



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 13, 2018 – 06:59 pm GMT

PDB ID : 1Y1R
Title : Crystal Structure of the Uridine Phosphorylase from Salmonella Typhimurium
in Complex with Inhibitor and Phosphate Ion at 2.11Å Resolution
Authors : Dontsova, M.V.; Gabdoulkhakov, A.G.; Kachalova, G.S.; Betzel, C.; Ealick,
S.E.; Mikhailov, A.M.
Deposited on : 2004-11-19
Resolution : 2.11 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix)	:	1.13
EDS	:	trunk31020
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk31020

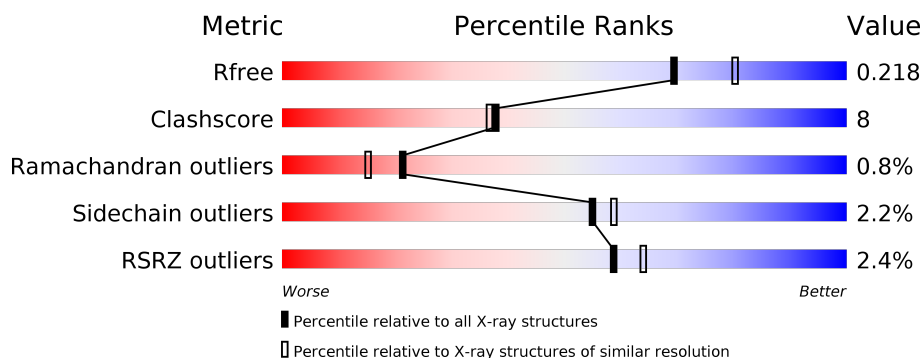
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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}	111664	5449 (2.14-2.10)
Clashscore	122126	5972 (2.14-2.10)
Ramachandran outliers	120053	5914 (2.14-2.10)
Sidechain outliers	120020	5915 (2.14-2.10)
RSRZ outliers	108989	5321 (2.14-2.10)

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.

Mol	Chain	Length	Quality of chain
1	A	253	<div> <div>3%</div> <div> <div></div> <div>79%</div> <div>18%</div> <div>• •</div> </div> </div>
1	B	253	<div> <div>80%</div> <div>14%</div> <div>• •</div> </div>
1	C	253	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>• •</div> </div> </div>
1	D	253	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>20%</div> <div>• 5%</div> </div> </div>
1	E	253	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>5%</div> </div> </div>
1	F	253	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>16%</div> <div>• 5%</div> </div> </div>

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
3	BJE	D	706	-	-	-	X

2 Entry composition [i](#)

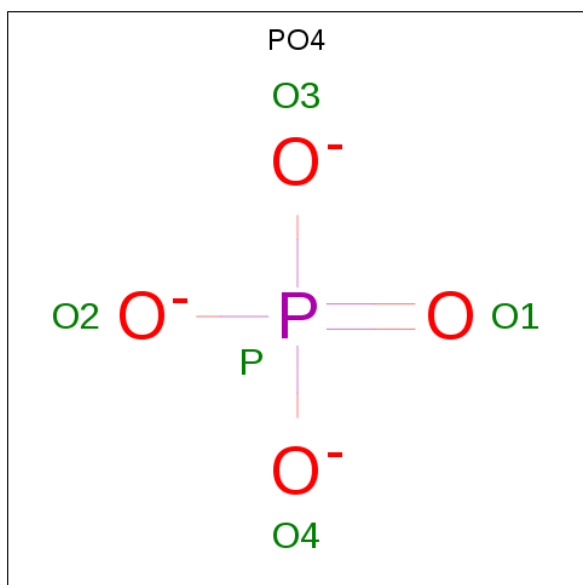
There are 4 unique types of molecules in this entry. The entry contains 11625 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Uridine phosphorylase.

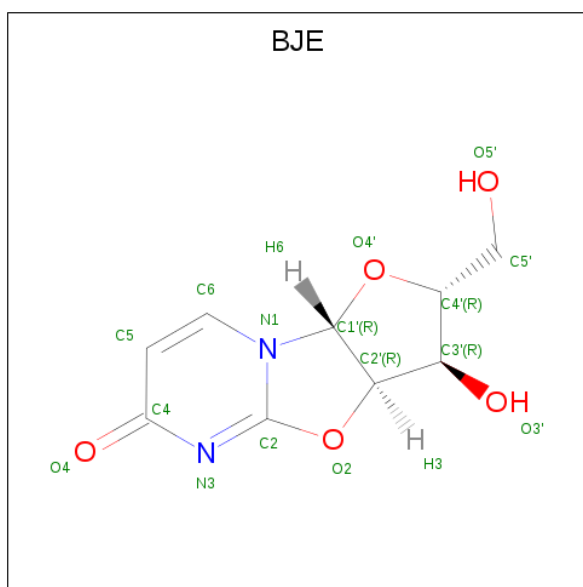
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	0	0
			1886	1180	332	362	12			
1	C	243	Total	C	N	O	S	0	0	0
			1824	1143	322	347	12			
1	D	241	Total	C	N	O	S	0	0	0
			1806	1132	318	344	12			
1	F	240	Total	C	N	O	S	0	0	0
			1800	1128	318	343	11			
1	E	241	Total	C	N	O	S	0	0	0
			1807	1132	319	345	11			
1	B	242	Total	C	N	O	S	0	0	0
			1814	1137	320	346	11			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total O P 5 4 1	0	0
2	F	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0

- Molecule 3 is 2,2'-Anhydro-(1-beta-D-ribofuranosyl)uracil (three-letter code: BJE) (formula: $C_9H_{10}N_2O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total C N O 16 9 2 5	0	0
3	B	1	Total C N O 16 9 2 5	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	122	Total O 122 122	0	0
4	C	109	Total O 109 109	0	0
4	D	82	Total O 82 82	0	0
4	F	95	Total O 95 95	0	0

Continued on next page...

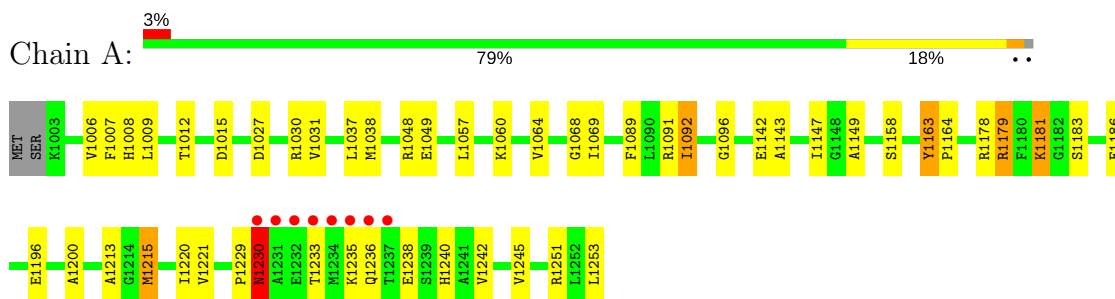
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	117	Total 117	O 117	0	0
4	B	116	Total 116	O 116	0	0

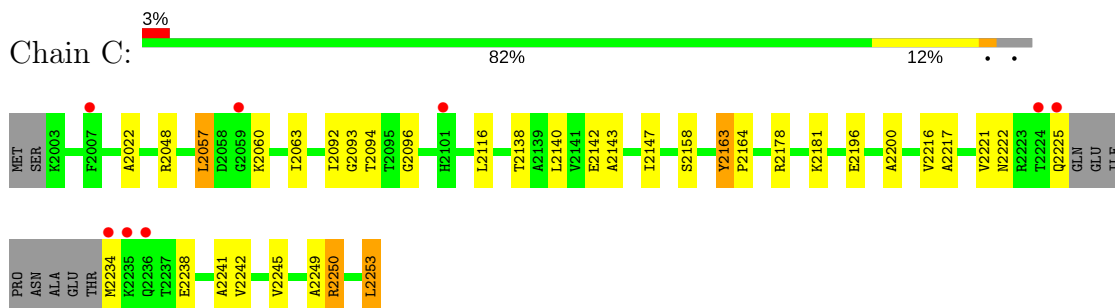
3 Residue-property plots

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

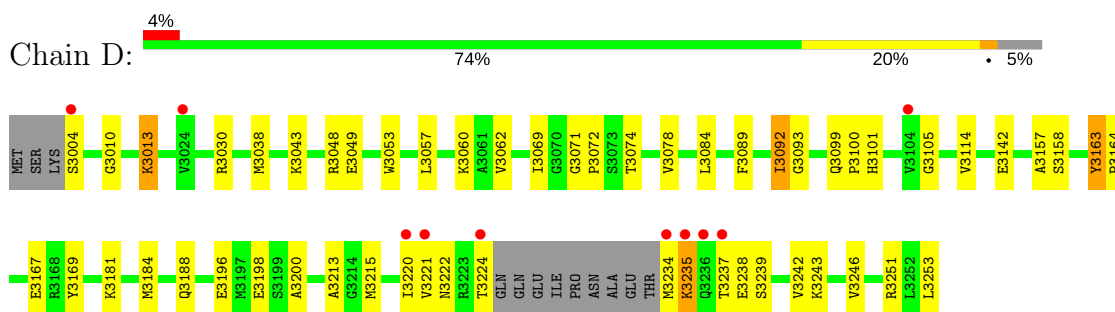
• Molecule 1: Uridine phosphorylase



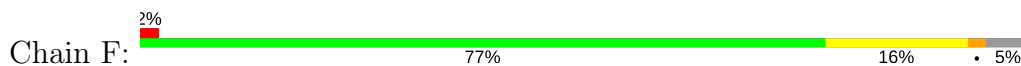
• Molecule 1: Uridine phosphorylase

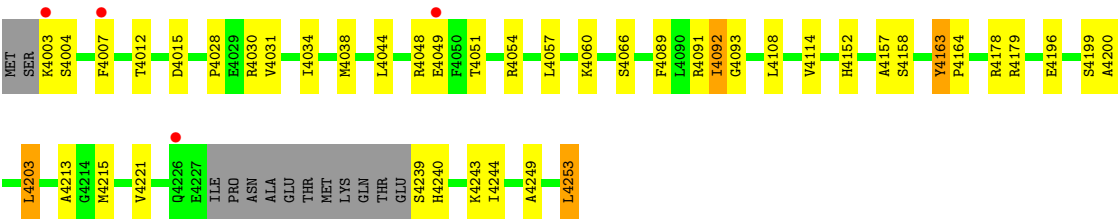


• Molecule 1: Uridine phosphorylase

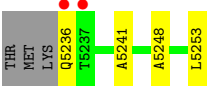
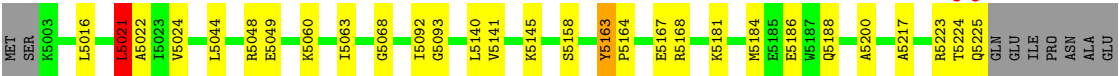
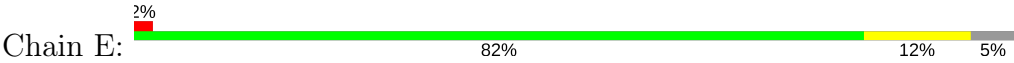


• Molecule 1: Uridine phosphorylase

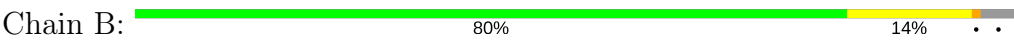




● Molecule 1: Uridine phosphorylase



● Molecule 1: Uridine phosphorylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.88Å 124.07Å 134.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.75 – 2.11 28.75 – 2.09	Depositor EDS
% Data completeness (in resolution range)	97.5 (28.75-2.11) 98.3 (28.75-2.09)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.180 , 0.219 0.180 , 0.218	Depositor DCC
R_{free} test set	4192 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	12.0	Xtriage
Anisotropy	0.386	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 60.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11625	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: PO4, BJE

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	0.39	0/1916	0.69	0/2595
1	B	0.40	0/1843	0.65	0/2497
1	C	0.38	0/1852	0.64	0/2505
1	D	0.37	0/1834	0.63	0/2482
1	E	0.39	0/1835	0.67	1/2484 (0.0%)
1	F	0.39	0/1828	0.65	0/2474
All	All	0.39	0/11108	0.66	1/15037 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	5021	LEU	CA-CB-CG	5.32	127.53	115.30

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	1886	0	1900	40	0
1	B	1814	0	1825	31	0
1	C	1824	0	1843	26	0
1	D	1806	0	1822	46	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1807	0	1821	24	0
1	F	1800	0	1814	32	0
2	B	5	0	0	0	0
2	D	5	0	0	0	0
2	F	5	0	0	0	0
3	B	16	0	0	1	0
3	D	16	0	0	2	0
4	A	122	0	0	2	0
4	B	116	0	0	2	0
4	C	109	0	0	1	0
4	D	82	0	0	0	0
4	E	117	0	0	1	0
4	F	95	0	0	1	0
All	All	11625	0	11025	183	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 8.

The worst 5 of 183 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:1229:PRO:HB2	1:A:1233:THR:HG22	1.21	1.13
1:D:3049:GLU:HG2	1:B:6049:GLU:HG2	1.27	1.08
1:D:3100:PRO:HG3	1:D:3224:THR:HG21	1.52	0.92
1:F:4091:ARG:HB3	1:F:4215:MET:HG3	1.53	0.89
1:A:1229:PRO:HB2	1:A:1233:THR:CG2	2.03	0.87

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	249/253 (98%)	237 (95%)	9 (4%)	3 (1%)	14	8
1	B	238/253 (94%)	232 (98%)	5 (2%)	1 (0%)	36	34
1	C	239/253 (94%)	234 (98%)	4 (2%)	1 (0%)	36	34
1	D	237/253 (94%)	231 (98%)	4 (2%)	2 (1%)	21	15
1	E	237/253 (94%)	231 (98%)	4 (2%)	2 (1%)	21	15
1	F	236/253 (93%)	226 (96%)	8 (3%)	2 (1%)	21	15
All	All	1436/1518 (95%)	1391 (97%)	34 (2%)	11 (1%)	21	15

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1230	ASN
1	A	1235	LYS
1	A	1163	TYR
1	D	3163	TYR
1	D	3235	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	200/202 (99%)	194 (97%)	6 (3%)	44	46
1	B	192/202 (95%)	187 (97%)	5 (3%)	49	52
1	C	193/202 (96%)	189 (98%)	4 (2%)	56	60
1	D	191/202 (95%)	187 (98%)	4 (2%)	56	60
1	E	191/202 (95%)	189 (99%)	2 (1%)	78	83
1	F	190/202 (94%)	186 (98%)	4 (2%)	56	60
All	All	1157/1212 (96%)	1132 (98%)	25 (2%)	55	59

5 of 25 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	3030	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	3215	MET
1	B	6196	GLU
1	D	3092	ILE
1	F	4092	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	5014	ASN
1	E	5020	GLN
1	B	6225	GLN
1	F	4226	GLN
1	B	6188	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

5 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	PO4	B	705	-	4,4,4	1.23	0	6,6,6	0.38	0
3	BJE	B	707	-	14,18,18	2.05	5 (35%)	15,27,27	4.37	8 (53%)
2	PO4	D	703	-	4,4,4	1.10	0	6,6,6	0.37	0
3	BJE	D	706	-	14,18,18	2.10	5 (35%)	15,27,27	4.63	8 (53%)
2	PO4	F	704	-	4,4,4	1.09	0	6,6,6	0.40	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	PO4	B	705	-	-	0/0/0/0	0/0/0/0
3	BJE	B	707	-	-	0/2/26/26	0/3/3/3
2	PO4	D	703	-	-	0/0/0/0	0/0/0/0
3	BJE	D	706	-	-	0/2/26/26	0/3/3/3
2	PO4	F	704	-	-	0/0/0/0	0/0/0/0

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	707	BJE	O2-C2'	-2.13	1.43	1.45
3	D	706	BJE	O2-C2'	-2.05	1.43	1.45
3	B	707	BJE	O4'-C1'	2.06	1.44	1.41
3	D	706	BJE	O4'-C1'	2.19	1.44	1.41
3	D	706	BJE	C6-N1	3.13	1.40	1.35

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	707	BJE	O2-C2'-C1'	-6.30	97.89	105.47
3	D	706	BJE	O2-C2'-C1'	-4.96	99.50	105.47
3	D	706	BJE	O4'-C1'-C2'	-4.65	98.51	106.60
3	B	707	BJE	O4'-C1'-C2'	-4.14	99.40	106.60
3	D	706	BJE	C5'-C4'-C3'	-3.48	106.62	115.06

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	707	BJE	1	0
3	D	706	BJE	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	251/253 (99%)	-0.01	8 (3%) 47 54	7, 15, 58, 99	0
1	B	242/253 (95%)	-0.20	1 (0%) 92 93	9, 15, 36, 55	0
1	C	243/253 (96%)	0.07	8 (3%) 46 53	8, 18, 42, 85	0
1	D	241/253 (95%)	0.06	10 (4%) 37 44	9, 17, 48, 89	0
1	E	241/253 (95%)	-0.13	4 (1%) 70 74	7, 15, 41, 61	0
1	F	240/253 (94%)	-0.05	4 (1%) 70 74	8, 18, 41, 70	0
All	All	1458/1518 (96%)	-0.04	35 (2%) 59 64	7, 16, 43, 99	0

The worst 5 of 35 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1234	MET	8.6
1	A	1231	ALA	8.4
1	D	3234	MET	7.9
1	A	1233	THR	6.3
1	D	3235	LYS	5.3

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. 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	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BJE	D	706	16/16	0.70	0.48	20,20,20,20	0
3	BJE	B	707	16/16	0.79	0.18	20,20,20,20	0
2	PO4	D	703	5/5	0.94	0.28	20,20,20,20	0
2	PO4	F	704	5/5	0.96	0.24	20,20,20,20	0
2	PO4	B	705	5/5	0.97	0.12	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.