



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 19, 2020 – 09:13 pm BST

PDB ID : 1XN1  
Title : Crystal Structure Of Lumazine Synthase From Brucella Abortus (Orthorhombic Form At 3.05 Angstroms)  
Authors : Klinke, S.; Zylberman, V.; Vega, D.R.; Guimaraes, B.G.; Braden, B.C.; Goldbaum, F.A.  
Deposited on : 2004-10-04  
Resolution : 3.05 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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	:	2.11
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)	:	2.11

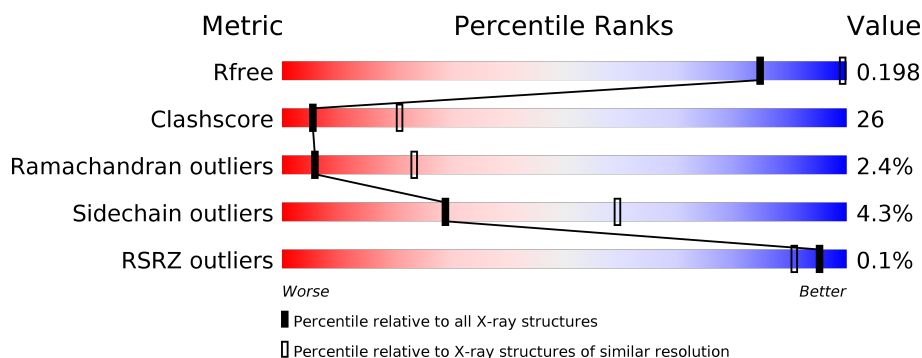
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.05 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	1754 (3.10-3.02)
Clashscore	141614	1864 (3.10-3.02)
Ramachandran outliers	138981	1794 (3.10-3.02)
Sidechain outliers	138945	1793 (3.10-3.02)
RSRZ outliers	127900	1713 (3.10-3.02)

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 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	158	<div> <div>%</div> <div> <div></div> <div>50%</div> <div>39%</div> <div>5%</div> <div>6%</div> </div> </div>
1	B	158	<div> <div>52%</div> <div>37%</div> <div>•</div> <div>7%</div> </div>
1	C	158	<div> <div>%</div> <div> <div></div> <div>48%</div> <div>40%</div> <div>5%</div> <div>7%</div> </div> </div>
1	D	158	<div> <div>48%</div> <div>41%</div> <div>•</div> <div>7%</div> </div>
1	E	158	<div> <div>53%</div> <div>39%</div> <div>•</div> <div>6%</div> </div>
1	F	158	<div> <div>47%</div> <div>44%</div> <div>••</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	158	
1	H	158	
1	I	158	
1	J	158	

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
4	SO4	D	211	-	-	-	X

## 2 Entry composition

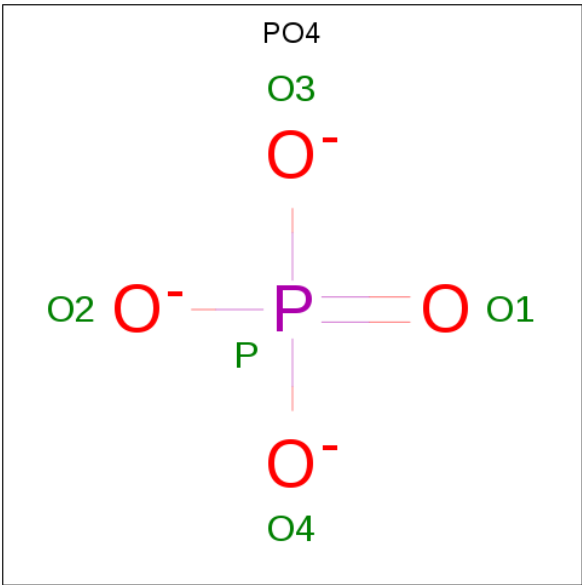
There are 5 unique types of molecules in this entry. The entry contains 11578 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 6,7-dimethyl-8-ribityllumazine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	148	Total	C	N	O	S	0	0	0
			1142	736	201	203	2			
1	B	147	Total	C	N	O	S	0	0	0
			1133	730	199	202	2			
1	C	147	Total	C	N	O	S	0	0	0
			1127	727	199	199	2			
1	D	147	Total	C	N	O	S	0	0	0
			1137	733	200	202	2			
1	E	148	Total	C	N	O	S	0	0	0
			1140	735	201	202	2			
1	F	148	Total	C	N	O	S	0	0	0
			1142	736	201	203	2			
1	G	148	Total	C	N	O	S	0	0	0
			1144	736	203	203	2			
1	H	146	Total	C	N	O	S	0	0	0
			1126	725	200	199	2			
1	I	148	Total	C	N	O	S	0	0	0
			1138	733	200	203	2			
1	J	147	Total	C	N	O	S	0	0	0
			1131	728	198	203	2			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		

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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Na	0	0
			1	1		
3	J	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		
3	A	2	Total	Na	0	0
			2	2		
3	F	1	Total	Na	0	0
			1	1		

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 5 is water.

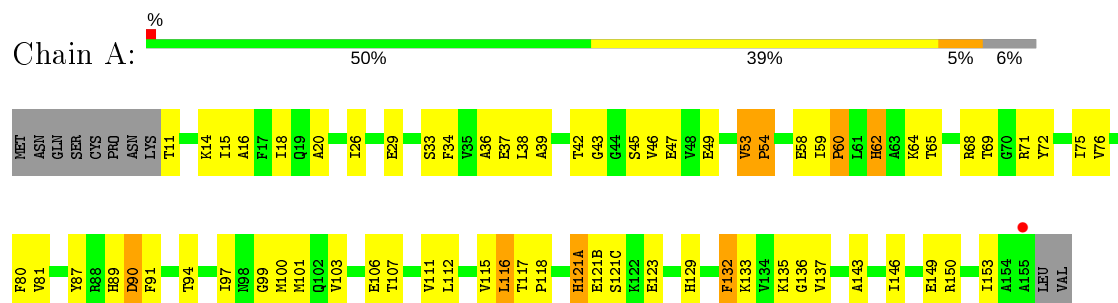
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	O	0	0
			4	4		
5	C	3	Total	O	0	0
			3	3		
5	D	3	Total	O	0	0
			3	3		
5	G	1	Total	O	0	0
			1	1		
5	H	1	Total	O	0	0
			1	1		
5	I	2	Total	O	0	0
			2	2		
5	J	2	Total	O	0	0
			2	2		



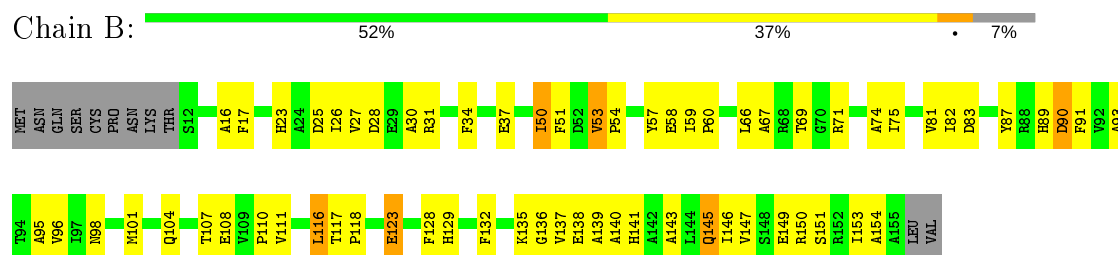
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

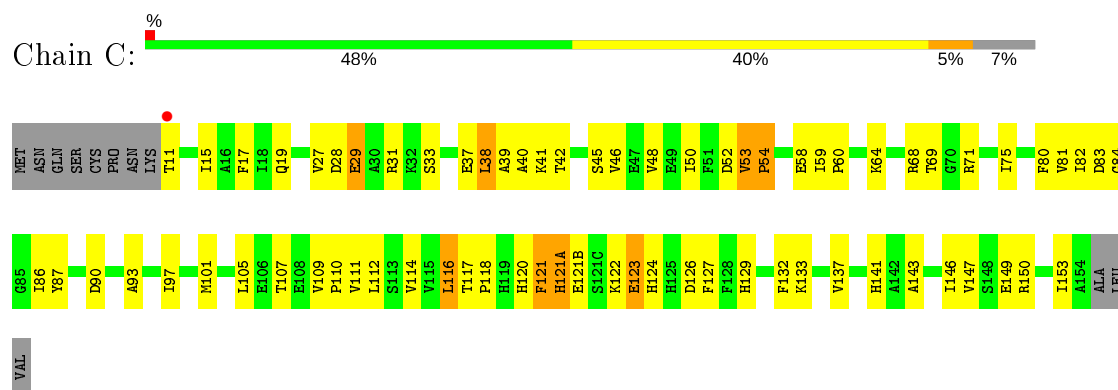
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



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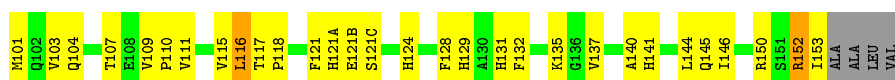
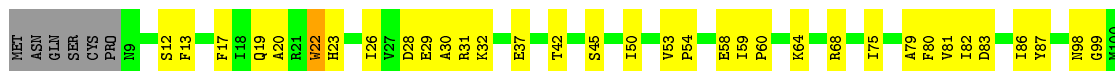
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase





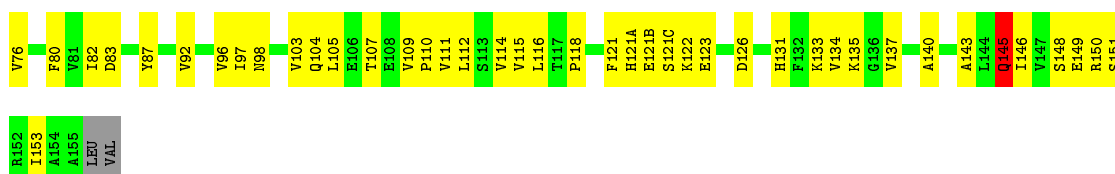
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain E: 53% 39% 6%



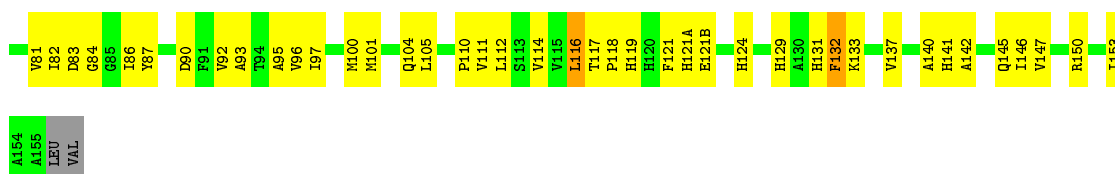
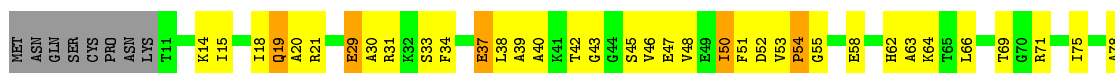
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain F: 47% 44% 6%



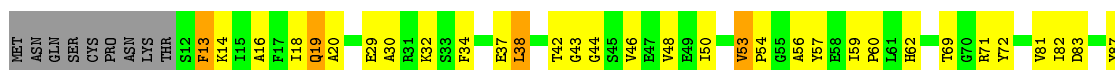
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

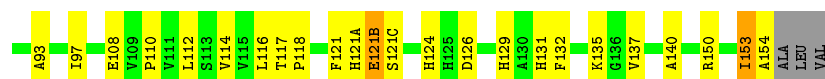
Chain G: 45% 44% 6%



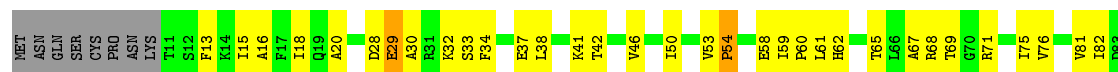
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase

Chain H: 57% 32% 8%

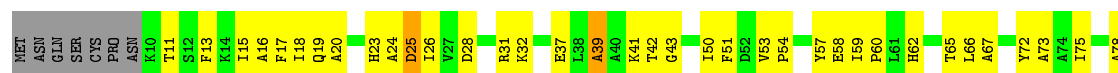




- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.55Å 120.24Å 159.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.00 – 3.05 56.25 – 3.05	Depositor EDS
% Data completeness (in resolution range)	97.6 (56.00-3.05) 99.0 (56.25-3.05)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.75 (at 3.07Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.199 , 0.256 0.202 , 0.198	Depositor DCC
$R_{free}$ test set	2162 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.1	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 38.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	11578	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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, 0.333 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: NA, PO4, 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	0.41	0/1171	0.61	0/1589
1	B	0.41	0/1162	0.62	0/1578
1	C	0.41	0/1156	0.61	0/1571
1	D	0.42	0/1166	0.64	0/1582
1	E	0.41	0/1169	0.64	0/1587
1	F	0.43	0/1171	0.62	0/1589
1	G	0.39	0/1173	0.61	0/1592
1	H	0.42	0/1155	0.63	0/1569
1	I	0.42	0/1167	0.62	0/1585
1	J	0.45	0/1160	0.65	0/1577
All	All	0.42	0/11650	0.63	0/15819

There are no bond length outliers.

There are no bond angle outliers.

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	1142	0	1120	60	0
1	B	1133	0	1107	61	0
1	C	1127	0	1101	68	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1137	0	1115	62	0
1	E	1140	0	1115	62	0
1	F	1142	0	1120	80	0
1	G	1144	0	1120	77	0
1	H	1126	0	1098	54	0
1	I	1138	0	1109	57	0
1	J	1131	0	1095	83	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	15	0	0	0	0
2	D	10	0	0	0	0
2	E	15	0	0	0	0
2	F	10	0	0	0	0
2	G	10	0	0	1	0
2	H	15	0	0	0	0
2	I	10	0	0	0	0
2	J	15	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	J	1	0	0	0	0
4	B	10	0	0	0	0
4	C	5	0	0	0	0
4	D	15	0	0	0	0
4	E	10	0	0	0	0
4	G	5	0	0	0	0
4	H	10	0	0	1	0
4	I	15	0	0	0	0
4	J	5	0	0	0	0
5	A	4	0	0	0	0
5	C	3	0	0	0	0
5	D	3	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
5	I	2	0	0	0	0
5	J	2	0	0	0	0
All	All	11578	0	11100	592	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 26.

The worst 5 of 592 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:I:37:GLU:HG3	1:I:137:VAL:HG13	1.33	1.11
1:C:31:ARG:HD2	1:C:50:ILE:HD13	1.37	1.05
1:I:33:SER:HB3	1:I:133:LYS:HE2	1.39	1.02
1:G:50:ILE:H	1:G:50:ILE:HD13	1.24	0.99
1:D:31:ARG:HD2	1:D:50:ILE:HD13	1.44	0.98

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	146/158 (92%)	128 (88%)	14 (10%)	4 (3%)	5	21
1	B	145/158 (92%)	127 (88%)	17 (12%)	1 (1%)	22	52
1	C	145/158 (92%)	121 (83%)	19 (13%)	5 (3%)	3	17
1	D	145/158 (92%)	125 (86%)	17 (12%)	3 (2%)	7	26
1	E	146/158 (92%)	131 (90%)	13 (9%)	2 (1%)	11	36
1	F	146/158 (92%)	128 (88%)	13 (9%)	5 (3%)	3	17
1	G	146/158 (92%)	130 (89%)	12 (8%)	4 (3%)	5	21
1	H	144/158 (91%)	127 (88%)	13 (9%)	4 (3%)	5	21
1	I	146/158 (92%)	127 (87%)	15 (10%)	4 (3%)	5	21
1	J	145/158 (92%)	128 (88%)	14 (10%)	3 (2%)	7	26
All	All	1454/1580 (92%)	1272 (88%)	147 (10%)	35 (2%)	6	23

5 of 35 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	121(B)	GLU

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Mol	Chain	Res	Type
1	D	85	GLY
1	D	121(B)	GLU
1	H	121(B)	GLU
1	I	121(B)	GLU

### 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	114/126 (90%)	108 (95%)	6 (5%)	22	51
1	B	113/126 (90%)	108 (96%)	5 (4%)	28	58
1	C	112/126 (89%)	106 (95%)	6 (5%)	22	51
1	D	114/126 (90%)	109 (96%)	5 (4%)	28	58
1	E	114/126 (90%)	111 (97%)	3 (3%)	46	72
1	F	114/126 (90%)	109 (96%)	5 (4%)	28	58
1	G	114/126 (90%)	108 (95%)	6 (5%)	22	51
1	H	112/126 (89%)	107 (96%)	5 (4%)	27	58
1	I	113/126 (90%)	112 (99%)	1 (1%)	78	90
1	J	113/126 (90%)	106 (94%)	7 (6%)	18	46
All	All	1133/1260 (90%)	1084 (96%)	49 (4%)	29	59

5 of 49 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	22	TRP
1	F	41	LYS
1	J	121(C)	SER
1	E	152	ARG
1	F	53	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 30 such sidechains are listed below:



Mol	Chain	Res	Type
1	F	98	ASN
1	F	129	HIS
1	J	121(A)	HIS
1	F	121(A)	HIS
1	F	145	GLN

### 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 46 ligands modelled in this entry, 7 are monoatomic - leaving 39 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	PO4	D	207	-	4,4,4	1.58	0	6,6,6	0.42	0
4	SO4	J	225	-	4,4,4	0.26	0	6,6,6	0.15	0
4	SO4	D	211	-	4,4,4	0.28	0	6,6,6	0.08	0
4	SO4	I	220	-	4,4,4	0.31	0	6,6,6	0.12	0
4	SO4	E	224	-	4,4,4	0.28	0	6,6,6	0.11	0
2	PO4	A	201	-	4,4,4	1.54	0	6,6,6	0.46	0
4	SO4	B	205	-	4,4,4	0.27	0	6,6,6	0.13	0
2	PO4	J	224	-	4,4,4	1.66	0	6,6,6	0.41	0
2	PO4	G	213	-	4,4,4	1.59	0	6,6,6	0.45	0
4	SO4	I	219	-	4,4,4	0.29	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	PO4	H	216	-	4,4,4	1.63	0	6,6,6	0.43	0
2	PO4	A	221	-	4,4,4	1.60	0	6,6,6	0.42	0
2	PO4	G	214	-	4,4,4	1.61	0	6,6,6	0.42	0
2	PO4	E	209	-	4,4,4	1.65	0	6,6,6	0.43	0
4	SO4	B	206	-	4,4,4	0.26	0	6,6,6	0.05	0
2	PO4	F	212	-	4,4,4	1.54	0	6,6,6	0.43	0
2	PO4	F	211	-	4,4,4	1.58	0	6,6,6	0.45	0
4	SO4	C	209	-	4,4,4	0.27	0	6,6,6	0.11	0
2	PO4	I	218	-	4,4,4	1.56	0	6,6,6	0.44	0
2	PO4	J	220	-	4,4,4	1.64	0	6,6,6	0.42	0
2	PO4	D	210	-	4,4,4	1.63	0	6,6,6	0.42	0
2	PO4	B	204	-	4,4,4	1.61	0	6,6,6	0.42	0
2	PO4	C	206	-	4,4,4	1.58	0	6,6,6	0.43	0
2	PO4	C	208	-	4,4,4	1.63	0	6,6,6	0.42	0
4	SO4	D	212	-	4,4,4	0.27	0	6,6,6	0.07	0
4	SO4	H	224	-	4,4,4	0.28	0	6,6,6	0.13	0
2	PO4	H	215	-	4,4,4	1.60	0	6,6,6	0.41	0
2	PO4	H	223	-	4,4,4	1.65	0	6,6,6	0.43	0
4	SO4	I	221	-	4,4,4	0.28	0	6,6,6	0.05	0
4	SO4	D	213	-	4,4,4	0.27	0	6,6,6	0.06	0
2	PO4	B	203	-	4,4,4	1.56	0	6,6,6	0.45	0
4	SO4	H	225	-	4,4,4	0.30	0	6,6,6	0.07	0
2	PO4	C	205	-	4,4,4	1.64	0	6,6,6	0.44	0
2	PO4	J	219	-	4,4,4	1.61	0	6,6,6	0.42	0
2	PO4	E	202	-	4,4,4	1.58	0	6,6,6	0.44	0
4	SO4	G	215	-	4,4,4	0.28	0	6,6,6	0.12	0
2	PO4	E	222	-	4,4,4	1.63	0	6,6,6	0.43	0
2	PO4	I	217	-	4,4,4	1.57	0	6,6,6	0.43	0
4	SO4	E	223	-	4,4,4	0.28	0	6,6,6	0.17	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	214	PO4	1	0
4	H	224	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 [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	148/158 (93%)	-0.45	1 (0%) 87 72	5, 20, 39, 52	0
1	B	147/158 (93%)	-0.35	0 100 100	7, 22, 43, 54	0
1	C	147/158 (93%)	-0.44	1 (0%) 87 72	4, 27, 45, 58	0
1	D	147/158 (93%)	-0.38	0 100 100	6, 20, 45, 57	0
1	E	148/158 (93%)	-0.55	0 100 100	6, 18, 35, 43	0
1	F	148/158 (93%)	-0.46	0 100 100	5, 19, 41, 52	0
1	G	148/158 (93%)	-0.27	0 100 100	7, 26, 47, 52	0
1	H	146/158 (92%)	-0.25	0 100 100	9, 23, 42, 56	0
1	I	148/158 (93%)	-0.30	0 100 100	4, 20, 41, 57	0
1	J	147/158 (93%)	-0.52	0 100 100	5, 16, 37, 50	0
All	All	1474/1580 (93%)	-0.40	2 (0%) 95 91	4, 21, 43, 58	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	11	THR	2.6
1	A	155	ALA	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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
4	SO4	H	225	5/5	0.77	0.35	45,45,47,47	5
4	SO4	C	209	5/5	0.78	0.34	25,25,26,27	5
4	SO4	D	211	5/5	0.78	0.54	42,43,43,44	5
3	NA	F	303	1/1	0.82	0.28	19,19,19,19	0
2	PO4	A	221	5/5	0.82	0.36	28,29,30,30	5
2	PO4	B	203	5/5	0.83	0.37	41,42,43,43	5
4	SO4	I	220	5/5	0.83	0.32	26,27,28,29	5
2	PO4	A	201	5/5	0.85	0.34	38,39,41,41	5
2	PO4	G	214	5/5	0.86	0.32	45,46,46,47	5
4	SO4	D	213	5/5	0.87	0.41	45,45,46,46	5
4	SO4	D	212	5/5	0.87	0.26	37,38,38,39	5
4	SO4	I	221	5/5	0.87	0.32	41,41,42,42	5
2	PO4	E	222	5/5	0.87	0.35	36,36,37,37	5
2	PO4	D	207	5/5	0.88	0.27	33,33,34,35	5
2	PO4	J	224	5/5	0.88	0.27	42,43,43,44	5
2	PO4	I	217	5/5	0.88	0.24	35,35,36,37	5
3	NA	G	304	1/1	0.89	0.20	21,21,21,21	0
4	SO4	E	223	5/5	0.89	0.21	39,41,42,43	5
4	SO4	G	215	5/5	0.90	0.24	21,21,22,23	5
4	SO4	B	206	5/5	0.90	0.47	31,31,32,32	5
2	PO4	H	216	5/5	0.90	0.25	31,31,32,32	5
2	PO4	D	210	5/5	0.90	0.25	34,34,35,35	5
2	PO4	C	206	5/5	0.91	0.37	21,22,23,23	5
3	NA	D	302	1/1	0.91	0.29	21,21,21,21	0
3	NA	A	301	1/1	0.91	0.20	2,2,2,2	0
2	PO4	E	209	5/5	0.91	0.27	34,35,36,37	5
2	PO4	F	212	5/5	0.93	0.21	17,17,19,19	5
2	PO4	H	223	5/5	0.93	0.23	16,17,18,18	5
2	PO4	E	202	5/5	0.93	0.22	16,16,18,19	5
2	PO4	G	213	5/5	0.93	0.23	31,31,32,32	5
2	PO4	B	204	5/5	0.93	0.23	20,20,21,21	5
4	SO4	J	225	5/5	0.93	0.23	19,19,20,21	5
2	PO4	C	208	5/5	0.93	0.20	24,24,25,25	5
3	NA	A	300	1/1	0.94	0.24	15,15,15,15	0
2	PO4	J	219	5/5	0.94	0.19	36,36,37,38	5
2	PO4	F	211	5/5	0.94	0.19	26,26,29,29	5
4	SO4	E	224	5/5	0.94	0.19	22,23,24,24	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	PO4	J	220	5/5	0.94	0.23	26,27,27,28	5
4	SO4	H	224	5/5	0.94	0.21	21,21,22,23	5
2	PO4	H	215	5/5	0.94	0.21	28,29,30,31	5
4	SO4	B	205	5/5	0.95	0.16	21,22,23,24	5
4	SO4	I	219	5/5	0.95	0.18	17,17,19,19	5
3	NA	B	306	1/1	0.95	0.39	21,21,21,21	0
2	PO4	C	205	5/5	0.95	0.20	32,34,34,35	5
2	PO4	I	218	5/5	0.96	0.17	16,16,18,18	5
3	NA	J	305	1/1	0.96	0.15	18,18,18,18	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.