



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 12, 2018 – 07:22 PM EST

PDB ID : 6AZN  
Title : Structural and biochemical characterization of a non-canonical biuret hydrolase (BiuH) from the cyanuric acid catabolism pathway of *Rhizobium leguminosorum* bv. *viciae* 3841  
Authors : Peat, T.S.; Esquirol, L.; Newman, J.; Scott, C.  
Deposited on : 2017-09-11  
Resolution : 1.75 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](mailto: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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030736

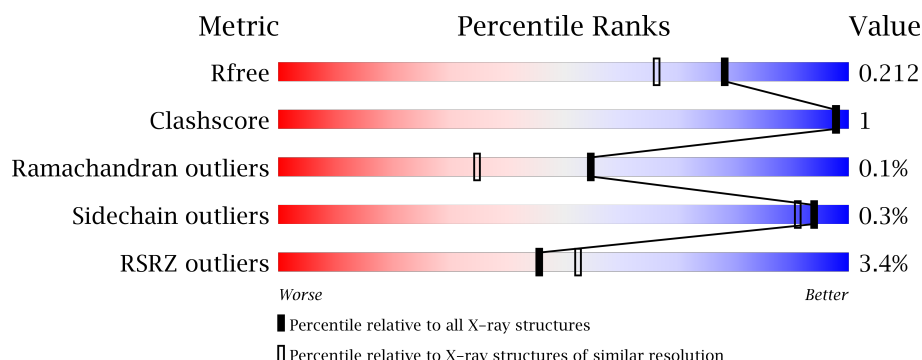
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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))
$R_{free}$	100719	1762 (1.76-1.76)
Clashscore	112137	1889 (1.76-1.76)
Ramachandran outliers	110173	1868 (1.76-1.76)
Sidechain outliers	110143	1868 (1.76-1.76)
RSRZ outliers	101464	1770 (1.76-1.76)

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  $\geq 3$ , 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	262	<div> <div>0.1%</div> <div>82%</div> <div>14%</div> </div>
1	B	262	<div> <div>5%</div> <div>80%</div> <div>15%</div> </div>
1	C	262	<div> <div>0.1%</div> <div>82%</div> <div>14%</div> </div>
1	D	262	<div> <div>3%</div> <div>82%</div> <div>14%</div> </div>
1	F	262	<div> <div>3%</div> <div>81%</div> <div>14%</div> </div>

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	G	262	
1	H	262	
1	I	262	

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
2	EDO	A	301	-	-	-	X
2	EDO	I	302	-	-	-	X
3	PO4	A	302	-	-	-	X
3	PO4	F	301	-	-	-	X
3	PO4	I	303	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15574 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 Putative amidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	5	0
			1780	1116	318	333	13			
1	B	223	Total	C	N	O	S	0	3	0
			1721	1079	303	327	12			
1	C	226	Total	C	N	O	S	0	4	0
			1766	1111	313	329	13			
1	D	226	Total	C	N	O	S	0	5	0
			1776	1115	316	332	13			
1	F	226	Total	C	N	O	S	0	4	0
			1768	1110	314	331	13			
1	G	226	Total	C	N	O	S	0	2	0
			1748	1099	309	327	13			
1	H	229	Total	C	N	O	S	0	2	0
			1766	1109	313	331	13			
1	I	227	Total	C	N	O	S	0	4	0
			1773	1115	312	333	13			

There are 240 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q1M7F4
A	-18	GLY	-	expression tag	UNP Q1M7F4
A	-17	SER	-	expression tag	UNP Q1M7F4
A	-16	SER	-	expression tag	UNP Q1M7F4
A	-15	HIS	-	expression tag	UNP Q1M7F4
A	-14	HIS	-	expression tag	UNP Q1M7F4
A	-13	HIS	-	expression tag	UNP Q1M7F4
A	-12	HIS	-	expression tag	UNP Q1M7F4
A	-11	HIS	-	expression tag	UNP Q1M7F4
A	-10	HIS	-	expression tag	UNP Q1M7F4
A	-9	SER	-	expression tag	UNP Q1M7F4
A	-8	SER	-	expression tag	UNP Q1M7F4
A	-7	GLY	-	expression tag	UNP Q1M7F4

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	expression tag	UNP Q1M7F4
A	-5	VAL	-	expression tag	UNP Q1M7F4
A	-4	PRO	-	expression tag	UNP Q1M7F4
A	-3	ARG	-	expression tag	UNP Q1M7F4
A	-2	GLY	-	expression tag	UNP Q1M7F4
A	-1	SER	-	expression tag	UNP Q1M7F4
A	0	HIS	-	expression tag	UNP Q1M7F4
A	175	SER	CYS	conflict	UNP Q1M7F4
A	234	GLY	-	expression tag	UNP Q1M7F4
A	235	LEU	-	expression tag	UNP Q1M7F4
A	236	VAL	-	expression tag	UNP Q1M7F4
A	237	PRO	-	expression tag	UNP Q1M7F4
A	238	ARG	-	expression tag	UNP Q1M7F4
A	239	GLY	-	expression tag	UNP Q1M7F4
A	240	SER	-	expression tag	UNP Q1M7F4
A	241	ILE	-	expression tag	UNP Q1M7F4
A	242	GLU	-	expression tag	UNP Q1M7F4
B	-19	MET	-	initiating methionine	UNP Q1M7F4
B	-18	GLY	-	expression tag	UNP Q1M7F4
B	-17	SER	-	expression tag	UNP Q1M7F4
B	-16	SER	-	expression tag	UNP Q1M7F4
B	-15	HIS	-	expression tag	UNP Q1M7F4
B	-14	HIS	-	expression tag	UNP Q1M7F4
B	-13	HIS	-	expression tag	UNP Q1M7F4
B	-12	HIS	-	expression tag	UNP Q1M7F4
B	-11	HIS	-	expression tag	UNP Q1M7F4
B	-10	HIS	-	expression tag	UNP Q1M7F4
B	-9	SER	-	expression tag	UNP Q1M7F4
B	-8	SER	-	expression tag	UNP Q1M7F4
B	-7	GLY	-	expression tag	UNP Q1M7F4
B	-6	LEU	-	expression tag	UNP Q1M7F4
B	-5	VAL	-	expression tag	UNP Q1M7F4
B	-4	PRO	-	expression tag	UNP Q1M7F4
B	-3	ARG	-	expression tag	UNP Q1M7F4
B	-2	GLY	-	expression tag	UNP Q1M7F4
B	-1	SER	-	expression tag	UNP Q1M7F4
B	0	HIS	-	expression tag	UNP Q1M7F4
B	175	SER	CYS	conflict	UNP Q1M7F4
B	234	GLY	-	expression tag	UNP Q1M7F4
B	235	LEU	-	expression tag	UNP Q1M7F4
B	236	VAL	-	expression tag	UNP Q1M7F4
B	237	PRO	-	expression tag	UNP Q1M7F4

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
B	238	ARG	-	expression tag	UNP Q1M7F4
B	239	GLY	-	expression tag	UNP Q1M7F4
B	240	SER	-	expression tag	UNP Q1M7F4
B	241	ILE	-	expression tag	UNP Q1M7F4
B	242	GLU	-	expression tag	UNP Q1M7F4
C	-19	MET	-	initiating methionine	UNP Q1M7F4
C	-18	GLY	-	expression tag	UNP Q1M7F4
C	-17	SER	-	expression tag	UNP Q1M7F4
C	-16	SER	-	expression tag	UNP Q1M7F4
C	-15	HIS	-	expression tag	UNP Q1M7F4
C	-14	HIS	-	expression tag	UNP Q1M7F4
C	-13	HIS	-	expression tag	UNP Q1M7F4
C	-12	HIS	-	expression tag	UNP Q1M7F4
C	-11	HIS	-	expression tag	UNP Q1M7F4
C	-10	HIS	-	expression tag	UNP Q1M7F4
C	-9	SER	-	expression tag	UNP Q1M7F4
C	-8	SER	-	expression tag	UNP Q1M7F4
C	-7	GLY	-	expression tag	UNP Q1M7F4
C	-6	LEU	-	expression tag	UNP Q1M7F4
C	-5	VAL	-	expression tag	UNP Q1M7F4
C	-4	PRO	-	expression tag	UNP Q1M7F4
C	-3	ARG	-	expression tag	UNP Q1M7F4
C	-2	GLY	-	expression tag	UNP Q1M7F4
C	-1	SER	-	expression tag	UNP Q1M7F4
C	0	HIS	-	expression tag	UNP Q1M7F4
C	175	SER	CYS	conflict	UNP Q1M7F4
C	234	GLY	-	expression tag	UNP Q1M7F4
C	235	LEU	-	expression tag	UNP Q1M7F4
C	236	VAL	-	expression tag	UNP Q1M7F4
C	237	PRO	-	expression tag	UNP Q1M7F4
C	238	ARG	-	expression tag	UNP Q1M7F4
C	239	GLY	-	expression tag	UNP Q1M7F4
C	240	SER	-	expression tag	UNP Q1M7F4
C	241	ILE	-	expression tag	UNP Q1M7F4
C	242	GLU	-	expression tag	UNP Q1M7F4
D	-19	MET	-	initiating methionine	UNP Q1M7F4
D	-18	GLY	-	expression tag	UNP Q1M7F4
D	-17	SER	-	expression tag	UNP Q1M7F4
D	-16	SER	-	expression tag	UNP Q1M7F4
D	-15	HIS	-	expression tag	UNP Q1M7F4
D	-14	HIS	-	expression tag	UNP Q1M7F4
D	-13	HIS	-	expression tag	UNP Q1M7F4

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
D	-12	HIS	-	expression tag	UNP Q1M7F4
D	-11	HIS	-	expression tag	UNP Q1M7F4
D	-10	HIS	-	expression tag	UNP Q1M7F4
D	-9	SER	-	expression tag	UNP Q1M7F4
D	-8	SER	-	expression tag	UNP Q1M7F4
D	-7	GLY	-	expression tag	UNP Q1M7F4
D	-6	LEU	-	expression tag	UNP Q1M7F4
D	-5	VAL	-	expression tag	UNP Q1M7F4
D	-4	PRO	-	expression tag	UNP Q1M7F4
D	-3	ARG	-	expression tag	UNP Q1M7F4
D	-2	GLY	-	expression tag	UNP Q1M7F4
D	-1	SER	-	expression tag	UNP Q1M7F4
D	0	HIS	-	expression tag	UNP Q1M7F4
D	175	SER	CYS	conflict	UNP Q1M7F4
D	234	GLY	-	expression tag	UNP Q1M7F4
D	235	LEU	-	expression tag	UNP Q1M7F4
D	236	VAL	-	expression tag	UNP Q1M7F4
D	237	PRO	-	expression tag	UNP Q1M7F4
D	238	ARG	-	expression tag	UNP Q1M7F4
D	239	GLY	-	expression tag	UNP Q1M7F4
D	240	SER	-	expression tag	UNP Q1M7F4
D	241	ILE	-	expression tag	UNP Q1M7F4
D	242	GLU	-	expression tag	UNP Q1M7F4
F	-19	MET	-	initiating methionine	UNP Q1M7F4
F	-18	GLY	-	expression tag	UNP Q1M7F4
F	-17	SER	-	expression tag	UNP Q1M7F4
F	-16	SER	-	expression tag	UNP Q1M7F4
F	-15	HIS	-	expression tag	UNP Q1M7F4
F	-14	HIS	-	expression tag	UNP Q1M7F4
F	-13	HIS	-	expression tag	UNP Q1M7F4
F	-12	HIS	-	expression tag	UNP Q1M7F4
F	-11	HIS	-	expression tag	UNP Q1M7F4
F	-10	HIS	-	expression tag	UNP Q1M7F4
F	-9	SER	-	expression tag	UNP Q1M7F4
F	-8	SER	-	expression tag	UNP Q1M7F4
F	-7	GLY	-	expression tag	UNP Q1M7F4
F	-6	LEU	-	expression tag	UNP Q1M7F4
F	-5	VAL	-	expression tag	UNP Q1M7F4
F	-4	PRO	-	expression tag	UNP Q1M7F4
F	-3	ARG	-	expression tag	UNP Q1M7F4
F	-2	GLY	-	expression tag	UNP Q1M7F4
F	-1	SER	-	expression tag	UNP Q1M7F4

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	HIS	-	expression tag	UNP Q1M7F4
F	175	SER	CYS	conflict	UNP Q1M7F4
F	234	GLY	-	expression tag	UNP Q1M7F4
F	235	LEU	-	expression tag	UNP Q1M7F4
F	236	VAL	-	expression tag	UNP Q1M7F4
F	237	PRO	-	expression tag	UNP Q1M7F4
F	238	ARG	-	expression tag	UNP Q1M7F4
F	239	GLY	-	expression tag	UNP Q1M7F4
F	240	SER	-	expression tag	UNP Q1M7F4
F	241	ILE	-	expression tag	UNP Q1M7F4
F	242	GLU	-	expression tag	UNP Q1M7F4
G	-19	MET	-	initiating methionine	UNP Q1M7F4
G	-18	GLY	-	expression tag	UNP Q1M7F4
G	-17	SER	-	expression tag	UNP Q1M7F4
G	-16	SER	-	expression tag	UNP Q1M7F4
G	-15	HIS	-	expression tag	UNP Q1M7F4
G	-14	HIS	-	expression tag	UNP Q1M7F4
G	-13	HIS	-	expression tag	UNP Q1M7F4
G	-12	HIS	-	expression tag	UNP Q1M7F4
G	-11	HIS	-	expression tag	UNP Q1M7F4
G	-10	HIS	-	expression tag	UNP Q1M7F4
G	-9	SER	-	expression tag	UNP Q1M7F4
G	-8	SER	-	expression tag	UNP Q1M7F4
G	-7	GLY	-	expression tag	UNP Q1M7F4
G	-6	LEU	-	expression tag	UNP Q1M7F4
G	-5	VAL	-	expression tag	UNP Q1M7F4
G	-4	PRO	-	expression tag	UNP Q1M7F4
G	-3	ARG	-	expression tag	UNP Q1M7F4
G	-2	GLY	-	expression tag	UNP Q1M7F4
G	-1	SER	-	expression tag	UNP Q1M7F4
G	0	HIS	-	expression tag	UNP Q1M7F4
G	175	SER	CYS	conflict	UNP Q1M7F4
G	234	GLY	-	expression tag	UNP Q1M7F4
G	235	LEU	-	expression tag	UNP Q1M7F4
G	236	VAL	-	expression tag	UNP Q1M7F4
G	237	PRO	-	expression tag	UNP Q1M7F4
G	238	ARG	-	expression tag	UNP Q1M7F4
G	239	GLY	-	expression tag	UNP Q1M7F4
G	240	SER	-	expression tag	UNP Q1M7F4
G	241	ILE	-	expression tag	UNP Q1M7F4
G	242	GLU	-	expression tag	UNP Q1M7F4
H	-19	MET	-	initiating methionine	UNP Q1M7F4

*Continued on next page...*



*Continued from previous page...*

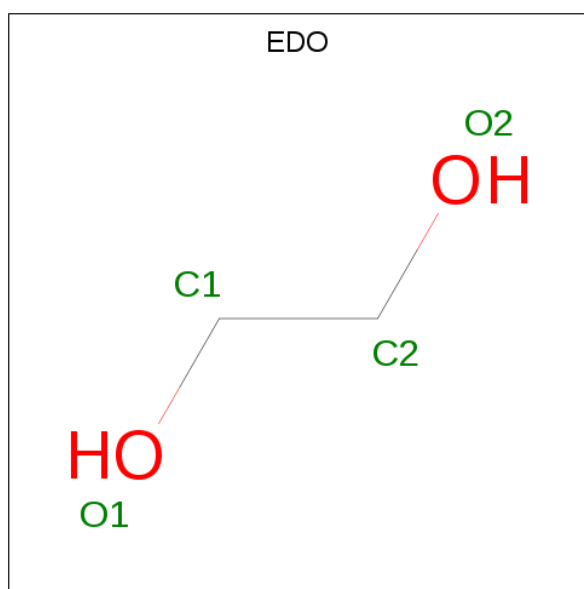
Chain	Residue	Modelled	Actual	Comment	Reference
H	-18	GLY	-	expression tag	UNP Q1M7F4
H	-17	SER	-	expression tag	UNP Q1M7F4
H	-16	SER	-	expression tag	UNP Q1M7F4
H	-15	HIS	-	expression tag	UNP Q1M7F4
H	-14	HIS	-	expression tag	UNP Q1M7F4
H	-13	HIS	-	expression tag	UNP Q1M7F4
H	-12	HIS	-	expression tag	UNP Q1M7F4
H	-11	HIS	-	expression tag	UNP Q1M7F4
H	-10	HIS	-	expression tag	UNP Q1M7F4
H	-9	SER	-	expression tag	UNP Q1M7F4
H	-8	SER	-	expression tag	UNP Q1M7F4
H	-7	GLY	-	expression tag	UNP Q1M7F4
H	-6	LEU	-	expression tag	UNP Q1M7F4
H	-5	VAL	-	expression tag	UNP Q1M7F4
H	-4	PRO	-	expression tag	UNP Q1M7F4
H	-3	ARG	-	expression tag	UNP Q1M7F4
H	-2	GLY	-	expression tag	UNP Q1M7F4
H	-1	SER	-	expression tag	UNP Q1M7F4
H	0	HIS	-	expression tag	UNP Q1M7F4
H	175	SER	CYS	conflict	UNP Q1M7F4
H	234	GLY	-	expression tag	UNP Q1M7F4
H	235	LEU	-	expression tag	UNP Q1M7F4
H	236	VAL	-	expression tag	UNP Q1M7F4
H	237	PRO	-	expression tag	UNP Q1M7F4
H	238	ARG	-	expression tag	UNP Q1M7F4
H	239	GLY	-	expression tag	UNP Q1M7F4
H	240	SER	-	expression tag	UNP Q1M7F4
H	241	ILE	-	expression tag	UNP Q1M7F4
H	242	GLU	-	expression tag	UNP Q1M7F4
I	-19	MET	-	initiating methionine	UNP Q1M7F4
I	-18	GLY	-	expression tag	UNP Q1M7F4
I	-17	SER	-	expression tag	UNP Q1M7F4
I	-16	SER	-	expression tag	UNP Q1M7F4
I	-15	HIS	-	expression tag	UNP Q1M7F4
I	-14	HIS	-	expression tag	UNP Q1M7F4
I	-13	HIS	-	expression tag	UNP Q1M7F4
I	-12	HIS	-	expression tag	UNP Q1M7F4
I	-11	HIS	-	expression tag	UNP Q1M7F4
I	-10	HIS	-	expression tag	UNP Q1M7F4
I	-9	SER	-	expression tag	UNP Q1M7F4
I	-8	SER	-	expression tag	UNP Q1M7F4
I	-7	GLY	-	expression tag	UNP Q1M7F4

*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
I	-6	LEU	-	expression tag	UNP Q1M7F4
I	-5	VAL	-	expression tag	UNP Q1M7F4
I	-4	PRO	-	expression tag	UNP Q1M7F4
I	-3	ARG	-	expression tag	UNP Q1M7F4
I	-2	GLY	-	expression tag	UNP Q1M7F4
I	-1	SER	-	expression tag	UNP Q1M7F4
I	0	HIS	-	expression tag	UNP Q1M7F4
I	175	SER	CYS	conflict	UNP Q1M7F4
I	234	GLY	-	expression tag	UNP Q1M7F4
I	235	LEU	-	expression tag	UNP Q1M7F4
I	236	VAL	-	expression tag	UNP Q1M7F4
I	237	PRO	-	expression tag	UNP Q1M7F4
I	238	ARG	-	expression tag	UNP Q1M7F4
I	239	GLY	-	expression tag	UNP Q1M7F4
I	240	SER	-	expression tag	UNP Q1M7F4
I	241	ILE	-	expression tag	UNP Q1M7F4
I	242	GLU	-	expression tag	UNP Q1M7F4

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



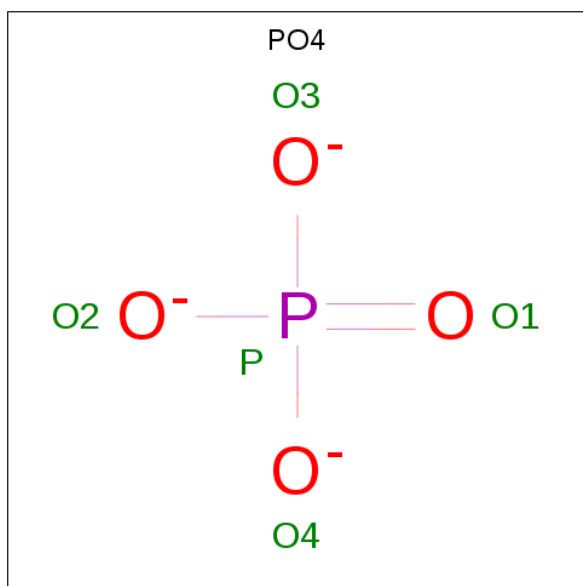
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	I	1	Total	C	O	0	0
			4	2	2		

Continued on next page...

*Continued from previous page...*

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	I	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	F	1	Total	O	P	0	0
			5	4	1		
3	I	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	200	Total	O	0	0
			200	200		
4	B	187	Total	O	0	0
			187	187		
4	C	197	Total	O	0	0
			197	197		
4	D	194	Total	O	0	0
			194	194		
4	F	189	Total	O	0	0
			189	189		

*Continued on next page...*

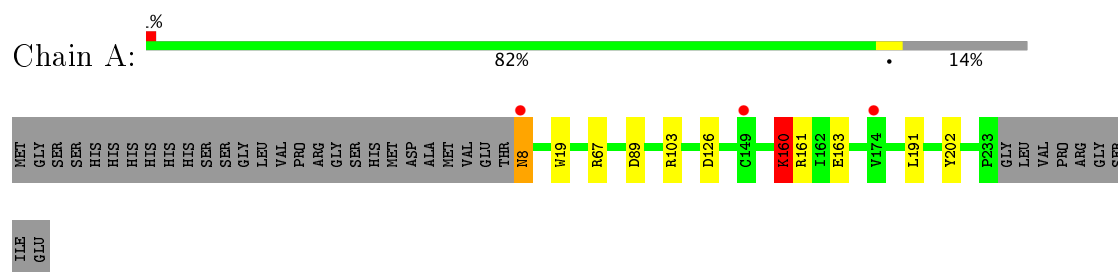
*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	163	Total 163	O 163	0	0
4	H	112	Total 112	O 112	0	0
4	I	203	Total 203	O 203	0	0

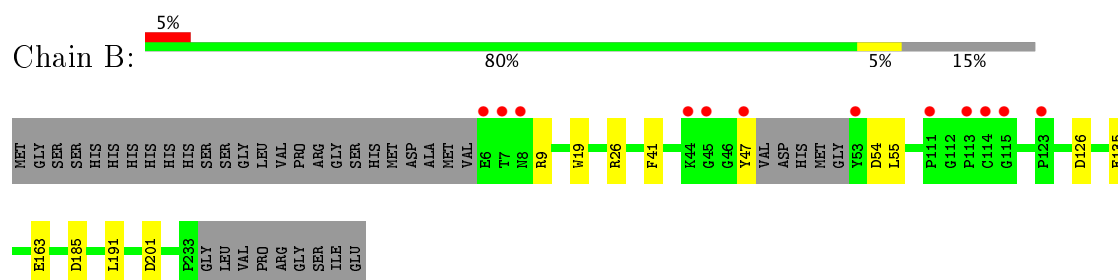
### 3 Residue-property plots [i](#)

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

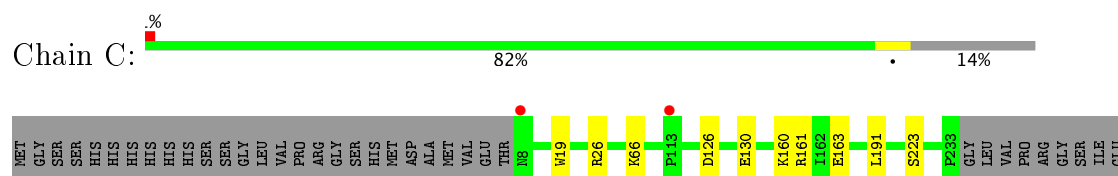
- Molecule 1: Putative amidase



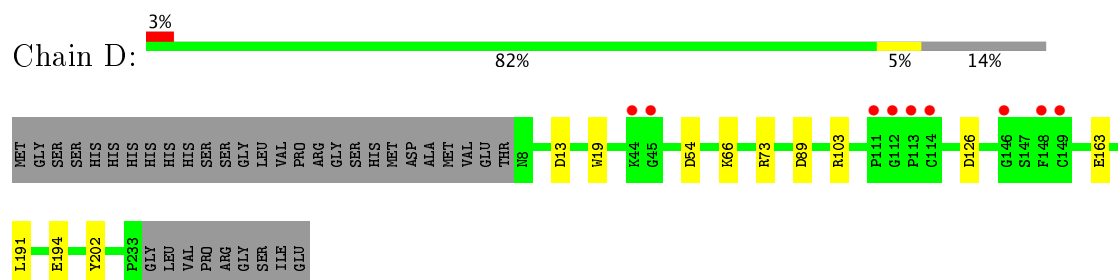
- Molecule 1: Putative amidase



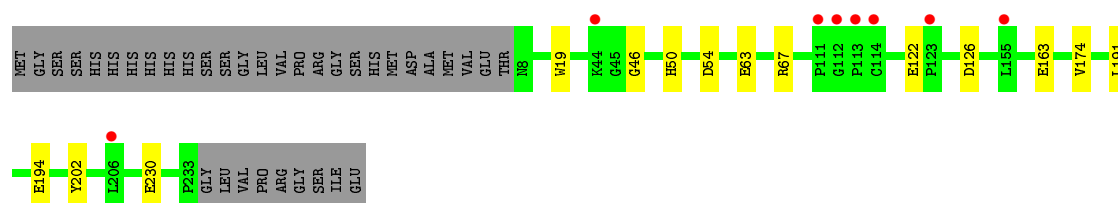
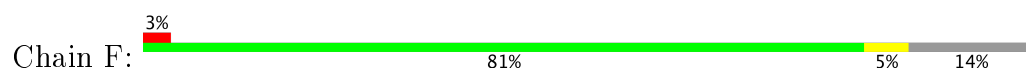
- Molecule 1: Putative amidase



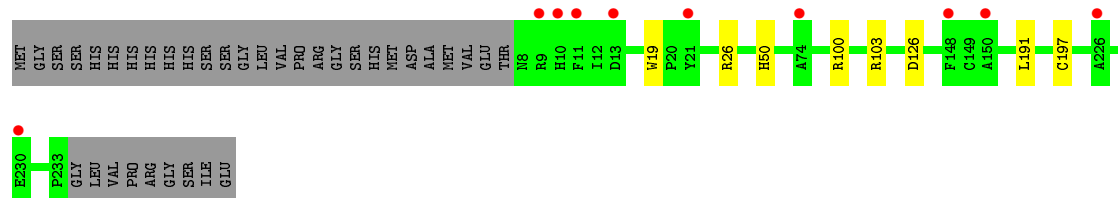
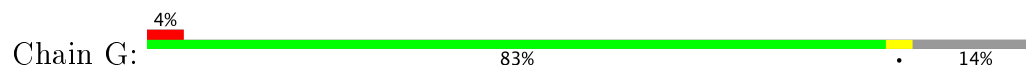
- Molecule 1: Putative amidase



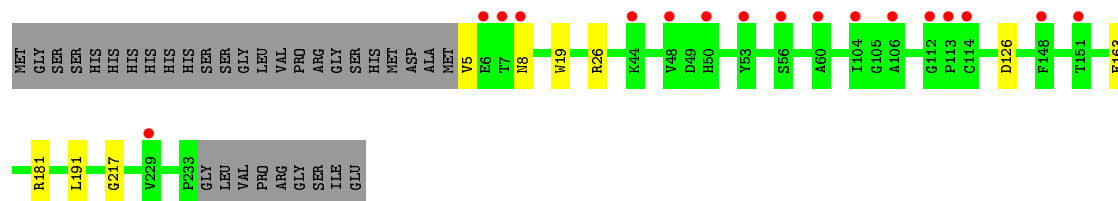
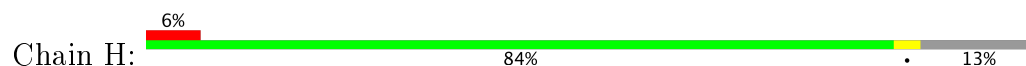
- Molecule 1: Putative amidase



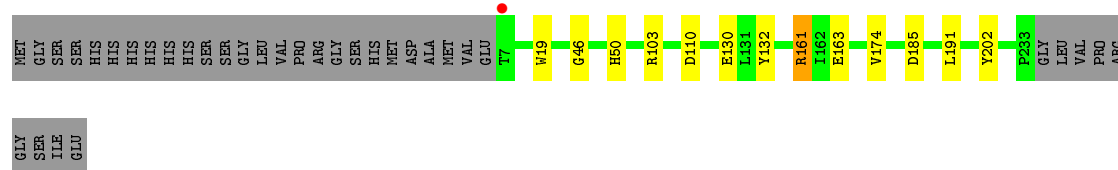
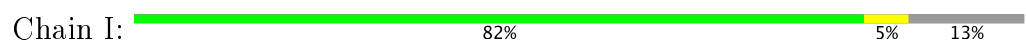
- Molecule 1: Putative amidase



- Molecule 1: Putative amidase



- Molecule 1: Putative amidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.18Å 86.92Å 342.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.48 – 1.75 43.58 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.1 (44.48-1.75) 99.1 (43.58-1.75)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.80 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.174 , 0.204 0.185 , 0.212	Depositor DCC
$R_{free}$ test set	11028 reflections (5.23%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.6	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 38.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	15574	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, EDO

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	0.94	1/1820 (0.1%)	1.01	11/2470 (0.4%)
1	B	1.02	2/1756 (0.1%)	1.01	7/2384 (0.3%)
1	C	0.93	3/1806 (0.2%)	0.91	3/2452 (0.1%)
1	D	0.90	1/1815 (0.1%)	0.95	9/2462 (0.4%)
1	F	0.90	3/1807 (0.2%)	0.94	5/2452 (0.2%)
1	G	0.87	2/1787 (0.1%)	0.87	3/2426 (0.1%)
1	H	0.80	0/1805	0.89	6/2451 (0.2%)
1	I	0.93	4/1813 (0.2%)	1.02	10/2463 (0.4%)
All	All	0.91	16/14409 (0.1%)	0.95	54/19560 (0.3%)

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	163	GLU	CD-OE1	-7.07	1.17	1.25
1	F	163	GLU	CD-OE2	-6.79	1.18	1.25
1	I	130	GLU	CD-OE2	-6.23	1.18	1.25
1	A	163	GLU	CD-OE1	-6.21	1.18	1.25
1	I	132[A]	TYR	CB-CG	6.06	1.60	1.51

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	161	ARG	NE-CZ-NH2	12.60	126.60	120.30
1	I	132[A]	TYR	CB-CG-CD2	-9.58	115.25	121.00
1	I	132[B]	TYR	CB-CG-CD2	-9.58	115.25	121.00
1	G	103	ARG	NE-CZ-NH2	8.39	124.49	120.30
1	I	132[A]	TYR	CB-CG-CD1	8.30	125.98	121.00

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	1780	0	1748	5	0
1	B	1721	0	1684	2	0
1	C	1766	0	1739	3	0
1	D	1776	0	1750	3	0
1	F	1768	0	1740	3	0
1	G	1748	0	1723	3	0
1	H	1766	0	1733	3	0
1	I	1773	0	1737	3	0
2	A	4	0	6	0	0
2	C	4	0	6	0	0
2	I	8	0	12	1	0
3	A	5	0	0	1	0
3	F	5	0	0	0	0
3	I	5	0	0	0	0
4	A	200	0	0	1	0
4	B	187	0	0	1	0
4	C	197	0	0	1	0
4	D	194	0	0	2	0
4	F	189	0	0	1	0
4	G	163	0	0	1	0
4	H	112	0	0	1	0
4	I	203	0	0	0	0
All	All	15574	0	13878	24	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 1.

The worst 5 of 24 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:F:194[A]:GLU:OE2	4:F:401:HOH:O	1.99	0.79
1:H:5:VAL:CB	4:H:405:HOH:O	2.33	0.76
1:B:135[B]:GLU:OE2	4:B:301:HOH:O	2.15	0.63
1:A:8:ASN:ND2	1:A:8:ASN:C	2.57	0.58
1:I:161:ARG:NH2	2:I:302:EDO:O1	2.41	0.54

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	229/262 (87%)	225 (98%)	4 (2%)	0	100	100
1	B	222/262 (85%)	215 (97%)	7 (3%)	0	100	100
1	C	228/262 (87%)	223 (98%)	5 (2%)	0	100	100
1	D	229/262 (87%)	224 (98%)	5 (2%)	0	100	100
1	F	228/262 (87%)	224 (98%)	3 (1%)	1 (0%)	38	18
1	G	226/262 (86%)	223 (99%)	3 (1%)	0	100	100
1	H	229/262 (87%)	223 (97%)	6 (3%)	0	100	100
1	I	229/262 (87%)	225 (98%)	3 (1%)	1 (0%)	38	18
All	All	1820/2096 (87%)	1782 (98%)	36 (2%)	2 (0%)	55	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	174	VAL
1	I	174	VAL

### 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	189/214 (88%)	187 (99%)	2 (1%)	78	64

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	181/214 (85%)	180 (99%)	1 (1%)	89	82
1	C	187/214 (87%)	187 (100%)	0	100	100
1	D	188/214 (88%)	186 (99%)	2 (1%)	78	64
1	F	187/214 (87%)	186 (100%)	1 (0%)	91	86
1	G	185/214 (86%)	185 (100%)	0	100	100
1	H	186/214 (87%)	186 (100%)	0	100	100
1	I	187/214 (87%)	187 (100%)	0	100	100
All	All	1490/1712 (87%)	1484 (100%)	6 (0%)	94	89

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	47	TYR
1	F	230	GLU
1	D	66[A]	LYS
1	A	160	LYS
1	D	66[B]	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN

### 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

7 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	EDO	A	301	-	3,3,3	0.77	0	2,2,2	0.86	0
3	PO4	A	302	-	4,4,4	1.38	1 (25%)	6,6,6	2.22	2 (33%)
2	EDO	C	301	-	3,3,3	0.63	0	2,2,2	1.42	0
3	PO4	F	301	-	4,4,4	1.31	0	6,6,6	1.33	0
2	EDO	I	301	-	3,3,3	0.54	0	2,2,2	1.20	0
2	EDO	I	302	-	3,3,3	0.28	0	2,2,2	1.11	0
3	PO4	I	303	-	4,4,4	0.81	0	6,6,6	1.48	1 (16%)

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	EDO	A	301	-	-	0/1/1/1	0/0/0/0
3	PO4	A	302	-	-	0/0/0/0	0/0/0/0
2	EDO	C	301	-	-	0/1/1/1	0/0/0/0
3	PO4	F	301	-	-	0/0/0/0	0/0/0/0
2	EDO	I	301	-	-	0/1/1/1	0/0/0/0
2	EDO	I	302	-	-	0/1/1/1	0/0/0/0
3	PO4	I	303	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	PO4	P-O3	-2.66	1.45	1.54

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	302	PO4	O4-P-O3	-3.44	95.27	107.90
3	I	303	PO4	O4-P-O3	2.23	116.11	107.90
3	A	302	PO4	O4-P-O1	3.08	124.08	110.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	302	PO4	1	0
2	I	302	EDO	1	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 ⓘ

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	226/262 (86%)	-0.11	3 (1%) 77 84	15, 23, 36, 64	0
1	B	223/262 (85%)	0.23	12 (5%) 26 33	15, 24, 52, 75	0
1	C	226/262 (86%)	-0.07	2 (0%) 84 89	17, 27, 39, 63	0
1	D	226/262 (86%)	0.32	9 (3%) 39 46	17, 26, 41, 66	0
1	F	226/262 (86%)	0.18	8 (3%) 44 51	19, 27, 40, 60	0
1	G	226/262 (86%)	0.16	10 (4%) 35 42	20, 31, 47, 67	0
1	H	229/262 (87%)	0.51	17 (7%) 15 21	21, 36, 62, 81	0
1	I	227/262 (86%)	-0.19	1 (0%) 92 95	19, 24, 33, 75	0
All	All	1809/2096 (86%)	0.13	62 (3%) 46 52	15, 27, 48, 81	0

The worst 5 of 62 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	7	THR	4.8
1	B	45	GLY	4.8
1	B	53	TYR	4.4
1	B	111	PRO	4.3
1	D	114	CYS	4.1

### 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 [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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	PO4	A	302	5/5	0.80	0.35	8.96	43,45,48,61	5
2	EDO	I	302	4/4	0.90	0.34	6.70	50,52,56,59	0
3	PO4	I	303	5/5	0.83	0.27	4.85	44,49,56,60	5
3	PO4	F	301	5/5	0.75	0.20	3.99	49,50,55,69	5
2	EDO	A	301	4/4	0.83	0.23	3.10	39,46,47,51	0
2	EDO	C	301	4/4	0.83	0.16	1.33	43,53,53,55	0
2	EDO	I	301	4/4	0.94	0.07	0.25	34,35,36,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.