



# wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 26, 2023 – 03:28 AM EDT

PDB ID : 6BFZ  
Title : Crystal structure of enolase from E. coli with a mixture of apo form, substrate, and product form  
Authors : Erlandsen, H.; Wright, D.; Krucinska, J.  
Deposited on : 2017-10-27  
Resolution : 2.21 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

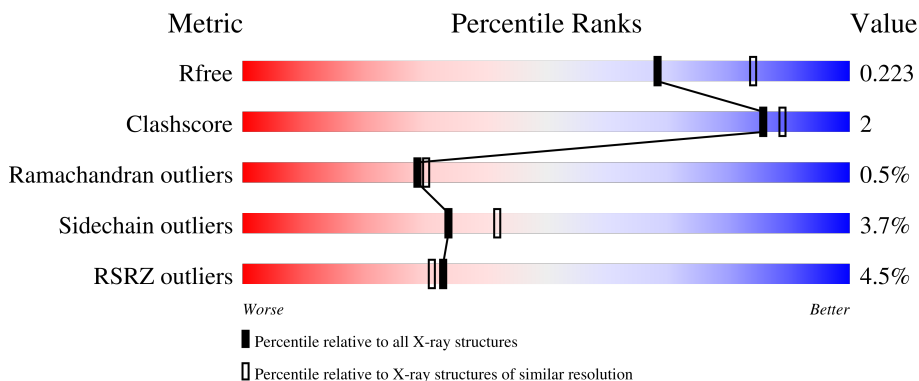
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.21 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.



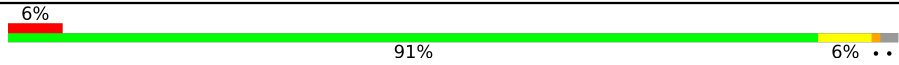
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5912 (2.24-2.20)
Clashscore	141614	6646 (2.24-2.20)
Ramachandran outliers	138981	6543 (2.24-2.20)
Sidechain outliers	138945	6544 (2.24-2.20)
RSRZ outliers	127900	5797 (2.24-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	449	 89% 8% ..
1	B	449	 90% 6% ..
1	C	449	 87% 7% ..
1	D	449	 88% 9% ..
1	E	449	 91% 6% ..

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Mol	Chain	Length	Quality of chain
1	F	449	

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
5	PEP	E	602	-	X	X	-

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 20287 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 Enolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	439	Total 3246	C 2030	N 557	O 644	S 15	0	0	0
1	B	439	Total 3254	C 2036	N 559	O 644	S 15	0	0	0
1	E	439	Total 3254	C 2036	N 559	O 644	S 15	0	0	0
1	F	438	Total 3254	C 2036	N 559	O 644	S 15	0	1	0
1	D	438	Total 3253	C 2034	N 559	O 645	S 15	0	1	0
1	C	429	Total 3183	C 1996	N 544	O 628	S 15	0	0	0

There are 102 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	GLY	-	expression tag	UNP B7MLA0
A	-16	SER	-	expression tag	UNP B7MLA0
A	-15	HIS	-	expression tag	UNP B7MLA0
A	-14	MET	-	expression tag	UNP B7MLA0
A	-13	ALA	-	expression tag	UNP B7MLA0
A	-12	SER	-	expression tag	UNP B7MLA0
A	-11	MET	-	expression tag	UNP B7MLA0
A	-10	THR	-	expression tag	UNP B7MLA0
A	-9	GLY	-	expression tag	UNP B7MLA0
A	-8	GLY	-	expression tag	UNP B7MLA0
A	-7	GLN	-	expression tag	UNP B7MLA0
A	-6	GLN	-	expression tag	UNP B7MLA0
A	-5	MET	-	expression tag	UNP B7MLA0
A	-4	GLY	-	expression tag	UNP B7MLA0
A	-3	ARG	-	expression tag	UNP B7MLA0
A	-2	GLY	-	expression tag	UNP B7MLA0
A	-1	SER	-	expression tag	UNP B7MLA0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	GLY	-	expression tag	UNP B7MLA0
B	-16	SER	-	expression tag	UNP B7MLA0
B	-15	HIS	-	expression tag	UNP B7MLA0
B	-14	MET	-	expression tag	UNP B7MLA0
B	-13	ALA	-	expression tag	UNP B7MLA0
B	-12	SER	-	expression tag	UNP B7MLA0
B	-11	MET	-	expression tag	UNP B7MLA0
B	-10	THR	-	expression tag	UNP B7MLA0
B	-9	GLY	-	expression tag	UNP B7MLA0
B	-8	GLY	-	expression tag	UNP B7MLA0
B	-7	GLN	-	expression tag	UNP B7MLA0
B	-6	GLN	-	expression tag	UNP B7MLA0
B	-5	MET	-	expression tag	UNP B7MLA0
B	-4	GLY	-	expression tag	UNP B7MLA0
B	-3	ARG	-	expression tag	UNP B7MLA0
B	-2	GLY	-	expression tag	UNP B7MLA0
B	-1	SER	-	expression tag	UNP B7MLA0
E	-17	GLY	-	expression tag	UNP B7MLA0
E	-16	SER	-	expression tag	UNP B7MLA0
E	-15	HIS	-	expression tag	UNP B7MLA0
E	-14	MET	-	expression tag	UNP B7MLA0
E	-13	ALA	-	expression tag	UNP B7MLA0
E	-12	SER	-	expression tag	UNP B7MLA0
E	-11	MET	-	expression tag	UNP B7MLA0
E	-10	THR	-	expression tag	UNP B7MLA0
E	-9	GLY	-	expression tag	UNP B7MLA0
E	-8	GLY	-	expression tag	UNP B7MLA0
E	-7	GLN	-	expression tag	UNP B7MLA0
E	-6	GLN	-	expression tag	UNP B7MLA0
E	-5	MET	-	expression tag	UNP B7MLA0
E	-4	GLY	-	expression tag	UNP B7MLA0
E	-3	ARG	-	expression tag	UNP B7MLA0
E	-2	GLY	-	expression tag	UNP B7MLA0
E	-1	SER	-	expression tag	UNP B7MLA0
F	-17	GLY	-	expression tag	UNP B7MLA0
F	-16	SER	-	expression tag	UNP B7MLA0
F	-15	HIS	-	expression tag	UNP B7MLA0
F	-14	MET	-	expression tag	UNP B7MLA0
F	-13	ALA	-	expression tag	UNP B7MLA0
F	-12	SER	-	expression tag	UNP B7MLA0
F	-11	MET	-	expression tag	UNP B7MLA0
F	-10	THR	-	expression tag	UNP B7MLA0

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-9	GLY	-	expression tag	UNP B7MLA0
F	-8	GLY	-	expression tag	UNP B7MLA0
F	-7	GLN	-	expression tag	UNP B7MLA0
F	-6	GLN	-	expression tag	UNP B7MLA0
F	-5	MET	-	expression tag	UNP B7MLA0
F	-4	GLY	-	expression tag	UNP B7MLA0
F	-3	ARG	-	expression tag	UNP B7MLA0
F	-2	GLY	-	expression tag	UNP B7MLA0
F	-1	SER	-	expression tag	UNP B7MLA0
D	-17	GLY	-	expression tag	UNP B7MLA0
D	-16	SER	-	expression tag	UNP B7MLA0
D	-15	HIS	-	expression tag	UNP B7MLA0
D	-14	MET	-	expression tag	UNP B7MLA0
D	-13	ALA	-	expression tag	UNP B7MLA0
D	-12	SER	-	expression tag	UNP B7MLA0
D	-11	MET	-	expression tag	UNP B7MLA0
D	-10	THR	-	expression tag	UNP B7MLA0
D	-9	GLY	-	expression tag	UNP B7MLA0
D	-8	GLY	-	expression tag	UNP B7MLA0
D	-7	GLN	-	expression tag	UNP B7MLA0
D	-6	GLN	-	expression tag	UNP B7MLA0
D	-5	MET	-	expression tag	UNP B7MLA0
D	-4	GLY	-	expression tag	UNP B7MLA0
D	-3	ARG	-	expression tag	UNP B7MLA0
D	-2	GLY	-	expression tag	UNP B7MLA0
D	-1	SER	-	expression tag	UNP B7MLA0
C	-17	GLY	-	expression tag	UNP B7MLA0
C	-16	SER	-	expression tag	UNP B7MLA0
C	-15	HIS	-	expression tag	UNP B7MLA0
C	-14	MET	-	expression tag	UNP B7MLA0
C	-13	ALA	-	expression tag	UNP B7MLA0
C	-12	SER	-	expression tag	UNP B7MLA0
C	-11	MET	-	expression tag	UNP B7MLA0
C	-10	THR	-	expression tag	UNP B7MLA0
C	-9	GLY	-	expression tag	UNP B7MLA0
C	-8	GLY	-	expression tag	UNP B7MLA0
C	-7	GLN	-	expression tag	UNP B7MLA0
C	-6	GLN	-	expression tag	UNP B7MLA0
C	-5	MET	-	expression tag	UNP B7MLA0
C	-4	GLY	-	expression tag	UNP B7MLA0
C	-3	ARG	-	expression tag	UNP B7MLA0
C	-2	GLY	-	expression tag	UNP B7MLA0

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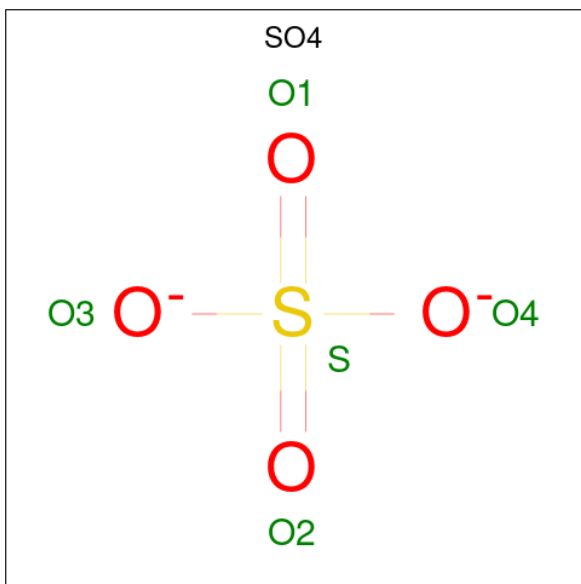
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Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	SER	-	expression tag	UNP B7MLA0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0
2	E	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0
2	D	1	Total Mg 1 1	0	0
2	C	1	Total Mg 1 1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



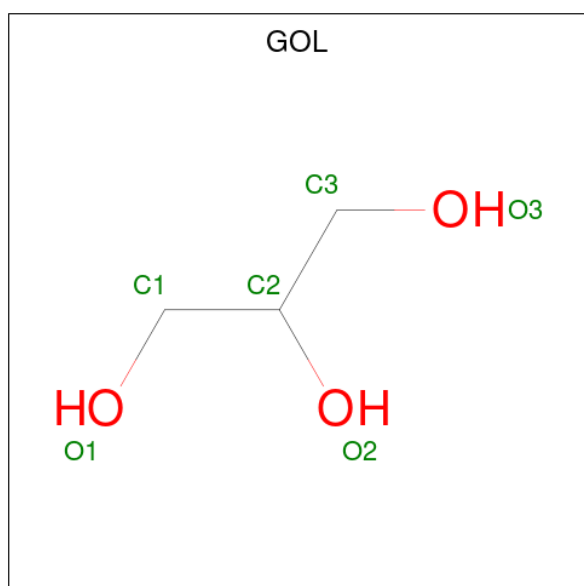
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

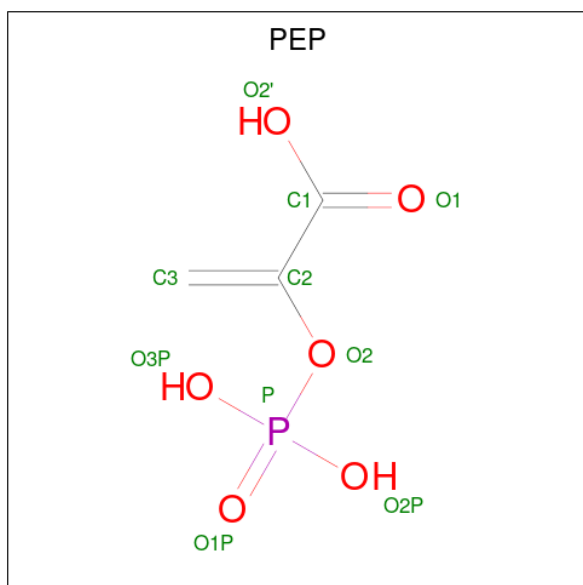
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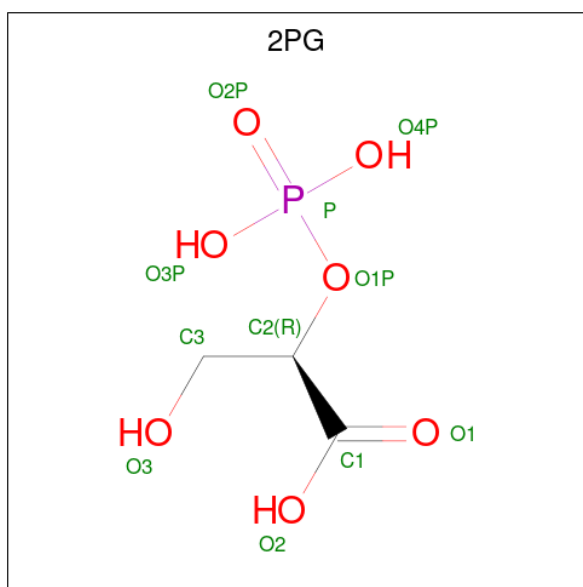
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	F	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is PHOSPHOENOLPYRUVATE (three-letter code: PEP) (formula:  $C_3H_5O_6P$ ).



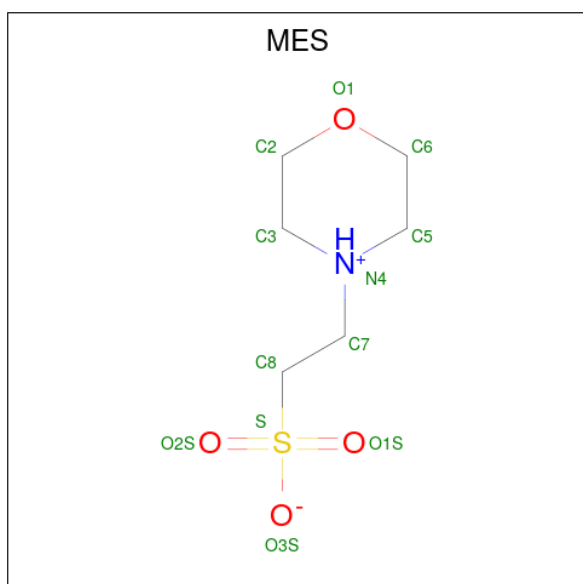
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	E	1	Total	C	O	P	0	0
			10	3	6	1		

- Molecule 6 is 2-PHOSPHOGLYCERIC ACID (three-letter code: 2PG) (formula:  $C_3H_7O_7P$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
6	E	1	11	3	7	1	0	0
6	F	1	11	3	7	1	0	0
6	D	1	11	3	7	1	0	0
6	C	1	11	3	7	1	0	0

- Molecule 7 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	E	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
7	D	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

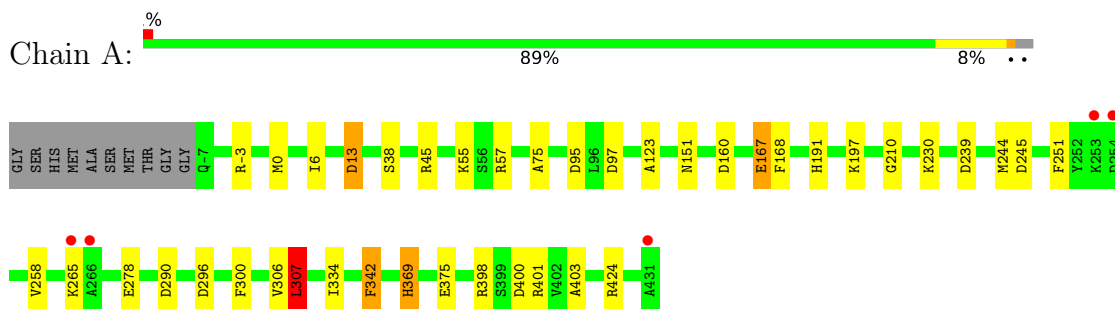
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	133	Total	O	0	0
			133	133		
8	B	110	Total	O	0	0
			110	110		
8	E	133	Total	O	0	0
			133	133		
8	F	97	Total	O	0	0
			97	97		
8	D	113	Total	O	0	0
			113	113		
8	C	90	Total	O	0	0
			90	90		

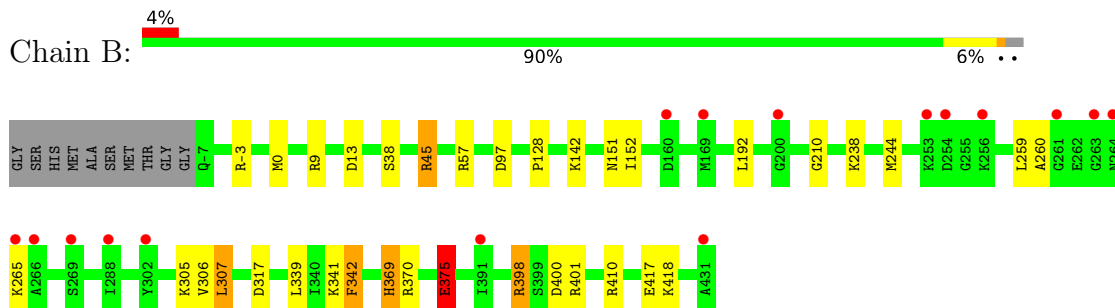
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

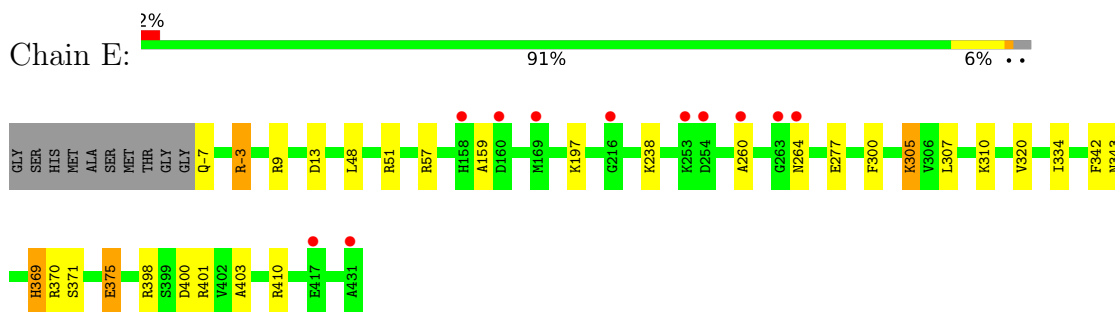
- Molecule 1: Enolase



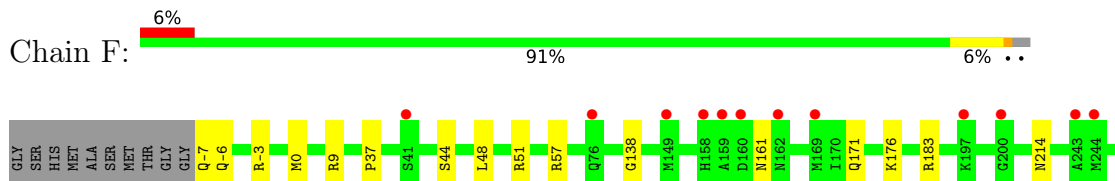
- Molecule 1: Enolase

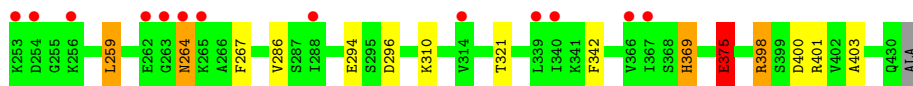


- Molecule 1: Enolase

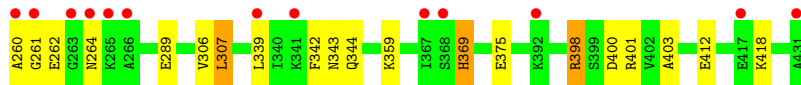
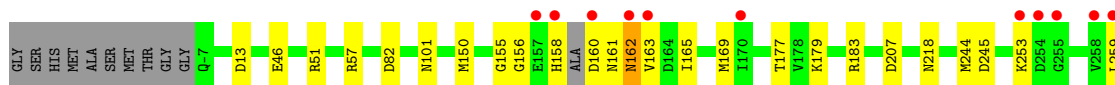
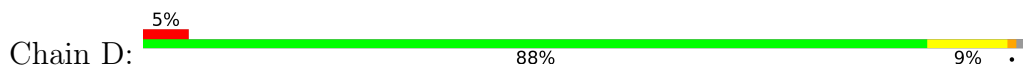


- Molecule 1: Enolase

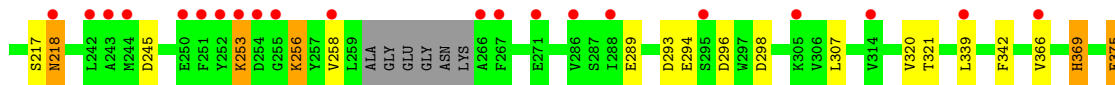
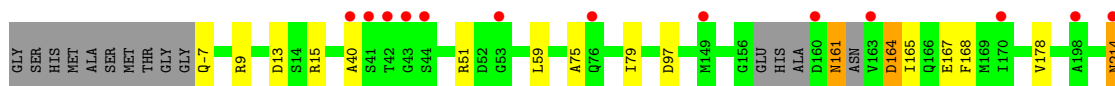
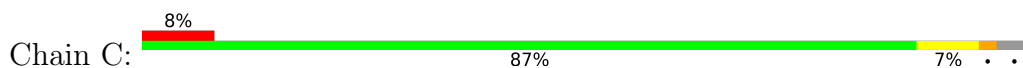




● Molecule 1: Enolase



● Molecule 1: Enolase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.26Å 143.29Å 207.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.69 – 2.21 29.69 – 2.21	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.69-2.21) 99.4 (29.69-2.21)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.55 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.171 , 0.218 0.180 , 0.223	Depositor DCC
$R_{free}$ test set	8374 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.0	Xtrriage
Anisotropy	0.042	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 31.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	20287	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.68% 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 i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: 2PG, GOL, MG, SO4, PEP, MES

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	0.90	1/3288 (0.0%)	0.96	13/4427 (0.3%)
1	B	0.90	2/3296 (0.1%)	0.95	11/4435 (0.2%)
1	C	0.93	0/3221	0.95	8/4330 (0.2%)
1	D	0.92	0/3294	0.93	10/4432 (0.2%)
1	E	0.90	1/3296 (0.0%)	0.92	9/4435 (0.2%)
1	F	0.88	0/3299	0.94	16/4439 (0.4%)
All	All	0.90	4/19694 (0.0%)	0.94	67/26498 (0.3%)

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	C	0	1
1	D	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	38	SER	CB-OG	7.53	1.52	1.42
1	E	371	SER	CB-OG	-5.72	1.34	1.42
1	A	38	SER	CB-OG	5.20	1.49	1.42
1	B	417	GLU	CG-CD	5.15	1.59	1.51

The worst 5 of 67 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	ARG	NE-CZ-NH1	11.78	126.19	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45	ARG	NE-CZ-NH1	11.04	125.82	120.30
1	A	45	ARG	NE-CZ-NH2	-10.15	115.23	120.30
1	B	45	ARG	NE-CZ-NH2	-10.08	115.26	120.30
1	D	13	ASP	CB-CG-OD1	8.41	125.87	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	40	ALA	Peptide
1	D	156	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	3246	0	3257	13	0
1	B	3254	0	3279	10	0
1	C	3183	0	3216	15	0
1	D	3253	0	3267	25	0
1	E	3254	0	3279	10	0
1	F	3254	0	3280	11	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	20	0	0	0	0
3	B	10	0	0	1	0
3	C	10	0	0	0	0
3	D	5	0	0	0	0
3	E	10	0	0	0	0
3	F	10	0	0	0	0
4	B	6	0	8	1	0
4	C	6	0	8	2	0
4	F	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	10	0	1	6	0
6	C	11	0	4	0	0
6	D	11	0	4	0	0
6	E	11	0	1	3	0
6	F	11	0	4	0	0
7	D	12	0	13	0	0
7	E	12	0	13	0	0
8	A	133	0	0	3	0
8	B	110	0	0	0	0
8	C	90	0	0	1	0
8	D	113	0	0	2	0
8	E	133	0	0	2	0
8	F	97	0	0	0	0
All	All	20287	0	19642	85	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 2.

The worst 5 of 85 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:D:101[B]:ASN:H	1:D:101[B]:ASN:HD22	1.22	0.87
5:E:602:PEP:O2'	6:E:603:2PG:O2P	1.94	0.85
1:D:101[B]:ASN:ND2	8:D:601:HOH:O	2.17	0.77
1:E:-7:GLN:OE1	8:E:702:HOH:O	2.05	0.75
1:F:369:HIS:CD2	1:F:401:ARG:HH11	2.06	0.73

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	437/449 (97%)	423 (97%)	13 (3%)	1 (0%)	47	54
1	B	437/449 (97%)	423 (97%)	12 (3%)	2 (0%)	29	30
1	C	421/449 (94%)	397 (94%)	21 (5%)	3 (1%)	22	21
1	D	435/449 (97%)	417 (96%)	15 (3%)	3 (1%)	22	21
1	E	437/449 (97%)	418 (96%)	15 (3%)	4 (1%)	17	15
1	F	437/449 (97%)	422 (97%)	14 (3%)	1 (0%)	47	54
All	All	2604/2694 (97%)	2500 (96%)	90 (4%)	14 (0%)	29	30

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	262	GLU
1	C	164	ASP
1	D	261	GLY
1	B	260	ALA
1	E	159	ALA

### 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	329/337 (98%)	317 (96%)	12 (4%)	35	43
1	B	331/337 (98%)	317 (96%)	14 (4%)	30	36
1	C	325/337 (96%)	311 (96%)	14 (4%)	29	35
1	D	331/337 (98%)	322 (97%)	9 (3%)	44	55
1	E	331/337 (98%)	320 (97%)	11 (3%)	38	47
1	F	332/337 (98%)	319 (96%)	13 (4%)	32	40
All	All	1979/2022 (98%)	1906 (96%)	73 (4%)	34	42

5 of 73 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	400	ASP

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Mol	Chain	Res	Type
1	C	369	HIS
1	C	97	ASP
1	C	253	LYS
1	B	400	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 38 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	171	GLN
1	C	214	ASN
1	D	264	ASN
1	D	369	HIS
1	C	369	HIS

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 6 are monoatomic - leaving 23 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	SO4	F	503	-	4,4,4	0.43	0	6,6,6	0.56	0
6	2PG	E	603	2,5	9,10,10	1.31	1 (11%)	11,14,14	2.57	6 (54%)
6	2PG	C	501	2	9,10,10	1.30	1 (11%)	11,14,14	1.98	4 (36%)
7	MES	D	504	-	12,12,12	1.83	1 (8%)	14,16,16	2.27	6 (42%)
3	SO4	E	605	-	4,4,4	0.40	0	6,6,6	0.46	0
3	SO4	A	604	-	4,4,4	0.45	0	6,6,6	0.48	0
6	2PG	F	501	2	9,10,10	1.39	1 (11%)	11,14,14	2.85	7 (63%)
3	SO4	A	603	-	4,4,4	0.28	0	6,6,6	0.50	0
3	SO4	A	605	-	4,4,4	0.63	0	6,6,6	0.16	0
7	MES	E	606	-	12,12,12	2.16	2 (16%)	14,16,16	2.45	7 (50%)
5	PEP	E	602	6,2	9,9,9	1.34	1 (11%)	11,13,13	2.77	6 (54%)
6	2PG	D	501	2	9,10,10	1.67	4 (44%)	11,14,14	2.51	4 (36%)
3	SO4	C	504	-	4,4,4	0.74	0	6,6,6	0.43	0
3	SO4	D	503	-	4,4,4	0.40	0	6,6,6	0.50	0
3	SO4	F	504	-	4,4,4	0.84	0	6,6,6	0.44	0
3	SO4	C	503	-	4,4,4	0.62	0	6,6,6	0.53	0
3	SO4	B	602	-	4,4,4	0.31	0	6,6,6	1.46	1 (16%)
4	GOL	F	505	-	5,5,5	0.55	0	5,5,5	1.02	0
4	GOL	B	604	-	5,5,5	0.49	0	5,5,5	1.86	1 (20%)
3	SO4	E	604	-	4,4,4	0.69	0	6,6,6	0.55	0
4	GOL	C	505	-	5,5,5	0.56	0	5,5,5	1.44	1 (20%)
3	SO4	A	602	-	4,4,4	0.71	0	6,6,6	1.40	1 (16%)
3	SO4	B	603	-	4,4,4	0.59	0	6,6,6	2.23	3 (50%)

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
6	2PG	F	501	2	-	2/11/11/11	-
4	GOL	F	505	-	-	2/4/4/4	-
5	PEP	E	602	6,2	-	6/9/9/9	-
6	2PG	D	501	2	-	0/11/11/11	-
6	2PG	E	603	2,5	-	5/11/11/11	-
6	2PG	C	501	2	-	0/11/11/11	-
7	MES	E	606	-	-	4/6/14/14	0/1/1/1
4	GOL	B	604	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	MES	D	504	-	-	0/6/14/14	0/1/1/1
4	GOL	C	505	-	-	2/4/4/4	-

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	E	606	MES	C8-S	-6.62	1.68	1.77
7	D	504	MES	C8-S	-5.17	1.70	1.77
5	E	602	PEP	C2-C1	-2.76	1.46	1.49
6	D	501	2PG	P-O2P	2.63	1.59	1.50
6	F	501	2PG	C2-C1	2.51	1.54	1.52

The worst 5 of 47 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	F	501	2PG	O4P-P-O2P	5.52	132.29	110.68
6	D	501	2PG	O3P-P-O2P	5.17	130.91	110.68
7	E	606	MES	C6-C5-N4	4.93	117.58	110.10
5	E	602	PEP	O2-C2-C3	-4.65	115.83	124.79
6	E	603	2PG	O1P-P-O2P	-4.60	91.63	109.39

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	604	GOL	O1-C1-C2-C3
4	C	505	GOL	C1-C2-C3-O3
4	C	505	GOL	O2-C2-C3-O3
5	E	602	PEP	O1-C1-C2-C3
5	E	602	PEP	O2'-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	603	2PG	3	0
5	E	602	PEP	6	0
3	B	602	SO4	1	0
4	B	604	GOL	1	0
4	C	505	GOL	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, 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	439/449 (97%)	-0.30	5 (1%) 80 79	25, 34, 56, 84	0
1	B	439/449 (97%)	-0.15	16 (3%) 42 40	24, 36, 65, 97	0
1	C	429/449 (95%)	0.26	36 (8%) 11 9	26, 40, 82, 107	0
1	D	438/449 (97%)	-0.04	24 (5%) 25 23	26, 37, 67, 103	0
1	E	439/449 (97%)	-0.16	11 (2%) 57 55	25, 34, 61, 104	0
1	F	438/449 (97%)	-0.03	25 (5%) 23 22	27, 37, 64, 104	0
All	All	2622/2694 (97%)	-0.07	117 (4%) 33 31	24, 36, 67, 107	0

The worst 5 of 117 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	431	ALA	7.0
1	E	263	GLY	6.7
1	D	157	GLU	5.7
1	C	266	ALA	5.6
1	C	253	LYS	5.2

### 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 monosaccharides 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	SO4	F	503	5/5	0.77	0.31	87,103,110,116	0
7	MES	D	504	12/12	0.81	0.25	62,68,91,92	0
7	MES	E	606	12/12	0.85	0.15	52,56,74,77	0
4	GOL	F	505	6/6	0.85	0.23	42,51,61,76	0
4	GOL	B	604	6/6	0.86	0.20	51,55,57,72	0
2	MG	C	502	1/1	0.86	0.06	50,50,50,50	0
4	GOL	C	505	6/6	0.87	0.19	41,49,60,73	0
2	MG	B	601	1/1	0.93	0.17	53,53,53,53	0
3	SO4	F	504	5/5	0.93	0.11	52,53,60,77	0
3	SO4	D	503	5/5	0.93	0.27	88,91,98,107	0
3	SO4	E	605	5/5	0.93	0.22	92,94,105,108	0
6	2PG	C	501	11/11	0.94	0.11	45,55,63,64	0
3	SO4	E	604	5/5	0.95	0.14	62,63,64,77	0
2	MG	A	601	1/1	0.95	0.10	54,54,54,54	0
6	2PG	F	501	11/11	0.95	0.10	33,38,45,58	0
6	2PG	D	501	11/11	0.96	0.11	35,40,46,53	0
2	MG	D	502	1/1	0.96	0.06	41,41,41,41	0
6	2PG	E	603	11/11	0.96	0.11	17,25,28,36	11
3	SO4	C	503	5/5	0.96	0.12	53,60,70,76	0
3	SO4	B	603	5/5	0.97	0.10	39,45,48,55	0
2	MG	E	601	1/1	0.97	0.08	35,35,35,35	0
3	SO4	A	602	5/5	0.97	0.09	41,46,53,55	0
3	SO4	A	604	5/5	0.97	0.08	46,47,49,57	0
3	SO4	A	605	5/5	0.97	0.12	52,56,60,68	0
5	PEP	E	602	10/10	0.97	0.12	36,46,56,63	10
3	SO4	B	602	5/5	0.98	0.10	44,49,52,54	0
3	SO4	C	504	5/5	0.98	0.10	51,51,59,64	0
3	SO4	A	603	5/5	0.99	0.09	41,43,45,50	0
2	MG	F	502	1/1	0.99	0.08	39,39,39,39	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.