

Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	7S6T
Title	:	Complex structure of Methane monooxygenase hydroxylase and regulatory
		subunit H33A
Authors	:	Johns, J.C.; Banerjee, R.; Semonis, M.M.; Shi, K.; Aihara, H.; Lipscomb, J.D.
Deposited on	:	2021-09-14
Resolution	:	1.82 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org* A user guide is available at 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	:	2.35.1
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 (i)

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

The reported resolution of this entry is 1.82 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	Similar resolution
	$(\# { m Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$
R_{free}	130704	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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	
1	٨	F1F	3%	
1	A	515	94%	6%
1	Ε	515	93%	7%
	D	202	% •	
2	В	392	97%	•
		202	.%	
2	F,	392	95%	5%
	-		.% ■	
3	С	168	97%	•



Mol	Chain	Length	Quality of chain	
3	G	168	% 99%	·
4	D	136	90%	7% •
4	Н	136	94%	6%



2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 22073 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 Methane monooxygenase component A alpha chain.

Mol	Chain	Residues		At	oms		ZeroOcc	AltConf	Trace	
1	А	515	Total 4217	C 2700	N 737	O 768	S 12	0	4	0
1	Е	515	Total 4246	C 2715	N 738	0 780	S 13	0	8	0

• Molecule 2 is a protein called Methane monooxygenase beta chain.

Mol	Chain	Residues		Ate	oms		ZeroOcc	AltConf	Trace	
2	В	D 202	Total	С	Ν	Ο	S	0	1	0
	D	592	3190	2034	557	594	5	0	L	0
9	Б	202	Total	С	Ν	0	S	0	2	0
2	Г	392	3197	2038	558	596	5	0	Δ	0

• Molecule 3 is a protein called Methane monooxygenase gamma chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	С	168	Total 1371	C 880	N 236	0 254	S 1	0	1	0
3	G	168	Total 1371	C 880	N 236	0 254	S 1	0	1	0

• Molecule 4 is a protein called Methane monooxygenase regulatory protein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	а	D 121	Total	С	Ν	0	S	0	0	0
4	D	101	991	633	159	196	3	0		
4	ц	136	Total	С	Ν	0	S	0	0	0
4	4 H		1027	656	164	204	3	0	0	0

There are 2 discrepancies between the modelled and reference sequences:



Chain	Residue Modelled Actual		Comment	Reference	
D	33	ALA	HIS	engineered mutation	UNP A0A2D2D0T8
Н	33	ALA	HIS	engineered mutation	UNP A0A2D2D0T8

• Molecule 5 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	2	Total Fe 2 2	0	0
5	Е	2	Total Fe 2 2	0	0



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	С	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	С	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	С	1	$\begin{array}{c c} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	D	1	$\begin{array}{ccc} \overline{\text{Total}} & C & O \\ 4 & 2 & 2 \end{array}$	0	0
6	D	1	$\begin{array}{c cc} \overline{\text{Total}} & C & O \\ 4 & 2 & 2 \end{array}$	0	0
6	D	1	$\begin{array}{c ccc} \hline \text{Total} & \text{C} & \text{O} \\ \hline 4 & 2 & 2 \end{array}$	0	0
6	Е	1	$\begin{array}{c cc} \hline \text{Total} & \text{C} & \text{O} \\ \hline 4 & 2 & 2 \end{array}$	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	Е	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	Е	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 4 2 2 \end{array}$	0	0
6	Е	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 4 2 2 \end{array}$	0	0
6	Е	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 4 2 2 \end{array}$	0	0
6	Е	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	Ε	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	Ε	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	Ε	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 4 2 2 \end{array}$	0	0
6	F	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 4 2 2 \end{array}$	0	0
6	G	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	G	1	$\begin{array}{ccc} \overline{\text{Total}} & \mathcal{C} & \mathcal{O} \\ 4 & 2 & 2 \end{array}$	0	0
6	G	1	$\begin{array}{c ccc} \hline \text{Total} & \text{C} & \text{O} \\ \hline 4 & 2 & 2 \end{array}$	0	0
6	G	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 4 & 2 & 2 \end{array}$	0	0



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
6	Н	1	Total 4	${ m C} 2$	O 2	0	0

• Molecule 7 is BENZOIC ACID (three-letter code: BEZ) (formula: $C_7H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 9 & 7 & 2 \end{array}$	0	0
7	Е	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 9 & 7 & 2 \end{array}$	0	0

• Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	А	445	Total O 446 446	0	1
8	В	420	Total O 421 421	0	1
8	С	198	Total O 198 198	0	0
8	D	64	$\begin{array}{cc} \text{Total} & \text{O} \\ 64 & 64 \end{array}$	0	0
8	Ε	428	Total O 432 432	0	4
8	F	410	Total O 410 410	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	G	200	Total O 200 200	0	0
8	Н	70	Total O 70 70	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: Methane monooxygenase component A alpha chain



• Molecule 1: Methane monooxygenase component A alpha chain



• Molecule 2: Methane monooxygenase beta chain



• Molecule 3: Methane monooxygenase gamma chain





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	102.52Å 105.60Å 299.98Å	Deperitor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
$\mathbf{P}_{\text{assolution}}(\hat{\mathbf{A}})$	48.51 - 1.82	Depositor
Resolution (A)	48.51 - 1.82	EDS
% Data completeness	99.4 (48.51-1.82)	Depositor
(in resolution range)	99.4 (48.51-1.82)	EDS
R _{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.73 (at 1.82 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
P. P.	0.146 , 0.172	Depositor
n, n_{free}	0.146 , 0.172	DCC
R_{free} test set	14182 reflections (4.90%)	wwPDB-VP
Wilson B-factor $(Å^2)$	28.6	Xtriage
Anisotropy	0.539	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.33 , 49.2	EDS
L-test for twinning ²	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	22073	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 2.05% 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, BEZ, 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).

Mal	Chain	Bond	lengths	Bond angles	
MOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.55	0/4347	0.66	0/5903
1	Е	0.53	0/4376	0.66	1/5944~(0.0%)
2	В	0.56	0/3283	0.66	0/4464
2	F	0.54	0/3290	0.65	0/4476
3	С	0.47	0/1397	0.62	0/1888
3	G	0.50	0/1397	0.62	0/1888
4	D	0.47	0/1006	0.58	0/1360
4	Н	0.45	0/1042	0.57	0/1410
All	All	0.53	0/20138	0.64	1/27333~(0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1
1	Е	0	2
All	All	0	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	Е	187	VAL	CG1-CB-CG2	5.31	119.40	110.90

There are no chirality outliers.

All (3) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	А	489	ARG	Sidechain
1	Ε	186	ARG	Sidechain
1	Е	489	ARG	Sidechain

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	4217	0	4020	17	0
1	Е	4246	0	4029	28	0
2	В	3190	0	3019	10	0
2	F	3197	0	3023	17	0
3	С	1371	0	1412	5	0
3	G	1371	0	1412	2	0
4	D	991	0	988	6	0
4	Н	1027	0	1026	4	0
5	А	2	0	0	0	0
5	Е	2	0	0	0	0
6	А	56	0	84	9	0
6	В	28	0	42	1	0
6	С	12	0	18	0	0
6	D	12	0	18	3	0
6	Ε	36	0	54	4	0
6	F	36	0	54	2	0
6	G	16	0	24	0	0
6	Н	4	0	6	0	0
7	А	9	0	5	0	0
7	Ε	9	0	5	0	0
8	А	446	0	0	5	0
8	В	421	0	0	2	0
8	С	198	0	0	3	0
8	D	64	0	0	1	0
8	E	432	0	0	9	0
8	F	410	0	0	7	0
8	G	200	0	0	0	0
8	Н	70	0	0	0	0
All	All	22073	0	19239	81	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.

All (81) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:192:PHE:O	8:A:701:HOH:O	2.02	0.77
1:E:192:PHE:O	8:E:701:HOH:O	2.03	0.75
6:A:609:EDO:O2	8:A:702:HOH:O	2.03	0.74
6:F:409:EDO:O1	8:F:501:HOH:O	2.08	0.71
4:D:3:SER:N	8:D:301:HOH:O	2.23	0.70
2:B:321[A]:ASN:ND2	8:B:502:HOH:O	2.26	0.68
1:E:333:ARG:NH2	4:H:27:GLU:OE2	2.28	0.67
1:A:493[A]:LYS:NZ	8:A:704:HOH:O	2.28	0.66
2:F:321[A]:ASN:ND2	8:F:503:HOH:O	2.28	0.65
1:A:504:ASP:OD2	1:A:505:ASN:ND2	2.27	0.65
1:A:333:ARG:NH2	4:D:27:GLU:OE2	2.28	0.65
1:E:187:VAL:HG13	1:E:281:TYR:HB3	1.79	0.64
6:F:409:EDO:O2	8:F:502:HOH:O	2.15	0.64
6:A:617:EDO:H12	8:E:842:HOH:O	1.97	0.63
1:A:49:LYS:NZ	1:A:270:ASP:OD2	2.33	0.61
2:B:80:HIS:CD2	3:C:141:MET:HG2	2.35	0.61
2:F:80:HIS:CD2	3:G:141:MET:HG2	2.36	0.60
2:F:295:ARG:NH2	8:F:504:HOH:O	2.33	0.60
1:E:480[A]:GLU:OE2	8:E:702:HOH:O	2.17	0.59
1:A:511:ASP:OD1	1:A:514[B]:ARG:NH2	2.36	0.57
1:E:138:LEU:HD22	2:F:163:PHE:CE1	2.42	0.55
1:A:333:ARG:HD2	6:A:610:EDO:H22	1.88	0.55
3:C:36:PHE:HA	8:C:462:HOH:O	2.07	0.54
1:E:70[B]:MET:HE2	4:H:102:TYR:HB3	1.91	0.53
1:E:406:GLN:NE2	8:E:706:HOH:O	2.31	0.53
1:A:200:CYS:HA	1:A:203:ASN:OD1	2.10	0.52
6:A:609:EDO:H11	8:A:823:HOH:O	2.10	0.52
6:A:608:EDO:H11	8:A:963:HOH:O	2.08	0.51
1:E:200:CYS:HA	1:E:203:ASN:OD1	2.12	0.50
1:A:354:TRP:CG	1:A:355:PRO:HD3	2.47	0.49
1:E:354:TRP:CG	1:E:355:PRO:HD3	2.48	0.49
1:E:133:GLN:NE2	8:E:720:HOH:O	2.46	0.49
1:E:361:LEU:HD21	6:E:607:EDO:H11	1.96	0.47
1:A:240:GLU:HG3	4:D:109:SER:O	2.14	0.47
2:B:325:GLU:OE2	8:B:501:HOH:O	2.20	0.47
2:F:189:LYS:HA	2:F:189:LYS:HD3	1.70	0.47
6:A:617:EDO:H21	1:E:83:LEU:HD12	1.96	0.47
6:E:609:EDO:H22	8:E:751:HOH:O	2.13	0.47



		Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:138:LEU:HD22	2:B:163:PHE:CE1	2.50	0.46
1:E:240:GLU:HG3	4:H:109:SER:O	2.14	0.46
1:E:245[B]:ARG:NH1	8:E:708:HOH:O	2.37	0.45
1:A:466:CYS:HB2	2:B:76:THR:HA	1.99	0.45
6:A:617:EDO:H11	1:E:79:PHE:HB3	1.99	0.45
1:E:333:ARG:NH1	8:E:725:HOH:O	2.50	0.45
1:E:187:VAL:CG1	1:E:281:TYR:HB3	2.45	0.44
2:B:189:LYS:HA	2:B:189:LYS:HD3	1.73	0.44
6:E:605:EDO:H22	8:E:970:HOH:O	2.18	0.44
6:A:608:EDO:C2	2:B:77:GLN:HE22	2.30	0.44
6:D:202:EDO:H22	1:E:88:ARG:O	2.18	0.44
2:B:121:ARG:HD3	2:F:118:GLU:OE2	2.18	0.43
4:D:99:PHE:HE1	6:D:201:EDO:H11	1.83	0.43
4:H:5:HIS:HE1	4:H:81:ASP:OD1	2.00	0.43
4:D:5:HIS:HE1	4:D:81:ASP:OD1	2.01	0.43
1:A:440:GLU:O	6:A:616:EDO:H22	2.19	0.43
1:E:352:ALA:HA	1:E:404:PRO:HB2	2.01	0.43
2:F:57:CYS:HB3	8:F:839:HOH:O	2.19	0.43
1:E:207:VAL:O	1:E:211:CYS:HB3	2.19	0.43
1:E:296:PHE:HA	6:E:608:EDO:H12	2.00	0.42
3:C:121:PRO:HD3	3:C:129:PHE:CG	2.54	0.42
3:C:36:PHE:HB2	8:C:467:HOH:O	2.17	0.42
1:A:44:THR:HB	1:A:127:SER:HA	2.01	0.42
1:E:123:MET:HG3	2:F:171:ARG:HD3	2.01	0.42
1:E:184:MET:HA	1:E:187:VAL:CG2	2.50	0.42
1:E:504:ASP:OD2	1:E:505:ASN:ND2	2.52	0.42
1:E:138:LEU:HD22	2:F:163:PHE:HE1	1.82	0.42
4:D:99:PHE:CE1	6:D:201:EDO:H11	2.55	0.41
1:A:276:TRP:CE3	1:A:331:SER:HB2	2.55	0.41
2:B:191:ASP:O	2:B:195:MET:HG2	2.21	0.41
2:F:79:PHE:HB2	2:F:83:ARG:HB3	2.01	0.41
1:E:466:CYS:HB2	2:F:76:THR:HA	2.02	0.41
1:E:490[A]:SER:OG	2:F:35:GLN:HB2	2.20	0.41
2:B:45:ARG:HD3	6:B:404:EDO:H22	2.03	0.41
2:F:100:ARG:NH1	8:F:521:HOH:O	2.53	0.41
2:F:358:SER:O	2:F:362:GLN:HG3	2.21	0.41
1:A:49:LYS:HB2	1:A:49:LYS:HE3	1.86	0.40
1:A:185:LYS:O	1:A:189:ALA:HB3	2.21	0.40
2:F:191:ASP:O	2:F:195:MET:HG2	2.20	0.40
1:E:323:LYS:HA	1:E:323:LYS:HD2	1.91	0.40
2:F:37:LYS:NZ	8:F:510:HOH:O	2.42	0.40



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:40:TYR:HE2	8:C:462:HOH:O	2.05	0.40
2:F:80:HIS:CG	3:G:141:MET:HG2	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	entiles
1	А	517/515~(100%)	499~(96%)	18 (4%)	0	100	100
1	Е	521/515~(101%)	506 (97%)	15 (3%)	0	100	100
2	В	391/392~(100%)	381~(97%)	10 (3%)	0	100	100
2	F	392/392~(100%)	382 (97%)	10 (3%)	0	100	100
3	С	167/168~(99%)	165~(99%)	2(1%)	0	100	100
3	G	167/168~(99%)	165~(99%)	2(1%)	0	100	100
4	D	129/136~(95%)	124 (96%)	4 (3%)	1 (1%)	19	7
4	Н	134/136~(98%)	129 (96%)	4 (3%)	1 (1%)	22	10
All	All	2418/2422~(100%)	2351 (97%)	65 (3%)	2(0%)	51	37

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	73	ALA
4	Н	73	ALA

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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.



Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	430/426~(101%)	425~(99%)	5(1%)	71 64
1	Ε	434/426~(102%)	428~(99%)	6 (1%)	67 58
2	В	324/324~(100%)	321~(99%)	3~(1%)	78 74
2	F	325/324~(100%)	321~(99%)	4 (1%)	71 64
3	С	146/145~(101%)	146 (100%)	0	100 100
3	G	146/145~(101%)	145~(99%)	1 (1%)	84 80
4	D	103/107~(96%)	101~(98%)	2(2%)	57 45
4	Н	107/107~(100%)	105 (98%)	2 (2%)	57 45
All	All	2015/2004~(100%)	1992 (99%)	23 (1%)	71 67

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	32	LEU
1	А	52	MET
1	А	112	VAL
1	А	307	ARG
1	А	437	ARG
2	В	163	PHE
2	В	176	ASP
2	В	270	ARG
4	D	114	ARG
4	D	129	MET
1	Е	52	MET
1	Е	112	VAL
1	Е	187	VAL
1	Е	244	LEU
1	Е	307	ARG
1	Е	437	ARG
2	F	163	PHE
2	F	176	ASP
2	F	206	LEU
2	F	270	ARG
3	G	31	LYS
4	Н	57	LEU
4	Н	114	ARG



Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
4	D	5	HIS
4	Н	5	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 56 ligands modelled in this entry, 4 are monoatomic - leaving 52 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).

			T • 1	B	ond leng	gths	Bond angles			
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	F	409	-	3,3,3	0.38	0	2,2,2	0.53	0
6	EDO	В	407	-	3,3,3	0.59	0	2,2,2	0.30	0
6	EDO	А	612	-	3,3,3	0.58	0	2,2,2	0.17	0
6	EDO	Е	603	-	3,3,3	0.47	0	2,2,2	0.36	0
6	EDO	А	610	-	3,3,3	0.46	0	2,2,2	0.35	0
6	EDO	В	402	-	3,3,3	0.35	0	2,2,2	0.54	0
6	EDO	А	608	-	3,3,3	0.43	0	2,2,2	0.59	0
6	EDO	F	408	-	3,3,3	0.39	0	2,2,2	0.50	0
6	EDO	А	617	-	3,3,3	0.51	0	2,2,2	0.23	0
6	EDO	А	616	-	3,3,3	0.49	0	2,2,2	0.23	0



	T a	Chain	Dag	T : 1-	B	ond leng	$_{ m gths}$	В	ond ang	les
NIOI	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	В	403	-	3,3,3	0.41	0	$2,\!2,\!2$	0.54	0
6	EDO	В	405	-	3,3,3	0.51	0	$2,\!2,\!2$	0.26	0
6	EDO	Е	604	-	3,3,3	0.50	0	$2,\!2,\!2$	0.35	0
6	EDO	В	401	-	3,3,3	0.52	0	$2,\!2,\!2$	0.35	0
6	EDO	А	605	-	3,3,3	0.60	0	2,2,2	0.41	0
6	EDO	С	201	-	3,3,3	0.50	0	$2,\!2,\!2$	0.27	0
6	EDO	А	603	-	3,3,3	0.46	0	$2,\!2,\!2$	0.30	0
6	EDO	F	407	-	3,3,3	0.43	0	$2,\!2,\!2$	0.63	0
6	EDO	D	203	-	3,3,3	0.46	0	$2,\!2,\!2$	0.48	0
6	EDO	Е	607	-	3,3,3	0.65	0	$2,\!2,\!2$	0.02	0
6	EDO	G	204	-	3,3,3	0.43	0	$2,\!2,\!2$	0.27	0
6	EDO	А	613	-	3,3,3	0.65	0	$2,\!2,\!2$	0.16	0
6	EDO	G	201	-	3,3,3	0.51	0	$2,\!2,\!2$	0.35	0
6	EDO	F	405	-	3,3,3	0.44	0	2,2,2	0.46	0
6	EDO	Е	608	-	3,3,3	0.59	0	$2,\!2,\!2$	0.21	0
6	EDO	Н	201	-	3,3,3	0.52	0	2,2,2	0.21	0
6	EDO	С	203	-	3,3,3	0.48	0	2,2,2	0.32	0
6	EDO	D	201	-	3,3,3	0.51	0	2,2,2	0.19	0
6	EDO	В	406	-	3,3,3	0.72	0	2,2,2	0.55	0
6	EDO	А	614	-	3,3,3	0.49	0	2,2,2	0.29	0
6	EDO	F	402	-	3,3,3	0.53	0	2,2,2	0.33	0
7	BEZ	Е	606	5	9,9,9	0.89	0	11,11,11	0.92	0
6	EDO	Е	609	-	3,3,3	0.36	0	$2,\!2,\!2$	0.55	0
6	EDO	F	406	-	3,3,3	0.53	0	2,2,2	0.19	0
6	EDO	G	203	-	3,3,3	0.46	0	2,2,2	0.35	0
6	EDO	С	202	-	3,3,3	0.39	0	$2,\!2,\!2$	0.49	0
7	BEZ	А	607	5	9,9,9	0.82	0	11,11,11	1.02	0
6	EDO	F	403	-	3,3,3	0.43	0	$2,\!2,\!2$	0.61	0
6	EDO	Е	610	-	3,3,3	0.49	0	$2,\!2,\!2$	0.30	0
6	EDO	А	609	-	3,3,3	0.60	0	$2,\!2,\!2$	0.21	0
6	EDO	Е	611	-	3,3,3	0.57	0	$2,\!2,\!2$	0.15	0
6	EDO	А	615	-	3,3,3	0.44	0	2,2,2	0.50	0
6	EDO	D	202	-	3,3,3	0.58	0	$2,\!2,\!2$	0.11	0
6	EDO	Е	605	-	3,3,3	0.55	0	2,2,2	0.24	0
6	EDO	F	401	-	3,3,3	0.44	0	$2,\!2,\!2$	0.50	0
6	EDO	Е	612	-	3,3,3	0.51	0	$2,\!2,\!2$	0.30	0
6	EDO	G	202	-	3,3,3	0.40	0	$2,\!2,\!2$	0.50	0
6	EDO	F	404	-	3,3,3	0.41	0	2,2,2	0.50	0
6	EDO	А	606	-	3,3,3	0.55	0	$2,\!2,\!2$	0.27	0
6	EDO	А	604	-	3,3,3	0.52	0	2,2,2	0.74	0
6	EDO	В	404	-	3,3,3	0.49	0	$2,\!2,\!2$	0.26	0
6	EDO	A	611	-	3,3,3	0.52	0	$2,\!2,\!2$	0.31	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	EDO	F	409	-	-	1/1/1/1	-
6	EDO	В	407	-	-	0/1/1/1	-
6	EDO	А	612	-	-	0/1/1/1	-
6	EDO	Е	603	-	-	0/1/1/1	-
6	EDO	А	610	-	-	0/1/1/1	_
6	EDO	В	402	-	-	1/1/1/1	-
6	EDO	А	608	-	-	0/1/1/1	-
6	EDO	F	408	-	-	0/1/1/1	-
6	EDO	А	617	-	-	0/1/1/1	-
6	EDO	А	616	-	-	0/1/1/1	-
6	EDO	В	403	-	-	0/1/1/1	-
6	EDO	В	405	-	-	0/1/1/1	_
6	EDO	Е	604	-	_	1/1/1/1	_
6	EDO	В	401	-	-	0/1/1/1	_
6	EDO	А	605	-	_	0/1/1/1	-
6	EDO	С	201	-	-	0/1/1/1	-
6	EDO	А	603	-	-	0/1/1/1	_
6	EDO	F	407	-	-	1/1/1/1	-
6	EDO	D	203	-	_	1/1/1/1	_
6	EDO	Е	607	-	-	1/1/1/1	-
6	EDO	G	204	-	-	0/1/1/1	_
6	EDO	А	613	-	-	1/1/1/1	-
6	EDO	G	201	-	-	0/1/1/1	-
6	EDO	F	405	-	-	0/1/1/1	_
6	EDO	Е	608	-	-	1/1/1/1	-
6	EDO	Н	201	-	-	0/1/1/1	_
6	EDO	С	203	-	_	0/1/1/1	-
6	EDO	D	201	-	-	0/1/1/1	-
6	EDO	В	406	-	_	1/1/1/1	_
6	EDO	А	614	-	-	0/1/1/1	-
6	EDO	F	402	-	-	0/1/1/1	_
7	BEZ	Е	606	5	_	0/4/4/4	0/1/1/1
6	EDO	Е	609	-	-	0/1/1/1	-
6	EDO	F	406	-	-	0/1/1/1	_
6	EDO	G	203	-	-	0/1/1/1	-
6	EDO	С	202	-	-	1/1/1/1	-
7	BEZ	А	607	5	-	0/4/4/4	0/1/1/1
6	EDO	F	403	-	-	0/1/1/1	-
6	EDO	Е	610	-	-	0/1/1/1	-



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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	EDO	А	609	-	-	1/1/1/1	-
6	EDO	Е	611	-	-	0/1/1/1	-
6	EDO	А	615	-	-	0/1/1/1	-
6	EDO	D	202	-	-	0/1/1/1	-
6	EDO	Е	605	-	-	1/1/1/1	-
6	EDO	F	401	-	-	1/1/1/1	-
6	EDO	Е	612	-	-	0/1/1/1	-
6	EDO	G	202	-	-	0/1/1/1	-
6	EDO	F	404	-	-	0/1/1/1	-
6	EDO	А	606	-	-	0/1/1/1	-
6	EDO	А	604	-	-	0/1/1/1	-
6	EDO	В	404	-	-	0/1/1/1	-
6	EDO	А	611	-	-	0/1/1/1	_

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	401	EDO	O1-C1-C2-O2
6	F	409	EDO	O1-C1-C2-O2
6	А	609	EDO	O1-C1-C2-O2
6	Е	605	EDO	O1-C1-C2-O2
6	F	407	EDO	O1-C1-C2-O2
6	D	203	EDO	O1-C1-C2-O2
6	А	613	EDO	O1-C1-C2-O2
6	В	402	EDO	O1-C1-C2-O2
6	Ε	607	EDO	O1-C1-C2-O2
6	В	406	EDO	O1-C1-C2-O2
6	С	202	EDO	O1-C1-C2-O2
6	Ē	604	EDO	O1-C1-C2-O2
6	Е	608	EDO	O1-C1-C2-O2

There are no ring outliers.

13 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	F	409	EDO	2	0
6	А	610	EDO	1	0



Contre	Continuaca from pretious page										
Mol	Chain	Res	Type	Clashes	Symm-Clashes						
6	А	608	EDO	2	0						
6	А	617	EDO	3	0						
6	А	616	EDO	1	0						
6	Е	607	EDO	1	0						
6	Е	608	EDO	1	0						
6	D	201	EDO	2	0						
6	Е	609	EDO	1	0						
6	А	609	EDO	2	0						
6	D	202	EDO	1	0						
6	Е	605	EDO	1	0						
6	В	404	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 (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	0	$WAB(A^2)$	Q<0.9
1	А	515/515~(100%)	-0.30	13 (2%) 57 55	2 20	, 27, 41, 67	0
1	Ε	515/515~(100%)	-0.27	12 (2%) 60 5	3 21	, 27, 42, 68	0
2	В	392/392~(100%)	-0.36	5 (1%) 77 74	20	, 26, 37, 67	1 (0%)
2	F	392/392~(100%)	-0.36	3 (0%) 86 84	21	, 27, 37, 64	1 (0%)
3	С	168/168~(100%)	-0.48	2 (1%) 79 76	23	, 31, 43, 55	0
3	G	168/168~(100%)	-0.44	1 (0%) 89 88	24	, 31, 43, 53	0
4	D	131/136~(96%)	0.35	6 (4%) 32 27	27	, 34, 49, 81	0
4	Н	136/136~(100%)	0.51	15 (11%) 5 4	27	, 35, 56, 85	0
All	All	2417/2422 (99%)	-0.25	57 (2%) 59 54	1 20	, 28, 43, 85	2 (0%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	Н	135	LEU	6.8
4	Н	138	ILE	6.1
4	Н	133	ARG	5.7
4	Н	134	ALA	5.3
4	D	133	ARG	4.9
2	F	4	PRO	4.2
4	D	3	SER	4.2
2	В	4	PRO	3.7
2	F	395	ASN	3.6
1	А	193	ILE	3.3
1	А	53	ALA	3.2
1	Е	259	ASN	3.1
4	Н	136	THR	3.1
4	Н	118	LEU	3.1
1	A	187	VAL	3.0
1	Е	53	ALA	3.0



Mol	Chain	Res	Type	RSRZ
3	С	169	ALA	3.0
1	Е	141	VAL	3.0
3	С	2	ALA	2.8
4	D	4	ALA	2.8
4	D	62	ALA	2.8
1	А	118	ILE	2.7
1	Е	189	ALA	2.7
1	Е	112	VAL	2.7
3	G	169	ALA	2.6
1	Е	193	ILE	2.6
1	А	141	VAL	2.6
4	D	118	LEU	2.6
2	F	86	TRP	2.6
2	В	395	ASN	2.5
4	Н	62	ALA	2.5
1	А	259	ASN	2.5
1	Е	261	PRO	2.5
1	А	120	ALA	2.4
1	Е	118	ILE	2.4
1	Е	145	ILE	2.4
4	Н	67	ILE	2.4
1	А	145	ILE	2.4
4	Н	14	GLN	2.4
4	Н	4	ALA	2.4
2	В	86	TRP	2.3
4	D	59	GLY	2.3
4	Н	3	SER	2.3
4	Н	137	ASP	2.3
1	А	189	ALA	2.3
1	А	261	PRO	2.2
1	Е	119	ALA	2.2
1	А	142	LEU	2.2
1	Е	187	VAL	2.2
4	Н	119	GLY	2.1
2	В	169	VAL	2.1
4	Н	16	THR	2.1
4	Н	65	PRO	2.1
2	В	75	TRP	2.1
1	А	55	GLU	2.1
1	Е	120	ALA	2.0
1	А	323	LYS	2.0

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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^{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	$B-factors(Å^2)$	Q<0.9
6	EDO	А	610	4/4	0.61	0.37	98,98,98,100	0
6	EDO	А	617	4/4	0.61	0.23	55,58,60,63	0
6	EDO	Е	608	4/4	0.64	0.20	52,54,55,56	0
6	EDO	А	614	4/4	0.66	0.21	75,76,79,80	0
6	EDO	В	406	4/4	0.67	0.21	38,45,49,50	0
6	EDO	А	613	4/4	0.70	0.16	46,51,54,58	0
6	EDO	Е	611	4/4	0.72	0.16	52,55,59,60	0
6	EDO	В	405	4/4	0.73	0.15	69,69,70,73	0
6	EDO	А	609	4/4	0.73	0.24	53,56,57,61	0
6	EDO	F	409	4/4	0.75	0.24	68,68,68,70	0
6	EDO	С	202	4/4	0.78	0.18	62,63,66,70	0
6	EDO	G	203	4/4	0.81	0.15	67,68,68,69	0
6	EDO	F	404	4/4	0.83	0.11	60,61,63,65	0
6	EDO	F	402	4/4	0.84	0.19	56,60,61,63	0
6	EDO	А	616	4/4	0.84	0.15	56,59,63,67	0
6	EDO	D	201	4/4	0.85	0.18	51,53,56,58	0
6	EDO	D	203	4/4	0.85	0.25	67,68,69,70	0
6	EDO	В	403	4/4	0.86	0.22	61,63,63,67	0
6	EDO	А	606	4/4	0.86	0.11	45,48,49,57	0
6	EDO	А	608	4/4	0.86	0.38	51,53,54,55	0
6	EDO	D	202	4/4	0.88	0.15	42,45,46,48	0
6	EDO	В	401	4/4	0.88	0.18	40,43,47,51	0
6	EDO	Е	605	4/4	0.88	0.22	45,47,48,49	0
6	EDO	В	407	4/4	0.89	0.15	32,36,47,55	0
6	EDO	Е	609	4/4	0.90	0.27	59,60,60,66	0
6	EDO	В	402	4/4	0.90	0.17	45,46,46,49	0
6	EDO	G	204	4/4	0.90	0.19	70,70,71,71	0



Mol	Type	Chain	Res	Atoms	RSCC	RSR	B -factors($Å^2$)	Q<0.9
6	EDO	Е	607	4/4	0.91	0.16	44,50,55,60	0
6	EDO	Е	603	4/4	0.91	0.18	44,49,51,58	0
6	EDO	А	611	4/4	0.91	0.12	38,41,41,43	0
6	EDO	С	203	4/4	0.92	0.17	56, 56, 58, 61	0
6	EDO	G	202	4/4	0.92	0.29	56, 56, 56, 58	0
6	EDO	Е	604	4/4	0.92	0.14	40,40,41,41	0
6	EDO	В	404	4/4	0.92	0.12	44,49,51,54	0
6	EDO	А	604	4/4	0.93	0.09	35,39,40,41	0
6	EDO	F	406	4/4	0.93	0.09	45,46,50,50	0
6	EDO	F	403	4/4	0.93	0.14	36,38,41,42	0
6	EDO	А	612	4/4	0.95	0.14	40,46,51,51	0
6	EDO	А	615	4/4	0.95	0.14	$28,\!36,\!45,\!52$	0
6	EDO	F	405	4/4	0.95	0.11	61,62,66,69	0
6	EDO	А	605	4/4	0.95	0.09	32,33,34,40	0
6	EDO	Н	201	4/4	0.95	0.17	45,46,48,48	0
6	EDO	F	408	4/4	0.96	0.11	49,49,53,58	0
6	EDO	Е	610	4/4	0.96	0.12	$28,\!36,\!46,\!52$	0
6	EDO	F	407	4/4	0.96	0.20	$55,\!56,\!56,\!57$	0
6	EDO	А	603	4/4	0.97	0.23	41,43,43,43	0
6	EDO	Е	612	4/4	0.97	0.16	39,41,43,43	0
7	BEZ	А	607	9/9	0.97	0.13	$26,\!28,\!30,\!31$	0
6	EDO	G	201	4/4	0.98	0.07	31,32,33,33	0
6	EDO	С	201	4/4	0.98	0.06	$29,\!31,\!31,\!31$	0
6	EDO	F	401	4/4	0.98	0.09	38,39,42,50	0
7	BEZ	Е	606	9/9	0.98	0.14	29,31,32,33	0
5	FE	A	601	1/1	1.00	0.08	$25,\!25,\!25,\!25$	0
5	FE	A	602	1/1	1.00	0.08	28,28,28,28	0
5	FE	Е	601	1/1	1.00	0.09	27,27,27,27	0
5	FE	Е	602	1/1	1.00	0.07	$2\overline{8,28,28,28}$	0

6.5 Other polymers (i)

There are no such residues in this entry.

