



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:34 AM GMT

PDB ID : 2HR7  
Title : Insulin receptor (domains 1-3)  
Authors : Garrett, T.P.J.; Ward, C.W.  
Deposited on : 2006-07-19  
Resolution : 2.32 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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

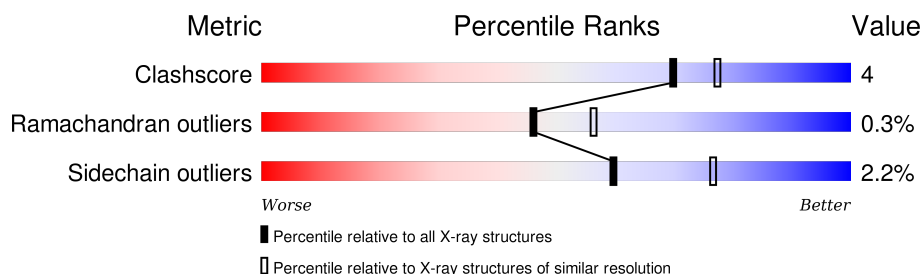
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.32 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	5057 (2.34-2.30)
Ramachandran outliers	100387	5008 (2.34-2.30)
Sidechain outliers	100360	5007 (2.34-2.30)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	486	
1	B	486	

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
2	NAG	A	501	X	-	-	-

## 2 Entry composition [i](#)

There are 12 unique types of molecules in this entry. The entry contains 8665 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.

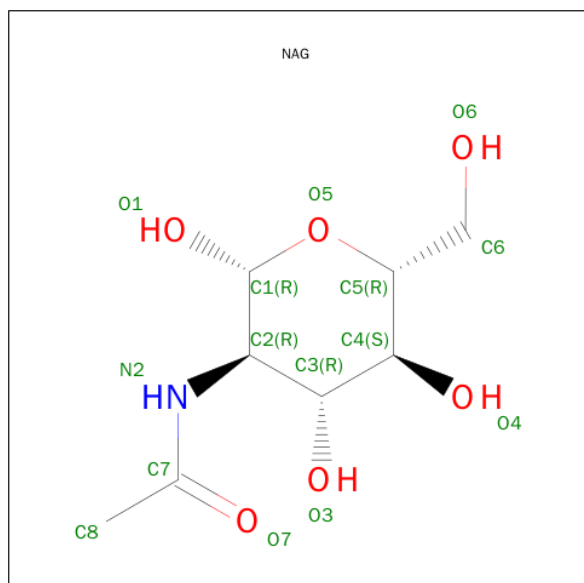
- Molecule 1 is a protein called Insulin receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	3	12	1
			3747	2348	670	687	42			
1	B	465	Total	C	N	O	S	0	22	1
			3770	2365	669	693	43			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	144	HIS	TYR	VARIANT	UNP P06213
A	421	ILE	THR	VARIANT	UNP P06213
A	486	SER	-	CLONING ARTIFACT	UNP P06213
B	144	HIS	TYR	VARIANT	UNP P06213
B	421	ILE	THR	VARIANT	UNP P06213
B	486	SER	-	CLONING ARTIFACT	UNP P06213

- 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	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		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	6	Total	C	N	O	0	0
			72	40	2	30		
4	B	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 5 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 6 is a polymer of unknown type called SUGAR (4-MER).

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

- Molecule 7 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	3	Total	C	N	O	0	0
			38	22	2	14		

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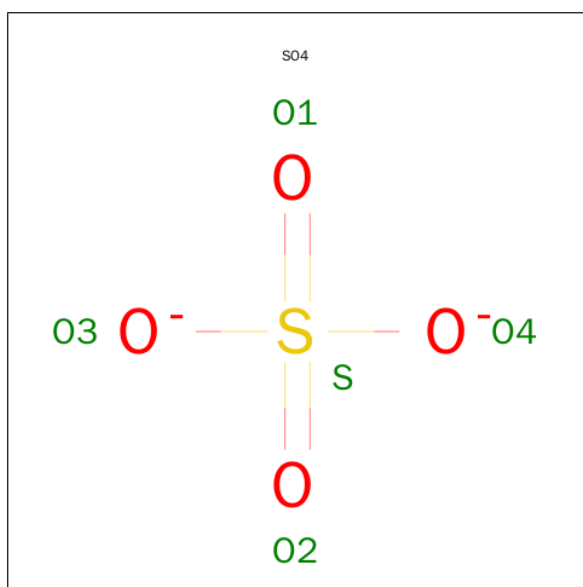
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	3	Total	C	N	O	0	0
			38	22	2	14		
7	B	3	Total	C	N	O	0	0
			38	22	2	14		
7	B	3	Total	C	N	O	0	0
			38	22	2	14		

- Molecule 8 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 9 is SULFATE ION (three-letter code: SO<sub>4</sub>) (formula: O<sub>4</sub>S).



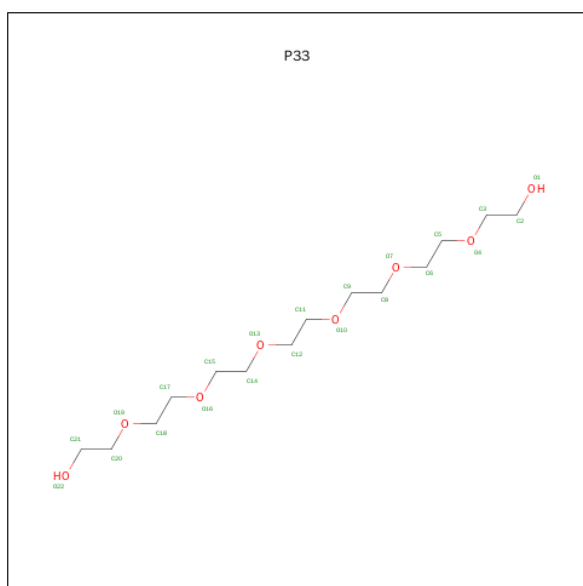
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	O	S	0	0
			5	4	1		
9	A	1	Total	O	S	0	0
			5	4	1		
9	A	1	Total	O	S	0	0
			5	4	1		
9	A	1	Total	O	S	0	0
			5	4	1		
9	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	O	S	0	0
			5	4	1		
9	B	1	Total	O	S	0	0
			5	4	1		
9	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 10 is 3,6,9,12,15,18-HEXAOXAIICOSANE-1,20-DIOL (three-letter code: P33) (formula:  $C_{14}H_{30}O_8$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			22	14	8		

- Molecule 11 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

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

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	192	Total	O	0	0
			192	192		
12	B	179	Total	O	0	1
			180	180		

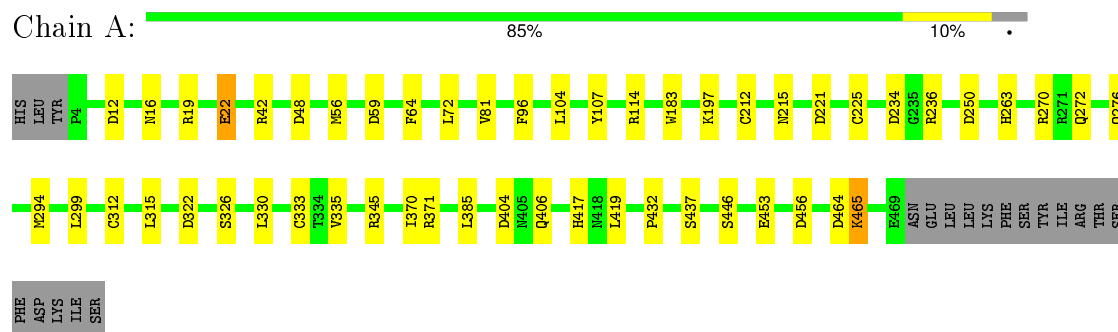


### 3 Residue-property plots

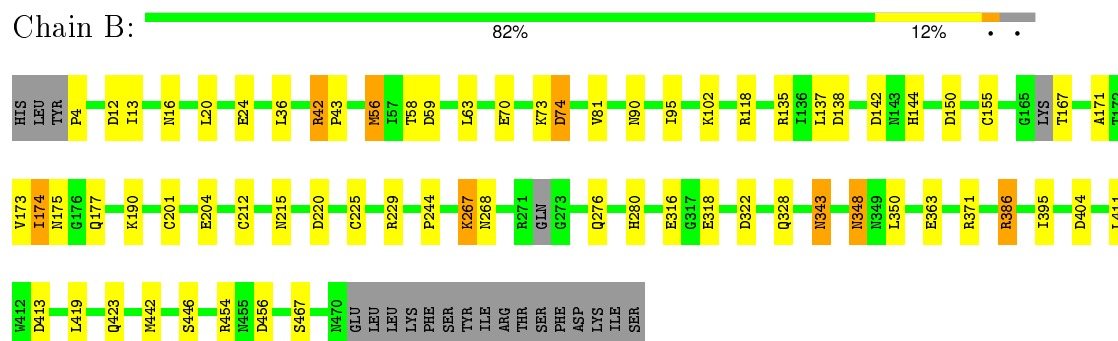
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

- Molecule 1: Insulin receptor



- Molecule 1: Insulin receptor



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.86Å 130.24Å 160.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.32	Depositor
% Data completeness (in resolution range)	100.0 (40.00-2.32)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC 5.2.0003	Depositor
R, $R_{free}$	0.195 , 0.231	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8665	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, BMA, NAG, FUC, P33, MAN, SO4

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	0.92	1/3875 (0.0%)	0.89	14/5236 (0.3%)
1	B	0.88	2/3913 (0.1%)	0.94	17/5288 (0.3%)
All	All	0.90	3/7788 (0.0%)	0.92	31/10524 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	270	ARG	CD-NE	-28.27	0.98	1.46
1	B	42	ARG	NE-CZ	16.32	1.54	1.33
1	B	42	ARG	CD-NE	6.33	1.57	1.46

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	42	ARG	CD-NE-CZ	-7.82	112.65	123.60
1	B	142	ASP	CB-CG-OD2	7.78	125.30	118.30
1	B	150	ASP	CB-CG-OD2	7.62	125.16	118.30
1	B	322	ASP	CB-CG-OD2	7.48	125.03	118.30
1	B	12	ASP	CB-CG-OD2	7.34	124.91	118.30
1	A	404	ASP	CB-CG-OD2	7.04	124.64	118.30
1	B	59	ASP	CB-CG-OD2	6.66	124.30	118.30
1	B	404	ASP	CB-CG-OD2	6.53	124.17	118.30
1	B	138	ASP	CB-CG-OD2	6.48	124.14	118.30
1	A	250	ASP	CB-CG-OD2	6.46	124.12	118.30
1	B	220	ASP	CB-CG-OD2	6.43	124.09	118.30
1	A	48	ASP	CB-CG-OD2	6.36	124.02	118.30
1	B	56	MET	CG-SD-CE	-6.33	90.07	100.20
1	A	322	ASP	CB-CG-OD2	6.12	123.81	118.30
1	A	12	ASP	CB-CG-OD1	-6.05	112.85	118.30
1	B	4	PRO	N-CA-CB	6.03	110.54	103.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	ASP	CB-CG-OD2	6.01	123.71	118.30
1	B	456	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	456	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	413	ASP	CB-CG-OD1	5.78	123.50	118.30
1	B	137	LEU	CA-CB-CG	5.67	128.34	115.30
1	A	16	ASN	CB-CA-C	-5.67	99.07	110.40
1	B	137	LEU	CB-CA-C	-5.53	99.69	110.20
1	A	221	ASP	CB-CG-OD1	5.38	123.14	118.30
1	A	59	ASP	CB-CG-OD2	5.36	123.12	118.30
1	A	464	ASP	CB-CG-OD2	5.36	123.12	118.30
1	B	454	ARG	NE-CZ-NH1	5.17	122.88	120.30
1	B	74	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	42[A]	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	A	42[B]	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	A	236	ARG	NE-CZ-NH2	-5.00	117.80	120.30

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	3747	0	3655	24	0
1	B	3770	0	3640	46	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
3	A	28	0	25	0	0
4	A	72	0	61	0	0
4	B	72	0	61	0	0
5	A	39	0	34	0	0
6	A	49	0	43	2	0
6	B	49	0	43	0	0
7	A	76	0	68	0	0
7	B	76	0	68	3	0
8	B	83	0	70	2	0
9	A	30	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	10	0	0	0	0
10	A	22	0	29	0	0
11	A	78	0	104	3	0
11	B	36	0	48	4	0
12	A	192	0	0	3	0
12	B	180	0	0	8	0
All	All	8665	0	8001	72	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 4.

All (72) 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:411:LEU:HD13	1:B:442:MET:CE	1.73	1.17
1:B:348[B]:ASN:HD21	1:B:350[B]:LEU:HD23	1.07	1.16
1:B:411:LEU:HD13	1:B:442:MET:HE3	1.35	1.05
1:B:348[B]:ASN:ND2	1:B:350[B]:LEU:HD23	1.76	0.99
6:A:553:NAG:O3	6:A:554:BMA:H2	1.71	0.89
1:B:386[B]:ARG:HD3	12:B:825[B]:HOH:O	1.78	0.83
1:B:316:GLU:OE2	12:B:849:HOH:O	2.05	0.73
1:A:453:GLU:O	12:A:781:HOH:O	2.09	0.70
1:B:386[B]:ARG:CD	12:B:825[B]:HOH:O	2.39	0.67
1:B:411:LEU:HD13	1:B:442:MET:HE2	1.76	0.66
1:A:345:ARG:NH2	9:A:602:SO4:O4	2.30	0.65
1:B:280:HIS:ND1	12:B:906:HOH:O	2.29	0.64
1:B:190:LYS:HD3	8:B:541:NAG:H82	1.82	0.62
1:A:56:MET:HG3	1:A:81:VAL:HB	1.84	0.60
1:A:263:HIS:CE1	1:A:276:GLN:HB3	2.36	0.60
1:A:312:CYS:SG	1:A:333[B]:CYS:HB2	2.41	0.59
1:B:56:MET:HG3	1:B:81:VAL:HB	1.83	0.59
1:B:386[A]:ARG:CD	7:B:592:FUC:H62	2.33	0.59
1:B:90:ASN:HB3	11:B:586:GOL:H12	1.87	0.56
1:B:174:ILE:HG13	1:B:175:ASN:H	1.71	0.54
1:A:114:ARG:HE	11:A:584:GOL:H2	1.73	0.54
1:B:190:LYS:CD	8:B:541:NAG:H82	2.37	0.53
1:A:333[B]:CYS:SG	1:A:335:VAL:O	2.67	0.53
1:B:411:LEU:CD1	1:B:442:MET:CE	2.67	0.53
6:A:553:NAG:C3	6:A:554:BMA:H2	2.40	0.52
1:B:371:ARG:HH12	11:B:588:GOL:C3	2.24	0.50
1:B:70:GLU:CD	1:B:102:LYS:HZ2	2.16	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:LYS:HB3	1:B:276:GLN:HG2	1.93	0.49
1:B:419:LEU:O	1:B:446:SER:HA	2.13	0.49
1:A:212[B]:CYS:SG	1:A:215:ASN:O	2.71	0.48
1:A:370:ILE:N	1:A:370:ILE:HD12	2.28	0.48
11:A:612:GOL:H12	12:A:777:HOH:O	2.13	0.47
1:A:315:LEU:HD21	1:B:167:THR:HG21	1.97	0.47
1:B:13:ILE:HD13	1:B:20:LEU:HA	1.96	0.47
1:B:348[B]:ASN:HD21	1:B:350[B]:LEU:CD2	1.99	0.47
1:B:411:LEU:HD22	1:B:442:MET:HE1	1.97	0.46
1:A:465:LYS:HA	1:A:465:LYS:HE3	1.98	0.46
1:B:328:GLN:NE2	12:B:904:HOH:O	2.49	0.46
1:B:386[A]:ARG:HD2	7:B:592:FUC:H62	1.97	0.46
1:B:423:GLN:NE2	12:B:686:HOH:O	2.48	0.46
1:B:386[A]:ARG:HD3	7:B:592:FUC:H62	1.98	0.45
1:B:90:ASN:HB3	11:B:586:GOL:C1	2.47	0.45
11:A:612:GOL:C1	12:A:777:HOH:O	2.64	0.45
1:B:36:LEU:N	1:B:36:LEU:HD23	2.32	0.45
1:A:212[B]:CYS:HB2	1:A:225:CYS:SG	2.57	0.45
1:A:385:LEU:O	1:A:417:HIS:NE2	2.41	0.45
1:B:56:MET:CE	1:B:58:THR:HG22	2.47	0.44
1:A:330:LEU:O	1:A:333[A]:CYS:SG	2.76	0.44
1:B:371:ARG:HH12	11:B:588:GOL:H31	1.81	0.44
1:B:63:LEU:HD12	1:B:95:ILE:CD1	2.48	0.43
1:A:197:LYS:HZ2	1:A:197:LYS:HG2	1.76	0.43
1:B:118:ARG:HD2	1:B:144:HIS:HB3	1.99	0.43
1:B:244:PRO:O	12:B:661:HOH:O	2.20	0.43
1:B:73:LYS:NZ	1:B:74:ASP:OD2	2.46	0.43
1:A:263:HIS:ND1	1:A:276:GLN:HB3	2.33	0.42
1:B:212[B]:CYS:HB2	1:B:225:CYS:SG	2.60	0.42
1:A:419:LEU:O	1:A:446:SER:HA	2.20	0.42
1:B:395:ILE:HG13	1:B:395:ILE:O	2.20	0.42
1:A:315:LEU:HD22	1:B:171:ALA:HB2	2.02	0.41
1:A:72:LEU:HB2	1:A:104:LEU:O	2.20	0.41
1:B:190:LYS:NZ	1:B:201:CYS:O	2.40	0.41
1:B:363:GLU:OE2	12:B:962:HOH:O	2.22	0.41
1:B:174:ILE:HG13	1:B:175:ASN:N	2.36	0.41
1:B:42:ARG:HB3	1:B:43:PRO:HD2	2.02	0.41
1:A:19:ARG:O	1:A:22:GLU:HB2	2.21	0.41
1:A:406:GLN:OE1	1:A:432:PRO:HG3	2.21	0.41
1:A:64:PHE:HA	1:A:96:PHE:O	2.21	0.40
1:A:294:MET:SD	1:A:299:LEU:HA	2.62	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212[B]:CYS:SG	1:B:215:ASN:O	2.79	0.40
1:A:107:TYR:HA	1:A:183:TRP:CD1	2.57	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	476/486 (98%)	455 (96%)	20 (4%)	1 (0%)	52	64
1	B	481/486 (99%)	458 (95%)	21 (4%)	2 (0%)	39	48
All	All	957/972 (98%)	913 (95%)	41 (4%)	3 (0%)	46	56

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	268	ASN
1	A	272	GLN
1	B	174	ILE

### 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	431/441 (98%)	426 (99%)	5 (1%)	78	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	431/441 (98%)	414 (96%)	17 (4%)	39	53
All	All	862/882 (98%)	840 (97%)	22 (3%)	60	71

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	326	SER
1	A	371	ARG
1	A	437	SER
1	A	465	LYS
1	B	16	ASN
1	B	135	ARG
1	B	155	CYS
1	B	173	VAL
1	B	177	GLN
1	B	204[A]	GLU
1	B	204[B]	GLU
1	B	229	ARG
1	B	267	LYS
1	B	318	GLU
1	B	343[A]	ASN
1	B	343[B]	ASN
1	B	348[A]	ASN
1	B	348[B]	ASN
1	B	386[A]	ARG
1	B	386[B]	ARG
1	B	467	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	281	ASN
1	A	328	GLN
1	B	34	GLN

### 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 ⓘ

44 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	511	1,3	14,14,15	0.55	0	15,19,21	1.33	2 (13%)
3	NAG	A	512	3	14,14,15	0.46	0	15,19,21	0.85	0
4	NAG	A	531	1,4	14,14,15	0.76	0	15,19,21	1.73	2 (13%)
4	NAG	A	532	4	14,14,15	0.72	0	15,19,21	1.47	5 (33%)
4	BMA	A	533	4	11,11,12	0.73	0	14,15,17	0.97	0
4	MAN	A	534	4	11,11,12	0.65	0	14,15,17	1.43	2 (14%)
4	MAN	A	535	4	11,11,12	0.59	0	14,15,17	1.25	1 (7%)
4	MAN	A	536	4	11,11,12	0.78	0	14,15,17	1.52	2 (14%)
5	NAG	A	541	1,5	14,14,15	0.90	1 (7%)	15,19,21	1.91	4 (26%)
5	NAG	A	542	5	14,14,15	0.76	1 (7%)	15,19,21	1.08	1 (6%)
5	BMA	A	543	5	11,11,12	0.76	0	14,15,17	1.55	2 (14%)
6	NAG	A	551	1,6	14,14,15	0.80	1 (7%)	15,19,21	1.20	1 (6%)
6	FUC	A	552	6	10,10,11	0.87	0	14,14,16	1.44	3 (21%)
6	NAG	A	553	6	14,14,15	0.93	1 (7%)	15,19,21	2.44	6 (40%)
6	BMA	A	554	6	11,11,12	0.66	0	14,15,17	2.19	4 (28%)
7	NAG	A	571	1,7	14,14,15	0.53	0	15,19,21	1.43	3 (20%)
7	FUC	A	572	7	10,10,11	0.60	0	14,14,16	1.06	1 (7%)
7	NAG	A	573	7	14,14,15	0.62	0	15,19,21	0.72	0
7	NAG	A	591	1,7	14,14,15	0.55	0	15,19,21	0.95	2 (13%)
7	FUC	A	592	7	10,10,11	0.70	0	14,14,16	0.99	0
7	NAG	A	593	7	14,14,15	0.55	0	15,19,21	0.76	0
4	NAG	B	531	1,4	14,14,15	0.81	0	15,19,21	1.93	3 (20%)
4	NAG	B	532	4	14,14,15	0.65	0	15,19,21	1.52	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	BMA	B	533	4	11,11,12	0.65	0	14,15,17	1.27	2 (14%)
4	MAN	B	534	4	11,11,12	0.65	0	14,15,17	1.32	2 (14%)
4	MAN	B	535	4	11,11,12	0.43	0	14,15,17	1.42	1 (7%)
4	MAN	B	536	4	11,11,12	0.54	0	14,15,17	0.83	0
8	NAG	B	541	1,8	14,14,15	1.03	1 (7%)	15,19,21	1.29	1 (6%)
8	NAG	B	542	8	14,14,15	0.77	1 (7%)	15,19,21	1.26	2 (13%)
8	BMA	B	543	8	11,11,12	0.72	0	14,15,17	1.03	1 (7%)
8	MAN	B	544	8	11,11,12	0.62	0	14,15,17	1.42	1 (7%)
8	MAN	B	547	8	11,11,12	0.79	0	14,15,17	1.63	3 (21%)
8	MAN	B	548	8	11,11,12	0.64	0	14,15,17	0.81	0
8	MAN	B	549	8	11,11,12	0.66	0	14,15,17	1.38	1 (7%)
6	NAG	B	551	1,6	14,14,15	0.72	0	15,19,21	1.20	1 (6%)
6	FUC	B	552	6	10,10,11	0.52	0	14,14,16	1.28	1 (7%)
6	NAG	B	553	6	14,14,15	0.50	0	15,19,21	1.03	0
6	BMA	B	554	6	11,11,12	0.80	0	14,15,17	1.30	2 (14%)
7	NAG	B	571	1,7	14,14,15	0.80	1 (7%)	15,19,21	1.01	0
7	FUC	B	572	7	10,10,11	1.10	1 (10%)	14,14,16	2.08	7 (50%)
7	NAG	B	573	7	14,14,15	0.56	0	15,19,21	1.07	1 (6%)
7	NAG	B	591	1,7	14,14,15	0.55	0	15,19,21	0.89	1 (6%)
7	FUC	B	592	7	10,10,11	0.91	0	14,14,16	1.47	4 (28%)
7	NAG	B	593	7	14,14,15	0.68	0	15,19,21	1.57	3 (20%)

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	511	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	512	3	-	0/6/23/26	0/1/1/1
4	NAG	A	531	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	532	4	-	0/6/23/26	0/1/1/1
4	BMA	A	533	4	-	0/2/19/22	0/1/1/1
4	MAN	A	534	4	-	0/2/19/22	0/1/1/1
4	MAN	A	535	4	-	0/2/19/22	0/1/1/1
4	MAN	A	536	4	-	0/2/19/22	0/1/1/1
5	NAG	A	541	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	542	5	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BMA	A	543	5	-	0/2/19/22	0/1/1/1
6	NAG	A	551	1,6	-	0/6/23/26	0/1/1/1
6	FUC	A	552	6	-	0/0/17/20	0/1/1/1
6	NAG	A	553	6	-	0/6/23/26	0/1/1/1
6	BMA	A	554	6	-	0/2/19/22	0/1/1/1
7	NAG	A	571	1,7	-	1/6/23/26	0/1/1/1
7	FUC	A	572	7	-	0/0/17/20	0/1/1/1
7	NAG	A	573	7	-	0/6/23/26	0/1/1/1
7	NAG	A	591	1,7	-	0/6/23/26	0/1/1/1
7	FUC	A	592	7	-	0/0/17/20	0/1/1/1
7	NAG	A	593	7	-	0/6/23/26	0/1/1/1
4	NAG	B	531	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	532	4	-	0/6/23/26	0/1/1/1
4	BMA	B	533	4	-	0/2/19/22	0/1/1/1
4	MAN	B	534	4	-	0/2/19/22	0/1/1/1
4	MAN	B	535	4	-	0/2/19/22	0/1/1/1
4	MAN	B	536	4	-	0/2/19/22	0/1/1/1
8	NAG	B	541	1,8	-	0/6/23/26	0/1/1/1
8	NAG	B	542	8	-	0/6/23/26	0/1/1/1
8	BMA	B	543	8	-	0/2/19/22	0/1/1/1
8	MAN	B	544	8	-	0/2/19/22	0/1/1/1
8	MAN	B	547	8	-	0/2/19/22	0/1/1/1
8	MAN	B	548	8	-	0/2/19/22	0/1/1/1
8	MAN	B	549	8	-	0/2/19/22	0/1/1/1
6	NAG	B	551	1,6	-	0/6/23/26	0/1/1/1
6	FUC	B	552	6	-	0/0/17/20	0/1/1/1
6	NAG	B	553	6	-	0/6/23/26	0/1/1/1
6	BMA	B	554	6	-	0/2/19/22	0/1/1/1
7	NAG	B	571	1,7	-	0/6/23/26	0/1/1/1
7	FUC	B	572	7	-	0/0/17/20	0/1/1/1
7	NAG	B	573	7	-	0/6/23/26	0/1/1/1
7	NAG	B	591	1,7	-	0/6/23/26	0/1/1/1
7	FUC	B	592	7	-	0/0/17/20	0/1/1/1
7	NAG	B	593	7	-	0/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	541	NAG	O5-C1	-2.45	1.39	1.43
6	A	551	NAG	O5-C5	-2.41	1.38	1.43
6	A	553	NAG	O5-C1	-2.25	1.40	1.43
8	B	542	NAG	O5-C1	-2.24	1.40	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	541	NAG	C2-N2	-2.16	1.42	1.46
5	A	542	NAG	C1-C2	2.02	1.55	1.52
7	B	572	FUC	C2-C3	2.04	1.55	1.52
7	B	571	NAG	C1-C2	2.41	1.55	1.52

All (80) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	531	NAG	C2-N2-C7	-3.97	117.94	123.04
5	A	541	NAG	C8-C7-N2	-3.89	108.67	116.11
7	B	572	FUC	O5-C1-C2	-3.78	104.72	110.86
4	B	532	NAG	C2-N2-C7	-3.71	118.27	123.04
8	B	541	NAG	C3-C2-N2	-3.66	101.80	110.56
4	B	532	NAG	C1-O5-C5	-3.27	108.09	112.25
8	B	542	NAG	C2-N2-C7	-3.21	118.91	123.04
6	A	553	NAG	C4-C3-C2	-3.19	106.27	111.23
6	B	551	NAG	C3-C2-N2	-2.92	103.56	110.56
6	A	551	NAG	C3-C2-N2	-2.86	103.70	110.56
4	A	534	MAN	C3-C4-C5	-2.85	105.24	110.20
7	B	572	FUC	C1-C2-C3	-2.80	106.23	109.54
4	A	532	NAG	C2-N2-C7	-2.72	119.55	123.04
6	A	553	NAG	O3-C3-C2	-2.69	103.78	109.11
4	B	533	BMA	O5-C1-C2	-2.64	106.58	110.86
7	B	593	NAG	C4-C3-C2	-2.57	107.23	111.23
4	A	532	NAG	O6-C6-C5	-2.56	102.89	111.33
6	A	553	NAG	C3-C4-C5	-2.47	105.89	110.20
4	B	531	NAG	O6-C6-C5	-2.43	103.30	111.33
7	A	571	NAG	C4-C3-C2	-2.37	107.54	111.23
5	A	542	NAG	C2-N2-C7	-2.35	120.02	123.04
3	A	511	NAG	O3-C3-C4	-2.28	105.19	110.34
6	A	553	NAG	O4-C4-C5	-2.28	103.19	109.24
3	A	511	NAG	C2-N2-C7	-2.24	120.16	123.04
8	B	547	MAN	O3-C3-C4	-2.24	105.29	110.34
7	B	591	NAG	C4-C3-C2	-2.22	107.78	111.23
4	A	532	NAG	O4-C4-C3	-2.21	105.37	110.34
4	A	532	NAG	C3-C4-C5	-2.18	106.39	110.20
8	B	543	BMA	O6-C6-C5	-2.08	104.46	111.33
5	A	541	NAG	O4-C4-C3	-2.06	105.69	110.34
6	A	554	BMA	O3-C3-C2	-2.05	106.30	110.00
7	A	591	NAG	C2-N2-C7	-2.02	120.44	123.04
4	A	532	NAG	C3-C2-N2	-2.00	105.77	110.56
4	B	534	MAN	O2-C2-C1	2.00	113.22	109.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	592	FUC	C2-C3-C4	2.00	114.44	111.04
4	B	533	BMA	C3-C4-C5	2.01	113.71	110.20
6	A	552	FUC	O5-C5-C6	2.03	109.49	106.13
4	A	534	MAN	O4-C4-C3	2.06	114.99	110.34
7	A	572	FUC	O5-C5-C6	2.07	109.55	106.13
6	A	554	BMA	O2-C2-C1	2.09	113.39	109.21
8	B	542	NAG	C1-O5-C5	2.09	114.90	112.25
7	A	571	NAG	O4-C4-C5	2.09	114.77	109.24
7	B	572	FUC	O2-C2-C3	2.14	114.42	110.12
8	B	547	MAN	O2-C2-C3	2.16	114.46	110.12
7	B	573	NAG	C2-N2-C7	2.17	125.82	123.04
7	A	591	NAG	C1-O5-C5	2.18	115.02	112.25
6	A	553	NAG	C2-N2-C7	2.19	125.85	123.04
7	B	592	FUC	C1-C2-C3	2.25	112.20	109.54
7	B	572	FUC	O2-C2-C1	2.27	113.77	109.21
7	B	592	FUC	O5-C5-C4	2.28	113.48	109.53
7	B	572	FUC	O5-C5-C6	2.29	109.91	106.13
7	B	572	FUC	O5-C5-C4	2.29	113.50	109.53
5	A	543	BMA	C1-O5-C5	2.45	115.36	112.25
6	B	554	BMA	O5-C5-C6	2.48	112.71	107.35
7	B	592	FUC	C3-C4-C5	2.60	114.10	109.72
4	A	535	MAN	C1-O5-C5	2.65	115.61	112.25
6	B	554	BMA	C1-C2-C3	2.66	112.69	109.54
4	A	536	MAN	C2-C3-C4	2.73	115.68	111.04
7	A	571	NAG	C1-O5-C5	2.78	115.78	112.25
4	A	531	NAG	O4-C4-C3	2.82	116.69	110.34
4	B	534	MAN	C1-O5-C5	3.15	116.24	112.25
6	A	552	FUC	C1-O5-C5	3.15	117.24	112.38
7	B	593	NAG	O3-C3-C2	3.24	115.54	109.11
4	A	536	MAN	C3-C4-C5	3.27	115.89	110.20
6	A	552	FUC	C1-C2-C3	3.28	113.43	109.54
5	A	541	NAG	O7-C7-N2	3.57	129.14	121.86
7	B	593	NAG	C1-O5-C5	3.64	116.87	112.25
6	A	554	BMA	C3-C4-C5	3.71	116.66	110.20
8	B	547	MAN	C2-C3-C4	3.71	117.35	111.04
4	B	531	NAG	O4-C4-C5	3.73	119.12	109.24
4	A	531	NAG	C1-O5-C5	3.74	117.00	112.25
7	B	572	FUC	C3-C4-C5	3.80	116.13	109.72
6	B	552	FUC	O5-C5-C6	3.85	112.49	106.13
5	A	543	BMA	C3-C4-C5	4.00	117.17	110.20
8	B	549	MAN	C1-O5-C5	4.23	117.61	112.25
5	A	541	NAG	C2-N2-C7	4.30	128.56	123.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	544	MAN	C1-C2-C3	4.48	114.83	109.54
4	B	535	MAN	C1-O5-C5	4.70	118.21	112.25
6	A	554	BMA	C1-O5-C5	6.17	120.08	112.25
6	A	553	NAG	O4-C4-C3	6.96	126.02	110.34

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	571	NAG	O7-C7-N2-C2

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	553	NAG	2	0
6	A	554	BMA	2	0
8	B	541	NAG	2	0
7	B	592	FUC	3	0

## 5.6 Ligand geometry [i](#)

32 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	501	1	14,14,15	0.70	0	15,19,21	1.20	2 (13%)
2	NAG	A	581	1	14,14,15	0.79	0	15,19,21	1.17	2 (13%)
11	GOL	A	582	-	5,5,5	0.44	0	5,5,5	0.86	0
11	GOL	A	583	-	5,5,5	0.41	0	5,5,5	0.49	0
11	GOL	A	584	-	5,5,5	0.27	0	5,5,5	0.38	0
11	GOL	A	585	-	5,5,5	0.31	0	5,5,5	0.63	0
11	GOL	A	587	-	5,5,5	0.29	0	5,5,5	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
11	GOL	A	594	-	5,5,5	0.34	0	5,5,5	0.69	0
11	GOL	A	595	-	5,5,5	0.48	0	5,5,5	0.55	0
11	GOL	A	596	-	5,5,5	0.39	0	5,5,5	0.33	0
11	GOL	A	597	-	5,5,5	0.87	0	5,5,5	1.01	0
11	GOL	A	599	-	5,5,5	0.47	0	5,5,5	0.33	0
9	SO4	A	601	-	4,4,4	0.54	0	6,6,6	0.48	0
9	SO4	A	602	-	4,4,4	0.77	0	6,6,6	1.29	1 (16%)
9	SO4	A	603	-	4,4,4	0.52	0	6,6,6	0.58	0
9	SO4	A	605	-	4,4,4	0.43	0	6,6,6	0.58	0
9	SO4	A	607	-	4,4,4	0.14	0	6,6,6	0.64	0
9	SO4	A	608	-	4,4,4	0.21	0	6,6,6	0.17	0
10	P33	A	609	-	21,21,21	1.02	1 (4%)	20,20,20	0.84	0
11	GOL	A	610	-	5,5,5	0.38	0	5,5,5	0.85	0
11	GOL	A	611	-	5,5,5	0.30	0	5,5,5	0.68	0
11	GOL	A	612	-	5,5,5	0.39	0	5,5,5	0.77	0
2	NAG	B	511	1	14,14,15	0.62	0	15,19,21	1.61	2 (13%)
11	GOL	B	580	-	5,5,5	0.48	0	5,5,5	0.37	0
2	NAG	B	581	1	14,14,15	0.45	0	15,19,21	1.48	2 (13%)
11	GOL	B	586	-	5,5,5	0.49	0	5,5,5	1.23	0
11	GOL	B	588	-	5,5,5	0.72	0	5,5,5	1.71	2 (40%)
11	GOL	B	589	-	5,5,5	0.30	0	5,5,5	0.64	0
11	GOL	B	598	-	5,5,5	0.67	0	5,5,5	0.77	0
9	SO4	B	604	-	4,4,4	0.36	0	6,6,6	0.89	1 (16%)
9	SO4	B	606	-	4,4,4	0.28	0	6,6,6	0.42	0
11	GOL	B	607	-	5,5,5	0.45	0	5,5,5	0.41	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	501	1	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	A	581	1	-	1/6/23/26	0/1/1/1
11	GOL	A	582	-	-	0/4/4/4	0/0/0/0
11	GOL	A	583	-	-	0/4/4/4	0/0/0/0
11	GOL	A	584	-	-	0/4/4/4	0/0/0/0
11	GOL	A	585	-	-	0/4/4/4	0/0/0/0
11	GOL	A	587	-	-	0/4/4/4	0/0/0/0
11	GOL	A	594	-	-	0/4/4/4	0/0/0/0
11	GOL	A	595	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	GOL	A	596	-	-	0/4/4/4	0/0/0/0
11	GOL	A	597	-	-	0/4/4/4	0/0/0/0
11	GOL	A	599	-	-	0/4/4/4	0/0/0/0
9	SO4	A	601	-	-	0/0/0/0	0/0/0/0
9	SO4	A	602	-	-	0/0/0/0	0/0/0/0
9	SO4	A	603	-	-	0/0/0/0	0/0/0/0
9	SO4	A	605	-	-	0/0/0/0	0/0/0/0
9	SO4	A	607	-	-	0/0/0/0	0/0/0/0
9	SO4	A	608	-	-	0/0/0/0	0/0/0/0
10	P33	A	609	-	-	0/19/19/19	0/0/0/0
11	GOL	A	610	-	-	0/4/4/4	0/0/0/0
11	GOL	A	611	-	-	0/4/4/4	0/0/0/0
11	GOL	A	612	-	-	0/4/4/4	0/0/0/0
2	NAG	B	511	1	-	0/6/23/26	0/1/1/1
11	GOL	B	580	-	-	0/4/4/4	0/0/0/0
2	NAG	B	581	1	-	0/6/23/26	0/1/1/1
11	GOL	B	586	-	-	0/4/4/4	0/0/0/0
11	GOL	B	588	-	-	0/4/4/4	0/0/0/0
11	GOL	B	589	-	-	0/4/4/4	0/0/0/0
11	GOL	B	598	-	-	0/4/4/4	0/0/0/0
9	SO4	B	604	-	-	0/0/0/0	0/0/0/0
9	SO4	B	606	-	-	0/0/0/0	0/0/0/0
11	GOL	B	607	-	-	0/4/4/4	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	A	609	P33	O22-C21	-3.36	1.23	1.42

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	581	NAG	C2-N2-C7	-3.18	118.96	123.04
2	B	511	NAG	C4-C3-C2	-2.96	106.62	111.23
9	A	602	SO4	O2-S-O1	-2.82	100.56	109.50
11	B	588	GOL	O2-C2-C3	-2.03	99.35	108.65
9	B	604	SO4	O2-S-O1	2.06	116.02	109.50
2	A	501	NAG	C3-C4-C5	2.31	114.22	110.20
11	B	588	GOL	O2-C2-C1	2.37	119.52	108.65
2	A	581	NAG	C2-N2-C7	2.49	126.24	123.04
2	A	581	NAG	C1-O5-C5	2.88	115.91	112.25
2	B	581	NAG	O5-C5-C6	3.07	114.00	107.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAG	C4-C3-C2	3.31	116.37	111.23
2	B	511	NAG	C1-O5-C5	4.09	117.44	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	501	NAG	C1

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	581	NAG	O7-C7-N2-C2

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	584	GOL	1	0
9	A	602	SO4	1	0
11	A	612	GOL	2	0
11	B	586	GOL	2	0
11	B	588	GOL	2	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](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.