



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 05:00 pm BST

PDB ID : 6MKY
Title : Human SDS22
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Deposited on : 2018-09-26
Resolution : 2.90 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 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

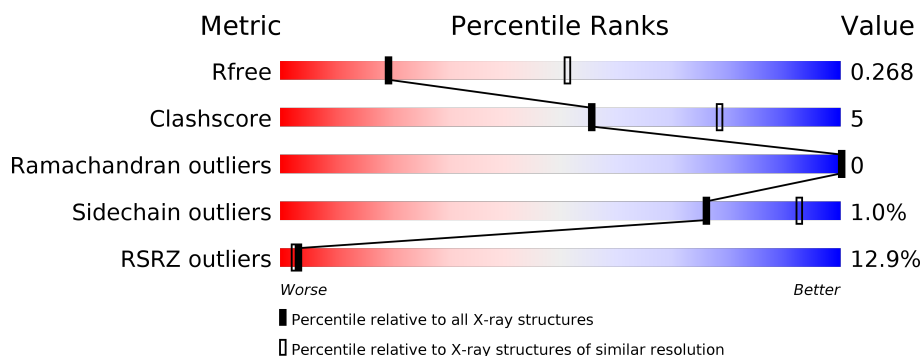
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 ≥ 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	261	<div> <div>8%</div> <div>84%</div> <div>15%</div> <div>..</div> </div>
1	B	261	<div> <div>18%</div> <div>88%</div> <div>12%</div> </div>

2 Entry composition [i](#)

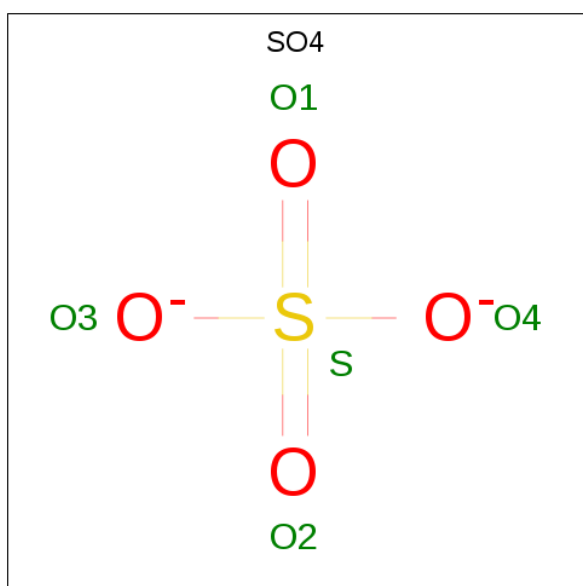
There are 3 unique types of molecules in this entry. The entry contains 4301 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein phosphatase 1 regulatory subunit 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	1	0
			2113	1332	368	405	8			
1	B	260	Total	C	N	O	S	0	0	0
			2121	1336	372	406	7			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

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

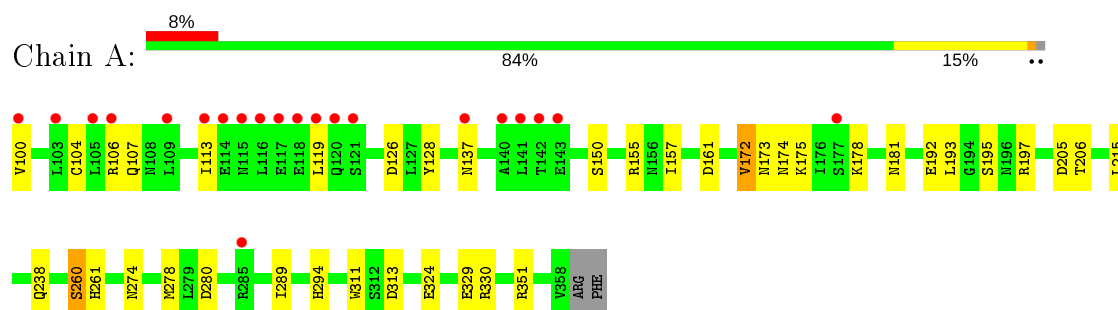
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	8	Total	O	0	0
			8	8		

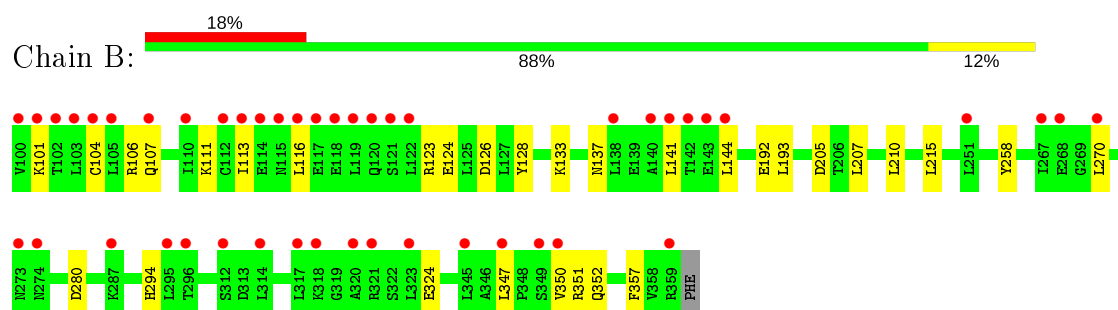
3 Residue-property plots [i](#)

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: Protein phosphatase 1 regulatory subunit 7



- Molecule 1: Protein phosphatase 1 regulatory subunit 7



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.10Å 76.55Å 128.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.55 – 2.90 38.27 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.2 (37.55-2.90) 99.3 (38.27-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.238 , 0.272 0.236 , 0.268	Depositor DCC
R_{free} test set	1738 reflections (10.31%)	wwPDB-VP
Wilson B-factor (Å ²)	73.2	Xtriage
Anisotropy	0.452	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4301	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.61% 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	0.23	0/2134	0.46	0/2876
1	B	0.23	0/2139	0.44	0/2882
All	All	0.23	0/4273	0.45	0/5758

There are no bond length outliers.

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	2113	0	2208	29	0
1	B	2121	0	2216	21	0
2	A	40	0	0	3	0
2	B	10	0	0	1	0
3	A	9	0	0	0	0
3	B	8	0	0	0	0
All	All	4301	0	4424	44	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 5.

All (44) 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:113:ILE:O	1:B:137:ASN:ND2	2.27	0.68
1:A:106:ARG:HG3	1:A:128:TYR:HD1	1.61	0.65
1:A:289:ILE:HB	1:A:313:ASP:HB3	1.81	0.63
1:A:238:GLN:NE2	2:A:401:SO4:O1	2.33	0.62
1:B:141:LEU:HB3	1:B:144:LEU:HD13	1.82	0.60
1:B:270:LEU:O	1:B:294:HIS:NE2	2.23	0.60
1:A:100:VAL:HG21	1:A:119:LEU:HB3	1.83	0.60
1:B:106:ARG:HG2	1:B:107:GLN:HG3	1.86	0.57
1:B:111:LYS:HD2	1:B:133:LYS:HE2	1.87	0.56
1:A:260:SER:OG	1:A:280:ASP:OD2	2.25	0.54
1:A:324:GLU:OE2	1:A:351:ARG:NH1	2.43	0.52
1:B:207:LEU:HB3	1:B:210:LEU:HD13	1.91	0.52
1:A:178:LYS:NZ	2:B:401:SO4:O3	2.40	0.52
1:B:193:LEU:HB2	1:B:215:LEU:HD23	1.92	0.51
1:A:106:ARG:HG2	1:A:107:GLN:HG3	1.93	0.50
1:A:113:ILE:O	1:A:137:ASN:ND2	2.36	0.50
1:A:324:GLU:HG3	1:A:351:ARG:HD3	1.94	0.49
1:A:193:LEU:HB2	1:A:215:LEU:HD23	1.94	0.49
1:B:106:ARG:HG3	1:B:128:TYR:HD2	1.79	0.48
1:B:101:LYS:HE2	1:B:123:ARG:HG3	1.94	0.48
1:A:197:ARG:HG2	1:B:128:TYR:CZ	2.49	0.48
1:A:205:ASP:OD2	1:B:352:GLN:NE2	2.40	0.48
1:A:206:THR:HG22	1:B:357:PHE:HZ	1.79	0.47
1:A:155:ARG:NH1	1:B:124:GLU:OE2	2.34	0.46
1:B:104:CYS:HA	1:B:126:ASP:HB3	1.98	0.46
1:A:311:TRP:N	2:A:402:SO4:O1	2.49	0.46
1:A:206:THR:HA	1:B:357:PHE:CZ	2.51	0.46
1:A:104[B]:CYS:HA	1:A:126:ASP:HB3	1.98	0.45
1:B:113:ILE:HG23	1:B:116:LEU:HD12	1.97	0.45
1:A:150:SER:OG	1:A:172:VAL:HG23	2.16	0.45
1:B:205:ASP:N	1:B:205:ASP:OD1	2.50	0.45
1:A:104[A]:CYS:HA	1:A:126:ASP:HB3	1.98	0.45
1:A:205:ASP:OD1	1:A:205:ASP:N	2.50	0.45
1:A:261:HIS:NE2	2:A:407:SO4:O3	2.48	0.45
1:A:157:ILE:O	1:A:181:ASN:ND2	2.49	0.44
1:A:172:VAL:HG22	1:A:192:GLU:HG2	1.99	0.44
1:A:174:ASN:O	1:B:106:ARG:NH1	2.50	0.44
1:A:329:GLU:OE2	1:A:330:ARG:NH2	2.52	0.43
1:A:173:ASN:HA	1:A:195:SER:O	2.19	0.43
1:A:274:ASN:OD1	1:A:294:HIS:HD2	2.01	0.43
1:B:258:TYR:HA	1:B:280:ASP:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ARG:HG2	1:A:175:LYS:HB2	2.03	0.41
1:B:324:GLU:HG3	1:B:351:ARG:HD3	2.02	0.41
1:B:347:LEU:O	1:B:350:VAL:HG22	2.21	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	258/261 (99%)	238 (92%)	20 (8%)	0	100	100
1	B	258/261 (99%)	239 (93%)	19 (7%)	0	100	100
All	All	516/522 (99%)	477 (92%)	39 (8%)	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	246/247 (100%)	242 (98%)	4 (2%)	62	86
1	B	246/247 (100%)	245 (100%)	1 (0%)	91	97
All	All	492/494 (100%)	487 (99%)	5 (1%)	76	92

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	ASP
1	A	172	VAL
1	A	260	SER
1	A	278	MET
1	B	192	GLU

Some 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

10 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	0.13	0	6,6,6	0.07	0
2	SO4	A	404	-	4,4,4	0.14	0	6,6,6	0.06	0
2	SO4	A	403	-	4,4,4	0.15	0	6,6,6	0.05	0
2	SO4	A	407	-	4,4,4	0.16	0	6,6,6	0.05	0
2	SO4	A	406	-	4,4,4	0.15	0	6,6,6	0.05	0
2	SO4	A	401	-	4,4,4	0.13	0	6,6,6	0.10	0
2	SO4	A	408	-	4,4,4	0.14	0	6,6,6	0.05	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	405	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	A	402	-	4,4,4	0.13	0	6,6,6	0.05	0
2	SO4	B	402	-	4,4,4	0.13	0	6,6,6	0.05	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.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	SO4	1	0
2	A	407	SO4	1	0
2	A	401	SO4	1	0
2	A	402	SO4	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, 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	259/261 (99%)	0.50	21 (8%)	12 9	38, 57, 98, 139	0
1	B	260/261 (99%)	1.05	46 (17%)	1 1	50, 87, 118, 153	0
All	All	519/522 (99%)	0.78	67 (12%)	3 2	38, 72, 115, 153	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	116	LEU	12.8
1	B	119	LEU	7.9
1	A	116	LEU	7.9
1	B	118	GLU	7.3
1	B	120	GLN	6.7
1	B	115	ASN	6.6
1	B	117	GLU	6.4
1	B	317	LEU	6.2
1	A	115	ASN	5.9
1	B	112	CYS	5.7
1	B	122	LEU	5.2
1	B	121	SER	5.0
1	A	117	GLU	4.6
1	B	140	ALA	4.2
1	B	295	LEU	4.2
1	A	103	LEU	4.1
1	B	312	SER	4.0
1	B	321	ARG	4.0
1	B	113	ILE	3.9
1	B	103	LEU	3.9
1	B	345	LEU	3.9
1	A	118	GLU	3.8
1	B	101	LYS	3.7
1	B	102	THR	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	110	ILE	3.7
1	A	140	ALA	3.5
1	B	141	LEU	3.5
1	B	143	GLU	3.4
1	A	120	GLN	3.3
1	A	109	LEU	3.2
1	A	141	LEU	3.2
1	B	107	GLN	3.1
1	A	142	THR	3.1
1	B	320	ALA	3.1
1	A	119	LEU	3.0
1	A	100	VAL	3.0
1	A	143	GLU	3.0
1	B	323	LEU	2.9
1	A	113	ILE	2.9
1	A	114	GLU	2.9
1	B	114	GLU	2.7
1	B	314	LEU	2.7
1	B	142	THR	2.7
1	A	177	SER	2.7
1	B	274	ASN	2.7
1	B	347	LEU	2.6
1	A	121	SER	2.6
1	B	138	LEU	2.5
1	B	296	THR	2.5
1	B	359	ARG	2.5
1	B	350	VAL	2.4
1	B	100	VAL	2.4
1	B	268	GLU	2.3
1	B	105	LEU	2.3
1	B	270	LEU	2.3
1	B	144	LEU	2.2
1	B	273	ASN	2.2
1	B	287	LYS	2.2
1	B	318	LYS	2.2
1	A	106	ARG	2.2
1	A	285	ARG	2.1
1	A	105	LEU	2.1
1	B	104	CYS	2.1
1	A	137	ASN	2.0
1	B	251	LEU	2.0
1	B	267	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	349	SER	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, 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
2	SO4	A	405	5/5	0.76	0.29	95,100,122,130	5
2	SO4	A	406	5/5	0.82	0.30	64,87,104,117	0
2	SO4	A	407	5/5	0.84	0.20	65,99,103,110	0
2	SO4	A	404	5/5	0.86	0.23	78,118,125,134	0
2	SO4	A	408	5/5	0.87	0.26	88,99,118,120	0
2	SO4	A	402	5/5	0.91	0.20	78,100,121,121	0
2	SO4	B	402	5/5	0.92	0.42	114,118,129,130	5
2	SO4	A	403	5/5	0.93	0.15	38,68,108,119	0
2	SO4	B	401	5/5	0.99	0.22	60,68,70,80	0
2	SO4	A	401	5/5	0.99	0.22	45,46,62,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.