



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 02:55 PM BST

PDB ID : 3GXF  
Title : Crystal structure of acid-beta-glucosidase with isofagomine at neutral pH  
Authors : Lieberman, R.L.  
Deposited on : 2009-04-02  
Resolution : 2.40 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.

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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 : 2.13.1  
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) : 2.13.1

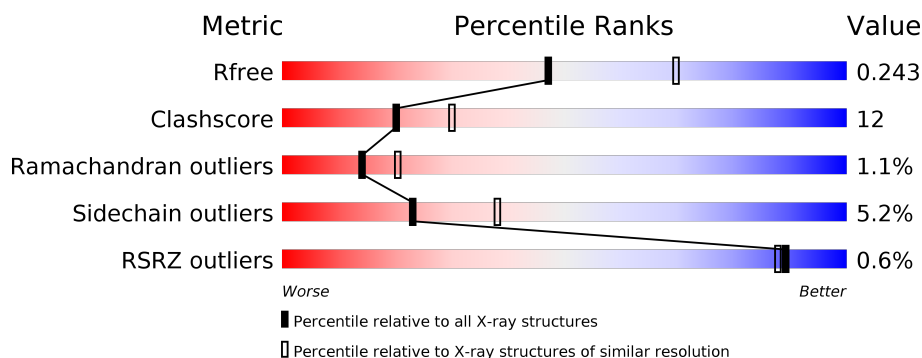
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	497	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>..</div> </div> </div>
1	B	497	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>20%</div> <div>..</div> </div> </div>
1	C	497	<div> <div></div> <div> <div></div> <div>76%</div> <div>20%</div> <div>..</div> </div> </div>
1	D	497	<div> <div></div> <div> <div></div> <div>78%</div> <div>19%</div> <div>..</div> </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	501	-	-	X	-
2	PO4	A	509	-	-	X	-
2	PO4	B	499	-	-	X	-
2	PO4	B	502	-	-	X	-
2	PO4	B	506	-	-	X	-
2	PO4	C	500	-	-	X	-
2	PO4	D	507	-	-	X	-
4	GOL	A	517	-	-	X	-
4	GOL	B	519	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17051 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.

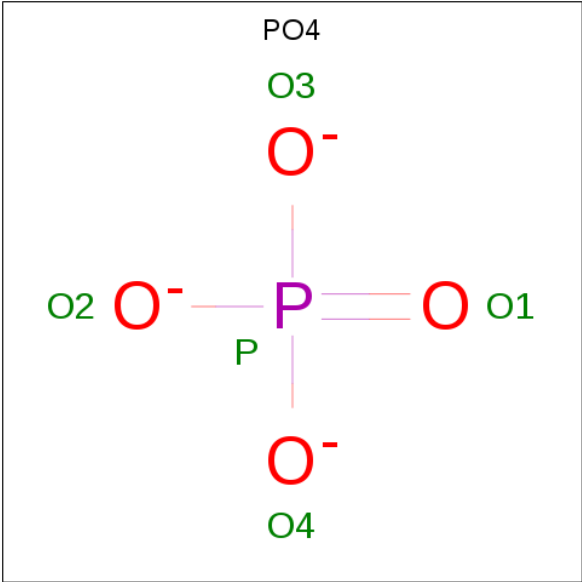
- Molecule 1 is a protein called Glucosylceramidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	0	0	0
			3930	2532	671	711	16			
1	B	497	Total	C	N	O	S	0	0	0
			3930	2532	671	711	16			
1	C	497	Total	C	N	O	S	0	0	0
			3930	2532	671	711	16			
1	D	497	Total	C	N	O	S	0	0	0
			3930	2532	671	711	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	495	HIS	ARG	variant	UNP P04062
B	495	HIS	ARG	variant	UNP P04062
C	495	HIS	ARG	variant	UNP P04062
D	495	HIS	ARG	variant	UNP P04062

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

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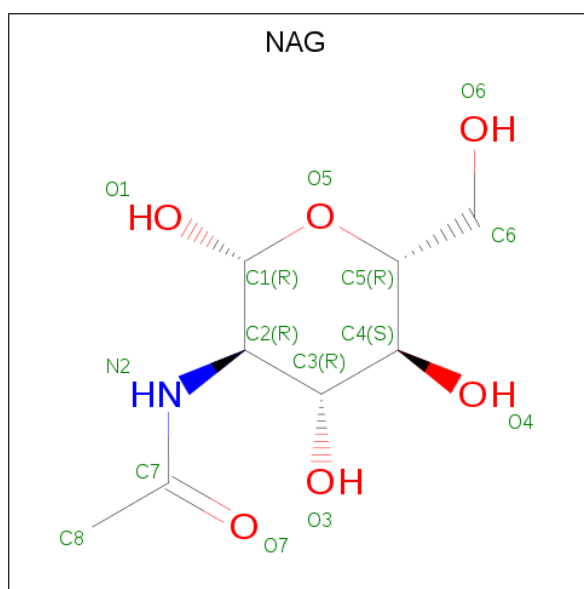
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		

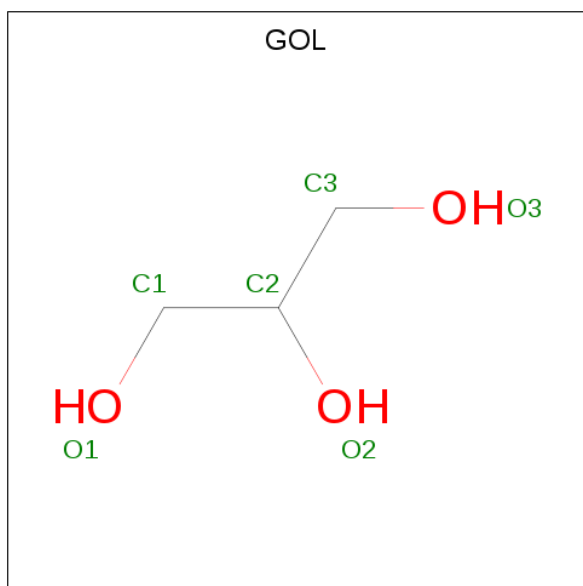
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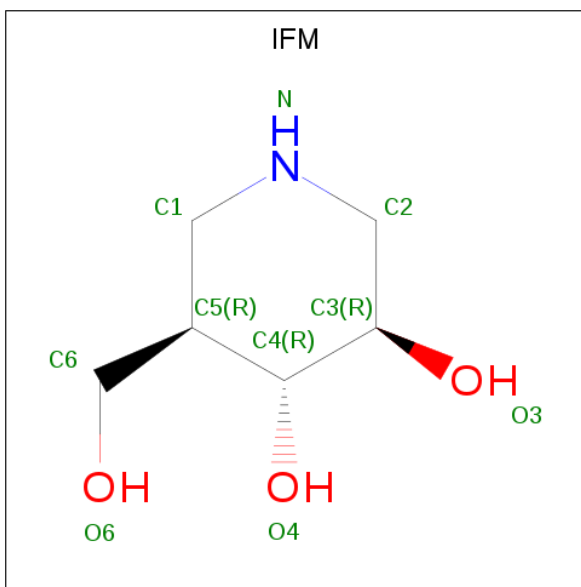
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is 5-HYDROXYMETHYL-3,4-DIHYDROXYPIPERIDINE (three-letter code: IFM) (formula:  $C_6H_{13}NO_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			10	6	1	3		
5	D	1	Total	C	N	O	0	0
			10	6	1	3		

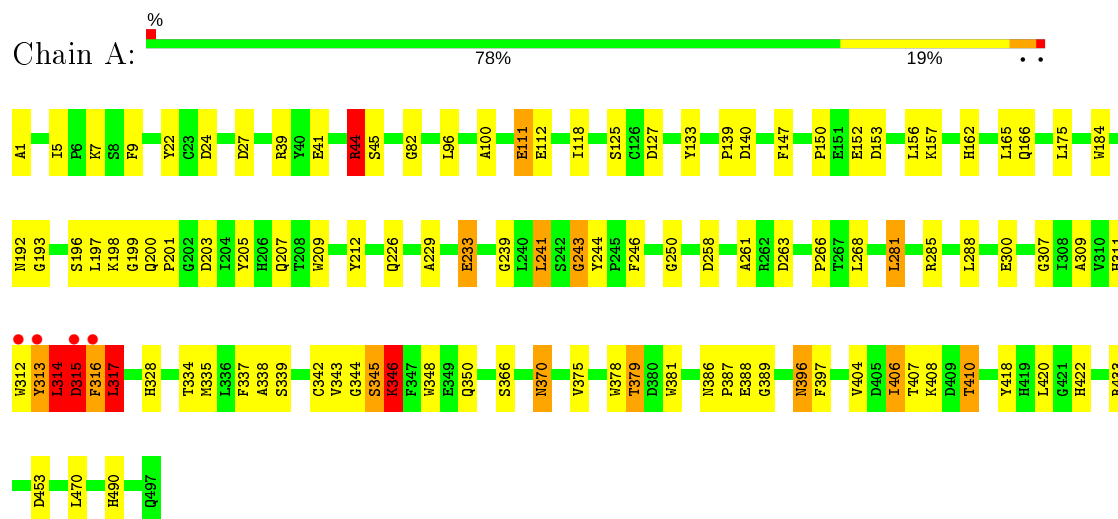
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	236	Total	O	0	0
			236	236		
6	B	243	Total	O	0	0
			243	243		
6	C	198	Total	O	0	0
			198	198		
6	D	228	Total	O	0	0
			228	228		

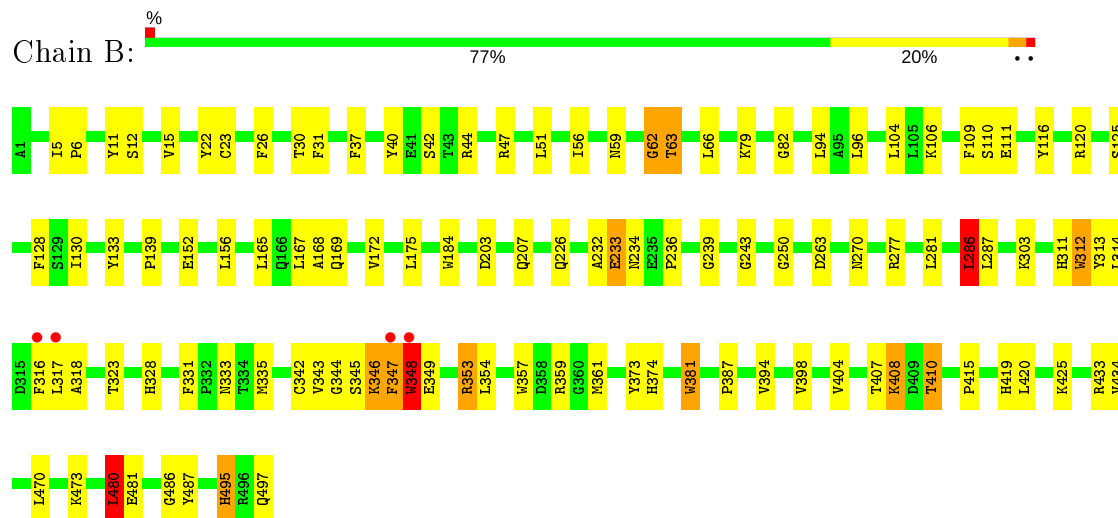
### 3 Residue-property plots

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: Glucosylceramidase

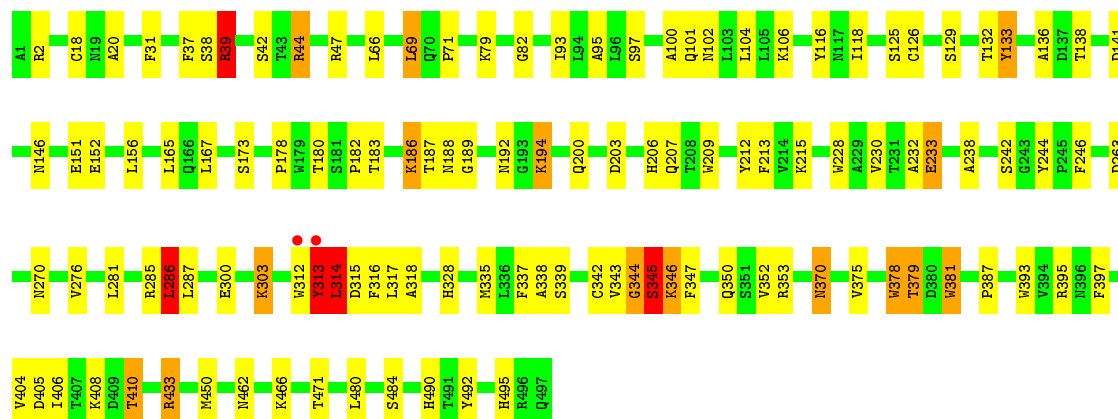


#### • Molecule 1: Glucosylceramidase



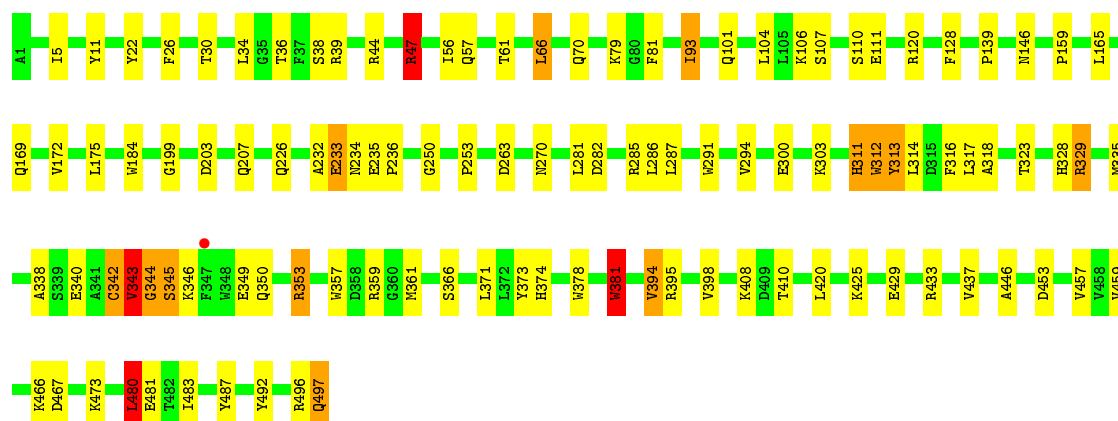
#### • Molecule 1: Glucosylceramidase





• Molecule 1: Glucosylceramidase

Chain D: 78% 19% 3%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.31Å 92.01Å 152.36Å 90.00° 111.21° 90.00°	Depositor
Resolution (Å)	32.12 – 2.40 32.12 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.2 (32.12-2.40) 93.1 (32.12-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.33 (at 2.39Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.178 , 0.245 0.178 , 0.243	Depositor DCC
$R_{free}$ test set	5209 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.5	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 15.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.460 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	17051	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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: GOL, PO4, NAG, IFM

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	0/4051	0.94	8/5523 (0.1%)
1	B	0.94	0/4051	0.91	7/5523 (0.1%)
1	C	0.97	1/4051 (0.0%)	0.91	5/5523 (0.1%)
1	D	0.93	0/4051	0.93	9/5523 (0.2%)
All	All	0.95	1/16204 (0.0%)	0.92	29/22092 (0.1%)

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	A	0	2
1	C	0	4
1	D	0	1
All	All	0	7

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	378	TRP	CG-CD1	5.50	1.44	1.36

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	39	ARG	NE-CZ-NH1	9.28	124.94	120.30
1	C	2	ARG	NE-CZ-NH2	-7.87	116.36	120.30
1	B	286	LEU	CA-CB-CG	7.71	133.03	115.30
1	D	39	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	A	317	LEU	CA-CB-CG	7.55	132.67	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	286	LEU	CA-CB-CG	7.49	132.51	115.30
1	A	243	GLY	N-CA-C	-7.30	94.86	113.10
1	B	62	GLY	N-CA-C	-7.14	95.26	113.10
1	D	343	VAL	N-CA-C	7.05	130.03	111.00
1	D	47	ARG	NE-CZ-NH1	6.98	123.79	120.30
1	B	317	LEU	CA-CB-CG	6.55	130.36	115.30
1	D	344	GLY	N-CA-C	-6.29	97.37	113.10
1	A	44	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	D	480	LEU	CA-CB-CG	5.84	128.74	115.30
1	A	313	TYR	C-N-CA	5.79	136.17	121.70
1	B	353	ARG	NE-CZ-NH2	-5.77	117.42	120.30
1	B	348	TRP	CB-CA-C	5.76	121.92	110.40
1	D	353	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	B	480	LEU	CA-CB-CG	5.58	128.13	115.30
1	A	44	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	C	313	TYR	C-N-CA	5.49	135.43	121.70
1	A	315	ASP	C-N-CA	5.39	135.17	121.70
1	B	347	PHE	N-CA-C	5.37	125.49	111.00
1	C	39	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	C	186	LYS	CD-CE-NZ	5.18	123.62	111.70
1	D	480	LEU	CB-CG-CD2	5.14	119.73	111.00
1	A	470	LEU	CB-CG-CD2	-5.14	102.27	111.00
1	A	127	ASP	CB-CG-OD1	5.12	122.91	118.30
1	D	453	ASP	CB-CG-OD1	5.06	122.85	118.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	241	LEU	Peptide
1	A	315	ASP	Peptide
1	C	313	TYR	Peptide
1	C	314	LEU	Peptide
1	C	316	PHE	Peptide
1	C	344	GLY	Peptide
1	D	342	CYS	Peptide

## 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	3930	0	3843	109	0
1	B	3930	0	3843	98	0
1	C	3930	0	3843	97	0
1	D	3930	0	3843	92	0
2	A	90	0	0	9	0
2	B	90	0	0	11	0
2	C	75	0	0	3	0
2	D	65	0	0	5	0
3	A	14	0	13	0	0
3	B	14	0	13	2	0
3	C	14	0	13	0	0
3	D	14	0	13	0	0
4	A	12	0	16	7	0
4	B	12	0	16	4	0
4	C	6	0	8	0	0
5	B	10	0	13	0	0
5	D	10	0	13	2	0
6	A	236	0	0	7	0
6	B	243	0	0	8	0
6	C	198	0	0	6	0
6	D	228	0	0	6	0
All	All	17051	0	15490	394	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 12.

All (394) 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:345:SER:CA	1:A:346:LYS:HB2	1.73	1.18
1:C:345:SER:HB3	1:C:346:LYS:HA	1.16	1.07
1:A:345:SER:HA	1:A:346:LYS:CB	1.81	1.05
1:C:39:ARG:HH11	1:C:39:ARG:HG2	1.21	1.04
1:D:286:LEU:HD21	1:D:314:LEU:CD1	1.89	1.03
1:A:313:TYR:HH	1:B:347:PHE:HZ	1.06	1.03
1:D:286:LEU:CD2	1:D:314:LEU:HD13	1.89	1.02
1:B:345:SER:HB2	6:B:669:HOH:O	1.59	1.02
1:A:44:ARG:CD	2:A:501:PO4:O2	2.13	0.96
1:D:329:ARG:HG2	1:D:329:ARG:HH11	1.28	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ARG:HD3	2:A:501:PO4:O2	1.66	0.95
1:D:286:LEU:CD2	1:D:314:LEU:CD1	2.44	0.95
1:D:270:ASN:HB3	6:D:679:HOH:O	1.67	0.94
1:C:44:ARG:HD2	2:C:500:PO4:O1	1.69	0.93
1:A:111:GLU:HB2	2:A:509:PO4:O4	1.68	0.92
1:C:312:TRP:CZ2	1:C:370:ASN:ND2	2.37	0.92
1:D:286:LEU:HD21	1:D:314:LEU:HD13	1.47	0.92
1:B:207:GLN:NE2	1:B:263:ASP:OD1	2.03	0.91
1:A:313:TYR:OH	1:B:347:PHE:HZ	1.55	0.89
1:D:487:TYR:HE1	2:D:507:PO4:O3	1.53	0.89
1:A:345:SER:HA	1:A:346:LYS:HB2	0.90	0.89
1:A:207:GLN:NE2	1:A:263:ASP:OD1	2.05	0.89
1:C:345:SER:CB	1:C:346:LYS:HA	2.04	0.88
1:A:285:ARG:HG2	1:A:317:LEU:HD22	1.56	0.87
1:D:344:GLY:HA2	1:D:345:SER:CB	2.05	0.87
1:A:370:ASN:OD1	1:A:375:VAL:HG11	1.78	0.83
1:D:487:TYR:CE1	2:D:507:PO4:O3	2.30	0.83
1:D:467:ASP:HB3	1:D:483:ILE:HD11	1.60	0.83
1:C:345:SER:HB3	1:C:346:LYS:CA	2.06	0.83
1:B:286:LEU:HD13	1:B:314:LEU:HD13	1.60	0.82
1:D:408:LYS:HB3	1:D:410:THR:HG23	1.63	0.80
1:D:226:GLN:OE1	2:D:501:PO4:O1	1.98	0.79
1:B:408:LYS:HB2	1:B:410:THR:HG23	1.63	0.78
1:B:62:GLY:HA2	1:B:63:THR:OG1	1.82	0.78
1:B:5:ILE:HD12	1:B:22:TYR:CE2	2.19	0.78
1:A:44:ARG:HD2	2:A:501:PO4:O2	1.83	0.78
1:A:111:GLU:HG2	1:A:112:GLU:OE2	1.84	0.78
1:A:312:TRP:HE1	1:A:370:ASN:HD21	1.32	0.77
1:C:192:ASN:HB2	1:C:244:TYR:O	1.84	0.77
1:C:312:TRP:CE2	1:C:370:ASN:ND2	2.53	0.77
1:D:235:GLU:OE2	1:D:311:HIS:HD2	1.68	0.77
1:D:329:ARG:NH1	1:D:329:ARG:HG2	1.99	0.76
1:D:343:VAL:HG21	1:D:359:ARG:HG2	1.68	0.76
1:D:286:LEU:HD23	1:D:314:LEU:HD13	1.68	0.75
1:B:333:ASN:N	2:B:506:PO4:O2	2.19	0.73
2:B:499:PO4:O4	6:B:522:HOH:O	2.06	0.73
1:D:286:LEU:HD21	1:D:314:LEU:HD12	1.68	0.73
1:B:96:LEU:HD21	1:B:404:VAL:HG13	1.70	0.72
1:A:312:TRP:NE1	1:A:370:ASN:ND2	2.38	0.72
1:A:226:GLN:NE2	6:A:662:HOH:O	2.22	0.71
1:A:328:HIS:HD2	6:A:627:HOH:O	1.74	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:LEU:HD23	1:D:314:LEU:CD1	2.21	0.70
1:D:344:GLY:HA2	1:D:345:SER:HB3	1.72	0.70
1:D:56:ILE:HG12	1:D:480:LEU:HD22	1.73	0.70
1:D:286:LEU:CD2	1:D:314:LEU:HD12	2.20	0.69
1:B:12:SER:OG	2:B:498:PO4:O1	2.09	0.69
1:A:1:ALA:HB2	1:A:27:ASP:OD1	1.93	0.69
1:A:397:PHE:HZ	4:A:517:GOL:H11	1.58	0.68
1:B:286:LEU:CD1	1:B:314:LEU:HD13	2.22	0.68
1:D:316:PHE:CE2	1:D:317:LEU:HD23	2.29	0.67
1:D:408:LYS:O	1:D:410:THR:HG23	1.94	0.67
1:A:397:PHE:HZ	4:A:517:GOL:C1	2.07	0.66
1:B:328:HIS:HD2	6:B:537:HOH:O	1.77	0.66
1:A:370:ASN:OD1	1:A:375:VAL:CG1	2.42	0.66
1:C:328:HIS:HD2	6:C:538:HOH:O	1.77	0.66
1:D:344:GLY:HA2	1:D:345:SER:HB2	1.76	0.66
1:C:31:PHE:O	6:C:592:HOH:O	2.14	0.66
1:A:152:GLU:HA	1:A:156:LEU:HD12	1.78	0.66
1:C:300:GLU:HA	1:C:303:LYS:HD3	1.78	0.65
1:C:285:ARG:HH11	1:C:317:LEU:HD23	1.62	0.65
1:A:312:TRP:HB3	1:A:314:LEU:HD23	1.78	0.65
1:B:56:ILE:HG12	1:B:480:LEU:HD22	1.78	0.65
1:D:44:ARG:NH2	1:D:487:TYR:CE1	2.63	0.65
1:B:152:GLU:HA	1:B:156:LEU:HD12	1.80	0.64
1:A:285:ARG:HH11	1:A:285:ARG:HG3	1.62	0.64
1:A:111:GLU:HB2	2:A:509:PO4:P	2.38	0.64
1:A:312:TRP:O	1:A:317:LEU:HG	1.98	0.63
1:A:346:LYS:HE3	1:A:348:TRP:CZ2	2.33	0.63
1:C:44:ARG:CD	2:C:500:PO4:O1	2.46	0.63
1:B:111:GLU:CG	1:B:169:GLN:HG3	2.29	0.62
1:B:286:LEU:HD13	1:B:314:LEU:CD1	2.29	0.62
1:B:111:GLU:HG3	1:B:169:GLN:HG3	1.80	0.62
1:A:39:ARG:HD2	6:A:643:HOH:O	1.99	0.62
1:B:346:LYS:HB3	1:B:348:TRP:CD1	2.35	0.62
1:C:125:SER:HB3	1:C:133:TYR:CE2	2.35	0.62
1:C:405:ASP:OD2	1:C:408:LYS:HE2	1.99	0.62
1:A:193:GLY:HA2	1:A:244:TYR:H	1.65	0.61
1:D:300:GLU:OE2	1:D:303:LYS:HE2	2.00	0.61
1:C:82:GLY:HA3	1:C:118:ILE:O	2.00	0.61
1:C:346:LYS:HE3	1:C:347:PHE:H	1.65	0.61
1:B:347:PHE:H	1:B:348:TRP:HB3	1.66	0.61
1:B:312:TRP:CD1	1:B:312:TRP:C	2.74	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:HIS:O	1:A:317:LEU:HD21	2.02	0.60
1:D:44:ARG:NH2	1:D:487:TYR:CD1	2.69	0.60
1:B:312:TRP:CD1	1:B:313:TYR:N	2.69	0.60
1:B:11:TYR:HD2	1:B:353:ARG:HH21	1.50	0.60
1:A:406:ILE:HG12	1:A:407:THR:N	2.16	0.60
1:A:45:SER:HB3	2:A:501:PO4:O4	2.02	0.60
1:D:316:PHE:CE1	1:D:346:LYS:HE2	2.37	0.59
1:D:496:ARG:O	1:D:497:GLN:HB3	2.02	0.59
1:A:313:TYR:OH	1:B:347:PHE:CZ	2.46	0.59
1:C:132:THR:O	1:C:133:TYR:HB3	2.01	0.59
1:B:44:ARG:NH2	1:B:487:TYR:CE1	2.68	0.58
1:C:138:THR:HG21	1:C:146:ASN:HD22	1.68	0.58
1:B:408:LYS:CB	1:B:410:THR:HG23	2.32	0.58
1:A:312:TRP:O	1:A:313:TYR:C	2.42	0.58
1:C:314:LEU:HD23	1:C:343:VAL:HG13	1.86	0.58
1:D:371:LEU:O	1:D:433:ARG:HD2	2.04	0.58
1:D:26:PHE:CD1	1:D:425:LYS:HE2	2.39	0.58
1:D:328:HIS:HD2	6:D:527:HOH:O	1.87	0.58
1:B:207:GLN:HE22	1:B:263:ASP:HA	1.68	0.57
1:A:193:GLY:HA2	1:A:244:TYR:N	2.19	0.57
1:D:165:LEU:HD22	1:D:172:VAL:HB	1.85	0.57
1:D:207:GLN:NE2	1:D:263:ASP:OD1	2.30	0.57
1:B:11:TYR:CD2	1:B:353:ARG:HB3	2.40	0.56
1:A:408:LYS:O	1:A:410:THR:HG22	2.05	0.56
1:A:313:TYR:HB2	6:A:592:HOH:O	2.05	0.56
1:B:331:PHE:HA	2:B:506:PO4:O1	2.05	0.56
1:C:136:ALA:O	1:C:182:PRO:HB3	2.05	0.56
1:A:350:GLN:OE1	1:A:350:GLN:HA	2.05	0.56
1:A:315:ASP:N	1:A:316:PHE:HB2	2.20	0.56
1:C:106:LYS:HG2	1:C:167:LEU:HD22	1.88	0.56
1:C:186:LYS:O	1:C:187:THR:C	2.44	0.56
1:C:93:ILE:HG12	1:C:104:LEU:HD22	1.88	0.56
1:D:79:LYS:O	1:D:429:GLU:HG3	2.05	0.56
1:B:109:PHE:O	1:B:168:ALA:HA	2.05	0.56
1:C:353:ARG:NH1	2:C:498:PO4:O1	2.39	0.56
1:A:345:SER:CB	1:A:346:LYS:HB2	2.35	0.55
1:D:312:TRP:CD1	1:D:313:TYR:N	2.74	0.55
1:A:198:LYS:HD3	1:A:205:TYR:CE2	2.41	0.55
1:B:5:ILE:HD12	1:B:22:TYR:CZ	2.42	0.55
1:C:338:ALA:O	1:C:379:THR:HG23	2.07	0.55
1:A:314:LEU:HB3	1:A:342:CYS:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:209:TRP:CZ3	1:C:212:TYR:CD2	2.95	0.55
1:C:408:LYS:O	1:C:410:THR:HG23	2.07	0.55
1:D:139:PRO:HA	1:D:184:TRP:CD1	2.42	0.55
1:A:285:ARG:NE	1:A:317:LEU:HD13	2.22	0.54
1:C:71:PRO:HB3	1:C:450:MET:CE	2.38	0.54
1:A:82:GLY:HA3	1:A:118:ILE:O	2.07	0.54
1:B:226:GLN:HG3	2:B:502:PO4:O3	2.08	0.54
1:B:130:ILE:HD12	1:C:393:TRP:HA	1.90	0.54
1:A:350:GLN:HG2	2:A:514:PO4:O2	2.08	0.54
1:B:243:GLY:N	2:B:499:PO4:O2	2.29	0.54
1:C:286:LEU:HD22	1:C:287:LEU:HG	1.88	0.54
1:B:31:PHE:CD1	1:B:495:HIS:CE1	2.96	0.54
1:C:203:ASP:C	1:C:203:ASP:OD1	2.46	0.54
1:A:343:VAL:HG22	1:A:344:GLY:HA2	1.90	0.54
1:B:128:PHE:HZ	1:B:398:VAL:HG22	1.73	0.54
1:C:178:PRO:HG3	1:C:213:PHE:CZ	2.42	0.54
1:B:165:LEU:HD22	1:B:172:VAL:HB	1.89	0.53
1:C:233:GLU:HB2	6:C:628:HOH:O	2.08	0.53
1:D:344:GLY:CA	1:D:345:SER:CB	2.84	0.53
1:B:318:ALA:HB1	1:B:323:THR:HG21	1.91	0.53
1:A:285:ARG:NH1	1:A:285:ARG:HG3	2.22	0.53
1:A:315:ASP:H	1:A:316:PHE:HB2	1.74	0.53
1:D:47:ARG:CG	1:D:47:ARG:HH11	2.22	0.53
1:A:203:ASP:C	1:A:203:ASP:OD1	2.45	0.53
1:A:397:PHE:CZ	4:A:517:GOL:C1	2.91	0.53
1:B:128:PHE:CZ	1:B:398:VAL:HG22	2.44	0.53
1:D:270:ASN:ND2	6:D:607:HOH:O	2.39	0.53
1:C:125:SER:HB3	1:C:133:TYR:CD2	2.44	0.53
1:B:313:TYR:HE2	1:B:342:CYS:SG	2.31	0.52
1:D:282:ASP:OD1	1:D:311:HIS:HE1	1.91	0.52
1:A:281:LEU:HD22	1:A:288:LEU:HD21	1.92	0.52
1:A:261:ALA:HB1	1:A:300:GLU:HG3	1.90	0.52
1:A:39:ARG:CD	6:A:643:HOH:O	2.55	0.52
1:A:312:TRP:HE1	1:A:370:ASN:ND2	1.99	0.52
1:C:313:TYR:HE2	1:C:342:CYS:SG	2.33	0.52
1:C:387:PRO:HD3	1:C:404:VAL:O	2.10	0.52
1:A:285:ARG:CZ	1:A:317:LEU:HD13	2.40	0.52
1:D:104:LEU:O	1:D:107:SER:HB2	2.09	0.52
1:C:285:ARG:NH2	1:C:318:ALA:O	2.43	0.51
1:C:207:GLN:NE2	1:C:263:ASP:OD1	2.32	0.51
1:A:315:ASP:OD2	1:A:346:LYS:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:373:TYR:O	1:B:374:HIS:HB2	2.09	0.51
1:D:234:ASN:O	1:D:235:GLU:C	2.48	0.51
1:A:197:LEU:HD11	1:A:209:TRP:CD1	2.46	0.51
1:B:270:ASN:ND2	6:B:877:HOH:O	2.35	0.51
1:C:313:TYR:CE2	1:C:342:CYS:SG	3.04	0.51
1:C:495:HIS:CD2	6:C:611:HOH:O	2.63	0.51
1:D:318:ALA:HB1	1:D:323:THR:HG21	1.93	0.51
1:A:209:TRP:CE3	1:A:209:TRP:HA	2.46	0.51
1:A:408:LYS:O	1:A:410:THR:CG2	2.59	0.51
1:B:11:TYR:HD2	1:B:353:ARG:HE	1.59	0.51
1:C:313:TYR:HH	1:C:397:PHE:HE2	1.55	0.51
1:A:339:SER:HA	1:A:379:THR:HG23	1.93	0.51
1:C:20:ALA:HA	1:C:100:ALA:HB2	1.93	0.51
1:D:357:TRP:O	1:D:361:MET:HG3	2.11	0.51
1:A:344:GLY:N	1:A:345:SER:HB2	2.25	0.50
1:D:286:LEU:HD12	1:D:287:LEU:HG	1.93	0.50
1:D:120:ARG:HD2	1:D:175:LEU:HD23	1.94	0.50
1:C:173:SER:HB3	1:C:228:TRP:CB	2.42	0.50
1:A:313:TYR:CA	1:A:314:LEU:HB2	2.41	0.50
1:C:194:LYS:HB2	1:C:242:SER:HA	1.92	0.50
1:C:39:ARG:NH1	1:C:39:ARG:HG2	1.99	0.50
1:A:162:HIS:O	1:A:166:GLN:HG3	2.11	0.50
1:A:24:ASP:HB2	4:A:518:GOL:H12	1.93	0.50
1:B:313:TYR:HE2	1:B:342:CYS:HG	1.57	0.50
1:A:386:ASN:HB2	1:A:387:PRO:CD	2.42	0.49
1:A:41:GLU:HG3	1:A:490:HIS:CD2	2.47	0.49
1:C:285:ARG:NH1	1:C:317:LEU:HB3	2.27	0.49
1:A:233:GLU:HB2	6:A:651:HOH:O	2.11	0.49
1:C:285:ARG:HH12	1:C:317:LEU:HB3	1.77	0.49
1:C:39:ARG:HH11	1:C:39:ARG:CG	2.09	0.49
1:C:433:ARG:NH1	6:C:630:HOH:O	2.34	0.49
1:B:333:ASN:HB2	2:B:506:PO4:O3	2.12	0.49
1:C:285:ARG:HG2	1:C:317:LEU:HD23	1.95	0.49
1:A:175:LEU:HA	1:A:229:ALA:O	2.13	0.49
1:A:418:TYR:O	1:A:422:HIS:HD2	1.96	0.49
1:B:12:SER:HB2	1:B:44:ARG:HD3	1.94	0.49
1:A:125:SER:HB3	1:A:133:TYR:CE2	2.48	0.49
1:C:141:ASP:OD2	1:C:146:ASN:N	2.32	0.49
1:D:287:LEU:HD22	1:D:291:TRP:NE1	2.27	0.49
1:B:37:PHE:CG	1:B:480:LEU:HD13	2.47	0.49
1:D:343:VAL:HG21	1:D:359:ARG:CG	2.40	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:GLY:H	1:A:345:SER:HB2	1.78	0.49
1:B:312:TRP:HD1	1:B:313:TYR:N	2.10	0.49
1:C:285:ARG:NH1	1:C:317:LEU:HD23	2.27	0.48
1:C:314:LEU:HG	1:C:314:LEU:O	2.11	0.48
1:C:338:ALA:HB3	1:C:378:TRP:HA	1.94	0.48
1:C:69:LEU:HD13	1:C:71:PRO:HD3	1.95	0.48
1:B:286:LEU:HD23	1:B:287:LEU:HG	1.96	0.48
1:A:370:ASN:OD1	1:A:375:VAL:CB	2.62	0.48
1:D:159:PRO:HA	2:D:503:PO4:O1	2.14	0.48
1:D:66:LEU:HD11	1:D:473:LYS:HB2	1.95	0.48
1:A:397:PHE:CZ	4:A:517:GOL:H11	2.44	0.48
1:A:312:TRP:CE2	1:A:370:ASN:ND2	2.82	0.48
1:D:338:ALA:HB1	6:D:784:HOH:O	2.13	0.48
1:D:496:ARG:O	1:D:497:GLN:CB	2.61	0.48
1:C:71:PRO:HB3	1:C:450:MET:HE1	1.96	0.48
1:D:282:ASP:OD1	1:D:311:HIS:CE1	2.67	0.47
1:C:344:GLY:C	1:C:345:SER:O	2.52	0.47
1:B:5:ILE:HD13	3:B:516:NAG:H62	1.96	0.47
1:A:198:LYS:HD3	1:A:205:TYR:CZ	2.50	0.47
1:B:236:PRO:O	1:B:250:GLY:HA2	2.14	0.47
1:B:234:ASN:OD1	1:B:311:HIS:HE1	1.97	0.47
1:B:51:LEU:H	4:B:519:GOL:C2	2.27	0.47
1:C:95:ALA:HB3	1:C:406:ILE:HG21	1.96	0.47
1:C:232:ALA:O	1:C:233:GLU:HB2	2.14	0.47
1:D:47:ARG:HG3	1:D:47:ARG:HH11	1.77	0.47
1:A:153:ASP:HA	1:A:157:LYS:HB2	1.96	0.47
1:C:314:LEU:HD23	1:C:343:VAL:CG1	2.44	0.47
1:D:408:LYS:HB3	1:D:410:THR:CG2	2.41	0.47
1:D:70:GLN:CD	1:D:437:VAL:HG23	2.35	0.47
1:D:93:ILE:O	1:D:101:GLN:HG2	2.15	0.47
1:D:207:GLN:HE22	1:D:263:ASP:HA	1.80	0.47
1:A:192:ASN:HB2	1:A:244:TYR:O	2.13	0.47
1:A:345:SER:OG	1:A:346:LYS:HB2	2.14	0.47
1:C:97:SER:O	1:C:101:GLN:HG3	2.15	0.47
1:A:209:TRP:HE3	1:A:209:TRP:HA	1.80	0.46
1:B:226:GLN:NE2	6:B:711:HOH:O	2.47	0.46
1:B:139:PRO:HA	1:B:184:TRP:CD1	2.50	0.46
1:B:94:LEU:HD12	1:B:156:LEU:HD23	1.97	0.46
1:C:186:LYS:HE2	1:C:188:ASN:OD1	2.15	0.46
1:B:31:PHE:CD1	1:B:495:HIS:HE1	2.33	0.46
1:D:26:PHE:CG	1:D:425:LYS:HE2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:81:PHE:HA	1:D:378:TRP:O	2.15	0.46
1:A:387:PRO:HD3	1:A:404:VAL:O	2.15	0.46
1:A:199:GLY:HA3	1:A:203:ASP:OD2	2.15	0.46
1:B:387:PRO:HD3	1:B:404:VAL:O	2.16	0.46
1:D:236:PRO:O	1:D:250:GLY:HA2	2.16	0.46
1:A:239:GLY:HA3	1:A:250:GLY:CA	2.46	0.46
1:B:106:LYS:HG2	1:B:167:LEU:HD22	1.97	0.46
1:A:338:ALA:HB3	1:A:378:TRP:HA	1.98	0.45
1:B:381:TRP:HA	1:B:381:TRP:CE3	2.52	0.45
1:C:370:ASN:OD1	1:C:375:VAL:HG11	2.16	0.45
1:D:57:GLN:HG3	1:D:481:GLU:OE1	2.15	0.45
1:D:11:TYR:HD2	1:D:353:ARG:HH21	1.60	0.45
1:B:433:ARG:HG2	1:B:434:VAL:N	2.31	0.45
1:B:66:LEU:HD11	1:B:473:LYS:HB2	1.99	0.45
1:C:395:ARG:O	1:C:397:PHE:HD1	2.00	0.45
1:C:270:ASN:ND2	6:C:563:HOH:O	2.49	0.45
1:C:183:THR:O	1:C:189:GLY:HA2	2.17	0.45
1:D:44:ARG:HG3	1:D:357:TRP:CG	2.51	0.45
1:A:328:HIS:HE1	2:A:504:PO4:O1	2.00	0.45
1:C:102:ASN:O	1:C:106:LYS:HG3	2.17	0.45
1:A:201:PRO:HG2	1:A:258:ASP:HB2	1.99	0.45
1:B:343:VAL:HG21	1:B:359:ARG:HG2	1.98	0.45
1:C:37:PHE:CD2	1:C:480:LEU:HG	2.52	0.45
1:B:207:GLN:HG3	2:B:510:PO4:O4	2.17	0.44
1:D:287:LEU:HD22	1:D:291:TRP:HE1	1.82	0.44
1:D:44:ARG:HG3	1:D:357:TRP:CB	2.47	0.44
1:B:104:LEU:C	1:B:104:LEU:HD23	2.38	0.44
1:D:253:PRO:HB2	1:D:294:VAL:HG11	2.00	0.44
1:A:7:LYS:HE3	1:A:9:PHE:CZ	2.53	0.44
1:C:314:LEU:CD2	1:C:343:VAL:HG13	2.47	0.44
1:D:285:ARG:NH2	1:D:318:ALA:O	2.44	0.44
1:B:381:TRP:HA	1:B:381:TRP:HE3	1.83	0.44
1:C:346:LYS:HE3	1:C:347:PHE:N	2.33	0.44
1:C:152:GLU:HA	1:C:156:LEU:HD12	1.98	0.44
1:C:462:ASN:HB2	1:C:484:SER:OG	2.17	0.44
1:A:209:TRP:CZ3	1:A:212:TYR:CD2	3.06	0.44
1:A:345:SER:CA	1:A:346:LYS:CB	2.62	0.44
1:C:79:LYS:HE2	1:C:228:TRP:CE2	2.52	0.44
1:C:408:LYS:O	1:C:410:THR:CG2	2.65	0.44
1:A:193:GLY:HA2	1:A:243:GLY:HA2	2.00	0.43
1:B:26:PHE:CD1	1:B:425:LYS:HE2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:CYS:O	1:C:129:SER:HB2	2.18	0.43
1:C:313:TYR:HE2	1:C:342:CYS:HG	1.58	0.43
1:B:44:ARG:NH2	1:B:487:TYR:CD1	2.68	0.43
1:D:146:ASN:ND2	6:D:529:HOH:O	2.50	0.43
1:B:232:ALA:O	1:B:233:GLU:HB2	2.18	0.43
1:B:344:GLY:HA2	1:B:349:GLU:HB3	1.99	0.43
1:B:316:PHE:CE1	1:B:346:LYS:NZ	2.86	0.43
1:B:37:PHE:CD2	1:B:480:LEU:HD13	2.54	0.43
1:A:313:TYR:N	1:A:314:LEU:HB2	2.33	0.43
1:B:120:ARG:HA	1:B:175:LEU:O	2.19	0.43
4:B:519:GOL:H31	6:B:659:HOH:O	2.19	0.43
1:C:337:PHE:CE2	1:C:379:THR:CG2	3.01	0.43
1:D:313:TYR:HE2	1:D:342:CYS:HG	1.65	0.43
1:A:24:ASP:OD2	4:A:518:GOL:C1	2.66	0.43
1:B:22:TYR:CD2	1:B:23:CYS:N	2.87	0.43
1:B:5:ILE:O	1:B:5:ILE:HG22	2.18	0.43
1:A:96:LEU:HB3	1:A:100:ALA:HB3	2.00	0.43
1:B:125:SER:HB3	1:B:133:TYR:CE2	2.54	0.43
1:B:345:SER:CB	6:B:669:HOH:O	2.38	0.43
1:B:40:TYR:CZ	1:B:51:LEU:HD13	2.54	0.43
1:B:5:ILE:HD13	3:B:516:NAG:C6	2.49	0.43
1:D:106:LYS:O	1:D:110:SER:HB3	2.19	0.43
1:D:287:LEU:HB3	1:D:291:TRP:CD1	2.54	0.43
1:D:44:ARG:NH1	2:D:502:PO4:O1	2.52	0.43
1:C:339:SER:HA	1:C:379:THR:HG23	2.00	0.42
1:C:71:PRO:HB3	1:C:450:MET:HE3	2.00	0.42
1:C:238:ALA:HB1	1:C:244:TYR:CZ	2.54	0.42
1:B:357:TRP:O	1:B:361:MET:HG3	2.19	0.42
1:D:235:GLU:OE2	1:D:311:HIS:CD2	2.60	0.42
1:D:395:ARG:NH1	6:D:625:HOH:O	2.46	0.42
1:A:300:GLU:HA	1:A:300:GLU:OE2	2.19	0.42
1:B:226:GLN:CG	2:B:502:PO4:O3	2.68	0.42
1:C:66:LEU:HD13	1:C:471:THR:HB	2.02	0.42
1:A:328:HIS:CD2	6:A:627:HOH:O	2.61	0.42
1:A:5:ILE:HG12	1:A:22:TYR:CE2	2.54	0.42
1:B:277:ARG:NH1	2:B:507:PO4:O3	2.53	0.42
1:B:51:LEU:H	4:B:519:GOL:H2	1.85	0.42
1:C:173:SER:HB3	1:C:228:TRP:HB2	2.02	0.42
1:D:457:VAL:HA	1:D:492:TYR:O	2.20	0.42
1:A:389:GLY:O	1:A:396:ASN:HB3	2.20	0.42
1:A:313:TYR:HA	1:A:314:LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:LYS:CB	1:A:410:THR:HG23	2.50	0.42
1:B:51:LEU:HB3	4:B:519:GOL:H2	2.01	0.42
1:C:82:GLY:HA2	1:C:116:TYR:CD1	2.55	0.42
1:D:128:PHE:CZ	1:D:398:VAL:HG22	2.55	0.42
1:A:147:PHE:C	1:A:147:PHE:CD2	2.93	0.42
1:C:230:VAL:HG13	1:C:276:VAL:HG11	2.01	0.42
1:B:415:PRO:O	1:B:419:HIS:HD2	2.03	0.42
1:B:82:GLY:HA2	1:B:116:TYR:CD1	2.55	0.42
1:D:313:TYR:CZ	5:D:512:IFM:H1C1	2.55	0.42
1:A:261:ALA:O	1:A:266:PRO:CG	2.68	0.41
1:A:345:SER:CB	1:A:397:PHE:O	2.68	0.41
1:A:111:GLU:CB	2:A:509:PO4:P	3.08	0.41
1:B:343:VAL:HG21	1:B:359:ARG:CG	2.50	0.41
1:C:165:LEU:HD23	1:C:165:LEU:N	2.36	0.41
1:C:18:CYS:O	1:C:410:THR:HB	2.20	0.41
1:D:232:ALA:O	1:D:233:GLU:HB2	2.18	0.41
1:D:234:ASN:ND2	1:D:235:GLU:HG3	2.35	0.41
1:B:6:PRO:HA	1:B:15:VAL:O	2.20	0.41
1:A:139:PRO:HA	1:A:184:TRP:CD1	2.55	0.41
1:D:5:ILE:HG12	1:D:22:TYR:CE2	2.55	0.41
1:D:47:ARG:NH1	1:D:47:ARG:CG	2.83	0.41
1:D:199:GLY:HA3	1:D:203:ASP:OD2	2.20	0.41
1:D:446:ALA:HA	1:D:459:VAL:O	2.20	0.41
1:A:307:GLY:HA2	1:A:334:THR:HG22	2.02	0.41
1:B:111:GLU:HG2	6:B:672:HOH:O	2.20	0.41
1:B:44:ARG:HG3	1:B:357:TRP:CG	2.56	0.41
1:C:38:SER:O	1:C:492:TYR:HA	2.20	0.41
1:C:93:ILE:O	1:C:101:GLN:HG2	2.20	0.41
1:B:79:LYS:NZ	2:B:500:PO4:O3	2.46	0.41
1:C:232:ALA:O	1:C:233:GLU:CB	2.69	0.41
1:C:350:GLN:HB3	1:C:352:VAL:O	2.20	0.41
1:C:337:PHE:CE2	1:C:379:THR:HG21	2.55	0.41
1:D:373:TYR:O	1:D:374:HIS:HB2	2.21	0.41
1:D:381:TRP:HA	1:D:381:TRP:HE3	1.86	0.41
1:B:354:LEU:HD23	1:B:354:LEU:HA	1.94	0.41
1:B:486:GLY:O	1:B:487:TYR:C	2.59	0.41
1:C:39:ARG:HD2	1:C:490:HIS:CE1	2.56	0.41
1:A:209:TRP:O	1:A:212:TYR:HB3	2.20	0.41
1:B:203:ASP:C	1:B:203:ASP:OD1	2.58	0.41
1:B:42:SER:HA	1:B:47:ARG:O	2.20	0.41
1:B:62:GLY:CA	1:B:63:THR:OG1	2.62	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:209:TRP:HA	1:C:209:TRP:CE3	2.56	0.41
1:C:206:HIS:HB3	1:C:263:ASP:OD2	2.21	0.41
1:D:381:TRP:HA	1:D:381:TRP:CE3	2.55	0.41
1:D:394:VAL:HG13	1:D:394:VAL:O	2.21	0.41
1:A:24:ASP:OD2	4:A:518:GOL:H12	2.21	0.41
1:D:313:TYR:HE2	1:D:342:CYS:SG	2.44	0.41
1:D:344:GLY:HA3	1:D:349:GLU:OE1	2.21	0.41
1:B:373:TYR:O	1:B:374:HIS:CB	2.70	0.40
1:C:39:ARG:NH1	1:C:39:ARG:CG	2.75	0.40
1:C:42:SER:HA	1:C:47:ARG:O	2.21	0.40
1:A:125:SER:HB3	1:A:133:TYR:CD2	2.55	0.40
1:A:453:ASP:OD1	1:A:453:ASP:C	2.59	0.40
1:A:309:ALA:HA	1:A:337:PHE:O	2.22	0.40
1:B:239:GLY:HA3	1:B:250:GLY:N	2.36	0.40
1:D:340:GLU:OE2	5:D:512:IFM:H2C1	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	495/497 (100%)	451 (91%)	37 (8%)	7 (1%)	11	15
1	B	495/497 (100%)	462 (93%)	29 (6%)	4 (1%)	19	29
1	C	495/497 (100%)	455 (92%)	34 (7%)	6 (1%)	13	19
1	D	495/497 (100%)	461 (93%)	29 (6%)	5 (1%)	15	23
All	All	1980/1988 (100%)	1829 (92%)	129 (6%)	22 (1%)	14	20

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	GLU
1	A	314	LEU
1	A	316	PHE
1	A	317	LEU
1	A	346	LYS
1	D	233	GLU
1	D	313	TYR
1	D	345	SER
1	B	233	GLU
1	C	233	GLU
1	C	345	SER
1	B	63	THR
1	A	281	LEU
1	B	281	LEU
1	B	348	TRP
1	C	281	LEU
1	C	314	LEU
1	C	381	TRP
1	D	281	LEU
1	D	381	TRP
1	C	133	TYR
1	A	150	PRO

### 5.3.2 Protein sidechains ⓘ

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	424/424 (100%)	400 (94%)	24 (6%)	20	33
1	B	424/424 (100%)	404 (95%)	20 (5%)	26	42
1	C	424/424 (100%)	403 (95%)	21 (5%)	24	40
1	D	424/424 (100%)	401 (95%)	23 (5%)	22	36
All	All	1696/1696 (100%)	1608 (95%)	88 (5%)	23	38

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ARG
1	A	111	GLU
1	A	140	ASP
1	A	165	LEU
1	A	196	SER
1	A	200	GLN
1	A	241	LEU
1	A	246	PHE
1	A	268	LEU
1	A	314	LEU
1	A	317	LEU
1	A	335	MET
1	A	345	SER
1	A	346	LYS
1	A	366	SER
1	A	370	ASN
1	A	379	THR
1	A	381	TRP
1	A	388	GLU
1	A	396	ASN
1	A	406	ILE
1	A	410	THR
1	A	420	LEU
1	A	433	ARG
1	B	30	THR
1	B	59	ASN
1	B	110	SER
1	B	286	LEU
1	B	303	LYS
1	B	312	TRP
1	B	335	MET
1	B	346	LYS
1	B	348	TRP
1	B	381	TRP
1	B	394	VAL
1	B	407	THR
1	B	408	LYS
1	B	410	THR
1	B	420	LEU
1	B	470	LEU
1	B	480	LEU
1	B	481	GLU
1	B	495	HIS

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Mol	Chain	Res	Type
1	B	497	GLN
1	C	39	ARG
1	C	44	ARG
1	C	69	LEU
1	C	151	GLU
1	C	180	THR
1	C	194	LYS
1	C	200	GLN
1	C	215	LYS
1	C	246	PHE
1	C	286	LEU
1	C	303	LYS
1	C	315	ASP
1	C	335	MET
1	C	345	SER
1	C	346	LYS
1	C	370	ASN
1	C	379	THR
1	C	381	TRP
1	C	410	THR
1	C	433	ARG
1	C	466	LYS
1	D	30	THR
1	D	34	LEU
1	D	36	THR
1	D	38	SER
1	D	47	ARG
1	D	61	THR
1	D	66	LEU
1	D	93	ILE
1	D	111	GLU
1	D	169	GLN
1	D	311	HIS
1	D	312	TRP
1	D	329	ARG
1	D	335	MET
1	D	343	VAL
1	D	350	GLN
1	D	366	SER
1	D	381	TRP
1	D	394	VAL
1	D	420	LEU

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Mol	Chain	Res	Type
1	D	466	LYS
1	D	480	LEU
1	D	497	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	ASN
1	A	146	ASN
1	A	223	HIS
1	A	328	HIS
1	A	495	HIS
1	B	59	ASN
1	B	143	GLN
1	B	200	GLN
1	B	207	GLN
1	B	226	GLN
1	B	328	HIS
1	B	365	HIS
1	B	396	ASN
1	B	495	HIS
1	C	57	GLN
1	C	146	ASN
1	C	270	ASN
1	C	328	HIS
1	C	362	GLN
1	C	365	HIS
1	C	495	HIS
1	D	146	ASN
1	D	226	GLN
1	D	311	HIS
1	D	328	HIS
1	D	365	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

75 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	PO4	D	509	-	4,4,4	0.75	0	6,6,6	0.74	0
2	PO4	B	505	-	4,4,4	0.81	0	6,6,6	0.94	0
2	PO4	C	498	-	4,4,4	1.46	0	6,6,6	0.81	0
2	PO4	C	505	-	4,4,4	0.79	0	6,6,6	0.42	0
2	PO4	B	509	-	4,4,4	0.78	0	6,6,6	0.75	0
2	PO4	B	510	-	4,4,4	0.76	0	6,6,6	0.67	0
3	NAG	C	513	1	14,14,15	0.92	0	17,19,21	1.92	5 (29%)
4	GOL	A	517	-	5,5,5	0.30	0	5,5,5	1.26	0
5	IFM	B	517	-	9,10,10	1.06	1 (11%)	9,13,13	4.41	6 (66%)
2	PO4	A	504	-	4,4,4	0.65	0	6,6,6	0.86	0
2	PO4	A	514	-	4,4,4	0.64	0	6,6,6	0.99	0
2	PO4	D	503	-	4,4,4	0.74	0	6,6,6	1.05	1 (16%)
2	PO4	D	508	-	4,4,4	0.78	0	6,6,6	0.56	0
2	PO4	A	508	-	4,4,4	1.07	0	6,6,6	0.61	0
2	PO4	B	515	-	4,4,4	0.92	0	6,6,6	0.60	0
2	PO4	C	509	-	4,4,4	0.75	0	6,6,6	0.89	0
2	PO4	A	498	-	4,4,4	1.47	0	6,6,6	1.76	2 (33%)
2	PO4	B	498	-	4,4,4	1.28	0	6,6,6	1.76	2 (33%)
2	PO4	C	511	-	4,4,4	0.59	0	6,6,6	0.52	0
2	PO4	A	512	-	4,4,4	0.74	0	6,6,6	0.65	0
3	NAG	A	516	1	14,14,15	0.93	0	17,19,21	1.63	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	518	-	5,5,5	0.34	0	5,5,5	0.64	0
2	PO4	A	502	-	4,4,4	0.82	0	6,6,6	0.80	0
2	PO4	D	504	-	4,4,4	0.82	0	6,6,6	0.44	0
2	PO4	A	499	-	4,4,4	0.93	0	6,6,6	1.35	1 (16%)
2	PO4	A	511	-	4,4,4	0.69	0	6,6,6	0.76	0
2	PO4	B	504	-	4,4,4	0.74	0	6,6,6	0.67	0
2	PO4	D	510	-	4,4,4	0.87	0	6,6,6	0.46	0
2	PO4	C	503	-	4,4,4	0.76	0	6,6,6	0.95	1 (16%)
4	GOL	C	514	-	5,5,5	0.42	0	5,5,5	0.74	0
2	PO4	B	508	-	4,4,4	0.71	0	6,6,6	0.75	0
2	PO4	C	510	-	4,4,4	0.65	0	6,6,6	0.85	0
2	PO4	B	514	-	4,4,4	0.79	0	6,6,6	0.71	0
2	PO4	A	510	-	4,4,4	0.78	0	6,6,6	1.13	0
2	PO4	C	500	-	4,4,4	0.90	0	6,6,6	1.40	1 (16%)
2	PO4	D	506	-	4,4,4	0.74	0	6,6,6	0.67	0
2	PO4	C	504	-	4,4,4	0.81	0	6,6,6	0.91	0
2	PO4	B	511	-	4,4,4	0.72	0	6,6,6	1.28	1 (16%)
2	PO4	B	499	-	4,4,4	0.94	0	6,6,6	1.26	0
2	PO4	D	502	-	4,4,4	1.01	0	6,6,6	1.19	1 (16%)
2	PO4	B	512	-	4,4,4	0.91	0	6,6,6	0.77	0
2	PO4	D	505	-	4,4,4	0.86	0	6,6,6	0.97	0
2	PO4	C	512	-	4,4,4	0.76	0	6,6,6	0.84	0
2	PO4	B	507	-	4,4,4	0.81	0	6,6,6	0.92	0
2	PO4	D	500	-	4,4,4	0.83	0	6,6,6	0.85	0
2	PO4	C	502	-	4,4,4	1.47	1 (25%)	6,6,6	0.93	0
2	PO4	A	500	-	4,4,4	1.15	0	6,6,6	1.11	0
2	PO4	C	508	-	4,4,4	0.82	0	6,6,6	0.55	0
2	PO4	A	513	-	4,4,4	0.90	0	6,6,6	0.51	0
2	PO4	A	503	-	4,4,4	0.85	0	6,6,6	0.61	0
2	PO4	B	502	-	4,4,4	0.85	0	6,6,6	1.10	1 (16%)
2	PO4	D	501	-	4,4,4	0.73	0	6,6,6	1.33	1 (16%)
2	PO4	D	498	-	4,4,4	1.01	0	6,6,6	1.89	2 (33%)
2	PO4	B	501	-	4,4,4	0.94	0	6,6,6	0.98	0
2	PO4	A	501	-	4,4,4	1.05	0	6,6,6	1.48	2 (33%)
2	PO4	C	506	-	4,4,4	1.06	0	6,6,6	0.93	0
2	PO4	C	501	-	4,4,4	0.93	0	6,6,6	1.42	1 (16%)
2	PO4	A	505	-	4,4,4	0.71	0	6,6,6	0.52	0
2	PO4	A	515	-	4,4,4	0.60	0	6,6,6	0.95	0
2	PO4	C	499	-	4,4,4	0.69	0	6,6,6	1.92	3 (50%)
2	PO4	B	500	-	4,4,4	0.82	0	6,6,6	1.26	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	PO4	A	506	-	4,4,4	0.64	0	6,6,6	1.02	0
5	IFM	D	512	-	9,10,10	1.18	1 (11%)	9,13,13	4.25	2 (22%)
3	NAG	B	516	1	14,14,15	0.70	0	17,19,21	1.87	6 (35%)
2	PO4	B	513	-	4,4,4	0.72	0	6,6,6	0.82	0
2	PO4	D	507	-	4,4,4	0.90	0	6,6,6	0.97	0
4	GOL	B	518	-	5,5,5	0.31	0	5,5,5	0.60	0
2	PO4	C	507	-	4,4,4	0.86	0	6,6,6	0.73	0
2	PO4	B	503	-	4,4,4	0.83	0	6,6,6	0.47	0
2	PO4	A	507	-	4,4,4	0.85	0	6,6,6	0.20	0
2	PO4	A	509	-	4,4,4	1.02	0	6,6,6	0.64	0
2	PO4	D	499	-	4,4,4	0.57	0	6,6,6	0.79	0
2	PO4	B	506	-	4,4,4	1.05	0	6,6,6	0.58	0
4	GOL	B	519	-	5,5,5	0.57	0	5,5,5	0.76	0
3	NAG	D	511	1	14,14,15	0.82	0	17,19,21	1.90	4 (23%)

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
5	IFM	D	512	-	-	0/2/16/16	0/1/1/1
4	GOL	B	519	-	-	4/4/4/4	-
3	NAG	B	516	1	-	0/6/23/26	0/1/1/1
3	NAG	C	513	1	-	0/6/23/26	0/1/1/1
3	NAG	D	511	1	-	0/6/23/26	0/1/1/1
4	GOL	A	517	-	-	1/4/4/4	-
4	GOL	B	518	-	-	4/4/4/4	-
4	GOL	C	514	-	-	4/4/4/4	-
3	NAG	A	516	1	-	0/6/23/26	0/1/1/1
4	GOL	A	518	-	-	0/4/4/4	-
5	IFM	B	517	-	-	0/2/16/16	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	512	IFM	C3-C4	2.53	1.56	1.52
5	B	517	IFM	C6-C5	2.33	1.57	1.52
2	C	502	PO4	P-O1	2.05	1.55	1.50

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	512	IFM	C1-N-C2	11.43	124.24	111.70
5	B	517	IFM	C1-N-C2	9.71	122.35	111.70
5	B	517	IFM	C2-C3-C4	6.07	117.45	110.33
3	D	511	NAG	C1-O5-C5	5.21	119.25	112.19
5	B	517	IFM	O4-C4-C5	-4.92	101.76	110.08
5	D	512	IFM	C2-C3-C4	4.78	115.94	110.33
3	B	516	NAG	C2-N2-C7	-4.11	117.05	122.90
3	C	513	NAG	O4-C4-C5	3.93	119.05	109.30
3	C	513	NAG	C1-O5-C5	3.86	117.43	112.19
3	B	516	NAG	C4-C3-C2	3.43	116.05	111.02
2	D	498	PO4	O2-P-O1	-3.21	99.13	110.89
3	D	511	NAG	C2-N2-C7	-3.11	118.48	122.90
3	A	516	NAG	C1-C2-N2	3.04	115.68	110.49
3	A	516	NAG	C1-O5-C5	2.98	116.22	112.19
3	B	516	NAG	C1-O5-C5	2.91	116.14	112.19
3	C	513	NAG	C8-C7-N2	2.91	121.02	116.10
3	B	516	NAG	O5-C1-C2	-2.83	106.81	111.29
2	C	499	PO4	O4-P-O3	2.83	117.04	107.97
2	A	498	PO4	O4-P-O2	2.78	116.89	107.97
2	B	498	PO4	O3-P-O1	-2.64	101.22	110.89
3	C	513	NAG	C1-C2-N2	2.64	115.00	110.49
3	D	511	NAG	O4-C4-C5	2.64	115.85	109.30
2	C	499	PO4	O4-P-O1	-2.59	101.40	110.89
3	B	516	NAG	O4-C4-C5	2.58	115.70	109.30
3	A	516	NAG	C4-C3-C2	2.48	114.65	111.02
3	C	513	NAG	O3-C3-C2	-2.47	104.35	109.47
2	A	501	PO4	O3-P-O1	2.44	119.84	110.89
5	B	517	IFM	O3-C3-C4	2.40	114.94	110.14
2	D	501	PO4	O4-P-O1	-2.36	102.25	110.89
2	B	498	PO4	O3-P-O2	2.34	115.48	107.97
5	B	517	IFM	O3-C3-C2	-2.34	105.05	109.61
3	B	516	NAG	O3-C3-C4	-2.31	105.00	110.35
2	A	499	PO4	O4-P-O3	2.30	115.35	107.97
2	C	501	PO4	O3-P-O1	-2.29	102.51	110.89
3	A	516	NAG	C8-C7-N2	2.28	119.95	116.10
2	C	499	PO4	O4-P-O2	2.25	115.20	107.97
2	D	498	PO4	O3-P-O1	2.24	119.09	110.89
3	A	516	NAG	O4-C4-C5	2.22	114.81	109.30
2	B	502	PO4	O3-P-O2	2.20	115.05	107.97
2	A	501	PO4	O4-P-O1	-2.20	102.85	110.89
2	B	511	PO4	O4-P-O3	2.18	114.97	107.97
5	B	517	IFM	O6-C6-C5	2.17	116.33	111.36
2	D	502	PO4	O4-P-O1	-2.17	102.97	110.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	PO4	O4-P-O2	2.12	114.78	107.97
2	C	503	PO4	O4-P-O3	2.06	114.57	107.97
2	A	498	PO4	O4-P-O3	-2.04	101.41	107.97
3	D	511	NAG	O5-C1-C2	-2.02	108.10	111.29
2	D	503	PO4	O4-P-O1	-2.00	103.57	110.89

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	519	GOL	O1-C1-C2-O2
4	B	519	GOL	O1-C1-C2-C3
4	B	519	GOL	C1-C2-C3-O3
4	C	514	GOL	O1-C1-C2-C3
4	B	518	GOL	C1-C2-C3-O3
4	B	519	GOL	O2-C2-C3-O3
4	B	518	GOL	O1-C1-C2-C3
4	B	518	GOL	O2-C2-C3-O3
4	A	517	GOL	O2-C2-C3-O3
4	B	518	GOL	O1-C1-C2-O2
4	C	514	GOL	O1-C1-C2-O2
4	C	514	GOL	C1-C2-C3-O3
4	C	514	GOL	O2-C2-C3-O3

There are no ring outliers.

22 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	498	PO4	1	0
2	B	510	PO4	1	0
4	A	517	GOL	4	0
2	A	504	PO4	1	0
2	A	514	PO4	1	0
2	D	503	PO4	1	0
2	B	498	PO4	1	0
4	A	518	GOL	3	0
2	C	500	PO4	2	0
2	B	499	PO4	2	0
2	D	502	PO4	1	0
2	B	507	PO4	1	0
2	B	502	PO4	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	PO4	1	0
2	A	501	PO4	4	0
2	B	500	PO4	1	0
5	D	512	IFM	2	0
3	B	516	NAG	2	0
2	D	507	PO4	2	0
2	A	509	PO4	3	0
2	B	506	PO4	3	0
4	B	519	GOL	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 [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<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	497/497 (100%)	-0.50	4 (0%) 86 84	14, 24, 44, 62	0
1	B	497/497 (100%)	-0.52	4 (0%) 86 84	13, 23, 43, 64	0
1	C	497/497 (100%)	-0.49	2 (0%) 92 91	14, 24, 43, 61	0
1	D	497/497 (100%)	-0.51	1 (0%) 95 94	13, 23, 42, 64	0
All	All	1988/1988 (100%)	-0.50	11 (0%) 89 88	13, 24, 43, 64	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	348	TRP	4.0
1	B	347	PHE	3.3
1	B	317	LEU	3.2
1	A	316	PHE	3.1
1	A	312	TRP	2.7
1	C	312	TRP	2.6
1	A	313	TYR	2.4
1	C	313	TYR	2.4
1	B	316	PHE	2.4
1	A	315	ASP	2.2
1	D	347	PHE	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](#)

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

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	PO4	C	508	5/5	0.68	0.24	107,107,109,109	0
2	PO4	C	506	5/5	0.78	0.30	110,110,111,112	0
2	PO4	B	510	5/5	0.80	0.27	116,117,117,117	0
2	PO4	A	513	5/5	0.80	0.30	105,106,106,106	0
2	PO4	C	502	5/5	0.80	0.39	118,118,118,119	0
2	PO4	D	506	5/5	0.81	0.28	101,102,103,103	0
2	PO4	D	505	5/5	0.81	0.33	95,96,96,97	0
2	PO4	C	511	5/5	0.81	0.35	98,98,99,99	0
2	PO4	C	509	5/5	0.82	0.18	95,95,96,96	0
2	PO4	B	509	5/5	0.82	0.30	81,82,82,83	0
2	PO4	A	510	5/5	0.83	0.19	81,82,83,83	0
2	PO4	D	507	5/5	0.83	0.18	72,73,75,77	0
2	PO4	D	510	5/5	0.84	0.30	116,117,117,117	0
2	PO4	B	504	5/5	0.85	0.32	89,89,90,91	0
2	PO4	B	508	5/5	0.85	0.17	113,113,113,113	0
2	PO4	A	503	5/5	0.86	0.36	100,100,101,101	0
2	PO4	C	507	5/5	0.87	0.17	78,80,81,81	0
2	PO4	A	509	5/5	0.87	0.24	98,100,100,101	0
2	PO4	B	507	5/5	0.88	0.20	91,92,94,94	0
2	PO4	A	507	5/5	0.88	0.17	103,104,104,105	0
2	PO4	B	502	5/5	0.88	0.26	77,79,80,80	0
2	PO4	C	510	5/5	0.89	0.20	77,80,80,81	0
2	PO4	B	514	5/5	0.89	0.16	95,95,95,96	0
2	PO4	D	509	5/5	0.90	0.19	83,85,85,85	0
2	PO4	D	503	5/5	0.90	0.19	93,93,95,95	0
4	GOL	B	519	6/6	0.90	0.25	35,40,41,42	0
3	NAG	B	516	14/15	0.90	0.14	39,41,44,45	0
2	PO4	B	506	5/5	0.90	0.20	107,107,108,108	0
2	PO4	D	508	5/5	0.91	0.30	85,86,87,87	0
2	PO4	D	501	5/5	0.91	0.24	49,53,55,58	0
2	PO4	A	511	5/5	0.91	0.21	76,77,78,78	0
2	PO4	A	515	5/5	0.91	0.25	71,72,73,74	0
2	PO4	B	515	5/5	0.91	0.17	67,67,70,70	0
3	NAG	D	511	14/15	0.91	0.14	24,39,42,42	0
2	PO4	C	505	5/5	0.92	0.33	103,104,104,105	0
2	PO4	C	504	5/5	0.92	0.26	77,78,79,80	0
2	PO4	D	504	5/5	0.92	0.15	78,79,79,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	PO4	B	503	5/5	0.92	0.37	105,106,107,107	0
2	PO4	A	506	5/5	0.93	0.09	67,69,70,71	0
2	PO4	A	502	5/5	0.93	0.11	74,74,76,77	0
2	PO4	B	513	5/5	0.93	0.16	61,64,65,66	0
2	PO4	A	505	5/5	0.93	0.30	84,85,85,85	0
2	PO4	B	505	5/5	0.93	0.23	72,72,73,75	0
4	GOL	A	517	6/6	0.94	0.15	36,42,42,43	0
2	PO4	A	504	5/5	0.94	0.11	60,63,65,65	0
2	PO4	A	501	5/5	0.94	0.12	49,50,52,56	0
2	PO4	A	508	5/5	0.94	0.16	61,61,62,62	0
2	PO4	A	512	5/5	0.95	0.12	61,63,64,64	0
4	GOL	A	518	6/6	0.95	0.22	54,57,57,57	0
2	PO4	C	500	5/5	0.95	0.14	51,53,54,54	0
4	GOL	B	518	6/6	0.95	0.14	38,41,43,44	0
5	IFM	D	512	10/10	0.95	0.17	24,30,34,34	0
2	PO4	D	502	5/5	0.96	0.16	49,55,56,57	0
5	IFM	B	517	10/10	0.96	0.14	22,24,30,31	0
4	GOL	C	514	6/6	0.96	0.16	38,39,40,41	0
3	NAG	A	516	14/15	0.96	0.13	25,30,32,33	0
2	PO4	B	511	5/5	0.96	0.13	66,66,67,68	0
2	PO4	A	514	5/5	0.97	0.10	62,63,64,64	0
3	NAG	C	513	14/15	0.97	0.11	25,30,31,34	0
2	PO4	C	512	5/5	0.97	0.16	69,69,70,71	0
2	PO4	C	501	5/5	0.97	0.11	37,38,42,42	0
2	PO4	C	503	5/5	0.97	0.14	61,61,63,63	0
2	PO4	D	500	5/5	0.97	0.10	33,33,39,39	0
2	PO4	B	512	5/5	0.98	0.19	51,53,54,54	0
2	PO4	B	499	5/5	0.98	0.11	35,37,38,41	0
2	PO4	B	501	5/5	0.98	0.09	42,43,48,49	0
2	PO4	A	499	5/5	0.98	0.11	40,41,42,43	0
2	PO4	D	499	5/5	0.98	0.09	37,39,40,41	0
2	PO4	C	499	5/5	0.98	0.11	35,36,38,42	0
2	PO4	B	500	5/5	0.98	0.12	26,34,37,38	0
2	PO4	B	498	5/5	0.99	0.10	24,25,30,31	0
2	PO4	C	498	5/5	0.99	0.14	25,27,30,30	0
2	PO4	D	498	5/5	0.99	0.11	21,24,25,26	0
2	PO4	A	500	5/5	0.99	0.12	41,41,43,45	0
2	PO4	A	498	5/5	0.99	0.13	24,25,32,32	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.