



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2017 – 06:35 AM EDT

PDB ID : 3DGH
Title : Crystal Structure of Drosophila Thioredoxin Reductase, C-terminal 8-residue truncation
Authors : Eckenroth, B.E.; Hondal, R.J.; Everse, S.J.
Deposited on : unknown
Resolution : 1.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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 : rb-20030345
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) : rb-20030345

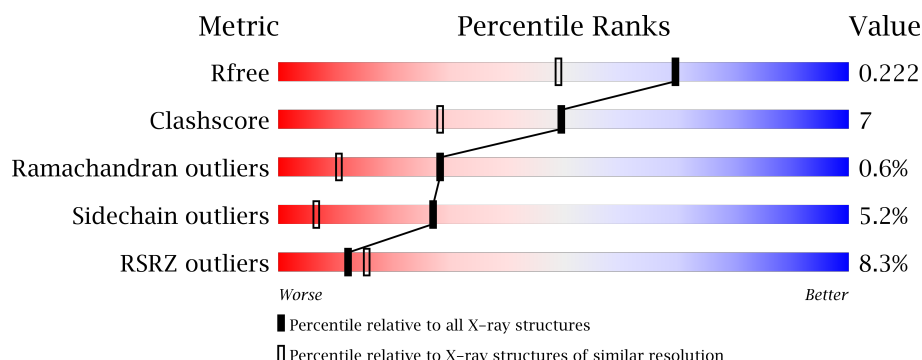
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	2694 (1.76-1.72)
Clashscore	112137	2854 (1.76-1.72)
Ramachandran outliers	110173	2824 (1.76-1.72)
Sidechain outliers	110143	2824 (1.76-1.72)
RSRZ outliers	101464	2705 (1.76-1.72)

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. 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	483	<div> <div>5%</div> <div>87%</div> <div>9%</div> <div>..</div> </div>
1	B	483	<div> <div>12%</div> <div>82%</div> <div>14%</div> <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
3	SO4	A	503	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8058 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 Thioredoxin reductase 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	477	Total	C	N	O	S	0	0	0
			3657	2333	621	691	12			
1	B	473	Total	C	N	O	S	0	0	0
			3631	2318	617	684	12			

There are 14 discrepancies between the modelled and reference sequences:

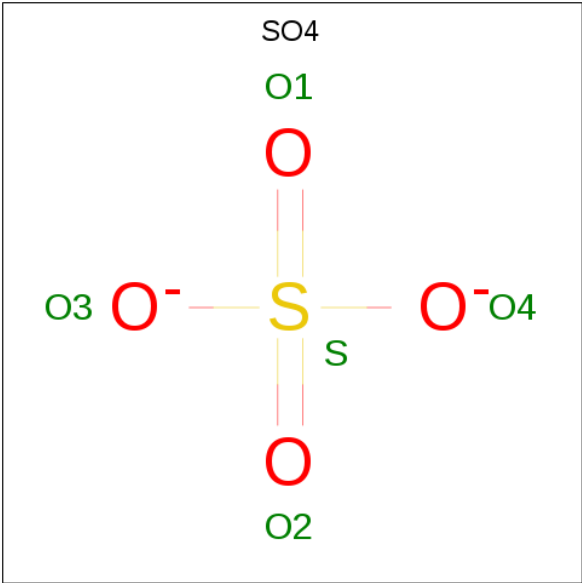
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP P91938
A	2	ALA	-	expression tag	UNP P91938
A	3	PRO	-	expression tag	UNP P91938
A	4	VAL	-	expression tag	UNP P91938
A	5	GLN	-	expression tag	UNP P91938
A	90	ASP	GLU	engineered	UNP P91938
A	274	SER	ALA	engineered	UNP P91938
B	1	MET	-	expression tag	UNP P91938
B	2	ALA	-	expression tag	UNP P91938
B	3	PRO	-	expression tag	UNP P91938
B	4	VAL	-	expression tag	UNP P91938
B	5	GLN	-	expression tag	UNP P91938
B	90	ASP	GLU	engineered	UNP P91938
B	274	SER	ALA	engineered	UNP P91938

- 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 53	C 27	N 9	O 15	P 2	0	0
2	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- 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	A	1	Total	O S	0	0
			5	4 1		

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

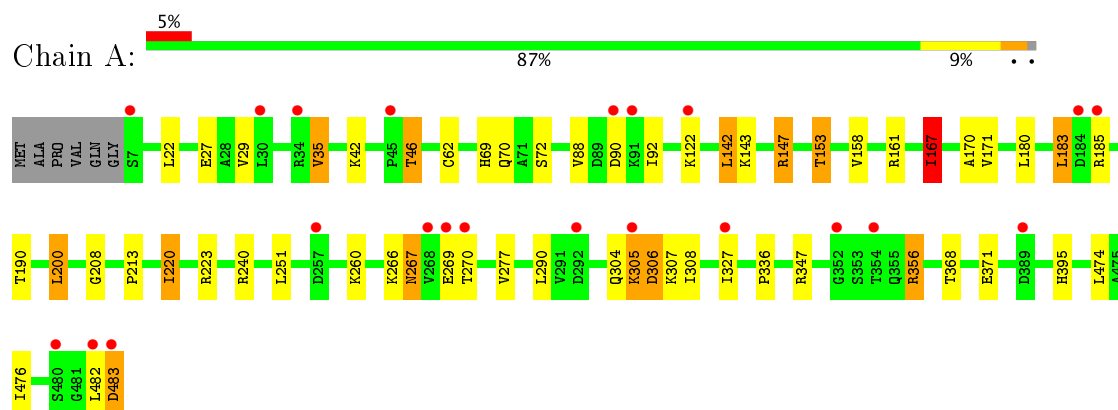
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	393	Total	O	0	0
			393	393		
4	B	251	Total	O	0	0
			251	251		

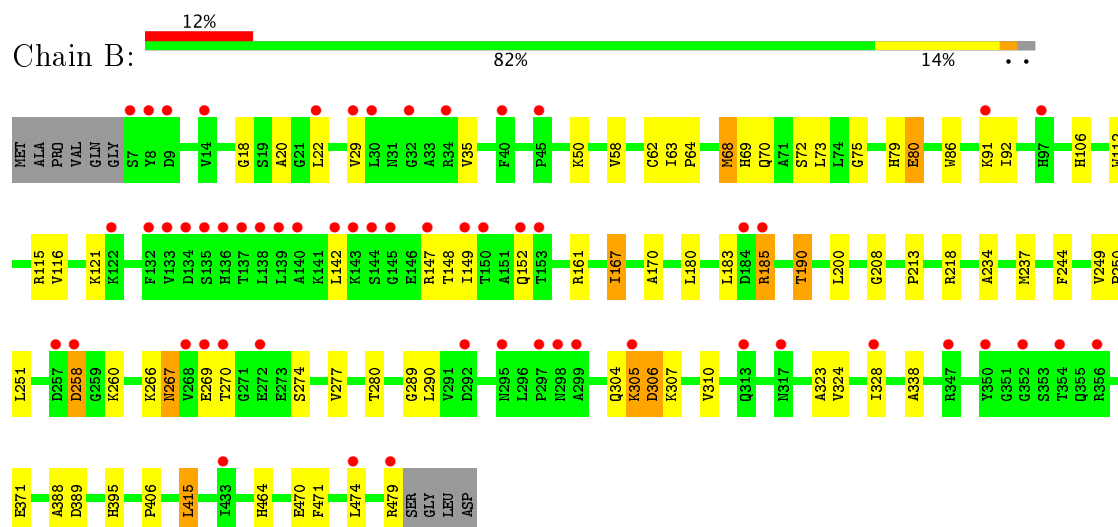
3 Residue-property plots [i](#)

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: Thioredoxin reductase 1, mitochondrial



- Molecule 1: Thioredoxin reductase 1, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	135.75Å 135.75Å 132.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.77 – 1.75 29.77 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.77-1.75) 99.9 (29.77-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 1.75Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.200 , 0.224 0.197 , 0.222	Depositor DCC
R_{free} test set	12552 reflections (10.03%)	DCC
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.009 for -h,-l,-k 0.004 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8058	wwPDB-VP
Average B, all atoms (Å ²)	6.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, FAD

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	0.66	0/3730	0.73	6/5063 (0.1%)
1	B	0.63	1/3704 (0.0%)	0.70	1/5028 (0.0%)
All	All	0.65	1/7434 (0.0%)	0.71	7/10091 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	80	GLU	CD-OE1	6.23	1.32	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	68	MET	CG-SD-CE	-8.03	87.35	100.20
1	A	147	ARG	NE-CZ-NH2	-6.29	117.15	120.30
1	A	142	LEU	CA-CB-CG	6.03	129.17	115.30
1	A	200	LEU	CB-CG-CD1	5.71	120.70	111.00
1	A	167	ILE	CG1-CB-CG2	5.50	123.50	111.40
1	A	90	ASP	N-CA-CB	-5.47	100.75	110.60
1	A	240	ARG	NE-CZ-NH2	-5.15	117.72	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3657	0	3678	42	0
1	B	3631	0	3655	60	0
2	A	53	0	31	1	0
2	B	53	0	31	3	0
3	A	10	0	0	2	0
3	B	10	0	0	0	0
4	A	393	0	0	3	0
4	B	251	0	0	1	1
All	All	8058	0	7395	99	1

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 7.

All (99) 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:A:88:VAL:HG12	1:B:92:ILE:HG21	1.34	1.08
1:B:415:LEU:HD11	1:B:474:LEU:HD11	1.30	1.06
1:A:158:VAL:HG22	1:A:327:ILE:HD12	1.43	0.99
1:B:415:LEU:CD1	1:B:474:LEU:HD11	2.02	0.89
1:A:158:VAL:CG2	1:A:327:ILE:HD12	2.03	0.88
1:B:324:VAL:HA	1:B:328:ILE:HD11	1.65	0.79
1:B:470:GLU:O	1:B:474:LEU:HD22	1.84	0.78
1:A:69:HIS:HE1	1:A:371:GLU:OE1	1.67	0.78
1:B:69:HIS:HE1	1:B:371:GLU:OE1	1.66	0.78
1:A:220:ILE:HD11	1:A:223:ARG:HB2	1.65	0.77
1:B:305:LYS:O	1:B:307:LYS:N	2.19	0.75
1:A:171:VAL:HG22	3:A:503:SO4:O4	1.87	0.74
1:A:171:VAL:HG22	3:A:503:SO4:S	2.27	0.74
1:A:356:ARG:CG	1:A:356:ARG:HH11	2.03	0.71
1:A:260:LYS:HB3	1:A:277:VAL:CG2	2.20	0.70
1:B:415:LEU:HD12	1:B:415:LEU:C	2.13	0.69
1:B:58:VAL:HA	1:B:63:ILE:HD13	1.76	0.67
1:B:304:GLN:O	1:B:305:LYS:O	2.13	0.66
1:A:482:LEU:O	1:A:483:ASP:OD1	2.13	0.66
1:A:88:VAL:HG12	1:B:92:ILE:CG2	2.20	0.65
1:B:324:VAL:HA	1:B:328:ILE:CD1	2.26	0.65
1:B:323:ALA:O	1:B:328:ILE:HD11	1.97	0.64
1:B:415:LEU:HD12	1:B:415:LEU:O	1.98	0.64
1:B:69:HIS:HD2	4:B:1154:HOH:O	1.80	0.64
1:A:27:GLU:OE2	1:A:347:ARG:HD2	1.97	0.64
1:A:476:ILE:CD1	1:A:482:LEU:HG	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:THR:HG23	4:A:912:HOH:O	1.99	0.62
1:B:470:GLU:O	1:B:474:LEU:CD2	2.47	0.62
1:B:471:PHE:HA	1:B:474:LEU:CD2	2.31	0.61
1:B:18:GLY:HA3	2:B:500:FAD:H51A	1.84	0.60
1:A:69:HIS:HD2	4:A:1230:HOH:O	1.84	0.60
1:B:415:LEU:HD11	1:B:474:LEU:CD1	2.20	0.60
1:B:75:GLY:O	1:B:79:HIS:HD2	1.86	0.59
1:B:237:MET:HE3	1:B:244:PHE:HZ	1.67	0.59
1:B:147:ARG:HG2	1:B:149:ILE:HG23	1.85	0.59
1:B:200:LEU:HD13	1:B:237:MET:HE1	1.84	0.58
1:B:260:LYS:HB3	1:B:277:VAL:CG1	2.34	0.58
1:B:72:SER:OG	1:B:208:GLY:HA3	2.03	0.58
1:B:18:GLY:CA	2:B:500:FAD:H51A	2.35	0.57
1:B:258:ASP:HB3	1:B:260:LYS:H	1.70	0.57
1:B:29:VAL:HG22	1:B:35:VAL:HG13	1.88	0.56
1:B:310:VAL:HG22	1:B:328:ILE:CG2	2.36	0.56
1:B:234:ALA:HA	1:B:237:MET:HE2	1.89	0.55
1:A:267:ASN:ND2	1:A:269:GLU:H	2.04	0.55
1:B:471:PHE:HA	1:B:474:LEU:HD21	1.88	0.55
1:A:167:ILE:HG22	1:A:170:ALA:HB2	1.90	0.54
1:B:415:LEU:CD1	1:B:415:LEU:C	2.76	0.54
1:B:190:THR:HB	1:B:280:THR:HB	1.91	0.53
1:B:161:ARG:HG2	1:B:290:LEU:HD11	1.91	0.53
1:A:308:ILE:HD12	1:A:327:ILE:HD13	1.91	0.53
1:B:310:VAL:HG22	1:B:328:ILE:HG21	1.91	0.53
1:A:356:ARG:HG3	1:A:356:ARG:HH11	1.72	0.52
1:B:260:LYS:HB3	1:B:277:VAL:HG13	1.92	0.52
1:B:185:ARG:H	1:B:185:ARG:HD3	1.75	0.52
1:A:476:ILE:HD13	1:A:482:LEU:HG	1.91	0.51
1:B:249:VAL:HG13	1:B:250:PRO:HD2	1.92	0.51
1:B:112:TRP:O	1:B:116:VAL:HG23	2.10	0.51
1:B:64:PRO:O	1:B:68:MET:HG2	2.11	0.51
1:B:167:ILE:HG22	1:B:170:ALA:HB2	1.92	0.50
1:B:471:PHE:O	1:B:474:LEU:HD23	2.11	0.50
1:A:356:ARG:HG2	1:A:356:ARG:HH11	1.77	0.50
1:A:29:VAL:HG22	1:A:35:VAL:HG13	1.94	0.50
1:B:305:LYS:C	1:B:307:LYS:H	2.15	0.49
1:A:183:LEU:HD22	1:A:185:ARG:H	1.77	0.49
1:B:388:ALA:O	1:B:479:ARG:NH2	2.46	0.49
1:B:18:GLY:N	2:B:500:FAD:H51A	2.27	0.48
1:A:183:LEU:HD22	1:A:185:ARG:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:VAL:CG2	1:A:35:VAL:HG13	2.43	0.48
1:B:267:ASN:ND2	1:B:270:THR:H	2.11	0.48
1:A:72:SER:OG	1:A:208:GLY:HA3	2.14	0.48
1:A:260:LYS:CB	1:A:277:VAL:CG2	2.90	0.48
1:A:305:LYS:O	1:A:307:LYS:N	2.46	0.48
1:A:251:LEU:HD11	1:A:266:LYS:HB2	1.95	0.47
1:A:161:ARG:HG2	1:A:290:LEU:HD11	1.97	0.47
1:B:29:VAL:CG2	1:B:35:VAL:HG13	2.45	0.47
1:A:42:LYS:NZ	1:A:42:LYS:HB2	2.31	0.46
1:A:190:THR:O	1:A:213:PRO:HA	2.17	0.45
1:B:29:VAL:HG11	1:B:121:LYS:O	2.16	0.45
1:A:336:PRO:HG3	1:B:464:HIS:HB2	1.99	0.45
1:B:251:LEU:HD11	1:B:266:LYS:HB3	1.99	0.45
1:B:70:GLN:HA	1:B:70:GLN:HE21	1.82	0.45
4:A:930:HOH:O	1:B:106:HIS:HE1	2.00	0.44
1:B:20:ALA:HB2	1:B:338:ALA:HB1	1.99	0.44
1:A:356:ARG:CG	1:A:356:ARG:NH1	2.72	0.44
1:A:305:LYS:O	1:A:306:ASP:OD2	2.36	0.44
1:A:267:ASN:HD22	1:A:269:GLU:H	1.65	0.44
1:B:80:GLU:HG3	1:B:406:PRO:HG3	2.00	0.44
1:B:267:ASN:HD22	1:B:270:THR:H	1.66	0.42
1:A:260:LYS:HB3	1:A:277:VAL:HG23	2.00	0.42
1:B:267:ASN:ND2	1:B:269:GLU:H	2.17	0.42
1:A:267:ASN:ND2	1:A:270:THR:H	2.18	0.42
1:B:251:LEU:HD11	1:B:266:LYS:CB	2.49	0.42
1:A:70:GLN:HE21	1:A:70:GLN:HA	1.85	0.41
1:B:190:THR:O	1:B:213:PRO:HA	2.19	0.41
1:A:356:ARG:HG3	1:A:356:ARG:NH1	2.36	0.41
2:A:500:FAD:H1'1	2:A:500:FAD:H9	1.90	0.40
1:A:92:ILE:HD12	1:B:86:TRP:CE3	2.57	0.40
1:A:304:GLN:O	1:A:305:LYS:O	2.40	0.40
1:B:289:GLY:HA3	1:B:306:ASP:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:961:HOH:O	4:B:961:HOH:O[7_555]	2.14	0.06

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	475/483 (98%)	463 (98%)	9 (2%)	3 (1%)	28	10
1	B	471/483 (98%)	455 (97%)	13 (3%)	3 (1%)	28	10
All	All	946/966 (98%)	918 (97%)	22 (2%)	6 (1%)	28	10

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	305	LYS
1	A	306	ASP
1	B	305	LYS
1	B	306	ASP
1	A	46	THR
1	B	148	THR

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	388/392 (99%)	368 (95%)	20 (5%)	27	7
1	B	385/392 (98%)	365 (95%)	20 (5%)	27	7
All	All	773/784 (99%)	733 (95%)	40 (5%)	27	7

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	35	VAL
1	A	46	THR
1	A	62	CYS
1	A	122	LYS
1	A	142	LEU
1	A	143	LYS
1	A	147	ARG
1	A	153	THR
1	A	167	ILE
1	A	180	LEU
1	A	183	LEU
1	A	200	LEU
1	A	220	ILE
1	A	267	ASN
1	A	356	ARG
1	A	368	THR
1	A	395	HIS
1	A	474	LEU
1	A	483	ASP
1	B	22	LEU
1	B	50	LYS
1	B	62	CYS
1	B	73	LEU
1	B	91	LYS
1	B	115	ARG
1	B	142	LEU
1	B	152	GLN
1	B	167	ILE
1	B	180	LEU
1	B	183	LEU
1	B	185	ARG
1	B	190	THR
1	B	218	ARG
1	B	258	ASP
1	B	267	ASN
1	B	274	SER
1	B	389	ASP
1	B	395	HIS
1	B	415	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	HIS
1	A	70	GLN
1	A	136	HIS
1	A	227	GLN
1	A	267	ASN
1	A	298	ASN
1	A	355	GLN
1	A	455	ASN
1	B	69	HIS
1	B	70	GLN
1	B	79	HIS
1	B	106	HIS
1	B	136	HIS
1	B	267	ASN
1	B	298	ASN
1	B	355	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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	1.37	6 (11%)	54,89,89	1.98	5 (9%)
3	SO4	A	501	-	4,4,4	0.16	0	6,6,6	0.58	0
3	SO4	A	503	-	4,4,4	0.47	0	6,6,6	0.45	0
2	FAD	B	500	-	51,58,58	1.38	7 (13%)	54,89,89	2.47	11 (20%)
3	SO4	B	502	-	4,4,4	0.15	0	6,6,6	0.27	0
3	SO4	B	504	-	4,4,4	0.15	0	6,6,6	0.30	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
3	SO4	A	503	-	-	0/0/0/0	0/0/0/0
2	FAD	B	500	-	-	0/28/50/50	0/6/6/6
3	SO4	B	502	-	-	0/0/0/0	0/0/0/0
3	SO4	B	504	-	-	0/0/0/0	0/0/0/0

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	FAD	C5X-N5	2.28	1.38	1.35
2	A	500	FAD	C5X-N5	2.47	1.39	1.35
2	B	500	FAD	C2A-N1A	2.78	1.39	1.33
2	A	500	FAD	C4-N3	3.08	1.38	1.33
2	B	500	FAD	C4-N3	3.09	1.38	1.33
2	B	500	FAD	C1'-N10	3.23	1.51	1.48
2	A	500	FAD	C4X-N5	3.34	1.38	1.33
2	A	500	FAD	C10-N1	3.44	1.38	1.33
2	B	500	FAD	C2A-N3A	3.56	1.38	1.32
2	A	500	FAD	C1'-N10	3.58	1.52	1.48
2	B	500	FAD	C10-N1	3.62	1.38	1.33
2	B	500	FAD	C4X-N5	3.80	1.38	1.33
2	A	500	FAD	C2A-N3A	4.10	1.39	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	N3A-C2A-N1A	-9.69	120.42	128.86
2	B	500	FAD	N3A-C2A-N1A	-9.39	120.68	128.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	FAD	C4B-O4B-C1B	-5.71	103.70	109.77
2	A	500	FAD	C4X-C4-N3	-3.66	118.27	123.48
2	B	500	FAD	C4X-C4-N3	-3.47	118.55	123.48
2	B	500	FAD	C5B-C4B-C3B	-2.60	105.37	115.29
2	B	500	FAD	O3B-C3B-C4B	-2.26	104.47	111.09
2	B	500	FAD	C5X-C9A-N10	2.77	119.72	117.66
2	A	500	FAD	C5X-C9A-N10	2.92	119.83	117.66
2	B	500	FAD	C4X-N5-C5X	3.48	120.43	116.76
2	A	500	FAD	C4X-N5-C5X	4.05	121.04	116.76
2	B	500	FAD	O5B-C5B-C4B	4.18	123.83	109.00
2	B	500	FAD	C1'-N10-C9A	4.74	122.69	118.35
2	A	500	FAD	C4-N3-C2	6.67	121.00	115.16
2	B	500	FAD	C4-N3-C2	6.80	121.10	115.16
2	B	500	FAD	O4B-C4B-C5B	6.90	132.70	109.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	FAD	1	0
3	A	503	SO4	2	0
2	B	500	FAD	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	477/483 (98%)	0.41	22 (4%) 33 39	2, 4, 12, 25	0
1	B	473/483 (97%)	0.75	57 (12%) 5 6	2, 7, 18, 28	0
All	All	950/966 (98%)	0.58	79 (8%) 12 15	2, 5, 16, 28	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	483	ASP	9.1
1	B	144	SER	8.9
1	A	482	LEU	6.2
1	A	7	SER	5.5
1	A	480	SER	5.0
1	B	7	SER	4.8
1	B	479	ARG	4.8
1	B	145	GLY	4.5
1	B	122	LYS	4.5
1	B	135	SER	4.5
1	B	143	LYS	4.3
1	B	34	ARG	4.2
1	B	268	VAL	4.0
1	B	140	ALA	4.0
1	B	152	GLN	4.0
1	B	32	GLY	4.0
1	B	132	PHE	3.8
1	A	305	LYS	3.7
1	B	133	VAL	3.7
1	B	30	LEU	3.6
1	B	257	ASP	3.6
1	B	134	ASP	3.5
1	B	270	THR	3.5
1	A	257	ASP	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	268	VAL	3.4
1	B	185	ARG	3.2
1	B	29	VAL	3.2
1	B	150	THR	3.2
1	B	354	THR	3.2
1	A	30	LEU	3.1
1	B	328	ILE	3.1
1	B	299	ALA	3.1
1	B	295	ASN	3.0
1	B	153	THR	3.0
1	B	147	ARG	3.0
1	B	8	TYR	3.0
1	B	142	LEU	3.0
1	A	269	GLU	2.8
1	A	270	THR	2.8
1	B	137	THR	2.8
1	B	269	GLU	2.8
1	A	185	ARG	2.7
1	B	149	ILE	2.7
1	B	258	ASP	2.7
1	A	352	GLY	2.7
1	B	305	LYS	2.6
1	B	272	GLU	2.5
1	A	184	ASP	2.5
1	A	122	LYS	2.5
1	B	474	LEU	2.5
1	B	139	LEU	2.5
1	A	327	ILE	2.4
1	B	292	ASP	2.4
1	B	313	GLN	2.3
1	B	356	ARG	2.3
1	A	354	THR	2.3
1	A	90	ASP	2.3
1	B	97	HIS	2.3
1	B	350	TYR	2.3
1	B	14	VAL	2.3
1	B	22	LEU	2.3
1	B	317	ASN	2.2
1	B	40	PHE	2.2
1	B	297	PRO	2.2
1	B	347	ARG	2.2
1	A	34	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	91	LYS	2.1
1	B	45	PRO	2.1
1	B	352	GLY	2.1
1	B	91	LYS	2.1
1	B	298	ASN	2.1
1	A	292	ASP	2.1
1	A	389	ASP	2.1
1	B	184	ASP	2.1
1	B	433	ILE	2.1
1	A	45	PRO	2.0
1	B	138	LEU	2.0
1	B	9	ASP	2.0
1	B	136	HIS	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, 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	A	501	5/5	0.97	0.12	0.92	20,20,21,23	0
3	SO4	A	503	5/5	0.98	0.10	-0.03	6,8,9,11	0
2	FAD	B	500	53/53	0.94	0.12	-0.07	2,3,6,12	0
2	FAD	A	500	53/53	0.97	0.12	-0.15	2,2,2,2	0
3	SO4	B	504	5/5	0.99	0.06	-1.66	15,16,17,18	0
3	SO4	B	502	5/5	0.98	0.07	-1.71	6,9,9,10	0

6.5 Other polymers

There are no such residues in this entry.