



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 11:22 am GMT

PDB ID : 3URH  
Title : Crystal structure of a dihydrolipoamide dehydrogenase from *Sinorhizobium meliloti* 1021  
Authors : Kumaran, D.; Chamala, S.; Evans, B.; Foti, R.; Gizzi, A.; Hillerich, B.; Kar, A.; LaFleur, J.; Seidel, R.; Villigas, G.; Zencheck, W.; Almo, S.C.; Swaminathan, S.; New York Structural Genomics Research Consortium (NYSGRG)  
Deposited on : 2011-11-22  
Resolution : 1.90 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

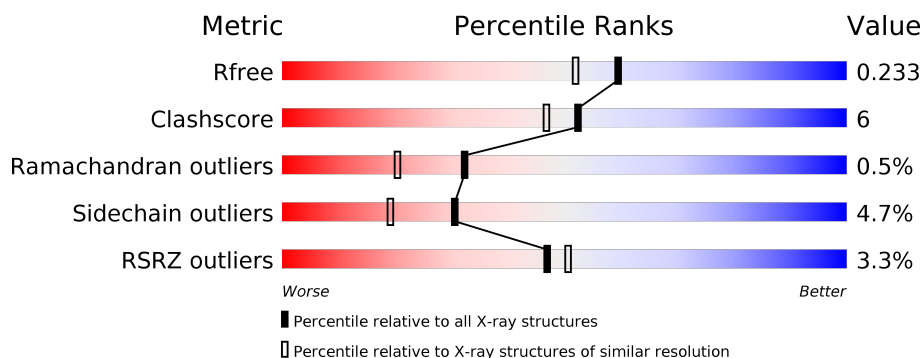
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
$R_{free}$	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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	491	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>6%</div> </div> </div>
1	B	491	<div> <div>3%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>5%</div> </div> </div>

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	FAD	B	500	-	-	-	X
3	SO4	B	501	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7292 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 Dihydrolipoyl dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	Se	0	0	0
			3365	2127	578	644	4	12			
1	B	465	Total	C	N	O	S	Se	0	0	0
			3399	2150	585	648	4	12			

There are 46 discrepancies between the modelled and reference sequences:

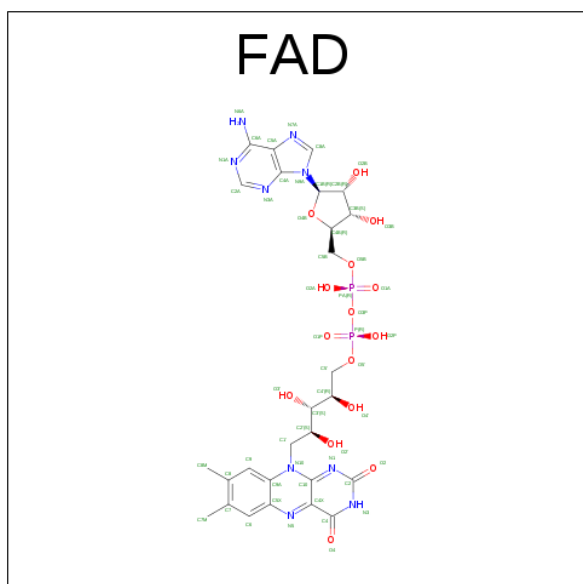
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MSE	-	EXPRESSION TAG	UNP Q92LK0
A	-21	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-20	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-19	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-18	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-17	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-16	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-15	SER	-	EXPRESSION TAG	UNP Q92LK0
A	-14	SER	-	EXPRESSION TAG	UNP Q92LK0
A	-13	GLY	-	EXPRESSION TAG	UNP Q92LK0
A	-12	VAL	-	EXPRESSION TAG	UNP Q92LK0
A	-11	ASP	-	EXPRESSION TAG	UNP Q92LK0
A	-10	LEU	-	EXPRESSION TAG	UNP Q92LK0
A	-9	GLY	-	EXPRESSION TAG	UNP Q92LK0
A	-8	THR	-	EXPRESSION TAG	UNP Q92LK0
A	-7	GLU	-	EXPRESSION TAG	UNP Q92LK0
A	-6	ASN	-	EXPRESSION TAG	UNP Q92LK0
A	-5	LEU	-	EXPRESSION TAG	UNP Q92LK0
A	-4	TYR	-	EXPRESSION TAG	UNP Q92LK0
A	-3	PHE	-	EXPRESSION TAG	UNP Q92LK0
A	-2	GLN	-	EXPRESSION TAG	UNP Q92LK0
A	-1	SER	-	EXPRESSION TAG	UNP Q92LK0
A	0	MSE	-	EXPRESSION TAG	UNP Q92LK0
B	-22	MSE	-	EXPRESSION TAG	UNP Q92LK0
B	-21	HIS	-	EXPRESSION TAG	UNP Q92LK0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-19	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-18	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-17	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-16	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-15	SER	-	EXPRESSION TAG	UNP Q92LK0
B	-14	SER	-	EXPRESSION TAG	UNP Q92LK0
B	-13	GLY	-	EXPRESSION TAG	UNP Q92LK0
B	-12	VAL	-	EXPRESSION TAG	UNP Q92LK0
B	-11	ASP	-	EXPRESSION TAG	UNP Q92LK0
B	-10	LEU	-	EXPRESSION TAG	UNP Q92LK0
B	-9	GLY	-	EXPRESSION TAG	UNP Q92LK0
B	-8	THR	-	EXPRESSION TAG	UNP Q92LK0
B	-7	GLU	-	EXPRESSION TAG	UNP Q92LK0
B	-6	ASN	-	EXPRESSION TAG	UNP Q92LK0
B	-5	LEU	-	EXPRESSION TAG	UNP Q92LK0
B	-4	TYR	-	EXPRESSION TAG	UNP Q92LK0
B	-3	PHE	-	EXPRESSION TAG	UNP Q92LK0
B	-2	GLN	-	EXPRESSION TAG	UNP Q92LK0
B	-1	SER	-	EXPRESSION TAG	UNP Q92LK0
B	0	MSE	-	EXPRESSION TAG	UNP Q92LK0

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		

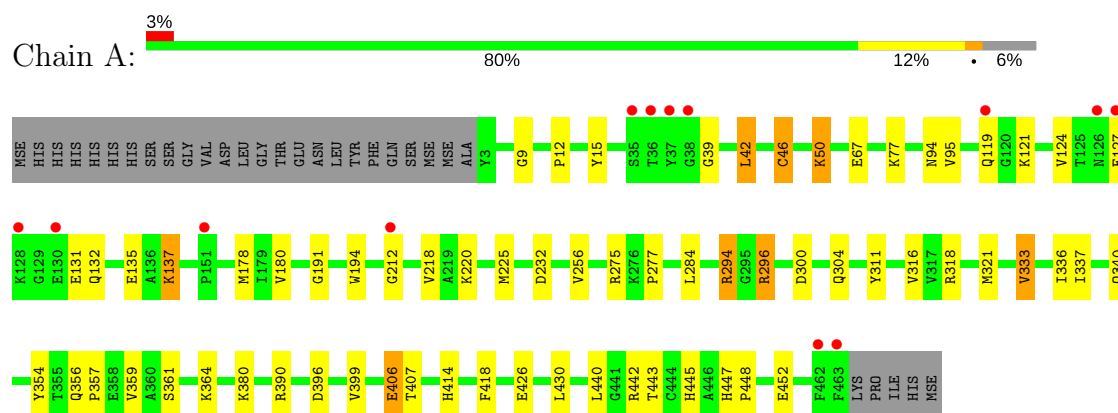
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	230	Total	O	0	0
			230	230		
5	B	178	Total	O	0	0
			178	178		

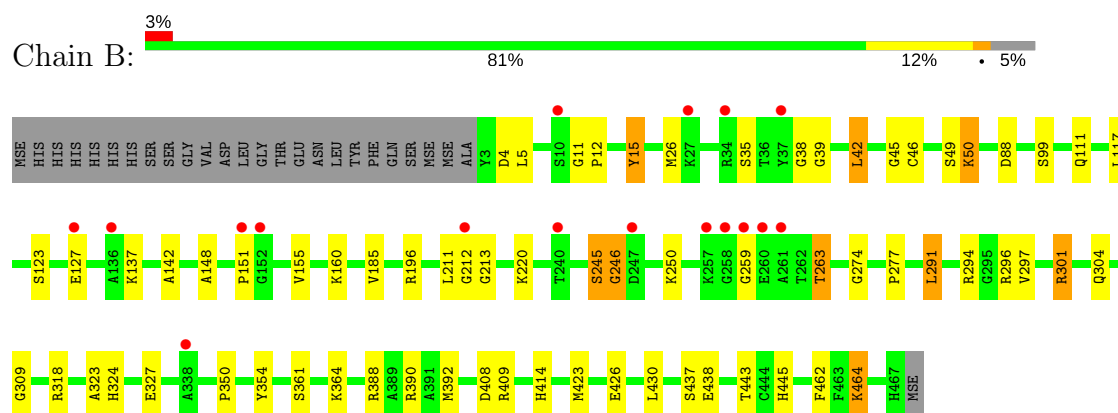
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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: Dihydrolipoyl dehydrogenase



#### • Molecule 1: Dihydrolipoyl dehydrogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.43Å 104.72Å 120.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.77 – 1.90 40.36 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.5 (42.77-1.90) 98.5 (40.36-1.90)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.79 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.184 , 0.227 0.192 , 0.233	Depositor DCC
$R_{free}$ test set	3721 reflections (5.33%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.9	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 41.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7292	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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: SO4, FAD, EDO

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	1.10	3/3400 (0.1%)	1.05	9/4567 (0.2%)
1	B	1.07	0/3436	1.03	7/4616 (0.2%)
All	All	1.08	3/6836 (0.0%)	1.04	16/9183 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	9	GLY	N-CA	6.97	1.56	1.46
1	A	354	TYR	CG-CD2	5.13	1.45	1.39
1	A	194	TRP	CD2-CE2	5.11	1.47	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	430	LEU	CB-CG-CD1	7.13	123.12	111.00
1	A	333	VAL	CG1-CB-CG2	-6.26	100.88	110.90
1	A	430	LEU	CA-CB-CG	-6.16	101.13	115.30
1	A	300	ASP	CB-CG-OD1	6.07	123.76	118.30
1	B	196	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	B	88	ASP	CB-CG-OD1	5.74	123.46	118.30
1	B	423	MSE	CA-CB-CG	-5.67	103.66	113.30
1	A	390	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	77	LYS	CD-CE-NZ	5.52	124.40	111.70
1	A	440	LEU	CB-CG-CD2	-5.48	101.69	111.00
1	B	245	SER	N-CA-C	-5.45	96.29	111.00
1	A	321	MSE	CA-CB-CG	-5.44	104.06	113.30
1	B	196	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	B	409	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	A	275	ARG	NE-CZ-NH1	5.29	122.94	120.30
1	A	284	LEU	CB-CG-CD1	-5.07	102.39	111.00

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	3365	0	3439	39	0
1	B	3399	0	3477	49	0
2	A	53	0	31	2	0
2	B	53	0	31	6	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	B	4	0	6	3	0
5	A	230	0	0	4	0
5	B	178	0	0	0	0
All	All	7292	0	6984	87	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 (87) 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:185:VAL:H	4:B:502:EDO:H21	1.19	1.07
1:A:232:ASP:HB2	5:A:640:HOH:O	1.61	0.99
1:B:304:GLN:HE21	1:B:309:GLY:H	0.97	0.92
1:B:4:ASP:HB3	1:B:26:MSE:HE3	1.49	0.91
1:B:185:VAL:N	4:B:502:EDO:H21	1.90	0.87
1:B:304:GLN:NE2	1:B:309:GLY:H	1.73	0.86
1:A:426:GLU:OE2	1:A:445:HIS:HE1	1.65	0.79
1:A:225:MSE:HE1	1:A:396:ASP:HB3	1.66	0.78
1:B:426:GLU:OE2	1:B:445:HIS:HE1	1.66	0.77
1:A:406:GLU:HG2	1:A:407:THR:HG23	1.66	0.76
1:B:185:VAL:H	4:B:502:EDO:C2	1.99	0.75
1:B:50:LYS:N	1:B:50:LYS:HD3	2.06	0.70
1:A:225:MSE:HE3	1:A:357:PRO:HG3	1.75	0.69
1:A:361:SER:OG	1:A:414:HIS:HD2	1.78	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:HIS:H	1:B:324:HIS:CD2	2.14	0.66
1:B:324:HIS:H	1:B:324:HIS:CD2	2.14	0.65
1:A:311:TYR:CE1	1:A:337:ILE:HD11	2.32	0.64
1:A:67:GLU:HG2	5:A:677:HOH:O	1.98	0.64
1:B:4:ASP:HB3	1:B:26:MSE:CE	2.25	0.63
1:A:12:PRO:HG2	2:A:500:FAD:O5'	1.99	0.63
1:A:232:ASP:CB	5:A:640:HOH:O	2.34	0.62
1:B:5:LEU:HB2	1:B:26:MSE:HE1	1.82	0.61
1:A:42:LEU:HD13	1:A:94:ASN:HB3	1.83	0.61
1:A:443:THR:O	1:A:445:HIS:HD2	1.84	0.61
1:A:311:TYR:CE1	1:A:337:ILE:CD1	2.84	0.61
1:A:426:GLU:OE2	1:A:445:HIS:CE1	2.52	0.60
1:B:250:LYS:HG3	1:B:263:THR:HG23	1.83	0.60
1:B:324:HIS:HE1	1:B:350:PRO:O	1.85	0.59
1:A:46:CYS:O	1:A:50:LYS:HD2	2.03	0.58
1:B:277:PRO:HG3	1:B:296:ARG:HG2	1.86	0.58
1:B:11:GLY:HA3	2:B:500:FAD:O5B	2.04	0.58
1:B:443:THR:O	1:B:445:HIS:HD2	1.87	0.57
1:B:4:ASP:CB	1:B:26:MSE:HE3	2.29	0.57
1:B:4:ASP:CB	1:B:26:MSE:CE	2.83	0.56
1:B:12:PRO:HD2	2:B:500:FAD:O5'	2.06	0.56
1:A:178:MSE:HE3	1:A:180:VAL:CG2	2.35	0.56
1:B:142:ALA:O	2:B:500:FAD:H52A	2.06	0.55
1:B:426:GLU:OE2	1:B:445:HIS:CE1	2.54	0.54
1:A:225:MSE:CE	1:A:396:ASP:HB3	2.36	0.54
1:B:12:PRO:HA	1:B:15:TYR:CD2	2.44	0.53
1:B:388:ARG:HG2	1:B:392:MSE:HE2	1.91	0.53
1:B:408:ASP:OD1	1:B:437:SER:OG	2.26	0.53
1:B:117:LEU:HD11	1:B:123:SER:HB3	1.91	0.52
1:A:333:VAL:O	1:A:337:ILE:HG12	2.09	0.52
1:A:304:GLN:HG2	1:A:311:TYR:CE2	2.45	0.51
1:B:45:GLY:O	1:B:49:SER:HB2	2.11	0.50
1:B:5:LEU:HB2	1:B:26:MSE:CE	2.41	0.50
1:B:291:LEU:HD12	1:B:297:VAL:HA	1.94	0.49
1:A:277:PRO:HG3	1:A:296:ARG:HG3	1.93	0.49
1:B:361:SER:OG	1:B:414:HIS:HD2	1.96	0.49
1:A:364:LYS:O	1:A:414:HIS:HE1	1.96	0.48
1:B:304:GLN:HE21	1:B:309:GLY:N	1.83	0.48
1:A:178:MSE:HE3	1:A:180:VAL:HG22	1.95	0.48
1:A:178:MSE:HE1	1:A:191:GLY:HA2	1.96	0.47
1:B:38:GLY:HA3	1:B:42:LEU:HD22	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:ALA:O	1:B:327:GLU:HG3	2.14	0.47
1:A:131:GLU:HG3	1:A:132:GLN:N	2.30	0.47
1:A:406:GLU:CG	1:A:407:THR:HG23	2.40	0.47
1:A:356:GLN:HE22	1:A:418:PHE:HA	1.79	0.46
1:A:442:ARG:HD3	5:A:669:HOH:O	2.15	0.46
1:B:50:LYS:HE3	1:B:354:TYR:CD2	2.51	0.46
1:A:39:GLY:HA2	2:A:500:FAD:O3B	2.16	0.46
1:B:464:LYS:HE3	1:B:464:LYS:HB3	1.88	0.46
1:A:137:LYS:HD3	1:A:137:LYS:H	1.81	0.46
1:B:12:PRO:HA	1:B:15:TYR:CE2	2.52	0.45
1:B:38:GLY:CA	1:B:42:LEU:HD22	2.46	0.45
1:A:294:ARG:HH11	1:A:294:ARG:CB	2.30	0.45
1:B:438:GLU:HB2	1:B:462:PHE:CD1	2.51	0.45
1:B:4:ASP:HB2	1:B:26:MSE:HE2	1.99	0.45
1:B:250:LYS:HE3	1:B:263:THR:HG21	1.99	0.44
2:B:500:FAD:H1'1	2:B:500:FAD:H9	1.83	0.44
1:A:218:VAL:HG11	1:A:359:VAL:HG13	2.00	0.43
1:A:94:ASN:ND2	1:B:390:ARG:NH1	2.66	0.43
1:A:336:ILE:HA	1:A:340:GLN:O	2.19	0.42
1:B:364:LYS:O	1:B:414:HIS:HE1	2.02	0.42
1:B:35:SER:HA	1:B:111:GLN:NE2	2.34	0.42
1:B:148:ALA:HB3	1:B:274:GLY:O	2.19	0.42
1:A:380:LYS:HA	1:A:399:VAL:O	2.20	0.41
1:B:301:ARG:HE	1:B:301:ARG:HB3	1.46	0.41
1:A:294:ARG:HB3	1:A:294:ARG:HH11	1.85	0.41
1:B:212:GLY:HA3	1:B:213:GLY:HA3	1.61	0.41
1:B:245:SER:O	1:B:246:GLY:C	2.59	0.41
1:B:39:GLY:HA2	2:B:500:FAD:O3B	2.20	0.41
1:B:50:LYS:CE	2:B:500:FAD:O4	2.69	0.41
1:A:448:PRO:HA	1:A:452:GLU:OE1	2.21	0.41
1:A:178:MSE:CE	1:A:191:GLY:HA2	2.51	0.41
1:A:124:VAL:O	1:A:131:GLU:HA	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries

of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	459/491 (94%)	443 (96%)	15 (3%)	1 (0%)	51	41
1	B	463/491 (94%)	442 (96%)	17 (4%)	4 (1%)	20	8
All	All	922/982 (94%)	885 (96%)	32 (4%)	5 (0%)	32	20

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	246	GLY
1	B	151	PRO
1	B	211	LEU
1	B	259	GLY
1	A	212	GLY

### 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	348/359 (97%)	331 (95%)	17 (5%)	29	17
1	B	352/359 (98%)	336 (96%)	16 (4%)	32	21
All	All	700/718 (98%)	667 (95%)	33 (5%)	30	19

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	TYR
1	A	42	LEU
1	A	46	CYS
1	A	50	LYS
1	A	95	VAL
1	A	119	GLN
1	A	121	LYS

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Mol	Chain	Res	Type
1	A	127	GLU
1	A	135	GLU
1	A	137	LYS
1	A	220	LYS
1	A	256	VAL
1	A	294	ARG
1	A	296	ARG
1	A	316	VAL
1	A	318	ARG
1	A	406	GLU
1	B	15	TYR
1	B	42	LEU
1	B	46	CYS
1	B	50	LYS
1	B	99	SER
1	B	127	GLU
1	B	137	LYS
1	B	155	VAL
1	B	160	LYS
1	B	220	LYS
1	B	263	THR
1	B	291	LEU
1	B	294	ARG
1	B	301	ARG
1	B	318	ARG
1	B	464	LYS

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

Mol	Chain	Res	Type
1	A	61	GLN
1	A	75	ASN
1	A	94	ASN
1	A	356	GLN
1	A	414	HIS
1	A	445	HIS
1	B	61	GLN
1	B	63	GLN
1	B	64	HIS
1	B	304	GLN
1	B	324	HIS
1	B	414	HIS

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Mol	Chain	Res	Type
1	B	445	HIS
1	B	467	HIS

### 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 ⓘ

5 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	FAD	A	500	-	51,58,58	2.09	10 (19%)	54,89,89	2.59	13 (24%)
3	SO4	A	501	-	4,4,4	0.60	0	6,6,6	0.60	0
2	FAD	B	500	-	51,58,58	1.97	11 (21%)	54,89,89	2.63	19 (35%)
3	SO4	B	501	-	4,4,4	0.78	0	6,6,6	0.81	0
4	EDO	B	502	-	3,3,3	0.67	0	2,2,2	1.47	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
2	FAD	A	500	-	-	0/28/50/50	0/6/6/6
3	SO4	A	501	-	-	0/0/0/0	0/0/0/0
2	FAD	B	500	-	-	0/28/50/50	0/6/6/6
3	SO4	B	501	-	-	0/0/0/0	0/0/0/0
4	EDO	B	502	-	-	0/1/1/1	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	FAD	O4B-C1B	2.38	1.44	1.41
2	A	500	FAD	C4A-N3A	2.48	1.39	1.35
2	B	500	FAD	C4X-N5	2.58	1.37	1.33
2	B	500	FAD	P-O1P	2.67	1.60	1.50
2	A	500	FAD	C4-C4X	2.72	1.46	1.41
2	A	500	FAD	C4-N3	2.97	1.38	1.33
2	B	500	FAD	C9A-N10	3.06	1.42	1.38
2	A	500	FAD	C4X-N5	3.32	1.38	1.33
2	B	500	FAD	C2A-N3A	3.37	1.37	1.32
2	A	500	FAD	C2A-N3A	3.41	1.37	1.32
2	B	500	FAD	C5A-C4A	3.72	1.48	1.40
2	A	500	FAD	C5A-C4A	3.82	1.49	1.40
2	A	500	FAD	C9A-C5X	3.98	1.50	1.42
2	B	500	FAD	C9A-C5X	4.00	1.50	1.42
2	A	500	FAD	C8-C7	4.26	1.51	1.41
2	B	500	FAD	C8-C7	4.27	1.51	1.41
2	B	500	FAD	C4-C4X	4.57	1.50	1.41
2	B	500	FAD	C4X-C10	4.88	1.49	1.41
2	B	500	FAD	C10-N1	5.03	1.40	1.33
2	A	500	FAD	C4X-C10	6.27	1.52	1.41
2	A	500	FAD	C10-N1	6.28	1.42	1.33

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	C4X-C10-N10	-8.02	114.95	120.52
2	B	500	FAD	N3A-C2A-N1A	-7.50	122.32	128.86
2	B	500	FAD	C4X-C10-N10	-7.17	115.54	120.52
2	A	500	FAD	N3A-C2A-N1A	-5.05	124.46	128.86
2	A	500	FAD	C4X-C4-N3	-5.03	116.32	123.48
2	A	500	FAD	C4-C4X-C10	-4.84	116.04	119.96
2	B	500	FAD	C4-C4X-C10	-4.78	116.09	119.96
2	B	500	FAD	C4X-C4-N3	-3.91	117.91	123.48
2	B	500	FAD	C4A-C5A-N7A	-3.64	105.89	109.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	O2B-C2B-C1B	-3.47	100.77	111.61
2	B	500	FAD	C4'-C3'-C2'	-3.29	106.32	113.41
2	B	500	FAD	O3B-C3B-C4B	-2.63	103.40	111.09
2	B	500	FAD	C9A-C5X-N5	-2.19	118.98	122.24
2	A	500	FAD	O3B-C3B-C4B	-2.16	104.77	111.09
2	B	500	FAD	O2P-P-O1P	-2.10	101.41	112.28
2	B	500	FAD	C6-C5X-C9A	2.01	121.61	119.00
2	B	500	FAD	C5X-C9A-N10	2.04	119.17	117.66
2	B	500	FAD	N6A-C6A-N1A	2.10	122.94	118.77
2	B	500	FAD	O5'-C5'-C4'	2.15	115.09	109.36
2	A	500	FAD	C4X-N5-C5X	2.21	119.10	116.76
2	B	500	FAD	C4-C4X-N5	2.26	121.16	118.68
2	A	500	FAD	C5X-C9A-N10	2.27	119.34	117.66
2	B	500	FAD	C2B-C3B-C4B	2.29	107.08	102.62
2	A	500	FAD	N6A-C6A-N1A	2.39	123.50	118.77
2	B	500	FAD	O5B-C5B-C4B	2.76	118.78	109.00
2	A	500	FAD	C10-C4X-N5	3.02	124.07	120.59
2	A	500	FAD	O2P-P-O1P	3.33	129.53	112.28
2	A	500	FAD	C1'-N10-C9A	3.51	121.56	118.35
2	B	500	FAD	O5'-P-O1P	4.59	127.78	109.25
2	B	500	FAD	C4X-N5-C5X	5.39	122.45	116.76
2	B	500	FAD	C4-N3-C2	8.39	122.50	115.16
2	A	500	FAD	C4-N3-C2	10.72	124.53	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	FAD	2	0
2	B	500	FAD	6	0
4	B	502	EDO	3	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	449/491 (91%)	0.01	13 (2%) 52 56	11, 22, 45, 80	0
1	B	453/491 (92%)	0.21	17 (3%) 41 45	11, 26, 51, 79	0
All	All	902/982 (91%)	0.11	30 (3%) 47 50	11, 24, 48, 80	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	128	LYS	4.5
1	A	463	PHE	4.2
1	B	257	LYS	4.1
1	A	462	PHE	4.0
1	B	151	PRO	4.0
1	B	258	GLY	3.9
1	A	127	GLU	3.5
1	A	126	ASN	3.2
1	B	34	ARG	3.2
1	A	38	GLY	3.1
1	A	36	THR	3.0
1	A	35	SER	3.0
1	B	10	SER	2.7
1	B	260	GLU	2.7
1	B	240	THR	2.7
1	A	130	GLU	2.6
1	B	127	GLU	2.6
1	B	212	GLY	2.5
1	A	37	TYR	2.5
1	B	261	ALA	2.4
1	B	152	GLY	2.4
1	B	247	ASP	2.4
1	B	136	ALA	2.4
1	A	151	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	338	ALA	2.3
1	B	37	TYR	2.3
1	A	212	GLY	2.2
1	B	27	LYS	2.2
1	B	259	GLY	2.1
1	A	119	GLN	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( $\text{\AA}^2$ )	Q<0.9
3	SO4	B	501	5/5	0.96	0.19	2.16	34,38,42,43	0
2	FAD	B	500	53/53	0.85	0.23	2.14	28,42,50,57	0
2	FAD	A	500	53/53	0.87	0.19	1.73	25,37,46,50	0
4	EDO	B	502	4/4	0.87	0.17	0.93	32,33,35,46	0
3	SO4	A	501	5/5	0.91	0.12	-	60,64,65,67	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.