



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 06:47 pm BST

PDB ID : 3E4O  
Title : Crystal structure of succinate bound state DctB  
Authors : Zhou, Y.F.; Nan, J.; Nan, B.Y.; Liang, Y.H.; Panjikar, S.; Su, X.D.  
Deposited on : 2008-08-12  
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

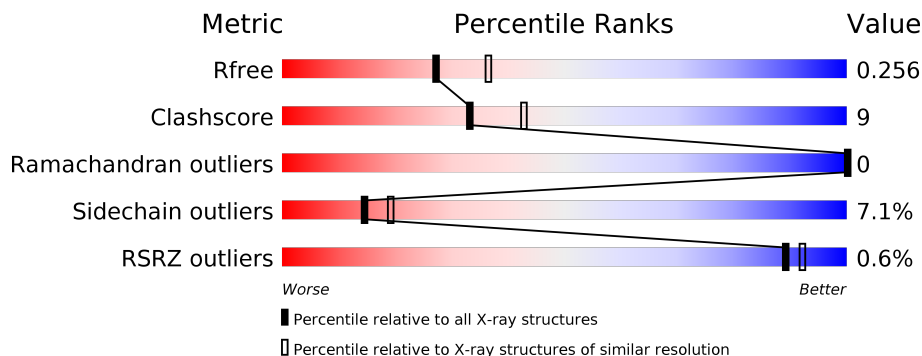
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	305	
1	B	305	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4107 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 C4-dicarboxylate transport sensor protein dctB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	254	1945	1232	349	360	4	0	0	0
1	B	254	1945	1232	349	360	4	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	MET	-	EXPRESSION TAG	UNP P13633
A	9	GLY	-	EXPRESSION TAG	UNP P13633
A	10	SER	-	EXPRESSION TAG	UNP P13633
A	11	SER	-	EXPRESSION TAG	UNP P13633
A	12	HIS	-	EXPRESSION TAG	UNP P13633
A	13	HIS	-	EXPRESSION TAG	UNP P13633
A	14	HIS	-	EXPRESSION TAG	UNP P13633
A	15	HIS	-	EXPRESSION TAG	UNP P13633
A	16	HIS	-	EXPRESSION TAG	UNP P13633
A	17	HIS	-	EXPRESSION TAG	UNP P13633
A	18	SER	-	EXPRESSION TAG	UNP P13633
A	19	SER	-	EXPRESSION TAG	UNP P13633
A	20	GLY	-	EXPRESSION TAG	UNP P13633
A	21	LEU	-	EXPRESSION TAG	UNP P13633
A	22	VAL	-	EXPRESSION TAG	UNP P13633
A	23	PRO	-	EXPRESSION TAG	UNP P13633
A	24	ARG	-	EXPRESSION TAG	UNP P13633
A	25	GLY	-	EXPRESSION TAG	UNP P13633
A	26	SER	-	EXPRESSION TAG	UNP P13633
A	27	HIS	-	EXPRESSION TAG	UNP P13633
A	28	MSE	-	EXPRESSION TAG	UNP P13633
A	29	ALA	-	EXPRESSION TAG	UNP P13633
A	30	SER	-	EXPRESSION TAG	UNP P13633
A	31	MSE	-	EXPRESSION TAG	UNP P13633
A	32	THR	-	EXPRESSION TAG	UNP P13633

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Chain	Residue	Modelled	Actual	Comment	Reference
A	33	GLY	-	EXPRESSION TAG	UNP P13633
A	34	GLY	-	EXPRESSION TAG	UNP P13633
A	35	GLU	-	EXPRESSION TAG	UNP P13633
A	36	GLU	-	EXPRESSION TAG	UNP P13633
A	37	MSE	-	EXPRESSION TAG	UNP P13633
A	38	GLY	-	EXPRESSION TAG	UNP P13633
A	39	ARG	-	EXPRESSION TAG	UNP P13633
A	40	GLY	-	EXPRESSION TAG	UNP P13633
A	41	SER	-	EXPRESSION TAG	UNP P13633
A	174	LYS	ASN	SEE REMARK 999	UNP P13633
A	309	ASN	LYS	SEE REMARK 999	UNP P13633
B	8	MET	-	EXPRESSION TAG	UNP P13633
B	9	GLY	-	EXPRESSION TAG	UNP P13633
B	10	SER	-	EXPRESSION TAG	UNP P13633
B	11	SER	-	EXPRESSION TAG	UNP P13633
B	12	HIS	-	EXPRESSION TAG	UNP P13633
B	13	HIS	-	EXPRESSION TAG	UNP P13633
B	14	HIS	-	EXPRESSION TAG	UNP P13633
B	15	HIS	-	EXPRESSION TAG	UNP P13633
B	16	HIS	-	EXPRESSION TAG	UNP P13633
B	17	HIS	-	EXPRESSION TAG	UNP P13633
B	18	SER	-	EXPRESSION TAG	UNP P13633
B	19	SER	-	EXPRESSION TAG	UNP P13633
B	20	GLY	-	EXPRESSION TAG	UNP P13633
B	21	LEU	-	EXPRESSION TAG	UNP P13633
B	22	VAL	-	EXPRESSION TAG	UNP P13633
B	23	PRO	-	EXPRESSION TAG	UNP P13633
B	24	ARG	-	EXPRESSION TAG	UNP P13633
B	25	GLY	-	EXPRESSION TAG	UNP P13633
B	26	SER	-	EXPRESSION TAG	UNP P13633
B	27	HIS	-	EXPRESSION TAG	UNP P13633
B	28	MSE	-	EXPRESSION TAG	UNP P13633
B	29	ALA	-	EXPRESSION TAG	UNP P13633
B	30	SER	-	EXPRESSION TAG	UNP P13633
B	31	MSE	-	EXPRESSION TAG	UNP P13633
B	32	THR	-	EXPRESSION TAG	UNP P13633
B	33	GLY	-	EXPRESSION TAG	UNP P13633
B	34	GLY	-	EXPRESSION TAG	UNP P13633
B	35	GLU	-	EXPRESSION TAG	UNP P13633
B	36	GLU	-	EXPRESSION TAG	UNP P13633
B	37	MSE	-	EXPRESSION TAG	UNP P13633
B	38	GLY	-	EXPRESSION TAG	UNP P13633

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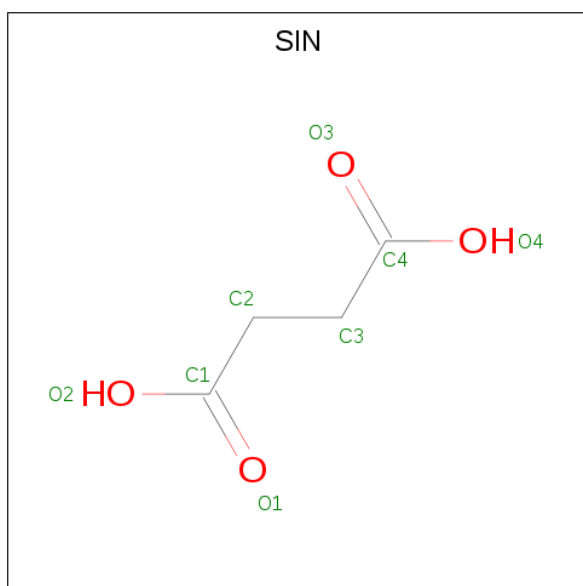
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Chain	Residue	Modelled	Actual	Comment	Reference
B	39	ARG	-	EXPRESSION TAG	UNP P13633
B	40	GLY	-	EXPRESSION TAG	UNP P13633
B	41	SER	-	EXPRESSION TAG	UNP P13633
B	174	LYS	ASN	SEE REMARK 999	UNP P13633
B	309	ASN	LYS	SEE REMARK 999	UNP P13633

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0

- Molecule 3 is SUCCINIC ACID (three-letter code: SIN) (formula: C<sub>4</sub>H<sub>6</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 8 4 4	0	0
3	A	1	Total C O 8 4 4	0	0

- Molecule 4 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	A	110	Total 110	O 110	0	0
4	B	89	Total 89	O 89	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.89Å 66.59Å 93.60Å 90.00° 94.71° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 48.79 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.30) 99.6 (48.79-2.30)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.58 (at 2.29Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.200 , 0.252 0.203 , 0.256	Depositor DCC
$R_{free}$ test set	1299 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.9	Xtrriage
Anisotropy	0.426	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 30.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4107	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.15% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, SIN

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	0.69	1/1981 (0.1%)	0.74	1/2687 (0.0%)
1	B	0.63	1/1981 (0.1%)	0.70	0/2687
All	All	0.66	2/3962 (0.1%)	0.72	1/5374 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	169	MSE	SE-CE	-5.87	1.60	1.95
1	A	169	MSE	SE-CE	-5.85	1.60	1.95

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	150	ALA	N-CA-C	-7.86	89.77	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149	TYR	Peptide

## 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	1945	0	1975	26	0
1	B	1945	0	1975	44	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	16	0	8	0	0
4	A	110	0	0	4	0
4	B	89	0	0	2	1
All	All	4107	0	3958	68	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 9.

All (68) 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:294:MSE:HG3	1:B:301:ARG:HH21	1.15	1.03
1:B:182:ARG:HH21	1:B:182:ARG:HG2	1.24	1.02
1:B:294:MSE:HG2	1:B:301:ARG:NH2	1.76	1.00
1:B:182:ARG:HH21	1:B:182:ARG:CG	1.92	0.83
1:B:294:MSE:HG2	1:B:301:ARG:CZ	2.09	0.80
1:A:239:ILE:CG2	1:A:244:LEU:HD13	2.13	0.79
1:B:169:MSE:HE1	1:B:174:LYS:O	1.83	0.77
1:B:241:GLU:O	1:B:241:GLU:HG3	1.84	0.77
1:A:294:MSE:HG3	1:B:301:ARG:NH2	1.97	0.77
1:A:73:ARG:HD3	1:A:77:GLU:OE1	1.89	0.73
1:A:292:GLU:OE2	1:A:301:ARG:NE	2.24	0.71
1:A:301:ARG:NH1	4:A:904:HOH:O	2.24	0.70
1:B:93:ARG:HD2	1:B:185:ASP:O	1.94	0.68
1:A:239:ILE:HG21	1:A:244:LEU:HD13	1.75	0.67
1:B:182:ARG:HG2	1:B:182:ARG:NH2	2.04	0.67
1:B:294:MSE:CG	1:B:301:ARG:NH2	2.55	0.66
1:B:169:MSE:HE1	1:B:174:LYS:C	2.16	0.64
1:B:245:ALA:HB3	1:B:246:PRO:HD3	1.79	0.63
1:B:68:LYS:NZ	1:B:304:GLN:HE21	1.98	0.62
1:B:102:PRO:O	1:B:106:ARG:HG3	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:ARG:HG3	1:A:296:PRO:O	2.03	0.59
1:B:294:MSE:SE	1:B:301:ARG:HH22	2.36	0.59
1:B:245:ALA:N	1:B:246:PRO:HD2	2.20	0.57
1:A:153:ASP:OD1	4:A:816:HOH:O	2.18	0.57
1:B:245:ALA:HB3	1:B:246:PRO:CD	2.34	0.56
1:B:108:ASN:HD21	1:B:134:ALA:HA	1.70	0.56
1:B:71:LEU:O	1:B:75:VAL:HG23	2.05	0.56
1:A:243:ARG:O	1:A:247:ILE:HG13	2.05	0.56
1:B:182:ARG:NH2	1:B:182:ARG:CG	2.62	0.54
1:B:295:VAL:HG22	1:B:298:THR:OG1	2.08	0.54
1:A:153:ASP:HB3	1:A:157:LEU:HD12	1.92	0.52
1:B:105:ASP:O	1:B:109:ARG:HG3	2.10	0.51
1:A:239:ILE:HG23	1:A:244:LEU:HD13	1.91	0.51
1:A:86:LEU:HD21	1:A:114:LEU:CD1	2.41	0.51
1:A:101:ARG:HB3	1:A:102:PRO:HD3	1.92	0.51
1:A:245:ALA:HB3	1:A:246:PRO:HD3	1.93	0.50
1:A:237:LYS:NZ	4:A:904:HOH:O	2.46	0.49
1:B:294:MSE:HG2	1:B:301:ARG:HH22	1.68	0.49
1:A:228:LEU:HD12	1:A:229:PRO:HD2	1.93	0.48
1:A:183:ARG:HD2	1:A:190:PRO:HB3	1.95	0.48
1:B:208:GLN:NE2	4:B:689:HOH:O	2.47	0.47
1:A:239:ILE:HG21	1:A:244:LEU:CD1	2.44	0.46
1:A:75:VAL:HG12	1:A:198:LEU:HD21	1.96	0.46
1:B:169:MSE:CE	1:B:174:LYS:C	2.84	0.46
1:B:207:TRP:HB2	1:B:227:SER:HB3	1.97	0.46
1:B:68:LYS:HZ2	1:B:304:GLN:HE21	1.63	0.46
1:B:241:GLU:HA	1:B:244:LEU:HB2	1.98	0.44
1:A:163:MSE:HE3	1:A:163:MSE:HB2	1.89	0.44
1:B:216:VAL:HG13	1:B:300:TRP:HB3	1.99	0.44
1:B:158:ALA:O	1:B:183:ARG:HB2	2.18	0.43
1:A:238:PRO:HD3	4:A:908:HOH:O	2.18	0.43
1:B:284:SER:O	1:B:285:THR:HG22	2.19	0.43
1:A:239:ILE:CG2	1:A:239:ILE:O	2.67	0.42
1:B:169:MSE:HE3	1:B:174:LYS:HA	2.02	0.42
1:B:97:LEU:HD23	1:B:191:LEU:HD21	2.03	0.41
1:A:275:THR:HA	1:A:288:PHE:O	2.20	0.41
1:B:245:ALA:N	1:B:246:PRO:CD	2.84	0.41
1:B:245:ALA:CB	1:B:246:PRO:CD	2.97	0.41
1:B:68:LYS:HZ3	1:B:304:GLN:HE21	1.67	0.41
1:B:108:ASN:HB3	1:B:138:TRP:CD1	2.56	0.41
1:A:82:LEU:HB3	1:A:83:PRO:HD3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:VAL:HG22	1:B:204:GLU:N	2.35	0.41
1:B:216:VAL:HA	1:B:301:ARG:O	2.21	0.40
1:B:292:GLU:HG3	1:B:303:GLU:HG3	2.03	0.40
1:B:219:ARG:HD2	1:B:219:ARG:O	2.21	0.40
1:B:128:ARG:HG2	4:B:642:HOH:O	2.21	0.40
1:B:165:GLU:HG2	1:B:178:LEU:HD11	2.03	0.40
1:A:207:TRP:HB2	1:A:227:SER:HB3	2.02	0.40

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:284:SER:OG	4:B:607:HOH:O[4_546]	2.19	0.01

## 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	252/305 (83%)	248 (98%)	4 (2%)	0	100	100
1	B	252/305 (83%)	247 (98%)	5 (2%)	0	100	100
All	All	504/610 (83%)	495 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	203/235 (86%)	193 (95%)	10 (5%)	25	35
1	B	203/235 (86%)	184 (91%)	19 (9%)	8	10
All	All	406/470 (86%)	377 (93%)	29 (7%)	14	19

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	SER
1	A	73	ARG
1	A	98	SER
1	A	156	ARG
1	A	163	MSE
1	A	201	ASP
1	A	208	GLN
1	A	239	ILE
1	A	297	SER
1	A	301	ARG
1	B	63	ILE
1	B	73	ARG
1	B	101	ARG
1	B	105	ASP
1	B	114	LEU
1	B	128	ARG
1	B	163	MSE
1	B	182	ARG
1	B	203	VAL
1	B	208	GLN
1	B	219	ARG
1	B	241	GLU
1	B	243	ARG
1	B	279	LEU
1	B	283	ASP
1	B	285	THR
1	B	295	VAL
1	B	301	ARG
1	B	308	LEU

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	208	GLN
1	B	79	GLN
1	B	108	ASN
1	B	304	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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	SIN	A	801	-	1,7,7	0.17	0	2,8,8	1.40	0
3	SIN	A	701	-	1,7,7	0.04	0	2,8,8	0.77	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	SIN	A	801	-	-	0/1/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SIN	A	701	-	-	0/1/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	250/305 (81%)	-0.12	2 (0%) 86 89	21, 34, 46, 54	0
1	B	250/305 (81%)	-0.00	1 (0%) 92 95	21, 35, 50, 58	2 (0%)
All	All	500/610 (81%)	-0.06	3 (0%) 89 92	21, 35, 49, 58	2 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	239	ILE	3.3
1	B	238	PRO	2.7
1	A	243	ARG	2.1

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

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

### 6.3 Carbohydrates [i](#)

There are no 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. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	B	601	1/1	0.97	0.05	38,38,38,38	0
2	MG	A	501	1/1	0.97	0.05	34,34,34,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	SIN	A	701	8/8	0.97	0.10	26,28,29,30	0
3	SIN	A	801	8/8	0.98	0.07	19,21,22,23	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.