



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 07:38 pm BST

PDB ID : 5ZXN  
Title : Crystal structure of CurA from *Vibrio vulnificus*  
Authors : Kim, M.-K.; Bae, D.-W.; Cha, S.-S.  
Deposited on : 2018-05-21  
Resolution : 1.85 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

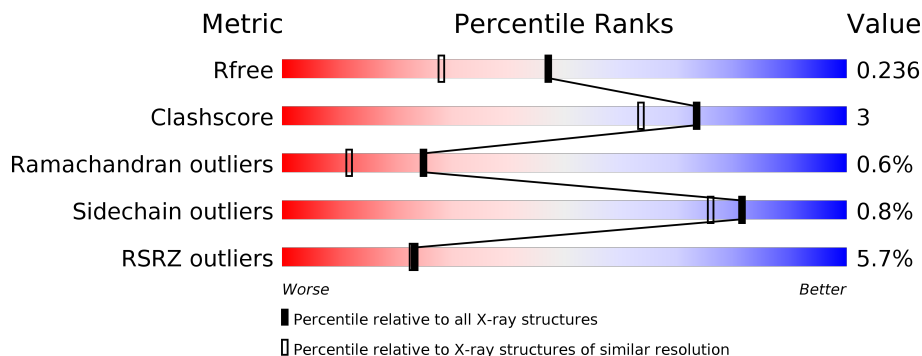
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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  $\geq 3$ , 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	340	
1	B	340	

## 2 Entry composition [i](#)

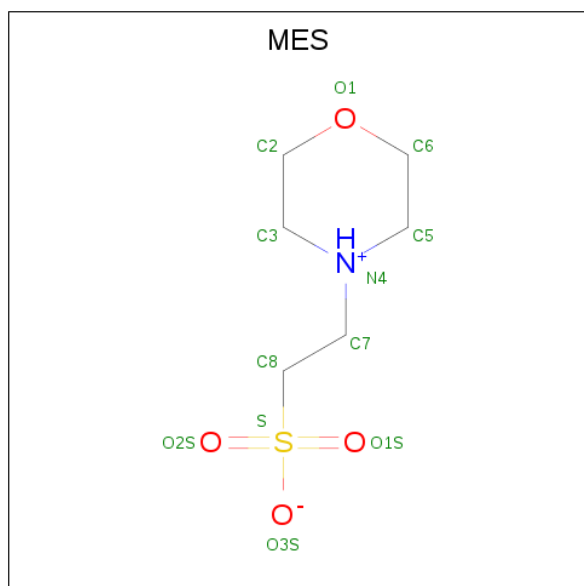
There are 4 unique types of molecules in this entry. The entry contains 5509 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NADP-dependent oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	S				Se
1	A	340	Total	C	N	O	S	Se	0	0	0
			2532	1605	431	475	6	15			
1	B	339	Total	C	N	O	S	Se	0	0	0
			2525	1600	430	474	6	15			

- Molecule 2 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
2	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	B	1	12	6	1	4	1	0	0

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	4	2	2	0	0

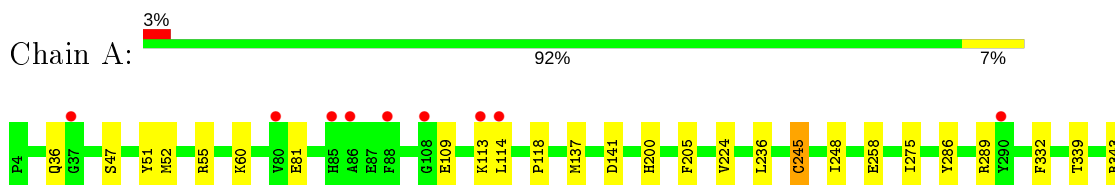
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	218	218	218	0	0
4	B	170	170	170	0	0

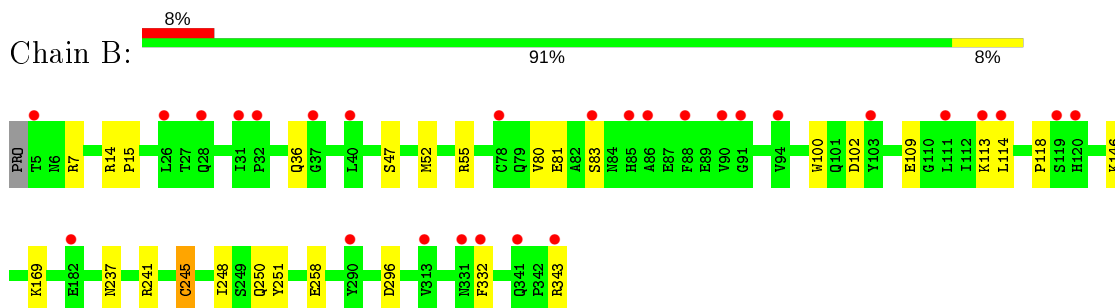
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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 ( $RSRZ > 2$ ). 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: NADP-dependent oxidoreductase



- Molecule 1: NADP-dependent oxidoreductase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.52Å 91.56Å 104.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 1.85 19.92 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.4 (19.92-1.85) 99.4 (19.92-1.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.03 (at 1.85Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.209 , 0.236 0.209 , 0.236	Depositor DCC
$R_{free}$ test set	3726 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.8	Xtrriage
Anisotropy	0.058	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 33.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.057 for k,h,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5509	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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: EDO, MES

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	1/2570 (0.0%)	0.54	0/3454
1	B	0.39	1/2562 (0.0%)	0.53	0/3443
All	All	0.39	2/5132 (0.0%)	0.54	0/6897

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	245	CYS	CB-SG	-8.09	1.68	1.82
1	A	245	CYS	CB-SG	-6.31	1.71	1.82

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	2532	0	2464	15	0
1	B	2525	0	2456	18	0
2	A	24	0	24	1	0
2	B	36	0	36	2	0
3	A	4	0	6	0	0
4	A	218	0	0	4	0
4	B	170	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5509	0	4986	33	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 3.

All (33) 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:146:LYS:NZ	4:B:502:HOH:O	2.16	0.78
1:B:55:ARG:NH1	4:B:506:HOH:O	2.27	0.67
1:A:224:VAL:HA	1:A:248:ILE:HG22	1.77	0.66
1:B:258:GLU:OE2	4:B:501:HOH:O	2.14	0.65
1:B:114:LEU:H	1:B:118:PRO:HB3	1.61	0.63
1:B:169:LYS:NZ	4:B:508:HOH:O	2.30	0.62
1:A:60:LYS:NZ	4:A:502:HOH:O	2.19	0.62
1:A:114:LEU:H	1:A:118:PRO:HB3	1.64	0.61
1:A:258:GLU:OE1	4:A:501:HOH:O	2.17	0.58
1:A:55:ARG:NH1	4:A:509:HOH:O	2.38	0.57
1:A:51:TYR:HB3	2:A:401:MES:H72	1.86	0.56
1:A:47:SER:O	1:A:52:MSE:HE1	2.09	0.52
1:B:248:ILE:HA	1:B:251:TYR:CD2	2.46	0.51
1:B:47:SER:O	1:B:52:MSE:HE1	2.12	0.49
1:B:113:LYS:HA	1:B:114:LEU:HA	1.58	0.48
1:B:80:VAL:HG11	1:B:83:SER:HB2	1.96	0.47
1:B:7:ARG:NE	1:B:102:ASP:OD2	2.32	0.46
1:B:250:GLN:HB3	4:B:636:HOH:O	2.15	0.46
1:A:339:THR:O	4:A:503:HOH:O	2.21	0.46
1:A:81:GLU:CD	1:A:343:ARG:HE	2.20	0.45
1:A:113:LYS:HA	1:A:114:LEU:HA	1.51	0.45
1:B:237:ASN:OD1	2:B:402:MES:H32	2.16	0.45
1:B:258:GLU:OE1	4:B:503:HOH:O	2.21	0.45
1:B:296:ASP:OD2	2:B:403:MES:H21	2.18	0.44
1:A:236:LEU:HD13	1:A:275:ILE:HD13	1.99	0.43
1:B:81:GLU:CD	1:B:343:ARG:HE	2.22	0.43
1:A:141:ASP:OD2	1:A:289:ARG:NH1	2.43	0.42
1:A:200:HIS:HA	1:A:205:PHE:CD1	2.54	0.42
1:A:137:MSE:HE2	1:A:286:TYR:CD1	2.54	0.42
1:B:100:TRP:N	1:B:100:TRP:CD1	2.87	0.42
1:B:241:ARG:HD3	1:B:241:ARG:N	2.36	0.41
1:A:137:MSE:O	1:A:141:ASP:HB2	2.21	0.41
1:B:14:ARG:HA	1:B:15:PRO:HD3	1.97	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	338/340 (99%)	327 (97%)	9 (3%)	2 (1%)	25	12
1	B	337/340 (99%)	324 (96%)	11 (3%)	2 (1%)	25	12
All	All	675/680 (99%)	651 (96%)	20 (3%)	4 (1%)	25	12

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	36	GLN
1	A	36	GLN
1	A	109	GLU
1	B	109	GLU

### 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	255/253 (101%)	253 (99%)	2 (1%)	81	76
1	B	254/253 (100%)	252 (99%)	2 (1%)	81	76
All	All	509/506 (101%)	505 (99%)	4 (1%)	81	76

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	245	CYS
1	A	332	PHE
1	B	245	CYS
1	B	332	PHE

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](#)

6 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	MES	A	402	-	12,12,12	2.02	1 (8%)	14,16,16	1.74	5 (35%)
2	MES	B	401	-	12,12,12	1.89	1 (8%)	14,16,16	2.01	4 (28%)
2	MES	B	403	-	12,12,12	2.25	1 (8%)	14,16,16	1.92	2 (14%)
3	EDO	A	403	-	3,3,3	0.56	0	2,2,2	0.15	0
2	MES	A	401	-	12,12,12	1.87	1 (8%)	14,16,16	2.31	6 (42%)
2	MES	B	402	-	12,12,12	2.21	1 (8%)	14,16,16	2.74	7 (50%)

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	MES	A	402	-	-	2/6/14/14	0/1/1/1
2	MES	B	401	-	-	0/6/14/14	0/1/1/1
2	MES	B	403	-	-	2/6/14/14	0/1/1/1
3	EDO	A	403	-	-	1/1/1/1	-
2	MES	A	401	-	-	1/6/14/14	0/1/1/1
2	MES	B	402	-	-	1/6/14/14	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	403	MES	C8-S	-7.52	1.66	1.77
2	B	402	MES	C8-S	-7.35	1.67	1.77
2	A	402	MES	C8-S	-6.59	1.68	1.77
2	B	401	MES	C8-S	-6.23	1.68	1.77
2	A	401	MES	C8-S	-6.13	1.68	1.77

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	MES	C5-N4-C3	7.24	125.14	108.83
2	B	401	MES	C5-N4-C3	5.35	120.86	108.83
2	B	403	MES	C5-N4-C3	5.24	120.63	108.83
2	A	401	MES	C5-N4-C3	4.94	119.96	108.83
2	B	402	MES	O1S-S-C8	4.31	112.11	106.92
2	A	401	MES	C7-N4-C5	4.30	122.24	111.23
2	B	401	MES	O1S-S-C8	3.63	111.28	106.92
2	A	402	MES	C5-N4-C3	3.54	116.80	108.83
2	B	403	MES	O3S-S-C8	3.31	111.12	105.77
2	B	402	MES	O2S-S-C8	2.87	110.37	106.92
2	A	402	MES	C7-N4-C5	2.59	117.85	111.23
2	A	401	MES	C6-C5-N4	-2.59	106.18	110.10
2	B	402	MES	C7-N4-C3	2.59	117.85	111.23
2	A	401	MES	O2S-S-C8	2.58	110.02	106.92
2	A	401	MES	C7-N4-C3	2.57	117.80	111.23
2	A	402	MES	O1S-S-C8	2.53	109.97	106.92
2	B	402	MES	C2-C3-N4	2.37	113.69	110.10
2	B	402	MES	O3S-S-O1S	-2.23	105.82	111.27
2	A	401	MES	O3S-S-C8	2.21	109.34	105.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	MES	O3S-S-C8	2.21	109.33	105.77
2	A	402	MES	O3S-S-C8	2.15	109.24	105.77
2	B	402	MES	C7-N4-C5	2.08	116.56	111.23
2	A	402	MES	O2S-S-C8	2.02	109.34	106.92
2	B	401	MES	C7-N4-C3	2.01	116.37	111.23

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	MES	C8-C7-N4-C5
2	A	401	MES	C8-C7-N4-C5
2	B	402	MES	C8-C7-N4-C3
2	B	403	MES	N4-C7-C8-S
2	B	403	MES	C8-C7-N4-C3
3	A	403	EDO	O1-C1-C2-O2
2	A	402	MES	C7-C8-S-O2S

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	403	MES	1	0
2	A	401	MES	1	0
2	B	402	MES	1	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, 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	325/340 (95%)	0.01	9 (2%) 53 52	18, 35, 69, 83	0
1	B	324/340 (95%)	0.36	28 (8%) 10 10	18, 42, 83, 95	0
All	All	649/680 (95%)	0.18	37 (5%) 23 23	18, 38, 80, 95	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	LEU	12.8
1	A	114	LEU	7.0
1	B	332	PHE	6.3
1	B	86	ALA	4.6
1	B	40	LEU	4.5
1	A	85	HIS	4.5
1	B	341	GLN	4.2
1	A	108	GLY	4.0
1	B	290	TYR	3.7
1	A	290	TYR	3.7
1	B	119	SER	3.6
1	B	31	ILE	3.6
1	A	86	ALA	3.5
1	B	88	PHE	3.5
1	B	182	GLU	3.5
1	B	37	GLY	3.4
1	B	5	THR	3.4
1	B	26	LEU	3.4
1	B	113	LYS	3.0
1	B	83	SER	3.0
1	B	103	TYR	2.9
1	B	91	GLY	2.9
1	B	343	ARG	2.9
1	B	313	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	37	GLY	2.7
1	B	78	CYS	2.7
1	B	28	GLN	2.6
1	B	85	HIS	2.6
1	B	331	ASN	2.6
1	A	113	LYS	2.4
1	A	88	PHE	2.4
1	B	111	LEU	2.3
1	B	90	VAL	2.3
1	B	120	HIS	2.2
1	B	32	PRO	2.2
1	A	80	VAL	2.0
1	B	94	VAL	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, 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
3	EDO	A	403	4/4	0.76	0.14	41,43,43,46	0
2	MES	B	402	12/12	0.78	0.29	46,59,66,76	0
2	MES	B	403	12/12	0.87	0.19	48,56,72,81	0
2	MES	A	402	12/12	0.93	0.20	33,43,56,56	0
2	MES	A	401	12/12	0.98	0.08	31,32,33,38	0
2	MES	B	401	12/12	0.98	0.09	29,32,34,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.