



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 05:27 PM UTC

PDB ID : 2VR4 / pdb\_00002vr4  
Title : Transition-state mimicry in mannoside hydrolysis: characterisation of twenty six inhibitors and insight into binding from linear free energy relationships and 3-D structure  
Authors : Tailford, L.E.; Offen, W.A.; Smith, N.L.; Dumon, C.; Moreland, C.; Gratien, J.; Heck, M.P.; Stick, R.V.; Bleriot, Y.; Vasella, A.; Gilbert, H.J.; Davies, G.J.  
Deposited on : 2008-03-25  
Resolution : 1.80 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 : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

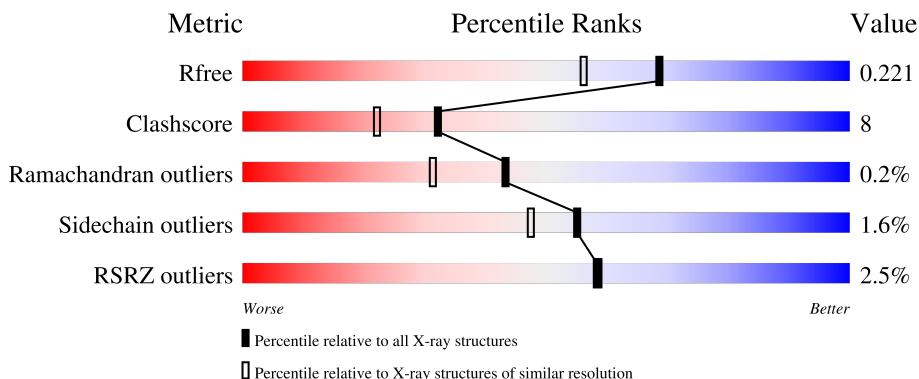
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


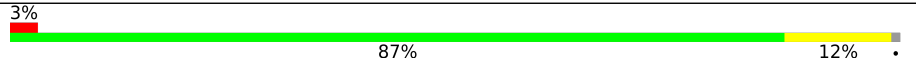
The reported resolution of this entry is 1.80 Å.

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))
$R_{free}$	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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  $\geq 3$ , 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	846	
1	B	846	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	CL	A	1890	-	-	X	-

## 2 Entry composition [i](#)

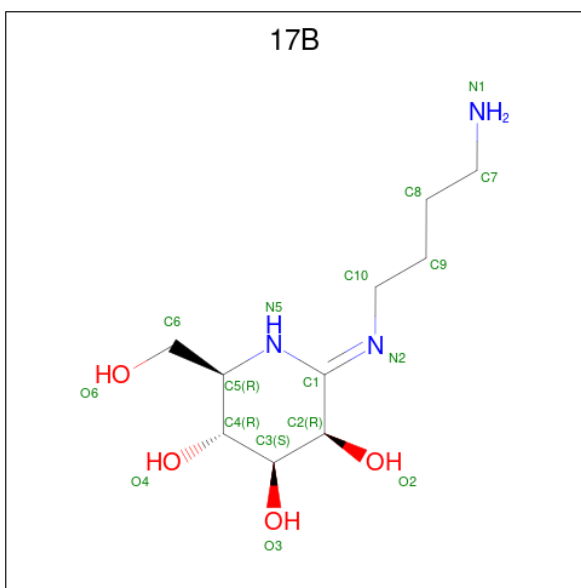
There are 6 unique types of molecules in this entry. The entry contains 15439 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BETA-MANNOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	836	Total 7093	C 4545	N 1193	O 1320	S 35	0	60	0
1	B	841	Total 6943	C 4451	N 1168	O 1289	S 35	0	33	0

- Molecule 2 is (2Z,3R,4S,5R,6R)-2-[(4-aminobutyl)imino]-6-(hydroxymethyl)piperidine-3,4,5-triol (CCD ID: 17B) (formula: C<sub>10</sub>H<sub>21</sub>N<sub>3</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total 17	C 10	N 3	O 4	0	0
2	B	1	Total 17	C 10	N 3	O 4	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: 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 O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0

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

- Molecule 4 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	6	Total Br 6 6	0	0
4	B	2	Total Br 2 2	0	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	5	Total Cl 5 5	0	0
5	B	6	Total Cl 6 6	0	0

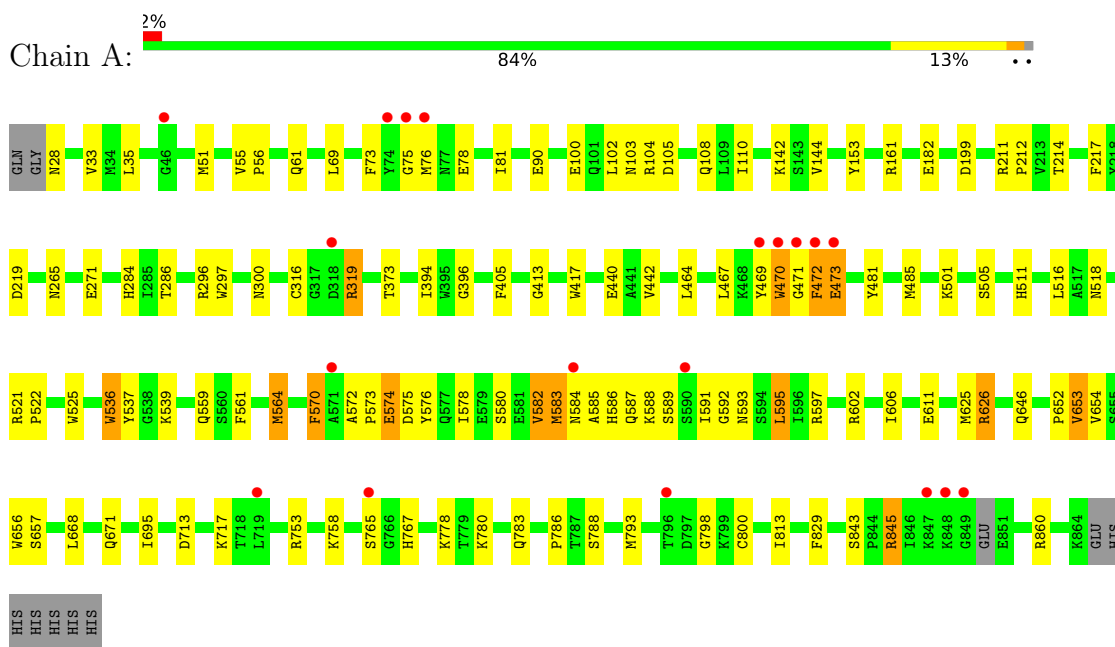
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	570	Total O 571 571	0	1
6	B	579	Total O 579 579	0	0

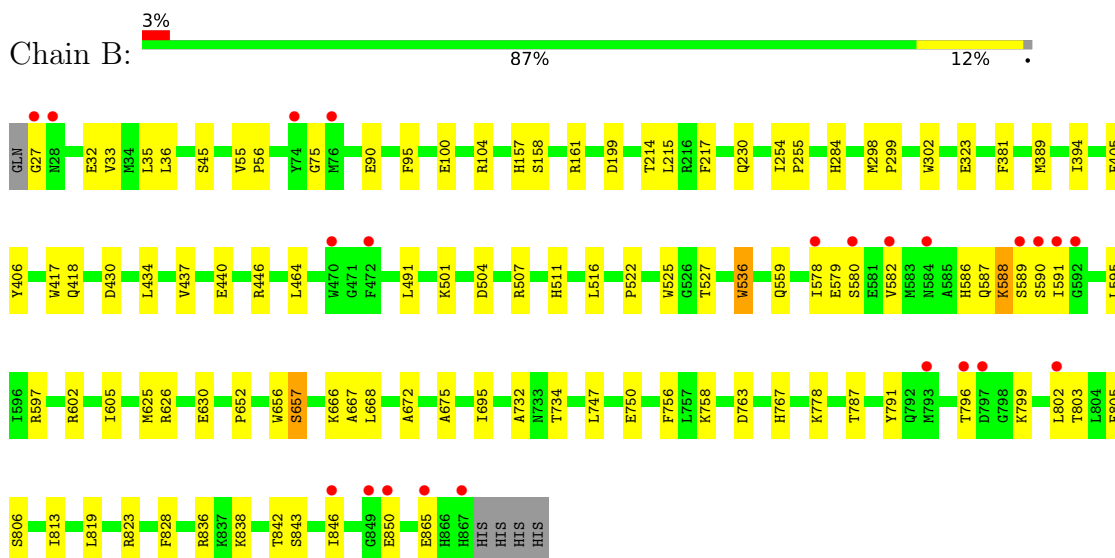
### 3 Residue-property plots [i](#)

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: BETA-MANNOSIDASE



#### • Molecule 1: BETA-MANNOSIDASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.19Å 114.74Å 99.17Å 90.00° 113.03° 90.00°	Depositor
Resolution (Å)	65.37 – 1.80 65.37 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.3 (65.37-1.80) 99.3 (65.37-1.80)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.42 (at 1.80Å)	Xtrriage
Refinement program	REFMAC 5.4.0065	Depositor
R, $R_{free}$	0.177 , 0.222 0.176 , 0.221	Depositor DCC
$R_{free}$ test set	8536 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	21.2	Xtrriage
Anisotropy	0.064	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	15439	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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: BR, 17B, CL, 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).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.96	1/7345 (0.0%)	0.95	5/9981 (0.1%)
1	B	1.01	1/7175 (0.0%)	0.97	8/9750 (0.1%)
All	All	0.99	2/14520 (0.0%)	0.96	13/19731 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	605	ILE	CA-CB	5.30	1.60	1.54
1	A	653	VAL	CA-CB	5.25	1.61	1.55

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	588[A]	LYS	N-CA-C	7.36	123.98	113.02
1	B	588[B]	LYS	N-CA-C	7.36	123.98	113.02
1	A	396	GLY	N-CA-C	6.42	123.27	115.31
1	A	470	TRP	N-CA-C	6.30	121.03	113.22
1	A	564	MET	N-CA-C	6.08	117.91	111.28
1	B	672	ALA	N-CA-C	-5.96	104.86	111.36
1	B	381	PHE	N-CA-C	5.77	117.56	111.28
1	B	430	ASP	CA-C-N	5.50	124.96	119.24
1	B	430	ASP	C-N-CA	5.50	124.96	119.24
1	A	144	VAL	N-CA-CB	-5.41	104.89	112.35
1	B	657	SER	N-CA-C	5.30	117.56	110.35
1	B	389	MET	N-CA-C	5.24	117.89	110.50
1	A	296	ARG	N-CA-C	5.15	117.90	110.28

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts

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	7093	0	6776	139	0
1	B	6943	0	6641	82	0
2	A	17	0	21	0	0
2	B	17	0	21	0	0
3	A	84	0	126	13	0
3	B	116	0	174	17	0
4	A	6	0	0	2	0
4	B	2	0	0	0	0
5	A	5	0	0	4	0
5	B	6	0	0	0	0
6	A	571	0	0	10	0
6	B	579	0	0	6	0
All	All	15439	0	13759	224	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 8.

All (224) 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:574[A]:GLU:N	1:A:574[A]:GLU:OE2	1.58	1.36
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:CG2	1.73	1.36
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:CG1	1.83	1.23
1:A:845:ARG:HG3	1:A:845:ARG:HH11	1.12	1.12
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:CB	1.99	1.11
1:A:467:LEU:HB3	1:A:485:MET:HE1	1.29	1.09
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:HG13	1.47	1.07
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:HG22	1.33	1.07
1:A:578[A]:ILE:CG2	1:A:578[A]:ILE:HA	1.87	1.03
1:A:470:TRP:N	1:A:471[A]:GLY:HA2	1.73	1.03
1:A:573[B]:PRO:HD2	1:A:574[B]:GLU:OE1	1.62	1.00
1:A:467:LEU:HD22	1:A:485:MET:HE2	1.46	0.97
1:A:578[A]:ILE:CG2	1:A:578[A]:ILE:CA	2.43	0.97
1:A:575[A]:ASP:CG	1:A:582[A]:VAL:HG13	1.90	0.96
1:A:481:TYR:CZ	1:A:485:MET:HE3	2.06	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:587:GLN:OE1	1:B:589[A]:SER:HB3	1.72	0.89
1:B:446:ARG:HD2	3:B:1896:EDO:H22	1.55	0.87
1:B:45:SER:HB2	1:B:90:GLU:OE2	1.74	0.87
1:A:214[A]:THR:HG22	6:A:2046:HOH:O	1.73	0.86
1:B:33:VAL:HG13	1:B:214[A]:THR:CG2	2.06	0.86
1:A:572[B]:ALA:HB1	1:A:574[B]:GLU:OE1	1.77	0.85
3:A:1886:EDO:H21	1:B:597:ARG:HH22	1.46	0.80
1:A:574[B]:GLU:H	1:A:574[B]:GLU:CD	1.88	0.80
1:B:230:GLN:HE21	3:B:1889:EDO:H21	1.45	0.80
1:A:758:LYS:HG3	6:A:2501:HOH:O	1.81	0.80
1:A:583[A]:MET:O	1:A:593[A]:ASN:ND2	2.17	0.77
1:A:69:LEU:HD22	1:A:81:ILE:HG22	1.67	0.77
1:A:35:LEU:HD23	1:A:214[B]:THR:HG22	1.65	0.77
1:A:373:THR:HG22	1:B:796:THR:OG1	1.85	0.77
1:A:573[A]:PRO:C	1:A:574[A]:GLU:OE2	2.29	0.76
1:B:695:ILE:HG23	3:B:1877:EDO:H22	1.68	0.76
1:B:158:SER:HB3	1:B:161:ARG:HG2	1.68	0.75
1:A:182[B]:GLU:H	1:A:182[B]:GLU:CD	1.98	0.71
1:B:464:LEU:HD22	1:B:511:HIS:HE1	1.55	0.71
1:A:561:PHE:HE1	1:A:583[A]:MET:HG3	1.56	0.71
1:A:574[A]:GLU:N	1:A:574[A]:GLU:CD	2.49	0.70
1:A:470:TRP:N	1:A:471[B]:GLY:HA2	2.07	0.70
1:B:33:VAL:CG1	1:B:214[A]:THR:CG2	2.70	0.70
1:A:472[A]:PHE:O	1:A:473[A]:GLU:HB3	1.89	0.70
3:A:1885:EDO:H22	6:A:2312:HOH:O	1.91	0.69
1:B:756:PHE:HE1	1:B:758:LYS:HG3	1.58	0.68
1:A:626:ARG:HG3	6:A:2458:HOH:O	1.93	0.68
1:A:845:ARG:HG3	1:A:845:ARG:NH1	1.92	0.66
1:A:104:ARG:HG3	1:A:217:PHE:HB3	1.77	0.66
1:A:537:TYR:CE1	5:A:1890:CL:CL	2.86	0.66
1:A:580[A]:SER:OG	1:A:583[A]:MET:HB2	1.96	0.65
1:A:564:MET:HG3	1:A:611[A]:GLU:HG2	1.78	0.65
1:A:469:TYR:C	1:A:471[A]:GLY:HA2	2.22	0.65
1:A:110:ILE:HB	1:A:214[A]:THR:HG23	1.78	0.64
1:A:786:PRO:HA	1:A:860[B]:ARG:NE	2.12	0.64
1:A:467:LEU:HD22	1:A:485:MET:CE	2.23	0.64
1:A:695:ILE:HG23	3:A:1881:EDO:H11	1.79	0.64
1:A:211:ARG:HG2	3:A:1889:EDO:H21	1.80	0.63
1:B:803:THR:HG21	1:B:836[B]:ARG:HD3	1.80	0.63
1:A:573[B]:PRO:CD	1:A:574[B]:GLU:OE1	2.43	0.62
1:A:561:PHE:CE1	1:A:583[A]:MET:HG3	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406[B]:TYR:OH	1:B:418:GLN:OE1	2.17	0.61
1:B:33:VAL:HG13	1:B:214[A]:THR:HG21	1.80	0.61
1:A:90:GLU:OE1	1:A:153[A]:TYR:OH	2.13	0.60
1:A:574[A]:GLU:CD	1:A:574[A]:GLU:H	2.03	0.60
5:A:1887:CL:CL	6:A:2464:HOH:O	2.54	0.60
3:B:1899:EDO:H11	6:B:2523:HOH:O	2.01	0.60
1:A:373:THR:CG2	1:B:796:THR:OG1	2.49	0.60
1:A:76:MET:HE3	1:A:585[B]:ALA:HB2	1.84	0.59
1:A:470:TRP:N	1:A:471[A]:GLY:CA	2.60	0.59
1:A:442:VAL:HG12	3:A:1885:EDO:H11	1.83	0.59
1:B:36:LEU:HA	3:B:1906:EDO:H21	1.85	0.59
1:B:104:ARG:HG3	1:B:217:PHE:HB3	1.85	0.58
1:A:481:TYR:CE1	1:A:485:MET:HE3	2.39	0.58
1:A:199:ASP:C	1:A:588[A]:LYS:HB2	2.29	0.58
1:A:525:TRP:CE2	3:A:1867:EDO:H22	2.39	0.57
1:B:298[A]:MET:HE3	1:B:299:PRO:HD2	1.86	0.57
1:A:481:TYR:OH	1:A:485:MET:HE3	2.04	0.56
1:A:110:ILE:HB	1:A:214[A]:THR:CG2	2.36	0.56
1:A:559:GLN:OE1	1:A:587[A]:GLN:NE2	2.24	0.56
3:B:1897:EDO:H11	6:B:2571:HOH:O	2.06	0.56
1:A:578[A]:ILE:CG2	1:A:578[A]:ILE:N	2.69	0.56
1:B:819:LEU:HD11	1:B:846:ILE:HD13	1.89	0.55
1:B:525:TRP:CE2	3:B:1870:EDO:H22	2.41	0.55
1:A:316:CYS:O	1:A:319[A]:ARG:HD3	2.07	0.55
1:A:539:LYS:HG2	1:A:595:LEU:HD11	1.88	0.54
1:A:467:LEU:CB	1:A:485:MET:HE1	2.21	0.54
1:A:765:SER:O	3:A:1896:EDO:H21	2.07	0.54
1:A:580[A]:SER:O	1:A:584[A]:ASN:HB2	2.07	0.54
1:A:845:ARG:HH11	1:A:845:ARG:CG	1.99	0.54
1:A:574[A]:GLU:OE2	1:A:574[A]:GLU:CA	2.50	0.53
1:B:33:VAL:CG1	1:B:214[A]:THR:HG21	2.34	0.53
1:A:472[A]:PHE:O	1:A:473[A]:GLU:CB	2.57	0.53
1:A:570[A]:PHE:CD1	1:A:570[A]:PHE:C	2.87	0.52
1:A:589[A]:SER:HB2	1:A:656:TRP:CH2	2.44	0.52
1:A:793:MET:SD	1:A:800[A]:CYS:SG	3.07	0.52
1:A:199:ASP:O	1:A:588[A]:LYS:HB2	2.10	0.52
1:A:472[A]:PHE:N	1:A:472[A]:PHE:CD1	2.73	0.52
1:A:591[A]:ILE:HB	1:A:595:LEU:HD22	1.91	0.52
1:A:575[A]:ASP:CG	1:A:582[A]:VAL:HG22	2.28	0.52
1:B:747:LEU:O	3:B:1886:EDO:H22	2.10	0.52
1:A:300:ASN:O	4:A:1882:BR:BR	2.83	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:787:THR:OG1	1:B:806:SER:OG	2.27	0.51
1:B:284:HIS:HD2	3:B:1903:EDO:H12	1.75	0.51
1:B:626[C]:ARG:HG3	1:B:675:ALA:HB2	1.93	0.51
1:A:271:GLU:HG2	6:A:2185:HOH:O	2.11	0.51
1:A:394:ILE:HD11	1:A:405:PHE:CE2	2.46	0.51
1:A:574[B]:GLU:CD	1:A:574[B]:GLU:N	2.64	0.50
1:B:394:ILE:HD11	1:B:405:PHE:CE2	2.46	0.50
1:B:586:HIS:HD2	1:B:652:PRO:O	1.94	0.50
1:B:33:VAL:CG1	1:B:214[A]:THR:HG23	2.41	0.50
1:B:799:LYS:HG3	1:B:842:THR:HG22	1.93	0.50
1:A:284:HIS:HB3	3:A:1893:EDO:H21	1.94	0.49
1:B:446:ARG:HH11	3:B:1896:EDO:C2	2.25	0.49
1:B:803:THR:CG2	1:B:836[B]:ARG:HD3	2.42	0.49
1:A:467:LEU:HB3	1:A:485:MET:CE	2.21	0.49
1:A:586[B]:HIS:HD2	1:A:652:PRO:O	1.95	0.49
1:B:591[A]:ILE:HD12	1:B:595:LEU:HD22	1.94	0.49
1:A:102:LEU:O	1:A:142:LYS:HE3	2.12	0.49
1:B:75:GLY:HA2	1:B:582:VAL:HG13	1.95	0.49
1:B:626[A]:ARG:CZ	1:B:630:GLU:OE2	2.61	0.49
1:B:763:ASP:OD2	1:B:767:HIS:HB3	2.12	0.48
1:A:753:ARG:NH1	1:A:783[B]:GLN:OE1	2.46	0.48
1:A:575[A]:ASP:OD2	1:A:582[A]:VAL:HG21	1.96	0.48
1:B:796:THR:CG2	1:B:799:LYS:HB3	2.43	0.48
1:A:582[A]:VAL:HA	1:A:585[A]:ALA:HB3	1.96	0.48
1:A:76:MET:HE1	6:A:2398:HOH:O	2.13	0.48
3:A:1886:EDO:H21	1:B:597:ARG:NH2	2.23	0.48
1:A:537:TYR:HE1	5:A:1890:CL:CL	2.30	0.48
1:A:559:GLN:CD	1:A:587[A]:GLN:HE22	2.17	0.47
1:A:76:MET:CE	6:A:2398:HOH:O	2.62	0.47
1:A:33:VAL:HG13	1:A:214[A]:THR:OG1	2.14	0.47
1:A:73:PHE:O	1:A:586[B]:HIS:NE2	2.39	0.47
1:B:55:VAL:HA	1:B:56:PRO:C	2.39	0.47
1:B:464:LEU:HD22	1:B:511:HIS:CE1	2.43	0.47
1:B:522:PRO:HA	1:B:525:TRP:CE2	2.50	0.47
1:A:561:PHE:HE1	1:A:583[A]:MET:CG	2.25	0.47
1:B:778:LYS:HE3	6:B:2422:HOH:O	2.15	0.46
1:B:850:GLU:HB3	3:B:1907:EDO:H21	1.97	0.46
1:A:78:GLU:OE1	1:A:588[B]:LYS:NZ	2.33	0.46
1:B:579:GLU:C	1:B:580:SER:O	2.57	0.46
1:B:589[B]:SER:HB2	1:B:656:TRP:CH2	2.50	0.46
1:B:667:ALA:HB2	1:B:828:PHE:CD1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:TRP:CZ2	1:A:592[B]:GLY:HA2	2.50	0.46
1:A:55:VAL:HA	1:A:56:PRO:C	2.40	0.46
1:B:501:LYS:HG3	6:B:2148:HOH:O	2.15	0.46
1:B:587:GLN:C	1:B:589[A]:SER:H	2.13	0.46
1:B:732:ALA:O	1:B:734:THR:HG23	2.16	0.46
1:B:823:ARG:HB3	3:B:1881:EDO:H11	1.98	0.46
1:A:572[B]:ALA:HB3	1:A:574[B]:GLU:HG2	1.98	0.46
1:B:298[A]:MET:HG3	1:B:302:TRP:O	2.15	0.46
1:A:286:THR:OG1	3:A:1893:EDO:H11	2.16	0.46
1:A:582[A]:VAL:C	1:A:584[A]:ASN:H	2.24	0.46
1:A:845:ARG:NH1	1:A:845:ARG:CG	2.67	0.46
1:B:657:SER:O	1:B:666:LYS:HD2	2.16	0.46
1:B:803:THR:CG2	1:B:836[B]:ARG:HG2	2.47	0.45
1:B:157:HIS:CE1	3:B:1900:EDO:C2	2.99	0.45
1:B:437:VAL:HG11	1:B:491:LEU:HD11	1.98	0.45
1:B:35:LEU:HD23	1:B:214[B]:THR:HG22	1.99	0.45
1:A:75:GLY:HA2	1:A:582[B]:VAL:HG13	1.98	0.45
1:B:802:LEU:O	1:B:838:LYS:HA	2.17	0.45
3:B:1878:EDO:H11	6:B:2563:HOH:O	2.15	0.45
1:B:695:ILE:CG2	3:B:1877:EDO:H22	2.44	0.45
1:B:791:TYR:HA	1:B:803:THR:O	2.17	0.45
1:B:254:ILE:HA	1:B:255:PRO:HD2	1.89	0.45
1:B:516[A]:LEU:HD23	1:B:527:THR:HG21	1.97	0.45
1:A:199:ASP:HA	1:A:588[A]:LYS:HB2	1.99	0.44
1:A:501:LYS:O	1:A:505:SER:HB3	2.17	0.44
1:A:576[A]:TYR:N	1:A:576[A]:TYR:CD1	2.85	0.44
1:A:695:ILE:CG2	3:A:1881:EDO:H11	2.47	0.44
1:A:573[B]:PRO:HA	1:A:576[B]:TYR:CE1	2.53	0.44
1:B:756:PHE:CE1	1:B:758:LYS:HG3	2.44	0.44
1:B:805:PHE:HD1	1:B:836[A]:ARG:HG3	1.83	0.44
1:A:212:PRO:HD3	3:A:1889:EDO:H11	2.00	0.44
1:A:265[B]:ASN:ND2	6:A:2179:HOH:O	2.48	0.44
1:A:578[A]:ILE:HG22	1:A:597[A]:ARG:HH21	1.83	0.43
1:A:373:THR:HG22	1:B:796:THR:HG1	1.80	0.43
1:A:813:ILE:HD11	1:A:829:PHE:CZ	2.53	0.43
3:A:1872:EDO:H12	3:A:1894:EDO:H22	2.00	0.43
1:A:575[A]:ASP:OD1	1:A:582[A]:VAL:HG13	2.16	0.43
1:B:625:MET:HB2	1:B:668:LEU:HD13	1.99	0.43
1:B:819:LEU:CD1	1:B:846:ILE:HD13	2.48	0.43
1:B:525:TRP:CD2	3:B:1870:EDO:H22	2.53	0.43
1:A:589[A]:SER:HB2	1:A:656:TRP:HH2	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:27:GLY:HA3	1:B:323:GLU:OE1	2.19	0.43
1:B:95:PHE:CE1	1:B:215:LEU:HD11	2.54	0.43
1:A:653:VAL:HG22	1:A:654:VAL:N	2.34	0.42
1:B:199:ASP:O	1:B:588[B]:LYS:HB2	2.19	0.42
1:A:587[A]:GLN:OE1	1:A:589[A]:SER:HB3	2.19	0.42
1:A:646:GLN:HG2	1:A:657:SER:HA	2.00	0.42
1:A:518:ASN:OD1	1:A:521[B]:ARG:NH1	2.48	0.42
1:B:75:GLY:O	1:B:586:HIS:HE1	2.02	0.42
1:A:767[A]:HIS:HE2	1:B:32:GLU:CD	2.26	0.42
1:A:108:GLN:HG3	4:A:1891:BR:BR	2.73	0.42
1:B:504:ASP:OD2	1:B:507:ARG:HD2	2.19	0.42
1:A:522:PRO:HA	1:A:525:TRP:CE2	2.55	0.42
1:B:589[B]:SER:HB2	1:B:656:TRP:HH2	1.84	0.42
1:A:798:GLY:HA2	1:A:843:SER:O	2.20	0.41
1:B:516[A]:LEU:HD12	1:B:516[A]:LEU:HA	1.85	0.41
1:B:843:SER:HB3	1:B:846:ILE:HG12	2.01	0.41
1:A:105:ASP:HB2	1:A:219:ASP:HA	2.02	0.41
1:B:602:ARG:O	1:B:778:LYS:HD3	2.20	0.41
1:B:796:THR:HG22	1:B:799:LYS:HB3	2.02	0.41
1:A:297:TRP:CZ2	1:A:413:GLY:HA2	2.56	0.41
1:A:464:LEU:HD22	1:A:511:HIS:HE1	1.85	0.41
1:A:573[B]:PRO:HA	1:A:576[B]:TYR:CD1	2.56	0.41
1:B:446:ARG:HH11	3:B:1896:EDO:H22	1.84	0.41
1:A:199:ASP:HA	1:A:588[A]:LYS:CB	2.51	0.41
1:A:564:MET:CG	1:A:611[A]:GLU:HG2	2.49	0.41
1:A:626:ARG:HD2	1:A:626:ARG:C	2.45	0.41
1:A:626:ARG:HB2	1:A:671:GLN:HB3	2.03	0.41
1:A:778:LYS:HA	5:A:1897:CL:CL	2.58	0.41
1:B:536:TRP:CH2	1:B:559:GLN:HB2	2.56	0.41
1:A:625:MET:HB2	1:A:668:LEU:HD13	2.03	0.41
1:B:434:LEU:HD23	1:B:434:LEU:HA	1.90	0.41
1:B:750:GLU:H	1:B:750:GLU:CD	2.28	0.41
1:A:575[A]:ASP:CG	1:A:582[A]:VAL:CG1	2.62	0.40
1:A:602:ARG:O	1:A:780:LYS:HE2	2.21	0.40
1:A:713:ASP:OD2	1:A:717:LYS:HB3	2.20	0.40
1:A:536:TRP:CH2	1:A:559:GLN:HB2	2.57	0.40
1:A:61:GLN:HB3	6:A:2019:HOH:O	2.21	0.40
1:A:51:MET:HE2	1:A:51:MET:HB3	1.95	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	Percentiles	
1	A	881/846 (104%)	859 (98%)	20 (2%)	2 (0%)	43	31
1	B	864/846 (102%)	835 (97%)	28 (3%)	1 (0%)	48	34
All	All	1745/1692 (103%)	1694 (97%)	48 (3%)	3 (0%)	43	31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	473[A]	GLU
1	A	536	TRP
1	B	536	TRP

### 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	768/754 (102%)	745 (97%)	23 (3%)	36	24
1	B	752/754 (100%)	744 (99%)	8 (1%)	65	60
All	All	1520/1508 (101%)	1489 (98%)	31 (2%)	55	38

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	103	ASN
1	A	161[B]	ARG

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Mol	Chain	Res	Type
1	A	161[D]	ARG
1	A	319[A]	ARG
1	A	319[B]	ARG
1	A	417	TRP
1	A	440	GLU
1	A	472[A]	PHE
1	A	472[B]	PHE
1	A	516	LEU
1	A	570[A]	PHE
1	A	570[B]	PHE
1	A	574[A]	GLU
1	A	574[B]	GLU
1	A	582[A]	VAL
1	A	582[B]	VAL
1	A	583[A]	MET
1	A	583[B]	MET
1	A	595	LEU
1	A	626	ARG
1	A	788	SER
1	A	845	ARG
1	B	100	GLU
1	B	417	TRP
1	B	440	GLU
1	B	578	ILE
1	B	590[A]	SER
1	B	590[B]	SER
1	B	813	ILE
1	B	865	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	162	GLN
1	A	635	ASN
1	B	230	GLN
1	B	274	GLN

### 5.3.3 RNA

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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 71 ligands modelled in this entry, 19 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	EDO	B	1892	-	3,3,3	0.41	0	2,2,2	0.59	0
3	EDO	B	1898	-	3,3,3	0.53	0	2,2,2	0.32	0
3	EDO	B	1873	-	3,3,3	0.51	0	2,2,2	0.30	0
3	EDO	A	1875	-	3,3,3	0.42	0	2,2,2	0.44	0
2	17B	B	1869	-	17,17,17	1.15	3 (17%)	18,22,22	1.19	2 (11%)
3	EDO	A	1873	-	3,3,3	0.35	0	2,2,2	0.88	0
3	EDO	B	1880	-	3,3,3	0.48	0	2,2,2	0.61	0
3	EDO	B	1872	-	3,3,3	0.62	0	2,2,2	0.16	0
3	EDO	A	1872	-	3,3,3	0.37	0	2,2,2	0.43	0
3	EDO	A	1895	-	3,3,3	0.51	0	2,2,2	0.41	0
3	EDO	B	1877	-	3,3,3	0.92	0	2,2,2	0.19	0
3	EDO	A	1874	-	3,3,3	0.84	0	2,2,2	0.39	0
3	EDO	A	1896	-	3,3,3	0.35	0	2,2,2	0.46	0
3	EDO	B	1881	-	3,3,3	0.45	0	2,2,2	0.59	0
2	17B	A	1865	-	17,17,17	1.42	4 (23%)	18,22,22	1.39	2 (11%)
3	EDO	A	1881	-	3,3,3	0.62	0	2,2,2	0.23	0
3	EDO	B	1882	-	3,3,3	0.88	0	2,2,2	0.33	0
3	EDO	B	1878	-	3,3,3	0.65	0	2,2,2	0.21	0
3	EDO	B	1871	-	3,3,3	0.64	0	2,2,2	0.33	0
3	EDO	A	1868	-	3,3,3	0.60	0	2,2,2	0.49	0
3	EDO	B	1874	-	3,3,3	0.68	0	2,2,2	0.40	0
3	EDO	B	1886	-	3,3,3	0.39	0	2,2,2	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	EDO	B	1887	-	3,3,3	0.43	0	2,2,2	0.51	0
3	EDO	B	1889	-	3,3,3	0.39	0	2,2,2	0.69	0
3	EDO	B	1897	-	3,3,3	0.28	0	2,2,2	0.65	0
3	EDO	A	1871	-	3,3,3	0.56	0	2,2,2	0.51	0
3	EDO	A	1885	-	3,3,3	0.53	0	2,2,2	0.14	0
3	EDO	A	1893	-	3,3,3	0.53	0	2,2,2	0.18	0
3	EDO	B	1875	-	3,3,3	0.55	0	2,2,2	0.23	0
3	EDO	B	1905	-	3,3,3	0.38	0	2,2,2	0.54	0
3	EDO	A	1889	-	3,3,3	0.51	0	2,2,2	0.47	0
3	EDO	A	1869	-	3,3,3	0.81	0	2,2,2	0.22	0
3	EDO	B	1896	-	3,3,3	0.33	0	2,2,2	0.33	0
3	EDO	B	1908	-	3,3,3	0.46	0	2,2,2	0.24	0
3	EDO	B	1910	-	3,3,3	0.42	0	2,2,2	0.57	0
3	EDO	A	1883	-	3,3,3	0.56	0	2,2,2	0.45	0
3	EDO	B	1870	-	3,3,3	0.55	0	2,2,2	1.01	0
3	EDO	A	1867	-	3,3,3	0.49	0	2,2,2	0.96	0
3	EDO	A	1898	-	3,3,3	0.51	0	2,2,2	0.31	0
3	EDO	B	1903	-	3,3,3	0.43	0	2,2,2	0.37	0
3	EDO	B	1907	-	3,3,3	0.53	0	2,2,2	0.43	0
3	EDO	A	1894	-	3,3,3	0.43	0	2,2,2	0.42	0
3	EDO	B	1900	-	3,3,3	0.30	0	2,2,2	0.24	0
3	EDO	B	1909	-	3,3,3	0.26	0	2,2,2	0.82	0
3	EDO	A	1886	-	3,3,3	0.35	0	2,2,2	0.59	0
3	EDO	B	1879	-	3,3,3	0.55	0	2,2,2	0.08	0
3	EDO	B	1876	-	3,3,3	0.99	0	2,2,2	0.43	0
3	EDO	A	1866	-	3,3,3	0.59	0	2,2,2	0.50	0
3	EDO	B	1899	-	3,3,3	0.58	0	2,2,2	0.18	0
3	EDO	A	1870	-	3,3,3	0.41	0	2,2,2	0.54	0
3	EDO	B	1906	-	3,3,3	0.38	0	2,2,2	0.74	0
3	EDO	A	1888	-	3,3,3	0.44	0	2,2,2	0.45	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	B	1892	-	-	1/1/1/1	-
3	EDO	B	1898	-	-	1/1/1/1	-
3	EDO	B	1873	-	-	0/1/1/1	-
3	EDO	A	1875	-	-	0/1/1/1	-
2	17B	B	1869	-	-	2/6/28/28	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	1873	-	-	1/1/1/1	-
3	EDO	B	1880	-	-	1/1/1/1	-
3	EDO	B	1872	-	-	0/1/1/1	-
3	EDO	A	1872	-	-	0/1/1/1	-
3	EDO	A	1895	-	-	1/1/1/1	-
3	EDO	B	1877	-	-	0/1/1/1	-
3	EDO	A	1874	-	-	0/1/1/1	-
3	EDO	A	1896	-	-	1/1/1/1	-
3	EDO	B	1881	-	-	1/1/1/1	-
2	17B	A	1865	-	-	1/6/28/28	0/1/1/1
3	EDO	A	1881	-	-	0/1/1/1	-
3	EDO	B	1882	-	-	0/1/1/1	-
3	EDO	B	1878	-	-	1/1/1/1	-
3	EDO	B	1871	-	-	0/1/1/1	-
3	EDO	A	1868	-	-	0/1/1/1	-
3	EDO	B	1874	-	-	0/1/1/1	-
3	EDO	B	1886	-	-	1/1/1/1	-
3	EDO	B	1887	-	-	1/1/1/1	-
3	EDO	B	1889	-	-	1/1/1/1	-
3	EDO	B	1897	-	-	0/1/1/1	-
3	EDO	A	1871	-	-	0/1/1/1	-
3	EDO	A	1885	-	-	0/1/1/1	-
3	EDO	A	1893	-	-	0/1/1/1	-
3	EDO	B	1875	-	-	0/1/1/1	-
3	EDO	B	1905	-	-	1/1/1/1	-
3	EDO	A	1889	-	-	1/1/1/1	-
3	EDO	A	1869	-	-	0/1/1/1	-
3	EDO	B	1896	-	-	0/1/1/1	-
3	EDO	B	1908	-	-	0/1/1/1	-
3	EDO	B	1910	-	-	1/1/1/1	-
3	EDO	A	1883	-	-	0/1/1/1	-
3	EDO	B	1870	-	-	1/1/1/1	-
3	EDO	A	1867	-	-	1/1/1/1	-
3	EDO	A	1898	-	-	1/1/1/1	-
3	EDO	B	1903	-	-	1/1/1/1	-
3	EDO	B	1907	-	-	0/1/1/1	-
3	EDO	A	1894	-	-	0/1/1/1	-
3	EDO	B	1900	-	-	1/1/1/1	-
3	EDO	B	1909	-	-	1/1/1/1	-
3	EDO	A	1886	-	-	1/1/1/1	-
3	EDO	B	1879	-	-	1/1/1/1	-
3	EDO	B	1876	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	1866	-	-	0/1/1/1	-
3	EDO	B	1899	-	-	1/1/1/1	-
3	EDO	A	1870	-	-	0/1/1/1	-
3	EDO	B	1906	-	-	1/1/1/1	-
3	EDO	A	1888	-	-	1/1/1/1	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1865	17B	C3-C2	2.95	1.57	1.53
2	B	1869	17B	C4-C5	2.50	1.57	1.52
2	B	1869	17B	C2-C1	2.33	1.58	1.50
2	A	1865	17B	C1-N2	2.23	1.33	1.28
2	A	1865	17B	C2-C1	2.10	1.57	1.50
2	A	1865	17B	C5-N5	-2.09	1.43	1.47
2	B	1869	17B	C1-N2	2.00	1.33	1.28

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1865	17B	O2-C2-C3	4.46	119.81	110.53
2	B	1869	17B	O2-C2-C3	3.18	117.14	110.53
2	A	1865	17B	C9-C10-N2	2.11	114.41	110.67
2	B	1869	17B	C6-C5-C4	2.01	119.26	112.03

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1889	EDO	O1-C1-C2-O2
3	B	1910	EDO	O1-C1-C2-O2
3	A	1889	EDO	O1-C1-C2-O2
3	B	1903	EDO	O1-C1-C2-O2
2	B	1869	17B	C9-C10-N2-C1
2	B	1869	17B	N1-C7-C8-C9
3	A	1898	EDO	O1-C1-C2-O2
2	A	1865	17B	N1-C7-C8-C9
3	A	1895	EDO	O1-C1-C2-O2
3	A	1896	EDO	O1-C1-C2-O2
3	B	1879	EDO	O1-C1-C2-O2
3	B	1909	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	1867	EDO	O1-C1-C2-O2
3	B	1886	EDO	O1-C1-C2-O2
3	A	1873	EDO	O1-C1-C2-O2
3	A	1888	EDO	O1-C1-C2-O2
3	B	1870	EDO	O1-C1-C2-O2
3	B	1887	EDO	O1-C1-C2-O2
3	B	1898	EDO	O1-C1-C2-O2
3	B	1899	EDO	O1-C1-C2-O2
3	B	1900	EDO	O1-C1-C2-O2
3	B	1906	EDO	O1-C1-C2-O2
3	B	1892	EDO	O1-C1-C2-O2
3	B	1905	EDO	O1-C1-C2-O2
3	A	1886	EDO	O1-C1-C2-O2
3	B	1878	EDO	O1-C1-C2-O2
3	B	1880	EDO	O1-C1-C2-O2
3	B	1881	EDO	O1-C1-C2-O2

There are no ring outliers.

22 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1872	EDO	1	0
3	B	1877	EDO	2	0
3	A	1896	EDO	1	0
3	B	1881	EDO	1	0
3	A	1881	EDO	2	0
3	B	1878	EDO	1	0
3	B	1886	EDO	1	0
3	B	1889	EDO	1	0
3	B	1897	EDO	1	0
3	A	1885	EDO	2	0
3	A	1893	EDO	2	0
3	A	1889	EDO	2	0
3	B	1896	EDO	3	0
3	B	1870	EDO	2	0
3	A	1867	EDO	1	0
3	B	1903	EDO	1	0
3	B	1907	EDO	1	0
3	A	1894	EDO	1	0
3	B	1900	EDO	1	0
3	A	1886	EDO	2	0
3	B	1899	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1906	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

### 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	836/846 (98%)	-0.03	19 (2%) 61 61	8, 22, 39, 53	59 (7%)
1	B	841/846 (99%)	-0.08	23 (2%) 56 56	4, 21, 39, 51	33 (3%)
All	All	1677/1692 (99%)	-0.06	42 (2%) 58 58	4, 21, 39, 53	92 (5%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	470	TRP	5.1
1	B	590[A]	SER	4.6
1	A	472[A]	PHE	3.8
1	B	578	ILE	3.7
1	A	471[A]	GLY	3.5
1	A	848	LYS	3.3
1	A	590[A]	SER	3.2
1	A	849	GLY	3.1
1	B	470	TRP	3.0
1	B	867	HIS	3.0
1	B	849	GLY	2.9
1	A	469	TYR	2.9
1	A	719	LEU	2.9
1	B	846	ILE	2.9
1	A	571[A]	ALA	2.8
1	B	797[A]	ASP	2.8
1	B	28[A]	ASN	2.8
1	A	46	GLY	2.8
1	B	591[A]	ILE	2.7
1	B	796	THR	2.7
1	B	850	GLU	2.7
1	A	847	LYS	2.6
1	A	765	SER	2.6
1	B	472	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	592[A]	GLY	2.6
1	B	76	MET	2.5
1	A	75	GLY	2.4
1	B	589[A]	SER	2.4
1	B	584	ASN	2.4
1	B	793	MET	2.4
1	B	74	TYR	2.3
1	A	584[A]	ASN	2.3
1	A	74	TYR	2.2
1	A	318	ASP	2.2
1	A	473[A]	GLU	2.2
1	B	580	SER	2.2
1	B	865	GLU	2.2
1	B	582	VAL	2.2
1	B	27	GLY	2.1
1	A	76	MET	2.1
1	B	802	LEU	2.1
1	A	796	THR	2.0

## 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 oligosaccharides 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
3	EDO	B	1906	4/4	0.71	0.17	53,53,54,55	0
3	EDO	A	1895	4/4	0.76	0.17	51,51,51,52	0
3	EDO	B	1908	4/4	0.80	0.15	48,49,50,51	0
3	EDO	B	1889	4/4	0.83	0.17	45,47,47,48	0
3	EDO	A	1894	4/4	0.83	0.17	59,59,59,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	EDO	A	1888	4/4	0.83	0.15	52,52,53,55	0
3	EDO	B	1909	4/4	0.84	0.15	47,47,48,52	0
3	EDO	B	1907	4/4	0.85	0.16	44,44,45,46	0
3	EDO	B	1898	4/4	0.85	0.16	40,42,42,44	0
3	EDO	B	1897	4/4	0.85	0.18	48,49,50,51	0
3	EDO	A	1893	4/4	0.86	0.14	43,44,46,49	0
3	EDO	A	1883	4/4	0.86	0.14	42,42,43,43	0
3	EDO	A	1886	4/4	0.87	0.14	47,48,49,49	0
3	EDO	A	1898	4/4	0.87	0.13	50,51,52,54	0
3	EDO	B	1903	4/4	0.87	0.12	39,44,47,50	0
3	EDO	B	1876	4/4	0.87	0.16	26,27,29,31	0
3	EDO	B	1877	4/4	0.87	0.15	26,28,28,30	0
3	EDO	A	1873	4/4	0.87	0.12	39,39,40,40	0
3	EDO	B	1896	4/4	0.87	0.14	33,35,36,42	0
3	EDO	B	1892	4/4	0.88	0.12	38,38,39,41	0
3	EDO	B	1879	4/4	0.88	0.16	50,51,51,52	0
3	EDO	A	1875	4/4	0.88	0.14	42,42,43,45	0
3	EDO	B	1872	4/4	0.89	0.10	34,36,39,41	0
3	EDO	B	1899	4/4	0.89	0.15	23,30,34,38	0
3	EDO	A	1896	4/4	0.89	0.10	44,44,45,47	0
3	EDO	B	1905	4/4	0.89	0.14	41,42,42,44	0
4	BR	A	1882	1/1	0.89	0.13	46,46,46,46	1
5	CL	B	1902	1/1	0.89	0.22	47,47,47,47	1
3	EDO	A	1881	4/4	0.90	0.13	31,33,36,36	0
3	EDO	A	1867	4/4	0.90	0.12	27,29,29,31	0
3	EDO	A	1885	4/4	0.90	0.15	26,30,31,42	0
5	CL	B	1891	1/1	0.90	0.15	54,54,54,54	0
5	CL	B	1894	1/1	0.90	0.15	43,43,43,43	1
3	EDO	A	1872	4/4	0.90	0.12	33,37,38,40	0
3	EDO	B	1881	4/4	0.91	0.10	41,42,43,43	0
3	EDO	B	1900	4/4	0.91	0.11	33,35,37,41	0
3	EDO	B	1871	4/4	0.92	0.10	24,25,26,27	0
3	EDO	B	1880	4/4	0.92	0.09	33,35,35,38	0
3	EDO	A	1870	4/4	0.93	0.08	26,27,29,29	0
5	CL	A	1892	1/1	0.93	0.15	63,63,63,63	0
3	EDO	A	1868	4/4	0.93	0.11	15,24,30,32	0
3	EDO	B	1886	4/4	0.93	0.09	42,44,45,46	0
3	EDO	B	1873	4/4	0.93	0.09	28,30,32,34	0
3	EDO	B	1870	4/4	0.94	0.07	20,23,23,27	0
5	CL	B	1901	1/1	0.94	0.21	36,36,36,36	1
3	EDO	A	1889	4/4	0.94	0.13	17,26,31,36	0
5	CL	A	1897	1/1	0.95	0.14	28,28,28,28	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	EDO	A	1874	4/4	0.95	0.11	16,22,26,27	0
3	EDO	B	1882	4/4	0.95	0.08	18,23,24,25	0
5	CL	A	1880	1/1	0.95	0.07	33,33,33,33	0
3	EDO	B	1874	4/4	0.95	0.07	19,24,26,26	0
3	EDO	B	1887	4/4	0.96	0.06	26,27,29,30	0
3	EDO	A	1871	4/4	0.96	0.07	21,22,22,28	0
3	EDO	A	1869	4/4	0.96	0.07	17,23,24,26	0
3	EDO	B	1878	4/4	0.96	0.08	14,24,29,29	0
3	EDO	B	1910	4/4	0.96	0.10	21,31,31,34	0
4	BR	A	1878	1/1	0.96	0.04	21,21,21,21	0
2	17B	B	1869	17/17	0.96	0.07	15,19,37,39	0
2	17B	A	1865	17/17	0.97	0.07	15,18,38,42	0
3	EDO	B	1875	4/4	0.97	0.05	15,17,18,19	0
3	EDO	A	1866	4/4	0.97	0.08	16,18,21,22	0
5	CL	B	1883	1/1	0.97	0.06	31,31,31,31	0
4	BR	A	1891	1/1	0.98	0.08	21,21,21,21	1
4	BR	A	1877	1/1	0.98	0.04	23,23,23,23	1
4	BR	A	1884	1/1	0.98	0.04	32,32,32,32	1
4	BR	A	1879	1/1	0.99	0.02	21,21,21,21	1
5	CL	A	1887	1/1	0.99	0.10	23,23,23,23	1
5	CL	A	1890	1/1	0.99	0.03	18,18,18,18	1
4	BR	B	1884	1/1	0.99	0.03	30,30,30,30	1
4	BR	B	1885	1/1	0.99	0.03	22,22,22,22	1
5	CL	B	1904	1/1	0.99	0.06	19,19,19,19	0

## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.