



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:41 PM GMT

PDB ID : 1BZL  
Title : CRYSTAL STRUCTURE OF TRYPANOSOMA CRUZI TRYPANOTHIONE REDUCTASE IN COMPLEX WITH TRYPANOTHIONE, AND THE STRUCTURE-BASED DISCOVERY OF NEW NATURAL PRODUCT INHIBITORS  
Authors : Bond, C.S.; Zhang, Y.; Berriman, M.; Cunningham, M.; Fairlamb, A.H.; Hunter, W.N.  
Deposited on : 1998-11-02  
Resolution : 2.40 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.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

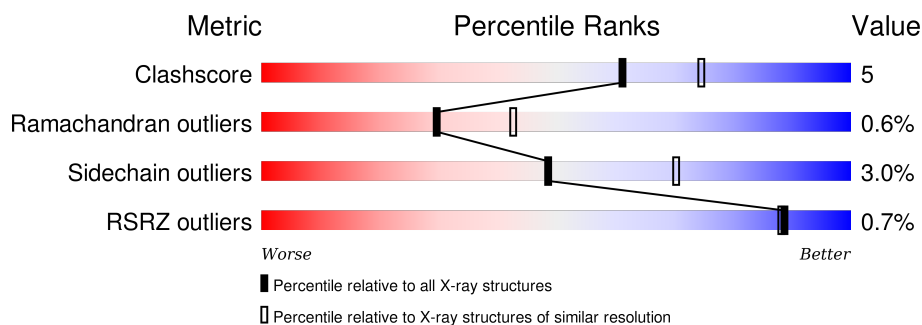
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	486	<div> <div></div> <div>85%13%.</div> </div>
1	B	486	<div> <div></div> <div>82%15%..</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	GCG	A	603	-	-	-	X
3	GCG	A	604	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8070 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 TRYPTANOTHIONE REDUCTASE (OXIDIZED FORM).

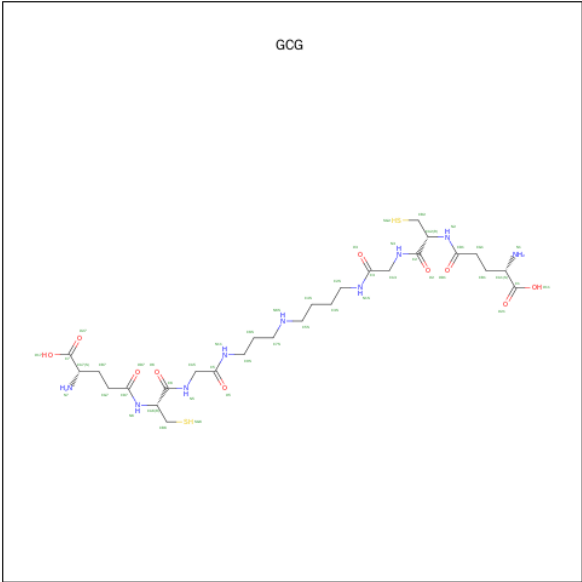
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	486	Total	C	N	O	S	0	0	0
			3741	2378	636	705	22			
1	B	483	Total	C	N	O	S	0	0	0
			3718	2364	632	701	21			

- 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 BIS(GAMMA-GLUTAMYL-CYSTEINYL-GLYCINYL)SPERMIDINE (three-letter code: GCG) (formula:  $C_{27}H_{49}N_9O_{10}S_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			48	27	9	10	2		
3	A	1	Total	C	N	O	S	0	0
			48	27	9	10	2		

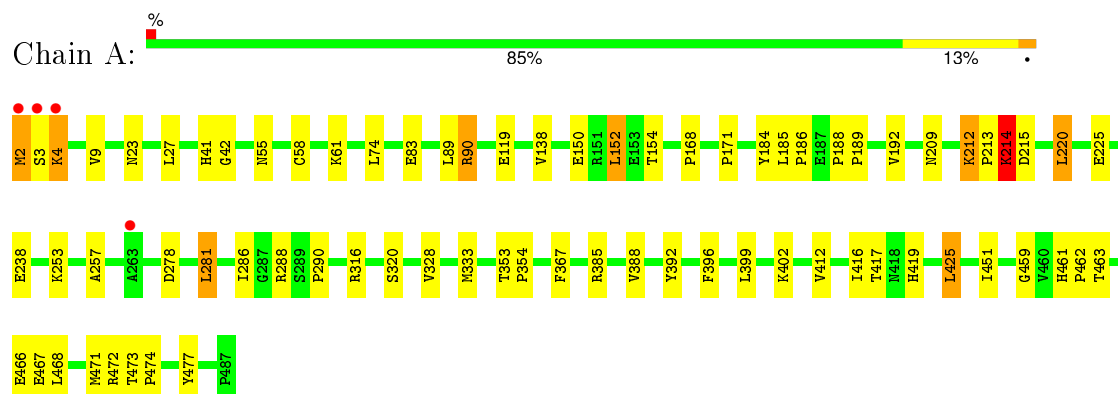
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	220	Total	O	0	0
			220	220		
4	B	189	Total	O	0	0
			189	189		

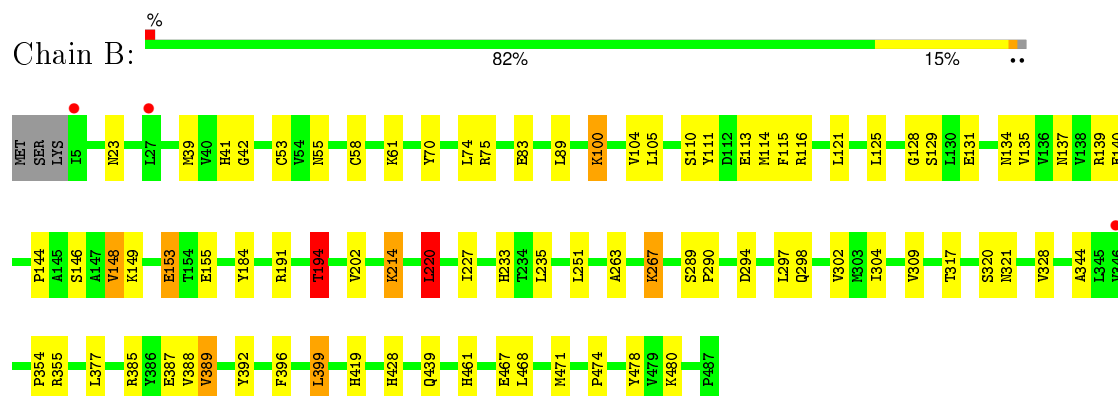
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: TRYPTANOTHIONE REDUCTASE (OXIDIZED FORM)



- Molecule 1: TRYPTANOTHIONE REDUCTASE (OXIDIZED FORM)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.33Å 93.33Å 157.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.40 19.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	87.4 (8.00-2.40) 94.8 (19.98-2.40)	Depositor EDS
$R_{merge}$	0.75	Depositor
$R_{sym}$	0.75	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.25 (at 2.41Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.209 , (Not available) 0.199 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	11.8	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.20 , 97.0	EDS
Estimated twinning fraction	0.209 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.31$ , $\langle L^2 \rangle = 0.15$	Xtriage
Outliers	0 of 49620 reflections	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	8070	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: FAD, GCG

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	0.61	0/3818	0.87	9/5170 (0.2%)
1	B	0.59	0/3795	0.89	15/5141 (0.3%)
All	All	0.60	0/7613	0.88	24/10311 (0.2%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	191	ARG	CB-CG-CD	-10.43	84.48	111.60
1	A	281	LEU	CA-CB-CG	9.35	136.80	115.30
1	B	220	LEU	CA-CB-CG	8.81	135.56	115.30
1	A	220	LEU	CA-CB-CG	8.45	134.74	115.30
1	A	152	LEU	CA-CB-CG	7.28	132.04	115.30
1	B	23	ASN	CA-CB-CG	6.43	127.55	113.40
1	B	399	LEU	CA-CB-CG	6.23	129.63	115.30
1	A	316	ARG	CB-CG-CD	-6.09	95.77	111.60
1	B	194	THR	N-CA-CB	-5.83	99.22	110.30
1	B	105	LEU	CA-CB-CG	5.83	128.70	115.30
1	B	214	LYS	CB-CG-CD	-5.77	96.59	111.60
1	B	75	ARG	CB-CG-CD	-5.66	96.89	111.60
1	B	148	VAL	CG1-CB-CG2	-5.62	101.90	110.90
1	B	194	THR	CB-CA-C	5.48	126.39	111.60
1	A	385	ARG	CB-CG-CD	-5.37	97.63	111.60
1	B	389	VAL	CG1-CB-CG2	-5.34	102.35	110.90
1	B	191	ARG	CG-CD-NE	5.29	122.90	111.80
1	B	354	PRO	N-CA-C	-5.28	98.38	112.10
1	A	9	VAL	CG1-CB-CG2	-5.27	102.46	110.90
1	A	3	SER	C-N-CA	5.20	134.71	121.70
1	A	425	LEU	CA-CB-CG	5.16	127.17	115.30
1	A	472	ARG	CB-CG-CD	5.06	124.76	111.60

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	385	ARG	CB-CG-CD	-5.03	98.52	111.60
1	B	202	VAL	CG1-CB-CG2	-5.00	102.90	110.90

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	3741	0	3752	37	0
1	B	3718	0	3725	46	0
2	A	53	0	31	0	0
2	B	53	0	31	1	0
3	A	96	0	90	8	0
4	A	220	0	0	3	0
4	B	189	0	0	8	0
All	All	8070	0	7629	84	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 5.

All (84) 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)
3:A:603:GCG:HB22	3:A:603:GCG:H5	1.53	0.73
1:B:233:HIS:HB2	4:B:631:HOH:O	1.93	0.68
1:A:225:GLU:HG2	1:A:253:LYS:HE3	1.77	0.66
3:A:603:GCG:SG6	1:B:461:HIS:CE1	2.91	0.63
1:B:317:THR:HB	4:B:686:HOH:O	1.98	0.62
3:A:604:GCG:HG11	4:A:766:HOH:O	2.00	0.61
1:B:113:GLU:HG3	1:B:116:ARG:HH21	1.66	0.61
1:A:214:LYS:HD2	1:A:215:ASP:H	1.65	0.61
1:B:290:PRO:HB3	1:B:328:VAL:HA	1.81	0.60
1:B:155:GLU:O	1:B:321:ASN:HB2	2.02	0.59
1:B:387:GLU:HG3	1:B:480:LYS:HG2	1.85	0.57

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:468:LEU:HA	1:A:471:MET:HE3	1.86	0.57
1:B:113:GLU:HG3	1:B:116:ARG:NH2	2.20	0.56
1:B:139:ARG:NH1	1:B:148:VAL:HG23	2.21	0.56
1:A:90:ARG:HG3	1:B:83:GLU:HG3	1.89	0.53
1:B:70:TYR:O	1:B:74:LEU:HG	2.08	0.53
1:B:131:GLU:HB3	1:B:135:VAL:HB	1.91	0.53
3:A:604:GCG:HB22	3:A:604:GCG:HA6	1.91	0.52
1:B:294:ASP:HB2	4:B:764:HOH:O	2.10	0.52
1:A:4:LYS:HG2	4:A:691:HOH:O	2.10	0.52
1:B:396:PHE:CZ	1:B:467:GLU:HG3	2.46	0.51
1:A:41:HIS:HB3	1:A:55:ASN:OD1	2.11	0.51
1:B:131:GLU:HB2	1:B:137:ASN:OD1	2.12	0.50
1:B:53:CYS:HB3	4:B:625:HOH:O	2.10	0.50
1:A:367:PHE:HE2	4:B:690:HOH:O	1.94	0.50
1:B:297:LEU:HB3	1:B:302:VAL:HG22	1.93	0.50
1:A:290:PRO:HB3	1:A:328:VAL:HA	1.94	0.49
1:B:389:VAL:HB	1:B:478:TYR:HB2	1.94	0.49
1:A:388:VAL:HG12	1:A:419:HIS:HB3	1.93	0.49
1:B:388:VAL:HG12	1:B:419:HIS:HB3	1.95	0.48
1:A:320:SER:HA	4:A:689:HOH:O	2.14	0.48
1:B:304:ILE:HD13	1:B:309:VAL:HA	1.96	0.47
1:B:115:PHE:CD1	1:B:121:LEU:HG	2.49	0.47
1:A:392:TYR:CZ	1:A:474:PRO:HG3	2.49	0.47
1:A:463:THR:O	1:A:466:GLU:HG2	2.15	0.46
1:A:2:MET:HA	1:A:2:MET:CE	2.44	0.46
1:A:396:PHE:CE1	1:A:467:GLU:HG3	2.50	0.46
1:B:392:TYR:CZ	1:B:474:PRO:HG3	2.51	0.46
1:A:4:LYS:HD3	1:A:154:THR:O	2.16	0.46
1:B:74:LEU:HD13	1:B:89:LEU:HD11	1.98	0.46
1:B:39:MET:HG3	1:B:125:LEU:HD13	1.98	0.46
1:B:41:HIS:HB3	1:B:55:ASN:OD1	2.17	0.45
1:B:468:LEU:HA	1:B:471:MET:HE3	1.98	0.45
1:B:297:LEU:HD22	1:B:302:VAL:HG21	1.97	0.45
1:B:134:ASN:O	1:B:153:GLU:HA	2.15	0.45
1:A:459:GLY:HA2	1:B:439:GLN:OE1	2.16	0.45
1:B:146:SER:HB2	4:B:776:HOH:O	2.17	0.45
1:A:288:ARG:HB2	1:A:333:MET:HE1	1.99	0.45
1:A:473:THR:HA	1:A:474:PRO:HD3	1.89	0.45
1:A:461:HIS:HA	1:A:462:PRO:HA	1.82	0.44
1:A:189:PRO:HB2	1:A:192:VAL:HB	1.99	0.44
1:B:42:GLY:HA2	1:B:184:TYR:CZ	2.52	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:LEU:HA	1:A:402:LYS:HD2	1.99	0.44
1:B:194:THR:HG22	1:B:220:LEU:HA	1.99	0.44
1:B:227:ILE:HG23	4:B:719:HOH:O	2.17	0.43
1:B:140:GLU:HG2	1:B:149:LYS:HD2	1.99	0.43
3:A:603:GCG:OD7	1:B:461:HIS:CE1	2.71	0.43
1:A:209:ASN:O	1:A:212:LYS:HD2	2.18	0.43
1:A:83:GLU:O	1:B:89:LEU:HA	2.18	0.43
1:B:399:LEU:HD12	4:B:679:HOH:O	2.18	0.43
3:A:603:GCG:SG6	1:B:461:HIS:HE1	2.40	0.43
1:A:23:ASN:O	1:A:27:LEU:HB2	2.19	0.42
1:A:214:LYS:CD	1:A:215:ASP:H	2.31	0.42
3:A:603:GCG:SG6	1:B:461:HIS:NE2	2.93	0.42
1:A:353:THR:HA	1:A:354:PRO:HD3	1.89	0.42
1:B:267:LYS:N	1:B:267:LYS:HD2	2.35	0.42
1:A:171:PRO:HD2	1:A:257:ALA:O	2.18	0.42
1:A:168:PRO:HD3	1:A:286:ILE:HD11	2.01	0.42
3:A:604:GCG:H9S1	1:B:110:SER:OG	2.20	0.42
1:B:128:GLY:HA2	1:B:137:ASN:O	2.20	0.42
1:A:238:GLU:HG3	1:A:412:VAL:HG21	2.01	0.42
1:B:344:ALA:HB1	1:B:355:ARG:O	2.20	0.41
1:A:188:PRO:HG2	1:A:213:PRO:HD3	2.03	0.41
1:A:74:LEU:HD13	1:A:89:LEU:HD11	2.02	0.41
1:B:235:LEU:HD11	1:B:428:HIS:HB3	2.02	0.41
1:B:111:TYR:O	1:B:114:MET:HB3	2.21	0.41
1:A:417:THR:HG21	1:A:451:ILE:HB	2.03	0.40
2:B:602:FAD:H9	2:B:602:FAD:H1'1	1.86	0.40
1:A:477:TYR:N	1:A:477:TYR:CD1	2.89	0.40
1:A:185:LEU:HA	1:A:186:PRO:HD3	1.87	0.40
1:A:192:VAL:HG23	1:A:281:LEU:HD12	2.03	0.40
1:B:100:LYS:O	1:B:104:VAL:HG23	2.22	0.40
1:A:42:GLY:HA2	1:A:184:TYR:CZ	2.56	0.40
1:A:138:VAL:HB	1:A:150:GLU:HB2	2.02	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	484/486 (100%)	468 (97%)	14 (3%)	2 (0%)	39	56
1	B	481/486 (99%)	460 (96%)	17 (4%)	4 (1%)	24	35
All	All	965/972 (99%)	928 (96%)	31 (3%)	6 (1%)	30	43

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	263	ALA
1	B	320	SER
1	A	4	LYS
1	A	214	LYS
1	B	214	LYS
1	B	144	PRO

### 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	407/407 (100%)	395 (97%)	12 (3%)	50	71
1	B	404/407 (99%)	392 (97%)	12 (3%)	48	70
All	All	811/814 (100%)	787 (97%)	24 (3%)	48	70

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	MET
1	A	58	CYS
1	A	61	LYS
1	A	90	ARG
1	A	119	GLU
1	A	152	LEU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	212	LYS
1	A	214	LYS
1	A	220	LEU
1	A	278	ASP
1	A	416	ILE
1	A	425	LEU
1	B	58	CYS
1	B	61	LYS
1	B	100	LYS
1	B	129	SER
1	B	153	GLU
1	B	194	THR
1	B	220	LEU
1	B	251	LEU
1	B	267	LYS
1	B	289	SER
1	B	298	GLN
1	B	377	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	GLN
1	A	106	ASN
1	A	137	ASN
1	A	156	HIS
1	A	180	ASN
1	A	209	ASN
1	A	217	GLN
1	A	306	ASN
1	A	433	ASN
1	B	106	ASN
1	B	134	ASN
1	B	180	ASN
1	B	249	GLN
1	B	298	GLN
1	B	299	ASN
1	B	433	ASN

### 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 ⓘ

4 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	601	-	48,58,58	1.40	6 (12%)	54,89,89	2.01	10 (18%)
3	GCG	A	603	-	41,47,47	0.60	0	46,58,58	0.98	4 (8%)
3	GCG	A	604	-	41,47,47	0.75	1 (2%)	46,58,58	1.20	4 (8%)
2	FAD	B	602	-	48,58,58	1.42	4 (8%)	54,89,89	2.06	7 (12%)

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	601	-	-	0/30/50/50	0/6/6/6
3	GCG	A	603	-	-	0/53/61/61	0/0/0/0
3	GCG	A	604	-	-	0/53/61/61	0/0/0/0
2	FAD	B	602	-	-	0/30/50/50	0/6/6/6

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	604	GCG	CB6-CA6	-2.80	1.49	1.53
2	A	601	FAD	C10-N10	-2.56	1.36	1.39
2	A	601	FAD	C8A-N7A	-2.22	1.30	1.34
2	B	602	FAD	C8A-N7A	-2.11	1.30	1.34
2	A	601	FAD	C2'-C3'	-2.02	1.49	1.53
2	B	602	FAD	O4B-C1B	2.07	1.43	1.41
2	A	601	FAD	C1'-N10	3.18	1.51	1.48
2	A	601	FAD	C4-N3	3.32	1.39	1.33
2	B	602	FAD	C4-C4X	5.07	1.51	1.41
2	A	601	FAD	C4-C4X	5.12	1.51	1.41
2	B	602	FAD	C1'-N10	5.18	1.53	1.48

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	602	FAD	C4X-C4-N3	-5.84	115.60	123.59
2	A	601	FAD	C4X-C4-N3	-5.77	115.69	123.59
3	A	604	GCG	C5-CA5-N5	-3.30	104.10	113.26
2	A	601	FAD	C4X-C10-N10	-2.94	118.79	120.52
2	B	602	FAD	O4'-C4'-C5'	-2.53	104.68	110.19
2	A	601	FAD	O2'-C2'-C1'	-2.40	104.04	109.94
2	A	601	FAD	O4'-C4'-C5'	-2.37	105.03	110.19
2	A	601	FAD	C4-C4X-C10	-2.16	118.56	119.94
3	A	603	GCG	C2-CA2-N2	-2.13	105.26	111.26
3	A	603	GCG	C8S-C9S-N11	2.01	118.08	112.19
2	A	601	FAD	P-O3P-PA	2.03	138.42	132.73
2	A	601	FAD	O3P-P-O5'	2.14	108.60	102.94
3	A	603	GCG	CA2-CB2-SG2	2.38	117.08	114.16
3	A	604	GCG	C4S-C3S-C2S	2.42	125.27	113.69
3	A	604	GCG	C3S-C2S-N1S	2.76	120.26	112.19
2	A	601	FAD	O3P-PA-O5B	2.80	110.37	102.94
3	A	603	GCG	CA6-N6-CD7	2.91	129.01	121.58
2	B	602	FAD	O3P-PA-O5B	3.16	111.33	102.94
3	A	604	GCG	C8S-C9S-N11	3.21	121.60	112.19
2	B	602	FAD	O4B-C1B-N9A	3.64	115.73	108.10
2	B	602	FAD	C4X-N5-C5X	4.01	121.38	116.76
2	B	602	FAD	O2A-PA-O3P	4.38	124.95	105.09
2	A	601	FAD	C4X-N5-C5X	4.43	121.86	116.76
2	B	602	FAD	C4-N3-C2	9.67	123.61	115.25
2	A	601	FAD	C4-N3-C2	9.84	123.75	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	603	GCG	5	0
3	A	604	GCG	3	0
2	B	602	FAD	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	486/486 (100%)	-0.65	4 (0%) 87 87	10, 25, 51, 87	0
1	B	483/486 (99%)	-0.43	3 (0%) 90 90	12, 33, 61, 72	0
All	All	969/972 (99%)	-0.54	7 (0%) 89 88	10, 28, 57, 87	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	MET	4.2
1	A	3	SER	3.5
1	B	346	VAL	3.3
1	A	4	LYS	3.2
1	B	27	LEU	2.4
1	B	5	ILE	2.3
1	A	263	ALA	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	GCG	A	603	48/48	0.74	0.31	7.92	53,71,78,84	0
3	GCG	A	604	48/48	0.72	0.30	5.48	62,69,75,77	0
2	FAD	A	601	53/53	0.97	0.09	-0.64	3,15,23,29	0
2	FAD	B	602	53/53	0.96	0.10	-0.78	13,21,39,44	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.