Full wwPDB X-ray Structure Validation Report (i)

Jan 3, 2024 – 09:32 am GMT

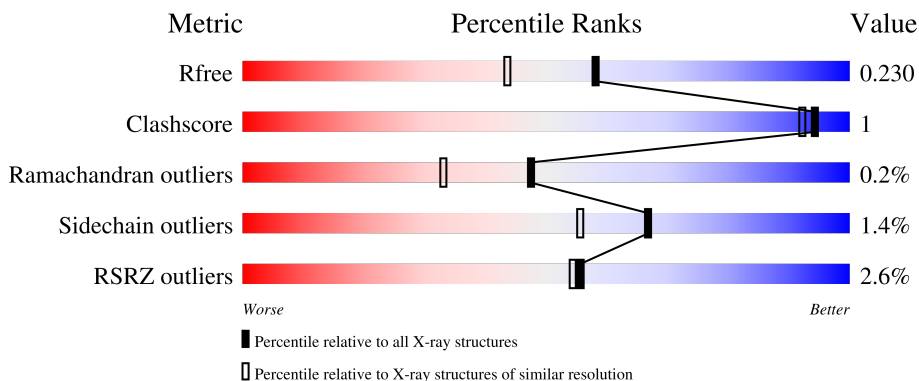|		
PDB ID	:	4YGM
Title	:	Vaccinia virus his-D4/A20(1-50) in complex with uracil
Authors	:	Tarbouriech, N.; Iseni, F.; Burmeister, W.P.
Deposited on	:	2015-02-26
Resolution	:	1.85 Å(reported)

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	Percentile Ranks	Value
Rfree		0.230
Clashscore		1
Ramachandran outliers		0.2%
Sidechain outliers		1.4%
RSRZ outliers		2.6%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

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 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	232	3% / 92% / • / 5%
1	B	232	% / 88% / 6% / 6%
2	C	52	4% / 90% / 6% / • / •
2	D	52	6% / 92% / • / • / •

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
4	URA	A	304	-	X	-	-
4	URA	B	305	-	X	-	-

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 4795 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 Uracil-DNA glycosylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	218	Total	C	N	O	S	0	0	0
			1773	1149	293	325	6			
1	A	220	Total	C	N	O	S	0	0	0
			1788	1158	295	329	6			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	MET	-	initiating methionine	UNP P20536
B	-12	GLY	-	expression tag	UNP P20536
B	-11	SER	-	expression tag	UNP P20536
B	-10	SER	-	expression tag	UNP P20536
B	-9	HIS	-	expression tag	UNP P20536
B	-8	HIS	-	expression tag	UNP P20536
B	-7	HIS	-	expression tag	UNP P20536
B	-6	HIS	-	expression tag	UNP P20536
B	-5	HIS	-	expression tag	UNP P20536
B	-4	HIS	-	expression tag	UNP P20536
B	-3	SER	-	expression tag	UNP P20536
B	-2	GLN	-	expression tag	UNP P20536
B	-1	ASP	-	expression tag	UNP P20536
B	0	PRO	-	expression tag	UNP P20536
A	-13	MET	-	initiating methionine	UNP P20536
A	-12	GLY	-	expression tag	UNP P20536
A	-11	SER	-	expression tag	UNP P20536
A	-10	SER	-	expression tag	UNP P20536
A	-9	HIS	-	expression tag	UNP P20536
A	-8	HIS	-	expression tag	UNP P20536
A	-7	HIS	-	expression tag	UNP P20536
A	-6	HIS	-	expression tag	UNP P20536
A	-5	HIS	-	expression tag	UNP P20536
A	-4	HIS	-	expression tag	UNP P20536
A	-3	SER	-	expression tag	UNP P20536

Continued from previous page...

| Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLN	-	expression tag	UNP P20536
A	-1	ASP	-	expression tag	UNP P20536
A	0	PRO	-	expression tag	UNP P20536

- Molecule 2 is a protein called DNA polymerase processivity factor component A20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	50	Total	C	N	O	S	0	0	0
			400	257	63	79	1			
2	C	51	Total	C	N	O	S	0	0	0
			405	260	64	80	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	GLY	-	expression tag	UNP P20995
D	0	ALA	-	expression tag	UNP P20995
C	-1	GLY	-	expression tag	UNP P20995
C	0	ALA	-	expression tag	UNP P20995

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).
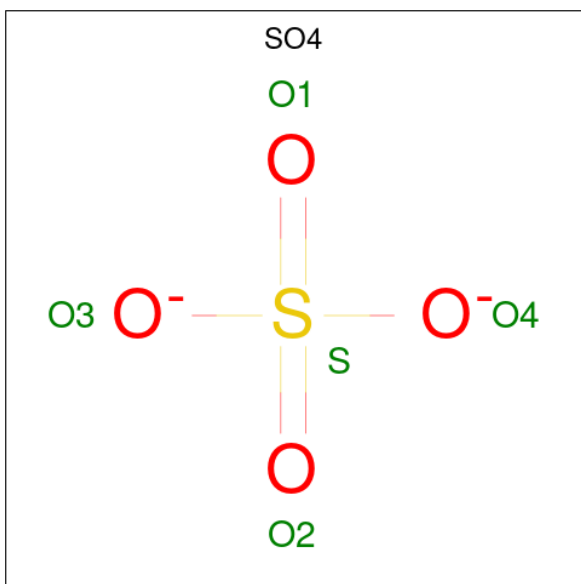
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		

Continued on next page...
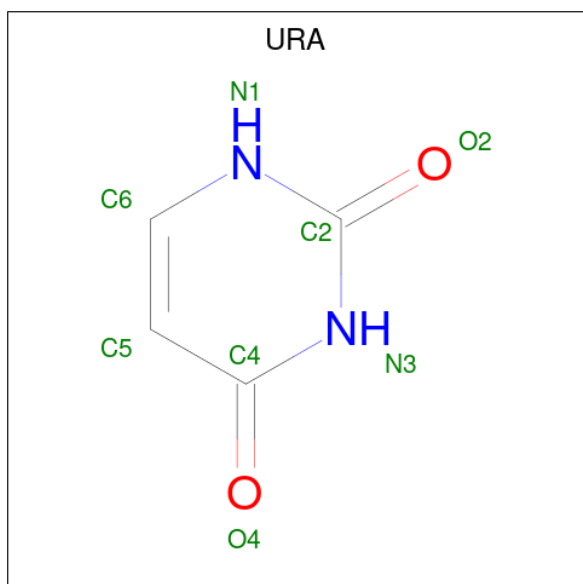
Continued from previous page...

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

- Molecule 4 is URACIL (three-letter code: URA) (formula: $C_4H_4N_2O_2$).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total 8	C 4	N 2	O 2	0	0
4	A	1	Total 8	C 4	N 2	O 2	0	0

- Molecule 5 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	163	Total 163	O 163	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	32	Total 32	O 32	0	0
5	A	153	Total 153	O 153	0	0
5	C	30	Total 30	O 30	0	0

3 Residue-property plots ⓘ

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.
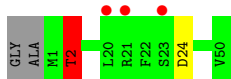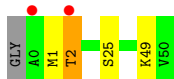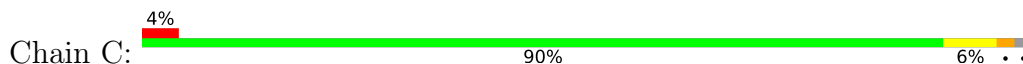
• Molecule 1: Uracil-DNA glycosylase
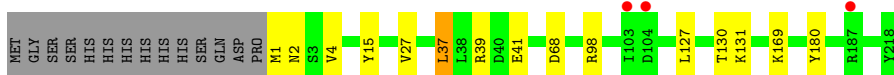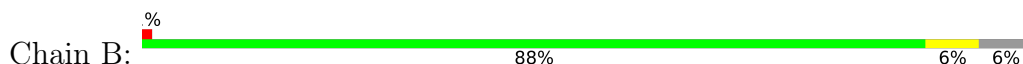
• Molecule 1: Uracil-DNA glycosylase

• Molecule 2: DNA polymerase processivity factor component A20

• Molecule 2: DNA polymerase processivity factor component A20

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	92.79Å 92.79Å 146.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	80.36 – 1.85 46.39 – 1.85	Depositor EDS
% Data completeness (in resolution range)	95.1 (80.36-1.85) 95.1 (46.39-1.85)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.26 (at 1.84Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.191 , 0.222 0.198 , 0.230	Depositor DCC
R_{free} test set	2992 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	29.0	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 38.7	EDS
L-test for twinning[2]	$< |L| > = 0.48, < L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	4795	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: SO4, URA

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	0.64	0/1838	0.81	1/2500 (0.0%)
1	B	0.68	0/1822	0.82	2/2477 (0.1%)
2	C	0.65	0/412	0.79	0/555
2	D	0.72	0/407	0.83	0/548
All	All	0.66	0/4479	0.81	3/6080 (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
2	C	0	2
2	D	0	1
All	All	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	193	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	B	98	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	39	ARG	NE-CZ-NH2	-5.10	117.75	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	1	MET	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
2	C	49	LYS	Peptide
2	D	2	THR	Peptide

5.2 Too-close contacts ⓘ

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	1788	0	1781	1	0
1	B	1773	0	1770	9	0
2	C	405	0	411	1	0
2	D	400	0	406	1	0
3	A	15	0	0	0	0
3	B	20	0	0	0	0
4	A	8	0	3	0	0
4	B	8	0	3	0	0
5	A	153	0	0	0	0
5	B	163	0	0	2	0
5	C	30	0	0	1	0
5	D	32	0	0	0	0
All	All	4795	0	4374	11	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 1.

All (11) 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:68:ASP:OD1	1:B:130:THR:HG23	2.09	0.53
1:B:180:TYR:CE2	2:D:2:THR:HG21	2.46	0.51
1:A:199:ASN:O	1:A:203:GLU:HG3	2.12	0.49
1:B:169:LYS:NZ	5:B:403:HOH:O	2.46	0.48
1:B:37:LEU:HD12	1:B:37:LEU:C	2.33	0.47
1:B:2:ASN:HB2	1:B:15:TYR:CZ	2.50	0.46
1:B:130:THR:HG22	1:B:131:LYS:HG3	1.97	0.46
1:B:41:GLU:HG2	1:B:127:LEU:HD22	1.98	0.46
1:B:4:VAL:HG21	1:B:27:VAL:HG21	1.97	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:N	5:B:404:HOH:O	2.50	0.44
2:C:2:THR:HG22	5:C:102:HOH:O	2.19	0.43

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	218/232 (94%)	212 (97%)	5 (2%)	1 (0%)	29	15
1	B	216/232 (93%)	208 (96%)	8 (4%)	0	100	100
2	C	49/52 (94%)	47 (96%)	2 (4%)	0	100	100
2	D	48/52 (92%)	45 (94%)	3 (6%)	0	100	100
All	All	531/568 (94%)	512 (96%)	18 (3%)	1 (0%)	47	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	138	ASP

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	202/213 (95%)	200 (99%)	2 (1%)	76	69

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	200/213 (94%)	199 (100%)	1 (0%)	88	86
2	C	45/45 (100%)	43 (96%)	2 (4%)	28	12
2	D	45/45 (100%)	43 (96%)	2 (4%)	28	12
All	All	492/516 (95%)	485 (99%)	7 (1%)	67	55

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	37	LEU
2	D	2	THR
2	D	24	ASP
1	A	103	ILE
1	A	117	ILE
2	C	2	THR
2	C	25	SER

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

9 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	$\#\|Z\| > 2$
3	SO4	A	301	-	4,4,4	0.40	0	6,6,6	0.38	0
3	SO4	A	302	-	4,4,4	0.23	0	6,6,6	0.20	0
4	URA	B	305	-	8,8,8	1.75	2 (25%)	9,10,10	3.42	6 (66%)
3	SO4	B	301	-	4,4,4	0.36	0	6,6,6	0.36	0
3	SO4	B	304	-	4,4,4	0.33	0	6,6,6	0.13	0
4	URA	A	304	-	8,8,8	1.59	3 (37%)	9,10,10	3.15	6 (66%)
3	SO4	A	303	-	4,4,4	0.42	0	6,6,6	0.19	0
3	SO4	B	303	-	4,4,4	0.56	0	6,6,6	0.39	0
3	SO4	B	302	-	4,4,4	0.20	0	6,6,6	0.42	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
4	URA	B	305	-	-	-	0/1/1/1
4	URA	A	304	-	-	-	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	305	URA	C4-N3	-3.28	1.32	1.38
4	B	305	URA	C6-C5	2.44	1.39	1.34
4	A	304	URA	C6-C5	2.29	1.39	1.34
4	A	304	URA	C4-N3	-2.13	1.34	1.38
4	A	304	URA	C5-C4	-2.10	1.39	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	305	URA	N1-C2-N3	5.62	121.49	115.13
4	B	305	URA	C6-N1-C2	-5.09	119.28	122.40
4	A	304	URA	N1-C2-N3	5.05	120.85	115.13
4	A	304	URA	C4-N3-C2	-4.46	121.28	125.70
4	B	305	URA	C4-N3-C2	-4.35	121.38	125.70
4	A	304	URA	C6-N1-C2	-3.92	119.99	122.40
4	A	304	URA	O4-C4-C5	-3.13	119.65	125.16

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	305	URA	O2-C2-N1	-2.92	119.57	122.79
4	B	305	URA	O4-C4-C5	-2.90	120.06	125.16
4	B	305	URA	C5-C4-N3	2.73	118.93	114.84
4	A	304	URA	O2-C2-N1	-2.60	119.93	122.79
4	A	304	URA	C5-C4-N3	2.46	118.52	114.84

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	220/232 (94%)	-0.16	6 (2%)	54	53	22, 31, 51, 71	0
1	B	218/232 (93%)	-0.09	3 (1%)	75	76	21, 31, 48, 66	0
2	C	51/52 (98%)	0.00	2 (3%)	39	38	25, 33, 53, 60	0
2	D	50/52 (96%)	0.15	3 (6%)	21	21	24, 34, 62, 80	0
All	All	539/568 (94%)	-0.09	14 (2%)	56	54	21, 32, 53, 80	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	36	TRP	4.8
1	A	185	ARG	3.8
1	B	103	ILE	3.6
1	A	186	ASP	3.4
2	D	20	LEU	2.7
1	A	104	ASP	2.6
2	C	2	THR	2.6
1	A	187	ARG	2.6
1	B	187	ARG	2.4
2	D	21	ARG	2.4
2	D	23	SER	2.3
2	C	0	ALA	2.2
1	A	39	ARG	2.2
1	B	104	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

WORLDWIDE
wwPDB
PROTEIN DATA BANK

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
4	URA	A	304	8/8	0.95	0.07	25,27,29,29	0
3	SO4	B	303	5/5	0.96	0.07	39,43,49,50	0
3	SO4	A	303	5/5	0.98	0.09	43,44,49,50	0
4	URA	B	305	8/8	0.98	0.08	21,22,23,24	0
3	SO4	B	304	5/5	0.98	0.19	48,49,50,50	5
3	SO4	A	302	5/5	0.99	0.14	52,54,55,59	0
3	SO4	B	301	5/5	0.99	0.09	42,45,46,48	0
3	SO4	B	302	5/5	0.99	0.11	48,49,51,55	0
3	SO4	A	301	5/5	0.99	0.15	41,42,43,45	0

6.5 Other polymers ⓘ

There are no such residues in this entry.