



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2021 – 11:47 PM EDT

PDB ID : 2H08
Title : crystal structure of human phosphoribosyl pyrophosphate synthetase 1 mutant Y146M
Authors : Li, S.; Peng, B.; Ding, J.
Deposited on : 2006-05-14
Resolution : 2.50 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 : 2.23.2
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) : 2.23.2

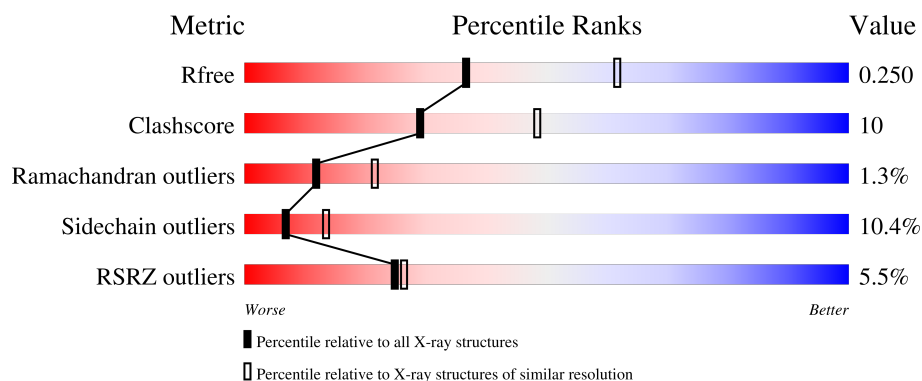
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4877 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 I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2332	1460	412	442	18			
1	B	308	Total	C	N	O	S	0	0	0
			2351	1473	414	446	18			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	146	MET	TYR	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	146	MET	TYR	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 (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	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		

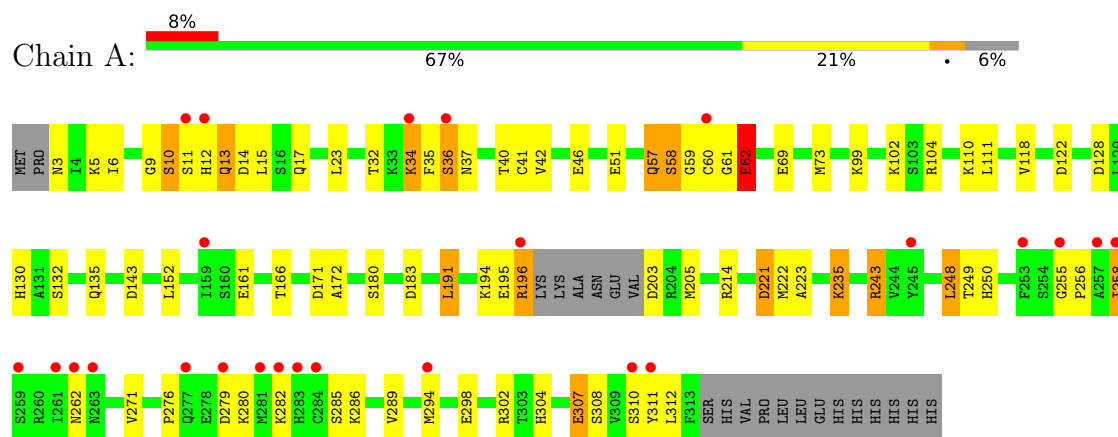
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	73	Total	O	0	0
			73	73		
3	B	91	Total	O	0	0
			91	91		

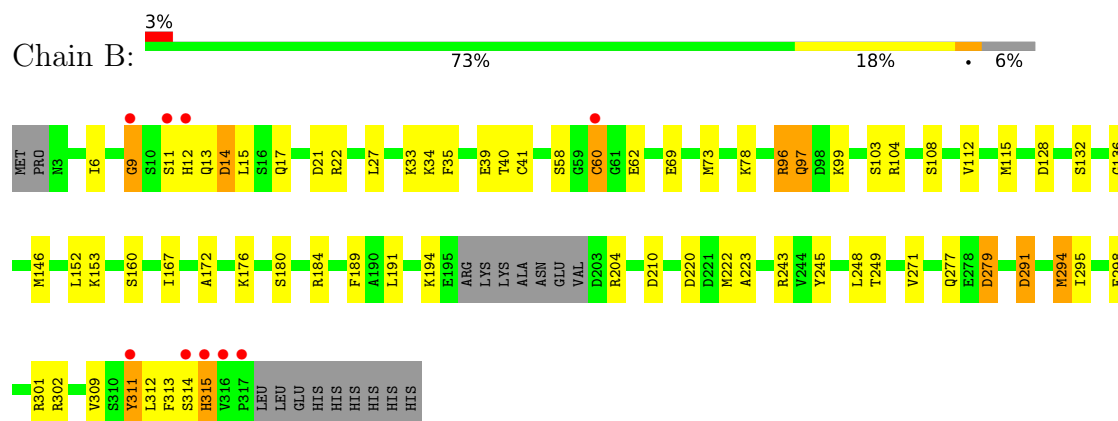
3 Residue-property plots [i](#)

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 I



• Molecule 1: Ribose-phosphate pyrophosphokinase I



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	169.91Å 169.91Å 61.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.05 – 2.50 49.05 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.05-2.50) 100.0 (49.05-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.207 , 0.257 0.199 , 0.250	Depositor DCC
R_{free} test set	1155 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4877	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.15% 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	0.44	0/2363	0.77	7/3192 (0.2%)
1	B	0.47	0/2384	0.75	4/3223 (0.1%)
All	All	0.45	0/4747	0.76	11/6415 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	171	ASP	CB-CG-OD2	6.28	123.96	118.30
1	B	210	ASP	CB-CG-OD2	6.21	123.89	118.30
1	B	128	ASP	CB-CG-OD2	5.81	123.53	118.30
1	A	143	ASP	CB-CG-OD2	5.80	123.52	118.30
1	A	221	ASP	CB-CG-OD2	5.77	123.50	118.30
1	A	183	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	279	ASP	CB-CG-OD2	5.32	123.08	118.30
1	A	122	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	128	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	291	ASP	CB-CG-OD2	5.15	122.93	118.30
1	B	279	ASP	CB-CG-OD2	5.06	122.85	118.30

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	2332	0	2373	50	0
1	B	2351	0	2388	44	0
2	A	15	0	0	2	0
2	B	15	0	0	0	0
3	A	73	0	0	5	0
3	B	91	0	0	0	0
All	All	4877	0	4761	91	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 10.

All (91) 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:172:ALA:HA	1:A:191:LEU:HD21	1.44	1.00
1:B:291:ASP:O	1:B:294:MET:HE3	1.68	0.93
1:A:10:SER:HB3	1:A:58:SER:OG	1.78	0.83
1:A:258:ILE:HG13	3:A:1075:HOH:O	1.83	0.78
1:B:291:ASP:O	1:B:294:MET:CE	2.31	0.78
1:B:298:GLU:O	1:B:302:ARG:HG2	1.84	0.77
1:A:102:LYS:HD3	1:B:146:MET:SD	2.31	0.71
1:A:110:LYS:HE3	1:B:136:GLY:O	1.96	0.66
1:A:130:HIS:HD2	1:A:221:ASP:OD1	1.79	0.65
1:A:10:SER:HB2	1:A:69:GLU:OE1	1.97	0.65
1:A:104:ARG:NH2	3:A:1022:HOH:O	2.32	0.63
1:B:220:ASP:HB3	1:B:248:LEU:HD12	1.80	0.63
1:B:220:ASP:HB3	1:B:248:LEU:CD1	2.28	0.63
1:A:34:LYS:HZ3	1:A:40:THR:HG23	1.62	0.62
1:A:42:VAL:HG11	1:A:73:MET:HG2	1.80	0.62
1:A:262:ASN:HA	1:A:286:LYS:HE3	1.81	0.61
1:B:96:ARG:HD3	1:B:222:MET:SD	2.40	0.61
1:A:307:GLU:HG3	1:A:308:SER:N	2.14	0.61
1:B:33:LYS:O	1:B:41:CYS:HB2	2.01	0.60
1:B:6:ILE:HG22	1:B:27:LEU:HD23	1.83	0.60
1:A:5:LYS:HE3	1:A:46:GLU:OE1	2.02	0.60
1:B:302:ARG:HB2	1:B:309:VAL:HG22	1.83	0.59
1:B:243:ARG:NH1	1:B:245:TYR:OH	2.31	0.59
1:B:9:GLY:HA2	1:B:73:MET:HE3	1.83	0.58
1:A:135:GLN:NE2	3:A:1020:HOH:O	2.37	0.57
1:A:12:HIS:CD2	1:A:60:CYS:HB3	2.41	0.55
1:A:223:ALA:HB2	1:A:248:LEU:HD22	1.89	0.55
1:A:32:THR:HG22	1:A:34:LYS:NZ	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:HIS:ND1	1:A:276:PRO:HG2	2.22	0.55
1:B:248:LEU:O	1:B:271:VAL:HA	2.08	0.54
1:A:172:ALA:CA	1:A:191:LEU:HD21	2.29	0.54
1:A:32:THR:HG22	1:A:34:LYS:HZ2	1.73	0.53
1:A:6:ILE:HD11	1:A:23:LEU:HD12	1.90	0.53
1:B:294:MET:SD	1:B:295:ILE:HG13	2.50	0.52
1:B:22:ARG:NE	1:B:294:MET:HB3	2.25	0.52
1:B:69:GLU:HG2	1:B:73:MET:HE2	1.90	0.52
1:A:248:LEU:O	1:A:271:VAL:HA	2.10	0.52
1:B:9:GLY:HA2	1:B:73:MET:CE	2.39	0.52
1:B:17:GLN:NE2	1:B:21:ASP:OD1	2.41	0.51
1:A:34:LYS:NZ	1:A:40:THR:HG23	2.26	0.51
1:A:9:GLY:HA2	1:A:73:MET:HE2	1.93	0.51
1:A:302:ARG:NH1	1:A:308:SER:O	2.38	0.51
1:B:12:HIS:CE1	1:B:60:CYS:HA	2.45	0.51
1:A:302:ARG:HG2	1:A:307:GLU:HG2	1.92	0.51
1:B:97:GLN:HB3	1:B:108:SER:HB2	1.92	0.51
1:A:59:GLY:HA2	3:A:1016:HOH:O	2.11	0.50
1:A:161:GLU:CD	1:A:243:ARG:HH12	2.16	0.49
1:A:3:ASN:HD22	1:A:304:HIS:HE1	1.59	0.49
1:A:196:ARG:NH2	3:A:1069:HOH:O	2.44	0.49
1:B:69:GLU:HG2	1:B:73:MET:CE	2.43	0.48
1:A:12:HIS:CE1	1:A:60:CYS:HA	2.48	0.48
1:B:172:ALA:HA	1:B:191:LEU:HD21	1.96	0.48
1:B:223:ALA:HB2	1:B:248:LEU:HG	1.95	0.47
1:B:152:LEU:HD11	1:B:184:ARG:HD2	1.95	0.47
1:A:161:GLU:OE2	1:A:243:ARG:NH1	2.48	0.47
1:A:34:LYS:HE2	1:A:40:THR:HA	1.97	0.46
1:A:99:LYS:HD3	1:B:132:SER:HB2	1.96	0.46
1:B:204:ARG:HG2	1:B:204:ARG:HH11	1.81	0.46
1:B:146:MET:CE	1:B:313:PHE:O	2.63	0.46
1:B:248:LEU:HB2	1:B:271:VAL:HG12	1.97	0.46
1:A:5:LYS:HG3	1:A:51:GLU:OE1	2.14	0.46
1:B:12:HIS:CD2	1:B:60:CYS:HA	2.51	0.45
1:A:12:HIS:HB2	1:A:59:GLY:H	1.81	0.45
1:B:12:HIS:NE2	1:B:60:CYS:HA	2.31	0.45
1:A:302:ARG:HD3	1:A:308:SER:O	2.17	0.45
1:A:3:ASN:ND2	1:A:304:HIS:HE1	2.15	0.44
1:A:205:MET:O	1:A:235:LYS:HE3	2.17	0.44
1:A:135:GLN:NE2	2:A:1006:SO4:O2	2.49	0.44
1:B:294:MET:H	1:B:294:MET:HG3	1.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:CYS:O	1:A:62:GLU:N	2.50	0.44
1:B:302:ARG:CB	1:B:309:VAL:HG22	2.47	0.44
1:B:13:GLN:O	1:B:14:ASP:C	2.56	0.44
1:A:130:HIS:CD2	1:A:221:ASP:OD1	2.64	0.44
1:B:35:PHE:HD1	1:B:39:GLU:HB3	1.84	0.43
1:B:311:TYR:O	1:B:315:HIS:HB2	2.18	0.43
1:A:298:GLU:HB3	1:A:312:LEU:HD11	2.01	0.43
1:A:9:GLY:HA2	1:A:73:MET:CE	2.48	0.42
1:B:298:GLU:O	1:B:302:ARG:CG	2.63	0.42
1:A:57:GLN:HG2	1:A:58:SER:O	2.19	0.42
2:A:1004:SO4:O3	1:B:104:ARG:NH1	2.52	0.42
1:B:298:GLU:OE1	1:B:301:ARG:NH1	2.53	0.42
1:B:112:VAL:HA	1:B:115:MET:HE2	2.02	0.41
1:B:298:GLU:HB3	1:B:312:LEU:HD21	2.01	0.41
1:B:167:ILE:O	1:B:189:PHE:HA	2.20	0.41
1:A:222:MET:HA	1:A:250:HIS:O	2.20	0.41
1:B:112:VAL:HA	1:B:115:MET:CE	2.50	0.41
1:B:34:LYS:HD3	1:B:40:THR:HG23	2.03	0.41
1:A:13:GLN:O	1:A:17:GLN:HB2	2.21	0.40
1:A:255:GLY:HA3	1:A:256:PRO:HD3	1.81	0.40
1:A:35:PHE:HB3	1:A:36:SER:H	1.52	0.40
1:A:271:VAL:O	1:A:289:VAL:HA	2.21	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	301/326 (92%)	280 (93%)	16 (5%)	5 (2%)	9	16
1	B	304/326 (93%)	286 (94%)	15 (5%)	3 (1%)	15	28
All	All	605/652 (93%)	566 (94%)	31 (5%)	8 (1%)	12	21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	SER
1	B	14	ASP
1	A	61	GLY
1	A	62	GLU
1	B	9	GLY
1	A	10	SER
1	B	11	SER
1	A	258	ILE

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	258/278 (93%)	225 (87%)	33 (13%)	4	8
1	B	261/278 (94%)	240 (92%)	21 (8%)	12	23
All	All	519/556 (93%)	465 (90%)	54 (10%)	7	13

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	13	GLN
1	A	14	ASP
1	A	15	LEU
1	A	34	LYS
1	A	37	ASN
1	A	41	CYS
1	A	57	GLN
1	A	58	SER
1	A	62	GLU
1	A	111	LEU
1	A	118	VAL
1	A	132	SER
1	A	152	LEU
1	A	166	THR

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Mol	Chain	Res	Type
1	A	180	SER
1	A	191	LEU
1	A	194	LYS
1	A	195	GLU
1	A	196	ARG
1	A	203	ASP
1	A	214	ARG
1	A	235	LYS
1	A	243	ARG
1	A	248	LEU
1	A	249	THR
1	A	280	LYS
1	A	282	LYS
1	A	285	SER
1	A	294	MET
1	A	307	GLU
1	A	310	SER
1	A	311	TYR
1	B	15	LEU
1	B	58	SER
1	B	60	CYS
1	B	62	GLU
1	B	78	LYS
1	B	96	ARG
1	B	97	GLN
1	B	99	LYS
1	B	103	SER
1	B	153	LYS
1	B	160	SER
1	B	176	LYS
1	B	180	SER
1	B	194	LYS
1	B	249	THR
1	B	277	GLN
1	B	279	ASP
1	B	294	MET
1	B	311	TYR
1	B	314	SER
1	B	315	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	37	ASN
1	A	57	GLN
1	A	130	HIS
1	A	135	GLN
1	A	158	ASN
1	A	277	GLN
1	B	38	GLN
1	B	64	ASN
1	B	263	ASN

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](#)

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	1005	-	4,4,4	0.09	0	6,6,6	0.21	0
2	SO4	B	1001	-	4,4,4	0.18	0	6,6,6	0.36	0
2	SO4	B	1003	-	4,4,4	0.14	0	6,6,6	0.20	0
2	SO4	A	1002	-	4,4,4	0.16	0	6,6,6	0.11	0
2	SO4	A	1004	-	4,4,4	0.15	0	6,6,6	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	1006	-	4,4,4	0.15	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1004	SO4	1	0
2	A	1006	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	305/326 (93%)	0.31	25 (8%) 11 11	21, 42, 75, 84	0
1	B	308/326 (94%)	-0.12	9 (2%) 51 55	20, 33, 63, 98	0
All	All	613/652 (94%)	0.09	34 (5%) 25 26	20, 37, 70, 98	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	258	ILE	5.3
1	B	60	CYS	5.3
1	B	316	VAL	5.2
1	A	262	ASN	4.5
1	A	284	CYS	4.4
1	A	263	ASN	4.3
1	B	317	PRO	4.3
1	A	60	CYS	4.1
1	A	282	LYS	3.8
1	A	311	TYR	3.6
1	B	314	SER	3.6
1	B	9	GLY	3.5
1	A	310	SER	3.5
1	B	315	HIS	3.4
1	A	259	SER	3.3
1	A	277	GLN	3.3
1	B	12	HIS	3.2
1	A	245	TYR	3.1
1	A	261	ILE	2.9
1	A	283	HIS	2.9
1	B	11	SER	2.8
1	A	159	ILE	2.8
1	A	257	ALA	2.7
1	A	34	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	281	MET	2.6
1	A	36	SER	2.5
1	A	294	MET	2.5
1	A	279	ASP	2.4
1	A	11	SER	2.4
1	A	255	GLY	2.3
1	B	311	TYR	2.2
1	A	196	ARG	2.2
1	A	12	HIS	2.0
1	A	253	PHE	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 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, 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(\AA^2)	Q<0.9
2	SO4	A	1006	5/5	0.93	0.30	66,66,67,67	5
2	SO4	B	1001	5/5	0.95	0.18	62,63,63,64	0
2	SO4	A	1002	5/5	0.96	0.13	78,79,79,79	0
2	SO4	B	1005	5/5	0.98	0.13	44,44,44,45	0
2	SO4	B	1003	5/5	0.99	0.10	32,33,33,34	0
2	SO4	A	1004	5/5	0.99	0.10	37,38,38,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.