



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 27, 2022 – 10:17 am BST

PDB ID : 7PR9  
Title : Crystal structure of Burkholderia pseudomallei heparanase in complex with covalent inhibitor VL166  
Authors : Wu, L.; Armstrong, Z.; Davies, G.J.  
Deposited on : 2021-09-21  
Resolution : 1.34 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](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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.29  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.29

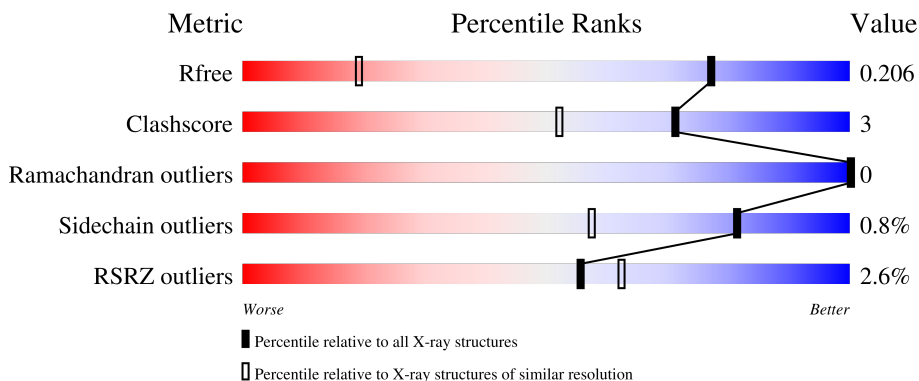
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.34 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	1385 (1.36-1.32)
Clashscore	141614	1417 (1.36-1.32)
Ramachandran outliers	138981	1397 (1.36-1.32)
Sidechain outliers	138945	1397 (1.36-1.32)
RSRZ outliers	127900	1369 (1.36-1.32)

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	438	
1	B	438	
2	C	2	
2	D	2	

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
3	EDO	A	502	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13984 atoms, of which 6419 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.

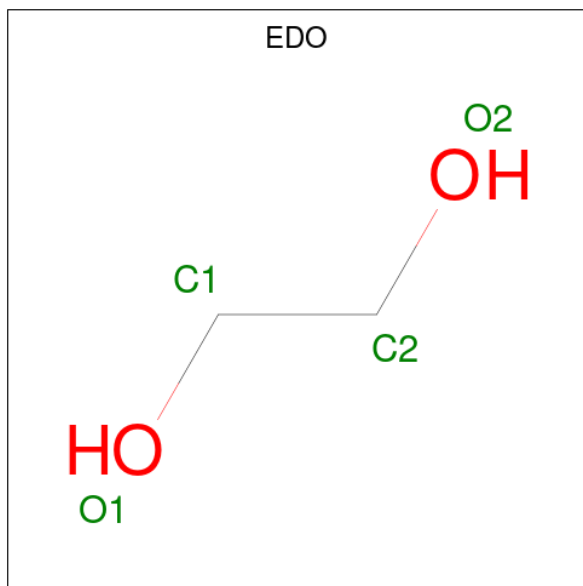
- Molecule 1 is a protein called Glyco\_hydro\_44 domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	437	Total 6498	C 2057	H 3209	N 590	O 633	S 9	93	18	0
1	B	437	Total 6334	C 2015	H 3128	N 567	O 616	S 8	90	7	0

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-(2R,3S,5R,6R)-2,3,4,5,6-pentakis(oxidanyl)cyclohexane-1-carboxylic acid.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	H	N	O			
2	C	2	Total 50	C 15	H 23	N 1	O 11	6	0	0
2	D	2	Total 50	C 15	H 23	N 1	O 11	6	0	0

- Molecule 3 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
3	A	1	Total	C	H	O	1	0
			10	2	6	2		
3	A	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		

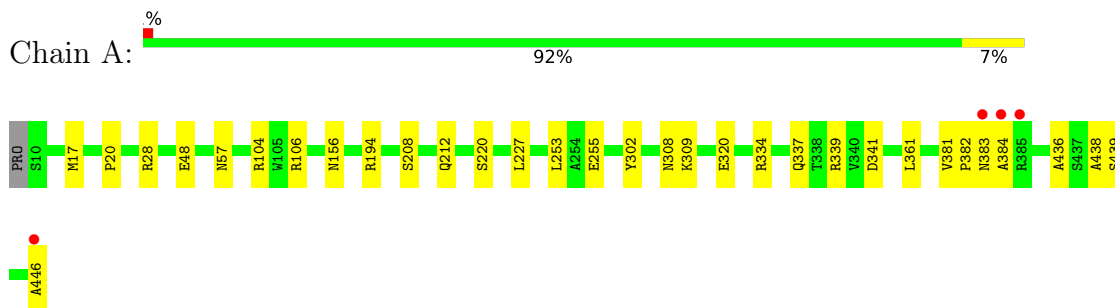
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	597	Total	O	0	0
			597	597		
4	B	395	Total	O	0	0
			395	395		

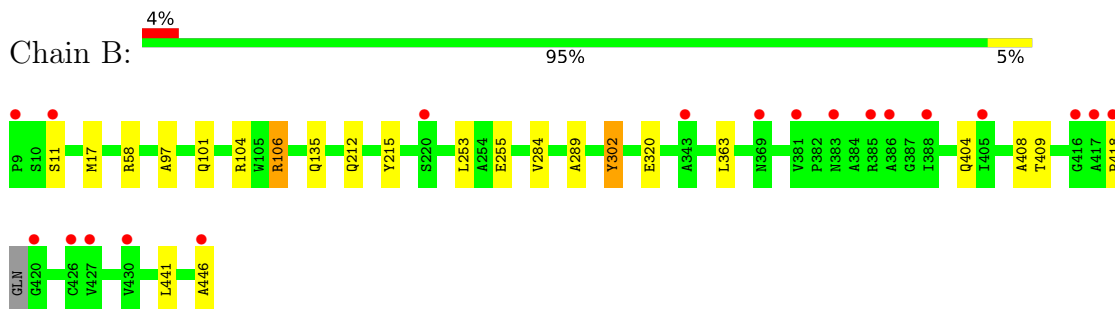
### 3 Residue-property plots [i](#)

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.

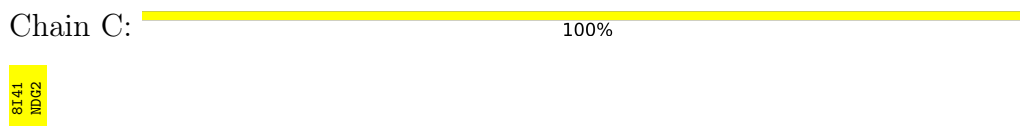
- Molecule 1: Glyco\_hydro\_44 domain-containing protein



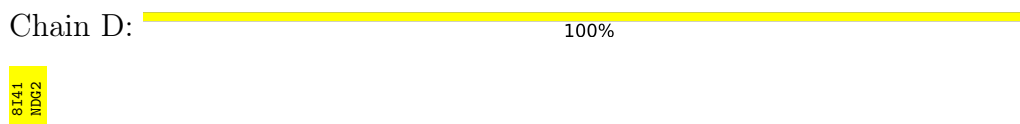
- Molecule 1: Glyco\_hydro\_44 domain-containing protein



- Molecule 2: 2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-(2R,3S,5R,6R)-2,3,4,5,6-pentakis (oxidanyl)cyclohexane-1-carboxylic acid



- Molecule 2: 2-acetamido-2-deoxy-alpha-D-glucopyranose-(1-4)-(2R,3S,5R,6R)-2,3,4,5,6-pentakis (oxidanyl)cyclohexane-1-carboxylic acid



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.87Å 104.14Å 113.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	77.00 – 1.34 76.88 – 1.34	Depositor EDS
% Data completeness (in resolution range)	98.7 (77.00-1.34) 98.2 (76.88-1.34)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.23 (at 1.34Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.176 , 0.199 0.183 , 0.206	Depositor DCC
$R_{free}$ test set	10538 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.1	Xtriage
Anisotropy	0.407	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	13984	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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: EDO, 8I4, NDG

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	0.83	4/3360 (0.1%)	0.93	2/4585 (0.0%)
1	B	0.82	1/3277 (0.0%)	0.89	4/4475 (0.1%)
All	All	0.82	5/6637 (0.1%)	0.91	6/9060 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	255	GLU	CD-OE2	10.25	1.36	1.25
1	A	255	GLU	CD-OE2	7.16	1.33	1.25
1	A	320	GLU	CD-OE1	-6.92	1.18	1.25
1	A	48	GLU	CD-OE2	-5.90	1.19	1.25
1	A	208	SER	CB-OG	-5.03	1.35	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	302	TYR	CB-CG-CD2	-5.67	117.60	121.00
1	A	28	ARG	NE-CZ-NH2	5.66	123.13	120.30
1	A	320	GLU	CG-CD-OE1	-5.63	107.03	118.30
1	B	106	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	B	320	GLU	OE1-CD-OE2	5.16	129.50	123.30
1	B	58	ARG	NE-CZ-NH2	-5.09	117.76	120.30

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	3289	3209	3186	38	0
1	B	3206	3128	3113	16	0
2	C	27	23	12	0	0
2	D	27	23	12	0	0
3	A	8	12	12	5	0
3	B	16	24	23	0	0
4	A	597	0	0	13	1
4	B	395	0	0	0	0
All	All	7565	6419	6358	45	1

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 3.

All (45) 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:384:ALA:O	4:A:601:HOH:O	1.78	1.01
1:A:339[B]:ARG:HH22	1:A:341:ASP:HB2	1.30	0.97
1:A:156:ASN:OD1	3:A:502:EDO:H11	1.75	0.87
1:A:156:ASN:OD1	3:A:502:EDO:C1	2.26	0.82
1:A:57:ASN:HB3	3:A:501:EDO:H12	1.65	0.78
1:A:17[B]:MET:HE3	4:A:860:HOH:O	1.88	0.73
1:A:104:ARG:NH2	1:B:446:ALA:HB1	2.04	0.72
1:A:446:ALA:OXT	1:B:104:ARG:NH1	2.24	0.70
1:A:227:LEU:O	4:A:602:HOH:O	2.13	0.67
1:B:404:GLN:HE21	1:B:409:THR:HG23	1.60	0.67
1:A:383[B]:ASN:HD21	1:B:97:ALA:HB1	1.60	0.66
1:A:337[A]:GLN:NE2	4:A:607:HOH:O	2.33	0.61
1:A:156:ASN:OD1	3:A:502:EDO:H12	2.02	0.58
1:A:104:ARG:HH22	1:B:446:ALA:HB1	1.69	0.55
1:A:438:ALA:HB3	4:A:620:HOH:O	2.06	0.54
1:A:104:ARG:CZ	1:B:446:ALA:HB1	2.38	0.54
1:A:17[B]:MET:HE1	1:A:361:LEU:HD22	1.90	0.53
1:A:220:SER:C	4:A:791:HOH:O	2.48	0.52
1:A:309:LYS:NZ	4:A:611:HOH:O	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106[A]:ARG:HD3	4:A:603:HOH:O	2.10	0.51
1:A:339[B]:ARG:NH2	1:A:341:ASP:HB2	2.13	0.50
1:B:363:LEU:HD12	1:B:441:LEU:HD11	1.94	0.50
1:A:334[A]:ARG:HH22	1:B:106:ARG:CZ	2.23	0.50
1:A:337[A]:GLN:NE2	4:A:608:HOH:O	2.38	0.49
1:A:382[A]:PRO:HG3	1:B:135:GLN:HG2	1.94	0.49
1:A:383[A]:ASN:CB	4:A:686:HOH:O	2.61	0.48
1:A:381[A]:VAL:HG11	1:A:384:ALA:HB2	1.95	0.48
1:B:212:GLN:O	1:B:253:LEU:HA	2.14	0.48
1:A:383[A]:ASN:HB3	4:A:686:HOH:O	2.15	0.47
1:A:20:PRO:HG3	1:A:337[A]:GLN:HG2	1.97	0.45
1:A:308:ASN:HB3	4:A:971:HOH:O	2.16	0.45
1:A:334[A]:ARG:HH22	1:B:106:ARG:NH2	2.15	0.44
1:A:382[A]:PRO:O	1:B:101:GLN:OE1	2.35	0.44
1:A:104:ARG:HD3	1:A:104:ARG:HA	1.70	0.44
1:A:212:GLN:O	1:A:253:LEU:HA	2.17	0.44
1:B:101:GLN:HA	1:B:104:ARG:NH2	2.33	0.44
1:A:383[A]:ASN:CG	4:A:686:HOH:O	2.57	0.42
1:A:17[B]:MET:CE	1:A:361:LEU:HD13	2.50	0.42
1:B:284:VAL:HG12	1:B:289:ALA:HB3	2.02	0.41
1:B:101:GLN:HA	1:B:104:ARG:HH21	1.85	0.41
1:A:339[B]:ARG:NH2	1:A:339[B]:ARG:HG2	2.36	0.41
1:A:194:ARG:NE	1:A:194:ARG:HA	2.35	0.41
1:B:408:ALA:CB	1:B:418:PRO:HD3	2.51	0.41
1:A:436:ALA:O	1:A:439[B]:SER:HB2	2.20	0.40
1:A:156:ASN:CG	3:A:502:EDO:H11	2.41	0.40

All (1) 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)
4:A:782:HOH:O	4:A:978:HOH:O[2_455]	2.13	0.07

## 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	452/438 (103%)	445 (98%)	7 (2%)	0	100	100
1	B	440/438 (100%)	433 (98%)	7 (2%)	0	100	100
All	All	892/876 (102%)	878 (98%)	14 (2%)	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 sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	326/310 (105%)	325 (100%)	1 (0%)	92	81
1	B	316/310 (102%)	311 (98%)	5 (2%)	62	29
All	All	642/620 (104%)	636 (99%)	6 (1%)	81	52

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	302	TYR
1	B	11	SER
1	B	17	MET
1	B	215[A]	TYR
1	B	215[B]	TYR
1	B	302	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	428	ASN
1	B	404	GLN
1	B	428	ASN

### 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](#)

4 monosaccharides 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	8I4	C	1	1,2	13,13,14	1.22	1 (7%)	17,19,21	1.44	2 (11%)
2	NDG	C	2	2	14,14,15	0.79	0	17,19,21	1.73	4 (23%)
2	8I4	D	1	1,2	13,13,14	1.40	3 (23%)	17,19,21	1.42	2 (11%)
2	NDG	D	2	2	14,14,15	0.84	0	17,19,21	1.90	2 (11%)

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	8I4	C	1	1,2	-	0/4/24/28	0/1/1/1
2	NDG	C	2	2	-	0/6/23/26	0/1/1/1
2	8I4	D	1	1,2	-	0/4/24/28	0/1/1/1
2	NDG	D	2	2	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	8I4	O6A-C6	2.42	1.29	1.22
2	C	1	8I4	C1-C7	-2.40	1.48	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	8I4	C5-C7	-2.22	1.50	1.54
2	D	1	8I4	C5-C4	-2.11	1.50	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	NDG	C1-O5-C5	5.99	120.31	112.19
2	C	2	NDG	C1-O5-C5	3.62	117.10	112.19
2	C	1	8I4	O4-C4-C5	-3.32	102.43	109.68
2	C	2	NDG	O7-C7-C8	2.88	127.41	122.06
2	C	2	NDG	C8-C7-N2	-2.74	111.45	116.10
2	D	1	8I4	C2-C3-C4	-2.68	106.26	110.89
2	C	2	NDG	C6-C5-C4	2.47	118.78	113.00
2	D	1	8I4	O6A-C6-C5	-2.13	117.11	122.78
2	D	2	NDG	C3-C4-C5	-2.08	106.53	110.24
2	C	1	8I4	C1-C7-C5	2.01	114.88	111.56

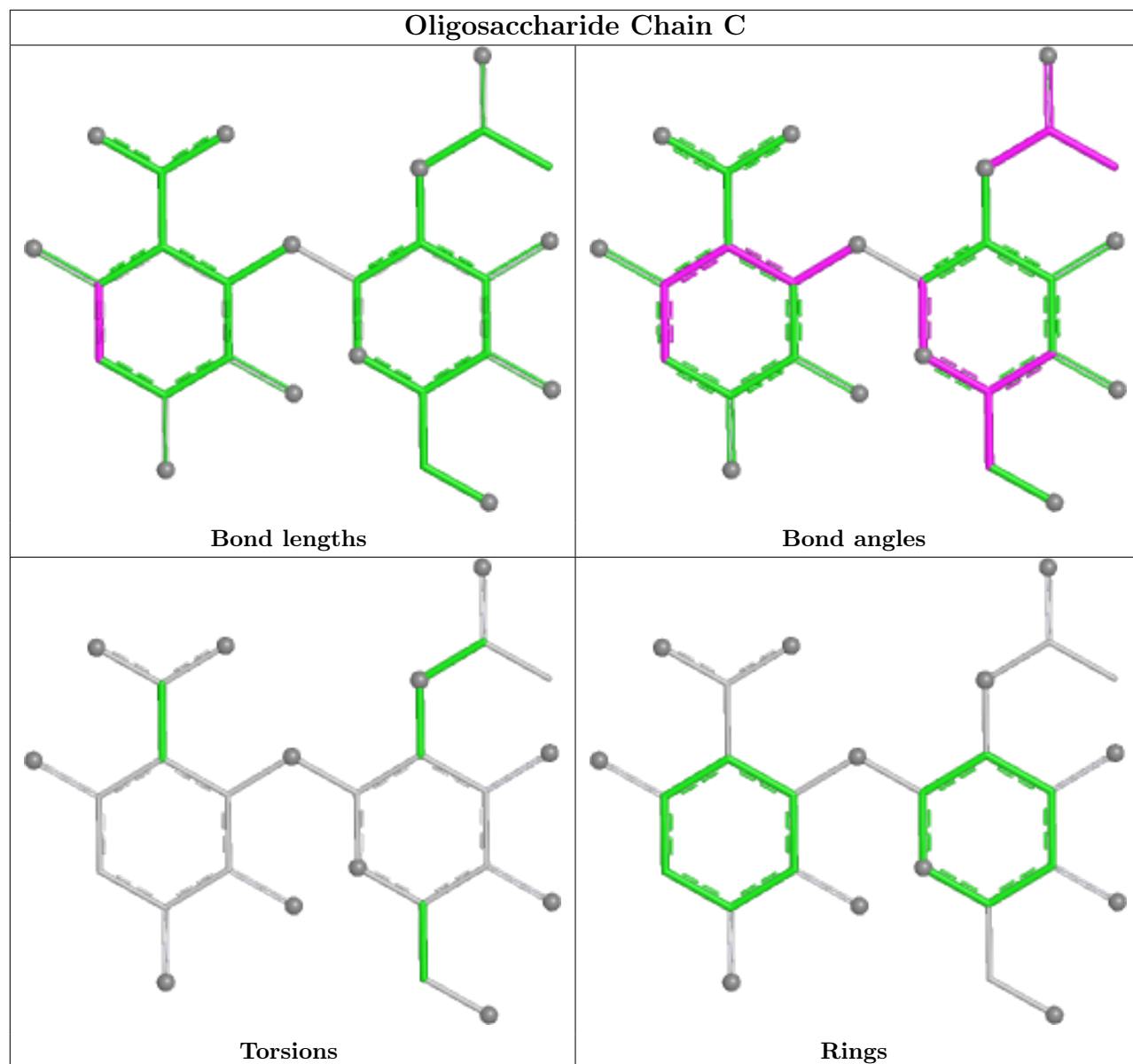
There are no chirality outliers.

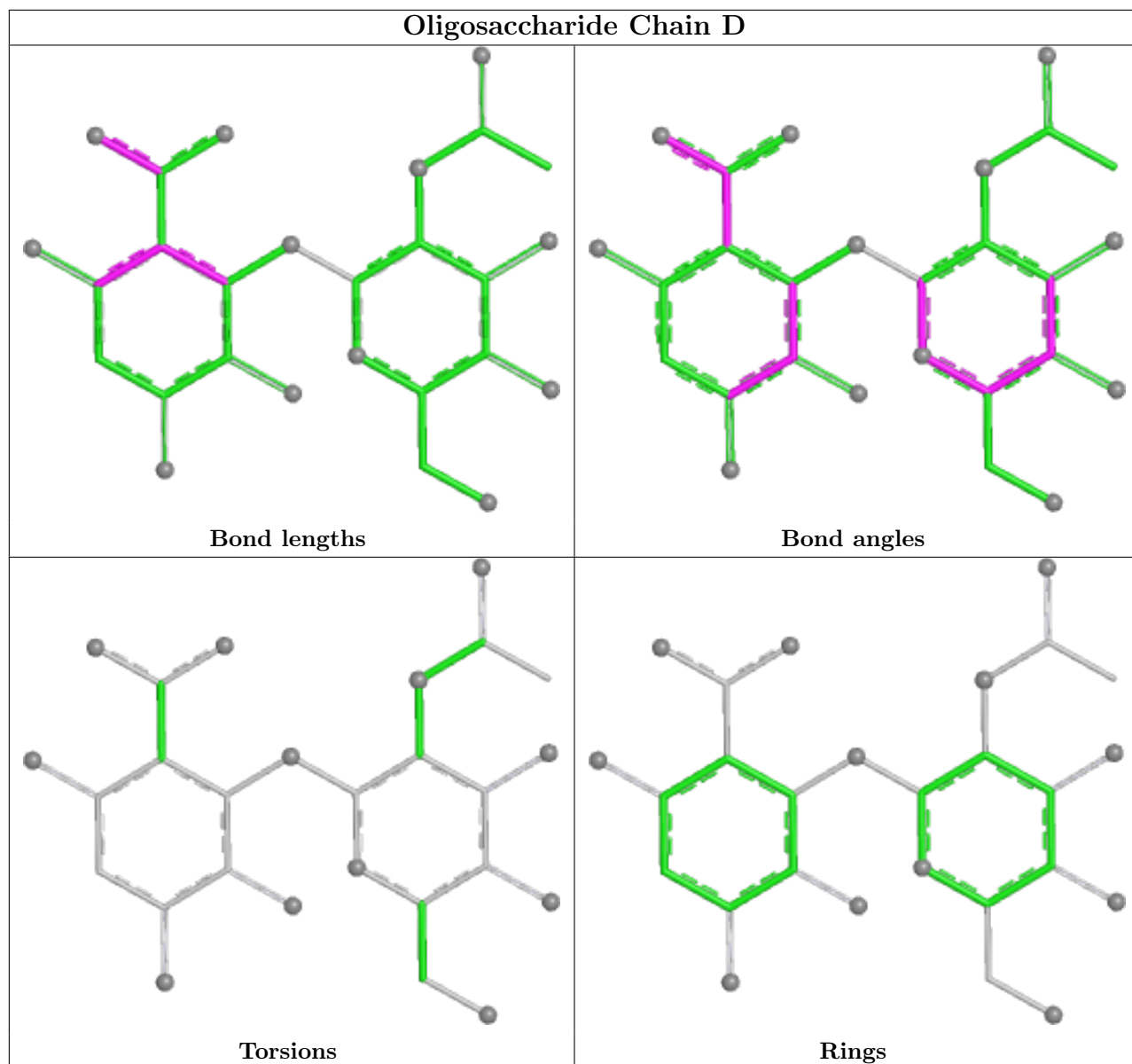
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





## 5.6 Ligand geometry [i](#)

6 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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	EDO	A	501	-	3,3,3	0.12	0	2,2,2	0.69	0
3	EDO	B	503	-	3,3,3	0.57	0	2,2,2	0.15	0
3	EDO	B	502	-	3,3,3	0.24	0	2,2,2	0.19	0
3	EDO	B	504	-	3,3,3	0.47	0	2,2,2	0.35	0
3	EDO	B	501	-	3,3,3	0.90	0	2,2,2	0.05	0
3	EDO	A	502	-	3,3,3	0.06	0	2,2,2	0.58	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
3	EDO	A	501	-	-	1/1/1/1	-
3	EDO	B	503	-	-	0/1/1/1	-
3	EDO	B	502	-	-	1/1/1/1	-
3	EDO	B	504	-	-	0/1/1/1	-
3	EDO	B	501	-	-	0/1/1/1	-
3	EDO	A	502	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	EDO	O1-C1-C2-O2
3	B	502	EDO	O1-C1-C2-O2
3	A	502	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	EDO	1	0
3	A	502	EDO	4	0



## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	437/438 (99%)	-0.21	4 (0%) 84 87	12, 16, 29, 50	0
1	B	437/438 (99%)	0.15	19 (4%) 35 41	13, 25, 48, 86	0
All	All	874/876 (99%)	-0.03	23 (2%) 56 62	12, 19, 40, 86	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	418	PRO	4.7
1	B	420	GLY	4.3
1	B	446	ALA	4.0
1	B	417	ALA	3.5
1	B	385	ARG	3.2
1	B	9	PRO	3.1
1	B	426	CYS	2.9
1	B	343	ALA	2.7
1	B	427	VAL	2.7
1	B	386	ALA	2.6
1	A	383[A]	ASN	2.4
1	A	446	ALA	2.4
1	B	405	ILE	2.4
1	B	220	SER	2.4
1	B	430	VAL	2.4
1	A	385	ARG	2.4
1	B	416	GLY	2.3
1	A	384	ALA	2.3
1	B	388	ILE	2.3
1	B	383	ASN	2.3
1	B	369	ASN	2.2
1	B	381	VAL	2.2
1	B	11	SER	2.1

## 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](#)

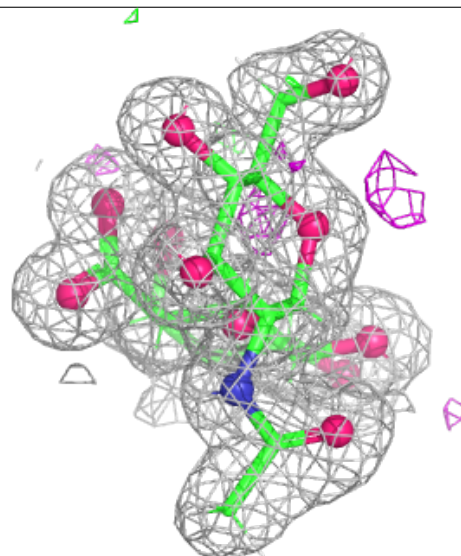
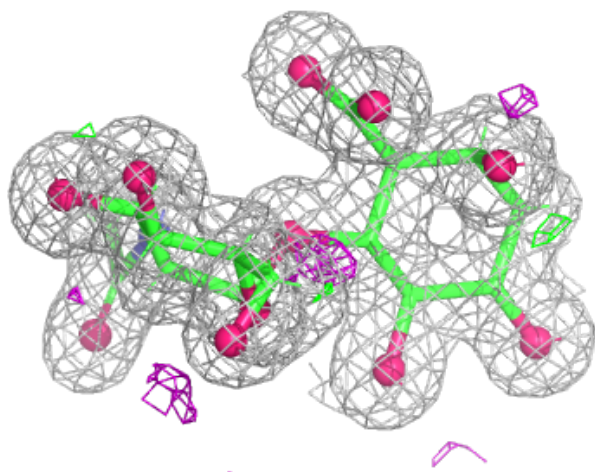
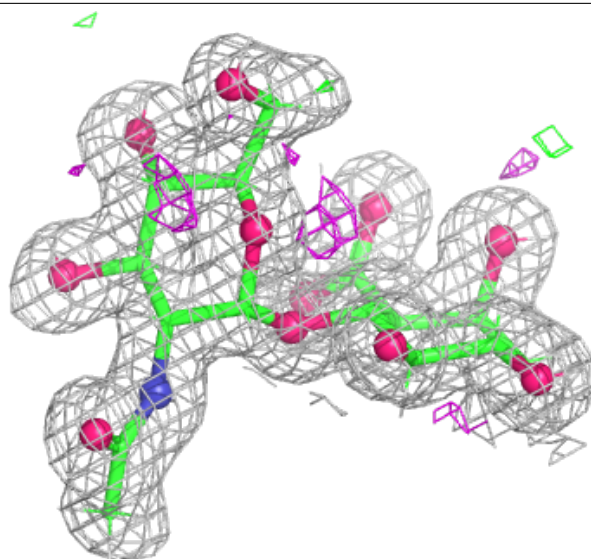
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

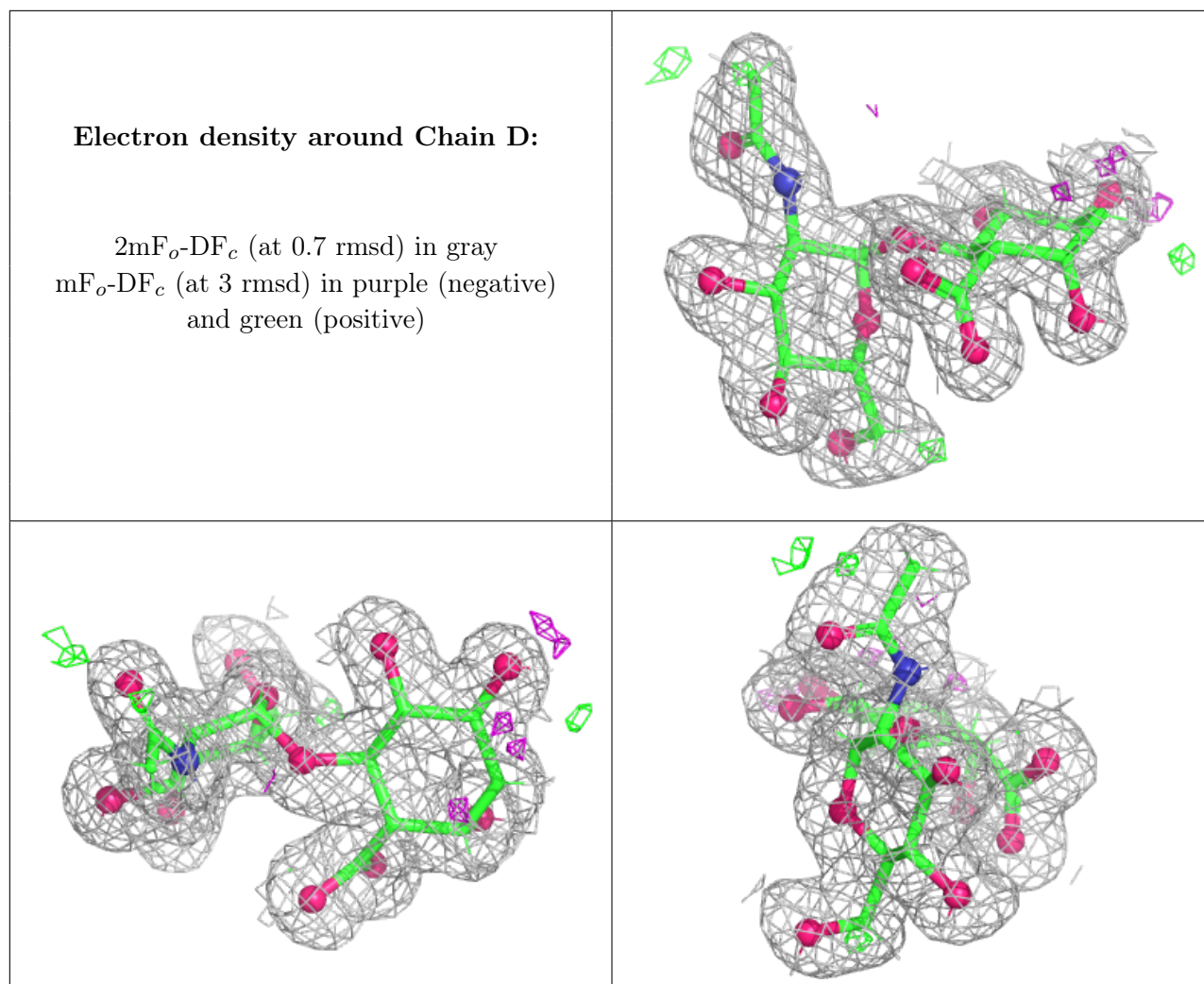
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NDG	D	2	14/15	0.94	0.07	21,25,28,33	3
2	8I4	D	1	13/14	0.95	0.07	18,19,20,22	3
2	NDG	C	2	14/15	0.96	0.06	16,20,25,29	3
2	8I4	C	1	13/14	0.98	0.06	14,15,16,16	3

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around Chain C:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 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	EDO	B	502	4/4	0.79	0.13	37,48,51,52	1
3	EDO	A	502	4/4	0.81	0.14	28,30,33,33	1
3	EDO	A	501	4/4	0.81	0.32	40,40,41,42	1
3	EDO	B	504	4/4	0.86	0.22	31,32,35,37	1
3	EDO	B	503	4/4	0.94	0.16	25,30,35,37	1
3	EDO	B	501	4/4	0.98	0.11	20,22,22,22	1

## 6.5 Other polymers

There are no such residues in this entry.