



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 01:12 PM GMT

PDB ID : 3T6V
Title : Crystal Structure of Laccase from *Steccherinum ochraceum*
Authors : Ferraroni, M.; Briganti, F.; Matera, I.; Kolomytseva, M.; Golovleva, L.; Scoz-
zafava, A.; Chernykh, A.M.
Deposited on : 2011-07-29
Resolution : 2.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at 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.

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

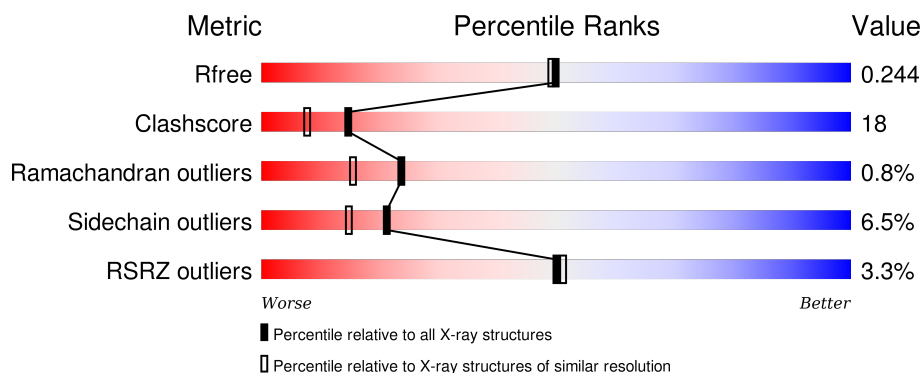
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 ≥ 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	495	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 75% 21% </div> </div>
1	B	495	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 75% 21% </div> </div>
1	C	495	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 75% 21% </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
3	SO4	A	505	-	-	-	X
3	SO4	A	506	-	-	-	X
3	SO4	A	507	-	-	-	X
3	SO4	C	505	-	-	X	-
3	SO4	C	506	-	-	X	X
4	GOL	A	509	-	-	X	X
4	GOL	B	506	-	-	X	X
4	GOL	B	508	-	-	-	X
4	GOL	B	509	-	-	-	X
4	GOL	C	507	-	-	X	X
4	GOL	C	508	-	-	-	X
5	CBS	B	511	-	-	-	X
5	CBS	C	509	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12647 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 Laccase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	495	Total	C	N	O	S	0	5	0
			3734	2358	625	731	20			
1	B	495	Total	C	N	O	S	0	6	0
			3739	2363	630	728	18			
1	C	495	Total	C	N	O	S	0	3	0
			3731	2356	629	729	17			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	4	Total	Cu	0	0
			4	4		
2	A	4	Total	Cu	0	0
			4	4		
2	C	4	Total	Cu	0	0
			4	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



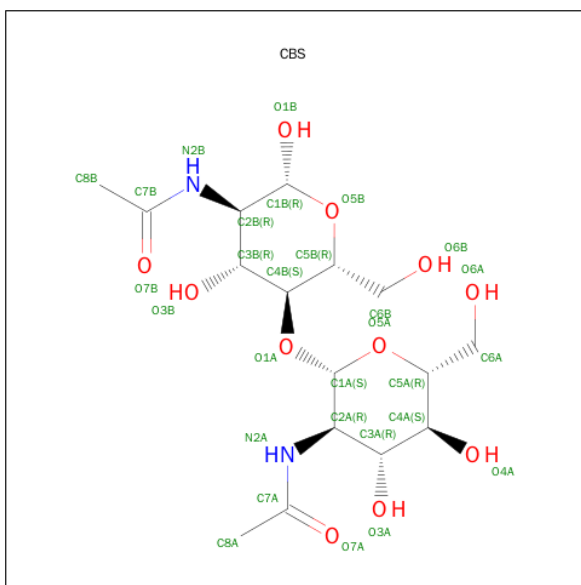
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 5 is SUGAR (DI(N-ACETYL-D-GLUCOSAMINE)) (three-letter code: CBS) (formula: C₁₆H₂₈N₂O₁₁).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			28	16	2	10		
5	A	1	Total	C	N	O	0	0
			28	16	2	10		
5	B	1	Total	C	N	O	0	0
			28	16	2	10		
5	B	1	Total	C	N	O	0	0
			28	16	2	10		
5	C	1	Total	C	N	O	0	0
			28	16	2	10		
5	C	1	Total	C	N	O	0	0
			14	8	1	5		

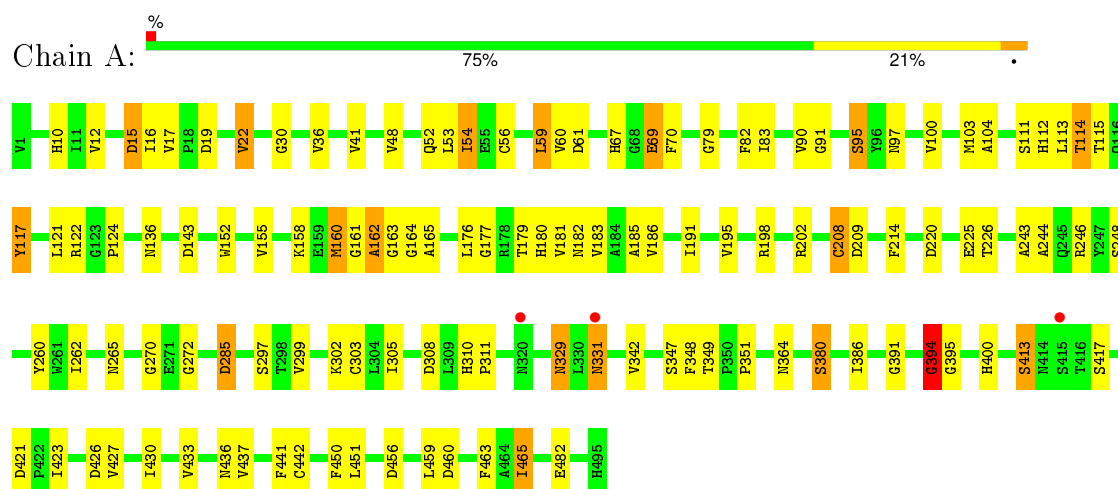
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	478	Total	O	0	0
			478	478		
6	B	376	Total	O	0	0
			376	376		
6	C	340	Total	O	0	0
			340	340		

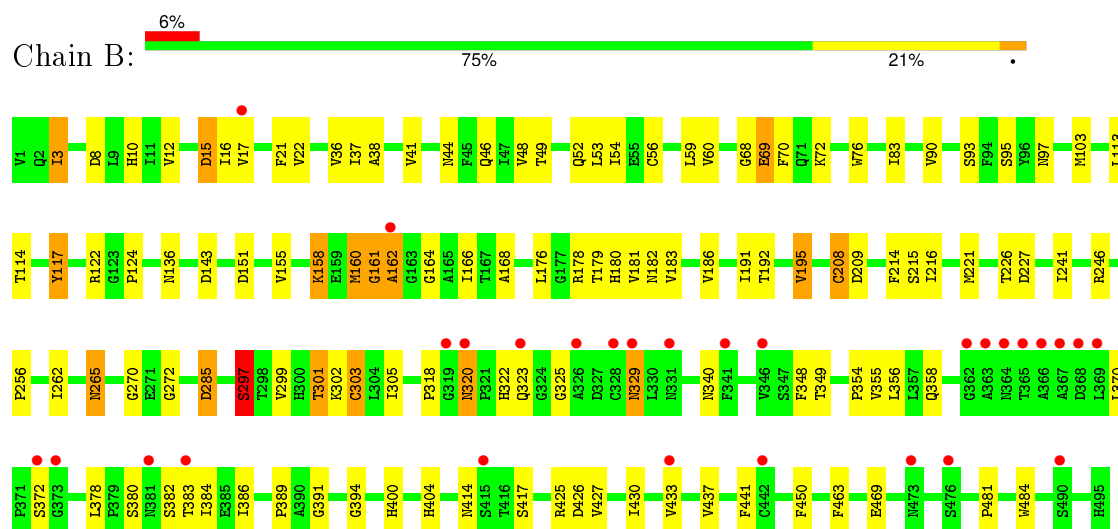
3 Residue-property plots

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.

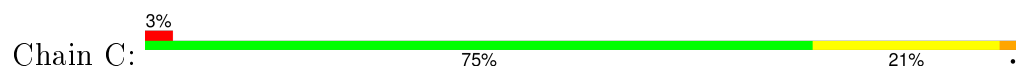
• Molecule 1: Laccase

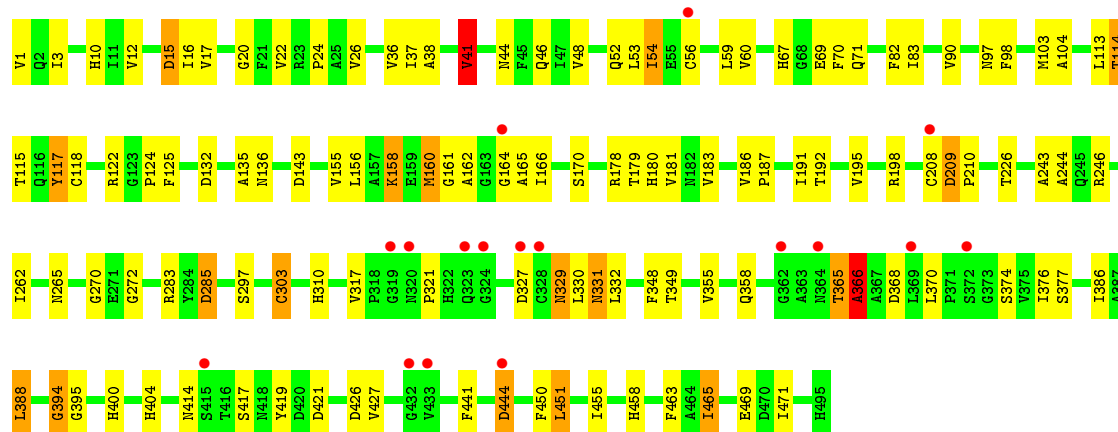


• Molecule 1: Laccase



• Molecule 1: Laccase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.53Å 140.66Å 174.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 70.33 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.3 (30.00-2.00) 98.3 (70.33-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.181 , 0.241 0.185 , 0.244	Depositor DCC
R_{free} test set	6068 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.485	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 62.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 120451 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12647	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, CU, CBS

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	1.12	12/3857 (0.3%)	0.96	5/5297 (0.1%)
1	B	1.03	7/3865 (0.2%)	0.95	5/5307 (0.1%)
1	C	1.01	1/3850 (0.0%)	0.96	9/5287 (0.2%)
All	All	1.05	20/11572 (0.2%)	0.95	19/15891 (0.1%)

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	285	ASP	CB-CG	5.92	1.64	1.51
1	A	208[A]	CYS	CB-SG	-5.89	1.72	1.81
1	A	208[B]	CYS	CB-SG	-5.89	1.72	1.81
1	B	69	GLU	CG-CD	-5.80	1.43	1.51
1	C	198	ARG	CG-CD	5.76	1.66	1.51

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	394	GLY	N-CA-C	-7.11	95.34	113.10
1	C	161	GLY	N-CA-C	-6.11	97.82	113.10
1	C	395	GLY	N-CA-C	-6.09	97.88	113.10
1	B	303	CYS	CA-CB-SG	-5.90	103.38	114.00
1	C	303	CYS	CA-CB-SG	-5.88	103.42	114.00

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	3734	0	3504	128	0
1	B	3739	0	3520	139	1
1	C	3731	0	3508	139	1
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
3	A	20	0	0	0	0
3	B	5	0	0	0	0
3	C	10	0	0	9	0
4	A	6	0	8	4	0
4	B	30	0	40	9	0
4	C	12	0	16	12	0
5	A	56	0	52	4	0
5	B	56	0	52	5	0
5	C	42	0	38	7	0
6	A	478	0	0	32	0
6	B	376	0	0	23	0
6	C	340	0	0	24	0
All	All	12647	0	10738	394	1

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 18.

The worst 5 of 394 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:414:ASN:ND2	5:C:509:CBS:C1B	1.68	1.53
1:B:414:ASN:HD21	5:B:511:CBS:C1B	1.18	1.52
1:A:436:ASN:ND2	5:A:511:CBS:C1B	1.68	1.51
1:B:414:ASN:ND2	5:B:511:CBS:C1B	1.75	1.43
5:C:509:CBS:H3B	5:C:509:CBS:C8B	1.41	1.42

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:ASN:ND2	1:C:136:ASN:ND2[3_554]	1.85	0.35

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	498/495 (101%)	476 (96%)	18 (4%)	4 (1%)	24	15
1	B	498/495 (101%)	473 (95%)	20 (4%)	5 (1%)	19	11
1	C	496/495 (100%)	470 (95%)	23 (5%)	3 (1%)	30	22
All	All	1492/1485 (100%)	1419 (95%)	61 (4%)	12 (1%)	24	15

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	ALA
1	B	162	ALA
1	C	209	ASP
1	C	366	ALA
1	A	209	ASP

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	410/406 (101%)	383 (93%)	27 (7%)	21	14
1	B	410/406 (101%)	385 (94%)	25 (6%)	23	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	408/406 (100%)	379 (93%)	29 (7%)	18	12
All	All	1228/1218 (101%)	1147 (93%)	81 (7%)	21	14

5 of 81 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	226	THR
1	B	349	THR
1	C	388	LEU
1	B	246	ARG
1	B	301	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 47 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	180	HIS
1	B	322	HIS
1	C	331	ASN
1	B	252	ASN
1	B	358	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 33 ligands modelled in this entry, 12 are monoatomic - leaving 21 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	SO4	A	505	-	4,4,4	0.45	0	6,6,6	0.74	0
3	SO4	A	506	-	4,4,4	0.23	0	6,6,6	0.64	0
3	SO4	A	507	-	4,4,4	1.54	0	6,6,6	0.99	0
3	SO4	A	508	-	4,4,4	0.62	0	6,6,6	0.44	0
4	GOL	A	509	-	5,5,5	0.46	0	5,5,5	0.62	0
5	CBS	A	510	1	29,29,30	0.57	0	36,41,43	1.21	2 (5%)
5	CBS	A	511	-	29,29,30	0.61	0	36,41,43	2.35	6 (16%)
3	SO4	B	505	-	4,4,4	0.40	0	6,6,6	0.65	0
4	GOL	B	506	-	5,5,5	0.70	0	5,5,5	1.85	2 (40%)
4	GOL	B	507	-	5,5,5	0.66	0	5,5,5	0.97	0
4	GOL	B	508	-	5,5,5	0.11	0	5,5,5	0.55	0
4	GOL	B	509	-	5,5,5	0.84	0	5,5,5	1.29	1 (20%)
4	GOL	B	510	-	5,5,5	0.44	0	5,5,5	0.54	0
5	CBS	B	511	-	29,29,30	0.50	0	36,41,43	1.23	2 (5%)
5	CBS	B	512	1	29,29,30	0.62	0	36,41,43	2.04	10 (27%)
3	SO4	C	505	-	4,4,4	0.81	0	6,6,6	0.78	0
3	SO4	C	506	-	4,4,4	0.17	0	6,6,6	0.43	0
4	GOL	C	507	-	5,5,5	0.41	0	5,5,5	1.45	1 (20%)
4	GOL	C	508	-	5,5,5	0.42	0	5,5,5	1.54	1 (20%)
5	CBS	C	509	-	29,29,30	0.53	0	36,41,43	1.08	3 (8%)
5	CBS	C	510	1	14,14,30	2.21	2 (14%)	15,19,43	2.82	5 (33%)

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	SO4	A	505	-	-	0/0/0/0	0/0/0/0
3	SO4	A	506	-	-	0/0/0/0	0/0/0/0
3	SO4	A	507	-	-	0/0/0/0	0/0/0/0
3	SO4	A	508	-	-	0/0/0/0	0/0/0/0
4	GOL	A	509	-	-	0/4/4/4	0/0/0/0
5	CBS	A	510	1	-	0/16/53/56	0/2/2/2
5	CBS	A	511	-	-	0/16/53/56	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	B	505	-	-	0/0/0/0	0/0/0/0
4	GOL	B	506	-	-	0/4/4/4	0/0/0/0
4	GOL	B	507	-	-	0/4/4/4	0/0/0/0
4	GOL	B	508	-	-	0/4/4/4	0/0/0/0
4	GOL	B	509	-	-	0/4/4/4	0/0/0/0
4	GOL	B	510	-	-	0/4/4/4	0/0/0/0
5	CBS	B	511	-	-	0/16/53/56	0/2/2/2
5	CBS	B	512	1	-	0/16/53/56	0/2/2/2
3	SO4	C	505	-	-	0/0/0/0	0/0/0/0
3	SO4	C	506	-	-	0/0/0/0	0/0/0/0
4	GOL	C	507	-	-	0/4/4/4	0/0/0/0
4	GOL	C	508	-	-	0/4/4/4	0/0/0/0
5	CBS	C	509	-	-	0/16/53/56	0/2/2/2
5	CBS	C	510	1	-	0/6/23/56	0/1/1/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	510	CBS	C7B-N2B	2.13	1.42	1.34
5	C	510	CBS	O7B-C7B	7.64	1.40	1.23

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	511	CBS	C4A-C3A-C2A	-9.49	97.27	110.43
5	C	510	CBS	C3B-C2B-N2B	-5.97	96.26	110.56
5	A	511	CBS	O5A-C1A-C2A	-4.47	101.05	110.78
5	B	512	CBS	O5A-C1A-C2A	-3.75	102.62	110.78
5	C	510	CBS	C3B-C4B-C5B	-3.71	103.73	110.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

11 monomers are involved in 50 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	509	GOL	4	0
5	A	511	CBS	4	0
4	B	506	GOL	6	0
4	B	507	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	510	GOL	2	0
5	B	511	CBS	3	0
5	B	512	CBS	2	0
3	C	505	SO4	4	0
3	C	506	SO4	5	0
4	C	507	GOL	12	0
5	C	509	CBS	7	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th 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 ²)	Q<0.9
1	A	495/495 (100%)	-0.13	3 (0%) 90 90	9, 19, 33, 46	0
1	B	495/495 (100%)	0.19	29 (5%) 26 27	10, 24, 45, 56	0
1	C	495/495 (100%)	0.16	17 (3%) 49 50	11, 26, 47, 62	0
All	All	1485/1485 (100%)	0.08	49 (3%) 50 51	9, 23, 43, 62	0

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	415	SER	6.9
1	B	328	CYS	6.1
1	C	324	GLY	5.2
1	B	320	ASN	4.4
1	B	415	SER	4.2

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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, 95th 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 ²)	Q<0.9
4	GOL	A	509	6/6	0.82	0.20	6.40	32,48,50,53	0
4	GOL	C	507	6/6	0.94	0.25	5.66	26,34,41,47	0
4	GOL	B	508	6/6	0.93	0.20	5.63	46,52,53,55	0
4	GOL	B	506	6/6	0.90	0.23	4.95	22,32,35,35	0
3	SO4	A	505	5/5	0.98	0.14	4.71	31,36,40,41	0
5	CBS	B	511	28/29	0.78	0.34	4.07	48,62,69,73	0
3	SO4	C	506	5/5	0.61	0.32	3.83	46,47,49,49	5
3	SO4	A	507	5/5	0.99	0.14	3.39	19,22,26,28	0
4	GOL	B	509	6/6	0.90	0.16	3.02	20,26,32,40	0
5	CBS	C	509	28/29	0.76	0.27	2.85	60,68,72,85	0
3	SO4	A	506	5/5	0.96	0.15	2.79	60,60,61,61	0
4	GOL	C	508	6/6	0.92	0.16	2.52	34,38,41,46	0
5	CBS	A	510	28/29	0.91	0.19	1.55	29,42,57,60	0
5	CBS	B	512	28/29	0.84	0.30	1.17	45,63,83,95	0
5	CBS	A	511	28/29	0.85	0.23	1.00	28,45,58,61	0
4	GOL	B	510	6/6	0.92	0.14	0.47	23,28,30,30	0
4	GOL	B	507	6/6	0.95	0.13	0.34	23,30,34,35	0
5	CBS	C	510	14/29	0.79	0.20	-0.02	54,60,62,62	0
3	SO4	C	505	5/5	0.99	0.12	-0.19	21,26,28,29	0
2	CU	A	501	1/1	1.00	0.08	-1.46	19,19,19,19	0
2	CU	A	504	1/1	1.00	0.08	-1.67	18,18,18,18	0
2	CU	B	501	1/1	1.00	0.07	-1.68	20,20,20,20	0
2	CU	A	503	1/1	1.00	0.08	-1.88	19,19,19,19	0
2	CU	B	502	1/1	1.00	0.07	-1.99	19,19,19,19	0
2	CU	C	501	1/1	0.99	0.07	-2.27	31,31,31,31	0
2	CU	A	502	1/1	1.00	0.07	-2.59	15,15,15,15	0
2	CU	B	504	1/1	1.00	0.05	-2.80	26,26,26,26	0
2	CU	B	503	1/1	1.00	0.07	-2.96	23,23,23,23	0
2	CU	C	502	1/1	1.00	0.07	-3.40	22,22,22,22	0
2	CU	C	504	1/1	1.00	0.05	-5.84	28,28,28,28	0
2	CU	C	503	1/1	0.99	0.07	-7.93	26,26,26,26	0
3	SO4	B	505	5/5	0.88	0.16	-	29,31,32,33	5
3	SO4	A	508	5/5	0.96	0.16	-	19,22,23,26	5

6.5 Other polymers

There are no such residues in this entry.