



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:02 AM UTC

PDB ID : 2W53 / pdb_00002w53
Title : Structure of SmeT, the repressor of the *Stenotrophomonas maltophilia* multidrug efflux pump SmeDEF.
Authors : Mate, M.J.; Romero, A.; Hernandez, A.; Martinez, J.L.
Deposited on : 2008-12-04
Resolution : 2.00 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

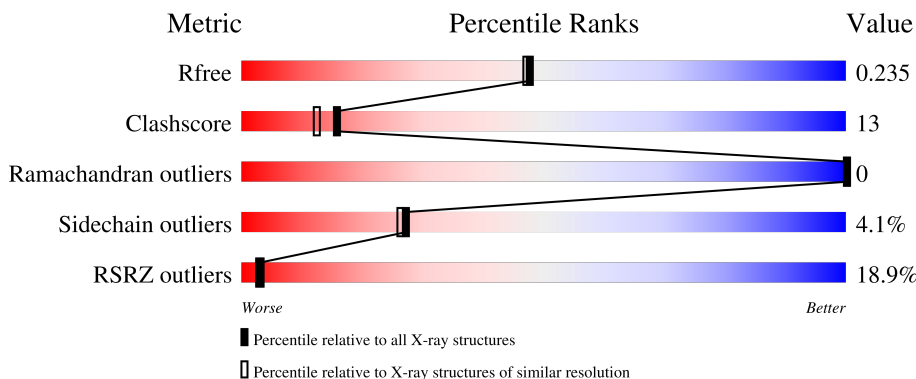
1 Overall quality at a glance

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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	219	
1	B	219	

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	SO4	A	1218	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3242 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.

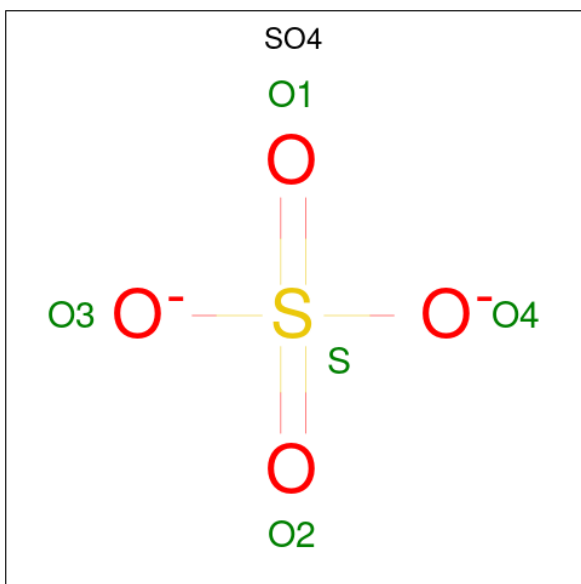
- Molecule 1 is a protein called REPRESSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	202	1615	1011	292	296	16	0	5	0
1	B	173	1348	848	243	245	12	0	5	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	LEU	GLN	engineered mutation	UNP Q8KLP4
B	166	LEU	GLN	engineered mutation	UNP Q8KLP4

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

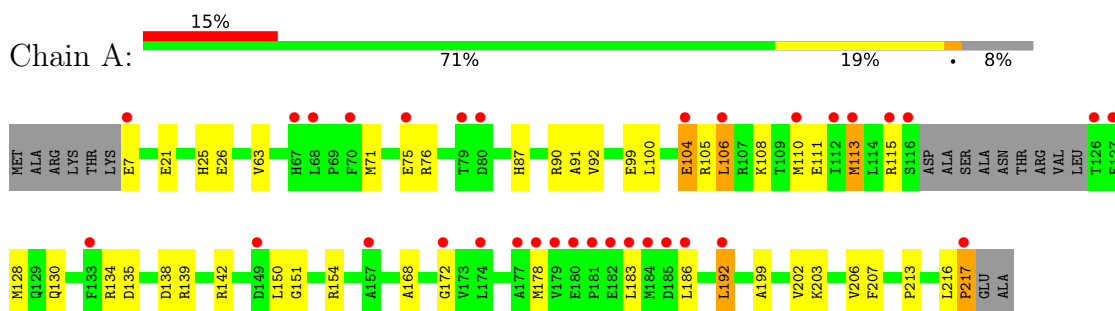
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	135	Total	O	0	0
			135	135		
3	B	134	Total	O	0	0
			134	134		

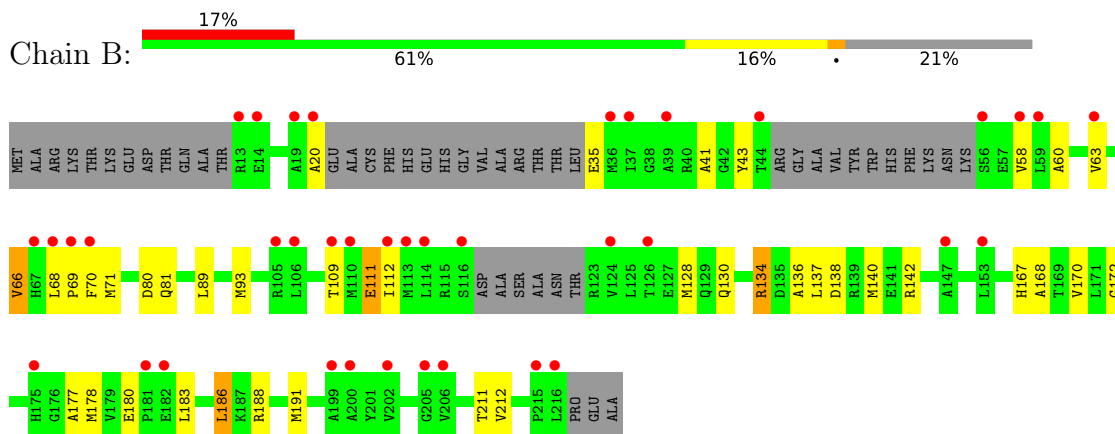
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: REPRESSOR



- Molecule 1: REPRESSOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.70Å 58.58Å 83.26Å 90.00° 103.14° 90.00°	Depositor
Resolution (Å)	19.86 – 2.00 19.86 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.0 (19.86-2.00) 96.8 (19.86-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.19 (at 2.01Å)	Xtrriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.194 , 0.236 0.194 , 0.235	Depositor DCC
R_{free} test set	1760 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtrriage
Anisotropy	0.095	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3242	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.98% 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: SO4

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	0.84	0/1655	0.98	1/2227 (0.0%)
1	B	0.94	0/1379	0.96	0/1855
All	All	0.89	0/3034	0.97	1/4082 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	217	PRO	N-CA-CB	6.15	109.76	103.00

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	1615	0	1619	43	0
1	B	1348	0	1346	41	0
2	A	5	0	0	2	0
2	B	5	0	0	0	0
3	A	135	0	0	1	0
3	B	134	0	0	4	0
All	All	3242	0	2965	79	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 13.

All (79) 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:115:ARG:HB3	1:A:178:MET:HE1	1.40	0.99
1:B:93[B]:MET:CE	1:B:170:VAL:HG21	1.96	0.96
1:A:111:GLU:O	1:A:115:ARG:HG3	1.74	0.88
1:A:71:MET:CE	1:A:139:ARG:HD2	2.05	0.86
1:A:154:ARG:NH1	1:A:199:ALA:O	2.10	0.83
1:A:104:GLU:CD	1:A:104:GLU:H	1.86	0.82
1:A:75:GLU:CD	1:A:142:ARG:HH22	1.89	0.80
1:B:35:GLU:OE2	1:B:35:GLU:HA	1.83	0.79
1:B:68:LEU:HB3	1:B:69:PRO:HD3	1.67	0.77
1:A:75:GLU:OE2	1:A:142:ARG:NH2	2.18	0.76
1:B:93[B]:MET:HE2	1:B:170:VAL:HG21	1.70	0.73
1:B:191:MET:HE3	3:B:2107:HOH:O	1.88	0.73
1:A:25:HIS:CD2	1:A:108:LYS:HG3	2.24	0.72
1:B:63:VAL:HG23	1:B:128:MET:HG2	1.70	0.72
1:B:93[B]:MET:CE	1:B:170:VAL:CG2	2.69	0.71
1:B:130:GLN:O	1:B:134[A]:ARG:HD3	1.90	0.71
1:B:140:MET:HE1	1:B:167:HIS:ND1	2.08	0.69
1:A:71:MET:HE3	1:A:139:ARG:CD	2.23	0.68
1:A:71:MET:HE2	1:A:135:ASP:HB3	1.75	0.68
1:B:138:ASP:OD2	1:B:142:ARG:NH2	2.24	0.68
1:B:93[B]:MET:HE2	1:B:170:VAL:CG2	2.25	0.67
1:A:87:HIS:NE2	1:A:206:VAL:O	2.29	0.65
1:B:93[B]:MET:HE1	1:B:170:VAL:HG21	1.78	0.65
1:B:63:VAL:CG2	1:B:128:MET:HG2	2.28	0.64
1:B:211:THR:OG1	3:B:2129:HOH:O	2.08	0.63
1:B:111:GLU:HG3	1:B:178:MET:SD	2.39	0.62
1:A:71:MET:CE	1:A:139:ARG:CD	2.76	0.62
1:B:80:ASP:OD1	1:B:81:GLN:NE2	2.29	0.61
1:A:71:MET:HE1	1:A:139:ARG:HD2	1.81	0.61
1:B:89:LEU:O	1:B:93[B]:MET:HG3	2.03	0.58
1:A:130:GLN:O	1:A:134:ARG:HG3	2.04	0.58
1:B:191:MET:HE1	1:B:212:VAL:HG22	1.86	0.58
1:B:68:LEU:HB3	1:B:69:PRO:CD	2.34	0.56
1:B:71:MET:HE1	1:B:136:ALA:HB2	1.86	0.56
1:A:76:ARG:HH22	1:A:217:PRO:C	2.14	0.56
1:B:177:ALA:CB	1:B:186:LEU:HD13	2.36	0.56
1:A:115:ARG:HG2	1:A:178:MET:SD	2.46	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:GLU:CD	1:A:142:ARG:NH2	2.60	0.55
1:A:99:GLU:HG2	1:A:106:LEU:HD22	1.90	0.52
1:A:110[A]:MET:HA	1:A:110[A]:MET:HE2	1.91	0.52
1:A:192:LEU:HD23	3:A:2060:HOH:O	2.09	0.52
1:B:191:MET:HE1	1:B:212:VAL:CG2	2.41	0.51
1:A:115:ARG:HB3	1:A:178:MET:CE	2.27	0.50
1:B:93[B]:MET:HE1	1:B:170:VAL:CG2	2.38	0.50
1:B:180:GLU:OE1	1:B:183:LEU:HG	2.12	0.50
1:A:154:ARG:HB2	2:A:1218:SO4:O1	2.12	0.50
1:B:20:ALA:HB1	1:B:58:VAL:HG11	1.94	0.49
1:B:109:THR:O	1:B:112:ILE:HG12	2.13	0.49
1:A:21:GLU:OE2	1:A:105:ARG:NH1	2.46	0.48
1:A:92:VAL:HG23	1:A:216:LEU:HD22	1.95	0.48
1:B:178:MET:HE2	3:B:2102:HOH:O	2.12	0.48
1:A:100:LEU:CD2	1:A:110[A]:MET:HG3	2.44	0.48
1:A:25:HIS:CG	1:A:108:LYS:HG3	2.48	0.48
1:A:63:VAL:HG13	1:A:113[B]:MET:HE2	1.94	0.47
1:A:25:HIS:CG	1:A:108:LYS:CG	2.97	0.47
1:A:134:ARG:HG2	1:B:183:LEU:HD11	1.94	0.47
1:B:60:ALA:HA	1:B:63:VAL:HG22	1.97	0.46
1:B:137:LEU:HA	1:B:140:MET:HE3	1.98	0.46
1:A:138:ASP:OD2	1:A:142:ARG:NH1	2.48	0.46
1:A:183:LEU:HD11	1:B:134[A]:ARG:HG3	1.99	0.45
1:B:70[B]:PHE:HD2	1:B:71:MET:HE2	1.81	0.45
1:A:154:ARG:HB2	2:A:1218:SO4:S	2.56	0.44
1:A:168:ALA:O	1:B:172:GLY:HA3	2.17	0.44
1:A:105:ARG:HG2	1:A:106:LEU:N	2.30	0.44
1:A:151:GLY:O	1:A:203:LYS:NZ	2.49	0.44
1:B:20:ALA:HB1	1:B:58:VAL:CG1	2.48	0.44
1:A:90:ARG:HD2	1:A:207:PHE:CD1	2.53	0.44
1:A:172:GLY:HA3	1:B:168:ALA:O	2.17	0.44
1:B:70[B]:PHE:CD2	1:B:71:MET:HE2	2.53	0.44
1:A:91:ALA:HA	1:A:213:PRO:HG2	2.00	0.43
1:A:92:VAL:CG2	1:A:216:LEU:HD22	2.49	0.43
1:A:128:MET:HE2	1:A:128:MET:HB3	1.66	0.43
1:B:35:GLU:OE2	1:B:35:GLU:CA	2.58	0.41
1:B:66:VAL:HG21	1:B:109:THR:HG21	2.03	0.41
1:A:154:ARG:NH1	1:A:202:VAL:O	2.53	0.41
1:A:183:LEU:CD1	1:B:134[A]:ARG:HG3	2.52	0.40
1:B:188:ARG:HD2	3:B:2105:HOH:O	2.21	0.40
1:B:41:ALA:HB1	1:B:43:TYR:CD2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:LEU:HD21	1:A:110[A]:MET:HG3	2.02	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	203/219 (93%)	202 (100%)	1 (0%)	0	100	100
1	B	170/219 (78%)	169 (99%)	1 (1%)	0	100	100
All	All	373/438 (85%)	371 (100%)	2 (0%)	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	168/179 (94%)	159 (95%)	9 (5%)	20	17
1	B	136/179 (76%)	131 (96%)	5 (4%)	30	30
All	All	304/358 (85%)	290 (95%)	14 (5%)	27	22

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

Mol	Chain	Res	Type
1	A	7	GLU
1	A	26	GLU
1	A	104	GLU
1	A	106	LEU
1	A	113[A]	MET
1	A	113[B]	MET
1	A	150	LEU
1	A	186	LEU
1	A	192	LEU
1	B	66	VAL
1	B	111	GLU
1	B	134[A]	ARG
1	B	134[B]	ARG
1	B	186	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	SO4	B	1217	-	4,4,4	0.33	0	6,6,6	0.50	0
2	SO4	A	1218	-	4,4,4	0.26	0	6,6,6	0.28	0

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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1218	SO4	2	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	202/219 (92%)	1.06	33 (16%) 4 4	25, 39, 47, 62	5 (2%)
1	B	173/219 (78%)	1.41	38 (21%) 2 2	22, 39, 57, 69	5 (2%)
All	All	375/438 (85%)	1.22	71 (18%) 3 3	22, 39, 52, 69	10 (2%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	116	SER	6.8
1	A	126	THR	5.7
1	A	116	SER	5.5
1	B	216	LEU	4.0
1	B	44	THR	3.9
1	A	104	GLU	3.6
1	B	182	GLU	3.6
1	A	217	PRO	3.5
1	A	192	LEU	3.4
1	B	126	THR	3.4
1	A	133	PHE	3.4
1	B	19	ALA	3.1
1	B	13	ARG	3.1
1	B	124	VAL	3.1
1	B	70[A]	PHE	3.1
1	B	215	PRO	3.1
1	B	181	PRO	3.0
1	B	63	VAL	3.0
1	B	58	VAL	3.0
1	A	178	MET	2.9
1	B	175	HIS	2.9
1	A	182	GLU	2.9
1	A	186	LEU	2.8
1	A	181	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	149	ASP	2.8
1	B	36	MET	2.7
1	B	67	HIS	2.7
1	B	39	ALA	2.7
1	B	105	ARG	2.7
1	B	109	THR	2.7
1	A	177	ALA	2.7
1	A	183	LEU	2.6
1	B	59	LEU	2.6
1	B	113	MET	2.6
1	A	157	ALA	2.6
1	B	20	ALA	2.6
1	B	147	ALA	2.6
1	B	68	LEU	2.6
1	B	200	ALA	2.6
1	B	56	SER	2.6
1	A	110[A]	MET	2.5
1	B	153	LEU	2.5
1	B	112	ILE	2.5
1	A	106	LEU	2.5
1	B	206	VAL	2.5
1	A	70[A]	PHE	2.5
1	A	115	ARG	2.5
1	A	112	ILE	2.5
1	B	202	VAL	2.4
1	B	106	LEU	2.4
1	A	80	ASP	2.4
1	A	7	GLU	2.4
1	A	184	MET	2.4
1	B	114	LEU	2.4
1	A	67	HIS	2.3
1	A	127	GLU	2.3
1	B	110[A]	MET	2.3
1	A	179	VAL	2.2
1	B	69	PRO	2.2
1	A	75	GLU	2.2
1	B	37	ILE	2.2
1	A	113[A]	MET	2.2
1	A	172	GLY	2.2
1	B	205	GLY	2.1
1	B	14	GLU	2.1
1	B	199	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	180	GLU	2.1
1	A	174	LEU	2.0
1	A	79	THR	2.0
1	A	68	LEU	2.0
1	A	185	ASP	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 oligosaccharides 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
2	SO4	A	1218	5/5	0.75	0.14	82,83,83,84	0
2	SO4	B	1217	5/5	0.92	0.19	52,54,56,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.