	330	352	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	52	SER	-	EXPRESSION TAG	UNP P76141
B	52	SER	-	EXPRESSION TAG	UNP P76141

- Molecule 2 is (2S)-2,3,3-TRIHYDROXY-4-OXOPENTYL DIHYDROGEN PHOSPHATE (three-letter code: D5X) (formula: $C_5H_{11}O_8P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			14	5	8	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total C O P 14 5 8 1	0	0

- Molecule 3 is water.

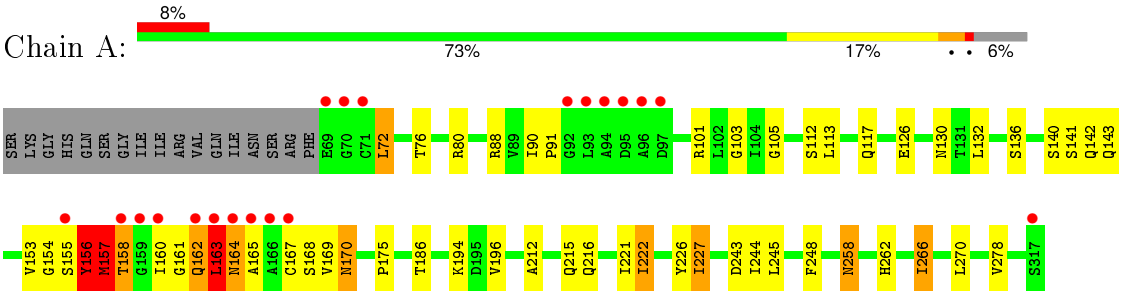
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	74	Total O 74 74	0	0
3	B	99	Total O 99 99	0	0

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.

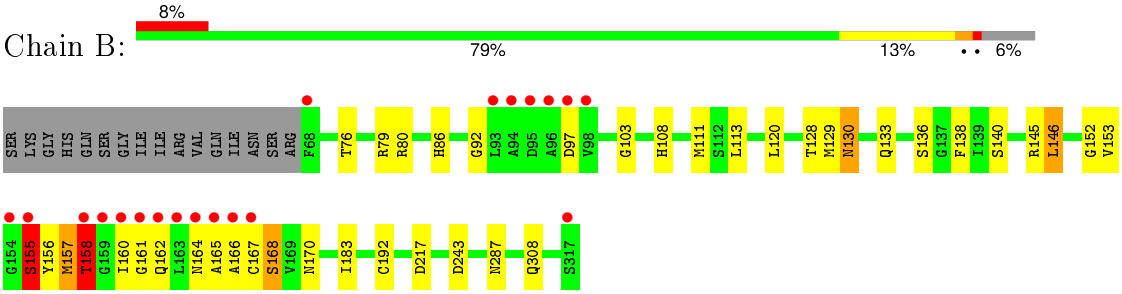
- Molecule 1: Transcriptional regulator LsrR

Chain A:



- Molecule 1: Transcriptional regulator LsrR

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α, β, γ	117.69Å 117.69Å 79.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.68 – 2.30 34.66 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.5 (34.68-2.30) 96.5 (34.66-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.47 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.172 , 0.233 0.180 , 0.238	Depositor DCC
R_{free} test set	1400 reflections (5.49%)	DCC
Wilson B-factor (Å^2)	32.3	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 45.4	EDS
Estimated twinning fraction	0.057 for h,-h-k,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 26859 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3877	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.60% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: D5X

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	0.76	0/1838	0.92	1/2483 (0.0%)
1	B	0.78	0/1887	0.90	0/2547
All	All	0.77	0/3725	0.91	1/5030 (0.0%)

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	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	157	MET	N-CA-C	-9.00	86.71	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	156	TYR	Peptide
1	A	157	MET	Peptide
1	B	152	GLY	Peptide

5.2 Too-close contacts (i)

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	1820	0	1874	40	0
1	B	1856	0	1921	32	0
2	A	14	0	11	0	0
2	B	14	0	11	0	0
3	A	74	0	0	2	0
3	B	99	0	0	4	0
All	All	3877	0	3817	70	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 9.

All (70) 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:258:ASN:HD22	1:A:258:ASN:H	1.09	0.99
1:B:164:ASN:HB2	1:B:165:ALA:HA	1.53	0.89
1:B:164:ASN:HB2	1:B:167:CYS:SG	2.15	0.86
1:B:157:MET:HA	1:B:160:ILE:HG13	1.63	0.81
1:A:258:ASN:ND2	1:A:258:ASN:H	1.77	0.79
1:B:164:ASN:CB	1:B:165:ALA:HA	2.12	0.78
1:B:153:VAL:O	1:B:156:TYR:HB2	1.86	0.74
1:B:129:MET:HB3	1:B:158:THR:HG21	1.72	0.71
1:A:153:VAL:O	1:A:156:TYR:HB2	1.93	0.68
1:A:143:GLN:NE2	1:A:165:ALA:O	2.27	0.68
1:B:133:GLN:HE21	1:B:158:THR:HG23	1.59	0.67
1:A:143:GLN:HA	1:A:167:CYS:SG	2.35	0.66
1:A:226:TYR:OH	3:A:635:HOH:O	2.10	0.66
1:B:108:HIS:HB2	3:B:574:HOH:O	1.96	0.65
1:B:165:ALA:HA	1:B:167:CYS:SG	2.42	0.60
1:A:258:ASN:N	1:A:258:ASN:ND2	2.47	0.60
1:A:221:ILE:HG23	1:A:227:ILE:CD1	2.32	0.59
1:B:157:MET:CA	1:B:160:ILE:HG13	2.31	0.59
1:A:162:GLN:CD	1:A:163:LEU:H	2.06	0.58
1:A:154:GLY:O	1:A:156:TYR:O	2.22	0.57
1:B:155:SER:H	1:B:157:MET:HG2	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:HIS:CD2	3:B:539:HOH:O	2.57	0.56
1:B:164:ASN:CB	1:B:165:ALA:CA	2.84	0.56
1:B:111:MET:HG2	1:B:138:PHE:CD2	2.41	0.56
1:A:155:SER:C	1:A:156:TYR:O	2.43	0.55
1:A:162:GLN:OE1	1:A:163:LEU:O	2.25	0.55
1:A:258:ASN:HD22	1:A:258:ASN:N	1.83	0.55
1:B:164:ASN:HB3	1:B:166:ALA:N	2.23	0.54
1:A:154:GLY:O	1:A:158:THR:HG21	2.08	0.54
1:A:170:ASN:OD1	1:A:196:VAL:HG22	2.08	0.54
1:B:160:ILE:HD12	1:B:160:ILE:C	2.29	0.53
1:A:76:THR:O	1:A:80:ARG:HG3	2.09	0.52
1:B:156:TYR:O	1:B:157:MET:HG2	2.09	0.52
1:A:212:ALA:H	1:A:215:GLN:HE21	1.58	0.52
1:A:164:ASN:O	1:A:165:ALA:HB3	2.09	0.51
1:B:103:GLY:HA3	1:B:130[A]:ASN:HD22	1.76	0.51
1:A:163:LEU:HD11	1:A:169:VAL:HG23	1.94	0.50
1:B:92:GLY:H	1:B:308:GLN:HE21	1.60	0.50
1:A:154:GLY:HA3	3:B:536:HOH:O	2.11	0.49
1:A:155:SER:HA	3:A:657:HOH:O	2.11	0.49
1:B:86:HIS:CD2	1:B:113:LEU:HD21	2.47	0.49
1:B:161:GLY:O	1:B:162:GLN:HG2	2.13	0.48
1:A:113:LEU:HD13	1:A:278:VAL:HG11	1.94	0.48
1:A:162:GLN:O	1:A:163:LEU:HB3	2.14	0.48
1:A:163:LEU:HG	1:A:164:ASN:N	2.29	0.48
1:B:133:GLN:HE21	1:B:158:THR:CG2	2.26	0.47
1:A:248:PHE:H	1:A:262:HIS:CE1	2.32	0.47
1:A:175:PRO:HB3	1:B:192:CYS:SG	2.56	0.46
1:B:76:THR:O	1:B:80:ARG:HG3	2.16	0.46
1:B:164:ASN:CB	1:B:167:CYS:SG	2.97	0.46
1:A:163:LEU:O	1:A:164:ASN:HB2	2.14	0.46
1:B:133:GLN:NE2	1:B:158:THR:HG23	2.30	0.46
1:A:244:ILE:HG22	1:A:245:LEU:HG	1.97	0.46
1:B:160:ILE:HD12	1:B:161:GLY:N	2.31	0.46
1:B:145:ARG:HG2	1:B:168:SER:HB3	1.98	0.45
1:A:72:LEU:HD13	1:A:72:LEU:C	2.36	0.45
1:A:103:GLY:HA3	1:A:130:ASN:ND2	2.31	0.45
1:A:222:ILE:HD12	1:A:227:ILE:CD1	2.47	0.45
1:A:117:GLN:HA	1:A:143:GLN:O	2.17	0.45
1:A:186:THR:HG21	1:B:183:ILE:HG12	2.00	0.44
1:B:76:THR:HA	1:B:79:ARG:NH1	2.33	0.43
1:A:222:ILE:HD12	1:A:227:ILE:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:GLN:O	1:A:216:GLN:C	2.56	0.43
1:A:266:ILE:N	1:A:266:ILE:CD1	2.83	0.42
1:B:120:LEU:HB3	1:B:146:LEU:HD12	2.01	0.42
1:A:161:GLY:O	1:A:162:GLN:CB	2.67	0.42
1:A:163:LEU:HD12	1:A:167:CYS:HB3	2.03	0.41
1:A:221:ILE:HG23	1:A:227:ILE:HD11	2.03	0.41
1:B:128:THR:HG23	3:B:508:HOH:O	2.20	0.40
1:A:90:ILE:HD11	1:A:105:GLY:HA3	2.03	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	247/266 (93%)	230 (93%)	10 (4%)	7 (3%)	6	4
1	B	252/266 (95%)	234 (93%)	13 (5%)	5 (2%)	9	7
All	All	499/532 (94%)	464 (93%)	23 (5%)	12 (2%)	7	5

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	156	TYR
1	A	160	ILE
1	A	163	LEU
1	A	164	ASN
1	B	155	SER
1	B	157	MET
1	A	162	GLN
1	B	158	THR
1	A	158	THR
1	B	168	SER

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Mol	Chain	Res	Type
1	B	97	ASP
1	A	91	PRO

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	188/203 (93%)	167 (89%)	21 (11%)	7 8
1	B	193/203 (95%)	182 (94%)	11 (6%)	25 34
All	All	381/406 (94%)	349 (92%)	32 (8%)	14 16

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	LEU
1	A	88	ARG
1	A	101	ARG
1	A	112	SER
1	A	126	GLU
1	A	132	LEU
1	A	136	SER
1	A	140	SER
1	A	141	SER
1	A	142	GLN
1	A	157	MET
1	A	163	LEU
1	A	168	SER
1	A	170	ASN
1	A	194	LYS
1	A	222	ILE
1	A	227	ILE
1	A	243	ASP
1	A	258	ASN
1	A	266	ILE
1	A	270	LEU

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Mol	Chain	Res	Type
1	B	130[A]	ASN
1	B	130[B]	ASN
1	B	136	SER
1	B	140	SER
1	B	146	LEU
1	B	155	SER
1	B	158	THR
1	B	170	ASN
1	B	217	ASP
1	B	243	ASP
1	B	287	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	GLN
1	A	130	ASN
1	A	142	GLN
1	A	191	ASN
1	A	201	GLN
1	A	215	GLN
1	A	258	ASN
1	A	262	HIS
1	B	133	GLN
1	B	142	GLN
1	B	164	ASN
1	B	191	ASN
1	B	287	ASN
1	B	308	GLN

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 (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

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	D5X	A	501	-	8,13,13	1.60	3 (37%)	11,20,20	2.28	4 (36%)
2	D5X	B	401	-	8,13,13	1.27	1 (12%)	11,20,20	1.68	3 (27%)

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	D5X	A	501	-	-	0/6/18/18	0/0/0/0
2	D5X	B	401	-	-	0/6/18/18	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	D5X	P-O5	-2.85	1.50	1.60
2	A	501	D5X	C1-C2	2.15	1.56	1.49
2	A	501	D5X	P-O3P	2.22	1.62	1.54
2	B	401	D5X	P-O3P	2.30	1.63	1.54

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	D5X	C5-C4-C3	-2.95	107.14	112.06
2	B	401	D5X	O5-P-O2P	-2.44	100.93	107.14
2	B	401	D5X	O1P-P-O3P	2.50	116.89	107.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	D5X	O4-C4-C3	2.96	115.27	110.20
2	A	501	D5X	O3P-P-O5	3.59	116.89	106.56
2	B	401	D5X	O2-C2-C3	3.65	123.52	119.44
2	A	501	D5X	O2-C2-C3	4.46	124.42	119.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 95^{th} 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	249/266 (93%)	0.18	20 (8%)	15	21	22, 33, 102, 126	0
1	B	250/266 (93%)	0.19	20 (8%)	15	21	19, 32, 99, 139	0
All	All	499/532 (93%)	0.18	40 (8%)	15	21	19, 33, 102, 139	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	93	LEU	13.2
1	B	158	THR	11.3
1	B	95	ASP	11.3
1	B	94	ALA	10.9
1	A	165	ALA	10.6
1	A	96	ALA	9.6
1	B	96	ALA	8.1
1	A	95	ASP	6.3
1	B	159	GLY	6.1
1	B	162	GLN	5.8
1	B	166	ALA	5.8
1	A	94	ALA	5.6
1	B	167	CYS	5.6
1	B	160	ILE	5.3
1	B	317	SER	5.3
1	A	160	ILE	5.2
1	A	92	GLY	5.2
1	B	165	ALA	4.7
1	B	97	ASP	4.3
1	A	70	GLY	4.3
1	B	154	GLY	4.2
1	A	167	CYS	4.1
1	A	163	LEU	4.1
1	A	166	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	97	ASP	4.0
1	A	317	SER	3.8
1	B	68	PHE	3.7
1	A	164	ASN	3.6
1	B	164	ASN	3.2
1	A	159	GLY	3.1
1	A	71	CYS	3.0
1	B	155	SER	2.8
1	B	98	VAL	2.7
1	A	162	GLN	2.7
1	A	69	GLU	2.7
1	B	93	LEU	2.7
1	A	155	SER	2.5
1	B	163	LEU	2.5
1	B	161	GLY	2.3
1	A	158	THR	2.1

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 carbohydrates 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. 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, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
2	D5X	B	401	14/14	0.98	0.14	-0.15	23,40,57,59	0
2	D5X	A	501	14/14	0.99	0.12	-0.46	26,38,58,61	0

6.5 Other polymers (i)

There are no such residues in this entry.