



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:34 AM GMT

PDB ID : 2HRG  
Title : Crystal Structure of Blue Laccase from *Trametes trogii* complexed with p-methylbenzoate  
Authors : Matera, I.; Gullotto, A.; Ferraroni, M.; Tilli, S.; Briganti, F.; Scozzafava, A.  
Deposited on : 2006-07-20  
Resolution : 1.58 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

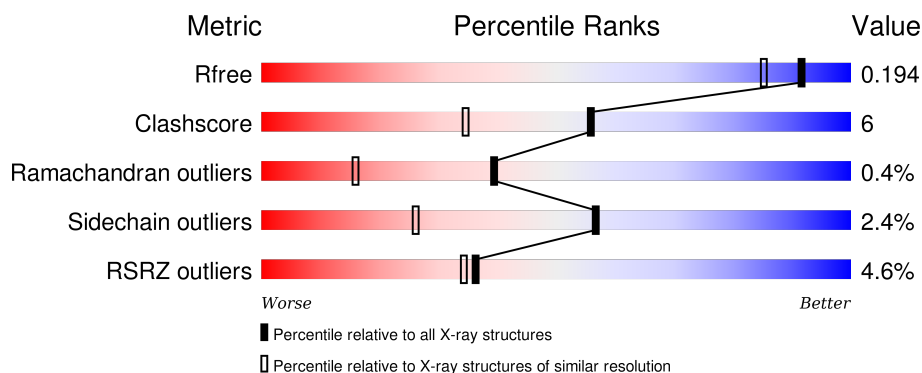
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.58 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	3815 (1.60-1.56)
Clashscore	102246	4131 (1.60-1.56)
Ramachandran outliers	100387	4021 (1.60-1.56)
Sidechain outliers	100360	4018 (1.60-1.56)
RSRZ outliers	91569	3823 (1.60-1.56)

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	496	<div> <div>5%</div> <div>91%</div> <div>8%</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
2	CBS	A	497	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CBS	A	1006	-	-	-	X
5	CA	A	1001	-	-	-	X
5	CA	A	1002	-	-	-	X
6	4MA	A	3001	-	-	-	X
7	GOL	A	2001	-	-	-	X
7	GOL	A	2002	-	-	-	X

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 4478 atoms, of which 2 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.

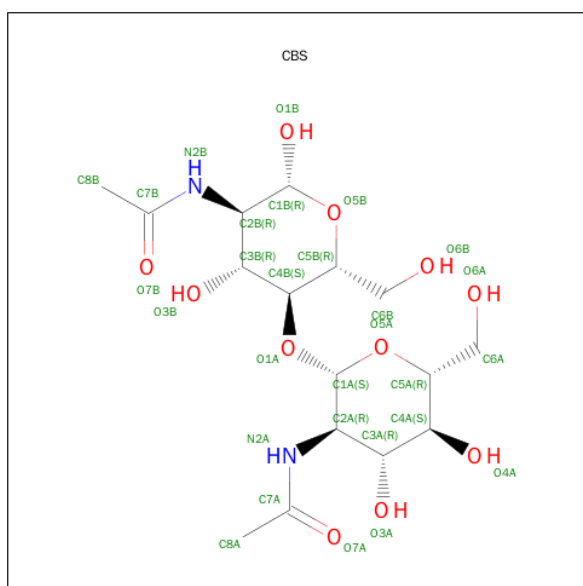
- Molecule 1 is a protein called Laccase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	496	Total	C	H	N	O	S	0	20	0
			3817	2425	2	641	738	11			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	3	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 3 is SUGAR (DI(N-ACETYL-D-GLUCOSAMINE)) (three-letter code: CBS) (formula: C<sub>16</sub>H<sub>28</sub>N<sub>2</sub>O<sub>11</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			28	16	2	10		

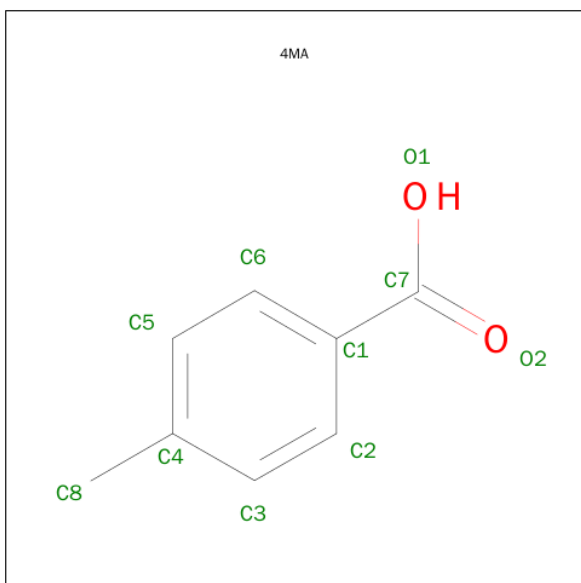
- Molecule 4 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	Cu	0	0
			4	4		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Ca	0	0
			3	3		

- Molecule 6 is 4-METHYLBENZOIC ACID (three-letter code: 4MA) (formula: C<sub>8</sub>H<sub>8</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	8	2		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

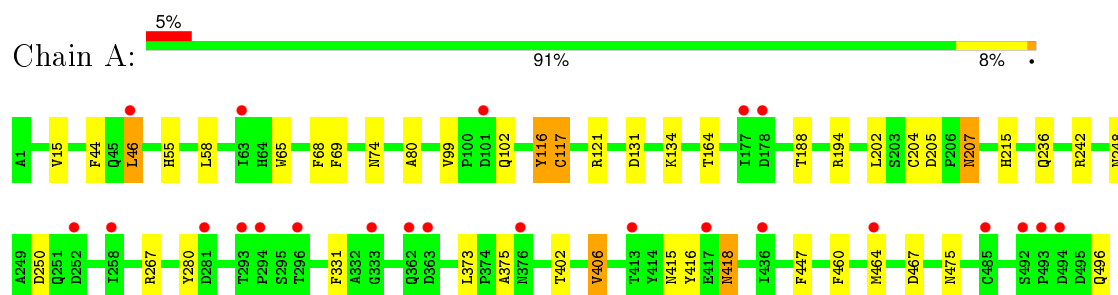
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	548	Total	O	0	0
			548	548		

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

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.

• Molecule 1: Laccase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.39Å 85.13Å 108.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.33 – 1.58 22.33 – 1.58	Depositor EDS
% Data completeness (in resolution range)	99.5 (22.33-1.58) 99.5 (22.33-1.58)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.73 (at 1.58Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.174 , 0.192 0.176 , 0.194	Depositor DCC
$R_{free}$ test set	5332 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.1	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.43 , 60.7	EDS
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 106932 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4478	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: GOL, BMA, 4MA, CA, CBS, CU, MAN

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	0.39	0/4020	0.58	0/5520

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
2	A	1	0

There are no bond length outliers.

There are no bond angle outliers.

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	497	CBS	C1B

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	3815	2	3630	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	50	0	44	1	0
3	A	28	0	26	1	0
4	A	4	0	0	0	0
5	A	3	0	0	0	0
6	A	10	0	7	1	0
7	A	18	0	24	2	0
8	A	548	0	0	5	0
All	All	4476	2	3731	48	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 6.

All (48) 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:373:LEU:HD11	1:A:464:MET:CE	1.66	1.25
1:A:373:LEU:HD21	1:A:464:MET:CE	1.81	1.09
1:A:373:LEU:HD21	1:A:464:MET:HE3	1.30	1.07
1:A:373:LEU:CD2	1:A:464:MET:HE3	1.85	1.07
1:A:373:LEU:HD11	1:A:464:MET:HE2	1.12	1.05
1:A:373:LEU:CG	1:A:464:MET:HE3	1.89	1.03
1:A:373:LEU:CD1	1:A:464:MET:CE	2.46	0.92
1:A:373:LEU:HD11	1:A:464:MET:HE3	1.63	0.81
1:A:250:ASP:HB3	8:A:3497:HOH:O	1.82	0.79
1:A:373:LEU:CD1	1:A:464:MET:HE3	2.09	0.78
1:A:68[B]:PHE:CE1	1:A:99[B]:VAL:HG21	2.20	0.77
1:A:373:LEU:HD21	1:A:464:MET:HE1	1.67	0.76
1:A:68[B]:PHE:HE1	1:A:99[B]:VAL:HG21	1.53	0.72
1:A:121[A]:ARG:NH1	1:A:204[A]:CYS:SG	2.64	0.71
1:A:116:TYR:CE2	1:A:204[B]:CYS:SG	2.89	0.66
1:A:215:HIS:HE1	1:A:280:TYR:OH	1.78	0.66
1:A:248:ASN:HB3	8:A:3497:HOH:O	1.96	0.65
1:A:116:TYR:HD2	1:A:121[A]:ARG:HD3	1.66	0.61
1:A:373:LEU:CD2	1:A:464:MET:CE	2.56	0.60
1:A:331:PHE:CZ	6:A:3001:4MA:H83	2.37	0.60
1:A:207:ASN:ND2	1:A:236:GLN:HE21	1.99	0.60
1:A:55:HIS:HD2	1:A:496[B]:GLN:O	1.87	0.57
1:A:80:ALA:HB3	7:A:2002:GOL:H32	1.86	0.57
1:A:373:LEU:HG	1:A:464:MET:HE3	1.83	0.56
1:A:69:PHE:H	1:A:102:GLN:HE22	1.53	0.54
1:A:194:ARG:HG2	1:A:248:ASN:ND2	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:PHE:HB3	1:A:46[B]:LEU:CD2	2.38	0.54
1:A:121[B]:ARG:HD2	1:A:202:LEU:HB3	1.90	0.53
1:A:267:ARG:NH2	8:A:3248:HOH:O	2.43	0.52
1:A:406[A]:VAL:HG21	1:A:416:TYR:CE2	2.45	0.51
1:A:74:ASN:H	1:A:475:ASN:ND2	2.10	0.50
1:A:68[B]:PHE:CD1	1:A:99[B]:VAL:HG22	2.47	0.50
3:A:1006:CBS:H1A	3:A:1006:CBS:H61B	1.94	0.49
1:A:68[B]:PHE:CE1	1:A:99[B]:VAL:CG2	2.93	0.49
1:A:117[A]:CYS:SG	8:A:3197:HOH:O	2.59	0.49
1:A:415:ASN:HD21	1:A:418:ASN:ND2	2.11	0.48
1:A:131:ASP:HB3	1:A:134:LYS:HB3	1.97	0.46
1:A:99[A]:VAL:CG2	1:A:102:GLN:HB2	2.46	0.46
1:A:68[B]:PHE:HD1	1:A:99[B]:VAL:HG22	1.83	0.44
1:A:46[A]:LEU:HD23	1:A:65:TRP:CZ3	2.53	0.44
1:A:68[B]:PHE:CD1	1:A:99[B]:VAL:CG2	3.01	0.44
1:A:74:ASN:H	1:A:475:ASN:HD22	1.65	0.44
1:A:375:ALA:HB2	1:A:467:ASP:CG	2.40	0.42
2:A:497:CBS:H61B	2:A:497:CBS:C7A	2.49	0.42
1:A:164:THR:HG23	8:A:3327:HOH:O	2.20	0.42
1:A:69:PHE:HE1	1:A:402[A]:THR:HG23	1.85	0.42
1:A:99[B]:VAL:CG1	1:A:102:GLN:HB2	2.50	0.41
1:A:15:VAL:HA	7:A:2001:GOL:H12	2.03	0.41

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	514/496 (104%)	502 (98%)	10 (2%)	2 (0%)	39 16

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	205	ASP
1	A	58	LEU

### 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	430/409 (105%)	416 (97%)	14 (3%)	45	16

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46[A]	LEU
1	A	46[B]	LEU
1	A	116	TYR
1	A	117[A]	CYS
1	A	117[B]	CYS
1	A	188[A]	THR
1	A	188[B]	THR
1	A	207	ASN
1	A	242	ARG
1	A	406[A]	VAL
1	A	406[B]	VAL
1	A	418	ASN
1	A	447	PHE
1	A	460	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	45	GLN
1	A	55	HIS
1	A	102	GLN
1	A	207	ASN
1	A	215	HIS

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Mol	Chain	Res	Type
1	A	248	ASN
1	A	362	GLN
1	A	418	ASN
1	A	475	ASN
1	A	479	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 ⓘ

3 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$
2	CBS	A	497	1,2	29,29,30	2.28	4 (13%)	36,41,43	3.71	25 (69%)
2	BMA	A	498	2	11,11,12	0.54	0	14,15,17	0.94	1 (7%)
2	MAN	A	499	2	11,11,12	0.56	0	14,15,17	0.93	2 (14%)

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	CBS	A	497	1,2	1/1/11/14	0/16/53/56	0/2/2/2
2	BMA	A	498	2	-	0/2/19/22	0/1/1/1
2	MAN	A	499	2	-	0/2/19/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	497	CBS	C7A-N2A	2.83	1.45	1.34
2	A	497	CBS	C7B-N2B	2.87	1.45	1.34
2	A	497	CBS	O7B-C7B	7.91	1.41	1.23
2	A	497	CBS	O7A-C7A	8.12	1.42	1.23

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	497	CBS	O7B-C7B-C8B	-6.86	109.48	122.06
2	A	497	CBS	O7A-C7A-C8A	-6.40	110.31	122.06
2	A	497	CBS	O7B-C7B-N2B	-5.31	111.04	121.86
2	A	497	CBS	O7A-C7A-N2A	-4.51	112.66	121.86
2	A	497	CBS	C2B-N2B-C7B	-4.46	117.31	123.04
2	A	497	CBS	C8B-C7B-N2B	-3.20	109.99	116.11
2	A	497	CBS	C1A-O1A-C4B	-3.04	110.07	118.01
2	A	497	CBS	O6A-C6A-C5A	-2.69	102.46	111.33
2	A	497	CBS	C8A-C7A-N2A	-2.27	111.76	116.11
2	A	497	CBS	C2A-N2A-C7A	-2.27	117.28	123.10
2	A	497	CBS	O5B-C5B-C6B	2.10	111.89	107.35
2	A	499	MAN	C1-C2-C3	2.14	112.08	109.54
2	A	497	CBS	O3A-C3A-C2A	2.18	114.10	109.66
2	A	499	MAN	C1-O5-C5	2.18	115.02	112.25
2	A	497	CBS	O3A-C3A-C4A	2.54	116.06	110.34
2	A	498	BMA	C1-O5-C5	2.65	115.61	112.25
2	A	497	CBS	C3B-C4B-C5B	2.86	117.31	110.84
2	A	497	CBS	C4B-C3B-C2B	3.02	117.34	109.74
2	A	497	CBS	O4A-C4A-C5A	3.08	117.39	109.24
2	A	497	CBS	O3B-C3B-C4B	3.52	118.20	109.87
2	A	497	CBS	C3A-C4A-C5A	3.99	117.16	110.20
2	A	497	CBS	O5A-C1A-C2A	4.21	119.94	110.78
2	A	497	CBS	O5B-C5B-C4B	4.25	117.75	110.01
2	A	497	CBS	C1A-O5A-C5A	4.61	122.69	113.75
2	A	497	CBS	O1A-C4B-C3B	4.67	119.23	107.17
2	A	497	CBS	O5A-C5A-C4A	4.77	118.63	109.68
2	A	497	CBS	C4A-C3A-C2A	5.29	117.76	110.43
2	A	497	CBS	C1B-O5B-C5B	9.27	124.02	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	497	CBS	C1B

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	497	CBS	1	0

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 7 are monoatomic - leaving 5 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	CBS	A	1006	1	29,29,30	2.40	4 (13%)	36,41,43	3.55	21 (58%)
7	GOL	A	2001	-	5,5,5	0.45	0	5,5,5	0.28	0
7	GOL	A	2002	-	5,5,5	0.33	0	5,5,5	0.37	0
7	GOL	A	2003	-	5,5,5	0.42	0	5,5,5	0.17	0
6	4MA	A	3001	-	7,10,10	1.50	1 (14%)	10,13,13	0.96	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	CBS	A	1006	1	-	0/16/53/56	0/2/2/2
7	GOL	A	2001	-	-	0/4/4/4	0/0/0/0
7	GOL	A	2002	-	-	0/4/4/4	0/0/0/0
7	GOL	A	2003	-	-	0/4/4/4	0/0/0/0
6	4MA	A	3001	-	-	0/0/4/4	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	3001	4MA	C8-C4	-3.75	1.37	1.51

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1006	CBS	C7B-N2B	3.02	1.45	1.34
3	A	1006	CBS	C7A-N2A	3.07	1.46	1.34
3	A	1006	CBS	O7A-C7A	8.30	1.42	1.23
3	A	1006	CBS	O7B-C7B	8.37	1.42	1.23

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1006	CBS	O7A-C7A-C8A	-6.94	109.33	122.06
3	A	1006	CBS	O7B-C7B-C8B	-6.19	110.70	122.06
3	A	1006	CBS	O7A-C7A-N2A	-5.11	111.44	121.86
3	A	1006	CBS	O7B-C7B-N2B	-5.06	111.55	121.86
3	A	1006	CBS	C1A-O1A-C4B	-4.56	106.08	118.01
3	A	1006	CBS	C8B-C7B-N2B	-2.63	111.07	116.11
3	A	1006	CBS	C8A-C7A-N2A	-2.37	111.58	116.11
3	A	1006	CBS	C2B-N2B-C7B	-2.21	120.20	123.04
3	A	1006	CBS	O3B-C3B-C4B	2.05	114.74	109.87
3	A	1006	CBS	O5B-C5B-C6B	2.23	112.17	107.35
3	A	1006	CBS	C3A-C2A-N2A	2.47	115.77	110.66
3	A	1006	CBS	O1A-C4B-C5B	2.65	116.29	109.32
3	A	1006	CBS	C4B-C3B-C2B	3.14	117.63	109.74
3	A	1006	CBS	O5B-C5B-C4B	3.55	116.47	110.01
3	A	1006	CBS	C3B-C4B-C5B	4.14	120.20	110.84
3	A	1006	CBS	O5A-C1A-C2A	4.33	120.19	110.78
3	A	1006	CBS	O5A-C5A-C4A	4.75	118.60	109.68
3	A	1006	CBS	C3A-C4A-C5A	5.05	119.00	110.20
3	A	1006	CBS	C1A-O5A-C5A	5.24	123.91	113.75
3	A	1006	CBS	C4A-C3A-C2A	5.72	118.36	110.43
3	A	1006	CBS	C1B-O5B-C5B	8.45	122.97	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1006	CBS	1	0
7	A	2001	GOL	1	0
7	A	2002	GOL	1	0
6	A	3001	4MA	1	0



## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	496/496 (100%)	0.12	23 (4%) 36 34	9, 17, 28, 34	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	333	GLY	4.7
1	A	362	GLN	4.7
1	A	294	PRO	4.5
1	A	293	THR	3.8
1	A	494	ASP	3.6
1	A	281	ASP	3.2
1	A	493	PRO	2.7
1	A	178	ASP	2.6
1	A	46[A]	LEU	2.5
1	A	101	ASP	2.4
1	A	258	ILE	2.4
1	A	252	ASP	2.4
1	A	376	ASN	2.3
1	A	177	ILE	2.3
1	A	63[A]	ILE	2.3
1	A	413	THR	2.2
1	A	492	SER	2.2
1	A	296	THR	2.2
1	A	363	ASP	2.2
1	A	485	CYS	2.2
1	A	436	ILE	2.1
1	A	417	GLU	2.1
1	A	464	MET	2.1

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

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

## 6.3 Carbohydrates ⓘ

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(Å <sup>2</sup> )	Q<0.9
2	CBS	A	497	28/29	0.89	0.09	0.71	15,17,22,25	0
2	MAN	A	499	11/12	0.75	0.23	-	40,41,42,44	0
2	BMA	A	498	11/12	0.78	0.18	-	30,34,36,38	0

## 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. 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(Å <sup>2</sup> )	Q<0.9
5	CA	A	1002	1/1	0.85	0.38	20.68	48,48,48,48	0
3	CBS	A	1006	28/29	0.75	0.20	8.89	25,28,34,37	0
7	GOL	A	2001	6/6	0.77	0.29	4.83	42,43,44,45	0
5	CA	A	1001	1/1	0.78	0.28	4.18	106,106,106,106	0
7	GOL	A	2002	6/6	0.75	0.15	4.03	30,36,37,38	0
6	4MA	A	3001	10/10	0.85	0.16	3.21	28,33,34,35	0
7	GOL	A	2003	6/6	0.84	0.13	0.55	34,36,37,39	0
4	CU	A	1500	1/1	1.00	0.02	-3.59	14,14,14,14	0
4	CU	A	1497	1/1	1.00	0.03	-3.83	14,14,14,14	0
4	CU	A	1498	1/1	1.00	0.02	-4.86	13,13,13,13	0
5	CA	A	1003	1/1	0.75	0.10	-	48,48,48,48	0
4	CU	A	1499	1/1	0.99	0.17	-	22,22,22,22	0

## 6.5 Other polymers

There are no such residues in this entry.