



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:57 AM GMT

PDB ID : 2YWV  
Title : Crystal structure of SAICAR synthetase from *Geobacillus kaustophilus*  
Authors : Kanagawa, M.; Baba, S.; Kuramitsu, S.; Yokoyama, S.; Kawai, G.; Sampei, G.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2007-04-23  
Resolution : 1.75 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

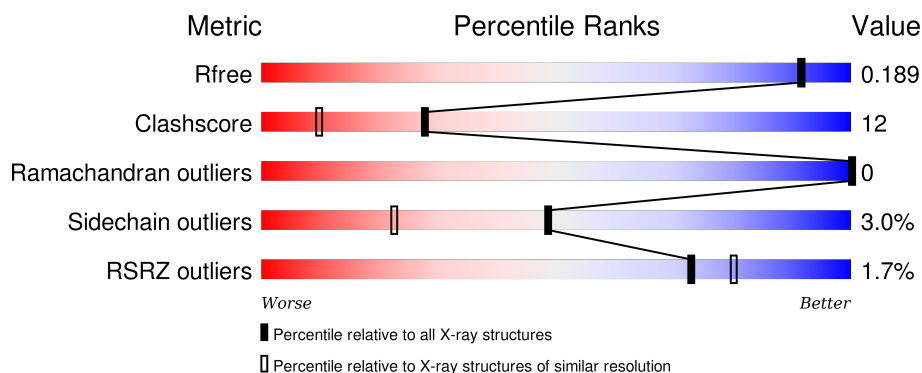
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

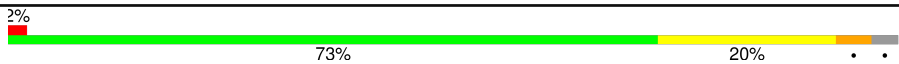
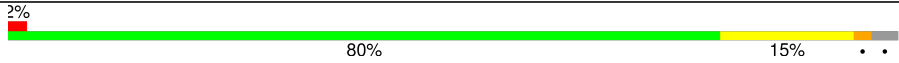
The reported resolution of this entry is 1.75 Å.

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))
$R_{free}$	91344	1609 (1.76-1.76)
Clashscore	102246	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)
RSRZ outliers	91569	1610 (1.76-1.76)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	244	
1	B	244	

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	545	-	-	-	X
2	SO4	B	548	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4370 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 Phosphoribosylaminoimidazole succinocarboxamide synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1877	1199	322	354	2			
1	B	237	Total	C	N	O	S	0	0	0
			1877	1199	322	354	2			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q5L3D5
A	0	HIS	-	EXPRESSION TAG	UNP Q5L3D5
B	-1	GLY	-	EXPRESSION TAG	UNP Q5L3D5
B	0	HIS	-	EXPRESSION TAG	UNP Q5L3D5

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).

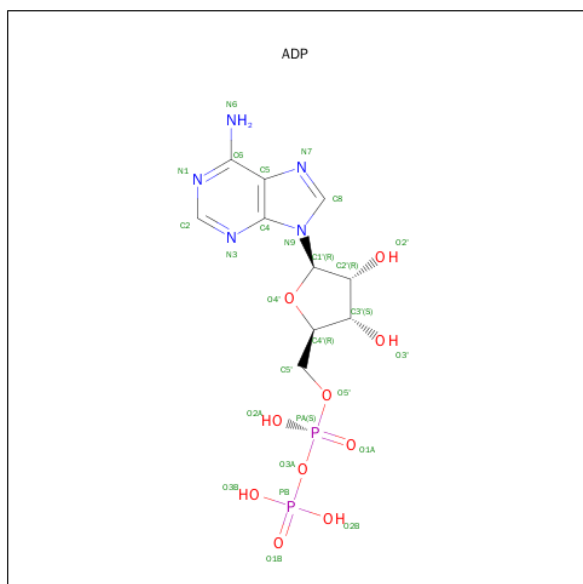


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

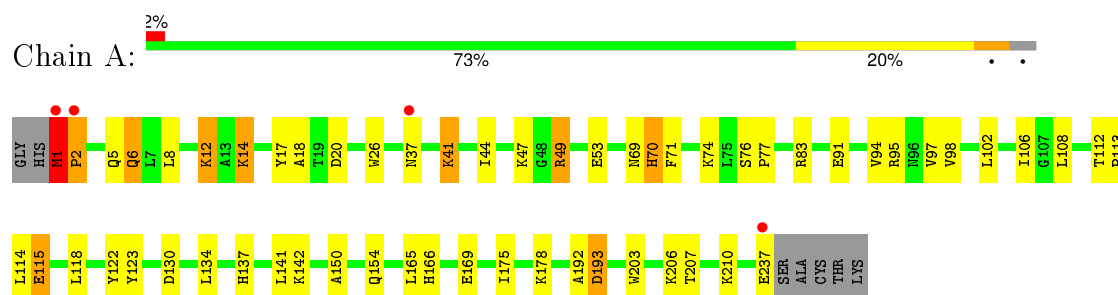
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	263	Total	O	0	0
			263	263		
5	B	262	Total	O	0	0
			262	262		

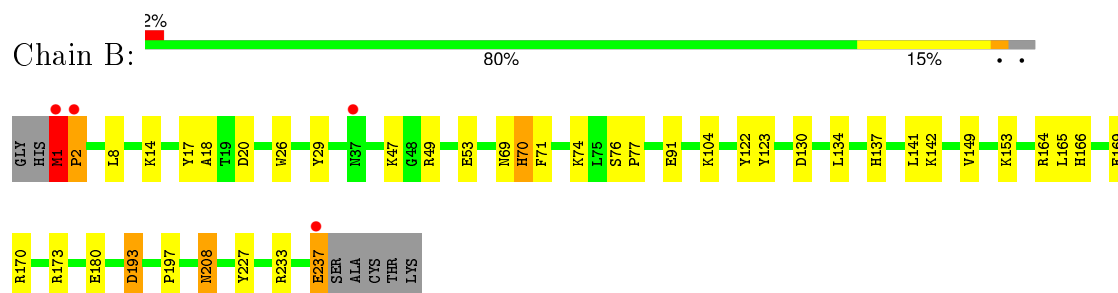
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Phosphoribosylaminoimidazole succinocarboxamide synthetase



- Molecule 1: Phosphoribosylaminoimidazole succinocarboxamide synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.89Å 56.40Å 56.41Å 76.39° 64.36° 64.35°	Depositor
Resolution (Å)	37.00 – 1.75 41.64 – 1.76	Depositor EDS
% Data completeness (in resolution range)	93.9 (37.00-1.75) 92.4 (41.64-1.76)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.31 (at 1.76Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.188 , 0.191 0.188 , 0.189	Depositor DCC
$R_{free}$ test set	4714 reflections (10.08%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.4	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
Estimated twinning fraction	0.020 for h,h-k,h-l 0.480 for -h,-l,-k 0.018 for -h,-h+l,-h+k	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 47381 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4370	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.30% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, SO4, ADP

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	0.29	0/1906	1.23	17/2576 (0.7%)
1	B	0.30	0/1906	0.69	4/2576 (0.2%)
All	All	0.30	0/3812	1.00	21/5152 (0.4%)

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	4

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	12	LYS	CA-CB-CG	26.54	171.78	113.40
1	A	41	LYS	CA-CB-CG	20.50	158.51	113.40
1	A	192	ALA	C-N-CA	17.88	166.39	121.70
1	A	192	ALA	O-C-N	-15.32	98.19	122.70
1	A	2	PRO	O-C-N	15.11	146.88	122.70
1	B	2	PRO	CB-CA-C	14.03	147.07	112.00
1	A	2	PRO	CA-C-N	-12.02	90.75	117.20
1	A	193	ASP	CB-CA-C	11.56	133.53	110.40
1	A	115	GLU	CA-CB-CG	-11.21	88.74	113.40
1	A	14	LYS	CG-CD-CE	10.72	144.07	111.90
1	A	192	ALA	CA-C-N	10.49	140.28	117.20
1	A	2	PRO	C-N-CA	-9.78	97.26	121.70
1	A	1	MET	O-C-N	-9.39	103.25	121.10
1	A	193	ASP	CA-C-N	-8.71	98.04	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	ASP	N-CA-C	-8.54	87.95	111.00
1	A	1	MET	CA-C-N	6.88	136.38	117.10
1	B	2	PRO	N-CA-C	-5.85	96.89	112.10
1	A	193	ASP	CA-C-O	5.69	132.06	120.10
1	B	1	MET	CB-CA-C	5.33	121.06	110.40
1	A	1	MET	CB-CA-C	5.33	121.06	110.40
1	B	193	ASP	O-C-N	5.20	131.02	122.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	MET	Mainchain,Peptide
1	A	193	ASP	Mainchain
1	A	83	ARG	Sidechain

## 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	1877	0	1927	49	0
1	B	1877	0	1928	41	0
2	A	15	0	0	0	0
2	B	20	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	27	0	12	3	0
4	B	27	0	12	3	0
5	A	263	0	0	16	0
5	B	262	0	0	11	0
All	All	4370	0	3879	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 12.

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:B:237:GLU:OE2	5:B:839:HOH:O	1.55	1.20
1:B:237:GLU:HG2	5:B:816:HOH:O	1.49	1.13
1:B:1:MET:H3	1:B:2:PRO:HD3	1.03	1.12
1:B:1:MET:N	1:B:2:PRO:HD3	1.61	1.06
1:A:74:LYS:NZ	5:A:910:HOH:O	1.89	1.05
1:B:53:GLU:OE1	5:B:855:HOH:O	1.80	0.99
1:B:1:MET:N	1:B:2:PRO:CD	2.27	0.97
1:B:74:LYS:NZ	5:B:828:HOH:O	1.98	0.95
2:B:546:SO4:O3	5:B:799:HOH:O	1.88	0.92
1:A:97:VAL:HA	5:A:818:HOH:O	1.70	0.91
1:A:115:GLU:O	1:A:115:GLU:HG3	1.65	0.87
1:A:6:GLN:NE2	1:A:6:GLN:H	1.72	0.87
1:B:70:HIS:HD2	4:B:601:ADP:HN61	1.18	0.86
1:A:53:GLU:OE2	1:A:237:GLU:HG2	1.78	0.83
1:A:70:HIS:HD2	4:A:600:ADP:HN61	1.23	0.83
1:B:2:PRO:HA	1:B:20:ASP:OD1	1.80	0.81
1:A:76:SER:HB2	1:A:77:PRO:HD2	1.71	0.73
1:A:154:GLN:NE2	5:A:730:HOH:O	2.23	0.71
1:A:70:HIS:CD2	4:A:600:ADP:HN61	2.09	0.70
1:B:53:GLU:CD	5:B:855:HOH:O	2.26	0.69
1:B:70:HIS:CD2	4:B:601:ADP:HN61	2.06	0.68
1:A:49:ARG:HG3	1:A:77:PRO:HB2	1.75	0.67
1:B:76:SER:HB2	1:B:77:PRO:HD2	1.75	0.67
1:A:203:TRP:NE1	5:A:827:HOH:O	2.28	0.67
1:A:150:ALA:O	1:A:154:GLN:HG2	1.95	0.66
1:B:53:GLU:CD	1:B:237:GLU:HB3	2.17	0.65
1:A:6:GLN:H	1:A:6:GLN:HE21	1.44	0.65
1:B:153:LYS:HB3	5:B:797:HOH:O	1.96	0.64
1:B:69:ASN:HD22	1:B:71:PHE:H	1.44	0.64
1:A:206:LYS:HE3	1:A:207:THR:HG23	1.81	0.63
1:A:154:GLN:CD	5:A:792:HOH:O	2.36	0.63
1:A:69:ASN:HD22	1:A:71:PHE:H	1.46	0.63
1:B:165:LEU:O	1:B:169:GLU:HG3	2.00	0.62
1:B:237:GLU:CG	5:B:816:HOH:O	2.24	0.62
1:B:149:VAL:O	1:B:153:LYS:HG3	2.00	0.62
1:B:1:MET:H2	1:B:2:PRO:CD	2.10	0.61
1:B:49:ARG:HH11	1:B:49:ARG:HB3	1.64	0.61
1:B:164:ARG:NH1	5:B:962:HOH:O	2.12	0.61
1:A:47:LYS:HD3	5:A:813:HOH:O	2.00	0.61
1:A:53:GLU:CD	1:A:237:GLU:HG2	2.21	0.61
1:A:134:LEU:H	1:A:137:HIS:CD2	2.19	0.61
1:B:53:GLU:OE1	1:B:237:GLU:HB3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:GLN:HB3	1:A:6:GLN:NE2	2.17	0.60
1:B:134:LEU:H	1:B:137:HIS:CD2	2.20	0.59
1:B:173:ARG:HH11	1:B:208:ASN:HD21	1.51	0.58
1:B:1:MET:H3	1:B:2:PRO:CD	1.90	0.58
1:A:49:ARG:O	1:A:53:GLU:HG3	2.05	0.57
1:A:8:LEU:HD11	1:A:18:ALA:HB2	1.87	0.57
1:A:94:VAL:HG13	5:A:860:HOH:O	2.04	0.56
1:A:210:LYS:NZ	5:A:827:HOH:O	2.29	0.55
1:B:70:HIS:HD2	4:B:601:ADP:N6	1.99	0.55
1:A:106:ILE:HD12	1:A:108:LEU:HD22	1.90	0.54
1:B:14:LYS:HD3	1:B:29:TYR:CZ	2.44	0.53
1:A:113:PRO:HA	5:A:818:HOH:O	2.09	0.53
1:A:114:LEU:N	5:A:818:HOH:O	2.42	0.52
1:A:91:GLU:HB2	1:A:122:TYR:HB2	1.91	0.51
1:B:91:GLU:HB2	1:B:122:TYR:HB2	1.91	0.51
1:B:166:HIS:HE1	5:B:718:HOH:O	1.93	0.51
1:B:8:LEU:HD11	1:B:18:ALA:HB2	1.92	0.51
1:A:118:LEU:N	5:A:860:HOH:O	2.45	0.50
1:A:165:LEU:O	1:A:169:GLU:HG3	2.11	0.50
1:A:134:LEU:H	1:A:137:HIS:HD2	1.60	0.50
1:A:44:ILE:HB	5:A:813:HOH:O	2.11	0.50
1:A:17:TYR:HB2	1:A:26:TRP:HB3	1.92	0.49
1:A:70:HIS:HD2	4:A:600:ADP:N6	2.01	0.49
1:A:166:HIS:HE1	5:A:719:HOH:O	1.95	0.49
1:A:6:GLN:N	1:A:6:GLN:HE21	2.11	0.49
1:B:14:LYS:HD3	1:B:29:TYR:CE2	2.48	0.48
1:B:134:LEU:H	1:B:137:HIS:HD2	1.61	0.47
1:A:98:VAL:HB	1:A:112:THR:HB	1.97	0.47
1:B:104:LYS:HB2	1:B:104:LYS:NZ	2.29	0.47
1:A:97:VAL:HB	1:A:175:ILE:HG23	1.96	0.46
1:B:123:TYR:O	1:B:130:ASP:HA	2.16	0.46
1:A:154:GLN:NE2	5:A:792:HOH:O	2.48	0.46
1:A:6:GLN:NE2	1:A:6:GLN:N	2.53	0.46
1:B:47:LYS:HE3	5:B:829:HOH:O	2.16	0.45
1:B:180:GLU:HB2	1:B:193:ASP:HB2	1.99	0.45
1:B:208:ASN:HA	1:B:208:ASN:HD22	1.63	0.44
1:A:178:LYS:NZ	5:A:812:HOH:O	2.49	0.43
1:A:123:TYR:O	1:A:130:ASP:HA	2.18	0.43
1:A:91:GLU:OE1	1:A:178:LYS:HD2	2.18	0.43
1:B:17:TYR:HB2	1:B:26:TRP:HB3	1.99	0.43
1:B:170:ARG:HD3	1:B:233:ARG:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:PRO:HA	1:A:20:ASP:OD1	2.20	0.42
1:B:197:PRO:HB2	1:B:227:TYR:CZ	2.55	0.42
1:A:95:ARG:HD2	1:A:102:LEU:HD22	2.02	0.41
1:A:102:LEU:O	1:A:106:ILE:HG12	2.20	0.41
1:A:114:LEU:HG	5:A:818:HOH:O	2.20	0.41
1:B:141:LEU:O	1:B:142:LYS:HB2	2.21	0.41
1:A:141:LEU:O	1:A:142:LYS:HB2	2.21	0.40
1:A:53:GLU:CD	1:A:237:GLU:CG	2.90	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	235/244 (96%)	226 (96%)	9 (4%)	0	100	100
1	B	235/244 (96%)	225 (96%)	10 (4%)	0	100	100
All	All	470/488 (96%)	451 (96%)	19 (4%)	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	200/205 (98%)	192 (96%)	8 (4%)	38	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	200/205 (98%)	196 (98%)	4 (2%)	63	39
All	All	400/410 (98%)	388 (97%)	12 (3%)	48	22

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	6	GLN
1	A	12	LYS
1	A	14	LYS
1	A	37	ASN
1	A	41	LYS
1	A	49	ARG
1	A	70	HIS
1	B	1	MET
1	B	70	HIS
1	B	208	ASN
1	B	237	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	37	ASN
1	A	69	ASN
1	A	70	HIS
1	A	80	GLN
1	A	137	HIS
1	A	154	GLN
1	A	166	HIS
1	A	171	ASN
1	B	37	ASN
1	B	69	ASN
1	B	70	HIS
1	B	80	GLN
1	B	137	HIS
1	B	166	HIS
1	B	171	ASN
1	B	208	ASN

### 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 ⓘ

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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	A	543	-	4,4,4	1.52	0	6,6,6	0.93	0
2	SO4	A	544	-	4,4,4	1.49	0	6,6,6	0.97	0
2	SO4	A	545	-	4,4,4	3.39	2 (50%)	6,6,6	0.94	0
4	ADP	A	600	-	22,29,29	1.35	1 (4%)	27,45,45	3.48	11 (40%)
2	SO4	B	541	-	4,4,4	1.46	0	6,6,6	0.93	0
2	SO4	B	542	-	4,4,4	1.50	0	6,6,6	0.95	0
2	SO4	B	546	-	4,4,4	3.40	2 (50%)	6,6,6	0.95	0
2	SO4	B	548	-	4,4,4	3.40	2 (50%)	6,6,6	0.94	0
4	ADP	B	601	-	22,29,29	1.38	1 (4%)	27,45,45	3.44	11 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	543	-	-	0/0/0/0	0/0/0/0
2	SO4	A	544	-	-	0/0/0/0	0/0/0/0
2	SO4	A	545	-	-	0/0/0/0	0/0/0/0
4	ADP	A	600	-	-	0/12/32/32	0/3/3/3
2	SO4	B	541	-	-	0/0/0/0	0/0/0/0
2	SO4	B	542	-	-	0/0/0/0	0/0/0/0
2	SO4	B	546	-	-	0/0/0/0	0/0/0/0
2	SO4	B	548	-	-	0/0/0/0	0/0/0/0
4	ADP	B	601	-	-	0/12/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	548	SO4	O3-S	-4.88	1.29	1.47
2	B	546	SO4	O3-S	-4.88	1.29	1.47
2	A	545	SO4	O3-S	-4.88	1.29	1.47
4	A	600	ADP	O4'-C1'	4.53	1.46	1.41
4	B	601	ADP	O4'-C1'	4.58	1.47	1.41
2	B	548	SO4	O1-S	4.63	1.63	1.47
2	A	545	SO4	O1-S	4.63	1.63	1.47
2	B	546	SO4	O1-S	4.66	1.63	1.47

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	600	ADP	N3-C2-N1	-10.00	121.24	128.89
4	B	601	ADP	N3-C2-N1	-9.94	121.29	128.89
4	B	601	ADP	O3A-PA-O5'	-6.46	85.80	102.94
4	A	600	ADP	O3A-PA-O5'	-6.35	86.09	102.94
4	A	600	ADP	PA-O3A-PB	-5.72	113.47	132.67
4	B	601	ADP	O5'-PA-O1A	-5.62	87.79	109.62
4	A	600	ADP	O5'-PA-O1A	-5.60	87.89	109.62
4	B	601	ADP	PA-O3A-PB	-5.49	114.27	132.67
4	A	600	ADP	O2A-PA-O5'	-5.25	82.00	108.46
4	B	601	ADP	O2A-PA-O5'	-5.22	82.13	108.46
4	A	600	ADP	O3B-PB-O3A	-4.60	84.24	105.09
4	B	601	ADP	O2B-PB-O3A	-4.50	84.70	105.09
4	A	600	ADP	O2B-PB-O3A	-4.15	86.27	105.09
4	B	601	ADP	O3B-PB-O3A	-3.60	88.76	105.09
4	B	601	ADP	C4-C5-N7	-2.56	107.13	109.48
4	A	600	ADP	C4-C5-N7	-2.55	107.13	109.48
4	A	600	ADP	O3B-PB-O1B	2.15	117.50	110.58
4	B	601	ADP	O3B-PB-O2B	2.70	117.67	107.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	ADP	O2A-PA-O3A	2.72	117.44	105.09
4	A	600	ADP	O2A-PA-O3A	2.78	117.69	105.09
4	B	601	ADP	O2B-PB-O1B	4.43	124.85	110.58
4	A	600	ADP	O2B-PB-O1B	4.53	125.17	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	600	ADP	3	0
2	B	546	SO4	1	0
4	B	601	ADP	3	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	237/244 (97%)	-0.12	4 (1%) 73 80	11, 18, 33, 50	0
1	B	237/244 (97%)	-0.15	4 (1%) 73 80	12, 18, 34, 48	0
All	All	474/488 (97%)	-0.14	8 (1%) 73 80	11, 18, 34, 50	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	3.6
1	B	37	ASN	3.1
1	A	2	PRO	2.9
1	A	37	ASN	2.5
1	B	2	PRO	2.2
1	A	237	GLU	2.2
1	A	1	MET	2.0
1	B	237	GLU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	SO4	A	545	5/5	0.92	0.21	10.46	66,66,67,67	0
2	SO4	B	548	5/5	0.92	0.30	9.72	93,93,94,94	0
2	SO4	B	541	5/5	0.98	0.08	0.28	29,30,31,32	0
2	SO4	A	543	5/5	0.97	0.09	0.05	30,31,32,34	0
4	ADP	B	601	27/27	0.96	0.09	-0.33	12,15,20,26	0
4	ADP	A	600	27/27	0.96	0.09	-0.35	11,14,20,25	0
3	MG	B	702	1/1	0.94	0.10	-0.42	22,22,22,22	0
2	SO4	B	542	5/5	0.97	0.07	-0.83	32,33,33,35	0
2	SO4	A	544	5/5	0.97	0.08	-0.96	36,37,37,38	0
2	SO4	B	546	5/5	0.87	0.27	-	72,72,73,73	0
3	MG	A	701	1/1	0.91	0.11	-	23,23,23,23	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.