

Full wwPDB X-ray Structure Validation Report (i)

Nov 6, 2023 – 12:35 PM EST

PDB ID	:	6BPR
Title	:	Crystal structure of cysteine, nitric oxide-bound ferrous form of the un-
		crosslinked F2-Tyr157 human cysteine dioxygenase
Authors	:	Liu, A.; Li, J.; Shin, I.
Deposited on		
Resolution	:	1.96 Å(reported)
Authors Deposited on	:	crosslinked F2-Tyr157 human cysteine dioxygenase Liu, A.; Li, J.; Shin, I.

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 https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

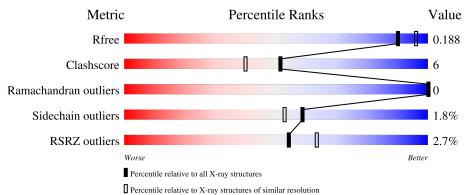
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.96 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 >=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 <=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		
			2%		
1	А	200	81%	11%	• 7%

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
6	GOL	А	307	-	-	-	Х

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Mo	l Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GOL	А	313	-	-	-	Х



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2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 1823 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 Cysteine dioxygenase type 1.

Mol	Chain	Residues		-	Ato	\mathbf{ms}			ZeroOcc	AltConf	Trace
1	А	186	Total 1565	C 988	F 2	N 278	O 284	S 13	0	9	0

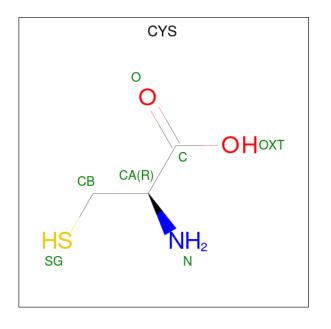
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	1	SER	-	expression tag	UNP Q16878
А	137	VAL	ILE	conflict	UNP Q16878

• Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

N	[o]	Chain	Residues	Aton	ns	ZeroOcc	AltConf
	2	А	1	Total 1	Fe 1	0	0

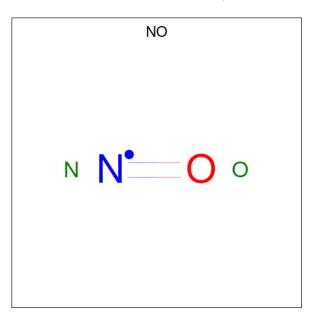
• Molecule 3 is CYSTEINE (three-letter code: CYS) (formula: C₃H₇NO₂S).



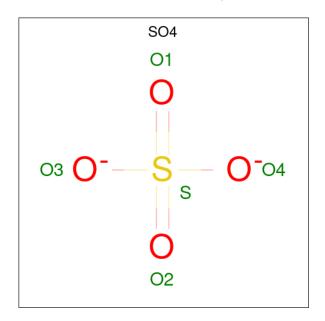


Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf
2	Λ	1	Total	С	Ν	0	S	0	0
0	A	1	7	3	1	2	1	0	0

 $\bullet\,$ Molecule 4 is NITRIC OXIDE (three-letter code: NO) (formula: NO).



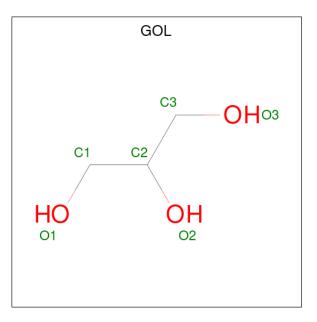
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	А	1	Total N 2 1	O 1	0	0





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
5	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
5	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 6 3 3 \end{array}$	0	0

• Molecule 7 is water.

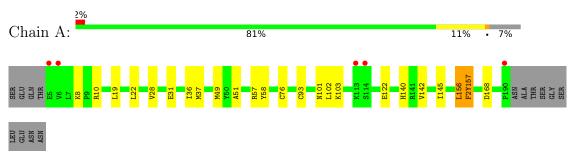


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	191	Total O 191 191	0	0



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: Cysteine dioxygenase type 1



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants	131.31Å 131.31Å 34.17Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.83 - 1.96	Depositor
Resolution (A)	32.83 - 1.96	EDS
% Data completeness	100.0 (32.83-1.96)	Depositor
(in resolution range)	94.0 (32.83-1.96)	EDS
R _{merge}	0.14	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.54 (at 1.97 \text{\AA})$	Xtriage
Refinement program	PHENIX	Depositor
D D.	0.161 , 0.188	Depositor
R, R_{free}	0.160 , 0.188	DCC
R_{free} test set	2004 reflections $(8.12%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	24.7	Xtriage
Anisotropy	0.332	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36 , 46.4	EDS
L-test for twinning ²	$< L > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	1823	wwPDB-VP
Average B, all atoms $(Å^2)$	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.87% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: FE, NO, GOL, SO4, F2Y

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).

Mal	Chain		lengths		angles
	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.39	0/1610	0.58	0/2168

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	А	1565	0	1536	18	0
2	А	1	0	0	0	0
3	А	7	0	3	0	0
4	А	2	0	0	0	0
5	А	15	0	0	0	0
6	А	42	0	56	3	0
7	А	191	0	0	5	2
All	All	1823	0	1595	18	2

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 6.

All (18) 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:93[B]:CYS:SG	1:A:157:F2Y:F1	2.24	0.85
1:A:10[A]:ARG:NH2	7:A:401:HOH:O	2.10	0.84
1:A:168:ASP:HB2	6:A:307:GOL:H31	1.74	0.70
1:A:57[A]:ARG:NH2	7:A:403:HOH:O	2.28	0.66
1:A:8:LYS:HE2	7:A:460:HOH:O	2.11	0.50
1:A:8:LYS:NZ	7:A:407:HOH:O	2.45	0.50
1:A:49[B]:MET:HE3	1:A:49[B]:MET:H	1.77	0.49
1:A:103:LYS:HE3	1:A:122:GLU:HG2	1.94	0.49
1:A:101:ASN:HB2	1:A:145:ILE:HG12	1.95	0.49
1:A:58:TYR:OH	6:A:308:GOL:H12	2.12	0.48
1:A:37[B]:MET:SD	1:A:156[B]:LEU:HD12	2.55	0.46
1:A:19:LEU:HD13	1:A:156[A]:LEU:HD13	1.96	0.46
1:A:31:GLU:HG3	7:A:525:HOH:O	2.15	0.46
1:A:51:ALA:HA	1:A:76:CYS:SG	2.58	0.43
1:A:8:LYS:NZ	6:A:313:GOL:H2	2.33	0.43
1:A:22:LEU:HD11	1:A:36:ILE:HD12	2.01	0.42
1:A:102:LEU:HD13	1:A:142:VAL:CG1	2.51	0.40
1:A:19:LEU:HD22	1:A:156[B]:LEU:HD11	2.04	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:552:HOH:O	7:A:564:HOH:O[1_554]	2.14	0.06
7:A:571:HOH:O	7:A:587:HOH:O[4_665]	2.16	0.04

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	Percentile	s
1	А	192/200~(96%)	190 (99%)	2(1%)	0	100 100	

There are no Ramachandran outliers to report.



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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed Rotameric		Outliers	Percentiles	
1	А	174/177~(98%)	170~(98%)	4 (2%)	50 42	

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	28	VAL
1	А	140	HIS
1	А	156[A]	LEU
1	А	156[B]	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

1 non-standard protein/DNA/RNA residue is 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	Bo	ond leng	ths	В	ond ang	les
	Type	Ullaili	nes	Link	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
1	F2Y	А	157	1	13,14,15	0.95	1 (7%)	$16,\!19,\!21$	2.08	5 (31%)

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
1	F2Y	А	157	1	-	2/5/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	157	F2Y	O-C	2.42	1.29	1.19

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	157	F2Y	CD2-CE2-CZ	-4.83	120.01	123.79
1	А	157	F2Y	F2-CE2-CZ	4.41	120.41	117.13
1	А	157	F2Y	CD1-CE1-CZ	-2.94	121.49	123.79
1	А	157	F2Y	CB-CA-C	-2.44	106.89	111.47
1	А	157	F2Y	CE2-CZ-CE1	2.05	119.60	116.68

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	А	157	F2Y	CA-CB-CG-CD1
1	А	157	F2Y	CA-CB-CG-CD2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	А	157	F2Y	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 for Mogul analysis.



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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).

Mal	Trune	Chain	Dec	Link	В	ond leng	gths	В	ond ang	gles
Mol	Type	Chain	Res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
6	GOL	А	310	-	$5,\!5,\!5$	1.25	1 (20%)	$5,\!5,\!5$	0.90	0
4	NO	А	303	2	0,1,1	-	-	-		
6	GOL	А	313	-	$5,\!5,\!5$	1.06	0	$5,\!5,\!5$	0.95	0
6	GOL	А	309	-	$5,\!5,\!5$	1.13	0	$5,\!5,\!5$	0.93	0
3	CYS	А	302	2	$5,\!6,\!6$	1.12	1 (20%)	5,7,7	1.56	2 (40%)
6	GOL	А	308	-	$5,\!5,\!5$	1.01	0	$5,\!5,\!5$	0.89	0
5	SO4	А	304	-	4,4,4	0.29	0	$6,\!6,\!6$	0.36	0
5	SO4	А	306	-	4,4,4	0.12	0	$6,\!6,\!6$	0.21	0
6	GOL	А	312	-	$5,\!5,\!5$	0.93	0	$5,\!5,\!5$	0.88	0
6	GOL	А	311	-	$5,\!5,\!5$	0.79	0	$5,\!5,\!5$	1.00	0
5	SO4	А	305	-	4,4,4	0.14	0	$6,\!6,\!6$	0.16	0
6	GOL	А	307	-	$5,\!5,\!5$	1.05	0	$5,\!5,\!5$	0.85	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
6	GOL	А	310	-	-	4/4/4/4	-
6	GOL	А	313	-	-	0/4/4/4	-
6	GOL	А	309	-	-	4/4/4/4	-
3	CYS	А	302	2	-	1/6/6/6	-
6	GOL	А	308	-	-	0/4/4/4	-
6	GOL	А	312	-	-	2/4/4/4	-
6	GOL	А	311	-	-	3/4/4/4	-
6	GOL	А	307	-	_	2/4/4/4	_

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
3	А	302	CYS	OXT-C	-2.36	1.22	1.30
6	А	310	GOL	C3-C2	2.23	1.60	1.51



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	302	CYS	OXT-C-O	-2.49	118.43	124.09
3	А	302	CYS	OXT-C-CA	2.03	120.30	113.38

All (2) bond angle outliers are listed below:

There are no chirality outliers.

 \mathbf{Mol} Chain Res Type Atoms 6 309 GOL C1-C2-C3-O3 А GOL 310 O1-C1-C2-C3 6 А GOL 6 311 O1-C1-C2-C3 А GOL 6 А 312 O1-C1-C2-C3 6 А 307 GOL O1-C1-C2-C3 GOL 6 А 309 O1-C1-C2-C3 6 А 310GOL C1-C2-C3-O3 GOL 01-C1-C2-O2 6 А 307 А GOL O1-C1-C2-O2 6 311 6309 GOL O1-C1-C2-O2 А 6 309 GOL <u>0</u>2-C2-C3-O3 А 6 А 310 GOL O2-C2-C3-O3 6 А 312 GOL O1-C1-C2-O2 6 А 311 GOL O2-C2-C3-O3 6 А 310GOL O1-C1-C2-O2 3 А 302 CYS OXT-C-CA-N

All (16) torsion outliers are listed below:

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	А	313	GOL	1	0
6	А	308	GOL	1	0
6	А	307	GOL	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 (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^{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		$\mathbf{OWAB}(\mathbf{A}^2)$	Q < 0.9	
1	А	185/200~(92%)	-0.16	5 (2%)	54	63	18, 26, 47, 81	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	5	GLU	4.5
1	А	113	LYS	3.7
1	А	190	PRO	2.9
1	А	6	VAL	2.6
1	А	114	SER	2.5

6.2 Non-standard residues in protein, DNA, RNA chains (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^{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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
1	F2Y	А	157	14/15	0.97	0.09	$20,\!27,\!39,\!39$	0

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,



Mol	Type	Chain	Res	Atoms	RSCC	RSR	B -factors($Å^2$)	Q < 0.9
6	GOL	А	313	6/6	0.64	0.64	$58,\!69,\!70,\!73$	0
6	GOL	А	309	6/6	0.65	0.29	52,54,59,62	0
6	GOL	А	307	6/6	0.73	0.41	$48,\!53,\!56,\!62$	0
6	GOL	А	312	6/6	0.81	0.23	52,59,62,67	0
6	GOL	А	310	6/6	0.84	0.21	35,45,49,50	0
6	GOL	А	308	6/6	0.85	0.13	42,45,51,51	0
6	GOL	А	311	6/6	0.87	0.38	$39,\!43,\!45,\!58$	0
3	CYS	А	302	7/7	0.96	0.08	$26,\!31,\!35,\!35$	0
5	SO4	А	306	5/5	0.98	0.32	$47,\!49,\!51,\!57$	0
5	SO4	А	305	5/5	0.98	0.18	47,53,60,62	0
5	SO4	А	304	5/5	0.99	0.08	22,35,44,44	0
2	FE	А	301	1/1	1.00	0.07	21,21,21,21	0
4	NO	А	303	2/2	1.00	0.07	29,29,29,31	0

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.

6.5 Other polymers (i)

There are no such residues in this entry.

