Full wwPDB X-ray Structure Validation Report ⓘ

Sep 20, 2023 – 11:00 PM EDT

PDB ID	:	5JW6
Title	:	Cystal structure of aspartate semialdehyde dehydrogenase from Aspergillus fumigatus
Authors	:	Dahal, G.P.; Viola, R.E.
Deposited on	:	2016-05-11
Resolution	:	2.39 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:

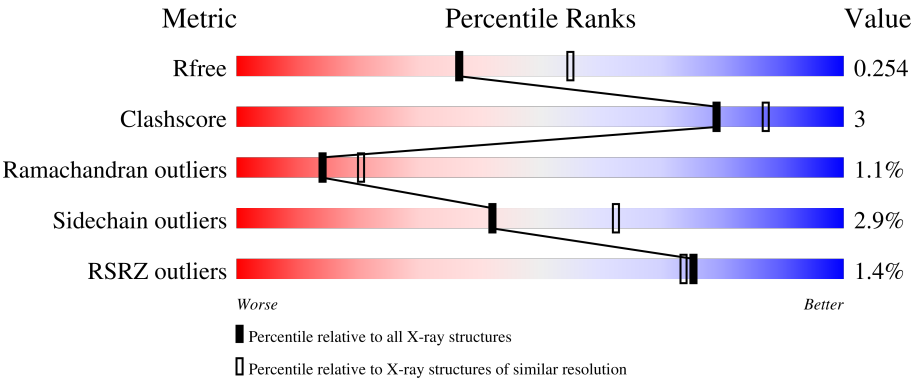
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

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

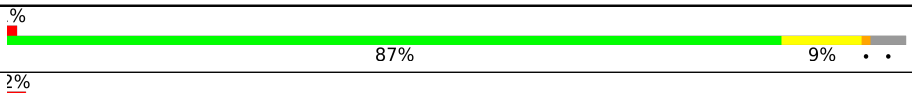
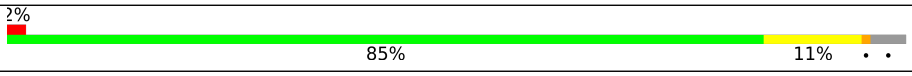
The reported resolution of this entry is 2.39 Å.

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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	371	
1	B	371	

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	402	-	-	X	-

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5598 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.

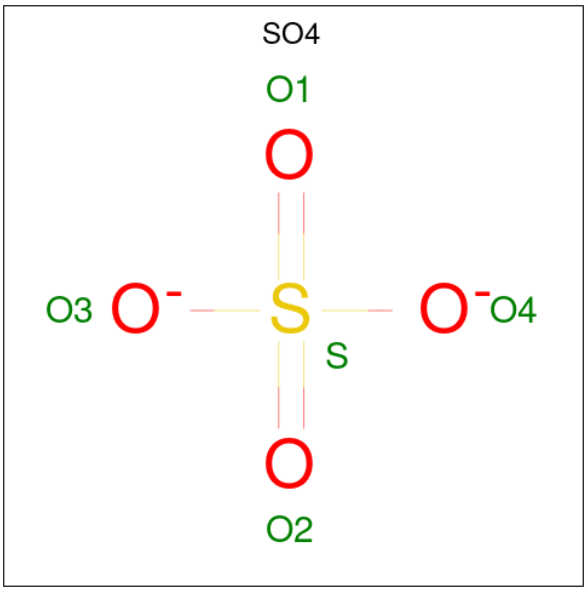
- Molecule 1 is a protein called Aspartate-semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	358	Total	C	N	O	S	0	1	0
			2733	1725	480	515	13			
1	B	358	Total	C	N	O	S	0	1	0
			2733	1725	480	515	13			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	364	LEU	-	expression tag	UNP Q4WWR8
A	365	GLU	-	expression tag	UNP Q4WWR8
A	366	HIS	-	expression tag	UNP Q4WWR8
A	367	HIS	-	expression tag	UNP Q4WWR8
A	368	HIS	-	expression tag	UNP Q4WWR8
A	369	HIS	-	expression tag	UNP Q4WWR8
A	370	HIS	-	expression tag	UNP Q4WWR8
A	371	HIS	-	expression tag	UNP Q4WWR8
B	364	LEU	-	expression tag	UNP Q4WWR8
B	365	GLU	-	expression tag	UNP Q4WWR8
B	366	HIS	-	expression tag	UNP Q4WWR8
B	367	HIS	-	expression tag	UNP Q4WWR8
B	368	HIS	-	expression tag	UNP Q4WWR8
B	369	HIS	-	expression tag	UNP Q4WWR8
B	370	HIS	-	expression tag	UNP Q4WWR8
B	371	HIS	-	expression tag	UNP Q4WWR8

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

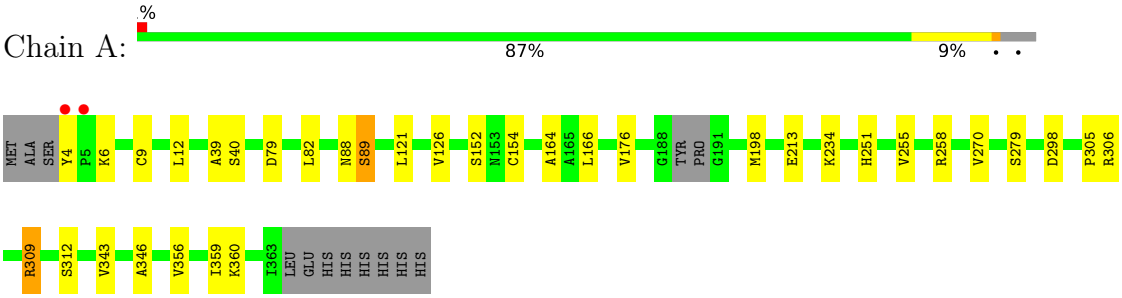
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	63	Total O 63 63	0	0
3	B	44	Total O 44 44	0	0

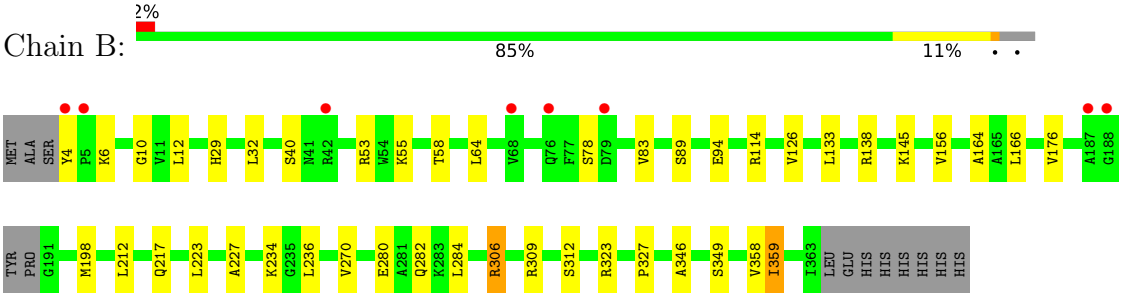
3 Residue-property plots (i)

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: Aspartate-semialdehyde dehydrogenase



• Molecule 1: Aspartate-semialdehyde dehydrogenase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	94.50Å 94.50Å 202.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	81.84 – 2.39 75.86 – 2.39	Depositor EDS
% Data completeness (in resolution range)	100.0 (81.84-2.39) 100.0 (75.86-2.39)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.90 (at 2.40Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.201 , 0.250 0.208 , 0.254	Depositor DCC
R_{free} test set	2153 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	34.1	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 29.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5598	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% 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

5.1 Standard geometry

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	0.89	1/2785 (0.0%)	0.99	10/3765 (0.3%)
1	B	0.88	1/2785 (0.0%)	0.99	7/3765 (0.2%)
All	All	0.88	2/5570 (0.0%)	0.99	17/7530 (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	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	213	GLU	CD-OE2	5.48	1.31	1.25
1	B	94	GLU	CG-CD	5.30	1.59	1.51

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	309	ARG	NE-CZ-NH1	10.87	125.74	120.30
1	B	309	ARG	NE-CZ-NH2	-10.32	115.14	120.30
1	A	198	MET	CG-SD-CE	-9.41	85.15	100.20
1	A	309	ARG	NE-CZ-NH2	-8.88	115.86	120.30
1	A	309	ARG	NE-CZ-NH1	8.13	124.37	120.30
1	A	258	ARG	NE-CZ-NH1	7.30	123.95	120.30
1	B	306	ARG	NE-CZ-NH1	-7.01	116.79	120.30
1	B	198	MET	CG-SD-CE	-6.59	89.66	100.20
1	B	306	ARG	NE-CZ-NH2	6.58	123.59	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	323	ARG	NE-CZ-NH1	6.47	123.54	120.30
1	A	258	ARG	NE-CZ-NH2	-6.24	117.18	120.30
1	A	306	ARG	NE-CZ-NH1	-5.90	117.35	120.30
1	A	79	ASP	CB-CG-OD1	5.67	123.40	118.30
1	B	114	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	A	306	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	A	298	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	309	ARG	CB-CG-CD	5.19	125.10	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	4	TYR	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	2733	0	2739	18	0
1	B	2733	0	2739	13	0
2	A	20	0	0	2	0
2	B	5	0	0	0	0
3	A	63	0	0	2	0
3	B	44	0	0	0	0
All	All	5598	0	5478	31	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 (31) 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:88:ASN:O	1:A:89:SER:HB3	1.83	0.78
1:A:88:ASN:O	1:A:89:SER:CB	2.41	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:GLN:NE2	1:B:236:LEU:O	2.32	0.62
1:A:309:ARG:HD2	3:A:523:HOH:O	2.03	0.58
1:A:166:LEU:HD21	1:A:270:VAL:HG13	1.91	0.53
1:A:279:SER:HB2	3:A:526:HOH:O	2.08	0.53
1:A:154[A]:CYS:SG	1:A:251:HIS:HE1	2.31	0.53
1:A:40:SER:HB3	2:A:402:SO4:O2	2.10	0.52
1:A:359:ILE:O	1:A:360:LYS:HB2	2.12	0.48
1:A:154[A]:CYS:SG	1:A:251:HIS:CE1	3.07	0.48
1:B:358:VAL:C	1:B:359:ILE:O	2.51	0.48
1:A:356:VAL:O	1:A:359:ILE:O	2.32	0.47
1:A:9:CYS:SG	1:A:82:LEU:HD11	2.55	0.47
1:B:223:LEU:HD22	1:B:227:ALA:HA	1.96	0.47
1:A:176:VAL:HG23	1:A:255:VAL:HG13	1.96	0.47
1:B:126:VAL:HG13	1:B:164:ALA:HB2	1.97	0.46
1:B:10:GLY:HA3	1:B:83:VAL:HG12	1.97	0.45
1:B:29:HIS:HB3	1:B:32:LEU:O	2.15	0.45
1:A:305:PRO:O	1:A:309:ARG:HG2	2.17	0.44
1:B:280:GLU:OE2	1:B:284:LEU:HD11	2.18	0.44
1:B:138:ARG:NH2	1:B:145:LYS:O	2.36	0.44
1:A:39:ALA:HB1	2:A:402:SO4:O4	2.19	0.42
1:B:64:LEU:HD12	1:B:64:LEU:O	2.19	0.42
1:A:126:VAL:HG13	1:A:164:ALA:HB2	2.02	0.42
1:B:166:LEU:HD21	1:B:270:VAL:HG13	2.01	0.42
1:B:234:LYS:HA	1:B:234:LYS:HE2	2.01	0.41
1:A:121:LEU:HD11	1:A:152:SER:HA	2.03	0.41
1:B:234:LYS:HA	1:B:234:LYS:CE	2.51	0.41
1:A:359:ILE:O	1:A:360:LYS:CB	2.69	0.40
1:B:156:VAL:HA	1:B:212:LEU:HD13	2.03	0.40
1:A:166:LEU:CD2	1:A:270:VAL:HG13	2.51	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	354/371 (95%)	337 (95%)	13 (4%)	4 (1%)	14 20
1	B	354/371 (95%)	333 (94%)	17 (5%)	4 (1%)	14 20
All	All	708/742 (95%)	670 (95%)	30 (4%)	8 (1%)	14 20

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	SER
1	B	89	SER
1	A	6	LYS
1	A	346	ALA
1	B	40	SER
1	B	346	ALA
1	B	359	ILE
1	A	343	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	293/305 (96%)	290 (99%)	3 (1%)	76 88
1	B	293/305 (96%)	279 (95%)	14 (5%)	25 41
All	All	586/610 (96%)	569 (97%)	17 (3%)	42 62

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	234	LYS
1	A	312	SER
1	B	4	TYR
1	B	6	LYS
1	B	12	LEU
1	B	53	ARG
1	B	55	LYS

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Mol	Chain	Res	Type
1	B	58	THR
1	B	78	SER
1	B	133	LEU
1	B	176	VAL
1	B	282	GLN
1	B	306	ARG
1	B	312	SER
1	B	327	PRO
1	B	349	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	ASN
1	A	251	HIS
1	B	41	ASN
1	B	151	ASN
1	B	214	ASN
1	B	251	HIS

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)

5 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	401	-	4,4,4	0.48	0	6,6,6	0.44	0
2	SO4	A	404	-	4,4,4	0.30	0	6,6,6	0.60	0
2	SO4	A	403	-	4,4,4	0.61	0	6,6,6	0.44	0
2	SO4	A	402	-	4,4,4	0.70	0	6,6,6	0.45	0
2	SO4	A	401	-	4,4,4	0.27	0	6,6,6	0.51	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	402	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

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^{th} 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	358/371 (96%)	-0.07	2 (0%)	89	88	18, 34, 56, 122	0
1	B	358/371 (96%)	0.13	8 (2%)	62	60	19, 39, 71, 115	0
All	All	716/742 (96%)	0.03	10 (1%)	75	73	18, 36, 66, 122	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	PRO	4.1
1	B	4	TYR	4.0
1	B	5	PRO	3.2
1	B	68	VAL	2.8
1	B	79	ASP	2.3
1	B	188	GLY	2.3
1	B	187	ALA	2.3
1	A	4	TYR	2.2
1	B	42	ARG	2.2
1	B	76	GLN	2.0

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

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

6.3 Carbohydrates

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, 95^{th} 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
2	SO4	A	402	5/5	0.94	0.15	59,59,66,73	0
2	SO4	A	403	5/5	0.95	0.15	44,49,54,62	0
2	SO4	A	404	5/5	0.96	0.13	55,64,69,79	0
2	SO4	A	401	5/5	0.99	0.13	23,25,27,27	0
2	SO4	B	401	5/5	0.99	0.13	22,24,29,32	0

6.5 Other polymers (i)

There are no such residues in this entry.