Full wwPDB X-ray Structure Validation Report (i)

Jan 4, 2024 – 08:01 am GMT

PDB ID	:	4UNM
Title	:	Structure of Galactose Oxidase homologue from Streptomyces lividans
Authors	:	Chaplin, A.K.; Hough, M.A.; Worrall, J.A.R.
Deposited on	:	2014-05-29
Resolution	:	1.77 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 (i) 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 (i)) were used in the production of this report:

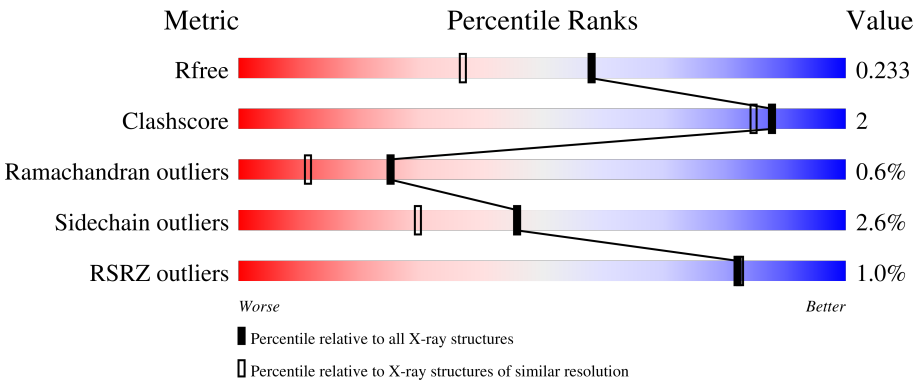
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.77 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 <=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	615	91% 6% • •
1	B	615	% 90% 7% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10084 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 SECRETED PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	599	4663	2938	811	903	11	50	3	0
1	B	609	4725	2979	820	915	11	79	1	0

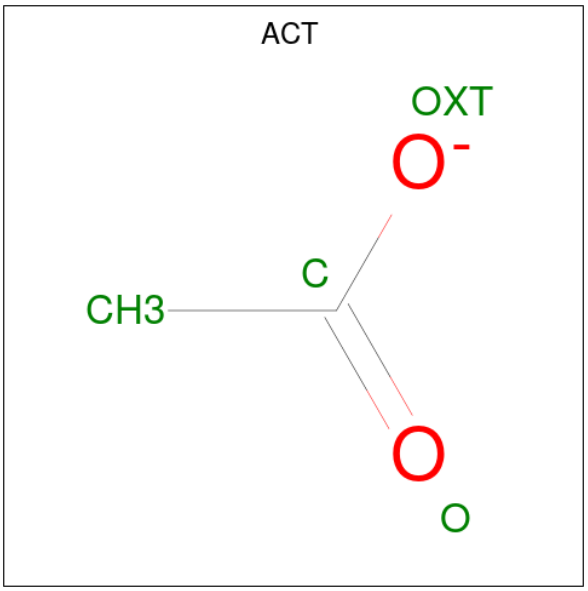
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	GLY	-	expression tag	UNP D6EWM0
A	32	SER	-	expression tag	UNP D6EWM0
A	33	HIS	-	expression tag	UNP D6EWM0
A	34	MET	-	expression tag	UNP D6EWM0
B	31	GLY	-	expression tag	UNP D6EWM0
B	32	SER	-	expression tag	UNP D6EWM0
B	33	HIS	-	expression tag	UNP D6EWM0
B	34	MET	-	expression tag	UNP D6EWM0

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

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

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 4 is water.

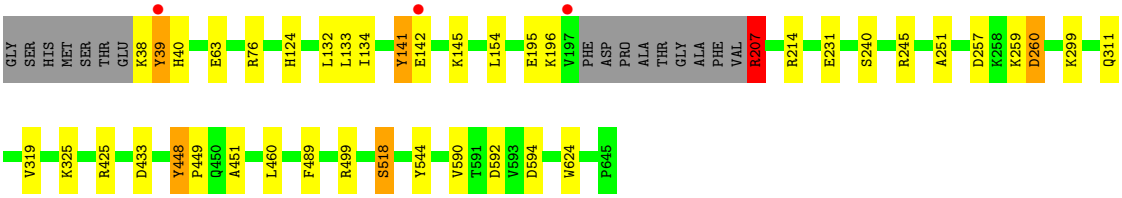
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	417	417	417	0	0
			Total	O		
4	B	261	261	261	0	0

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.

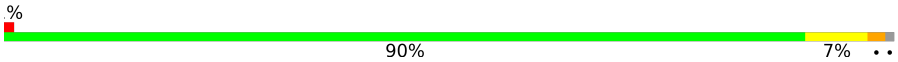
- Molecule 1: SECRETED PROTEIN

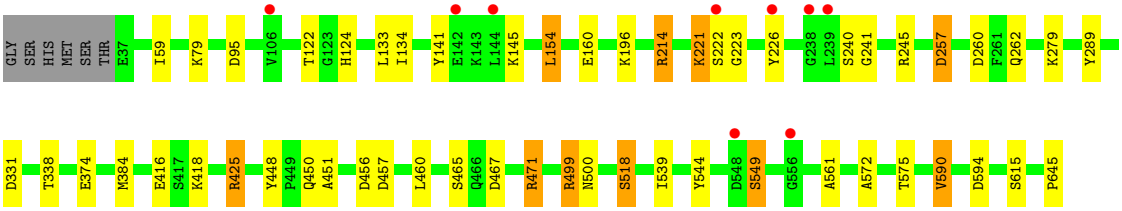
Chain A:



- Molecule 1: SECRETED PROTEIN

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	50.36Å 126.63Å 107.55Å 90.00° 91.10° 90.00°	Depositor
Resolution (Å)	53.50 – 1.77 63.31 – 1.77	Depositor EDS
% Data completeness (in resolution range)	99.7 (53.50-1.77) 99.7 (63.31-1.77)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.17 (at 1.77Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.188 , 0.228 0.195 , 0.233	Depositor DCC
R_{free} test set	6563 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	23.1	Xtriage
Anisotropy	0.604	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 41.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10084	wwPDB-VP
Average B, all atoms (Å^2)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: CU, ACT

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	3/4777 (0.1%)	0.82	6/6468 (0.1%)
1	B	0.77	7/4837 (0.1%)	0.95	16/6555 (0.2%)
All	All	0.73	10/9614 (0.1%)	0.89	22/13023 (0.2%)

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	4
1	B	0	4
All	All	0	8

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	221	LYS	CD-CE	25.56	2.15	1.51
1	B	245	ARG	CZ-NH1	-11.45	1.18	1.33
1	B	79	LYS	CE-NZ	-10.23	1.23	1.49
1	B	214	ARG	CZ-NH1	9.80	1.45	1.33
1	A	259	LYS	CE-NZ	9.79	1.73	1.49
1	B	549	SER	CB-OG	8.87	1.53	1.42
1	A	245	ARG	CZ-NH1	-8.15	1.22	1.33
1	B	374	GLU	CB-CG	-7.70	1.37	1.52
1	A	63	GLU	CD-OE1	-6.71	1.18	1.25
1	B	154	LEU	CG-CD1	-5.66	1.30	1.51

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	214	ARG	NE-CZ-NH1	-25.60	107.50	120.30
1	B	221	LYS	CD-CE-NZ	-22.44	60.08	111.70
1	B	221	LYS	CG-CD-CE	-18.89	55.22	111.90
1	B	425	ARG	NE-CZ-NH1	15.37	127.99	120.30
1	B	425	ARG	NE-CZ-NH2	-12.34	114.13	120.30
1	A	245	ARG	NE-CZ-NH1	11.33	125.96	120.30
1	B	499[A]	ARG	NE-CZ-NH1	-9.19	115.70	120.30
1	B	499[B]	ARG	NE-CZ-NH1	-9.19	115.70	120.30
1	A	433	ASP	CB-CG-OD2	9.17	126.55	118.30
1	B	154	LEU	CB-CG-CD2	-8.87	95.93	111.00
1	A	245	ARG	NH1-CZ-NH2	-8.00	110.59	119.40
1	B	499[A]	ARG	NE-CZ-NH2	6.89	123.74	120.30
1	B	499[B]	ARG	NE-CZ-NH2	6.89	123.74	120.30
1	B	214	ARG	NH1-CZ-NH2	-6.01	112.79	119.40
1	B	425	ARG	CD-NE-CZ	5.87	131.82	123.60
1	A	592	ASP	CB-CG-OD2	-5.72	113.15	118.30
1	B	95	ASP	CB-CG-OD1	5.72	123.45	118.30
1	B	471	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	A	141	TYR	CA-CB-CG	5.67	124.17	113.40
1	B	154	LEU	CD1-CG-CD2	5.53	127.08	110.50
1	B	331	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	592	ASP	CB-CG-OD1	5.03	122.82	118.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	195	GLU	Peptide
1	A	207	ARG	Peptide
1	A	257	ASP	Peptide
1	A	39	TYR	Peptide
1	B	214	ARG	Sidechain
1	B	223	GLY	Peptide
1	B	241	GLY	Peptide
1	B	257	ASP	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	4663	0	4608	15	0
1	B	4725	0	4656	20	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	12	0	9	0	0
3	B	4	0	3	1	0
4	A	417	0	0	2	0
4	B	261	0	0	2	0
All	All	10084	0	9276	36	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 2.

All (36) 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)
3:B:701:ACT:O	4:B:2220:HOH:O	1.63	1.13
1:B:384:MET:SD	1:B:425:ARG:HD2	2.29	0.73
1:A:325:LYS:NZ	4:A:2166:HOH:O	2.28	0.64
1:B:467:ASP:OD2	1:B:471:ARG:HD2	2.02	0.58
1:A:142:GLU:HB2	1:A:260:ASP:OD1	2.05	0.57
1:B:59:ILE:HD11	1:B:539:ILE:HD12	1.87	0.56
1:B:124:HIS:HA	1:B:133:LEU:O	2.08	0.54
1:B:572:ALA:O	1:B:575:THR:HG22	2.09	0.53
1:A:132:LEU:HD23	1:A:134:ILE:HD11	1.90	0.52
1:B:59:ILE:HD11	1:B:539:ILE:CD1	2.41	0.50
1:A:299:LYS:HB3	1:A:319:VAL:CG1	2.41	0.50
1:A:196:LYS:O	1:A:207:ARG:O	2.32	0.48
1:B:221:LYS:HB3	1:B:226:TYR:CE2	2.50	0.47
1:B:450:GLN:OE1	1:B:499[A]:ARG:NH1	2.48	0.46
1:B:499[B]:ARG:HA	1:B:518:SER:HB2	1.98	0.46
1:A:124:HIS:HA	1:A:133:LEU:O	2.16	0.46
1:B:499[A]:ARG:HA	1:B:518:SER:HB2	1.98	0.46
1:A:499[B]:ARG:HA	1:A:518:SER:HB2	1.99	0.45
1:A:311:GLN:NE2	4:A:2188:HOH:O	2.40	0.45
1:B:59:ILE:CD1	1:B:539:ILE:CD1	2.95	0.44
1:B:456:ASP:O	1:B:457:ASP:HB2	2.16	0.44
1:B:279:LYS:HE2	4:B:2080:HOH:O	2.17	0.44
1:B:122:THR:HG21	1:B:134:ILE:HD11	1.98	0.44
1:A:451:ALA:HA	1:A:460:LEU:O	2.18	0.43
1:B:289:TYR:HB3	1:B:338:THR:HB	2.00	0.43
1:B:257:ASP:HB3	1:B:416:GLU:OE2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:499[A]:ARG:HA	1:A:518:SER:HB2	2.00	0.42
1:A:448:TYR:N	1:A:449:PRO:CD	2.83	0.42
1:B:59:ILE:CD1	1:B:539:ILE:HD11	2.49	0.42
1:A:154:LEU:HD11	1:A:214:ARG:HB3	2.01	0.42
1:A:154:LEU:CD1	1:A:214:ARG:HB3	2.49	0.42
1:A:76:ARG:HD2	1:A:624:TRP:CD1	2.55	0.41
1:B:561:ALA:HA	1:B:645:PRO:HD2	2.01	0.41
1:B:465:SER:HB3	1:B:500:ASN:HB3	2.01	0.41
1:A:231:GLU:HA	1:A:251:ALA:O	2.20	0.41
1:B:451:ALA:HA	1:B:460:LEU:O	2.20	0.41

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	598/615 (97%)	578 (97%)	16 (3%)	4 (1%)	22 9
1	B	608/615 (99%)	581 (96%)	24 (4%)	3 (0%)	29 14
All	All	1206/1230 (98%)	1159 (96%)	40 (3%)	7 (1%)	25 11

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	518	SER
1	B	518	SER
1	A	40	HIS
1	A	448	TYR
1	B	448	TYR
1	A	590	VAL
1	B	590	VAL

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	501/510 (98%)	490 (98%)	11 (2%)	52	36
1	B	506/510 (99%)	491 (97%)	15 (3%)	41	24
All	All	1007/1020 (99%)	981 (97%)	26 (3%)	46	29

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	LYS
1	A	39	TYR
1	A	141	TYR
1	A	145	LYS
1	A	207	ARG
1	A	240	SER
1	A	260	ASP
1	A	425	ARG
1	A	489	PHE
1	A	544	TYR
1	A	594	ASP
1	B	141	TYR
1	B	145	LYS
1	B	154	LEU
1	B	160	GLU
1	B	196	LYS
1	B	222	SER
1	B	240	SER
1	B	260	ASP
1	B	262	GLN
1	B	418	LYS
1	B	544	TYR
1	B	549	SER
1	B	590	VAL
1	B	594	ASP
1	B	615	SER

Sometimes 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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	ACT	A	703	-	3,3,3	0.59	0	3,3,3	0.97	0
3	ACT	B	701	-	3,3,3	0.81	0	3,3,3	0.92	0
3	ACT	A	702	-	3,3,3	0.76	0	3,3,3	0.78	0
3	ACT	A	701	-	3,3,3	1.49	1 (33%)	3,3,3	0.95	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	ACT	O-C	2.44	1.33	1.22

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 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	701	ACT	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 (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	599/615 (97%)	-0.19	3 (0%)	91	91	14, 25, 41, 58	42 (7%)
1	B	609/615 (99%)	0.03	9 (1%)	73	73	18, 33, 53, 72	58 (9%)
All	All	1208/1230 (98%)	-0.08	12 (0%)	82	82	14, 28, 49, 72	100 (8%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	226	TYR	4.1
1	B	144	LEU	3.2
1	B	238	GLY	3.0
1	B	106	VAL	2.7
1	A	142	GLU	2.4
1	B	548	ASP	2.3
1	A	39	TYR	2.3
1	A	197	VAL	2.3
1	B	239	LEU	2.2
1	B	556	GLY	2.2
1	B	142	GLU	2.2
1	B	222	SER	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 monosaccharides 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(Å^2)	Q<0.9
3	ACT	A	701	4/4	0.68	0.20	26,32,35,39	0
3	ACT	B	701	4/4	0.90	0.20	25,27,27,31	4
3	ACT	A	703	4/4	0.93	0.16	37,43,44,48	0
3	ACT	A	702	4/4	0.95	0.16	17,18,18,19	4
2	CU	A	700	1/1	0.99	0.08	22,22,22,22	0
2	CU	B	700	1/1	0.99	0.08	30,30,30,30	0

6.5 Other polymers (i)

There are no such residues in this entry.