



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:22 AM UTC

PDB ID : 4F8E / pdb_00004f8e
Title : Crystal structure of human PRS1 D52H mutant
Authors : Chen, P.; Teng, M.; Li, X.
Deposited on : 2012-05-17
Resolution : 2.27 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

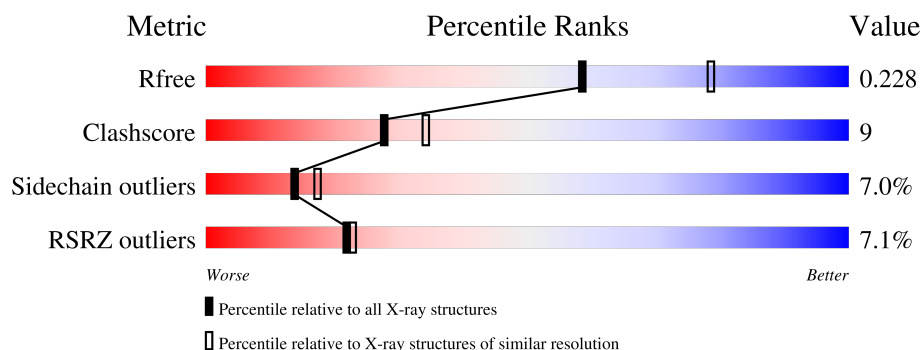
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	145	 10% 82% 14% ..
1	B	145	 6% 81% 17% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4713 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 Ribose-phosphate pyrophosphokinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	308	Total	C	N	O	S	0	0	0
			2287	1436	398	438	15			
1	B	308	Total	C	N	O	S	0	0	0
			2321	1460	403	441	17			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	52	HIS	ASP	engineered mutation	UNP P60891
A	319	LEU	-	expression tag	UNP P60891
A	320	GLU	-	expression tag	UNP P60891
A	321	HIS	-	expression tag	UNP P60891
A	322	HIS	-	expression tag	UNP P60891
A	323	HIS	-	expression tag	UNP P60891
A	324	HIS	-	expression tag	UNP P60891
A	325	HIS	-	expression tag	UNP P60891
A	326	HIS	-	expression tag	UNP P60891
B	52	HIS	ASP	engineered mutation	UNP P60891
B	319	LEU	-	expression tag	UNP P60891
B	320	GLU	-	expression tag	UNP P60891
B	321	HIS	-	expression tag	UNP P60891
B	322	HIS	-	expression tag	UNP P60891
B	323	HIS	-	expression tag	UNP P60891
B	324	HIS	-	expression tag	UNP P60891
B	325	HIS	-	expression tag	UNP P60891
B	326	HIS	-	expression tag	UNP P60891

- Molecule 2 is SULFATE ION (CCD ID: 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	B	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 MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	29	Total	O	0	0
			29	29		

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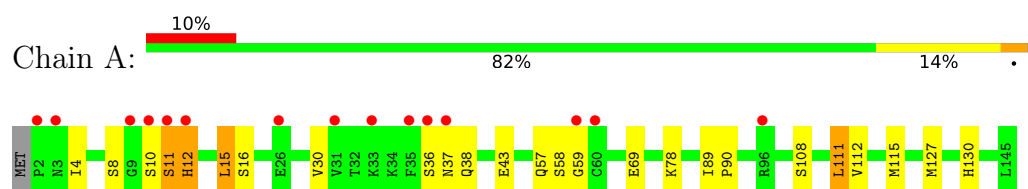
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	44	Total	O	0	0
			44	44		

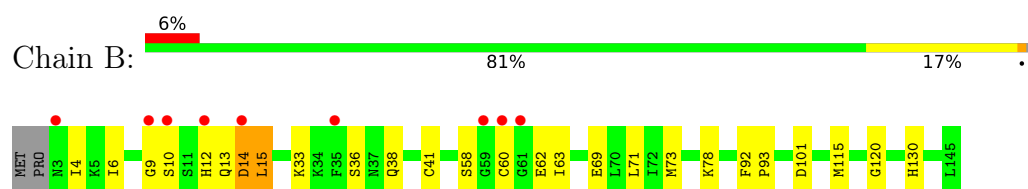
3 Residue-property plots [i](#)

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: Ribose-phosphate pyrophosphokinase 1



- Molecule 1: Ribose-phosphate pyrophosphokinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	170.31Å 170.31Å 61.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.27 50.00 – 2.27	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.27) 99.8 (50.00-2.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.27Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.178 , 0.233 (Not available) , 0.228	Depositor DCC
R_{free} test set	1555 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.417	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.019 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4713	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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: MG, 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	0.50	0/2321	0.88	3/3153 (0.1%)
1	B	0.54	0/2356	0.90	4/3196 (0.1%)
All	All	0.52	0/4677	0.89	7/6349 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15	LEU	N-CA-C	-8.88	102.36	113.55
1	B	14	ASP	N-CA-C	8.02	127.88	110.80
1	B	14	ASP	CB-CA-C	-7.06	96.36	110.42
1	B	13	GLN	N-CA-C	-6.08	106.02	113.50
1	A	169	SER	CA-C-N	5.94	125.56	119.56
1	A	169	SER	C-N-CA	5.94	125.56	119.56
1	A	4	ILE	CB-CA-C	-5.07	104.24	111.19

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	2287	0	2224	47	0
1	B	2321	0	2298	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	15	0	0	1	0
2	B	15	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	29	0	0	0	0
4	B	44	0	0	1	0
All	All	4713	0	4522	82	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 9.

All (82) 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:9:GLY:HA2	1:B:73:MET:HE1	1.39	1.01
1:B:10:SER:HB3	1:B:69:GLU:OE1	1.73	0.88
1:A:222:MET:HE2	1:A:252:ILE:HD11	1.59	0.82
1:B:298:GLU:O	1:B:302:ARG:HG2	1.83	0.78
1:A:10:SER:OG	1:A:58:SER:HB2	1.84	0.76
1:B:223:ALA:HB2	1:B:248:LEU:HG	1.70	0.73
1:B:221:ASP:O	1:B:249:THR:HG23	1.91	0.71
1:A:57:GLN:HG2	1:A:89:ILE:HA	1.77	0.65
1:A:253:PHE:O	1:A:280:LYS:HE2	1.97	0.64
1:B:10:SER:N	1:B:58:SER:OG	2.31	0.64
1:A:12:HIS:ND1	1:A:12:HIS:O	2.30	0.64
1:A:221:ASP:O	1:A:249:THR:HG23	1.98	0.63
1:A:256:PRO:O	1:A:259:SER:HB3	1.98	0.63
1:A:223:ALA:HB3	1:A:248:LEU:HD13	1.80	0.63
1:A:284:CYS:SG	1:A:287:ILE:HG13	2.40	0.61
1:B:224:ASP:HA	1:B:252:ILE:HB	1.81	0.61
1:B:278:GLU:H	1:B:278:GLU:CD	2.07	0.61
1:B:36:SER:C	1:B:38:GLN:H	2.09	0.60
1:A:177:ARG:CZ	1:A:221:ASP:HB3	2.32	0.59
1:A:148:GLU:HB3	1:A:149:PRO:HD3	1.83	0.59
1:B:71:LEU:HD11	1:B:115:MET:HE2	1.84	0.58
1:B:220:ASP:HB3	1:B:248:LEU:HD12	1.86	0.57
1:A:111:LEU:HD13	1:A:115:MET:HE2	1.86	0.56
1:B:78:LYS:HE3	1:B:120:GLY:O	2.06	0.56
1:B:130:HIS:HD2	1:B:221:ASP:OD2	1.89	0.56
1:A:192:ILE:HD12	1:A:232:ALA:HB1	1.90	0.53
1:A:172:ALA:HA	1:A:191:LEU:HD21	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:VAL:O	1:A:289:VAL:HA	2.09	0.52
1:B:223:ALA:CB	1:B:248:LEU:HG	2.40	0.52
1:A:57:GLN:HG2	1:A:90:PRO:HD2	1.91	0.51
1:B:33:LYS:O	1:B:41:CYS:HB2	2.11	0.51
1:A:8:SER:HB2	1:A:16:SER:CB	2.41	0.51
1:A:248:LEU:O	1:A:271:VAL:HA	2.11	0.50
1:B:12:HIS:O	1:B:12:HIS:ND1	2.45	0.50
1:A:57:GLN:O	1:A:57:GLN:HG3	2.10	0.50
1:B:169:SER:HB3	1:B:191:LEU:HD23	1.94	0.49
1:A:165:CYS:O	1:A:187:VAL:HB	2.12	0.49
1:A:207:LEU:C	1:A:207:LEU:HD23	2.39	0.47
1:A:12:HIS:HB3	1:A:59:GLY:H	1.80	0.47
1:B:12:HIS:CE1	4:B:1118:HOH:O	2.68	0.46
1:B:71:LEU:CD1	1:B:115:MET:HE2	2.46	0.46
1:A:37:ASN:O	1:A:38:GLN:HB3	2.15	0.46
1:B:221:ASP:HA	1:B:249:THR:CG2	2.45	0.46
1:B:258:ILE:HG23	1:B:284:CYS:HB2	1.97	0.46
1:B:292:ILE:C	1:B:292:ILE:HD12	2.41	0.46
1:A:130:HIS:HD2	1:A:221:ASP:OD1	1.98	0.46
1:B:9:GLY:O	1:B:69:GLU:CD	2.59	0.45
1:A:38:GLN:O	1:A:38:GLN:HG3	2.16	0.45
1:B:204:ARG:NE	1:B:204:ARG:HA	2.32	0.45
1:A:211:VAL:O	1:A:214:ARG:HG2	2.17	0.45
1:B:314:SER:HG	1:B:315:HIS:CE1	2.35	0.45
1:B:237:LEU:HD12	1:B:265:CYS:SG	2.57	0.45
1:A:57:GLN:CG	1:A:90:PRO:HD2	2.47	0.44
1:A:10:SER:CB	1:A:58:SER:HB2	2.48	0.44
1:B:148:GLU:HB3	1:B:149:PRO:HD3	1.98	0.44
1:A:150:ALA:HB1	1:A:290:ILE:HG21	2.00	0.44
1:A:224:ASP:HB3	2:A:401:SO4:O2	2.18	0.44
1:A:36:SER:OG	1:A:37:ASN:N	2.50	0.44
1:A:10:SER:OG	1:A:11:SER:N	2.47	0.43
1:A:237:LEU:HA	1:A:237:LEU:HD23	1.80	0.43
1:A:181:ILE:CD1	1:A:219:VAL:HG11	2.49	0.43
1:A:237:LEU:HD12	1:A:265:CYS:SG	2.59	0.43
1:B:314:SER:OG	1:B:315:HIS:CE1	2.71	0.43
1:A:10:SER:HB3	1:A:58:SER:CB	2.49	0.42
1:B:248:LEU:O	1:B:271:VAL:HA	2.20	0.42
1:A:108:SER:O	1:A:112:VAL:HG23	2.19	0.42
1:B:277:GLN:O	1:B:281:MET:HG3	2.20	0.42
1:A:223:ALA:CB	1:A:248:LEU:HD13	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:SER:H	1:A:69:GLU:CD	2.29	0.41
1:A:127:MET:HE3	1:A:127:MET:HB2	1.81	0.41
1:B:12:HIS:CD2	1:B:14:ASP:HB3	2.56	0.41
1:A:111:LEU:O	1:A:115:MET:HG3	2.20	0.41
1:A:152:LEU:HD12	1:A:152:LEU:HA	1.85	0.41
1:A:242:THR:HG22	1:A:243:ARG:HG3	2.02	0.41
1:B:309:VAL:O	1:B:309:VAL:HG12	2.21	0.41
1:B:92:PHE:HA	1:B:93:PRO:HD3	1.78	0.41
1:A:167:ILE:HG22	1:A:178:VAL:HB	2.03	0.40
1:A:172:ALA:HA	1:A:191:LEU:HD11	2.03	0.40
1:B:152:LEU:HD21	1:B:181:ILE:HG23	2.03	0.40
1:A:12:HIS:O	1:A:12:HIS:CG	2.73	0.40
1:A:12:HIS:NE2	1:A:15:LEU:HB2	2.36	0.40
1:B:167:ILE:O	1:B:189:PHE:HA	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	207/124 (167%)	189 (91%)	18 (9%)	9	11
1	B	248/124 (200%)	234 (94%)	14 (6%)	19	25
All	All	455/248 (184%)	423 (93%)	32 (7%)	14	17

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	12	HIS
1	A	15	LEU
1	A	30	VAL
1	A	43	GLU
1	A	78	LYS
1	A	111	LEU
1	A	152	LEU
1	A	195	GLU
1	A	228	THR
1	A	244	VAL
1	A	248	LEU
1	A	249	THR
1	A	278	GLU
1	A	303	THR
1	A	307	GLU
1	A	311	TYR
1	A	316	VAL
1	B	4	ILE
1	B	6	ILE
1	B	15	LEU
1	B	60	CYS
1	B	62	GLU
1	B	63	ILE
1	B	101	ASP
1	B	192	ILE
1	B	249	THR
1	B	252	ILE
1	B	260	ARG
1	B	294	MET
1	B	307	GLU
1	B	316	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	130	HIS
1	A	277	GLN
1	B	64	ASN
1	B	130	HIS
1	B	263	ASN
1	B	277	GLN

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Mol	Chain	Res	Type
1	B	283	HIS
1	B	315	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	$\# Z > 2$
2	SO4	B	1002	-	4,4,4	0.17	0	6,6,6	0.17	0
2	SO4	A	402	-	4,4,4	0.26	0	6,6,6	0.25	0
2	SO4	A	401	-	4,4,4	0.26	0	6,6,6	0.12	0
2	SO4	B	1003	-	4,4,4	0.12	0	6,6,6	0.25	0
2	SO4	A	403	-	4,4,4	0.26	0	6,6,6	0.11	0
2	SO4	B	1001	-	4,4,4	0.26	0	6,6,6	0.15	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 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	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	144/145 (99%)	0.22	15 (10%) 11 12	21, 35, 74, 119	10 (6%)
1	B	143/145 (98%)	-0.02	9 (6%) 26 27	14, 32, 76, 110	6 (4%)
All	All	287/290 (98%)	0.10	24 (8%) 22 17	14, 33, 76, 119	16 (5%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	35	PHE	8.0
1	A	60	CYS	6.4
1	B	60	CYS	4.8
1	B	61	GLY	4.7
1	A	9	GLY	4.6
1	A	3	ASN	4.4
1	A	12	HIS	3.7
1	A	96	ARG	3.4
1	B	12	HIS	3.1
1	A	2	PRO	3.0
1	B	59	GLY	2.8
1	A	37	ASN	2.8
1	A	59	GLY	2.8
1	A	26	GLU	2.8
1	B	14	ASP	2.6
1	A	36	SER	2.4
1	A	10	SER	2.4
1	A	33	LYS	2.3
1	A	31	VAL	2.3
1	B	35	PHE	2.3
1	B	9	GLY	2.2
1	B	3	ASN	2.1
1	A	11	SER	2.1
1	B	10	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 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($\text{\AA}^2$)	Q<0.9
2	SO4	A	403	5/5	0.85	0.11	63,86,92,95	0
2	SO4	A	401	5/5	0.89	0.08	83,83,86,87	0
3	MG	B	1004	1/1	0.90	0.25	45,45,45,45	0
3	MG	A	404	1/1	0.91	0.34	54,54,54,54	0
2	SO4	B	1003	5/5	0.91	0.15	49,60,68,68	0
2	SO4	B	1001	5/5	0.93	0.07	69,72,79,87	0
2	SO4	B	1002	5/5	0.99	0.05	33,34,35,37	0
2	SO4	A	402	5/5	0.99	0.06	32,38,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.