



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2017 – 09:22 AM EDT

PDB ID : 2EF6  
Title : Canavalia gladiata lectin complexed with Man1-3Man-OMe  
Authors : Moreno, F.B.M.B.; Bezerra, G.A.; Oliveira, T.M.; de Souza, E.P.; da Rocha, B.A.M.; Benevides, R.G.; Delatorre, P.; Cavada, B.S.; de Azevedo Jr., W.F.  
Deposited on : unknown  
Resolution : 2.10 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 : 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

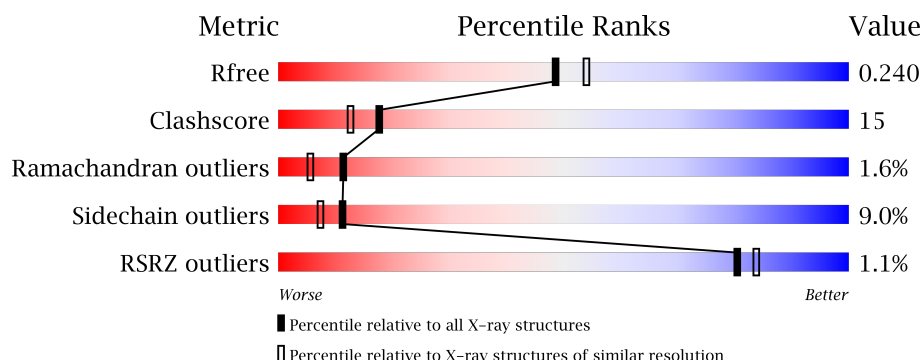
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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	237	<div> <div>75%</div> <div>21%</div> <div>...</div> </div>
1	B	237	<div> <div>68%</div> <div>23%</div> <div>5%</div> <div>..</div> </div>
1	C	237	<div> <div>60%</div> <div>34%</div> <div>..</div> </div>
1	D	237	<div> <div>62%</div> <div>31%</div> <div>5%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MDM	A	1002	X	-	-	-
2	MDM	B	1004	X	-	-	-
2	MDM	C	1003	X	-	-	-
2	MDM	D	1001	X	-	-	-

## 2 Entry composition [i](#)

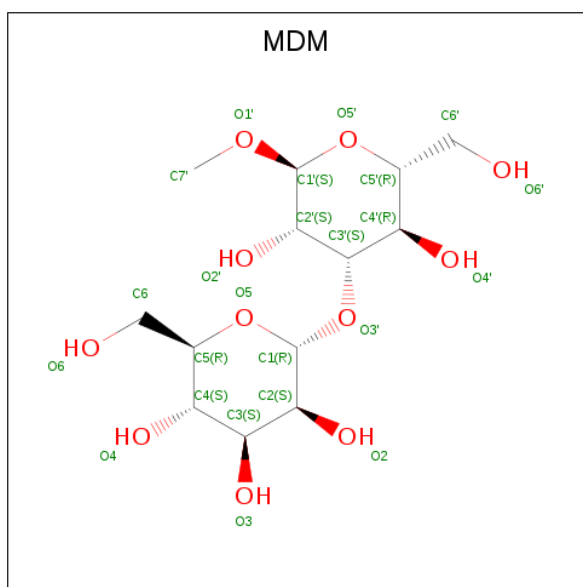
There are 5 unique types of molecules in this entry. The entry contains 7633 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Concanavalin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1779	1123	298	356	2			
1	B	234	Total	C	N	O	S	0	0	0
			1782	1125	298	357	2			
1	C	234	Total	C	N	O	S	0	0	0
			1779	1123	298	356	2			
1	D	234	Total	C	N	O	S	0	0	0
			1782	1125	298	357	2			

- Molecule 2 is METHYL-O3-(ALPHA-D-MANNOSE)-ALPHA-D-MANNOSE (three-letter code: MDM) (formula:  $C_{13}H_{24}O_{11}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			24	13	11		
2	B	1	Total	C	O	0	0
			24	13	11		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			24	13	11		
2	D	1	Total	C	O	0	0
			24	13	11		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		
3	D	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mn	0	0
			1	1		
4	A	1	Total	Mn	0	0
			1	1		
4	D	1	Total	Mn	0	0
			1	1		
4	C	1	Total	Mn	0	0
			1	1		

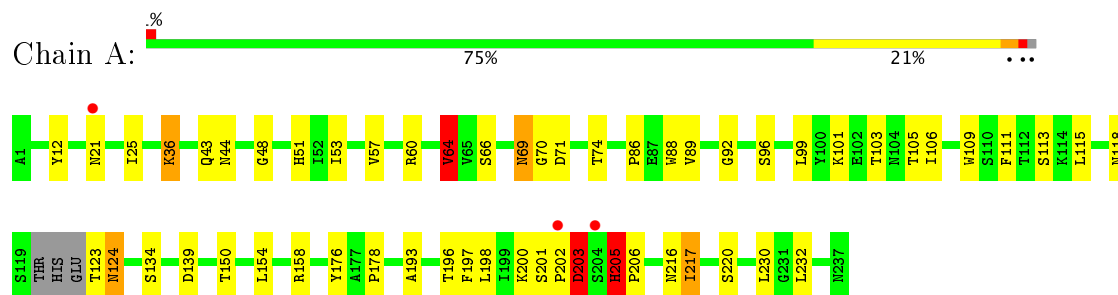
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	117	Total	O	0	0
			117	117		
5	B	137	Total	O	0	0
			137	137		
5	C	78	Total	O	0	0
			78	78		
5	D	75	Total	O	0	0
			75	75		

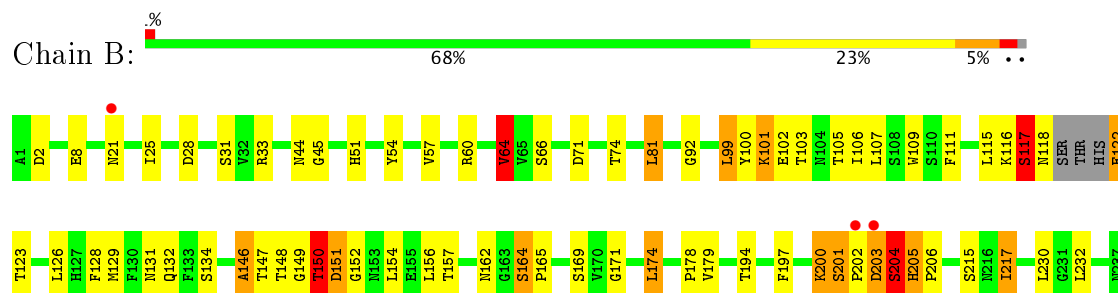
### 3 Residue-property plots [i](#)

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.

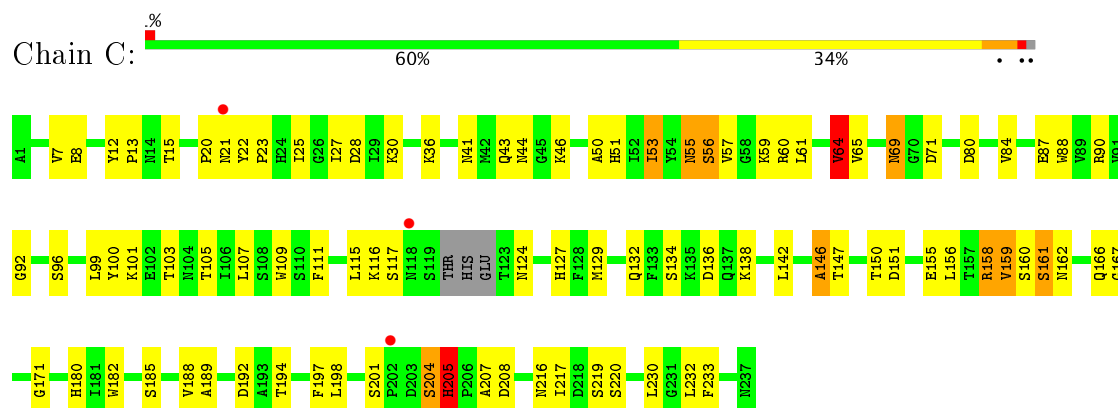
#### • Molecule 1: Concanavalin A



#### • Molecule 1: Concanavalin A

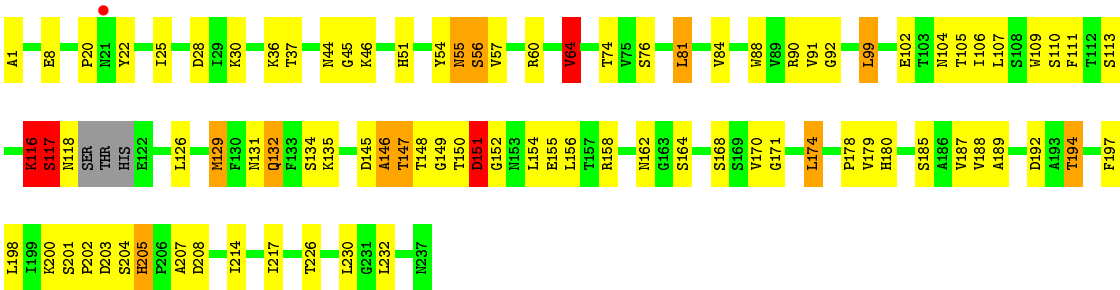


#### • Molecule 1: Concanavalin A



#### • Molecule 1: Concanavalin A





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.39 Å 69.39 Å 161.30 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.92 – 2.10 33.92 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.9 (33.92-2.10) 98.9 (33.92-2.10)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 2.10 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.189 , 0.245 0.187 , 0.240	Depositor DCC
$R_{free}$ test set	2524 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.2	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 20.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l 0.477 for h,-h-k,-l 0.022 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7633	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, MN, MDM

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	1.06	2/1819 (0.1%)	1.01	2/2476 (0.1%)
1	B	1.05	2/1822 (0.1%)	1.11	9/2480 (0.4%)
1	C	0.92	0/1819	0.96	5/2476 (0.2%)
1	D	0.93	0/1822	1.03	8/2480 (0.3%)
All	All	0.99	4/7282 (0.1%)	1.03	24/9912 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	6
1	C	0	3
1	D	0	5
All	All	0	14

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	21	ASN	CB-CG	6.91	1.67	1.51
1	B	21	ASN	CB-CG	6.29	1.65	1.51
1	A	36	LYS	CE-NZ	5.79	1.63	1.49
1	B	8	GLU	CD-OE2	-5.24	1.19	1.25

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	205	HIS	N-CA-C	9.60	136.91	111.00
1	B	64	VAL	CB-CA-C	-9.09	94.13	111.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	148	THR	N-CA-C	8.69	134.45	111.00
1	B	151	ASP	N-CA-C	8.21	133.18	111.00
1	A	64	VAL	CB-CA-C	-7.57	97.03	111.40
1	C	64	VAL	CB-CA-C	-7.46	97.23	111.40
1	B	148	THR	N-CA-C	7.21	130.46	111.00
1	D	205	HIS	N-CA-CB	-7.18	97.68	110.60
1	D	147	THR	N-CA-C	7.01	129.92	111.00
1	D	174	LEU	CA-CB-CG	7.00	131.39	115.30
1	D	64	VAL	CB-CA-C	-6.65	98.76	111.40
1	D	117	SER	N-CA-C	-6.53	93.36	111.00
1	B	174	LEU	CA-CB-CG	6.49	130.22	115.30
1	C	147	THR	N-CA-C	6.34	128.11	111.00
1	B	33	ARG	NE-CZ-NH2	-6.15	117.22	120.30
1	D	205	HIS	N-CA-C	5.99	127.16	111.00
1	B	2	ASP	CB-CG-OD2	-5.80	113.08	118.30
1	C	64	VAL	CG1-CB-CG2	5.67	119.97	110.90
1	A	205	HIS	N-CA-CB	5.60	120.68	110.60
1	B	150	THR	CB-CA-C	-5.58	96.53	111.60
1	B	204	SER	C-N-CA	5.54	135.56	121.70
1	B	146	ALA	N-CA-C	-5.34	96.57	111.00
1	D	174	LEU	CB-CG-CD2	5.19	119.83	111.00
1	C	55	ASN	N-CA-CB	-5.06	101.50	110.60

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	116	LYS	Peptide
1	B	117	SER	Peptide
1	B	146	ALA	Peptide
1	B	147	THR	Peptide
1	B	150	THR	Peptide
1	B	204	SER	Peptide
1	C	146	ALA	Peptide
1	C	160	SER	Peptide
1	C	204	SER	Peptide
1	D	116	LYS	Peptide
1	D	117	SER	Peptide
1	D	146	ALA	Peptide
1	D	147	THR	Peptide
1	D	204	SER	Peptide

## 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	1779	0	1730	39	0
1	B	1782	0	1731	60	0
1	C	1779	0	1730	70	0
1	D	1782	0	1731	67	0
2	A	24	0	24	1	0
2	B	24	0	24	1	0
2	C	24	0	23	2	0
2	D	24	0	24	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	117	0	0	6	0
5	B	137	0	0	11	0
5	C	78	0	0	13	0
5	D	75	0	0	7	0
All	All	7633	0	7017	218	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 15.

All (218) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:ARG:HH11	1:D:60:ARG:NH1	1.01	1.49
1:B:215:SER:HB2	5:B:1090:HOH:O	1.22	1.32
1:B:60:ARG:NH1	1:D:60:ARG:HH11	1.29	1.30
1:B:60:ARG:NH1	1:D:60:ARG:NH1	1.81	1.27
1:B:71:ASP:HB2	5:B:1084:HOH:O	1.43	1.16
1:A:60:ARG:HH21	1:C:60:ARG:NH2	1.50	1.10
1:D:150:THR:HA	1:D:152:GLY:N	1.69	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:SER:HB2	1:B:118:ASN:HB2	1.39	1.01
1:C:27:ILE:HG13	5:C:1066:HOH:O	1.62	1.00
1:B:202:PRO:O	5:B:1092:HOH:O	1.80	0.99
1:D:150:THR:CA	1:D:152:GLY:H	1.75	0.99
1:A:205:HIS:HD2	1:A:206:PRO:O	1.45	0.98
1:D:150:THR:HA	1:D:152:GLY:H	0.83	0.97
1:C:88:TRP:HB3	1:C:217:ILE:HD11	1.49	0.94
1:D:44:ASN:HD21	1:D:201:SER:H	1.05	0.93
1:A:60:ARG:HH21	1:C:60:ARG:HH21	0.93	0.92
1:B:64:VAL:HG13	1:B:74:THR:HG22	1.53	0.90
1:C:15:THR:HG21	1:C:21:ASN:HD22	1.38	0.89
1:B:204:SER:HB3	1:B:205:HIS:HB2	1.53	0.89
1:B:165:PRO:HD2	1:B:202:PRO:HG2	1.55	0.86
1:A:44:ASN:HD21	1:A:201:SER:H	1.22	0.86
1:B:150:THR:HA	1:B:152:GLY:H	1.41	0.85
1:C:205:HIS:HB3	5:C:1022:HOH:O	1.75	0.84
1:A:205:HIS:CD2	1:A:206:PRO:O	2.30	0.84
1:A:60:ARG:NH2	1:C:60:ARG:HH21	1.75	0.83
1:C:44:ASN:HD21	1:C:201:SER:H	1.29	0.80
1:A:60:ARG:NH2	1:C:60:ARG:NH2	2.29	0.80
1:C:156:LEU:O	1:C:171:GLY:HA3	1.81	0.80
1:A:25:ILE:HB	5:A:1108:HOH:O	1.82	0.79
1:B:118:ASN:O	5:B:1109:HOH:O	2.01	0.79
1:C:232:LEU:HD12	5:C:1068:HOH:O	1.83	0.79
1:B:117:SER:CB	1:B:118:ASN:HB2	2.12	0.79
1:A:124:ASN:HD21	1:B:132:GLN:H	1.30	0.78
1:D:44:ASN:ND2	1:D:201:SER:H	1.81	0.78
1:C:232:LEU:CD1	5:C:1068:HOH:O	2.33	0.77
1:D:149:GLY:HA2	1:D:150:THR:O	1.88	0.74
1:A:44:ASN:HD21	1:A:201:SER:N	1.86	0.73
1:C:53:ILE:C	1:C:53:ILE:HD13	2.10	0.72
1:C:59:LYS:HE2	5:C:1076:HOH:O	1.88	0.72
1:D:44:ASN:HD21	1:D:201:SER:N	1.85	0.71
1:B:45:GLY:HA2	1:B:200:LYS:HG3	1.74	0.69
1:C:61:LEU:O	5:C:1066:HOH:O	2.10	0.69
1:D:149:GLY:CA	1:D:150:THR:O	2.40	0.69
1:B:44:ASN:HD21	1:B:201:SER:H	1.40	0.69
1:D:60:ARG:HD3	1:D:76:SER:HB3	1.74	0.69
1:C:23:PRO:HA	5:C:1054:HOH:O	1.93	0.69
1:A:43:GLN:HE22	1:A:69:ASN:HD21	1.41	0.68
1:B:162:ASN:OD1	1:B:164:SER:HB2	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ASP:OD2	5:A:1040:HOH:O	2.13	0.67
1:C:90:ARG:NH1	1:C:217:ILE:HG23	2.09	0.67
1:B:205:HIS:H	1:B:206:PRO:HD3	1.59	0.67
1:D:162:ASN:HB2	5:D:1054:HOH:O	1.95	0.65
1:A:193:ALA:HB1	5:A:1098:HOH:O	1.95	0.65
1:A:118:ASN:O	5:A:1051:HOH:O	2.14	0.64
1:C:53:ILE:O	1:C:53:ILE:HD13	1.99	0.62
1:A:66:SER:HB2	5:A:1032:HOH:O	2.00	0.61
1:C:99:LEU:HB2	2:C:1003:MDM:H1	1.83	0.60
1:C:232:LEU:HG	5:C:1068:HOH:O	2.00	0.60
1:B:215:SER:CB	5:B:1090:HOH:O	2.03	0.59
1:D:45:GLY:HA2	1:D:198:LEU:HD21	1.84	0.59
1:C:87:GLU:HG3	1:C:180:HIS:CD2	2.38	0.59
1:B:203:ASP:O	1:B:205:HIS:HB2	2.03	0.58
1:C:232:LEU:CG	5:C:1068:HOH:O	2.51	0.58
1:B:117:SER:CB	1:B:118:ASN:CB	2.82	0.57
1:C:99:LEU:HD23	2:C:1003:MDM:H1'	1.85	0.57
1:D:74:THR:HG23	5:D:1018:HOH:O	2.02	0.57
1:C:124:ASN:ND2	1:D:131:ASN:H	2.02	0.57
5:B:1095:HOH:O	1:D:194:THR:HG21	2.04	0.57
1:D:99:LEU:HB2	2:D:1001:MDM:H1	1.85	0.57
1:C:27:ILE:CG1	5:C:1066:HOH:O	2.36	0.57
1:C:166:GLN:HG2	1:C:167:GLY:O	2.05	0.56
1:B:204:SER:HB3	1:B:205:HIS:CB	2.31	0.56
1:C:69:ASN:HD21	1:C:71:ASP:HB2	1.71	0.56
1:C:27:ILE:CD1	5:C:1066:HOH:O	2.52	0.56
1:A:111:PHE:CE2	1:A:113:SER:HB2	2.40	0.55
1:D:111:PHE:CE2	1:D:113:SER:HB2	2.41	0.55
1:C:43:GLN:HE22	1:C:69:ASN:ND2	2.04	0.54
1:C:56:SER:CB	1:C:189:ALA:H	2.20	0.54
1:D:170:VAL:HB	1:D:226:THR:HG22	1.89	0.54
1:B:44:ASN:HD21	1:B:201:SER:N	2.05	0.54
1:A:103:THR:OG1	1:A:200:LYS:HB2	2.08	0.54
1:D:107:LEU:O	5:D:1035:HOH:O	2.18	0.54
1:B:71:ASP:OD1	5:B:1045:HOH:O	2.18	0.54
1:C:158:ARG:HB2	1:C:166:GLN:HB3	1.89	0.54
1:C:116:LYS:HG3	1:C:188:VAL:HB	1.90	0.54
1:B:60:ARG:HH11	1:D:60:ARG:HH11	0.55	0.53
1:C:103:THR:HG22	1:C:159:VAL:CG1	2.38	0.53
1:C:20:PRO:HB2	1:C:22:TYR:CZ	2.43	0.53
1:C:8:GLU:OE1	1:C:28:ASP:OD2	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:VAL:HG22	1:D:57:VAL:CG2	2.37	0.53
1:B:215:SER:CA	5:B:1090:HOH:O	2.50	0.53
1:C:59:LYS:HD3	1:C:80:ASP:HB2	1.91	0.53
1:A:43:GLN:HE22	1:A:69:ASN:ND2	2.03	0.53
1:B:122:GLU:OE1	1:B:123:THR:OG1	2.27	0.53
1:A:57:VAL:HG21	1:C:64:VAL:HG22	1.90	0.53
1:D:99:LEU:HD12	2:D:1001:MDM:H1'	1.91	0.52
1:C:103:THR:HG22	1:C:159:VAL:HG11	1.91	0.52
1:D:36:LYS:HE2	5:D:1006:HOH:O	2.08	0.52
1:A:92:GLY:HA2	1:A:109:TRP:CH2	2.45	0.52
1:A:12:TYR:HD1	1:A:205:HIS:CD2	2.28	0.52
1:B:99:LEU:HD13	1:B:100:TYR:CE2	2.44	0.51
1:D:51:HIS:HB2	1:D:64:VAL:HG23	1.92	0.51
1:D:117:SER:OG	1:D:118:ASN:HB2	2.11	0.51
1:B:107:LEU:N	1:B:107:LEU:HD12	2.26	0.51
1:C:43:GLN:HE22	1:C:69:ASN:HD21	1.59	0.51
1:D:44:ASN:ND2	1:D:200:LYS:HA	2.26	0.51
1:D:118:ASN:HB2	1:D:187:VAL:HG21	1.93	0.50
1:B:165:PRO:CD	1:B:202:PRO:HG2	2.36	0.50
1:C:87:GLU:HG3	1:C:180:HIS:NE2	2.27	0.50
1:A:115:LEU:O	1:A:123:THR:HA	2.11	0.50
1:D:90:ARG:NH1	1:D:217:ILE:HG23	2.26	0.49
1:B:101:LYS:HG3	1:B:165:PRO:O	2.12	0.49
1:D:150:THR:HB	1:D:151:ASP:HA	1.93	0.49
1:D:55:ASN:ND2	5:D:1021:HOH:O	2.24	0.49
1:B:99:LEU:HD12	2:B:1004:MDM:H1'	1.94	0.49
1:C:155:GLU:CD	1:C:159:VAL:HG21	2.33	0.49
1:C:136:ASP:OD2	1:C:138:LYS:NZ	2.40	0.49
1:B:117:SER:HB3	1:B:118:ASN:HA	1.93	0.49
1:B:54:TYR:CE1	1:B:81:LEU:HD22	2.47	0.49
1:B:105:THR:O	1:B:197:PHE:HA	2.13	0.49
1:D:145:ASP:HB3	1:D:158:ARG:HG2	1.93	0.49
1:B:115:LEU:O	1:B:123:THR:HA	2.12	0.49
1:A:74:THR:HG23	1:C:57:VAL:HG13	1.95	0.49
1:A:178:PRO:HB3	1:A:217:ILE:HD11	1.94	0.48
1:B:178:PRO:HB3	1:B:217:ILE:HD11	1.94	0.48
1:C:142:LEU:HD22	1:C:146:ALA:HB1	1.95	0.48
1:C:107:LEU:HD11	1:C:198:LEU:HB2	1.96	0.48
1:A:106:ILE:HB	1:A:154:LEU:HB3	1.95	0.48
1:D:30:LYS:HE2	1:D:84:VAL:HG13	1.95	0.48
1:D:126:LEU:CD2	1:D:179:VAL:HG22	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:PHE:HB3	1:B:128:PHE:CZ	2.49	0.47
1:B:99:LEU:HD13	1:B:100:TYR:CZ	2.49	0.47
1:A:69:ASN:HD22	1:A:70:GLY:N	2.13	0.47
1:A:74:THR:HG23	1:C:57:VAL:CG1	2.44	0.47
1:B:117:SER:HB3	1:B:118:ASN:CA	2.45	0.47
1:C:55:ASN:HA	5:C:1017:HOH:O	2.15	0.47
1:A:88:TRP:HB3	1:A:217:ILE:HD11	1.96	0.47
1:C:88:TRP:HB3	1:C:217:ILE:CD1	2.34	0.47
1:D:8:GLU:OE1	1:D:28:ASP:OD2	2.32	0.47
1:C:51:HIS:HB2	1:C:64:VAL:HG23	1.97	0.47
1:C:216:ASN:O	1:C:219:SER:HB2	2.15	0.47
1:D:116:LYS:O	1:D:187:VAL:N	2.48	0.47
1:D:156:LEU:O	1:D:171:GLY:HA3	2.15	0.47
1:D:91:VAL:HG12	1:D:214:ILE:HG12	1.96	0.46
1:A:44:ASN:ND2	1:A:201:SER:H	2.00	0.46
1:C:87:GLU:CG	1:C:180:HIS:HE2	2.28	0.46
1:D:132:GLN:HG3	1:D:132:GLN:O	2.10	0.46
1:B:64:VAL:HG13	1:B:74:THR:CG2	2.37	0.46
1:D:88:TRP:CH2	1:D:180:HIS:HB2	2.50	0.46
1:B:156:LEU:O	1:B:171:GLY:HA3	2.15	0.46
1:D:25:ILE:O	1:D:37:THR:HA	2.15	0.46
1:B:157:THR:HB	1:B:169:SER:HB3	1.98	0.45
1:C:7:VAL:HG22	1:C:27:ILE:HD13	1.98	0.45
1:C:44:ASN:HD21	1:C:201:SER:N	2.04	0.45
1:D:102:GLU:OE2	1:D:104:ASN:OD1	2.35	0.45
1:D:116:LYS:HG3	1:D:188:VAL:HB	1.98	0.45
1:B:92:GLY:HA2	1:B:109:TRP:CH2	2.52	0.45
1:C:25:ILE:HG21	1:C:65:VAL:HG21	1.99	0.45
1:D:1:ALA:N	5:D:1067:HOH:O	2.50	0.45
1:D:46:LYS:HA	1:D:46:LYS:HD3	1.68	0.45
1:D:56:SER:HB3	1:D:189:ALA:H	1.82	0.45
1:A:96:SER:OG	1:A:230:LEU:HA	2.17	0.45
1:B:126:LEU:CD2	1:B:179:VAL:HG22	2.47	0.45
1:C:92:GLY:HA2	1:C:109:TRP:CH2	2.52	0.44
1:C:115:LEU:HD21	1:C:182:TRP:HA	1.98	0.44
1:C:232:LEU:HB3	1:C:233:PHE:CD2	2.53	0.44
1:D:54:TYR:CE1	1:D:81:LEU:HD22	2.52	0.44
1:B:118:ASN:HB3	5:B:1109:HOH:O	2.16	0.44
1:D:201:SER:OG	1:D:203:ASP:HB2	2.17	0.44
1:B:28:ASP:HB3	1:B:31:SER:O	2.16	0.44
1:B:57:VAL:CG2	1:D:64:VAL:HG22	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:HD23	2:A:1002:MDM:H7'3	2.00	0.44
1:C:207:ALA:HA	1:C:208:ASP:HA	1.82	0.44
1:A:51:HIS:HB2	1:A:64:VAL:HG23	2.00	0.44
1:B:64:VAL:HG22	1:D:57:VAL:HG21	1.98	0.44
1:B:204:SER:CB	1:B:205:HIS:HB2	2.37	0.43
1:D:117:SER:HB2	5:D:1042:HOH:O	2.17	0.43
1:D:105:THR:O	1:D:197:PHE:HA	2.18	0.43
1:B:126:LEU:HD23	1:B:179:VAL:HG22	2.01	0.43
1:C:12:TYR:HA	1:C:13:PRO:HD3	1.91	0.43
1:B:117:SER:CB	1:B:118:ASN:CA	2.96	0.43
1:B:204:SER:HB3	1:B:205:HIS:ND1	2.34	0.43
1:C:43:GLN:HB3	1:C:46:LYS:HG3	1.99	0.43
1:A:88:TRP:HB3	1:A:217:ILE:CD1	2.49	0.43
1:C:100:TYR:HB2	1:C:207:ALA:HB3	2.01	0.43
1:C:30:LYS:HD2	1:C:84:VAL:HG13	2.01	0.43
1:B:106:ILE:HB	1:B:154:LEU:HB3	1.99	0.43
1:B:25:ILE:HB	5:B:1093:HOH:O	2.18	0.43
1:C:87:GLU:CD	1:C:180:HIS:HE2	2.21	0.43
1:A:105:THR:O	1:A:197:PHE:HA	2.19	0.42
1:D:20:PRO:HB2	1:D:22:TYR:CZ	2.55	0.42
1:A:86:PRO:HG2	1:A:89:VAL:HG12	2.01	0.42
1:A:124:ASN:ND2	1:B:131:ASN:H	2.18	0.42
1:D:110:SER:HB3	1:D:129:MET:CE	2.50	0.42
1:C:124:ASN:HD21	1:D:132:GLN:H	1.67	0.42
1:B:66:SER:HB2	5:B:1122:HOH:O	2.18	0.42
1:C:50:ALA:O	1:C:194:THR:HA	2.20	0.42
1:C:30:LYS:HA	5:C:1012:HOH:O	2.20	0.42
1:A:48:GLY:O	1:A:196:THR:HA	2.19	0.42
1:D:92:GLY:HA2	1:D:109:TRP:CH2	2.55	0.42
1:C:7:VAL:HG22	1:C:27:ILE:CD1	2.50	0.41
1:D:106:ILE:HB	1:D:154:LEU:HB3	2.02	0.41
1:A:109:TRP:CD1	5:A:1098:HOH:O	2.72	0.41
1:D:126:LEU:HD23	1:D:179:VAL:HG22	2.01	0.41
1:C:96:SER:OG	1:C:230:LEU:HA	2.21	0.41
1:D:146:ALA:HA	1:D:155:GLU:O	2.21	0.41
1:C:111:PHE:O	1:C:127:HIS:HA	2.20	0.41
1:D:202:PRO:CD	1:D:203:ASP:H	2.34	0.41
1:D:178:PRO:HB3	1:D:217:ILE:CD1	2.51	0.41
1:B:51:HIS:HB2	1:B:64:VAL:HG23	2.03	0.41
1:D:202:PRO:HD2	1:D:203:ASP:H	1.86	0.41
1:C:105:THR:O	1:C:197:PHE:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:202:PRO:CD	1:D:203:ASP:N	2.84	0.41
1:D:149:GLY:CA	1:D:150:THR:C	2.89	0.41
1:D:178:PRO:HB3	1:D:217:ILE:HD11	2.02	0.41
1:A:139:ASP:C	1:A:176:TYR:HB2	2.42	0.40
1:B:102:GLU:HG3	1:B:103:THR:O	2.22	0.40
1:D:207:ALA:HA	1:D:208:ASP:HA	1.89	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	230/237 (97%)	219 (95%)	8 (4%)	3 (1%)	14	8
1	B	230/237 (97%)	216 (94%)	7 (3%)	7 (3%)	5	1
1	C	230/237 (97%)	218 (95%)	9 (4%)	3 (1%)	14	8
1	D	230/237 (97%)	219 (95%)	9 (4%)	2 (1%)	20	14
All	All	920/948 (97%)	872 (95%)	33 (4%)	15 (2%)	11	5

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	ASP
1	A	205	HIS
1	B	151	ASP
1	B	204	SER
1	B	205	HIS
1	C	161	SER
1	D	117	SER
1	D	151	ASP
1	B	149	GLY

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Mol	Chain	Res	Type
1	B	150	THR
1	B	117	SER
1	C	205	HIS
1	B	203	ASP
1	C	204	SER
1	A	202	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	199/202 (98%)	183 (92%)	16 (8%)	14	10
1	B	199/202 (98%)	183 (92%)	16 (8%)	14	10
1	C	199/202 (98%)	179 (90%)	20 (10%)	9	5
1	D	199/202 (98%)	179 (90%)	20 (10%)	9	5
All	All	796/808 (98%)	724 (91%)	72 (9%)	11	7

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LYS
1	A	53	ILE
1	A	64	VAL
1	A	69	ASN
1	A	71	ASP
1	A	101	LYS
1	A	124	ASN
1	A	134	SER
1	A	150	THR
1	A	158	ARG
1	A	198	LEU
1	A	203	ASP
1	A	216	ASN
1	A	217	ILE
1	A	220	SER

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Mol	Chain	Res	Type
1	A	232	LEU
1	B	64	VAL
1	B	81	LEU
1	B	99	LEU
1	B	101	LYS
1	B	122	GLU
1	B	129	MET
1	B	134	SER
1	B	150	THR
1	B	164	SER
1	B	174	LEU
1	B	194	THR
1	B	200	LYS
1	B	201	SER
1	B	217	ILE
1	B	230	LEU
1	B	232	LEU
1	C	36	LYS
1	C	41	ASN
1	C	53	ILE
1	C	56	SER
1	C	64	VAL
1	C	69	ASN
1	C	101	LYS
1	C	117	SER
1	C	129	MET
1	C	132	GLN
1	C	134	SER
1	C	150	THR
1	C	151	ASP
1	C	158	ARG
1	C	159	VAL
1	C	161	SER
1	C	162	ASN
1	C	185	SER
1	C	192	ASP
1	C	220	SER
1	D	55	ASN
1	D	56	SER
1	D	64	VAL
1	D	81	LEU
1	D	99	LEU

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Mol	Chain	Res	Type
1	D	116	LYS
1	D	129	MET
1	D	132	GLN
1	D	134	SER
1	D	135	LYS
1	D	151	ASP
1	D	164	SER
1	D	168	SER
1	D	174	LEU
1	D	185	SER
1	D	192	ASP
1	D	194	THR
1	D	205	HIS
1	D	230	LEU
1	D	232	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	44	ASN
1	A	69	ASN
1	A	83	ASN
1	A	104	ASN
1	A	124	ASN
1	A	166	GLN
1	A	205	HIS
1	B	41	ASN
1	B	44	ASN
1	B	51	HIS
1	B	104	ASN
1	B	132	GLN
1	B	166	GLN
1	C	21	ASN
1	C	44	ASN
1	C	69	ASN
1	C	104	ASN
1	C	124	ASN
1	C	132	GLN
1	C	166	GLN
1	D	44	ASN
1	D	104	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 ⓘ

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 for Mogul analysis.

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	MDM	A	1002	-	25,25,25	0.64	0	36,36,36	2.35	8 (22%)
2	MDM	B	1004	-	25,25,25	0.70	0	36,36,36	2.54	9 (25%)
2	MDM	C	1003	-	25,25,25	0.79	0	36,36,36	3.03	11 (30%)
2	MDM	D	1001	-	25,25,25	0.76	1 (4%)	36,36,36	2.73	14 (38%)

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	MDM	A	1002	-	1/1/10/10	0/10/50/50	0/2/2/2
2	MDM	B	1004	-	1/1/10/10	0/10/50/50	0/2/2/2
2	MDM	C	1003	-	2/2/10/10	0/10/50/50	0/2/2/2
2	MDM	D	1001	-	2/2/10/10	0/10/50/50	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1001	MDM	C4'-C3'	2.06	1.58	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1004	MDM	C1'-C2'-C3'	-7.82	96.44	110.04
2	A	1002	MDM	C1'-C2'-C3'	-7.00	97.86	110.04
2	C	1003	MDM	O3'-C1-C2	-4.97	96.92	108.11
2	D	1001	MDM	O3'-C1-C2	-4.90	97.07	108.11
2	A	1002	MDM	C1'-O5'-C5'	-4.67	104.92	113.72
2	A	1002	MDM	O2-C2-C1	-4.28	101.08	110.03
2	B	1004	MDM	C1'-O5'-C5'	-4.22	105.76	113.72
2	C	1003	MDM	O2'-C2'-C3'	-4.00	100.77	109.87
2	B	1004	MDM	O2-C2-C1	-3.88	101.91	110.03
2	D	1001	MDM	O2'-C2'-C3'	-3.59	101.69	109.87
2	C	1003	MDM	C7'-O1'-C1'	-3.28	108.16	113.29
2	D	1001	MDM	C1'-O5'-C5'	-3.07	107.94	113.72
2	C	1003	MDM	O5'-C1'-C2'	-2.71	105.07	110.30
2	B	1004	MDM	O5-C1-C2	-2.64	105.21	110.30
2	C	1003	MDM	C1'-O5'-C5'	-2.62	108.78	113.72
2	B	1004	MDM	O2'-C2'-C3'	-2.60	103.95	109.87
2	D	1001	MDM	O5'-C1'-C2'	-2.28	105.89	110.30
2	A	1002	MDM	C1-O3'-C3'	-2.28	112.45	118.00
2	D	1001	MDM	C1'-C2'-C3'	-2.20	106.22	110.04
2	B	1004	MDM	O3'-C1-C2	2.02	112.66	108.11
2	D	1001	MDM	O3-C3-C4	2.04	114.80	110.36
2	D	1001	MDM	O2-C2-C1	2.10	114.42	110.03
2	D	1001	MDM	O2'-C2'-C1'	2.22	114.68	110.03
2	D	1001	MDM	O3'-C3'-C2'	2.41	113.00	107.19
2	D	1001	MDM	C3'-C4'-C5'	2.42	114.86	109.67
2	A	1002	MDM	O2-C2-C3	2.43	115.65	110.36
2	A	1002	MDM	O2'-C2'-C1'	2.54	115.34	110.03
2	C	1003	MDM	O2'-C2'-C1'	2.74	115.77	110.03
2	B	1004	MDM	O5'-C5'-C6'	2.83	113.19	106.41
2	C	1003	MDM	O4'-C4'-C3'	3.18	117.11	109.87
2	A	1002	MDM	O5'-C5'-C6'	3.21	114.09	106.41
2	B	1004	MDM	O2'-C2'-C1'	4.33	119.08	110.03
2	D	1001	MDM	O3'-C3'-C4'	4.49	118.00	107.19
2	C	1003	MDM	O3'-C3'-C4'	5.57	120.59	107.19
2	D	1001	MDM	C1-O3'-C3'	6.57	133.99	118.00
2	D	1001	MDM	C4'-C3'-C2'	6.66	120.42	110.83
2	C	1003	MDM	C1-O3'-C3'	6.90	134.80	118.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1002	MDM	O1'-C1'-C2'	7.43	116.97	108.14
2	C	1003	MDM	C4'-C3'-C2'	7.43	121.53	110.83
2	D	1001	MDM	O1'-C1'-C2'	7.63	117.21	108.14
2	B	1004	MDM	O1'-C1'-C2'	8.44	118.18	108.14
2	C	1003	MDM	O1'-C1'-C2'	9.28	119.18	108.14

All (6) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1003	MDM	C3'
2	C	1003	MDM	C1'
2	D	1001	MDM	C3'
2	D	1001	MDM	C1'
2	A	1002	MDM	C1'
2	B	1004	MDM	C1'

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1002	MDM	1	0
2	B	1004	MDM	1	0
2	C	1003	MDM	2	0
2	D	1001	MDM	2	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	234/237 (98%)	-0.41	3 (1%) 77 81	11, 18, 33, 51	6 (2%)
1	B	234/237 (98%)	-0.38	3 (1%) 77 81	12, 18, 34, 47	4 (1%)
1	C	234/237 (98%)	-0.18	3 (1%) 77 81	15, 24, 40, 60	7 (2%)
1	D	234/237 (98%)	-0.23	1 (0%) 92 93	15, 24, 39, 51	6 (2%)
All	All	936/948 (98%)	-0.30	10 (1%) 80 84	11, 21, 37, 60	23 (2%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	203	ASP	3.6
1	C	118	ASN	3.3
1	A	21	ASN	3.1
1	A	202	PRO	2.9
1	B	202	PRO	2.6
1	B	21	ASN	2.6
1	D	21	ASN	2.6
1	C	21	ASN	2.5
1	A	204	SER	2.3
1	C	202	PRO	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(Å <sup>2</sup> )	Q<0.9
4	MN	B	239	1/1	1.00	0.10	1.65	17,17,17,17	0
2	MDM	B	1004	24/24	0.96	0.11	0.59	14,22,34,40	0
2	MDM	C	1003	24/24	0.95	0.14	0.44	19,31,41,44	0
2	MDM	A	1002	24/24	0.95	0.11	0.31	14,24,35,38	0
4	MN	A	239	1/1	1.00	0.09	-0.20	17,17,17,17	0
2	MDM	D	1001	24/24	0.95	0.11	-0.34	22,33,42,44	0
4	MN	C	239	1/1	0.99	0.08	-0.98	22,22,22,22	0
3	CA	A	238	1/1	1.00	0.07	-1.49	18,18,18,18	0
4	MN	D	239	1/1	0.99	0.07	-1.65	23,23,23,23	0
3	CA	C	238	1/1	0.97	0.07	-1.69	20,20,20,20	0
3	CA	B	238	1/1	1.00	0.07	-1.79	17,17,17,17	0
3	CA	D	238	1/1	0.99	0.05	-3.71	21,21,21,21	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.