



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2020 – 02:06 PM EST

PDB ID : 6EBC
Title : OhrB (Organic Hydroperoxide Resistance protein) wild type from *Chromobacterium violaceum* and reduced by DTT
Authors : Domingos, R.M.; Teixeira, R.D.; Alegria, T.G.P.; Vieira, P.S.; Murakami, M.T.; Netto, L.E.S.
Deposited on : 2018-08-06
Resolution : 1.87 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/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 : 1.8.0 (224370), CSD as540be (2019)
Xtriage (Phenix) : 1.13
EDS : 2.4
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.4

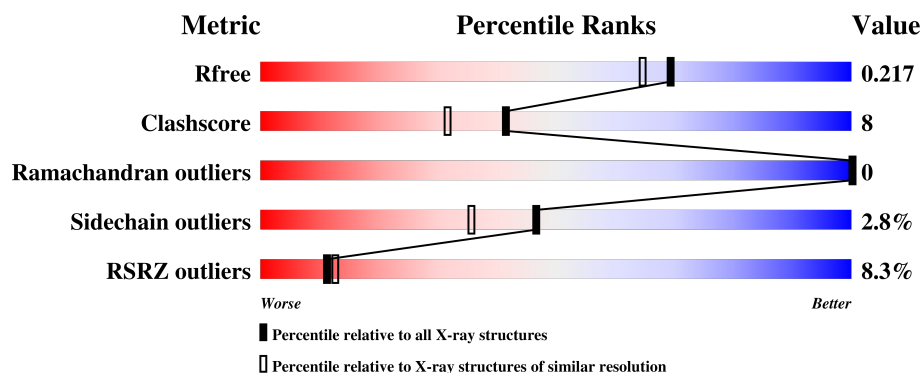
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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}	111664	8255 (1.90-1.86)
Clashscore	122126	9028 (1.90-1.86)
Ramachandran outliers	120053	8930 (1.90-1.86)
Sidechain outliers	120020	8930 (1.90-1.86)
RSRZ outliers	108989	8087 (1.90-1.86)

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	161	<div> <div>6%</div> <div>73%14%13%</div> </div>
1	B	161	<div> <div>9%</div> <div>72%16%11%</div> </div>
1	C	161	<div> <div>6%</div> <div>73%14%13%</div> </div>
1	D	161	<div> <div>9%</div> <div>71%17%12%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4394 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 Organic hydroperoxide resistance protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	140	Total	C	N	O	S	0	0	0
			1020	632	182	203	3			
1	B	143	Total	C	N	O	S	0	0	0
			1044	646	187	207	4			
1	C	140	Total	C	N	O	S	0	0	0
			1020	632	182	203	3			
1	D	142	Total	C	N	O	S	0	0	0
			1038	643	186	205	4			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A202B6V5
A	-18	GLY	-	expression tag	UNP A0A202B6V5
A	-17	SER	-	expression tag	UNP A0A202B6V5
A	-16	SER	-	expression tag	UNP A0A202B6V5
A	-15	HIS	-	expression tag	UNP A0A202B6V5
A	-14	HIS	-	expression tag	UNP A0A202B6V5
A	-13	HIS	-	expression tag	UNP A0A202B6V5
A	-12	HIS	-	expression tag	UNP A0A202B6V5
A	-11	HIS	-	expression tag	UNP A0A202B6V5
A	-10	HIS	-	expression tag	UNP A0A202B6V5
A	-9	SER	-	expression tag	UNP A0A202B6V5
A	-8	SER	-	expression tag	UNP A0A202B6V5
A	-7	GLY	-	expression tag	UNP A0A202B6V5
A	-6	LEU	-	expression tag	UNP A0A202B6V5
A	-5	VAL	-	expression tag	UNP A0A202B6V5
A	-4	PRO	-	expression tag	UNP A0A202B6V5
A	-3	ARG	-	expression tag	UNP A0A202B6V5
A	-2	GLY	-	expression tag	UNP A0A202B6V5
A	-1	SER	-	expression tag	UNP A0A202B6V5
A	0	HIS	-	expression tag	UNP A0A202B6V5
B	-19	MET	-	initiating methionine	UNP A0A202B6V5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP A0A202B6V5
B	-17	SER	-	expression tag	UNP A0A202B6V5
B	-16	SER	-	expression tag	UNP A0A202B6V5
B	-15	HIS	-	expression tag	UNP A0A202B6V5
B	-14	HIS	-	expression tag	UNP A0A202B6V5
B	-13	HIS	-	expression tag	UNP A0A202B6V5
B	-12	HIS	-	expression tag	UNP A0A202B6V5
B	-11	HIS	-	expression tag	UNP A0A202B6V5
B	-10	HIS	-	expression tag	UNP A0A202B6V5
B	-9	SER	-	expression tag	UNP A0A202B6V5
B	-8	SER	-	expression tag	UNP A0A202B6V5
B	-7	GLY	-	expression tag	UNP A0A202B6V5
B	-6	LEU	-	expression tag	UNP A0A202B6V5
B	-5	VAL	-	expression tag	UNP A0A202B6V5
B	-4	PRO	-	expression tag	UNP A0A202B6V5
B	-3	ARG	-	expression tag	UNP A0A202B6V5
B	-2	GLY	-	expression tag	UNP A0A202B6V5
B	-1	SER	-	expression tag	UNP A0A202B6V5
B	0	HIS	-	expression tag	UNP A0A202B6V5
C	-19	MET	-	initiating methionine	UNP A0A202B6V5
C	-18	GLY	-	expression tag	UNP A0A202B6V5
C	-17	SER	-	expression tag	UNP A0A202B6V5
C	-16	SER	-	expression tag	UNP A0A202B6V5
C	-15	HIS	-	expression tag	UNP A0A202B6V5
C	-14	HIS	-	expression tag	UNP A0A202B6V5
C	-13	HIS	-	expression tag	UNP A0A202B6V5
C	-12	HIS	-	expression tag	UNP A0A202B6V5
C	-11	HIS	-	expression tag	UNP A0A202B6V5
C	-10	HIS	-	expression tag	UNP A0A202B6V5
C	-9	SER	-	expression tag	UNP A0A202B6V5
C	-8	SER	-	expression tag	UNP A0A202B6V5
C	-7	GLY	-	expression tag	UNP A0A202B6V5
C	-6	LEU	-	expression tag	UNP A0A202B6V5
C	-5	VAL	-	expression tag	UNP A0A202B6V5
C	-4	PRO	-	expression tag	UNP A0A202B6V5
C	-3	ARG	-	expression tag	UNP A0A202B6V5
C	-2	GLY	-	expression tag	UNP A0A202B6V5
C	-1	SER	-	expression tag	UNP A0A202B6V5
C	0	HIS	-	expression tag	UNP A0A202B6V5
D	-19	MET	-	initiating methionine	UNP A0A202B6V5
D	-18	GLY	-	expression tag	UNP A0A202B6V5
D	-17	SER	-	expression tag	UNP A0A202B6V5

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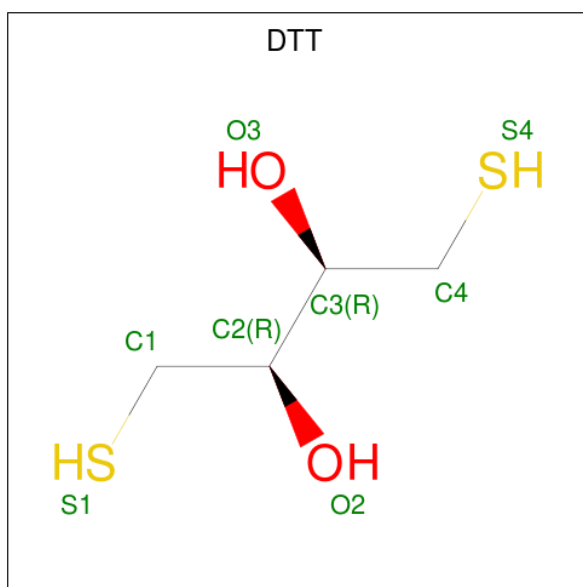
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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP A0A202B6V5
D	-15	HIS	-	expression tag	UNP A0A202B6V5
D	-14	HIS	-	expression tag	UNP A0A202B6V5
D	-13	HIS	-	expression tag	UNP A0A202B6V5
D	-12	HIS	-	expression tag	UNP A0A202B6V5
D	-11	HIS	-	expression tag	UNP A0A202B6V5
D	-10	HIS	-	expression tag	UNP A0A202B6V5
D	-9	SER	-	expression tag	UNP A0A202B6V5
D	-8	SER	-	expression tag	UNP A0A202B6V5
D	-7	GLY	-	expression tag	UNP A0A202B6V5
D	-6	LEU	-	expression tag	UNP A0A202B6V5
D	-5	VAL	-	expression tag	UNP A0A202B6V5
D	-4	PRO	-	expression tag	UNP A0A202B6V5
D	-3	ARG	-	expression tag	UNP A0A202B6V5
D	-2	GLY	-	expression tag	UNP A0A202B6V5
D	-1	SER	-	expression tag	UNP A0A202B6V5
D	0	HIS	-	expression tag	UNP A0A202B6V5

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Cl 2 2	0	0
2	A	2	Total Cl 2 2	0	0
2	D	1	Total Cl 1 1	0	0
2	C	2	Total Cl 2 2	0	0

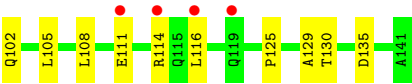
- Molecule 3 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (three-letter code: DTT) (formula: C₄H₁₀O₂S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	O	S	0	0
			8	4	2	2		
3	C	1	Total	C	O	S	0	0
			8	4	2	2		
3	D	1	Total	C	O	S	0	0
			8	4	2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	56	Total	O	0	0
			56	56		
4	B	68	Total	O	0	0
			68	68		
4	C	68	Total	O	0	0
			68	68		
4	D	49	Total	O	0	0
			49	49		



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	88.16Å 88.16Å 169.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.00 – 1.87 34.82 – 1.87	Depositor EDS
% Data completeness (in resolution range)	100.0 (35.00-1.87) 100.0 (34.82-1.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.193 , 0.214 0.199 , 0.217	Depositor DCC
R_{free} test set	3029 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 41.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.045 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4394	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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, 0.333 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: CL, DTT

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	0.76	0/1033	0.98	0/1397
1	B	0.79	1/1058 (0.1%)	1.00	0/1430
1	C	0.80	0/1033	1.00	0/1397
1	D	0.78	0/1052	0.95	0/1422
All	All	0.78	1/4176 (0.0%)	0.98	0/5646

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	98	GLU	CD-OE1	-5.19	1.20	1.25

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1020	0	1023	17	0
1	B	1044	0	1047	22	0
1	C	1020	0	1023	19	0
1	D	1038	0	1042	23	0
2	A	2	0	0	1	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	2	0	0	1	0
2	D	1	0	0	1	0
3	B	8	0	10	3	0
3	C	8	0	10	2	0
3	D	8	0	10	3	0
4	A	56	0	0	0	0
4	B	68	0	0	0	0
4	C	68	0	0	1	0
4	D	49	0	0	1	0
All	All	4394	0	4165	63	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 8.

All (63) 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:83:SER:HB2	1:B:102:GLN:HG3	1.51	0.93
1:C:104:ARG:HG2	1:C:141:ALA:HB3	1.59	0.85
1:C:104:ARG:HG2	1:C:141:ALA:CB	2.25	0.64
1:A:104:ARG:HG2	1:A:141:ALA:HB3	1.81	0.63
1:C:139:ARG:HD3	4:C:334:HOH:O	2.02	0.59
1:A:25:SER:OG	2:A:202:CL:CL	2.61	0.56
1:D:108:LEU:HD11	1:D:116:LEU:HD12	1.89	0.53
1:B:104:ARG:HG2	1:B:141:ALA:HB3	1.91	0.51
1:B:89:GLY:HA3	1:D:3:ILE:HG21	1.91	0.51
1:B:122:ILE:HG21	1:C:26:ASP:HA	1.92	0.51
1:B:130:THR:HA	1:D:129:ALA:O	2.11	0.51
1:B:68:VAL:HG11	1:B:119:GLN:HB3	1.93	0.51
1:A:113:ALA:HB1	1:A:140:LEU:HD21	1.93	0.50
1:B:14:VAL:HG11	1:D:59:ALA:HA	1.93	0.50
1:A:59:ALA:HA	1:C:14:VAL:HG11	1.93	0.50
1:B:69:ALA:HB2	1:B:116:LEU:HD21	1.94	0.50
1:C:60:CYS:SG	3:C:201:DTT:C4	3.00	0.50
1:B:95:PHE:CE2	3:D:201:DTT:H41	2.46	0.49
1:A:3:ILE:CG2	1:C:89:GLY:HA3	2.42	0.49
1:B:121:HIS:CD2	1:B:138:LEU:HD11	2.49	0.48
1:D:1:MET:HG3	1:D:2:SER:O	2.14	0.48
1:A:130:THR:HA	1:C:129:ALA:O	2.13	0.47
1:A:84:CYS:O	1:C:11:GLN:HA	2.14	0.47
1:B:97:LEU:HG	1:D:125:PRO:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:HIS:CG	1:B:138:LEU:HD11	2.50	0.47
1:D:25:SER:OG	2:D:202:CL:CL	2.69	0.47
1:C:121:HIS:CG	1:C:138:LEU:HD11	2.50	0.47
1:A:3:ILE:HG21	1:C:89:GLY:HA3	1.97	0.46
1:D:60:CYS:SG	3:D:201:DTT:H12	2.55	0.46
1:B:60:CYS:HG	3:B:201:DTT:HS2	1.61	0.46
1:D:108:LEU:CD1	1:D:116:LEU:HD12	2.45	0.46
1:B:129:ALA:O	1:D:130:THR:HA	2.16	0.46
1:A:100:GLU:OE1	1:A:139:ARG:HD3	2.16	0.46
1:A:60:CYS:SG	1:A:125:PRO:HG2	2.57	0.45
1:A:95:PHE:CE2	3:C:201:DTT:H12	2.51	0.45
1:D:6:ILE:HG21	1:D:9:ARG:CG	2.47	0.45
1:A:77:SER:HB3	1:A:79:ASP:OD1	2.16	0.44
1:B:0:HIS:NE2	1:D:135:ASP:OD1	2.43	0.44
1:B:38:GLU:HB3	1:D:7:LEU:O	2.18	0.44
1:B:89:GLY:HA3	1:D:3:ILE:CG2	2.48	0.44
1:B:95:PHE:CD2	3:D:201:DTT:H41	2.53	0.44
1:D:6:ILE:HG21	1:D:9:ARG:HG3	1.99	0.44
1:B:84:CYS:O	1:D:11:GLN:HA	2.18	0.43
1:D:102:GLN:HG3	4:D:325:HOH:O	2.19	0.43
1:C:6:ILE:HG21	1:C:9:ARG:CG	2.49	0.42
1:D:60:CYS:SG	1:D:125:PRO:HD2	2.59	0.42
1:B:14:VAL:HA	1:B:21:ASN:O	2.20	0.42
1:D:11:GLN:O	1:D:24:SER:HA	2.20	0.41
1:D:105:LEU:HD13	1:D:108:LEU:HD12	2.01	0.41
1:A:3:ILE:HG21	1:A:6:ILE:HD13	2.01	0.41
1:B:60:CYS:SG	1:B:125:PRO:HG2	2.60	0.41
1:B:122:ILE:CG2	1:C:26:ASP:HA	2.51	0.41
3:B:201:DTT:O3	3:B:201:DTT:S1	2.78	0.41
3:B:201:DTT:H11	1:D:95:PHE:CE2	2.56	0.41
1:C:23:GLU:HA	1:C:29:LEU:O	2.21	0.41
1:D:87:GLY:O	1:D:97:LEU:HA	2.21	0.41
1:A:129:ALA:O	1:C:130:THR:HA	2.21	0.41
1:A:7:LEU:O	1:C:38:GLU:HB3	2.21	0.41
1:A:14:VAL:HA	1:A:21:ASN:O	2.21	0.41
1:C:11:GLN:NE2	2:C:203:CL:CL	2.90	0.41
1:D:111:GLU:HA	1:D:114:ARG:NH1	2.36	0.41
1:A:88:ILE:HD12	1:C:125:PRO:HG3	2.04	0.40
1:C:108:LEU:HB3	1:C:112:GLU:HG3	2.03	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	138/161 (86%)	136 (99%)	2 (1%)	0	100	100
1	B	141/161 (88%)	137 (97%)	4 (3%)	0	100	100
1	C	138/161 (86%)	132 (96%)	6 (4%)	0	100	100
1	D	140/161 (87%)	138 (99%)	2 (1%)	0	100	100
All	All	557/644 (86%)	543 (98%)	14 (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	106/124 (86%)	105 (99%)	1 (1%)	81	79
1	B	109/124 (88%)	105 (96%)	4 (4%)	37	24
1	C	106/124 (86%)	103 (97%)	3 (3%)	47	36
1	D	108/124 (87%)	104 (96%)	4 (4%)	37	24
All	All	429/496 (86%)	417 (97%)	12 (3%)	47	36

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	TYR
1	B	5	THR

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Mol	Chain	Res	Type
1	B	57	TYR
1	B	77	SER
1	B	102	GLN
1	C	57	TYR
1	C	72	ARG
1	C	77	SER
1	D	4	GLU
1	D	19	GLU
1	D	57	TYR
1	D	77	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 10 ligands modelled in this entry, 7 are monoatomic - leaving 3 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	DTT	B	201	-	7,7,7	0.17	0	4,8,8	1.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	DTT	C	201	-	7,7,7	0.67	0	4,8,8	0.99	0
3	DTT	D	201	-	7,7,7	0.35	0	4,8,8	1.31	1 (25%)

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	DTT	B	201	-	-	4/8/8/8	-
3	DTT	C	201	-	-	5/8/8/8	-
3	DTT	D	201	-	-	3/8/8/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	201	DTT	C3-C4-S4	-2.16	107.92	114.45

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	201	DTT	C1-C2-C3-O3
3	B	201	DTT	C1-C2-C3-C4
3	B	201	DTT	O2-C2-C3-O3
3	B	201	DTT	O2-C2-C3-C4
3	C	201	DTT	S1-C1-C2-C3
3	C	201	DTT	C2-C3-C4-S4
3	C	201	DTT	O3-C3-C4-S4
3	D	201	DTT	S1-C1-C2-O2
3	D	201	DTT	S1-C1-C2-C3
3	C	201	DTT	S1-C1-C2-O2
3	C	201	DTT	O2-C2-C3-O3
3	D	201	DTT	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	201	DTT	3	0
3	C	201	DTT	2	0
3	D	201	DTT	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	140/161 (86%)	0.42	9 (6%) 19 21	23, 40, 69, 88	0
1	B	143/161 (88%)	0.57	14 (9%) 7 8	27, 39, 69, 85	0
1	C	140/161 (86%)	0.34	9 (6%) 19 21	24, 37, 69, 92	0
1	D	142/161 (88%)	0.64	15 (10%) 6 7	27, 43, 81, 103	0
All	All	565/644 (87%)	0.49	47 (8%) 11 13	23, 40, 75, 103	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	16	GLY	8.1
1	D	18	ARG	7.4
1	A	18	ARG	6.6
1	A	16	GLY	6.4
1	C	18	ARG	6.2
1	D	1	MET	5.2
1	D	19	GLU	5.2
1	B	75	THR	5.1
1	D	17	GLY	4.6
1	A	2	SER	4.5
1	B	18	ARG	4.2
1	A	4	GLU	4.1
1	B	4	GLU	4.1
1	C	73	LYS	4.1
1	A	19	GLU	4.0
1	B	16	GLY	4.0
1	D	4	GLU	3.9
1	B	1	MET	3.8
1	D	42	ALA	3.4
1	D	70	ALA	3.2
1	B	-1	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	92	PRO	2.9
1	B	76	LEU	2.8
1	B	19	GLU	2.8
1	D	114	ARG	2.8
1	C	17	GLY	2.7
1	B	57	TYR	2.7
1	D	116	LEU	2.6
1	A	132	GLY	2.5
1	C	72	ARG	2.5
1	A	5	THR	2.5
1	B	126	TYR	2.5
1	D	72	ARG	2.5
1	C	93	SER	2.5
1	D	71	LYS	2.4
1	B	55	ALA	2.4
1	D	73	LYS	2.4
1	D	111	GLU	2.3
1	A	134	ILE	2.2
1	B	5	THR	2.2
1	B	52	LEU	2.1
1	C	16	GLY	2.1
1	A	141	ALA	2.1
1	B	73	LYS	2.1
1	C	75	THR	2.1
1	C	2	SER	2.1
1	D	119	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

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, 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	B-factors(\AA^2)	Q<0.9
3	DTT	B	201	8/8	0.72	0.21	73,77,79,81	0
3	DTT	D	201	8/8	0.82	0.14	56,70,74,79	0
3	DTT	C	201	8/8	0.86	0.23	52,61,67,70	0
2	CL	A	201	1/1	0.95	0.07	39,39,39,39	0
2	CL	B	202	1/1	0.96	0.10	51,51,51,51	0
2	CL	C	202	1/1	0.97	0.05	45,45,45,45	0
2	CL	B	203	1/1	0.98	0.04	55,55,55,55	0
2	CL	C	203	1/1	0.98	0.06	45,45,45,45	0
2	CL	D	202	1/1	0.99	0.06	41,41,41,41	0
2	CL	A	202	1/1	0.99	0.06	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.