



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:05 am GMT

PDB ID : 2GIA  
Title : Crystal structures of trypanosoma brucei MRP1/MRP2  
Authors : Schumacher, M.A.; Karamooz, E.; Zikova, A.; Trantirek, L.; Lukes, J.  
Deposited on : 2006-03-28  
Resolution : 1.89 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

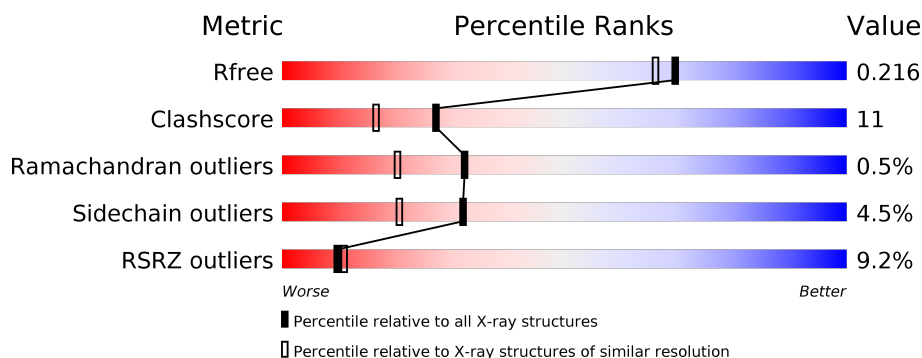
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 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	195	<div> <div>11%</div> <div> <div></div> <div>60%</div> <div>15%</div> <div>•</div> <div>21%</div> </div> </div>
1	G	195	<div> <div>13%</div> <div> <div></div> <div>60%</div> <div>17%</div> <div>•</div> <div>21%</div> </div> </div>
2	B	187	<div> <div>2%</div> <div> <div></div> <div>67%</div> <div>10%</div> <div>•</div> <div>22%</div> </div> </div>
2	D	187	<div> <div>2%</div> <div> <div></div> <div>65%</div> <div>13%</div> <div>•</div> <div>22%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACY	G	7402	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5509 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 mitochondrial RNA-binding protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1267	802	234	227	4			
1	G	155	Total	C	N	O	S	0	0	0
			1276	808	236	228	4			

- Molecule 2 is a protein called mitochondrial RNA-binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1195	754	220	217	4			
2	D	146	Total	C	N	O	S	0	0	0
			1195	754	220	217	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	43	GLU	LEU	CONFLICT	UNP P90629
D	43	GLU	LEU	CONFLICT	UNP P90629

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: C<sub>2</sub>H<sub>4</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		

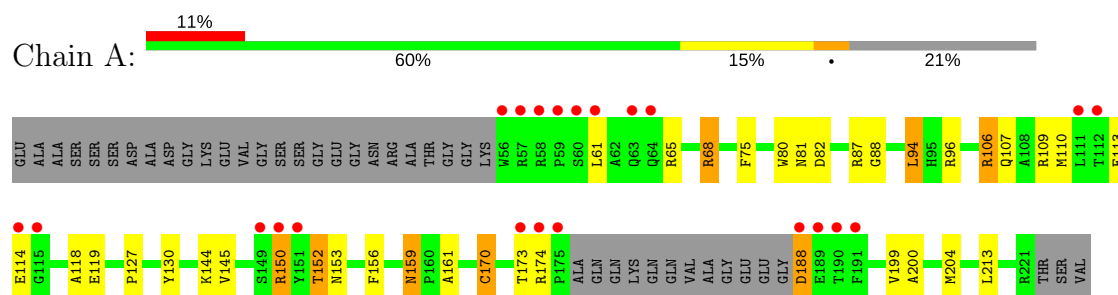
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	142	Total	O	0	0
			142	142		
4	B	167	Total	O	0	0
			167	167		
4	D	141	Total	O	0	0
			141	141		
4	G	118	Total	O	0	0
			118	118		

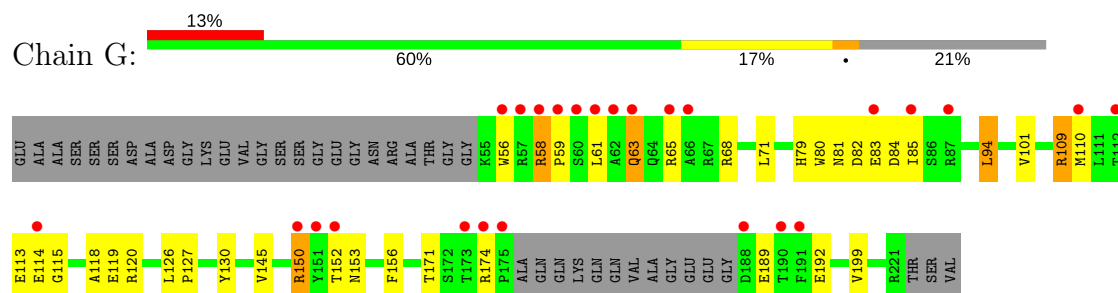
### 3 Residue-property plots [i](#)

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.

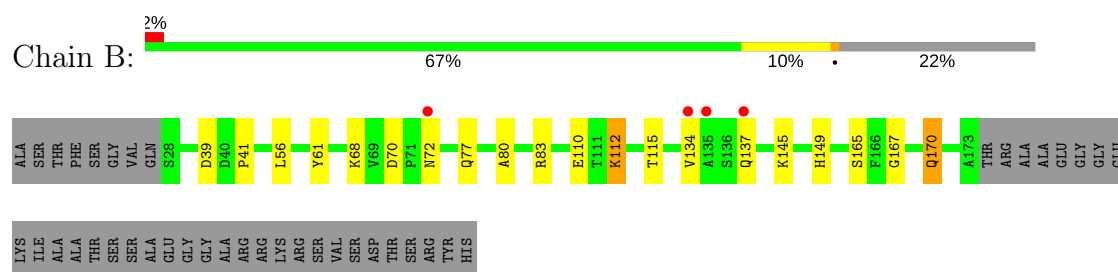
#### • Molecule 1: mitochondrial RNA-binding protein 2



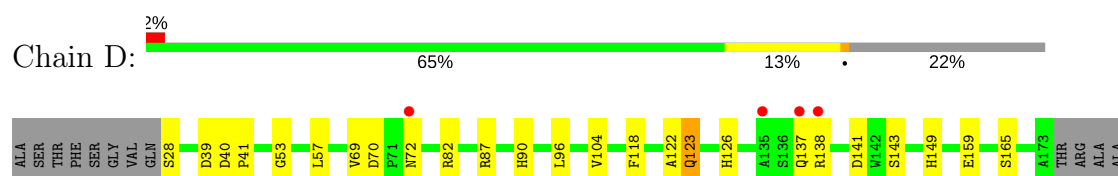
#### • Molecule 1: mitochondrial RNA-binding protein 2



#### • Molecule 2: mitochondrial RNA-binding protein 1



#### • Molecule 2: mitochondrial RNA-binding protein 1



GLU	GLY	GLY	GLY	LYS	ILE	ALA	ALA	THR	SER	SER	ALA	GLU	GLY	GLY	ALA	ARG	ARG	LYS	ARG	SER	VAL	SER	SER	ASP	THR	SER	ARG	TYR	HIS
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	60.27Å 85.70Å 86.86Å 90.00° 109.38° 90.00°	Depositor
Resolution (Å)	47.38 – 1.89 47.38 – 1.89	Depositor EDS
% Data completeness (in resolution range)	94.9 (47.38-1.89) 95.1 (47.38-1.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 1.90Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.179 , 0.216 0.179 , 0.216	Depositor DCC
$R_{free}$ test set	5784 reflections (9.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.8	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 60.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5509	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACY

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	0.72	1/1299 (0.1%)	0.80	1/1765 (0.1%)
1	G	0.63	0/1308	0.77	0/1776
2	B	0.77	0/1225	0.80	1/1654 (0.1%)
2	D	0.64	0/1225	0.77	0/1654
All	All	0.69	1/5057 (0.0%)	0.78	2/6849 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	170	CYS	CB-SG	5.60	1.91	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	LEU	CA-CB-CG	5.35	127.61	115.30
2	B	56	LEU	CA-CB-CG	5.25	127.37	115.30

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	1267	0	1238	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1276	0	1251	36	0
2	B	1195	0	1164	20	0
2	D	1195	0	1164	25	0
3	A	4	0	3	1	0
3	G	4	0	3	3	0
4	A	142	0	0	9	0
4	B	167	0	0	1	0
4	D	141	0	0	5	0
4	G	118	0	0	5	0
All	All	5509	0	4823	111	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 11.

All (111) 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:G:171:THR:HG22	1:G:192:GLU:HG2	1.41	1.02
1:G:152:THR:HG22	4:G:7508:HOH:O	1.69	0.91
1:G:79:HIS:HB3	4:G:7510:HOH:O	1.78	0.83
1:A:153:ASN:OD1	3:A:7401:ACY:H3	1.80	0.82
2:B:68:LYS:H	2:B:77:GLN:HE22	1.29	0.79
1:G:171:THR:HG23	3:G:7402:ACY:H1	1.63	0.79
1:G:56:TRP:HB2	1:G:58:ARG:HE	1.47	0.77
1:A:68:ARG:NH1	4:A:7497:HOH:O	2.18	0.76
1:A:152:THR:HG22	1:A:170:CYS:SG	2.26	0.76
2:D:149:HIS:HD2	4:D:318:HOH:O	1.67	0.76
1:A:68:ARG:NH2	2:D:165:SER:O	2.19	0.76
1:G:58:ARG:HB3	1:G:61:LEU:HB2	1.69	0.75
2:B:167:GLY:HA2	2:B:170:GLN:HE21	1.51	0.75
1:A:68:ARG:HH12	2:D:53:GLY:HA2	1.54	0.71
2:D:122:ALA:HB3	2:D:123:GLN:OE1	1.91	0.70
1:G:81:ASN:ND2	1:G:84:ASP:HB3	2.05	0.70
2:D:39:ASP:O	2:D:41:PRO:HD3	1.93	0.68
1:A:87:ARG:HD3	4:A:7491:HOH:O	1.92	0.68
1:G:150:ARG:HA	1:G:150:ARG:NE	2.09	0.68
1:G:81:ASN:HD21	1:G:84:ASP:HB3	1.58	0.66
1:A:152:THR:HG23	4:A:7518:HOH:O	1.97	0.65
1:G:56:TRP:HB2	1:G:58:ARG:NE	2.12	0.64
1:G:174:ARG:HG3	1:G:189:GLU:HB2	1.80	0.63
1:A:173:THR:HG23	4:A:7433:HOH:O	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:58:ARG:HG2	1:G:58:ARG:HH11	1.64	0.62
1:A:106:ARG:NH1	1:A:119:GLU:OE2	2.32	0.61
1:G:171:THR:CG2	3:G:7402:ACY:H1	2.30	0.61
1:A:127:PRO:HD2	1:A:130:TYR:CD2	2.35	0.61
1:G:153:ASN:HB2	4:G:7484:HOH:O	2.00	0.61
2:B:61:TYR:CD1	2:B:80:ALA:HA	2.37	0.59
1:G:127:PRO:HD2	1:G:130:TYR:HD2	1.67	0.59
2:D:57:LEU:HD22	2:D:87:ARG:NH1	2.18	0.59
1:G:171:THR:HG23	3:G:7402:ACY:CH3	2.32	0.58
1:A:152:THR:CG2	1:A:170:CYS:SG	2.92	0.57
1:A:174:ARG:O	1:A:188:ASP:HA	2.04	0.57
2:D:57:LEU:CD2	2:D:87:ARG:HG3	2.35	0.56
1:G:127:PRO:HD2	1:G:130:TYR:CD2	2.41	0.56
2:B:61:TYR:HD1	2:B:80:ALA:HA	1.71	0.56
2:D:57:LEU:HD22	2:D:87:ARG:HH11	1.70	0.56
2:B:83:ARG:CZ	2:B:134:VAL:HG11	2.36	0.55
1:G:109:ARG:HG3	1:G:113:GLU:HG3	1.88	0.55
2:D:82:ARG:NH2	4:D:338:HOH:O	2.32	0.55
1:G:145:VAL:CG1	1:G:156:PHE:HB3	2.35	0.55
1:A:75:PHE:CD1	1:A:204:MET:HE2	2.41	0.55
1:G:80:TRP:HH2	1:G:118:ALA:HB2	1.72	0.55
1:G:109:ARG:HB2	1:G:113:GLU:CD	2.26	0.55
1:A:127:PRO:HD2	1:A:130:TYR:HD2	1.71	0.55
2:B:167:GLY:CA	2:B:170:GLN:HE21	2.19	0.54
2:D:39:ASP:O	2:D:41:PRO:CD	2.55	0.54
2:B:170:GLN:CD	2:B:170:GLN:H	2.11	0.54
1:A:68:ARG:NH1	2:D:53:GLY:HA2	2.22	0.54
1:G:56:TRP:CB	1:G:58:ARG:HE	2.19	0.53
2:D:90:HIS:HE1	2:D:165:SER:OG	1.91	0.53
1:G:85:ILE:HD11	1:G:110:MET:HG3	1.89	0.53
1:A:199:VAL:HG11	2:D:104:VAL:HG11	1.91	0.53
2:B:68:LYS:H	2:B:77:GLN:NE2	2.03	0.53
4:A:7535:HOH:O	2:B:149:HIS:HD2	1.91	0.52
1:G:109:ARG:CZ	1:G:115:GLY:HA3	2.38	0.52
1:G:58:ARG:NH1	1:G:58:ARG:HG2	2.24	0.52
2:B:167:GLY:HA2	2:B:170:GLN:NE2	2.22	0.52
1:A:87:ARG:NE	4:A:7515:HOH:O	2.35	0.52
4:A:7410:HOH:O	2:D:90:HIS:HD2	1.92	0.52
1:G:174:ARG:CG	1:G:189:GLU:HB2	2.41	0.51
1:G:150:ARG:HE	1:G:150:ARG:HA	1.75	0.50
2:D:28:SER:HA	4:D:308:HOH:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:137:GLN:HA	2:B:137:GLN:HE21	1.76	0.49
1:A:159:ASN:HD22	1:A:159:ASN:C	2.14	0.49
1:G:120:ARG:HD3	4:G:7512:HOH:O	2.12	0.49
1:A:88:GLY:HA3	1:A:107:GLN:OE1	2.13	0.49
2:D:96:LEU:HD22	2:D:118:PHE:CD1	2.47	0.49
2:D:57:LEU:HD23	2:D:87:ARG:HG3	1.94	0.49
2:D:123:GLN:CD	2:D:123:GLN:H	2.17	0.48
2:D:159:GLU:OE2	4:D:298:HOH:O	2.19	0.48
2:B:137:GLN:HA	2:B:137:GLN:NE2	2.29	0.48
1:G:63:GLN:OE1	1:G:63:GLN:HA	2.12	0.48
1:A:68:ARG:HG2	4:A:7458:HOH:O	2.14	0.47
1:A:61:LEU:HD11	1:A:65:ARG:NE	2.29	0.47
2:B:145:LYS:HE2	4:B:329:HOH:O	2.14	0.47
1:A:80:TRP:HH2	1:A:118:ALA:HB2	1.80	0.47
1:A:144:LYS:HG2	1:A:145:VAL:N	2.28	0.47
2:B:68:LYS:N	2:B:77:GLN:HE22	2.05	0.47
2:B:110:GLU:HG2	2:B:115:THR:HG23	1.97	0.46
1:G:145:VAL:HG12	1:G:156:PHE:HB3	1.98	0.46
2:D:57:LEU:CD2	2:D:87:ARG:HH11	2.29	0.46
1:A:61:LEU:HD11	1:A:65:ARG:NH2	2.31	0.46
1:G:94:LEU:HG	1:G:101:VAL:HB	1.97	0.45
1:A:145:VAL:CG1	1:A:156:PHE:HB3	2.45	0.45
1:A:61:LEU:HD11	1:A:65:ARG:HE	1.82	0.45
2:D:126:HIS:NE2	2:D:143:SER:OG	2.41	0.45
1:A:150:ARG:HE	1:A:150:ARG:HA	1.82	0.44
2:D:72:ASN:ND2	4:D:305:HOH:O	2.49	0.44
2:B:165:SER:O	1:G:68:ARG:NH1	2.51	0.44
2:D:123:GLN:CD	2:D:123:GLN:N	2.71	0.44
2:B:112:LYS:HG2	2:B:112:LYS:O	2.18	0.43
2:D:69:VAL:HG23	2:D:70:ASP:N	2.34	0.43
1:A:109:ARG:HB2	1:A:113:GLU:CD	2.39	0.43
1:A:159:ASN:ND2	1:A:161:ALA:H	2.16	0.43
2:D:123:GLN:OE1	2:D:123:GLN:N	2.52	0.43
1:A:96:ARG:NH1	4:A:7459:HOH:O	2.51	0.43
1:G:58:ARG:HD2	1:G:58:ARG:N	2.34	0.43
1:A:61:LEU:HD11	1:A:65:ARG:HH21	1.85	0.42
1:G:126:LEU:HA	1:G:127:PRO:HD3	1.91	0.42
1:G:199:VAL:HG23	4:G:7446:HOH:O	2.19	0.41
2:B:137:GLN:HE21	2:B:137:GLN:CA	2.31	0.41
1:G:82:ASP:O	1:G:83:GLU:C	2.59	0.41
1:A:200:ALA:O	1:A:204:MET:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ARG:HH11	1:A:87:ARG:HG3	1.86	0.41
1:A:109:ARG:HB3	1:A:109:ARG:NH1	2.36	0.40
1:A:82:ASP:OD1	1:A:110:MET:HA	2.21	0.40
2:B:39:ASP:O	2:B:41:PRO:HD3	2.21	0.40
2:B:70:ASP:OD2	2:B:72:ASN:HB3	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	150/195 (77%)	146 (97%)	4 (3%)	0	100	100
1	G	151/195 (77%)	143 (95%)	7 (5%)	1 (1%)	25	13
2	B	144/187 (77%)	141 (98%)	3 (2%)	0	100	100
2	D	144/187 (77%)	137 (95%)	5 (4%)	2 (1%)	13	4
All	All	589/764 (77%)	567 (96%)	19 (3%)	3 (0%)	32	20

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	138	ARG
1	G	59	PRO
2	D	40	ASP

### 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	135/162 (83%)	125 (93%)	10 (7%)	16	7
1	G	136/162 (84%)	127 (93%)	9 (7%)	19	9
2	B	129/158 (82%)	127 (98%)	2 (2%)	68	65
2	D	129/158 (82%)	126 (98%)	3 (2%)	56	49
All	All	529/640 (83%)	505 (96%)	24 (4%)	32	21

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

Mol	Chain	Res	Type
1	A	68	ARG
1	A	81	ASN
1	A	94	LEU
1	A	106	ARG
1	A	114	GLU
1	A	150	ARG
1	A	152	THR
1	A	159	ASN
1	A	188	ASP
1	A	213	LEU
2	B	112	LYS
2	B	170	GLN
2	D	123	GLN
2	D	137	GLN
2	D	141	ASP
1	G	58	ARG
1	G	63	GLN
1	G	65	ARG
1	G	71	LEU
1	G	94	LEU
1	G	109	ARG
1	G	114	GLU
1	G	119	GLU
1	G	150	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN

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Mol	Chain	Res	Type
1	A	116	ASN
1	A	159	ASN
2	B	77	GLN
2	B	137	GLN
2	B	170	GLN
2	D	90	HIS
2	D	137	GLN
1	G	70	GLN
1	G	79	HIS
1	G	81	ASN
1	G	105	HIS
1	G	116	ASN

### 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](#)

2 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	$\# Z  > 2$
3	ACY	A	7401	-	1,3,3	2.37	1 (100%)	0,3,3	0.00	-
3	ACY	G	7402	-	1,3,3	0.66	0	0,3,3	0.00	-

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
3	ACY	A	7401	-	-	0/0/0/0	0/0/0/0
3	ACY	G	7402	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	7401	ACY	CH3-C	2.37	1.51	1.48

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	7401	ACY	1	0
3	G	7402	ACY	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, 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	154/195 (78%)	0.39	22 (14%) 3 3	14, 31, 71, 81	0
1	G	155/195 (79%)	0.70	25 (16%) 2 2	17, 38, 81, 88	0
2	B	146/187 (78%)	-0.05	4 (2%) 55 59	15, 25, 57, 70	0
2	D	146/187 (78%)	-0.02	4 (2%) 55 59	16, 33, 59, 78	0
All	All	601/764 (78%)	0.26	55 (9%) 10 11	14, 32, 72, 88	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	150	ARG	7.0
1	G	151	TYR	6.9
1	G	61	LEU	6.4
1	G	59	PRO	5.5
1	A	150	ARG	5.3
1	A	188	ASP	5.0
2	D	135	ALA	4.9
1	A	56	TRP	4.9
1	A	57	ARG	4.9
1	A	151	TYR	4.6
1	G	188	ASP	4.5
1	G	60	SER	4.3
1	G	191	PHE	4.2
1	G	83	GLU	4.2
1	A	61	LEU	4.1
2	D	72	ASN	3.9
1	G	174	ARG	3.9
1	G	58	ARG	3.8
1	G	175	PRO	3.8
1	G	114	GLU	3.8
1	G	190	THR	3.7

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Mol	Chain	Res	Type	RSRZ
1	G	65	ARG	3.7
1	A	59	PRO	3.6
1	G	87	ARG	3.6
2	B	135	ALA	3.5
1	A	58	ARG	3.5
1	A	114	GLU	3.5
2	B	72	ASN	3.4
1	G	112	THR	3.4
1	G	56	TRP	3.3
1	G	57	ARG	3.3
1	G	173	THR	3.3
2	D	138	ARG	3.3
1	G	85	ILE	3.2
1	G	63	GLN	3.2
2	D	137	GLN	2.8
1	A	189	GLU	2.8
2	B	137	GLN	2.8
1	A	191	PHE	2.8
1	A	63	GLN	2.7
1	A	173	THR	2.7
1	A	112	THR	2.7
1	A	60	SER	2.6
2	B	134	VAL	2.5
1	A	111	LEU	2.4
1	G	62	ALA	2.4
1	A	174	ARG	2.3
1	A	190	THR	2.3
1	G	66	ALA	2.3
1	A	149	SER	2.3
1	A	115	GLY	2.2
1	G	152	THR	2.2
1	A	175	PRO	2.2
1	A	64	GLN	2.2
1	G	110	MET	2.0

## 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 [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	ACY	A	7401	4/4	0.85	0.16	0.52	48,57,65,65	0
3	ACY	G	7402	4/4	0.97	0.17	0.28	49,58,67,72	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.