



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 8, 2021 – 06:22 PM EST

PDB ID : 6VR0  
Title : Agrobacterium Tumefaciens ADP-glucose pyrophosphorylase W106A  
Authors : Mascarenhas, R.N.; Liu, D.; Ballicora, M.; Iglesias, A.; Asencion, M.; Figueroa, C.  
Deposited on : 2020-02-06  
Resolution : 2.20 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.17  
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) : 2.17

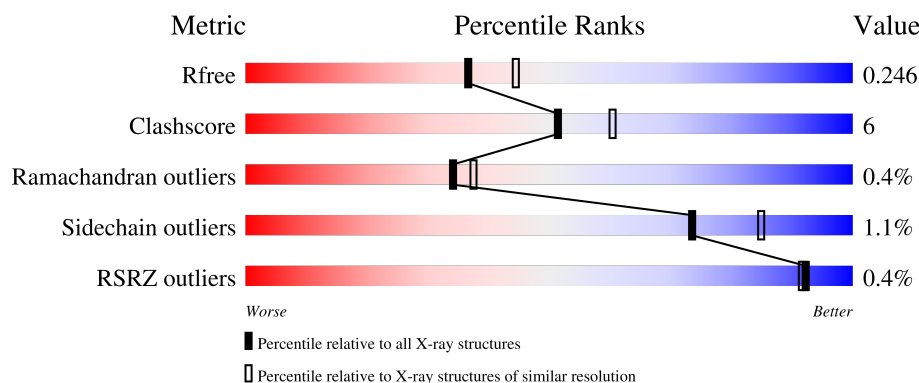
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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  $\geq 3$ , 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	413	
1	B	413	
1	C	413	
1	D	413	
1	E	413	

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Mol	Chain	Length	Quality of chain
1	F	413	
1	G	413	
1	H	413	
1	I	413	
1	J	413	

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	H	501	-	-	X	-
2	SO4	H	506	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 32581 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 Glucose-1-phosphate adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3174	2014	551	596	13			
1	B	403	Total	C	N	O	S	0	0	0
			3169	2011	550	595	13			
1	C	404	Total	C	N	O	S	0	0	0
			3167	2008	551	595	13			
1	D	404	Total	C	N	O	S	0	0	0
			3167	2008	551	595	13			
1	E	406	Total	C	N	O	S	0	0	0
			3186	2022	554	597	13			
1	F	403	Total	C	N	O	S	0	0	0
			3164	2008	548	595	13			
1	G	404	Total	C	N	O	S	0	0	0
			3174	2014	551	596	13			
1	H	404	Total	C	N	O	S	0	0	0
			3174	2014	551	596	13			
1	I	404	Total	C	N	O	S	0	0	0
			3174	2014	551	596	13			
1	J	404	Total	C	N	O	S	0	0	0
			3155	1999	549	594	13			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
B	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
C	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
D	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
E	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
F	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
G	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
H	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5
I	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5

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Chain	Residue	Modelled	Actual	Comment	Reference
J	67	ALA	GLN	engineered mutation	UNP A0A083ZTG5

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



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	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		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
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		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	I	1	Total	C	O	0	0
			6	3	3		

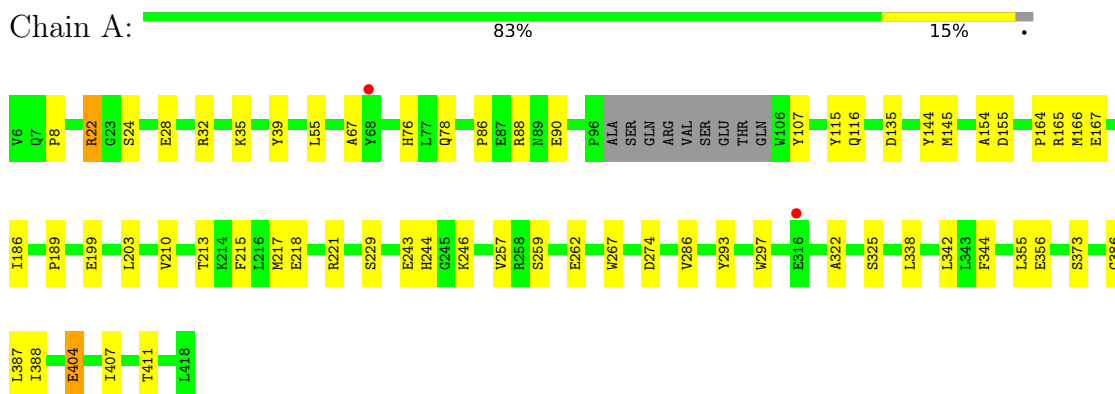
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	32	Total	O	0	0
			32	32		
4	B	23	Total	O	0	0
			23	23		
4	C	41	Total	O	0	0
			41	41		
4	D	20	Total	O	0	0
			20	20		
4	E	44	Total	O	0	0
			44	44		
4	F	42	Total	O	0	0
			42	42		
4	G	57	Total	O	0	0
			57	57		
4	H	45	Total	O	0	0
			45	45		
4	I	70	Total	O	0	0
			70	70		
4	J	52	Total	O	0	0
			52	52		

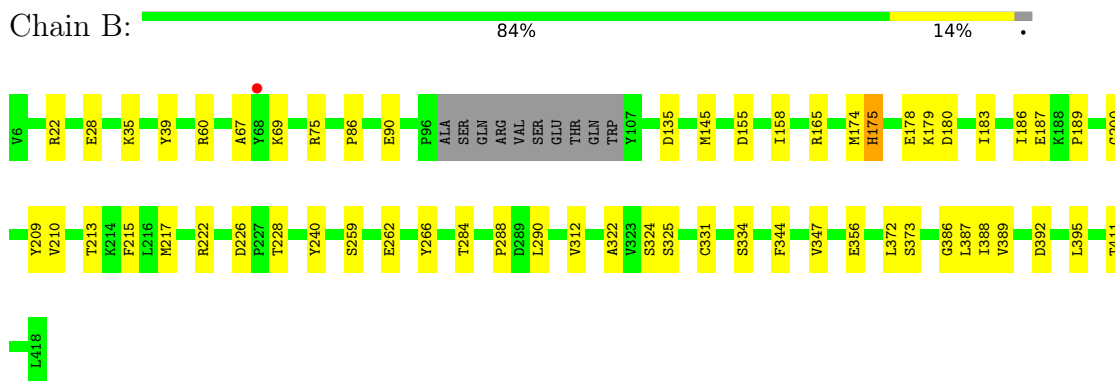
### 3 Residue-property plots [i](#)

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.

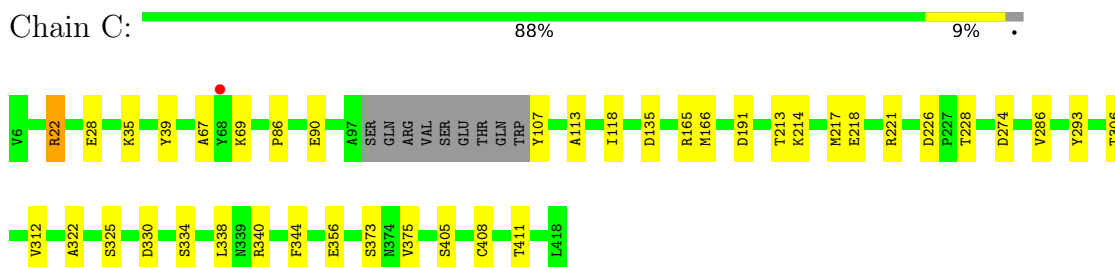
- Molecule 1: Glucose-1-phosphate adenylyltransferase



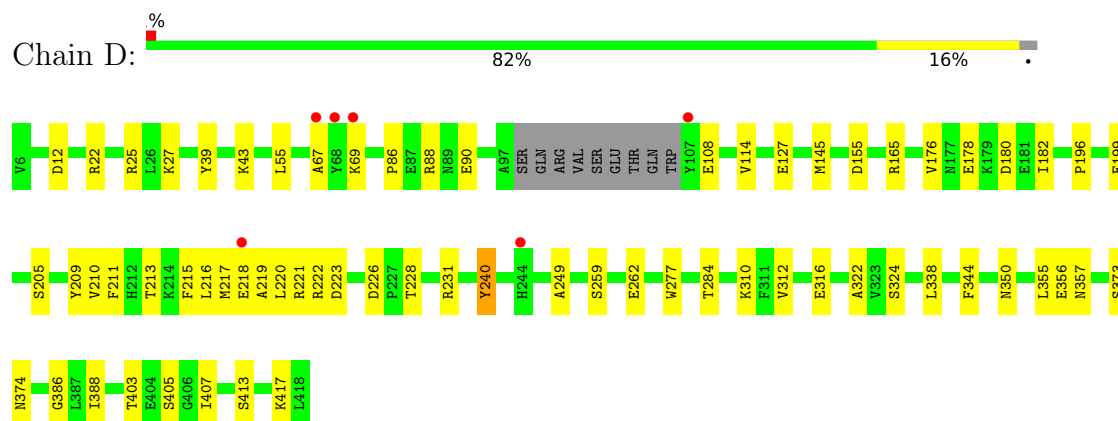
- Molecule 1: Glucose-1-phosphate adenylyltransferase



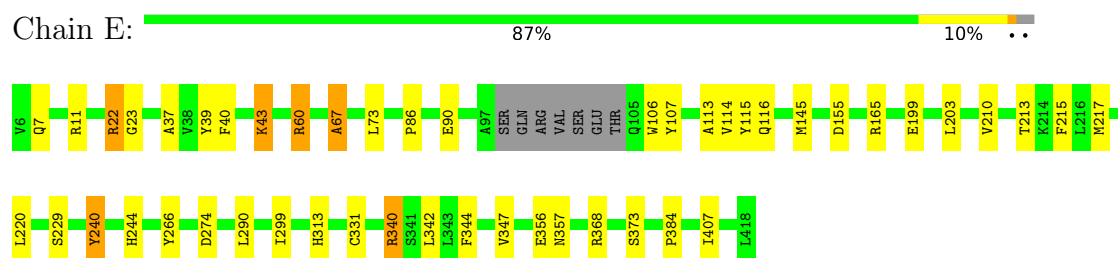
- Molecule 1: Glucose-1-phosphate adenylyltransferase



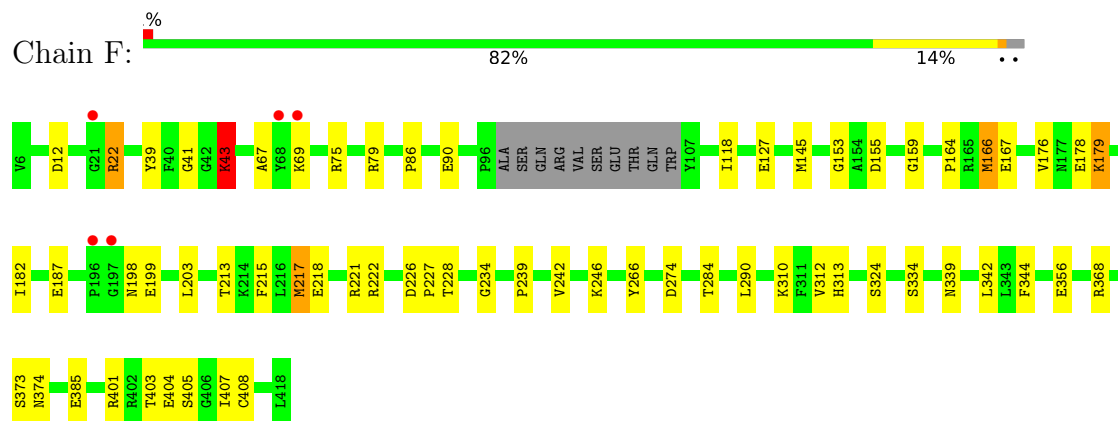
- Molecule 1: Glucose-1-phosphate adenylyltransferase



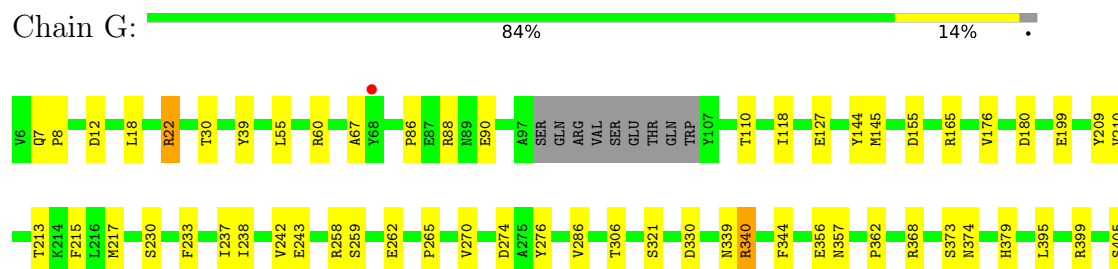
- Molecule 1: Glucose-1-phosphate adenylyltransferase



- Molecule 1: Glucose-1-phosphate adenylyltransferase



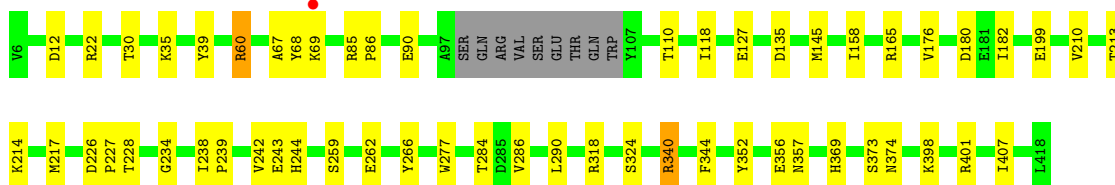
- Molecule 1: Glucose-1-phosphate adenylyltransferase





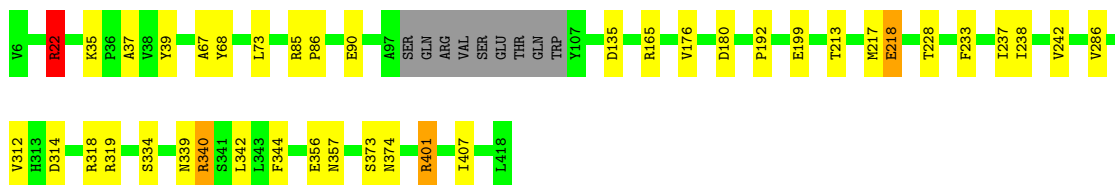
- Molecule 1: Glucose-1-phosphate adenylyltransferase

Chain H: 84% 13% .



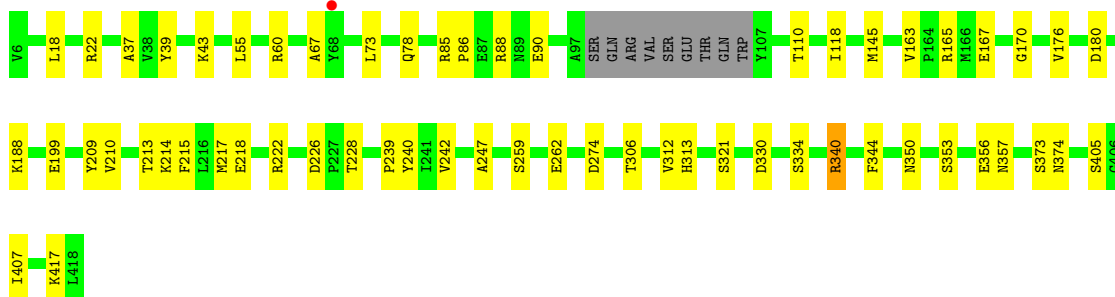
- Molecule 1: Glucose-1-phosphate adenylyltransferase

Chain I: 88% 9% ..



- Molecule 1: Glucose-1-phosphate adenylyltransferase

Chain J: 84% 14% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	433.32Å 141.04Å 92.51Å 90.00° 102.30° 90.00°	Depositor
Resolution (Å)	50.63 – 2.20 50.63 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (50.63-2.20) 89.4 (50.63-2.20)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.88 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, $R_{free}$	0.214 , 0.243 0.219 , 0.246	Depositor DCC
$R_{free}$ test set	13608 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.8	Xtriage
Anisotropy	0.639	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 10.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.478 for -h-2*1,-k,l	Xtriage
Reported twinning fraction	0.500 for -h-2*1,-k,l	Depositor
Outliers	4 of 272606 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	32581	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.58 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.6798e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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, 0.333 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, GOL

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	0/3248	0.52	2/4408 (0.0%)
1	B	0.31	1/3243 (0.0%)	0.52	1/4401 (0.0%)
1	C	0.27	0/3240	0.49	0/4397
1	D	0.35	1/3240 (0.0%)	0.58	3/4397 (0.1%)
1	E	0.28	0/3261	0.49	1/4427 (0.0%)
1	F	0.39	3/3237 (0.1%)	0.60	4/4393 (0.1%)
1	G	0.30	1/3248 (0.0%)	0.50	0/4408
1	H	0.27	0/3248	0.48	0/4408
1	I	0.32	0/3248	0.60	4/4408 (0.1%)
1	J	0.27	0/3226	0.48	0/4378
All	All	0.31	6/32439 (0.0%)	0.53	15/44025 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	43	LYS	CE-NZ	8.36	1.70	1.49
1	F	43	LYS	CD-CE	8.08	1.71	1.51
1	F	43	LYS	CB-CG	6.88	1.71	1.52
1	D	316	GLU	CB-CG	5.79	1.63	1.52
1	G	243	GLU	CB-CG	-5.50	1.41	1.52
1	B	187	GLU	CB-CG	5.13	1.61	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	218	GLU	CA-CB-CG	-13.65	83.37	113.40
1	F	43	LYS	CD-CE-NZ	-11.85	84.44	111.70
1	I	22	ARG	NE-CZ-NH2	-11.74	114.43	120.30
1	F	217	MET	CA-CB-CG	-7.31	100.87	113.30
1	I	22	ARG	CA-CB-CG	7.22	129.28	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	217	MET	N-CA-CB	-6.88	98.22	110.60
1	D	25	ARG	NE-CZ-NH1	6.81	123.70	120.30
1	I	22	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	F	217	MET	CB-CA-C	6.28	122.97	110.40
1	A	404	GLU	CA-CB-CG	-6.13	99.92	113.40
1	A	218	GLU	CB-CG-CD	-5.97	98.08	114.20
1	E	60	ARG	CG-CD-NE	-5.87	99.48	111.80
1	D	27	LYS	CD-CE-NZ	-5.79	98.38	111.70
1	D	43	LYS	CD-CE-NZ	-5.65	98.71	111.70
1	B	178	GLU	CB-CA-C	-5.45	99.51	110.40

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	3174	0	3100	39	0
1	B	3169	0	3098	32	0
1	C	3167	0	3096	23	0
1	D	3167	0	3096	49	0
1	E	3186	0	3108	34	0
1	F	3164	0	3093	54	0
1	G	3174	0	3103	37	0
1	H	3174	0	3103	41	0
1	I	3174	0	3103	26	0
1	J	3155	0	3084	42	0
2	A	45	0	0	1	0
2	B	25	0	0	2	0
2	C	45	0	0	0	0
2	D	40	0	0	2	0
2	E	50	0	0	1	0
2	F	35	0	0	2	0
2	G	50	0	0	3	0
2	H	45	0	0	6	0
2	I	60	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	50	0	0	1	0
3	I	6	0	8	1	0
4	A	32	0	0	1	0
4	B	23	0	0	0	0
4	C	41	0	0	2	0
4	D	20	0	0	1	0
4	E	44	0	0	3	0
4	F	42	0	0	2	0
4	G	57	0	0	0	0
4	H	45	0	0	0	0
4	I	70	0	0	2	0
4	J	52	0	0	0	0
All	All	32581	0	30992	355	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 6.

All (355) 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:F:43:LYS:NZ	1:F:43:LYS:CE	1.70	1.53
1:F:118:ILE:HG23	1:F:217:MET:HE1	1.42	1.01
1:F:43:LYS:NZ	1:F:43:LYS:CD	2.29	0.96
1:I:22:ARG:HD3	1:I:68:TYR:CZ	2.04	0.92
1:E:368:ARG:NH2	2:E:510:SO4:O2	2.05	0.88
1:A:154:ALA:HA	1:A:246:LYS:HG2	1.58	0.86
1:A:274:ASP:OD2	1:A:407:ILE:HD12	1.74	0.86
1:G:60:ARG:NH2	1:G:88:ARG:O	2.09	0.85
1:G:395:LEU:HD21	1:G:399:ARG:HH21	1.40	0.84
1:E:67:ALA:HB2	1:E:106:TRP:HE3	1.42	0.83
1:D:182:ILE:HD11	1:D:249:ALA:HB2	1.61	0.83
1:D:215:PHE:CE1	1:D:218:GLU:OE2	2.33	0.82
1:G:12:ASP:HB3	1:G:127:GLU:HB2	1.63	0.79
1:B:312:VAL:HG12	1:B:334:SER:HA	1.65	0.78
1:C:213:THR:O	1:C:217:MET:HG2	1.85	0.77
1:F:218:GLU:OE2	1:F:222:ARG:NH2	2.18	0.77
1:F:118:ILE:CG2	1:F:217:MET:HE1	2.14	0.77
1:A:165:ARG:HD2	1:A:199:GLU:O	1.84	0.77
1:D:215:PHE:CD1	1:D:218:GLU:OE2	2.39	0.76
1:A:274:ASP:OD2	1:A:407:ILE:CD1	2.34	0.75
1:G:165:ARG:HD2	1:G:199:GLU:O	1.87	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:165:ARG:HD2	1:D:199:GLU:O	1.87	0.74
1:J:60:ARG:NH2	1:J:88:ARG:O	2.18	0.74
1:B:165:ARG:HH21	1:D:388:ILE:HD11	1.53	0.73
1:F:41:GLY:HA3	4:F:605:HOH:O	1.89	0.72
1:I:165:ARG:HD2	1:I:199:GLU:O	1.90	0.72
1:G:321:SER:OG	1:J:321:SER:OG	2.08	0.72
1:J:165:ARG:HD2	1:J:199:GLU:O	1.91	0.71
1:F:176:VAL:HG11	1:F:203:LEU:HD12	1.71	0.71
1:H:401:ARG:NH1	2:H:507:SO4:O2	2.21	0.71
1:B:155:ASP:HB3	1:B:215:PHE:CE2	2.26	0.71
1:H:165:ARG:HD2	1:H:199:GLU:O	1.92	0.70
1:E:165:ARG:HD2	1:E:199:GLU:O	1.92	0.69
1:B:372:LEU:HD12	1:B:389:VAL:HB	1.75	0.69
1:H:85:ARG:NH2	2:H:501:SO4:O1	2.26	0.69
1:J:213:THR:O	1:J:217:MET:HG2	1.93	0.69
1:E:299:ILE:O	4:E:601:HOH:O	2.10	0.69
1:F:118:ILE:HG23	1:F:217:MET:CE	2.21	0.68
1:A:386:GLY:O	1:A:388:ILE:HD12	1.92	0.68
1:E:22:ARG:H	1:E:22:ARG:HD3	1.59	0.67
1:I:22:ARG:HD3	1:I:68:TYR:CE1	2.28	0.67
1:B:209:TYR:OH	2:B:504:SO4:O1	2.08	0.67
1:H:228:THR:HA	1:J:228:THR:CG2	2.25	0.67
1:D:338:LEU:HD21	1:D:355:LEU:HD12	1.77	0.66
1:C:221:ARG:NH1	4:C:602:HOH:O	2.27	0.66
1:E:107:TYR:CD2	1:E:113:ALA:HA	2.30	0.66
1:C:69:LYS:O	4:C:601:HOH:O	2.14	0.65
1:E:40:PHE:O	4:E:602:HOH:O	2.13	0.65
1:A:386:GLY:O	1:A:388:ILE:CD1	2.44	0.65
1:D:218:GLU:HG2	1:D:219:ALA:N	2.10	0.65
1:J:209:TYR:OH	2:J:507:SO4:O1	2.15	0.64
1:F:12:ASP:HB3	1:F:127:GLU:HB2	1.79	0.64
1:D:211:PHE:CD2	1:D:216:LEU:HD12	2.33	0.64
1:D:226:ASP:CG	1:D:228:THR:HG22	2.19	0.63
1:F:118:ILE:HG12	1:F:217:MET:HE1	1.80	0.63
1:G:368:ARG:NH2	2:G:510:SO4:O3	2.31	0.63
1:J:118:ILE:HG23	1:J:217:MET:HE1	1.80	0.63
2:A:508:SO4:O1	1:G:144:TYR:OH	2.11	0.63
1:G:399:ARG:NH1	2:G:506:SO4:O1	2.31	0.63
1:H:12:ASP:HB3	1:H:127:GLU:HB3	1.81	0.62
1:D:350:ASN:ND2	2:D:501:SO4:O1	2.31	0.62
1:E:107:TYR:HE2	1:E:116:GLN:HB2	1.66	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:510:GOL:O2	4:I:601:HOH:O	1.98	0.60
1:D:413:SER:O	1:D:417:LYS:HG3	2.02	0.60
1:A:404:GLU:O	1:A:404:GLU:HG3	2.02	0.60
1:F:75:ARG:HE	1:F:79:ARG:NH2	2.00	0.59
1:A:76:HIS:ND1	4:A:602:HOH:O	2.31	0.59
1:D:155:ASP:HB3	1:D:215:PHE:CZ	2.37	0.59
1:D:209:TYR:OH	2:D:503:SO4:O2	2.17	0.59
1:J:215:PHE:O	1:J:218:GLU:HG2	2.03	0.59
1:A:388:ILE:HD12	1:A:388:ILE:N	2.18	0.59
1:D:155:ASP:HB3	1:D:215:PHE:CE2	2.37	0.59
1:E:60:ARG:HE	1:E:90:GLU:HG2	1.66	0.59
1:G:209:TYR:OH	2:G:503:SO4:O3	2.18	0.58
1:J:145:MET:HE2	1:J:210:VAL:HG23	1.86	0.58
1:J:226:ASP:OD1	1:J:228:THR:HB	2.04	0.58
1:D:213:THR:O	1:D:217:MET:HG2	2.03	0.58
1:B:200:GLY:HA2	1:D:388:ILE:HD13	1.86	0.58
1:C:22:ARG:O	1:C:22:ARG:HD3	2.04	0.58
1:E:7:GLN:HE21	1:E:11:ARG:HD2	1.68	0.58
1:A:387:LEU:C	1:A:388:ILE:HD12	2.24	0.57
1:D:114:VAL:HB	1:D:220:LEU:HD21	1.86	0.57
1:D:222:ARG:HD3	1:D:240:TYR:CE2	2.39	0.57
1:A:55:LEU:HD21	1:G:88:ARG:HD2	1.86	0.57
1:C:356:GLU:O	1:C:373:SER:HA	2.05	0.57
1:C:312:VAL:HG23	1:C:334:SER:HA	1.87	0.56
1:D:176:VAL:HG22	1:D:180:ASP:HA	1.86	0.56
1:F:167:GLU:OE1	1:F:167:GLU:N	2.34	0.56
1:J:176:VAL:HG22	1:J:180:ASP:HA	1.87	0.56
1:H:228:THR:HA	1:J:228:THR:HG21	1.86	0.56
1:D:55:LEU:HD21	1:J:88:ARG:HD2	1.86	0.56
1:F:213:THR:O	1:F:217:MET:HG3	2.06	0.56
1:G:176:VAL:HG22	1:G:180:ASP:HA	1.88	0.56
1:H:227:PRO:O	1:J:228:THR:HG23	2.04	0.56
1:I:176:VAL:CG2	1:I:180:ASP:HA	2.36	0.56
1:F:69:LYS:O	4:F:601:HOH:O	2.18	0.55
1:H:30:THR:O	1:H:68:TYR:OH	2.18	0.55
1:I:176:VAL:HG22	1:I:180:ASP:HA	1.89	0.55
1:F:164:PRO:HB2	1:F:167:GLU:OE1	2.06	0.55
1:D:12:ASP:HB3	1:D:127:GLU:HB3	1.89	0.55
1:D:176:VAL:CG2	1:D:180:ASP:HA	2.36	0.55
1:E:114:VAL:HB	1:E:220:LEU:HD21	1.89	0.55
1:G:176:VAL:CG2	1:G:180:ASP:HA	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:270:VAL:HG12	1:G:276:TYR:HA	1.89	0.55
1:B:175:HIS:O	1:B:183:ILE:HG12	2.07	0.55
1:J:259:SER:HB3	1:J:262:GLU:HG3	1.89	0.55
1:E:356:GLU:O	1:E:373:SER:HA	2.08	0.54
1:J:22:ARG:HH11	1:J:22:ARG:HG3	1.71	0.54
1:B:158:ILE:HG21	1:B:174:MET:HE3	1.88	0.54
1:D:108:GLU:OE1	1:D:231:ARG:NH1	2.41	0.54
1:A:8:PRO:HB3	1:G:8:PRO:HG3	1.90	0.54
1:D:226:ASP:OD1	1:D:228:THR:HG22	2.07	0.54
1:I:356:GLU:O	1:I:373:SER:HA	2.08	0.53
1:J:176:VAL:CG2	1:J:180:ASP:HA	2.37	0.53
1:H:176:VAL:HG22	1:H:180:ASP:HA	1.90	0.53
1:E:67:ALA:CB	1:E:106:TRP:HE3	2.19	0.53
1:F:356:GLU:O	1:F:373:SER:HA	2.09	0.53
1:B:186:ILE:HG21	1:B:189:PRO:HB3	1.91	0.53
1:C:118:ILE:HG23	1:C:217:MET:HE1	1.91	0.53
1:F:403:THR:CG2	1:F:407:ILE:H	2.22	0.53
1:A:356:GLU:O	1:A:373:SER:HA	2.09	0.53
1:D:226:ASP:OD2	1:D:228:THR:HG22	2.09	0.52
1:D:259:SER:HB3	1:D:262:GLU:HG3	1.91	0.52
1:H:176:VAL:CG2	1:H:180:ASP:HA	2.39	0.52
1:D:338:LEU:HD23	1:D:355:LEU:HB2	1.91	0.52
1:E:43:LYS:HD3	4:E:602:HOH:O	2.08	0.52
1:A:88:ARG:HD2	1:G:55:LEU:HD21	1.91	0.52
1:J:274:ASP:CG	1:J:405:SER:HG	2.13	0.52
1:I:22:ARG:NH1	1:I:68:TYR:OH	2.42	0.52
1:A:164:PRO:HB2	1:A:167:GLU:OE2	2.10	0.52
1:D:374:ASN:HB3	1:D:407:ILE:HD13	1.91	0.52
1:F:155:ASP:HB3	1:F:215:PHE:CZ	2.45	0.52
1:H:213:THR:O	1:H:217:MET:HG2	2.09	0.51
1:H:182:ILE:O	1:H:242:VAL:HG23	2.11	0.51
1:D:22:ARG:HH22	1:D:69:LYS:HG3	1.74	0.51
1:D:145:MET:HE2	1:D:210:VAL:HG23	1.92	0.51
1:E:107:TYR:CE2	1:E:116:GLN:HB2	2.45	0.51
1:B:356:GLU:O	1:B:373:SER:HA	2.11	0.51
1:F:166:MET:HG2	1:F:167:GLU:OE1	2.11	0.51
1:J:374:ASN:HB3	1:J:407:ILE:HD13	1.91	0.51
1:B:145:MET:HE2	1:B:210:VAL:HG23	1.92	0.51
1:H:286:VAL:N	2:H:505:SO4:O3	2.36	0.51
1:E:240:TYR:HE1	1:E:244:HIS:HD1	1.59	0.51
1:F:118:ILE:HG21	1:F:221:ARG:NH2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:356:GLU:O	1:G:373:SER:HA	2.11	0.51
1:F:374:ASN:HB3	1:F:407:ILE:HD13	1.92	0.50
1:F:403:THR:HG22	1:F:407:ILE:H	1.75	0.50
1:H:228:THR:HA	1:J:228:THR:HG23	1.93	0.50
1:B:22:ARG:HH22	1:B:69:LYS:HG3	1.76	0.50
1:J:350:ASN:O	1:J:353:SER:OG	2.24	0.50
1:C:107:TYR:CE1	1:C:113:ALA:HA	2.46	0.50
1:E:22:ARG:HH11	1:E:23:GLY:H	1.60	0.50
1:C:165:ARG:NH2	1:H:369:HIS:O	2.44	0.50
1:H:68:TYR:CE1	1:H:69:LYS:HD3	2.46	0.50
1:H:239:PRO:HA	1:H:242:VAL:HG12	1.93	0.50
1:A:78:GLN:HE21	1:J:78:GLN:HE21	1.60	0.49
1:H:374:ASN:HB3	1:H:407:ILE:HD13	1.94	0.49
1:E:331:CYS:SG	1:E:347:VAL:HB	2.52	0.49
1:I:318:ARG:NH2	2:I:508:SO4:O3	2.43	0.49
1:H:234:GLY:N	2:H:506:SO4:O1	2.45	0.49
1:C:286:VAL:HA	1:F:312:VAL:HG21	1.95	0.49
1:E:22:ARG:NH1	1:E:23:GLY:H	2.10	0.49
1:J:239:PRO:HA	1:J:242:VAL:HG22	1.95	0.48
1:D:108:GLU:CD	1:D:231:ARG:HG3	2.34	0.48
1:C:226:ASP:OD1	1:C:228:THR:HB	2.13	0.48
1:A:243:GLU:OE1	1:A:244:HIS:CE1	2.67	0.48
1:B:386:GLY:O	1:B:388:ILE:HD12	2.13	0.48
1:A:22:ARG:NH1	1:A:24:SER:HA	2.28	0.48
1:B:259:SER:HB3	1:B:262:GLU:HG3	1.95	0.48
1:C:214:LYS:O	1:C:218:GLU:HG3	2.14	0.48
1:H:86:PRO:HA	1:H:90:GLU:O	2.14	0.48
1:H:118:ILE:HD12	1:H:217:MET:SD	2.53	0.48
1:H:145:MET:HE2	1:H:210:VAL:HG23	1.95	0.48
1:G:213:THR:O	1:G:217:MET:HG2	2.13	0.48
1:H:259:SER:HB3	1:H:262:GLU:HG3	1.95	0.48
1:E:145:MET:HE2	1:E:210:VAL:HG23	1.96	0.48
1:G:286:VAL:HG22	1:J:312:VAL:CG2	2.44	0.47
1:J:222:ARG:CD	1:J:240:TYR:HE2	2.26	0.47
1:I:39:TYR:HB3	1:I:344:PHE:CE2	2.49	0.47
1:B:322:ALA:HB1	1:B:325:SER:HB2	1.96	0.47
1:D:88:ARG:HD2	1:J:55:LEU:HD21	1.96	0.47
1:A:115:TYR:HD2	1:A:116:GLN:NE2	2.13	0.47
1:I:401:ARG:HE	1:I:401:ARG:HB2	1.53	0.47
1:H:398:LYS:HA	1:H:398:LYS:HD3	1.66	0.47
1:F:215:PHE:O	1:F:218:GLU:HG3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:226:ASP:CG	1:F:228:THR:HG22	2.36	0.46
1:D:403:THR:OG1	1:D:405:SER:OG	2.23	0.46
1:C:39:TYR:HB3	1:C:344:PHE:CE2	2.51	0.46
1:B:372:LEU:HD11	1:B:387:LEU:HD23	1.97	0.46
1:D:178:GLU:HG3	1:D:196:PRO:HB2	1.97	0.46
1:F:153:GLY:H	1:F:246:LYS:NZ	2.14	0.46
1:J:163:VAL:CG1	1:J:167:GLU:HB3	2.45	0.46
1:D:386:GLY:O	1:D:388:ILE:HD12	2.15	0.46
1:E:7:GLN:HE21	1:E:11:ARG:CD	2.29	0.46
1:A:39:TYR:HB3	1:A:344:PHE:CE2	2.50	0.46
1:A:145:MET:HE2	1:A:210:VAL:HG23	1.98	0.46
1:G:18:LEU:HD13	1:G:110:THR:HB	1.97	0.46
1:J:39:TYR:HB3	1:J:344:PHE:CE2	2.50	0.46
1:E:39:TYR:HB3	1:E:344:PHE:CE2	2.51	0.46
1:E:340:ARG:HD3	1:E:357:ASN:OD1	2.16	0.46
1:F:43:LYS:NZ	1:F:43:LYS:HD3	2.27	0.46
1:H:226:ASP:OD1	1:H:228:THR:HB	2.16	0.46
1:G:259:SER:HB3	1:G:262:GLU:HG3	1.98	0.45
1:H:85:ARG:NH2	2:H:501:SO4:S	2.89	0.45
1:J:356:GLU:O	1:J:373:SER:HA	2.16	0.45
1:B:179:LYS:O	1:B:180:ASP:HB2	2.16	0.45
1:F:239:PRO:HA	1:F:242:VAL:HG12	1.99	0.45
1:H:158:ILE:HD11	1:H:238:ILE:HD12	1.99	0.45
1:C:293:TYR:OH	1:F:310:LYS:HB2	2.16	0.45
1:E:67:ALA:CB	1:E:106:TRP:HA	2.46	0.45
1:H:356:GLU:O	1:H:373:SER:HA	2.16	0.45
1:J:43:LYS:HB2	1:J:43:LYS:HE2	1.67	0.45
1:C:28:GLU:HB3	1:C:411:THR:HG21	1.99	0.45
1:G:238:ILE:O	1:G:242:VAL:HG23	2.16	0.45
1:B:22:ARG:NH2	1:B:69:LYS:HG3	2.31	0.45
1:F:118:ILE:HG12	1:F:217:MET:CE	2.45	0.45
1:F:404:GLU:O	1:F:404:GLU:HG3	2.16	0.45
1:G:340:ARG:HD3	1:G:357:ASN:OD1	2.16	0.45
1:D:222:ARG:HD3	1:D:240:TYR:HE2	1.81	0.45
1:F:182:ILE:O	1:F:242:VAL:HG23	2.17	0.45
1:I:37:ALA:HB2	1:I:73:LEU:HB2	1.98	0.45
1:D:221:ARG:NH1	4:D:602:HOH:O	2.27	0.45
1:B:226:ASP:OD1	1:B:228:THR:HB	2.18	0.44
1:B:266:TYR:CE2	1:B:290:LEU:HB2	2.52	0.44
1:A:144:TYR:OH	1:G:7:GLN:OE1	2.20	0.44
1:D:338:LEU:CD2	1:D:355:LEU:HB2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:TYR:HB3	1:B:344:PHE:CE2	2.52	0.44
1:D:86:PRO:HA	1:D:90:GLU:O	2.18	0.44
1:D:284:THR:HB	1:D:324:SER:O	2.18	0.44
1:G:233:PHE:HA	1:G:237:ILE:HB	1.99	0.44
1:I:86:PRO:HA	1:I:90:GLU:O	2.18	0.44
1:C:86:PRO:HA	1:C:90:GLU:O	2.17	0.44
1:D:356:GLU:O	1:D:373:SER:HA	2.17	0.44
1:E:115:TYR:HA	1:E:220:LEU:HD23	1.99	0.44
1:G:306:THR:HB	1:G:330:ASP:HB2	2.00	0.44
1:I:238:ILE:O	1:I:242:VAL:HG23	2.17	0.44
1:A:28:GLU:HB3	1:A:411:THR:HG21	2.00	0.44
1:B:28:GLU:HB3	1:B:411:THR:HG21	2.00	0.44
1:C:35:LYS:HD3	1:C:135:ASP:HB2	2.00	0.44
1:C:274:ASP:CG	1:C:405:SER:HG	2.21	0.44
1:D:216:LEU:C	1:D:216:LEU:HD23	2.37	0.44
1:A:186:ILE:HG21	1:A:189:PRO:HB3	1.99	0.44
1:B:222:ARG:HD2	1:B:240:TYR:HE2	1.82	0.44
1:G:86:PRO:HA	1:G:90:GLU:O	2.18	0.44
1:G:39:TYR:HB3	1:G:344:PHE:CE2	2.51	0.43
1:J:37:ALA:HB2	1:J:73:LEU:HB2	1.99	0.43
1:A:293:TYR:OH	1:D:310:LYS:HB2	2.17	0.43
1:E:7:GLN:HG2	1:E:11:ARG:HD2	1.99	0.43
1:F:39:TYR:HB3	1:F:344:PHE:CE2	2.53	0.43
1:G:374:ASN:HB3	1:G:407:ILE:HD13	2.00	0.43
1:F:178:GLU:O	1:F:179:LYS:HB2	2.18	0.43
1:E:37:ALA:HB2	1:E:73:LEU:HB2	2.00	0.43
1:G:118:ILE:HG23	1:G:217:MET:HE3	2.01	0.43
1:G:145:MET:HE2	1:G:210:VAL:HG23	2.00	0.43
1:A:286:VAL:HA	1:D:312:VAL:HG21	2.00	0.43
1:C:322:ALA:HB1	1:C:325:SER:HB2	2.01	0.43
1:G:258:ARG:HD2	1:G:265:PRO:N	2.34	0.43
1:I:374:ASN:HB3	1:I:407:ILE:HD13	2.00	0.43
1:F:215:PHE:O	1:F:218:GLU:CG	2.66	0.43
1:H:239:PRO:HA	1:H:242:VAL:CG1	2.49	0.43
1:J:85:ARG:HD2	1:J:85:ARG:HA	1.89	0.43
1:I:312:VAL:HG13	1:I:334:SER:HA	2.01	0.43
1:J:340:ARG:HD3	1:J:357:ASN:OD1	2.19	0.43
1:A:213:THR:O	1:A:217:MET:HG2	2.19	0.43
1:D:39:TYR:HB3	1:D:344:PHE:CE2	2.54	0.43
1:F:226:ASP:OD2	1:F:228:THR:HG22	2.18	0.43
1:F:312:VAL:HG12	1:F:334:SER:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:314:ASP:OD1	1:I:319:ARG:HD3	2.19	0.43
1:G:362:PRO:O	1:G:379:HIS:HD2	2.02	0.42
1:H:60:ARG:H	1:H:60:ARG:HG2	1.50	0.42
1:J:86:PRO:HA	1:J:90:GLU:O	2.19	0.42
1:J:222:ARG:HD2	1:J:240:TYR:HE2	1.84	0.42
1:G:395:LEU:HG	1:G:399:ARG:HE	1.85	0.42
1:B:288:PRO:HD2	2:B:505:SO4:O2	2.18	0.42
1:I:85:ARG:HD2	1:I:85:ARG:HA	1.85	0.42
1:E:86:PRO:HA	1:E:90:GLU:O	2.18	0.42
1:E:213:THR:O	1:E:217:MET:HG2	2.20	0.42
1:G:22:ARG:NH1	1:G:30:THR:O	2.53	0.42
1:I:35:LYS:HD3	1:I:135:ASP:HB2	2.02	0.42
1:I:213:THR:O	1:I:217:MET:HG2	2.19	0.42
1:I:340:ARG:HD3	1:I:357:ASN:OD1	2.18	0.42
1:A:274:ASP:OD2	1:A:407:ILE:HD13	2.19	0.42
1:A:338:LEU:HD22	1:A:355:LEU:HD12	2.02	0.42
1:B:86:PRO:HA	1:B:90:GLU:O	2.18	0.42
1:F:86:PRO:HA	1:F:90:GLU:O	2.19	0.42
1:F:368:ARG:NH2	2:F:501:SO4:O2	2.52	0.42
1:H:39:TYR:HB3	1:H:344:PHE:CE2	2.54	0.42
1:I:286:VAL:HG23	2:I:506:SO4:S	2.59	0.42
1:E:155:ASP:HB3	1:E:215:PHE:CZ	2.55	0.42
1:H:318:ARG:HB3	1:H:352:TYR:CE1	2.55	0.42
1:A:257:VAL:HG11	1:A:297:TRP:HB2	2.02	0.42
1:C:375:VAL:HG12	1:C:408:CYS:HB2	2.02	0.42
1:E:67:ALA:HB1	1:E:106:TRP:HA	2.02	0.42
1:J:274:ASP:OD1	1:J:405:SER:OG	2.36	0.42
1:B:372:LEU:HD12	1:B:389:VAL:CB	2.45	0.41
1:E:266:TYR:CE2	1:E:290:LEU:HB2	2.55	0.41
1:C:306:THR:HB	1:C:330:ASP:HB2	2.02	0.41
1:A:155:ASP:HB3	1:A:215:PHE:CZ	2.55	0.41
1:D:322:ALA:HA	1:D:338:LEU:HB2	2.02	0.41
1:A:217:MET:O	1:A:221:ARG:HG3	2.21	0.41
1:B:35:LYS:HD3	1:B:135:ASP:HB2	2.03	0.41
1:F:234:GLY:N	2:F:503:SO4:O3	2.51	0.41
1:I:165:ARG:O	1:I:192:PRO:HG2	2.20	0.41
1:B:213:THR:O	1:B:217:MET:HG2	2.20	0.41
1:B:222:ARG:HD2	1:B:240:TYR:CE2	2.55	0.41
1:B:284:THR:HB	1:B:324:SER:O	2.21	0.41
1:D:223:ASP:OD1	1:D:231:ARG:NE	2.46	0.41
1:E:203:LEU:HD23	1:E:203:LEU:HA	1.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:22:ARG:HD2	1:F:22:ARG:O	2.21	0.41
1:G:274:ASP:CG	1:G:405:SER:HG	2.23	0.41
1:H:340:ARG:HD3	1:H:357:ASN:OD1	2.21	0.41
1:J:18:LEU:HD13	1:J:110:THR:HB	2.01	0.41
1:A:166:MET:SD	1:F:385:GLU:HB3	2.60	0.41
1:C:191:ASP:HB2	1:H:352:TYR:CD1	2.56	0.41
1:D:182:ILE:N	1:D:182:ILE:HD12	2.35	0.41
1:F:187:GLU:OE1	1:F:187:GLU:HA	2.20	0.41
1:F:203:LEU:HD23	1:F:203:LEU:HA	1.87	0.41
1:F:284:THR:HB	1:F:324:SER:O	2.21	0.41
1:I:339:ASN:O	1:I:356:GLU:HA	2.21	0.41
1:J:242:VAL:HG12	1:J:247:ALA:HB3	2.02	0.41
1:J:312:VAL:HG13	1:J:334:SER:HA	2.02	0.41
1:J:417:LYS:HB2	1:J:417:LYS:HE3	1.93	0.41
1:A:259:SER:HB3	1:A:262:GLU:HG3	2.03	0.41
1:B:331:CYS:SG	1:B:347:VAL:HB	2.61	0.41
1:C:322:ALA:HA	1:C:338:LEU:HB2	2.03	0.41
1:D:277:TRP:HZ2	1:D:357:ASN:HB3	1.85	0.41
1:F:43:LYS:CD	1:F:43:LYS:HZ2	2.25	0.41
1:F:339:ASN:O	1:F:356:GLU:HA	2.21	0.41
1:J:306:THR:HB	1:J:330:ASP:HB2	2.03	0.41
1:F:145:MET:HE3	1:F:159:GLY:N	2.36	0.41
1:F:199:GLU:OE1	1:F:199:GLU:N	2.53	0.41
1:H:35:LYS:HD3	1:H:135:ASP:HB2	2.02	0.41
1:I:233:PHE:HA	1:I:237:ILE:HB	2.03	0.41
1:A:322:ALA:HB2	1:A:338:LEU:HD12	2.03	0.41
1:E:274:ASP:OD1	1:E:407:ILE:HD13	2.21	0.41
1:F:227:PRO:O	1:I:228:THR:HB	2.21	0.41
1:F:274:ASP:CG	1:F:405:SER:HG	2.24	0.41
1:F:401:ARG:O	1:F:408:CYS:HA	2.21	0.41
1:H:243:GLU:HB3	1:H:244:HIS:ND1	2.36	0.41
1:H:266:TYR:CE2	1:H:290:LEU:HB2	2.56	0.41
1:H:277:TRP:HZ2	1:H:357:ASN:HB3	1.86	0.41
1:A:35:LYS:HD3	1:A:135:ASP:HB2	2.03	0.41
1:A:322:ALA:HB1	1:A:325:SER:HB2	2.02	0.41
1:F:266:TYR:CE2	1:F:290:LEU:HB2	2.55	0.41
1:F:274:ASP:OD1	1:F:403:THR:HG21	2.21	0.41
1:A:86:PRO:HA	1:A:90:GLU:O	2.21	0.40
1:B:392:ASP:HB3	1:B:395:LEU:HB3	2.04	0.40
1:G:155:ASP:HB3	1:G:215:PHE:CZ	2.56	0.40
1:J:170:GLY:O	1:J:188:LYS:NZ	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:339:ASN:O	1:G:356:GLU:HA	2.20	0.40
1:H:110:THR:HG21	2:H:506:SO4:O3	2.21	0.40
1:A:203:LEU:HD23	1:A:203:LEU:HA	1.95	0.40
1:H:12:ASP:HB3	1:H:127:GLU:CB	2.51	0.40
1:H:284:THR:HB	1:H:324:SER:O	2.20	0.40
1:I:319:ARG:NH1	4:I:614:HOH:O	2.54	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	400/413 (97%)	383 (96%)	14 (4%)	3 (1%)	19	19
1	B	399/413 (97%)	383 (96%)	15 (4%)	1 (0%)	41	46
1	C	400/413 (97%)	387 (97%)	12 (3%)	1 (0%)	41	46
1	D	400/413 (97%)	383 (96%)	16 (4%)	1 (0%)	41	46
1	E	402/413 (97%)	385 (96%)	15 (4%)	2 (0%)	29	31
1	F	399/413 (97%)	384 (96%)	12 (3%)	3 (1%)	19	19
1	G	400/413 (97%)	388 (97%)	11 (3%)	1 (0%)	41	46
1	H	400/413 (97%)	384 (96%)	15 (4%)	1 (0%)	41	46
1	I	400/413 (97%)	386 (96%)	13 (3%)	1 (0%)	41	46
1	J	400/413 (97%)	384 (96%)	14 (4%)	2 (0%)	29	31
All	All	4000/4130 (97%)	3847 (96%)	137 (3%)	16 (0%)	34	37

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	TRP

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Mol	Chain	Res	Type
1	A	67	ALA
1	E	67	ALA
1	F	179	LYS
1	A	107	TYR
1	B	67	ALA
1	C	67	ALA
1	D	67	ALA
1	F	67	ALA
1	G	67	ALA
1	H	67	ALA
1	I	67	ALA
1	J	67	ALA
1	E	313	HIS
1	F	313	HIS
1	J	313	HIS

### 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	335/344 (97%)	331 (99%)	4 (1%)	71	83
1	B	335/344 (97%)	332 (99%)	3 (1%)	78	88
1	C	334/344 (97%)	331 (99%)	3 (1%)	78	88
1	D	334/344 (97%)	332 (99%)	2 (1%)	86	93
1	E	335/344 (97%)	328 (98%)	7 (2%)	53	67
1	F	334/344 (97%)	329 (98%)	5 (2%)	65	78
1	G	335/344 (97%)	332 (99%)	3 (1%)	78	88
1	H	335/344 (97%)	331 (99%)	4 (1%)	71	83
1	I	335/344 (97%)	330 (98%)	5 (2%)	65	78
1	J	332/344 (96%)	330 (99%)	2 (1%)	86	93
All	All	3344/3440 (97%)	3306 (99%)	38 (1%)	73	85

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ARG
1	A	32	ARG
1	A	229	SER
1	A	342	LEU
1	B	60	ARG
1	B	75	ARG
1	B	175	HIS
1	C	22	ARG
1	C	166	MET
1	C	340	ARG
1	D	205	SER
1	D	240	TYR
1	E	22	ARG
1	E	43	LYS
1	E	229	SER
1	E	240	TYR
1	E	340	ARG
1	E	342	LEU
1	E	384	PRO
1	F	22	ARG
1	F	43	LYS
1	F	166	MET
1	F	198	ASN
1	F	342	LEU
1	G	22	ARG
1	G	230	SER
1	G	340	ARG
1	H	22	ARG
1	H	60	ARG
1	H	214	LYS
1	H	340	ARG
1	I	22	ARG
1	I	218	GLU
1	I	340	ARG
1	I	342	LEU
1	I	401	ARG
1	J	214	LYS
1	J	340	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN

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Mol	Chain	Res	Type
1	B	313	HIS
1	D	244	HIS
1	E	7	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

90 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	A	504	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	J	501	-	4,4,4	0.14	0	6,6,6	0.17	0
2	SO4	J	509	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	G	503	-	4,4,4	0.13	0	6,6,6	0.10	0
2	SO4	C	509	-	4,4,4	0.15	0	6,6,6	0.10	0
2	SO4	H	508	-	4,4,4	0.15	0	6,6,6	0.12	0
2	SO4	A	508	-	4,4,4	0.13	0	6,6,6	0.11	0
2	SO4	B	501	-	4,4,4	0.15	0	6,6,6	0.09	0
2	SO4	C	502	-	4,4,4	0.15	0	6,6,6	0.11	0
2	SO4	E	510	-	4,4,4	0.13	0	6,6,6	0.19	0
2	SO4	F	504	-	4,4,4	0.14	0	6,6,6	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	G	505	-	4,4,4	0.12	0	6,6,6	0.16	0
2	SO4	B	505	-	4,4,4	0.13	0	6,6,6	0.17	0
2	SO4	G	504	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	B	503	-	4,4,4	0.14	0	6,6,6	0.13	0
2	SO4	G	501	-	4,4,4	0.14	0	6,6,6	0.06	0
2	SO4	H	507	-	4,4,4	0.14	0	6,6,6	0.18	0
2	SO4	I	512	-	4,4,4	0.16	0	6,6,6	0.06	0
2	SO4	D	504	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	H	509	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	G	509	-	4,4,4	0.15	0	6,6,6	0.07	0
2	SO4	E	502	-	4,4,4	0.14	0	6,6,6	0.17	0
2	SO4	D	505	-	4,4,4	0.15	0	6,6,6	0.12	0
2	SO4	G	507	-	4,4,4	0.16	0	6,6,6	0.07	0
2	SO4	F	505	-	4,4,4	0.14	0	6,6,6	0.08	0
2	SO4	C	503	-	4,4,4	0.14	0	6,6,6	0.12	0
2	SO4	C	501	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	J	504	-	4,4,4	0.16	0	6,6,6	0.09	0
2	SO4	I	511	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	H	503	-	4,4,4	0.15	0	6,6,6	0.10	0
2	SO4	A	503	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	E	505	-	4,4,4	0.15	0	6,6,6	0.11	0
2	SO4	E	501	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	I	508	-	4,4,4	0.15	0	6,6,6	0.10	0
2	SO4	A	501	-	4,4,4	0.13	0	6,6,6	0.12	0
2	SO4	E	508	-	4,4,4	0.15	0	6,6,6	0.09	0
2	SO4	D	507	-	4,4,4	0.14	0	6,6,6	0.05	0
2	SO4	F	501	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	H	505	-	4,4,4	0.15	0	6,6,6	0.18	0
2	SO4	D	506	-	4,4,4	0.15	0	6,6,6	0.07	0
2	SO4	H	501	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	J	506	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	C	504	-	4,4,4	0.16	0	6,6,6	0.11	0
2	SO4	E	509	-	4,4,4	0.16	0	6,6,6	0.08	0
2	SO4	A	509	-	4,4,4	0.13	0	6,6,6	0.06	0
2	SO4	H	502	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	E	506	-	4,4,4	0.15	0	6,6,6	0.11	0
2	SO4	D	502	-	4,4,4	0.13	0	6,6,6	0.14	0
2	SO4	J	507	-	4,4,4	0.14	0	6,6,6	0.08	0
2	SO4	I	506	-	4,4,4	0.15	0	6,6,6	0.08	0
2	SO4	A	502	-	4,4,4	0.15	0	6,6,6	0.14	0
2	SO4	G	510	-	4,4,4	0.15	0	6,6,6	0.12	0
2	SO4	H	506	-	4,4,4	0.15	0	6,6,6	0.09	0
2	SO4	G	506	-	4,4,4	0.12	0	6,6,6	0.11	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	J	503	-	4,4,4	0.13	0	6,6,6	0.15	0
2	SO4	A	506	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	I	509	-	4,4,4	0.16	0	6,6,6	0.19	0
2	SO4	J	505	-	4,4,4	0.15	0	6,6,6	0.10	0
2	SO4	F	503	-	4,4,4	0.13	0	6,6,6	0.09	0
2	SO4	A	505	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	F	506	-	4,4,4	0.14	0	6,6,6	0.17	0
2	SO4	C	507	-	4,4,4	0.15	0	6,6,6	0.07	0
2	SO4	B	504	-	4,4,4	0.14	0	6,6,6	0.07	0
2	SO4	J	508	-	4,4,4	0.16	0	6,6,6	0.07	0
2	SO4	J	502	-	4,4,4	0.14	0	6,6,6	0.08	0
2	SO4	I	504	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	G	502	-	4,4,4	0.14	0	6,6,6	0.06	0
3	GOL	I	510	-	5,5,5	2.61	2 (40%)	5,5,5	1.51	1 (20%)
2	SO4	D	508	-	4,4,4	0.13	0	6,6,6	0.10	0
2	SO4	I	502	-	4,4,4	0.13	0	6,6,6	0.11	0
2	SO4	A	507	-	4,4,4	0.14	0	6,6,6	0.13	0
2	SO4	I	507	-	4,4,4	0.15	0	6,6,6	0.12	0
2	SO4	I	503	-	4,4,4	0.12	0	6,6,6	0.09	0
2	SO4	I	505	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	I	513	-	4,4,4	0.15	0	6,6,6	0.05	0
2	SO4	F	507	-	4,4,4	0.14	0	6,6,6	0.09	0
2	SO4	G	508	-	4,4,4	0.15	0	6,6,6	0.14	0
2	SO4	C	505	-	4,4,4	0.13	0	6,6,6	0.11	0
2	SO4	C	508	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	B	502	-	4,4,4	0.13	0	6,6,6	0.17	0
2	SO4	I	501	-	4,4,4	0.15	0	6,6,6	0.07	0
2	SO4	E	503	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	F	502	-	4,4,4	0.13	0	6,6,6	0.11	0
2	SO4	D	501	-	4,4,4	0.14	0	6,6,6	0.10	0
2	SO4	E	504	-	4,4,4	0.15	0	6,6,6	0.11	0
2	SO4	C	506	-	4,4,4	0.15	0	6,6,6	0.09	0
2	SO4	E	507	-	4,4,4	0.15	0	6,6,6	0.06	0
2	SO4	D	503	-	4,4,4	0.15	0	6,6,6	0.14	0
2	SO4	H	504	-	4,4,4	0.14	0	6,6,6	0.11	0
2	SO4	J	510	-	4,4,4	0.13	0	6,6,6	0.09	0

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
3	GOL	I	510	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	510	GOL	C1-C2	4.61	1.70	1.51
3	I	510	GOL	C3-C2	3.01	1.64	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	510	GOL	C3-C2-C1	2.34	120.81	111.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

19 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	503	SO4	1	0
2	A	508	SO4	1	0
2	E	510	SO4	1	0
2	B	505	SO4	1	0
2	H	507	SO4	1	0
2	I	508	SO4	1	0
2	F	501	SO4	1	0
2	H	505	SO4	1	0
2	H	501	SO4	2	0
2	J	507	SO4	1	0
2	I	506	SO4	1	0
2	G	510	SO4	1	0
2	H	506	SO4	2	0
2	G	506	SO4	1	0
2	F	503	SO4	1	0
2	B	504	SO4	1	0
3	I	510	GOL	1	0
2	D	501	SO4	1	0
2	D	503	SO4	1	0



## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 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	404/413 (97%)	-0.23	2 (0%) 91 90	18, 32, 44, 68	0
1	B	403/413 (97%)	-0.14	1 (0%) 95 94	19, 36, 48, 57	0
1	C	404/413 (97%)	-0.30	1 (0%) 95 94	16, 28, 41, 53	0
1	D	404/413 (97%)	-0.18	6 (1%) 73 72	19, 34, 48, 60	0
1	E	406/413 (98%)	-0.33	0 100 100	14, 26, 40, 49	0
1	F	403/413 (97%)	-0.20	5 (1%) 79 77	17, 32, 47, 62	0
1	G	404/413 (97%)	-0.32	1 (0%) 95 94	15, 25, 41, 52	0
1	H	404/413 (97%)	-0.31	1 (0%) 95 94	15, 28, 43, 55	0
1	I	404/413 (97%)	-0.32	0 100 100	16, 25, 42, 52	0
1	J	404/413 (97%)	-0.36	1 (0%) 95 94	15, 25, 39, 48	0
All	All	4040/4130 (97%)	-0.27	18 (0%) 92 91	14, 29, 45, 68	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	68	TYR	4.1
1	F	68	TYR	3.4
1	C	68	TYR	3.2
1	F	197	GLY	3.1
1	A	68	TYR	3.0
1	D	67	ALA	2.7
1	J	68	TYR	2.7
1	D	244	HIS	2.5
1	B	68	TYR	2.5
1	F	69	LYS	2.3
1	D	107	TYR	2.2
1	D	69	LYS	2.2
1	F	196	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	218	GLU	2.2
1	G	68	TYR	2.1
1	A	316	GLU	2.1
1	F	21	GLY	2.0
1	H	69	LYS	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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	I	510	6/6	0.78	0.27	33,44,49,51	0
2	SO4	F	507	5/5	0.94	0.12	36,38,48,54	0
2	SO4	I	506	5/5	0.94	0.12	35,35,35,44	0
2	SO4	J	510	5/5	0.94	0.22	44,49,58,63	0
2	SO4	C	508	5/5	0.94	0.14	38,46,58,66	0
2	SO4	H	507	5/5	0.95	0.10	34,37,49,51	0
2	SO4	F	504	5/5	0.95	0.11	34,36,47,49	0
2	SO4	I	511	5/5	0.95	0.16	30,32,39,39	0
2	SO4	H	501	5/5	0.95	0.12	25,42,49,66	0
2	SO4	H	505	5/5	0.95	0.14	27,30,47,52	0
2	SO4	E	506	5/5	0.96	0.12	37,40,42,51	0
2	SO4	F	503	5/5	0.96	0.21	40,42,56,61	0
2	SO4	C	506	5/5	0.96	0.22	46,46,59,61	0
2	SO4	C	507	5/5	0.96	0.13	40,48,57,58	0
2	SO4	J	505	5/5	0.96	0.11	22,26,29,30	0
2	SO4	J	509	5/5	0.96	0.12	36,37,51,58	0
2	SO4	G	505	5/5	0.96	0.11	20,27,39,50	0
2	SO4	B	502	5/5	0.96	0.10	34,35,48,51	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	F	506	5/5	0.97	0.10	35,36,41,48	0
2	SO4	C	502	5/5	0.97	0.10	25,38,44,48	0
2	SO4	C	509	5/5	0.97	0.11	23,23,39,46	0
2	SO4	D	505	5/5	0.97	0.09	33,36,48,49	0
2	SO4	D	507	5/5	0.97	0.13	44,47,52,62	0
2	SO4	C	505	5/5	0.97	0.13	23,35,40,52	0
2	SO4	I	505	5/5	0.97	0.12	27,28,42,44	0
2	SO4	E	508	5/5	0.97	0.10	32,41,48,49	0
2	SO4	I	509	5/5	0.97	0.06	35,37,50,55	0
2	SO4	E	509	5/5	0.97	0.14	30,38,39,40	0
2	SO4	I	512	5/5	0.97	0.09	36,38,46,50	0
2	SO4	I	513	5/5	0.97	0.07	40,41,54,55	0
2	SO4	E	510	5/5	0.97	0.10	33,34,46,50	0
2	SO4	A	509	5/5	0.97	0.07	39,43,52,54	0
2	SO4	A	507	5/5	0.97	0.07	32,37,44,54	0
2	SO4	F	505	5/5	0.97	0.15	35,35,51,59	0
2	SO4	A	506	5/5	0.98	0.09	41,45,51,54	0
2	SO4	A	501	5/5	0.98	0.09	23,33,36,40	0
2	SO4	A	508	5/5	0.98	0.08	26,38,41,46	0
2	SO4	A	503	5/5	0.98	0.10	32,33,49,52	0
2	SO4	A	504	5/5	0.98	0.11	33,41,47,49	0
2	SO4	G	503	5/5	0.98	0.11	24,35,38,41	0
2	SO4	G	504	5/5	0.98	0.10	30,43,49,49	0
2	SO4	D	503	5/5	0.98	0.12	41,43,44,53	0
2	SO4	G	507	5/5	0.98	0.13	39,45,61,63	0
2	SO4	G	508	5/5	0.98	0.08	25,27,41,48	0
2	SO4	G	510	5/5	0.98	0.13	28,30,38,41	0
2	SO4	D	504	5/5	0.98	0.11	30,33,42,45	0
2	SO4	H	503	5/5	0.98	0.07	21,28,36,39	0
2	SO4	B	503	5/5	0.98	0.13	29,39,51,52	0
2	SO4	H	506	5/5	0.98	0.17	29,37,40,44	0
2	SO4	D	506	5/5	0.98	0.08	31,50,55,57	0
2	SO4	H	508	5/5	0.98	0.08	34,43,45,45	0
2	SO4	H	509	5/5	0.98	0.13	33,35,43,48	0
2	SO4	I	502	5/5	0.98	0.10	31,34,41,48	0
2	SO4	I	503	5/5	0.98	0.08	24,31,34,38	0
2	SO4	I	504	5/5	0.98	0.11	24,24,38,46	0
2	SO4	B	504	5/5	0.98	0.19	40,50,51,54	0
2	SO4	D	508	5/5	0.98	0.09	32,36,41,47	0
2	SO4	I	507	5/5	0.98	0.08	29,31,36,42	0
2	SO4	I	508	5/5	0.98	0.08	31,33,39,40	0
2	SO4	E	504	5/5	0.98	0.11	33,45,47,48	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	E	505	5/5	0.98	0.12	31,35,42,45	0
2	SO4	B	505	5/5	0.98	0.07	30,30,36,40	0
2	SO4	E	507	5/5	0.98	0.08	34,46,50,52	0
2	SO4	A	505	5/5	0.98	0.07	32,38,49,52	0
2	SO4	J	508	5/5	0.98	0.07	31,34,35,39	0
2	SO4	C	503	5/5	0.98	0.09	27,28,42,51	0
2	SO4	C	504	5/5	0.98	0.09	25,27,43,55	0
2	SO4	F	502	5/5	0.98	0.08	30,34,40,40	0
2	SO4	F	501	5/5	0.99	0.10	27,27,29,32	0
2	SO4	G	506	5/5	0.99	0.09	31,37,41,46	0
2	SO4	E	503	5/5	0.99	0.08	19,25,32,41	0
2	SO4	A	502	5/5	0.99	0.09	21,33,38,42	0
2	SO4	G	509	5/5	0.99	0.08	33,37,47,52	0
2	SO4	B	501	5/5	0.99	0.09	19,31,35,36	0
2	SO4	D	501	5/5	0.99	0.09	21,32,38,44	0
2	SO4	H	502	5/5	0.99	0.09	18,24,27,31	0
2	SO4	D	502	5/5	0.99	0.05	28,31,40,43	0
2	SO4	H	504	5/5	0.99	0.11	25,31,38,40	0
2	SO4	C	501	5/5	0.99	0.10	24,29,34,35	0
2	SO4	J	501	5/5	0.99	0.10	24,33,36,43	0
2	SO4	J	502	5/5	0.99	0.10	19,22,27,30	0
2	SO4	J	503	5/5	0.99	0.11	23,30,36,37	0
2	SO4	J	504	5/5	0.99	0.07	22,23,33,34	0
2	SO4	G	501	5/5	0.99	0.12	18,20,25,25	0
2	SO4	J	506	5/5	0.99	0.08	24,31,37,49	0
2	SO4	J	507	5/5	0.99	0.10	32,37,43,44	0
2	SO4	G	502	5/5	0.99	0.09	18,27,31,32	0
2	SO4	E	501	5/5	0.99	0.11	19,20,24,29	0
2	SO4	E	502	5/5	0.99	0.07	21,26,32,35	0
2	SO4	I	501	5/5	0.99	0.12	16,18,26,31	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.