



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 12:48 PM 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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

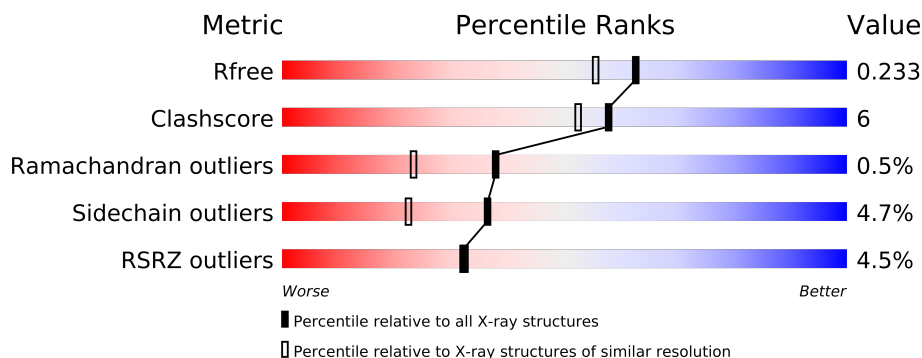
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	491	
1	B	491	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
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 hydrogen and 0 are deuterium.

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:

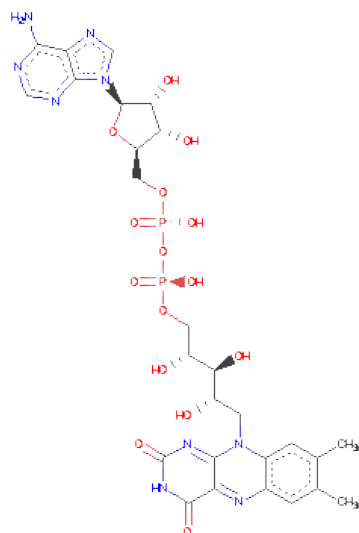
Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MSE	-	EXPRESSION TAG	UNP Q92LK0
A	-21	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-20	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-19	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-18	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-17	HIS	-	EXPRESSION TAG	UNP Q92LK0
A	-16	HIS	-	EXPRESSION TAG	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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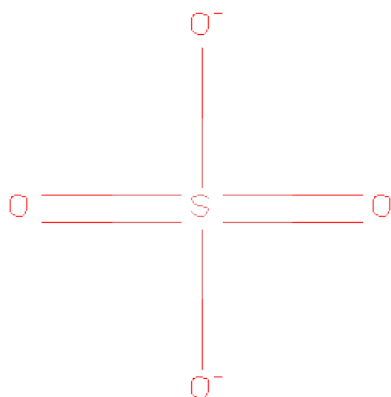
Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-19	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-18	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-17	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-16	HIS	-	EXPRESSION TAG	UNP Q92LK0
B	-15	SER	-	EXPRESSION TAG	UNP Q92LK0
B	-14	SER	-	EXPRESSION TAG	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: $C_{27}H_{33}N_9O_{15}P_2$).



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: O₄S).



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: C₂H₆O₂).



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

- Molecule 5 is water.

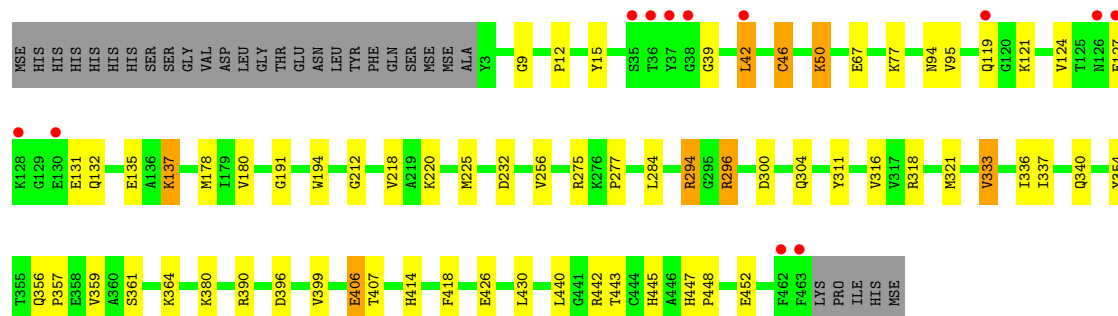
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	230	Total	O	0	0
			230	230		
5	B	178	Total	O	0	0
			178	178		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

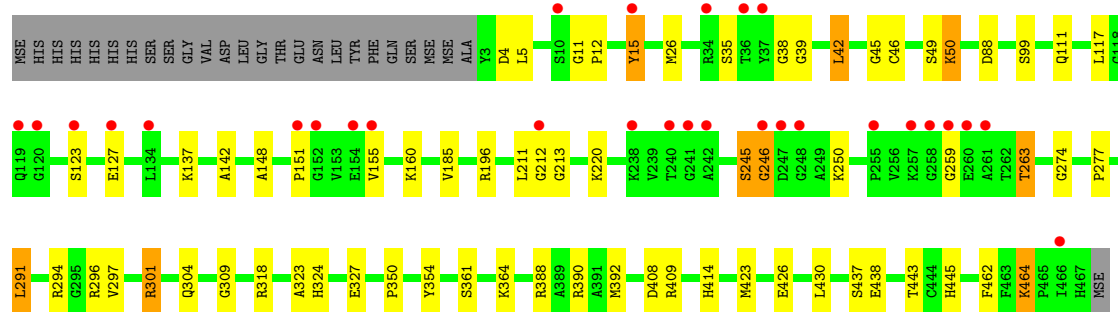
- Molecule 1: Dihydrolipoyl dehydrogenase

Chain A: 



- Molecule 1: Dihydrolipoyl dehydrogenase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	4.79 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.184 , 0.227 0.193 , 0.233	Depositor DCC
R_{free} test set	3721 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 73503 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7292	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (87) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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	Distance(Å)	Clash(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:B:361:SER:OG	1:B:414:HIS:HD2	1.96	0.49
1:A:277:PRO:HG3	1:A:296:ARG:HG3	1.93	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
1:B:323:ALA:O	1:B:327:GLU:HG3	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:406:GLU:CG	1:A:407:THR:HG23	2.40	0.47
1:A:131:GLU:HG3	1:A:132:GLN:N	2.30	0.47
1:A:442:ARG:HD3	5:A:669:HOH:O	2.15	0.46
1:A:356:GLN:HE22	1:A:418:PHE:HA	1.79	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:B:364:LYS:O	1:B:414:HIS:HE1	2.02	0.42
1:A:336:ILE:HA	1:A:340:GLN:O	2.19	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:HH11	1:A:294:ARG:HB3	1.85	0.41
1:B:245:SER:O	1:B:246:GLY:C	2.59	0.41
1:B:212:GLY:HA3	1:B:213:GLY:HA3	1.61	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%)	56	44
1	B	463/491 (94%)	442 (96%)	17 (4%)	4 (1%)	25	10
All	All	922/982 (94%)	885 (96%)	32 (4%)	5 (0%)	38	23

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%)	35	21
1	B	352/359 (98%)	336 (96%)	16 (4%)	38	24
All	All	700/718 (98%)	667 (95%)	33 (5%)	36	22

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
1	A	127	GLU
1	A	135	GLU
1	A	137	LYS
1	A	220	LYS
1	A	256	VAL

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

5.3.3 RNA ⓘ

There are no RNA chains 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	-	58,58,58	1.85	12 (20%)	85,89,89	2.16	21 (24%)
3	SO4	A	501	-	4,4,4	0.73	0	6,6,6	0.61	0
2	FAD	B	500	-	58,58,58	1.78	12 (20%)	85,89,89	2.26	18 (21%)
3	SO4	B	501	-	4,4,4	1.26	0	6,6,6	0.86	0
4	EDO	B	502	-	3,3,3	0.72	0	2,2,2	1.50	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	-	-	0/34/50/50	0/1/6/6
3	SO4	A	501	-	-	0/0/0/0	0/0/0/0
2	FAD	B	500	-	-	0/34/50/50	0/1/6/6
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 (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	FAD	C4X-C10	6.40	1.52	1.40
2	B	500	FAD	C4-C4X	5.41	1.50	1.41
2	B	500	FAD	C4X-C10	4.99	1.49	1.40
2	B	500	FAD	C9A-C5X	4.01	1.50	1.42
2	A	500	FAD	C9A-C5X	3.99	1.50	1.42
2	A	500	FAD	C5A-C4A	3.82	1.49	1.40
2	B	500	FAD	C8-C7	3.78	1.51	1.40
2	A	500	FAD	C8-C7	3.77	1.51	1.40
2	B	500	FAD	C5A-C4A	3.72	1.48	1.40
2	A	500	FAD	C10-N1	3.58	1.42	1.35
2	A	500	FAD	P-O3P	3.48	1.66	1.59
2	A	500	FAD	C4-C4X	3.25	1.46	1.41
2	B	500	FAD	PA-O3P	3.18	1.65	1.59
2	A	500	FAD	C2A-N3A	2.88	1.37	1.32
2	B	500	FAD	C2A-N3A	2.84	1.37	1.32
2	B	500	FAD	C9A-N10	2.65	1.42	1.38
2	B	500	FAD	C10-N1	2.60	1.40	1.35
2	A	500	FAD	C1'-C2'	2.50	1.54	1.51
2	B	500	FAD	P-O1P	2.50	1.60	1.51
2	A	500	FAD	O4-C4	2.44	1.29	1.24
2	A	500	FAD	C4A-N3A	2.36	1.39	1.35
2	A	500	FAD	PA-O3P	2.21	1.63	1.59
2	B	500	FAD	O4B-C1B	2.09	1.44	1.41
2	B	500	FAD	P-O3P	2.04	1.63	1.59

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	C4X-C10-N10	-11.14	114.95	120.51
2	B	500	FAD	C4X-C10-N10	-9.96	115.54	120.51
2	B	500	FAD	N3A-C2A-N1A	-7.64	122.32	128.71
2	B	500	FAD	N3A-C4A-N9A	6.34	136.87	125.43
2	A	500	FAD	N3A-C4A-N9A	6.14	136.52	125.43
2	B	500	FAD	C4X-N5-C5X	5.13	122.45	116.69
2	A	500	FAD	N3A-C2A-N1A	-5.09	124.46	128.71
2	B	500	FAD	O5'-P-O1P	4.70	127.78	109.37
2	A	500	FAD	C2'-C1'-N10	-4.36	106.67	112.45
2	A	500	FAD	O4B-C1B-C2B	-4.24	100.27	106.77
2	B	500	FAD	C4A-C5A-N7A	-4.24	105.89	109.52
2	B	500	FAD	C5A-C4A-N3A	-4.04	116.91	125.70
2	A	500	FAD	C2B-C1B-N9A	3.61	122.54	113.27
2	A	500	FAD	C5A-C4A-N3A	-3.59	117.88	125.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	O2B-C2B-C1B	-3.46	100.77	111.23
2	B	500	FAD	C2-N1-C10	3.19	118.20	114.98
2	A	500	FAD	O4B-C1B-N9A	-3.10	105.55	108.44
2	A	500	FAD	O2P-P-O1P	3.10	129.53	112.21
2	B	500	FAD	C4'-C3'-C2'	-3.07	106.32	113.25
2	B	500	FAD	N1-C10-N10	2.99	123.84	115.97
2	A	500	FAD	C10-C4X-N5	2.98	124.07	120.45
2	B	500	FAD	C2A-N3A-C4A	2.96	122.43	114.01
2	B	500	FAD	C2B-C1B-N9A	2.85	120.58	113.27
2	A	500	FAD	C1'-N10-C9A	2.76	121.56	118.87
2	B	500	FAD	O5B-C5B-C4B	2.68	118.78	108.94
2	B	500	FAD	O3B-C3B-C4B	-2.61	103.40	111.08
2	A	500	FAD	C5X-C9A-N10	2.58	119.34	116.80
2	A	500	FAD	C5'-C4'-C3'	-2.53	107.29	112.06
2	B	500	FAD	C5X-C9A-N10	2.41	119.17	116.80
2	A	500	FAD	C8A-N9A-C4A	2.35	108.69	106.90
2	A	500	FAD	N1-C10-N10	2.27	121.93	115.97
2	A	500	FAD	C2A-N3A-C4A	2.26	120.46	114.01
2	B	500	FAD	C2B-C3B-C4B	2.22	107.08	102.65
2	B	500	FAD	C9A-C5X-N5	-2.21	118.98	122.37
2	A	500	FAD	C4A-C5A-N7A	-2.18	107.66	109.52
2	A	500	FAD	C4X-N5-C5X	2.14	119.10	116.69
2	A	500	FAD	O3B-C3B-C4B	-2.14	104.77	111.08
2	B	500	FAD	C4X-C10-N1	-2.11	120.62	122.73
2	A	500	FAD	N6A-C6A-N1A	2.11	123.50	119.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	461/491 (93%)	0.13	12 (2%) 53 55	11, 22, 44, 80	0
1	B	465/491 (94%)	0.36	29 (6%) 20 20	11, 26, 50, 79	0
All	All	926/982 (94%)	0.25	41 (4%) 32 33	11, 24, 48, 80	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	258	GLY	6.2
1	A	463	PHE	6.0
1	A	36	THR	4.5
1	A	38	GLY	4.1
1	A	462	PHE	3.9
1	B	247	ASP	3.9
1	B	261	ALA	3.8
1	B	259	GLY	3.7
1	A	128	LYS	3.7
1	B	151	PRO	3.6
1	B	257	LYS	3.5
1	A	37	TYR	3.5
1	B	242	ALA	3.5
1	B	246	GLY	3.5
1	A	127	GLU	3.4
1	B	466	ILE	3.3
1	B	255	PRO	3.0
1	B	37	TYR	2.9
1	B	241	GLY	2.9
1	B	152	GLY	2.8
1	A	130	GLU	2.8
1	B	34	ARG	2.7
1	A	35	SER	2.7
1	B	240	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	10	SER	2.6
1	B	155	VAL	2.6
1	B	260	GLU	2.5
1	B	248	GLY	2.5
1	A	42	LEU	2.5
1	B	127	GLU	2.5
1	B	212	GLY	2.5
1	B	134	LEU	2.5
1	B	123	SER	2.4
1	B	120	GLY	2.3
1	B	238	LYS	2.3
1	A	119	GLN	2.3
1	B	154	GLU	2.2
1	B	119	GLN	2.1
1	B	15	TYR	2.1
1	B	36	THR	2.1
1	A	126	ASN	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

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. 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	B	501	5/5	0.17	2.35	34,38,42,43	0
2	FAD	B	500	53/53	0.19	1.30	28,42,50,57	0
2	FAD	A	500	53/53	0.18	0.93	25,37,46,50	0
4	EDO	B	502	4/4	0.16	0.51	32,33,35,46	0
3	SO4	A	501	5/5	0.10	-1.35	60,64,65,67	0

6.5 Other polymers ⓘ

There are no such residues in this entry.