



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 05:36 pm BST

PDB ID : 3MBS  
Title : Crystal structure of 8mer PNA  
Authors : Yeh, J.I.; Pohl, E.; Truan, D.; He, W.; Sheldrick, G.M.; Achim, C.  
Deposited on : 2010-03-26  
Resolution : 1.27 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.

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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 : 2.11  
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) : 2.11

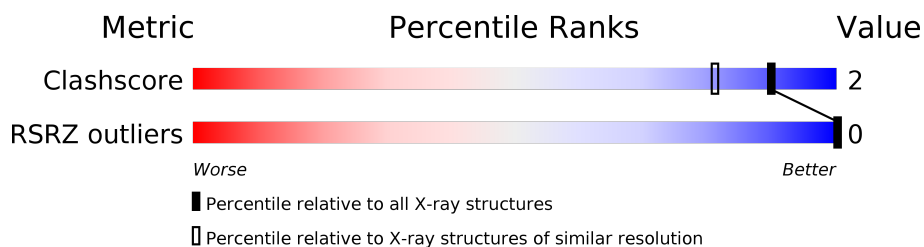
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 1.27 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))
Clashscore	141614	1926 (1.30-1.26)
RSRZ outliers	127900	1807 (1.30-1.26)

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 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	9	 89% 11%
1	B	9	 89% 11%
1	C	9	 89% 11%
1	D	9	 78% 11% 11%

## 2 Entry composition [i](#)

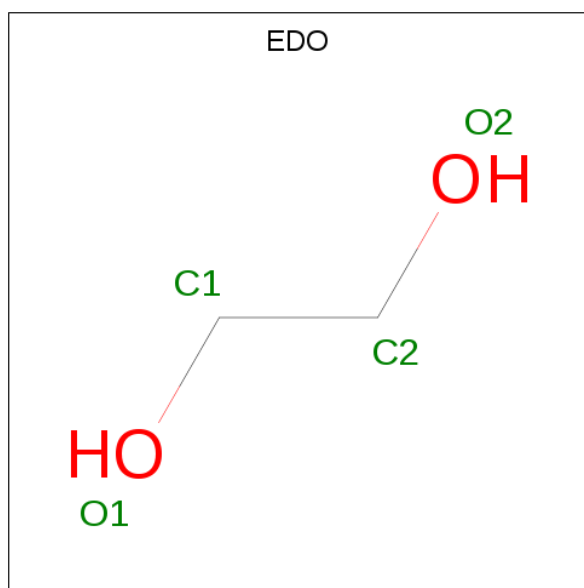
There are 3 unique types of molecules in this entry. The entry contains 892 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DNA chain called Peptide Nucleic Acid.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	9	Total	C	N	O	0	0	1
			157	85	48	24			
1	B	8	Total	C	N	O	0	0	0
			156	85	47	24			
1	C	9	Total	C	N	O	0	0	0
			159	87	48	24			
1	D	8	Total	C	N	O	0	0	0
			152	83	46	23			

- 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	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	61	Total	O	0	0
			61	61		
3	B	68	Total	O	0	0
			68	68		
3	C	60	Total	O	0	0
			60	60		
3	D	59	Total	O	0	0
			59	59		

### 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: Peptide Nucleic Acid

Chain A:  89% 11%




- Molecule 1: Peptide Nucleic Acid

Chain B:  89% 11%



- Molecule 1: Peptide Nucleic Acid

Chain C:  89% 11%



- Molecule 1: Peptide Nucleic Acid

Chain D:  78% 11% 11%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	18.87Å 28.92Å 54.97Å 88.10° 85.75° 79.56°	Depositor
Resolution (Å)	18.51 – 1.27 18.51 – 1.22	Depositor EDS
% Data completeness (in resolution range)	(Not available) (18.51-1.27) 82.6 (18.51-1.22)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.97 (at 1.22Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, $R_{free}$	0.179 , 0.220 0.184 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	10.9	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 91.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.59$ , $\langle L^2 \rangle = 0.44$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	892	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.18% 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: TPN, GPN, APN, EDO, CPN

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	C	7.76	1/2 (50.0%)	7.15	1/1 (100.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	3009	LYS	N-CA	-10.84	1.24	1.46

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	3009	LYS	N-CA-C	7.15	130.32	111.00

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	A	157	0	107	1	0
1	B	156	0	107	0	0
1	C	159	0	107	0	0
1	D	152	0	98	1	0
2	A	8	0	12	0	0
2	B	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	8	0	12	0	0
3	A	61	0	0	1	0
3	B	68	0	0	0	0
3	C	60	0	0	0	0
3	D	59	0	0	1	0
All	All	892	0	449	2	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 (2) 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:1009:LYS:N	3:A:9006:HOH:O	2.48	0.47
1:D:4004:APN:H5'1	3:D:9085:HOH:O	2.21	0.41

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

There are no protein molecules in this entry.

### 5.3.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

32 non-standard protein/DNA/RNA residues 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
1	TPN	A	1005	1	16,19,20	1.26	1 (6%)	15,25,27	3.52	3 (20%)
1	CPN	B	2008	1	15,18,19	1.18	2 (13%)	16,23,25	1.64	2 (12%)
1	GPN	A	1001	1	18,22,23	1.09	1 (5%)	17,30,32	3.08	10 (58%)
1	APN	D	4004	1	18,21,22	0.87	0	15,28,30	2.05	7 (46%)
1	GPN	A	1002	1	18,22,23	1.28	2 (11%)	17,30,32	3.00	6 (35%)
1	GPN	D	4001	1	15,20,23	1.26	1 (6%)	16,28,32	3.04	6 (37%)
1	TPN	B	2005	1	16,19,20	1.16	1 (6%)	15,25,27	3.43	2 (13%)
1	TPN	C	3005	1	16,19,20	0.96	1 (6%)	15,25,27	3.10	4 (26%)
1	CPN	D	4007	1	15,18,19	1.10	2 (13%)	16,23,25	1.70	2 (12%)
1	GPN	B	2002	1	18,22,23	1.39	2 (11%)	17,30,32	2.88	8 (47%)
1	APN	C	3004	1	18,21,22	0.80	0	15,28,30	1.98	5 (33%)
1	GPN	B	2006	1	18,22,23	1.48	1 (5%)	17,30,32	2.74	6 (35%)
1	CPN	D	4008	1	13,16,19	1.07	1 (7%)	15,21,25	1.24	2 (13%)
1	GPN	C	3001	1	18,22,23	1.29	2 (11%)	17,30,32	3.06	9 (52%)
1	CPN	A	1003	1	15,18,19	1.02	1 (6%)	16,23,25	1.86	5 (31%)
1	APN	A	1004	1	18,21,22	0.93	0	15,28,30	1.40	3 (20%)
1	CPN	B	2007	1	15,18,19	0.86	1 (6%)	16,23,25	1.58	3 (18%)
1	CPN	B	2003	1	15,18,19	1.10	1 (6%)	16,23,25	1.64	3 (18%)
1	GPN	A	1006	1	18,22,23	1.42	2 (11%)	17,30,32	2.73	9 (52%)
1	APN	B	2004	1	18,21,22	0.99	1 (5%)	15,28,30	1.85	6 (40%)
1	GPN	C	3006	1	18,22,23	1.19	1 (5%)	17,30,32	2.92	5 (29%)
1	GPN	D	4006	1	18,22,23	1.30	2 (11%)	17,30,32	3.63	12 (70%)
1	CPN	A	1007	1	15,18,19	1.09	2 (13%)	16,23,25	1.78	3 (18%)
1	CPN	C	3003	1	15,18,19	1.07	1 (6%)	16,23,25	2.05	4 (25%)
1	CPN	A	1008	1	15,18,19	1.07	1 (6%)	16,23,25	2.05	4 (25%)
1	CPN	C	3007	1	15,18,19	1.12	2 (13%)	16,23,25	1.78	5 (31%)
1	GPN	B	2001	1	18,22,23	1.26	2 (11%)	17,30,32	3.02	11 (64%)
1	GPN	C	3002	1	18,22,23	1.24	3 (16%)	17,30,32	2.56	6 (35%)
1	TPN	D	4005	1	16,19,20	1.40	2 (12%)	15,25,27	3.99	4 (26%)
1	CPN	D	4003	1	15,18,19	1.14	2 (13%)	16,23,25	1.66	2 (12%)
1	GPN	D	4002	1	18,22,23	1.40	2 (11%)	17,30,32	2.69	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	CPN	C	3008	1	15,18,19	1.48	3 (20%)	16,23,25	1.95	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	TPN	A	1005	1	-	0/13/14/15	0/1/1/1
1	CPN	B	2008	1	-	0/13/14/15	0/1/1/1
1	GPN	A	1001	1	-	1/13/14/15	0/2/2/2
1	APN	D	4004	1	-	0/13/14/15	0/2/2/2
1	GPN	A	1002	1	-	0/13/14/15	0/2/2/2
1	GPN	D	4001	1	-	1/10/11/15	0/2/2/2
1	TPN	B	2005	1	-	0/13/14/15	0/1/1/1
1	TPN	C	3005	1	-	0/13/14/15	0/1/1/1
1	CPN	D	4007	1	-	0/13/14/15	0/1/1/1
1	GPN	B	2002	1	-	0/13/14/15	0/2/2/2
1	APN	C	3004	1	-	0/13/14/15	0/2/2/2
1	GPN	B	2006	1	-	0/13/14/15	0/2/2/2
1	CPN	D	4008	1	-	2/11/11/15	0/1/1/1
1	GPN	C	3001	1	-	0/13/14/15	0/2/2/2
1	CPN	A	1003	1	-	0/13/14/15	0/1/1/1
1	APN	A	1004	1	-	0/13/14/15	0/2/2/2
1	CPN	B	2007	1	-	0/13/14/15	0/1/1/1
1	CPN	B	2003	1	-	0/13/14/15	0/1/1/1
1	GPN	A	1006	1	-	0/13/14/15	0/2/2/2
1	APN	B	2004	1	-	0/13/14/15	0/2/2/2
1	GPN	C	3006	1	-	0/13/14/15	0/2/2/2
1	GPN	D	4006	1	-	0/13/14/15	0/2/2/2
1	CPN	A	1007	1	-	0/13/14/15	0/1/1/1
1	CPN	C	3003	1	-	0/13/14/15	0/1/1/1
1	CPN	A	1008	1	-	0/13/14/15	0/1/1/1
1	CPN	C	3007	1	-	0/13/14/15	0/1/1/1
1	GPN	B	2001	1	-	0/13/14/15	0/2/2/2
1	GPN	C	3002	1	-	0/13/14/15	0/2/2/2
1	TPN	D	4005	1	-	0/13/14/15	0/1/1/1
1	CPN	D	4003	1	-	0/13/14/15	0/1/1/1
1	GPN	D	4002	1	-	0/13/14/15	0/2/2/2
1	CPN	C	3008	1	-	0/13/14/15	0/1/1/1

The worst 5 of 46 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2006	GP	C6-N1	5.74	1.43	1.33
1	D	4005	TP	C4-N3	4.37	1.40	1.33
1	B	2002	GP	C6-N1	4.33	1.40	1.33
1	D	4002	GP	C6-N1	4.22	1.40	1.33
1	A	1006	GP	C6-N1	4.17	1.40	1.33

The worst 5 of 167 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1005	TP	C4-N3-C2	12.77	125.92	115.14
1	B	2005	TP	C4-N3-C2	12.45	125.66	115.14
1	D	4005	TP	C4-N3-C2	11.44	124.80	115.14
1	C	3005	TP	C4-N3-C2	10.81	124.27	115.14
1	D	4005	TP	C2'-C3'-N4'	9.14	130.41	113.11

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1001	GP	C2'-C3'-N4'-C5'
1	D	4008	CP	C8'-C7'-N4'-C5'
1	D	4001	GP	O7'-C7'-N4'-C3'
1	D	4008	CP	O7'-C7'-N4'-C5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	4004	AP	1	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	EDO	A	5005	-	3,3,3	0.76	0	2,2,2	0.57	0
2	EDO	A	5003	-	3,3,3	0.75	0	2,2,2	0.51	0
2	EDO	D	5001	-	3,3,3	0.62	0	2,2,2	0.17	0
2	EDO	D	5002	-	3,3,3	0.55	0	2,2,2	0.29	0
2	EDO	B	5004	-	3,3,3	0.54	0	2,2,2	0.41	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	5005	-	-	0/1/1/1	-
2	EDO	A	5003	-	-	0/1/1/1	-
2	EDO	D	5001	-	-	0/1/1/1	-
2	EDO	D	5002	-	-	0/1/1/1	-
2	EDO	B	5004	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	1/9 (11%)	0.44	0 100 100	37, 37, 37, 37	0
1	B	0/9	-	-	-	-
1	C	1/9 (11%)	0.02	0 100 100	17, 17, 17, 17	0
1	D	0/9	-	-	-	-
All	All	2/36 (5%)	0.23	0 100 100	17, 17, 17, 37	0

There are no RSRZ outliers to report.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
1	CPN	A	1008	18/19	0.95	0.07	14,17,25,48	0
1	CPN	D	4007	18/19	0.96	0.08	13,15,22,23	0
1	GPN	C	3001	21/22	0.96	0.07	14,17,21,23	0
1	GPN	D	4006	21/22	0.96	0.08	11,14,20,26	0
1	TPN	A	1005	19/20	0.96	0.07	12,15,20,20	0
1	GPN	B	2001	21/22	0.96	0.07	15,18,19,20	0
1	TPN	D	4005	19/20	0.96	0.07	14,16,20,23	0
1	CPN	C	3008	18/19	0.96	0.07	8,10,13,21	0
1	APN	C	3004	20/21	0.97	0.06	11,14,19,22	0
1	GPN	B	2006	21/22	0.97	0.06	7,8,10,11	0
1	CPN	D	4008	16/19	0.97	0.06	14,16,24,24	0
1	GPN	A	1001	21/22	0.97	0.06	9,12,16,16	0
1	CPN	A	1003	18/19	0.97	0.06	8,10,15,18	0
1	APN	A	1004	20/21	0.97	0.06	10,11,18,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
1	GPN	A	1006	21/22	0.97	0.07	12,14,18,23	0
1	APN	B	2004	20/21	0.97	0.06	13,14,22,23	0
1	APN	D	4004	20/21	0.97	0.06	9,12,19,25	0
1	CPN	A	1007	18/19	0.97	0.07	13,15,22,32	0
1	CPN	C	3003	18/19	0.97	0.06	12,14,18,21	0
1	GPN	A	1002	21/22	0.97	0.06	8,10,12,12	0
1	GPN	D	4001	19/22	0.97	0.06	11,13,15,18	0
1	GPN	C	3002	21/22	0.97	0.06	13,15,17,18	0
1	CPN	B	2008	18/19	0.97	0.06	8,10,18,18	0
1	GPN	B	2002	21/22	0.97	0.05	13,16,17,18	0
1	TPN	C	3005	19/20	0.98	0.06	11,14,19,19	0
1	CPN	C	3007	18/19	0.98	0.06	6,8,11,11	0
1	GPN	C	3006	21/22	0.98	0.06	7,9,11,11	0
1	CPN	B	2007	18/19	0.98	0.06	7,8,10,11	0
1	CPN	B	2003	18/19	0.98	0.06	11,14,17,21	0
1	CPN	D	4003	18/19	0.98	0.06	9,10,12,13	0
1	GPN	D	4002	21/22	0.98	0.06	8,10,11,12	0
1	TPN	B	2005	19/20	0.98	0.05	11,13,19,19	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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( $\text{\AA}^2$ )	Q<0.9
2	EDO	B	5004	4/4	0.95	0.07	17,19,19,25	0
2	EDO	A	5005	4/4	0.96	0.08	16,22,27,27	0
2	EDO	D	5002	4/4	0.98	0.06	13,14,15,17	0
2	EDO	D	5001	4/4	0.98	0.06	11,13,14,14	0
2	EDO	A	5003	4/4	0.99	0.06	10,12,12,12	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.