



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 11:44 AM EDT

PDB ID : 4LZO / pdb_00004lzo
Title : Crystal structure of human PRS1 A87T mutant
Authors : Chen, P.; Teng, M.; Li, X.
Deposited on : 2013-07-31
Resolution : 3.31 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

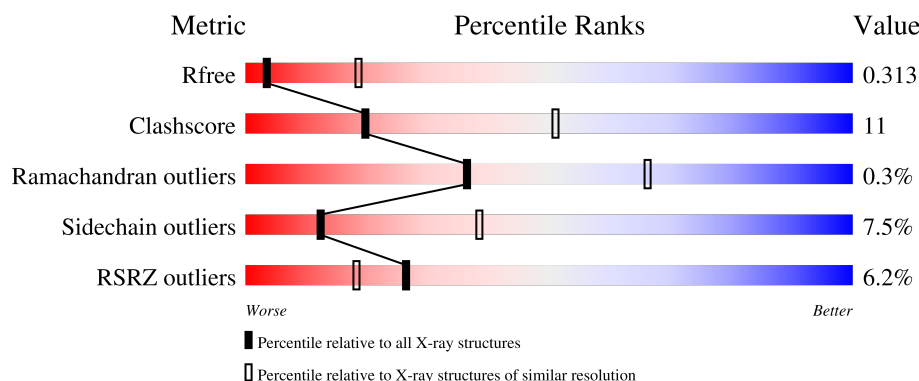
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.31 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	1153 (3.34-3.30)
Clashscore	190562	1193 (3.34-3.30)
Ramachandran outliers	187476	1172 (3.34-3.30)
Sidechain outliers	187428	1171 (3.34-3.30)
RSRZ outliers	180081	1153 (3.34-3.30)

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	326	
1	B	326	

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	B	401	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4714 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.

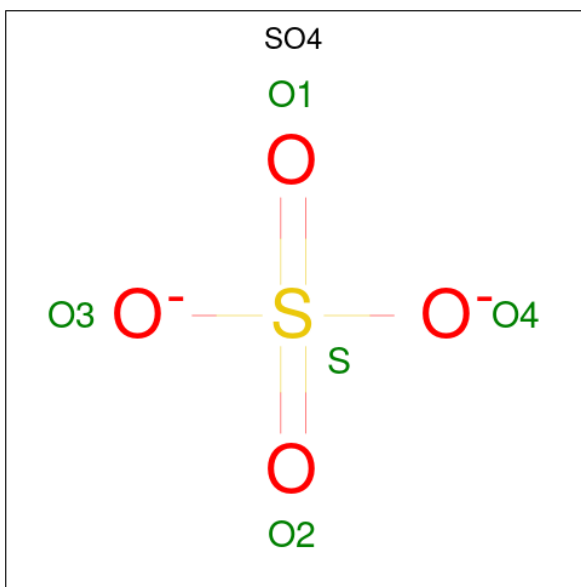
- Molecule 1 is a protein called Ribose-phosphate pyrophosphokinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2327	1459	408	443	17			
1	B	308	Total	C	N	O	S	0	0	0
			2357	1478	414	448	17			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	THR	ALA	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	87	THR	ALA	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).

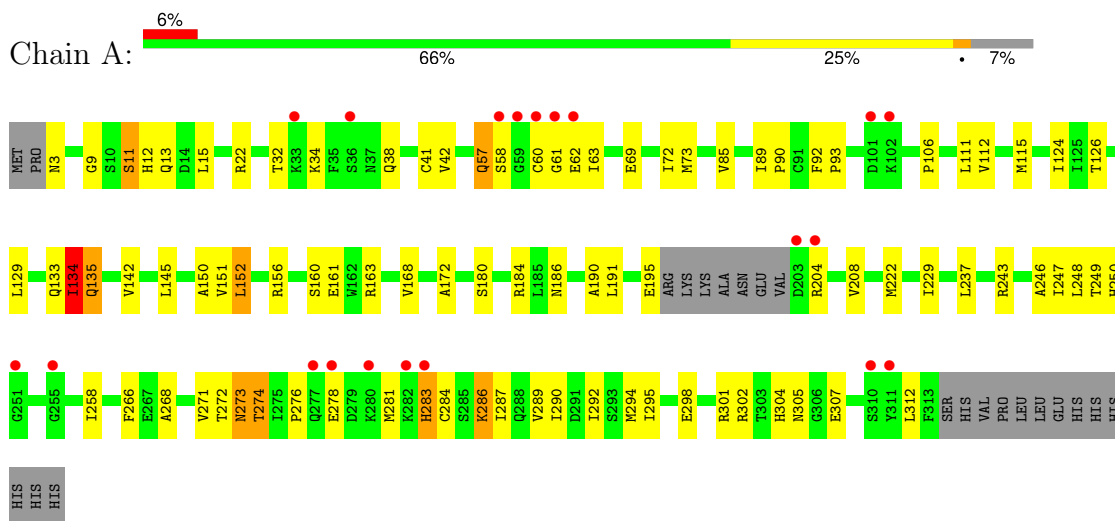


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

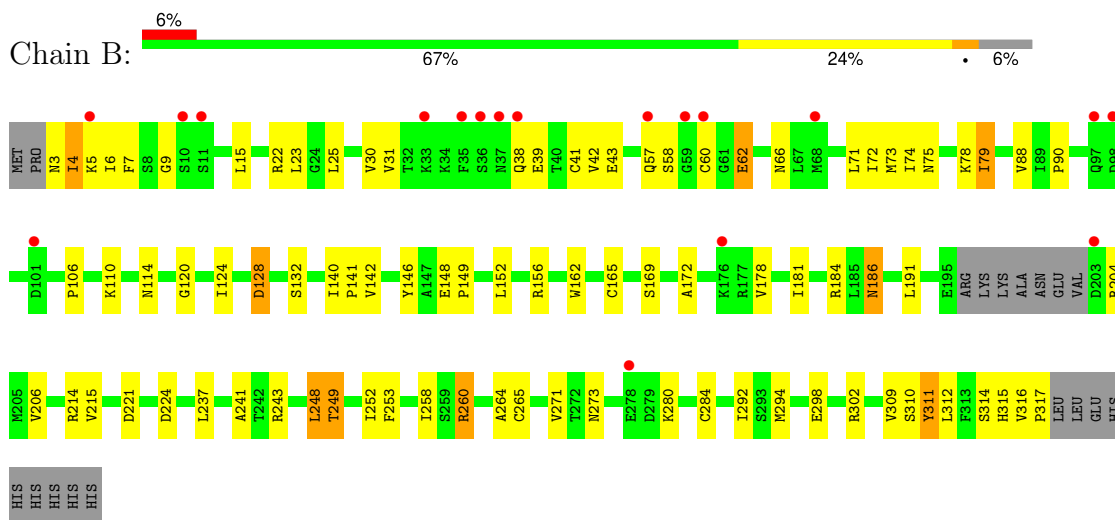
3 Residue-property plots [i](#)

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	170.40Å 170.40Å 62.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.18 – 3.31 34.18 – 3.31	Depositor EDS
% Data completeness (in resolution range)	99.8 (34.18-3.31) 99.9 (34.18-3.31)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.76 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.263 , 0.312 0.263 , 0.313	Depositor DCC
R_{free} test set	482 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	59.0	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.316 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4714	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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 ⓘ

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	0.51	0/2359	0.96	6/3189 (0.2%)
1	B	0.50	0/2391	0.85	1/3234 (0.0%)
All	All	0.51	0/4750	0.91	7/6423 (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	A	11	SER	N-CA-C	-19.35	81.36	108.07
1	A	135	GLN	N-CA-C	-9.99	102.16	114.75
1	B	39	GLU	N-CA-C	-6.22	100.77	110.42
1	A	13	GLN	N-CA-C	5.54	119.13	112.38
1	A	61	GLY	N-CA-C	-5.51	105.33	114.48
1	A	274	THR	N-CA-C	-5.46	106.13	112.89
1	A	134	ILE	CB-CA-C	-5.12	102.89	111.29

There are no chirality outliers.

There are no planarity outliers.

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	2327	0	2362	54	0
1	B	2357	0	2390	58	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	15	0	0	0	0
2	B	15	0	0	3	0
All	All	4714	0	4752	106	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 11.

All (106) 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:129:LEU:HD13	1:A:134:ILE:O	1.73	0.89
1:A:9:GLY:HA3	1:A:58:SER:HB2	1.55	0.89
1:A:12:HIS:CD2	1:A:276:PRO:HG3	2.09	0.88
1:A:12:HIS:CG	1:A:12:HIS:O	2.36	0.78
1:B:78:LYS:HE2	1:B:120:GLY:O	1.90	0.72
1:A:278:GLU:HA	1:A:281:MET:HE3	1.70	0.71
1:B:22:ARG:NH2	1:B:294:MET:HB3	2.08	0.69
1:B:148:GLU:O	1:B:152:LEU:HG	1.93	0.69
1:A:12:HIS:CD2	1:A:12:HIS:O	2.51	0.63
1:B:309:VAL:N	2:B:401:SO4:O3	2.32	0.62
1:A:172:ALA:HA	1:A:191:LEU:HD21	1.82	0.62
1:A:133:GLN:C	1:A:135:GLN:H	2.06	0.62
1:A:284:CYS:SG	1:A:287:ILE:HG13	2.40	0.61
1:A:112:VAL:HA	1:A:115:MET:HE2	1.82	0.61
1:B:9:GLY:O	1:B:58:SER:HB2	2.01	0.61
1:A:3:ASN:HD22	1:A:304:HIS:HE1	1.49	0.60
1:A:32:THR:HG22	1:A:42:VAL:HG22	1.84	0.59
1:B:124:ILE:HB	1:B:142:VAL:HG22	1.83	0.59
1:A:133:GLN:O	1:A:135:GLN:N	2.30	0.58
1:A:106:PRO:HG3	1:B:72:ILE:HG23	1.87	0.57
1:A:12:HIS:CD2	1:A:276:PRO:CG	2.84	0.55
1:B:6:ILE:HD11	1:B:23:LEU:CD1	2.35	0.55
1:A:12:HIS:HD2	1:A:276:PRO:HG3	1.66	0.55
1:B:149:PRO:HG3	1:B:314:SER:HA	1.88	0.55
1:A:126:THR:HG21	1:A:129:LEU:HD21	1.88	0.55
1:B:22:ARG:HH21	1:B:294:MET:HB3	1.72	0.54
1:A:145:LEU:HD13	1:A:295:ILE:HG22	1.90	0.54
1:B:294:MET:SD	1:B:294:MET:N	2.80	0.54
1:B:221:ASP:HA	1:B:249:THR:HG23	1.88	0.54
1:A:161:GLU:CD	1:A:243:ARG:HH21	2.17	0.53
1:A:298:GLU:HG2	1:A:312:LEU:HD21	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:GLN:C	1:A:135:GLN:N	2.67	0.53
1:B:6:ILE:HD11	1:B:23:LEU:HD12	1.91	0.52
1:B:273:ASN:OD1	1:B:292:ILE:HG13	2.09	0.52
1:B:310:SER:OG	2:B:401:SO4:O2	2.23	0.52
1:B:215:VAL:HG22	1:B:243:ARG:HG2	1.92	0.52
1:A:72:ILE:CG2	1:B:106:PRO:HG3	2.40	0.52
1:B:253:PHE:HB2	1:B:280:LYS:HZ3	1.75	0.52
1:B:152:LEU:HD21	1:B:181:ILE:HG23	1.93	0.51
1:A:161:GLU:OE1	1:A:243:ARG:NH2	2.44	0.51
1:B:156:ARG:HA	1:B:162:TRP:CD1	2.46	0.51
1:B:88:VAL:O	1:B:90:PRO:HD3	2.11	0.50
1:A:248:LEU:HB2	1:A:271:VAL:HG12	1.93	0.50
1:B:165:CYS:O	1:B:214:ARG:NH1	2.45	0.50
1:A:258:ILE:HG21	1:A:283:HIS:HB3	1.93	0.50
1:A:63:ILE:HG22	1:B:38:GLN:O	2.12	0.49
1:B:78:LYS:HG3	1:B:120:GLY:HA3	1.93	0.49
1:B:4:ILE:HG23	1:B:25:LEU:HD11	1.93	0.49
1:A:302:ARG:HA	1:A:307:GLU:HB3	1.95	0.48
1:B:258:ILE:HG23	1:B:284:CYS:HB2	1.95	0.48
1:A:124:ILE:HB	1:A:142:VAL:HG22	1.95	0.48
1:B:71:LEU:HA	1:B:74:ILE:HD12	1.96	0.48
1:A:222:MET:HB3	1:A:250:HIS:HB2	1.96	0.47
1:B:31:VAL:HB	1:B:43:GLU:HB3	1.96	0.47
1:B:148:GLU:OE2	1:B:184:ARG:NH1	2.48	0.47
1:B:298:GLU:O	1:B:302:ARG:HG2	2.14	0.47
1:B:140:ILE:HB	1:B:141:PRO:HD2	1.97	0.47
1:B:110:LYS:HE2	1:B:114:ASN:HD21	1.80	0.46
1:B:237:LEU:HD23	1:B:241:ALA:O	2.15	0.46
1:B:302:ARG:HB2	1:B:309:VAL:HG22	1.97	0.46
1:B:258:ILE:HD11	1:B:280:LYS:HB3	1.97	0.46
1:A:106:PRO:HG3	1:B:72:ILE:CG2	2.45	0.45
1:A:272:THR:OG1	1:A:274:THR:HG23	2.17	0.45
1:B:75:ASN:O	1:B:79:ILE:HG12	2.15	0.45
1:A:268:ALA:HA	1:A:286:LYS:HB2	1.98	0.45
1:B:7:PHE:HD2	1:B:30:VAL:HG23	1.79	0.45
1:A:151:VAL:HG13	1:A:247:ILE:HG21	1.99	0.45
1:A:57:GLN:NE2	1:A:90:PRO:HD2	2.32	0.44
1:B:152:LEU:CD2	1:B:181:ILE:HG23	2.47	0.44
1:B:237:LEU:HD12	1:B:265:CYS:SG	2.57	0.44
1:A:273:ASN:H	1:A:273:ASN:HD22	1.64	0.44
1:A:129:LEU:CD1	1:A:134:ILE:O	2.57	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:310:SER:C	1:B:312:LEU:H	2.26	0.44
1:A:302:ARG:HG2	1:A:307:GLU:HG2	1.99	0.44
1:A:72:ILE:HG23	1:B:106:PRO:HG3	1.99	0.44
1:A:246:ALA:HB2	1:A:266:PHE:CE2	2.53	0.44
1:A:152:LEU:HD21	1:A:184:ARG:HB3	1.99	0.44
1:A:152:LEU:HB3	1:A:156:ARG:HH12	1.83	0.43
1:B:42:VAL:HG11	1:B:73:MET:HG2	2.00	0.43
1:B:248:LEU:O	1:B:271:VAL:HA	2.19	0.43
1:A:22:ARG:HH11	1:A:294:MET:HE2	1.84	0.43
1:A:38:GLN:HG3	1:B:62:GLU:HA	2.00	0.43
1:A:271:VAL:O	1:A:289:VAL:HA	2.19	0.43
1:A:248:LEU:O	1:A:271:VAL:HA	2.19	0.42
1:A:301:ARG:HG2	1:A:305:ASN:ND2	2.35	0.42
1:B:146:TYR:OH	2:B:403:SO4:O3	2.33	0.42
1:B:253:PHE:HB2	1:B:280:LYS:NZ	2.32	0.42
1:A:69:GLU:O	1:A:73:MET:HG3	2.20	0.42
1:B:315:HIS:ND1	1:B:316:VAL:N	2.68	0.42
1:B:224:ASP:HA	1:B:252:ILE:HB	2.01	0.42
1:A:180:SER:O	1:A:184:ARG:HD3	2.20	0.42
1:B:110:LYS:HE2	1:B:114:ASN:ND2	2.35	0.42
1:B:224:ASP:O	1:B:253:PHE:HA	2.19	0.42
1:B:60:CYS:HB3	1:B:66:ASN:HD21	1.84	0.41
1:A:57:GLN:NE2	1:A:57:GLN:O	2.49	0.41
1:A:273:ASN:HD21	1:A:292:ILE:H	1.67	0.41
1:B:186:ASN:HD22	1:B:186:ASN:HA	1.69	0.41
1:A:34:LYS:HB3	1:A:34:LYS:HE2	1.86	0.41
1:A:168:VAL:HA	1:A:190:ALA:O	2.21	0.41
1:B:294:MET:HG2	1:B:317:PRO:HG2	2.02	0.41
1:A:150:ALA:HB1	1:A:290:ILE:HG21	2.02	0.41
1:B:128:ASP:OD2	1:B:146:TYR:HA	2.20	0.41
1:B:172:ALA:HA	1:B:191:LEU:HD21	2.02	0.41
1:B:260:ARG:O	1:B:264:ALA:HB2	2.21	0.41
1:A:92:PHE:HA	1:A:93:PRO:HD2	1.95	0.40
1:B:169:SER:HB2	1:B:178:VAL:HG21	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	300/326 (92%)	278 (93%)	21 (7%)	1 (0%)	36	66
1	B	304/326 (93%)	279 (92%)	24 (8%)	1 (0%)	36	66
All	All	604/652 (93%)	557 (92%)	45 (8%)	2 (0%)	36	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	ILE
1	B	311	TYR

5.3.2 Protein sidechains

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	258/279 (92%)	236 (92%)	22 (8%)	10	34
1	B	262/279 (94%)	245 (94%)	17 (6%)	15	43
All	All	520/558 (93%)	481 (92%)	39 (8%)	12	38

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	15	LEU
1	A	41	CYS
1	A	57	GLN

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Mol	Chain	Res	Type
1	A	60	CYS
1	A	62	GLU
1	A	85	VAL
1	A	89	ILE
1	A	111	LEU
1	A	152	LEU
1	A	160	SER
1	A	163	ARG
1	A	186	ASN
1	A	195	GLU
1	A	204	ARG
1	A	208	VAL
1	A	229	ILE
1	A	237	LEU
1	A	249	THR
1	A	273	ASN
1	A	283	HIS
1	A	286	LYS
1	B	3	ASN
1	B	4	ILE
1	B	5	LYS
1	B	15	LEU
1	B	41	CYS
1	B	57	GLN
1	B	62	GLU
1	B	79	ILE
1	B	128	ASP
1	B	132	SER
1	B	186	ASN
1	B	204	ARG
1	B	206	VAL
1	B	248	LEU
1	B	249	THR
1	B	260	ARG
1	B	311	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	13	GLN
1	A	57	GLN

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Mol	Chain	Res	Type
1	A	273	ASN
1	A	277	GLN
1	A	305	ASN
1	B	12	HIS
1	B	57	GLN
1	B	64	ASN
1	B	114	ASN
1	B	164	ASN
1	B	186	ASN
1	B	263	ASN
1	B	277	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	0.29	0	6,6,6	0.18	0
2	SO4	B	402	-	4,4,4	0.24	0	6,6,6	0.14	0
2	SO4	A	1002	-	4,4,4	0.25	0	6,6,6	0.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	1001	-	4,4,4	0.22	0	6,6,6	0.21	0
2	SO4	B	403	-	4,4,4	0.29	0	6,6,6	0.15	0
2	SO4	A	1003	-	4,4,4	0.25	0	6,6,6	0.17	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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	SO4	2	0
2	B	403	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	304/326 (93%)	-0.62	20 (6%) 24 16	0, 36, 56, 60	54 (17%)
1	B	308/326 (94%)	-0.70	18 (5%) 29 19	0, 32, 46, 57	49 (15%)
All	All	612/652 (93%)	-0.66	38 (6%) 26 18	0, 33, 52, 60	103 (16%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	CYS	9.4
1	B	36	SER	8.7
1	A	59	GLY	8.4
1	A	36	SER	8.4
1	B	101	ASP	7.5
1	A	278	GLU	7.1
1	B	203	ASP	6.8
1	A	203	ASP	6.7
1	A	310	SER	6.5
1	B	98	ASP	6.3
1	B	60	CYS	6.3
1	B	35	PHE	5.9
1	A	204	ARG	5.7
1	B	37	ASN	5.4
1	A	61	GLY	5.3
1	A	283	HIS	5.2
1	A	255	GLY	4.9
1	B	33	LYS	4.8
1	B	38	GLN	4.8
1	A	282	LYS	4.7
1	A	62	GLU	4.4
1	B	10	SER	4.4
1	A	101	ASP	4.0
1	B	97	GLN	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	311	TYR	3.9
1	A	58	SER	3.9
1	A	280	LYS	3.8
1	B	176	LYS	3.7
1	B	11	SER	3.5
1	B	68	MET	3.1
1	A	33	LYS	3.1
1	B	59	GLY	2.8
1	A	102	LYS	2.8
1	A	277	GLN	2.5
1	B	57	GLN	2.5
1	A	251	GLY	2.2
1	B	278	GLU	2.2
1	B	5	LYS	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	1002	5/5	0.96	0.10	34,34,35,35	5
2	SO4	A	1003	5/5	0.96	0.10	12,12,13,13	5
2	SO4	B	403	5/5	0.98	0.06	8,8,9,9	5
2	SO4	B	402	5/5	0.99	0.04	80,81,81,81	0
2	SO4	A	1001	5/5	1.00	0.02	30,30,30,30	0
2	SO4	B	401	5/5	1.00	0.02	23,23,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.