



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:20 PM GMT

PDB ID : 1OKE  
Title : Crystal structure of the dengue 2 virus envelope protein in complex with n-oc  
tyl-beta-D-glucoside  
Authors : Modis, Y.; Harrison, S.C.  
Deposited on : 2003-07-22  
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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

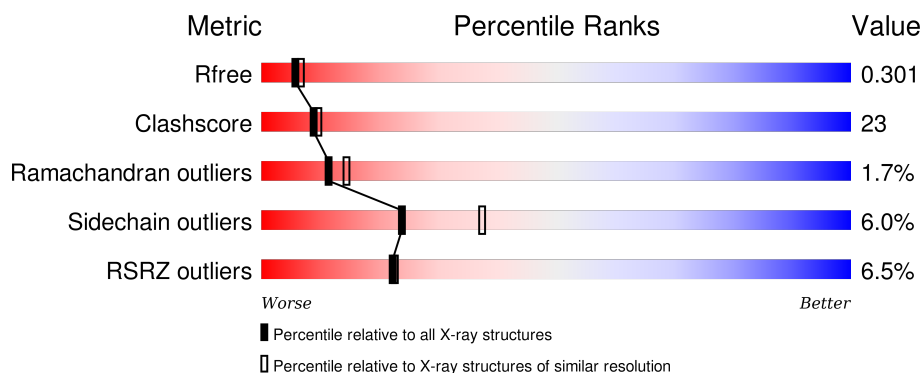
# 1 Overall quality at a glance

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))
$R_{free}$	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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. 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	394	<div> <div>11%</div> <div>61%</div> <div>35%</div> <div>• •</div> </div>
1	B	394	<div> <div>2%</div> <div>65%</div> <div>31%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	BOG	A	1400	-	-	-	X
4	BOG	B	1400	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6426 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.

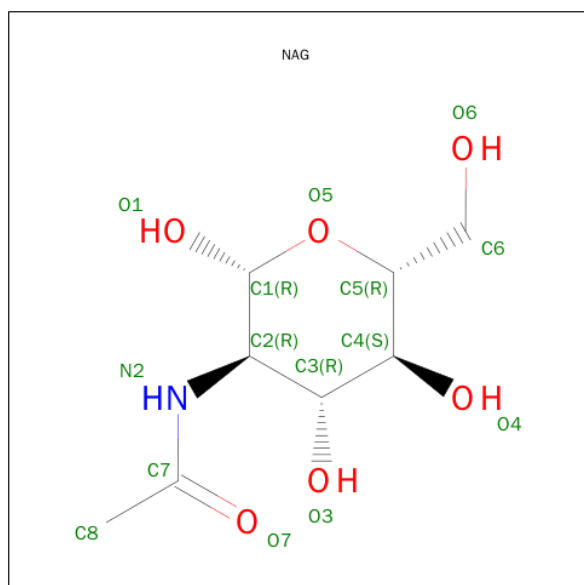
- Molecule 1 is a protein called MAJOR ENVELOPE PROTEIN E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			3062	1931	526	579	26			
1	B	394	Total	C	N	O	S	0	0	0
			3062	1931	526	579	26			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	71	GLU	ASP	CONFLICT	UNP P12823
A	390	ASN	ASP	CONFLICT	UNP P12823
B	71	GLU	ASP	CONFLICT	UNP P12823
B	390	ASN	ASP	CONFLICT	UNP P12823

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).

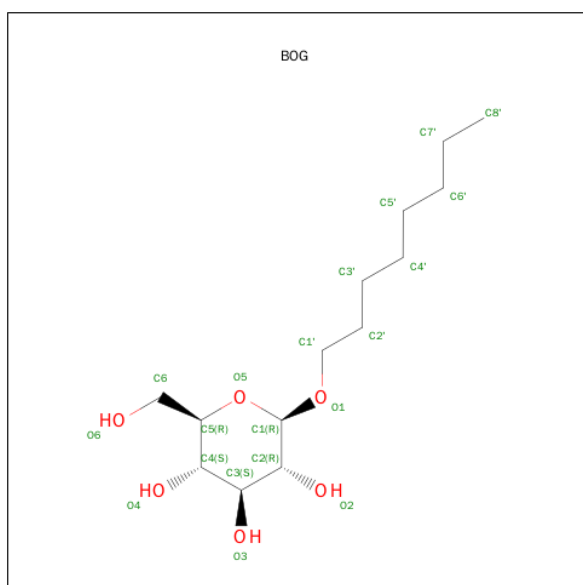


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	4	Total	C	N	O	0	0
			49	28	2	19		
3	B	4	Total	C	N	O	0	0
			49	28	2	19		

- Molecule 4 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula: C<sub>14</sub>H<sub>28</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			20	14	6		
4	B	1	Total	C	O	0	0
			20	14	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	49	Total	O	0	0
			49	49		

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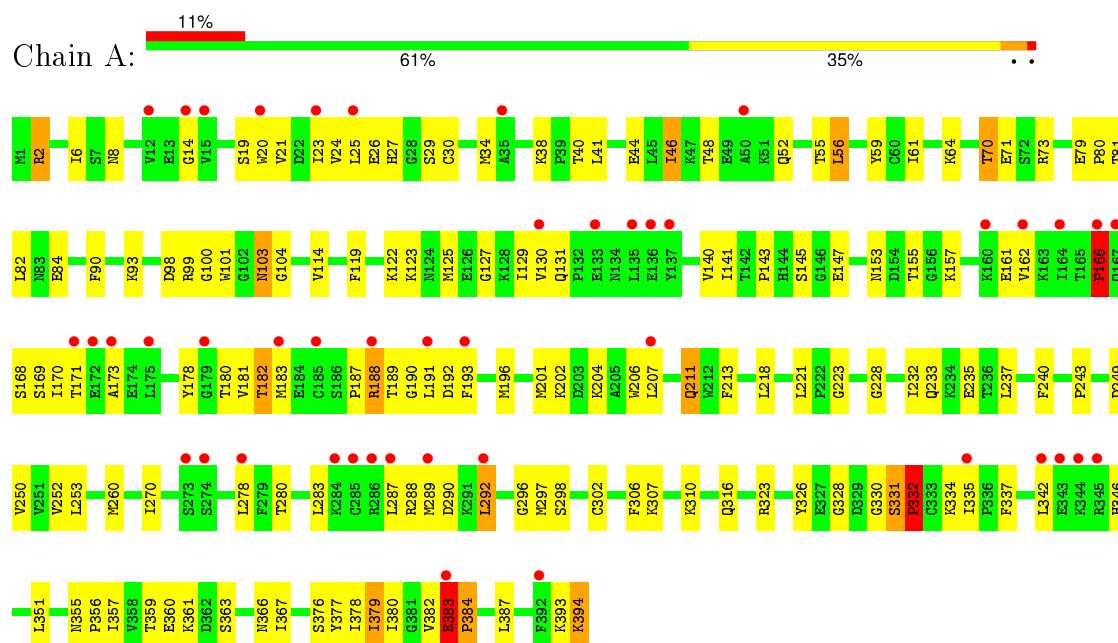
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	87	Total	O	0	0
			87	87		

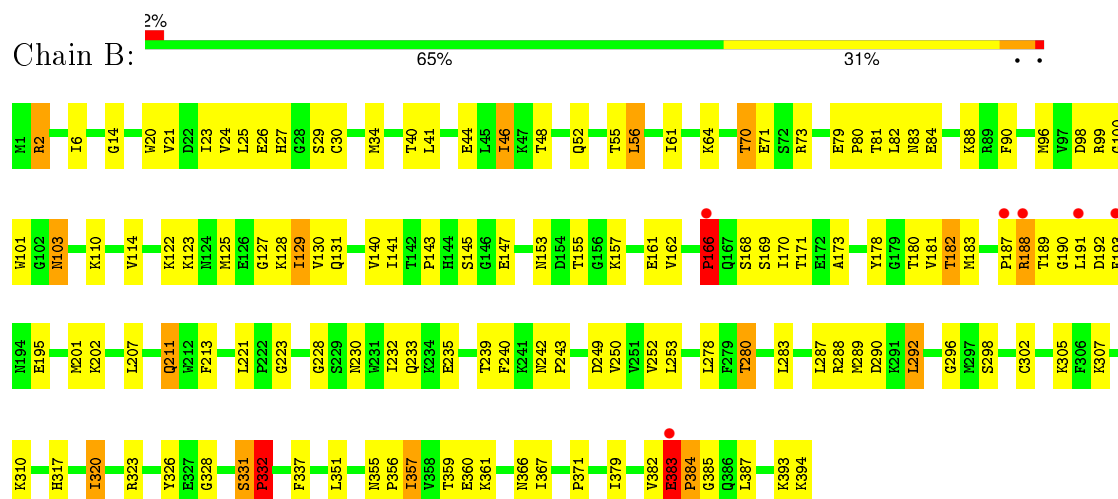
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: MAJOR ENVELOPE PROTEIN E



#### • Molecule 1: MAJOR ENVELOPE PROTEIN E



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.60Å 81.60Å 287.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.00 – 2.40 44.59 – 2.15	Depositor EDS
% Data completeness (in resolution range)	94.7 (45.00-2.40) 80.3 (44.59-2.15)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 2.16Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.263 , 0.294 0.271 , 0.301	Depositor DCC
$R_{free}$ test set	2130 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	55.3	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 47.5	EDS
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 49270 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6426	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.20% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: FUL, BMA, NAG, BOG

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	0.51	2/3124 (0.1%)	0.79	5/4219 (0.1%)
1	B	0.56	0/3124	0.81	3/4219 (0.1%)
All	All	0.53	2/6248 (0.0%)	0.80	8/8438 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	383	GLU	CG-CD	-7.70	1.40	1.51
1	A	383	GLU	CB-CG	-5.16	1.42	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	190	GLY	N-CA-C	-9.12	90.29	113.10
1	A	190	GLY	N-CA-C	-9.00	90.61	113.10
1	A	383	GLU	CA-CB-CG	-8.05	95.70	113.40
1	B	189	THR	N-CA-C	6.00	127.19	111.00
1	A	189	THR	N-CA-C	5.92	126.97	111.00
1	B	383	GLU	CA-CB-CG	5.75	126.05	113.40
1	A	383	GLU	N-CA-C	5.31	125.35	111.00
1	A	104	GLY	N-CA-C	5.19	126.07	113.10

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	3062	0	3065	154	0
1	B	3062	0	3065	140	0
2	A	14	0	13	0	0
2	B	14	0	13	1	0
3	A	49	0	43	3	0
3	B	49	0	43	2	0
4	A	20	0	28	1	0
4	B	20	0	28	1	0
5	A	49	0	0	15	0
5	B	87	0	0	11	0
All	All	6426	0	6298	294	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 23.

All (294) 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:B:166:PRO:HB3	1:B:187:PRO:CG	1.52	1.37
1:A:166:PRO:HB3	1:A:187:PRO:CG	1.52	1.36
1:A:166:PRO:CB	1:A:187:PRO:HG3	1.73	1.16
1:B:166:PRO:CB	1:B:187:PRO:HG3	1.74	1.16
1:A:289:MET:HA	5:A:2023:HOH:O	1.57	1.04
1:B:166:PRO:HB3	1:B:187:PRO:HG3	1.02	1.01
1:A:323:ARG:HE	1:A:366:ASN:HD21	1.04	1.01
1:A:166:PRO:HB3	1:A:187:PRO:HG3	1.02	0.99
1:B:323:ARG:HE	1:B:366:ASN:HD21	1.13	0.94
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.51	0.91
1:B:337:PHE:CD1	1:B:351:LEU:HD21	2.10	0.86
1:B:46:ILE:HG12	1:B:140:VAL:HG23	1.57	0.86
1:A:377:TYR:HD1	5:A:2043:HOH:O	1.58	0.86
1:A:201:MET:HB2	5:A:2015:HOH:O	1.76	0.85
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.58	0.83
1:B:359:THR:HG22	1:B:360:GLU:HG3	1.60	0.82
1:A:140:VAL:HG22	1:A:161:GLU:HG2	1.61	0.81
1:A:166:PRO:HB3	1:A:187:PRO:CD	2.11	0.80
1:B:140:VAL:HG22	1:B:161:GLU:HG2	1.62	0.80
1:A:323:ARG:HE	1:A:366:ASN:ND2	1.79	0.80
1:B:289:MET:HE1	5:B:2007:HOH:O	1.79	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:PRO:HB3	1:B:187:PRO:CD	2.12	0.79
1:A:46:ILE:HG12	1:A:140:VAL:HG23	1.64	0.79
1:A:178:TYR:HA	5:A:2012:HOH:O	1.83	0.79
1:A:323:ARG:NE	1:A:366:ASN:HD21	1.82	0.78
1:A:56:LEU:HG	1:A:129:ILE:HD11	1.66	0.78
1:A:61:ILE:HD11	1:A:123:LYS:HG2	1.64	0.78
1:A:337:PHE:CD1	1:A:351:LEU:HD21	2.19	0.77
1:A:359:THR:HG22	1:A:360:GLU:HG3	1.64	0.77
1:B:61:ILE:HD11	1:B:123:LYS:HG2	1.67	0.76
1:B:155:THR:HG22	1:B:157:LYS:H	1.52	0.74
1:B:191:LEU:HD12	1:B:192:ASP:H	1.50	0.74
1:B:323:ARG:HE	1:B:366:ASN:ND2	1.86	0.73
1:A:38:LYS:HE3	5:A:2003:HOH:O	1.89	0.73
1:A:331:SER:HB2	1:A:332:PRO:HD2	1.71	0.72
1:A:19:SER:HB2	5:A:2003:HOH:O	1.90	0.72
1:A:101:TRP:CE2	1:B:310:LYS:HE3	2.24	0.71
1:A:191:LEU:HD12	1:A:192:ASP:H	1.53	0.71
1:B:323:ARG:NE	1:B:366:ASN:HD21	1.89	0.71
1:A:48:THR:OG1	1:A:280:THR:HG21	1.90	0.71
1:A:166:PRO:HB3	1:A:187:PRO:HG2	1.68	0.70
1:B:166:PRO:HB3	1:B:187:PRO:HG2	1.69	0.70
1:A:155:THR:HG22	1:A:157:LYS:H	1.56	0.70
1:B:48:THR:OG1	1:B:280:THR:HG21	1.91	0.69
1:B:56:LEU:HG	1:B:129:ILE:HD11	1.73	0.69
1:A:14:GLY:HA3	1:A:21:VAL:HG12	1.74	0.68
1:B:14:GLY:HA3	1:B:21:VAL:HG12	1.76	0.68
1:A:166:PRO:CG	1:A:187:PRO:HG3	2.25	0.67
1:B:331:SER:HB2	1:B:332:PRO:HD2	1.77	0.67
1:A:243:PRO:HG3	1:B:278:LEU:HD22	1.75	0.66
1:B:166:PRO:CG	1:B:187:PRO:HG3	2.25	0.66
1:B:357:ILE:O	1:B:357:ILE:HG12	1.94	0.66
1:B:239:THR:HA	5:B:2056:HOH:O	1.93	0.66
1:A:55:THR:O	1:A:223:GLY:HA3	1.97	0.65
1:A:64:LYS:HB2	1:A:122:LYS:HD2	1.79	0.65
1:B:64:LYS:HB2	1:B:122:LYS:HD2	1.78	0.64
1:B:207:LEU:HD22	4:B:1400:BOG:H8'3	1.79	0.64
1:A:71:GLU:HG2	1:A:81:THR:H	1.63	0.64
1:B:141:ILE:N	1:B:141:ILE:HD12	2.13	0.64
1:A:141:ILE:HD12	1:A:141:ILE:N	2.13	0.64
3:A:1397:NAG:H4	3:A:1399:BMA:O2	1.98	0.63
1:A:331:SER:HB2	1:A:332:PRO:CD	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:PRO:CB	1:A:187:PRO:CG	2.45	0.63
1:B:240:PHE:CD2	1:B:250:VAL:HG22	2.33	0.63
1:B:320:ILE:HD13	1:B:371:PRO:HD3	1.80	0.63
1:A:206:TRP:HZ3	5:A:2015:HOH:O	1.81	0.63
1:B:280:THR:O	1:B:280:THR:HG22	1.99	0.62
1:B:331:SER:HB2	1:B:332:PRO:CD	2.29	0.62
1:B:52:GLN:NE2	1:B:52:GLN:HA	2.14	0.62
1:B:166:PRO:CB	1:B:187:PRO:CG	2.45	0.62
1:A:14:GLY:CA	1:A:21:VAL:HG12	2.30	0.62
3:A:1396:NAG:H62	3:A:1397:NAG:H82	1.81	0.62
1:B:178:TYR:O	1:B:292:LEU:HD23	1.99	0.62
3:B:1397:NAG:H4	3:B:1399:BMA:O2	1.99	0.62
1:A:288:ARG:HG2	1:A:288:ARG:HH11	1.64	0.62
1:A:130:VAL:O	1:A:193:PHE:HB3	1.99	0.61
1:A:178:TYR:O	1:A:292:LEU:HD23	2.00	0.61
1:A:100:GLY:H	1:A:103:ASN:HD21	1.48	0.61
3:B:1396:NAG:H62	3:B:1397:NAG:H82	1.83	0.60
1:A:204:LYS:HB2	5:A:2015:HOH:O	2.01	0.60
1:A:162:VAL:HG21	1:A:183:MET:CE	2.30	0.60
1:A:337:PHE:CD2	1:A:380:ILE:HD12	2.37	0.60
1:A:328:GLY:O	1:A:361:LYS:HE3	2.02	0.60
1:A:310:LYS:HE3	1:B:101:TRP:CE2	2.37	0.59
1:B:23:ILE:HD12	1:B:25:LEU:HD21	1.82	0.59
1:A:48:THR:HG21	1:A:280:THR:CG2	2.32	0.59
1:A:162:VAL:HG21	1:A:183:MET:HE1	1.83	0.59
1:B:14:GLY:CA	1:B:21:VAL:HG12	2.32	0.59
1:B:317:HIS:ND1	5:B:2070:HOH:O	2.32	0.59
1:A:307:LYS:NZ	1:A:307:LYS:HB3	2.18	0.59
1:B:100:GLY:H	1:B:103:ASN:HD21	1.50	0.59
1:A:280:THR:O	1:A:280:THR:HG22	2.03	0.59
1:B:288:ARG:HH11	1:B:288:ARG:HG2	1.66	0.59
1:A:316:GLN:HE22	1:B:110:LYS:HE2	1.67	0.58
1:B:162:VAL:HG21	1:B:183:MET:HE1	1.84	0.58
1:A:240:PHE:CD2	1:A:250:VAL:HG22	2.38	0.58
2:B:1395:NAG:O6	5:B:2084:HOH:O	2.17	0.58
1:B:328:GLY:O	1:B:361:LYS:HE3	2.03	0.58
1:A:170:ILE:HD12	1:A:170:ILE:H	1.69	0.58
1:B:356:PRO:HB3	1:B:367:ILE:HD12	1.86	0.58
1:B:71:GLU:OE2	1:B:73:ARG:NH1	2.37	0.57
1:A:23:ILE:HD12	1:A:25:LEU:HD21	1.85	0.57
1:A:232:ILE:HG22	1:A:233:GLN:HG3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:394:LYS:HD2	5:B:2082:HOH:O	2.04	0.57
1:B:170:ILE:H	1:B:170:ILE:HD12	1.70	0.56
1:B:128:LYS:NZ	5:B:2041:HOH:O	2.37	0.56
1:B:235:GLU:H	1:B:235:GLU:CD	2.08	0.56
1:A:41:LEU:HD23	1:A:143:PRO:HB3	1.87	0.56
1:B:130:VAL:O	1:B:193:PHE:HB3	2.05	0.56
1:B:162:VAL:HG21	1:B:183:MET:CE	2.35	0.56
1:B:305:LYS:HG2	5:B:2067:HOH:O	2.06	0.56
1:B:55:THR:O	1:B:223:GLY:HA3	2.05	0.56
1:A:235:GLU:H	1:A:235:GLU:CD	2.09	0.55
1:A:71:GLU:HG3	1:A:80:PRO:HB3	1.88	0.55
1:A:101:TRP:CD2	1:B:310:LYS:HE3	2.41	0.55
1:B:48:THR:HG21	1:B:280:THR:CG2	2.36	0.55
1:A:41:LEU:CD2	1:A:143:PRO:HB3	2.37	0.55
1:B:71:GLU:HG2	1:B:81:THR:H	1.70	0.55
1:B:331:SER:CB	1:B:332:PRO:CD	2.85	0.54
1:A:288:ARG:HG2	1:A:288:ARG:NH1	2.22	0.54
1:B:307:LYS:NZ	1:B:307:LYS:HB3	2.22	0.54
1:B:240:PHE:CE2	1:B:250:VAL:HG22	2.42	0.54
1:A:20:TRP:HA	1:A:287:LEU:O	2.07	0.54
1:B:41:LEU:HD23	1:B:143:PRO:HB3	1.89	0.54
1:B:337:PHE:CE1	1:B:351:LEU:HD21	2.43	0.54
1:A:71:GLU:OE2	1:A:73:ARG:NH1	2.41	0.54
1:B:355:ASN:N	1:B:356:PRO:HD3	2.22	0.54
1:B:232:ILE:HG22	1:B:233:GLN:HG3	1.87	0.54
1:B:383:GLU:HB3	1:B:384:PRO:CD	2.32	0.54
1:A:170:ILE:HD12	1:A:170:ILE:N	2.22	0.54
1:A:331:SER:CB	1:A:332:PRO:HD2	2.38	0.54
1:A:127:GLY:HA3	1:A:213:PHE:CZ	2.43	0.54
1:A:356:PRO:HB3	1:A:367:ILE:HD12	1.90	0.54
1:A:183:MET:HG2	1:A:287:LEU:HD23	1.89	0.53
1:A:357:ILE:O	1:A:357:ILE:CG1	2.54	0.53
1:A:331:SER:CB	1:A:332:PRO:CD	2.85	0.53
1:B:20:TRP:HA	1:B:287:LEU:O	2.09	0.53
1:B:41:LEU:CD2	1:B:143:PRO:HB3	2.39	0.53
1:B:288:ARG:NH1	1:B:288:ARG:HG2	2.24	0.53
1:A:183:MET:HG2	1:A:287:LEU:CD2	2.37	0.53
1:A:330:GLY:N	5:A:2034:HOH:O	2.41	0.53
1:A:2:ARG:CZ	1:A:6:ILE:HD11	2.38	0.53
1:B:383:GLU:CB	1:B:384:PRO:HD3	2.31	0.53
1:B:170:ILE:N	1:B:170:ILE:HD12	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:ILE:O	1:B:320:ILE:HG12	2.08	0.52
1:A:71:GLU:HG3	1:A:80:PRO:CB	2.38	0.52
1:B:356:PRO:HB3	1:B:367:ILE:CD1	2.40	0.52
1:A:355:ASN:N	1:A:356:PRO:HD3	2.24	0.52
1:B:393:LYS:HG3	5:B:2068:HOH:O	2.09	0.52
1:A:196:MET:HB2	5:A:2014:HOH:O	2.09	0.51
1:B:183:MET:HG2	1:B:287:LEU:HD23	1.91	0.51
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.37	0.51
1:B:27:HIS:HD2	1:B:280:THR:HB	1.74	0.51
1:A:378:ILE:CG2	1:A:380:ILE:HD11	2.41	0.51
1:A:24:VAL:C	1:A:25:LEU:HD23	2.31	0.51
1:B:331:SER:CB	1:B:332:PRO:HD2	2.40	0.51
1:B:127:GLY:HA3	1:B:213:PHE:CZ	2.45	0.51
1:B:191:LEU:HD12	1:B:192:ASP:N	2.24	0.50
1:B:182:THR:HG22	1:B:288:ARG:HG3	1.93	0.50
1:A:191:LEU:CD1	1:A:192:ASP:H	2.23	0.50
1:A:173:ALA:O	1:A:180:THR:HG23	2.10	0.50
1:A:173:ALA:HB3	1:A:181:VAL:HG12	1.94	0.50
1:A:182:THR:HG22	1:A:288:ARG:HG3	1.94	0.50
1:B:83:ASN:HB2	5:B:2016:HOH:O	2.11	0.50
1:A:181:VAL:HA	5:A:2023:HOH:O	2.10	0.50
1:A:27:HIS:HD2	1:A:280:THR:HB	1.76	0.50
1:A:61:ILE:CD1	1:A:123:LYS:HG2	2.39	0.50
1:A:240:PHE:CE2	1:A:250:VAL:HG22	2.47	0.50
1:B:183:MET:HG2	1:B:287:LEU:CD2	2.41	0.50
1:B:173:ALA:O	1:B:180:THR:HG23	2.11	0.50
1:A:80:PRO:HB2	1:A:114:VAL:HG12	1.92	0.49
1:A:191:LEU:HD12	1:A:192:ASP:N	2.26	0.49
1:A:48:THR:CB	1:A:280:THR:HG21	2.42	0.49
1:A:357:ILE:O	1:A:357:ILE:HG12	2.12	0.49
1:B:2:ARG:CZ	1:B:6:ILE:HD11	2.43	0.49
1:A:302:CYS:HB3	1:A:326:TYR:CZ	2.48	0.49
1:B:125:MET:HB3	1:B:201:MET:HG2	1.94	0.48
1:A:346:HIS:HA	5:A:2036:HOH:O	2.12	0.48
1:B:191:LEU:CD1	1:B:192:ASP:H	2.22	0.48
1:A:100:GLY:H	1:A:103:ASN:ND2	2.12	0.48
1:B:24:VAL:O	1:B:25:LEU:HD23	2.13	0.48
1:A:393:LYS:HG2	1:A:394:LYS:N	2.28	0.48
1:A:70:THR:HA	1:A:82:LEU:HD11	1.94	0.48
1:A:307:LYS:HZ3	1:A:307:LYS:HB3	1.79	0.48
1:A:323:ARG:HB2	1:A:366:ASN:ND2	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:ASN:OD1	1:A:155:THR:HB	2.14	0.48
1:B:100:GLY:H	1:B:103:ASN:ND2	2.12	0.48
1:B:173:ALA:HB3	1:B:181:VAL:HG12	1.96	0.48
1:B:24:VAL:C	1:B:25:LEU:HD23	2.34	0.48
1:A:48:THR:HG21	1:A:280:THR:HG23	1.95	0.47
1:A:383:GLU:CB	1:A:384:PRO:HD3	2.35	0.47
1:B:2:ARG:HD2	1:B:44:GLU:OE1	2.14	0.47
1:A:26:GLU:O	1:A:29:SER:HB3	2.14	0.47
1:B:71:GLU:HG3	1:B:80:PRO:HB3	1.96	0.47
1:B:80:PRO:HB2	1:B:114:VAL:HG12	1.95	0.47
1:B:46:ILE:HG12	1:B:140:VAL:CG2	2.37	0.47
1:B:202:LYS:HA	1:B:202:LYS:HE2	1.97	0.47
1:A:221:LEU:CD1	1:A:228:GLY:HA2	2.44	0.47
1:A:46:ILE:HG12	1:A:140:VAL:CG2	2.41	0.47
1:B:153:ASN:OD1	1:B:155:THR:HB	2.15	0.47
1:B:280:THR:CG2	1:B:280:THR:O	2.61	0.47
1:A:61:ILE:HD11	1:A:123:LYS:CG	2.40	0.47
1:A:337:PHE:CE1	1:A:351:LEU:HD21	2.50	0.47
1:B:48:THR:CB	1:B:280:THR:HG21	2.45	0.46
1:B:360:GLU:O	1:B:361:LYS:C	2.54	0.46
1:A:202:LYS:HA	1:A:202:LYS:HE2	1.97	0.46
1:B:310:LYS:HB3	1:B:323:ARG:HB3	1.96	0.46
1:A:8:ASN:HD22	1:A:29:SER:HA	1.79	0.46
1:B:71:GLU:HG3	1:B:80:PRO:CB	2.46	0.46
1:B:98:ASP:O	1:B:99:ARG:HD2	2.15	0.46
1:A:382:VAL:HG12	1:A:383:GLU:N	2.30	0.46
1:B:26:GLU:O	1:B:29:SER:HB3	2.15	0.46
1:B:188:ARG:HD2	1:B:188:ARG:HA	1.78	0.46
1:B:145:SER:OG	1:B:147:GLU:HB2	2.16	0.46
1:B:211:GLN:HE21	1:B:211:GLN:CA	2.28	0.46
1:A:24:VAL:O	1:A:25:LEU:HD23	2.16	0.45
1:A:2:ARG:HD2	1:A:44:GLU:OE1	2.17	0.45
1:A:125:MET:HB3	1:A:201:MET:HG2	1.98	0.45
1:A:48:THR:HG21	1:A:280:THR:HG21	1.98	0.45
1:A:188:ARG:HD2	1:A:188:ARG:HA	1.81	0.45
1:A:207:LEU:HA	1:A:207:LEU:HD12	1.85	0.45
3:A:1396:NAG:H61	3:A:1397:NAG:N2	2.32	0.45
1:B:296:GLY:C	1:B:298:SER:H	2.18	0.45
1:B:70:THR:HA	1:B:82:LEU:HD11	1.98	0.45
1:B:61:ILE:CD1	1:B:123:LYS:HG2	2.43	0.45
1:B:305:LYS:HA	5:B:2066:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:393:LYS:HG2	1:B:394:LYS:N	2.32	0.44
1:A:98:ASP:O	1:A:99:ARG:HD2	2.16	0.44
1:A:377:TYR:HB3	1:A:379:ILE:CD1	2.47	0.44
1:B:73:ARG:HD3	5:B:2010:HOH:O	2.17	0.44
1:A:211:GLN:CA	1:A:211:GLN:HE21	2.29	0.44
1:A:306:PHE:CZ	1:A:335:ILE:HG12	2.52	0.44
1:A:360:GLU:O	1:A:361:LYS:C	2.56	0.44
1:A:145:SER:OG	1:A:147:GLU:HB2	2.18	0.44
1:A:296:GLY:C	1:A:298:SER:H	2.21	0.44
1:B:71:GLU:CG	1:B:81:THR:H	2.31	0.44
1:A:378:ILE:HG22	1:A:380:ILE:HD11	1.99	0.44
1:B:240:PHE:CE2	1:B:250:VAL:CG2	3.00	0.44
1:A:337:PHE:CE2	1:A:380:ILE:HD12	2.52	0.43
1:B:71:GLU:HG2	1:B:81:THR:O	2.18	0.43
1:B:323:ARG:HB2	1:B:366:ASN:ND2	2.33	0.43
1:A:8:ASN:ND2	1:A:29:SER:HA	2.33	0.43
1:B:302:CYS:HB3	1:B:326:TYR:CZ	2.53	0.43
1:A:280:THR:CG2	1:A:280:THR:O	2.64	0.43
1:B:101:TRP:N	1:B:101:TRP:CD1	2.86	0.43
1:B:48:THR:HG21	1:B:280:THR:HG23	1.99	0.43
1:A:52:GLN:HA	1:A:52:GLN:NE2	2.34	0.43
1:B:48:THR:HG21	1:B:280:THR:HG21	2.01	0.43
1:B:96:MET:O	1:B:242:ASN:ND2	2.48	0.43
1:A:34:MET:HG2	1:A:40:THR:OG1	2.20	0.42
1:A:125:MET:HE2	1:A:260:MET:HG3	2.02	0.42
1:B:379:ILE:HD13	1:B:379:ILE:N	2.33	0.42
1:A:155:THR:HG22	1:A:157:LYS:HB2	2.01	0.42
1:A:278:LEU:HD22	1:B:243:PRO:HG3	2.01	0.42
1:A:46:ILE:O	1:A:46:ILE:HG22	2.19	0.42
1:A:71:GLU:CG	1:A:81:THR:H	2.31	0.42
1:B:52:GLN:HE21	1:B:52:GLN:HA	1.80	0.42
1:B:155:THR:HG22	1:B:157:LYS:HB2	2.02	0.42
1:A:168:SER:O	1:A:169:SER:OG	2.33	0.42
1:A:38:LYS:NZ	1:A:292:LEU:O	2.43	0.42
1:A:21:VAL:HG22	1:A:287:LEU:HB2	2.02	0.42
1:B:41:LEU:HD11	1:B:292:LEU:HD11	2.01	0.42
1:A:93:LYS:HB3	1:A:240:PHE:CE2	2.55	0.42
1:A:84:GLU:HG2	1:A:90:PHE:CE2	2.55	0.42
1:A:363:SER:HA	5:A:2039:HOH:O	2.19	0.42
1:B:34:MET:HG2	1:B:40:THR:OG1	2.20	0.42
1:B:195:GLU:HA	1:B:195:GLU:OE1	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:GLY:N	1:A:103:ASN:HD21	2.16	0.42
1:B:305:LYS:HE2	1:B:385:GLY:HA3	2.02	0.41
1:A:342:LEU:HA	5:A:2043:HOH:O	2.20	0.41
1:A:270:ILE:HD12	4:A:1400:BOG:H7'1	2.01	0.41
1:A:59:TYR:CD1	1:A:218:LEU:HB2	2.54	0.41
1:A:79:GLU:HA	1:A:80:PRO:HD3	1.87	0.41
1:A:170:ILE:HG22	1:A:171:THR:N	2.35	0.41
1:B:155:THR:CG2	1:B:157:LYS:HB2	2.50	0.41
1:A:155:THR:CG2	1:A:157:LYS:HB2	2.51	0.41
1:B:103:ASN:C	1:B:103:ASN:HD22	2.23	0.41
1:B:79:GLU:HA	1:B:80:PRO:HD3	1.85	0.41
1:A:297:MET:HA	1:A:334:LYS:NZ	2.34	0.41
1:A:103:ASN:C	1:A:103:ASN:HD22	2.24	0.41
1:A:237:LEU:HD23	1:A:237:LEU:HA	1.88	0.41
1:B:168:SER:O	1:B:169:SER:OG	2.35	0.41
1:B:84:GLU:HG2	1:B:90:PHE:CE2	2.56	0.41
1:B:88:LYS:HB2	1:B:230:ASN:ND2	2.36	0.41
1:A:155:THR:HG22	1:A:157:LYS:N	2.32	0.41
1:B:221:LEU:CD1	1:B:228:GLY:HA2	2.50	0.41
1:A:162:VAL:HG21	1:A:183:MET:HE3	2.03	0.40
1:B:382:VAL:HG12	1:B:383:GLU:N	2.36	0.40
1:B:21:VAL:HG22	1:B:287:LEU:HB2	2.03	0.40
1:B:170:ILE:HG22	1:B:171:THR:N	2.36	0.40
1:A:376:SER:C	5:A:2043:HOH:O	2.59	0.40
1:B:207:LEU:HD12	1:B:207:LEU:HA	1.82	0.40
1:A:307:LYS:NZ	1:A:307:LYS:CB	2.85	0.40
1:A:119:PHE:CD1	1:A:119:PHE:C	2.95	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	392/394 (100%)	354 (90%)	32 (8%)	6 (2%)	13	17
1	B	392/394 (100%)	352 (90%)	33 (8%)	7 (2%)	11	13
All	All	784/788 (100%)	706 (90%)	65 (8%)	13 (2%)	11	14

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	383	GLU
1	B	383	GLU
1	A	331	SER
1	B	331	SER
1	A	166	PRO
1	A	188	ARG
1	B	166	PRO
1	B	188	ARG
1	A	332	PRO
1	B	332	PRO
1	B	280	THR
1	A	384	PRO
1	B	384	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	342/342 (100%)	322 (94%)	20 (6%)	25	39
1	B	342/342 (100%)	321 (94%)	21 (6%)	23	36
All	All	684/684 (100%)	643 (94%)	41 (6%)	24	37

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	30	CYS
1	A	46	ILE

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Mol	Chain	Res	Type
1	A	56	LEU
1	A	70	THR
1	A	103	ASN
1	A	131	GLN
1	A	166	PRO
1	A	182	THR
1	A	211	GLN
1	A	249	ASP
1	A	252	VAL
1	A	253	LEU
1	A	283	LEU
1	A	290	ASP
1	A	292	LEU
1	A	332	PRO
1	A	379	ILE
1	A	387	LEU
1	A	394	LYS
1	B	2	ARG
1	B	30	CYS
1	B	46	ILE
1	B	56	LEU
1	B	70	THR
1	B	103	ASN
1	B	129	ILE
1	B	131	GLN
1	B	166	PRO
1	B	182	THR
1	B	211	GLN
1	B	249	ASP
1	B	252	VAL
1	B	253	LEU
1	B	283	LEU
1	B	290	ASP
1	B	292	LEU
1	B	320	ILE
1	B	332	PRO
1	B	357	ILE
1	B	387	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	52	GLN
1	A	103	ASN
1	A	131	GLN
1	A	211	GLN
1	A	256	GLN
1	A	316	GLN
1	A	325	GLN
1	A	366	ASN
1	B	8	ASN
1	B	52	GLN
1	B	103	ASN
1	B	131	GLN
1	B	211	GLN
1	B	248	GLN
1	B	256	GLN
1	B	276	ASN
1	B	316	GLN
1	B	325	GLN
1	B	366	ASN

### 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 ⓘ

8 carbohydrates 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
3	NAG	A	1396	1,3	14,14,15	0.65	0	15,19,21	0.92	1 (6%)
3	NAG	A	1397	3	14,14,15	0.74	0	15,19,21	1.04	1 (6%)
3	FUL	A	1398	3	10,10,11	0.42	0	14,14,16	0.92	2 (14%)
3	BMA	A	1399	3	11,11,12	0.57	0	14,15,17	0.62	0
3	NAG	B	1396	1,3	14,14,15	0.68	0	15,19,21	1.01	1 (6%)
3	NAG	B	1397	3	14,14,15	0.85	0	15,19,21	1.04	1 (6%)
3	FUL	B	1398	3	10,10,11	0.57	0	14,14,16	0.89	1 (7%)
3	BMA	B	1399	3	11,11,12	0.49	0	14,15,17	0.59	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	NAG	A	1396	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1397	3	-	0/6/23/26	0/1/1/1
3	FUL	A	1398	3	-	0/0/17/20	0/1/1/1
3	BMA	A	1399	3	-	0/2/19/22	0/1/1/1
3	NAG	B	1396	1,3	-	0/6/23/26	0/1/1/1
3	NAG	B	1397	3	-	0/6/23/26	0/1/1/1
3	FUL	B	1398	3	-	0/0/17/20	0/1/1/1
3	BMA	B	1399	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1396	NAG	C2-N2-C7	-2.87	119.36	123.04
3	B	1397	NAG	C2-N2-C7	-2.33	120.05	123.04
3	A	1396	NAG	C2-N2-C7	-2.29	120.09	123.04
3	A	1397	NAG	C2-N2-C7	-2.24	120.17	123.04
3	A	1398	FUL	C2-C3-C4	-2.08	107.52	111.04
3	B	1398	FUL	C1-O5-C5	2.15	115.70	112.38
3	A	1398	FUL	C1-O5-C5	2.38	116.06	112.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1396	NAG	2	0
3	A	1397	NAG	3	0
3	A	1399	BMA	1	0
3	B	1396	NAG	1	0
3	B	1397	NAG	2	0
3	B	1399	BMA	1	0

## 5.6 Ligand geometry

4 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	NAG	A	1395	1	14,14,15	0.71	0	15,19,21	0.68	0
4	BOG	A	1400	-	20,20,20	1.74	1 (5%)	25,25,25	1.11	2 (8%)
2	NAG	B	1395	1	14,14,15	0.55	0	15,19,21	0.75	1 (6%)
4	BOG	B	1400	-	20,20,20	1.60	1 (5%)	25,25,25	1.73	4 (16%)

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	NAG	A	1395	1	-	0/6/23/26	0/1/1/1
4	BOG	A	1400	-	-	0/11/31/31	0/1/1/1
2	NAG	B	1395	1	-	0/6/23/26	0/1/1/1
4	BOG	B	1400	-	-	0/11/31/31	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1400	BOG	O1-C1'	6.45	1.61	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1400	BOG	O1-C1'	6.72	1.61	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1400	BOG	O1-C1-C2	-6.39	99.97	108.04
4	A	1400	BOG	C1'-O1-C1	-3.38	108.03	113.94
4	A	1400	BOG	O1-C1-C2	-2.90	104.37	108.04
4	B	1400	BOG	O1-C1'-C2'	-2.58	99.63	109.88
4	B	1400	BOG	C1'-O1-C1	-2.53	109.53	113.94
2	B	1395	NAG	C2-N2-C7	-2.18	120.24	123.04
4	B	1400	BOG	O5-C1-O1	3.59	118.69	110.05

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1400	BOG	1	0
2	B	1395	NAG	1	0
4	B	1400	BOG	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	394/394 (100%)	0.49	45 (11%) 7 7	37, 80, 140, 164	0
1	B	394/394 (100%)	-0.11	6 (1%) 76 75	25, 60, 102, 150	0
All	All	788/788 (100%)	0.19	51 (6%) 22 22	25, 70, 133, 164	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	14	GLY	10.6
1	A	188	ARG	5.5
1	A	166	PRO	5.5
1	B	188	ARG	5.2
1	A	191	LEU	4.7
1	A	15	VAL	4.6
1	B	187	PRO	4.5
1	A	167	GLN	4.5
1	A	345	ARG	4.2
1	A	135	LEU	4.1
1	A	287	LEU	4.1
1	A	342	LEU	3.8
1	A	50	ALA	3.7
1	A	273	SER	3.6
1	A	23	ILE	3.6
1	A	274	SER	3.5
1	A	162	VAL	3.4
1	B	383	GLU	3.2
1	A	130	VAL	3.2
1	A	164	ILE	3.2
1	A	25	LEU	2.9
1	A	185	CYS	2.9
1	A	344	LYS	2.9
1	A	207	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	173	ALA	2.8
1	A	137	TYR	2.8
1	A	286	ARG	2.7
1	B	193	PHE	2.7
1	A	292	LEU	2.7
1	A	172	GLU	2.6
1	A	183	MET	2.6
1	A	171	THR	2.6
1	A	284	LYS	2.6
1	A	289	MET	2.5
1	A	136	GLU	2.5
1	A	343	GLU	2.5
1	A	35	ALA	2.5
1	A	285	CYS	2.4
1	B	191	LEU	2.4
1	A	278	LEU	2.4
1	A	20	TRP	2.3
1	A	335	ILE	2.2
1	A	383	GLU	2.2
1	A	160	LYS	2.2
1	A	179	GLY	2.1
1	B	166	PRO	2.1
1	A	133	GLU	2.1
1	A	12	VAL	2.0
1	A	392	PHE	2.0
1	A	175	LEU	2.0
1	A	193	PHE	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	B	1396	14/15	0.93	0.14	-0.01	63,70,78,90	0
3	NAG	A	1396	14/15	0.89	0.14	-0.14	75,78,90,92	0
3	BMA	B	1399	11/12	0.85	0.13	-	121,129,132,134	0
3	BMA	A	1399	11/12	0.71	0.15	-	120,127,131,132	0
3	FUL	B	1398	10/11	0.93	0.12	-	64,73,76,80	0
3	NAG	A	1397	14/15	0.88	0.13	-	85,103,109,116	0
3	NAG	B	1397	14/15	0.86	0.12	-	73,100,111,120	0
3	FUL	A	1398	10/11	0.93	0.17	-	75,81,89,94	0

## 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	BOG	A	1400	20/20	0.71	0.68	4.78	119,124,153,153	0
4	BOG	B	1400	20/20	0.79	0.36	3.55	89,119,122,124	0
2	NAG	B	1395	14/15	0.89	0.13	-	57,81,87,91	0
2	NAG	A	1395	14/15	0.89	0.13	-	96,108,115,117	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.