



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 8, 2021 – 02:06 PM EST

PDB ID : 7KJN  
Title : CRYSTAL STRUCTURE OF HUMAN MDMX IN COMPLEX WITH D-PEPTIDE INHIBITOR (DPMI-OMEGA)  
Authors : Tolbert, W.D.; Pazgier, M.  
Deposited on : 2020-10-26  
Resolution : 2.80 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

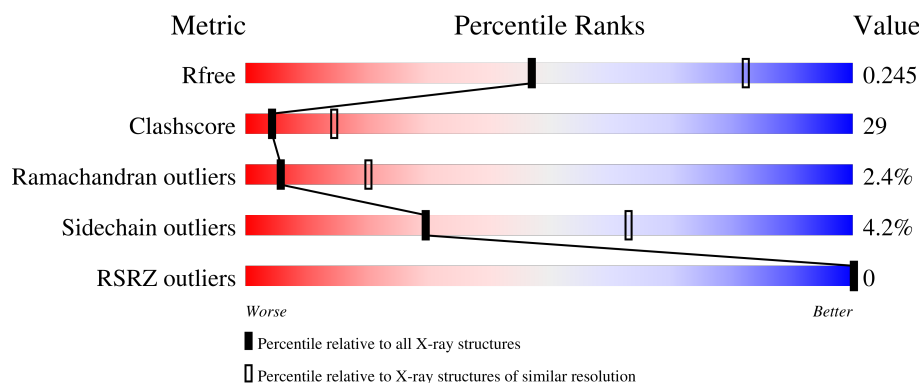
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


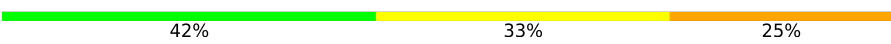

The reported resolution of this entry is 2.80 Å.

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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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	85	 61% 35% .
2	B	12	 42% 33% 25%
2	C	12	 50% 25% 25%

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
2	DAR	C	12	-	-	X	-

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 910 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 Protein Mdm4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	85	Total	C	N	O	S	0	0	0
			670	432	114	118	6			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	68	ALA	GLN	engineered mutation	UNP O15151
A	69	ALA	GLN	engineered mutation	UNP O15151
A	70	ALA	GLU	engineered mutation	UNP O15151

- Molecule 2 is a protein (with D amino acids) called D-PMI-omega.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	12	Total	C	Cl	N	O	0	0	0
			120	82	1	17	20			
2	C	12	Total	C	Cl	N	O	0	0	0
			120	82	1	17	20			

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

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

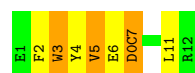
- Molecule 1: Protein Mdm4

Chain A: 



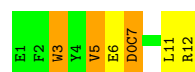
- Molecule 2: D-PMI-omega

Chain B: 



- Molecule 2: D-PMI-omega

Chain C: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	48.85Å 48.85Å 90.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.30 – 2.80 42.30 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.6 (42.30-2.80) 98.6 (42.30-2.80)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.41 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0258, PHENIX 1.17	Depositor
R, $R_{free}$	0.208 , 0.244 0.211 , 0.245	Depositor DCC
$R_{free}$ test set	184 reflections (5.59%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.9	Xtriage
Anisotropy	0.816	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 30.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.045 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	910	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.99% 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: DAR, DTR, DGL, DTY, DLY, DPN, DVA, DLE, D0C

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	0.70	0/683	1.00	0/919

There are no bond length outliers.

There are no bond angle outliers.

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	670	0	691	40	0
2	B	120	0	96	11	0
2	C	120	0	96	14	0
All	All	910	0	883	52	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 29.

All (52) 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:71:GLN:O	2:B:4:DTY:HE1	1.32	1.25

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:7:D0C:CE1	2:B:11:DLE:HD21	1.83	1.08
1:A:71:GLN:O	2:B:4:DTY:CE1	2.14	0.96
1:A:29:PRO:HG2	1:A:34:LEU:HD12	1.63	0.80
1:A:24:ILE:N	2:C:12:DAR:NH1	2.30	0.80
1:A:24:ILE:HD12	2:C:12:DAR:HB2	1.64	0.78
2:C:5:DVA:O	2:C:6:DGL:HB2	1.87	0.73
1:A:24:ILE:HA	2:C:12:DAR:HB3	1.72	0.71
2:C:12:DAR:O	2:C:12:DAR:HG3	1.91	0.70
1:A:59:TYR:OH	1:A:79:ASP:OD2	2.08	0.68
1:A:95:PRO:HG3	2:B:11:DLE:HD12	1.78	0.65
2:B:7:D0C:CD1	2:B:11:DLE:HD21	2.27	0.64
1:A:71:GLN:HB3	2:B:3:DTR:HD1	1.80	0.64
1:A:76:CYS:SG	1:A:76:CYS:O	2.58	0.62
2:B:5:DVA:HG23	2:B:6:DGL:HG2	1.83	0.61
1:A:24:ILE:N	2:C:12:DAR:HH11	1.98	0.59
2:B:7:D0C:CE1	2:B:11:DLE:CD2	2.71	0.58
1:A:96:SER:N	1:A:97:PRO:CD	2.68	0.57
1:A:53:MET:HE2	1:A:99:TYR:CE1	2.40	0.57
1:A:31:LEU:N	1:A:32:PRO:HD2	2.20	0.57
1:A:72:HIS:O	1:A:91:SER:HA	2.05	0.56
1:A:24:ILE:HD12	2:C:12:DAR:CB	2.33	0.56
1:A:53:MET:CE	1:A:99:TYR:CE1	2.89	0.56
2:B:7:D0C:CD1	2:B:11:DLE:CD2	2.84	0.55
1:A:28:ARG:HB3	1:A:107:VAL:HG22	1.93	0.50
1:A:72:HIS:HD1	1:A:91:SER:HG	1.57	0.50
1:A:24:ILE:HA	2:C:12:DAR:CB	2.42	0.49
1:A:85:LEU:HD23	1:A:101:MET:HB2	1.95	0.48
2:C:3:DTR:C	2:C:3:DTR:HE3	2.44	0.48
1:A:106:LEU:HD12	2:C:7:D0C:CL	2.51	0.48
2:C:3:DTR:C	2:C:3:DTR:CE3	2.92	0.47
1:A:72:HIS:ND1	1:A:91:SER:OG	2.42	0.46
2:C:3:DTR:HE3	2:C:3:DTR:O	2.16	0.45
1:A:63:LYS:HB2	1:A:65:LEU:HG	1.99	0.45
1:A:59:TYR:HA	1:A:62:VAL:HG22	2.00	0.44
1:A:71:GLN:HB3	2:B:3:DTR:CD1	2.47	0.44
1:A:29:PRO:HA	1:A:106:LEU:HD23	2.00	0.44
1:A:107:VAL:CG2	1:A:107:VAL:O	2.66	0.43
1:A:60:ILE:HD13	1:A:74:VAL:HG11	2.00	0.43
1:A:56:LEU:HA	1:A:56:LEU:HD23	1.76	0.43
1:A:31:LEU:N	1:A:32:PRO:CD	2.81	0.43
1:A:40:ALA:HB2	1:A:59:TYR:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:MET:HE1	1:A:99:TYR:CE1	2.54	0.42
1:A:84:LEU:HD12	1:A:84:LEU:HA	1.83	0.42
1:A:65:LEU:O	1:A:75:TYR:N	2.35	0.41
1:A:24:ILE:HA	2:C:12:DAR:HG2	2.02	0.41
1:A:72:HIS:O	1:A:91:SER:OG	2.33	0.41
1:A:85:LEU:HD23	1:A:85:LEU:HA	1.87	0.41
2:C:11:DLE:O	2:C:12:DAR:HB3	2.20	0.40
1:A:75:TYR:CD1	1:A:89:SER:HB3	2.57	0.40
1:A:107:VAL:O	1:A:107:VAL:HG23	2.20	0.40
2:B:5:DVA:HG23	2:B:6:DGL:N	2.36	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	83/85 (98%)	75 (90%)	6 (7%)	2 (2%)	<b>6</b> <b>20</b>

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	70	ALA
1	A	72	HIS

### 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	72/72 (100%)	69 (96%)	3 (4%)	30 63

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	VAL
1	A	34	LEU
1	A	84	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

24 non-standard protein/DNA/RNA residues 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	DPN	B	2	2	10,11,12	0.26	0	10,13,15	1.18	1 (10%)
2	DTY	B	4	2	11,12,13	0.40	0	12,15,17	0.61	0
2	DLE	C	11	2	5,7,8	0.82	0	5,8,10	0.35	0
2	DGL	B	8	2	4,8,9	0.71	0	1,9,11	1.48	0
2	DAR	C	12	2	7,11,11	0.38	0	6,13,13	0.29	0
2	DLE	C	10	2	5,7,8	0.29	0	5,8,10	0.56	0
2	DAR	B	12	2	7,11,11	0.45	0	6,13,13	0.58	0
2	D0C	C	7	2	11,12,13	1.06	0	12,15,17	1.11	2 (16%)
2	DTR	B	3	2	13,15,16	0.80	0	13,20,22	1.30	2 (15%)
2	DGL	B	6	2	4,8,9	0.54	0	1,9,11	0.50	0
2	DLY	C	9	2	7,8,9	0.29	0	3,8,10	0.06	0
2	D0C	B	7	2	11,12,13	2.01	3 (27%)	12,15,17	2.55	4 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	DVA	B	5	2	4,6,7	0.74	0	6,7,9	1.17	1 (16%)
2	DLE	B	10	2	5,7,8	0.65	0	5,8,10	0.21	0
2	DGL	C	8	2	4,8,9	0.70	0	1,9,11	0.25	0
2	DLY	B	9	2	7,8,9	0.38	0	3,8,10	0.29	0
2	DGL	C	6	2	4,8,9	0.49	0	1,9,11	1.22	0
2	DTY	C	4	2	11,12,13	0.41	0	12,15,17	0.55	0
2	DTR	C	3	2	13,15,16	0.76	0	13,20,22	1.86	3 (23%)
2	DGL	C	1	2	4,8,9	0.83	0	1,9,11	0.09	0
2	DVA	C	5	2	4,6,7	0.46	0	6,7,9	1.57	1 (16%)
2	DGL	B	1	2	4,8,9	0.59	0	1,9,11	0.12	0
2	DLE	B	11	2	5,7,8	0.69	0	5,8,10	0.51	0
2	DPN	C	2	2	10,11,12	0.47	0	10,13,15	0.87	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	DPN	B	2	2	-	2/5/6/8	0/1/1/1
2	DTY	B	4	2	-	1/5/6/8	0/1/1/1
2	DLE	C	11	2	-	0/5/6/8	-
2	DGL	B	8	2	-	3/4/7/9	-
2	DAR	C	12	2	-	4/7/11/11	-
2	DLE	C	10	2	-	0/5/6/8	-
2	DAR	B	12	2	-	5/7/11/11	-
2	D0C	C	7	2	-	0/5/6/8	0/1/1/1
2	DTR	B	3	2	-	0/4/6/8	0/2/2/2
2	DGL	B	6	2	-	1/4/7/9	-
2	DLY	C	9	2	-	4/6/7/9	-
2	D0C	B	7	2	-	0/5/6/8	0/1/1/1
2	DVA	B	5	2	-	0/5/6/8	-
2	DLE	B	10	2	-	2/5/6/8	-
2	DGL	C	8	2	-	3/4/7/9	-
2	DLY	B	9	2	-	1/6/7/9	-
2	DGL	C	6	2	-	2/4/7/9	-
2	DTY	C	4	2	-	0/5/6/8	0/1/1/1
2	DTR	C	3	2	-	2/4/6/8	0/2/2/2
2	DGL	C	1	2	-	3/4/7/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DVA	C	5	2	-	4/5/6/8	-
2	DGL	B	1	2	-	3/4/7/9	-
2	DLE	B	11	2	-	2/5/6/8	-
2	DPN	C	2	2	-	4/5/6/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	7	D0C	CZ-CL	-4.52	1.64	1.74
2	B	7	D0C	CE1-CZ	-3.01	1.32	1.38
2	B	7	D0C	CE1-CD1	2.80	1.43	1.38

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	7	D0C	CB-CA-C	-7.31	97.76	111.47
2	C	3	DTR	CG-CB-CA	4.13	120.91	114.53
2	B	2	DPN	CB-CA-C	3.58	118.17	111.47
2	C	5	DVA	CB-CA-C	-3.42	108.32	112.94
2	C	3	DTR	CB-CA-C	3.39	117.83	111.47
2	B	7	D0C	CG-CB-CA	-2.76	108.52	114.10
2	C	3	DTR	CB-CG-CD1	-2.75	124.57	127.97
2	B	3	DTR	CB-CG-CD1	-2.65	124.70	127.97
2	B	5	DVA	O-C-CA	-2.58	118.01	124.78
2	C	7	D0C	CG-CB-CA	-2.56	108.92	114.10
2	C	7	D0C	CB-CA-C	2.38	115.92	111.47
2	B	7	D0C	CE2-CD2-CG	-2.27	117.91	121.03
2	B	3	DTR	CB-CG-CD2	2.23	129.72	126.25
2	B	7	D0C	CD1-CE1-CZ	-2.04	117.09	119.24

There are no chirality outliers.

All (46) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	DGL	N-CA-CB-CG
2	B	1	DGL	C-CA-CB-CG
2	B	1	DGL	CA-CB-CG-CD
2	C	1	DGL	N-CA-CB-CG
2	C	1	DGL	CA-CB-CG-CD
2	B	2	DPN	C-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
2	C	2	DPN	N-CA-CB-CG
2	C	2	DPN	C-CA-CB-CG
2	C	3	DTR	N-CA-CB-CG
2	C	3	DTR	C-CA-CB-CG
2	B	4	DTY	O-C-CA-CB
2	C	5	DVA	N-CA-CB-CG1
2	C	5	DVA	C-CA-CB-CG1
2	C	5	DVA	C-CA-CB-CG2
2	C	6	DGL	N-CA-CB-CG
2	C	6	DGL	C-CA-CB-CG
2	B	8	DGL	O-C-CA-CB
2	C	8	DGL	N-CA-CB-CG
2	C	8	DGL	C-CA-CB-CG
2	C	9	DLY	N-CA-CB-CG
2	C	9	DLY	C-CA-CB-CG
2	B	12	DAR	N-CA-CB-CG
2	B	12	DAR	C-CA-CB-CG
2	B	12	DAR	NH1-CZ-NE-CD
2	B	12	DAR	NH2-CZ-NE-CD
2	C	12	DAR	CG-CD-NE-CZ
2	C	12	DAR	NH1-CZ-NE-CD
2	C	12	DAR	NH2-CZ-NE-CD
2	C	2	DPN	CA-CB-CG-CD1
2	C	2	DPN	CA-CB-CG-CD2
2	C	12	DAR	CA-CB-CG-CD
2	C	9	DLY	CE-CD-CG-CB
2	B	2	DPN	N-CA-CB-CG
2	B	9	DLY	CE-CD-CG-CB
2	C	1	DGL	C-CA-CB-CG
2	B	8	DGL	C-CA-CB-CG
2	B	11	DLE	CA-CB-CG-CD1
2	B	11	DLE	CA-CB-CG-CD2
2	B	10	DLE	CA-CB-CG-CD2
2	C	9	DLY	CG-CD-CE-NZ
2	B	12	DAR	NE-CD-CG-CB
2	C	5	DVA	N-CA-CB-CG2
2	B	10	DLE	CA-CB-CG-CD1
2	B	8	DGL	N-CA-CB-CG
2	B	6	DGL	CA-CB-CG-CD
2	C	8	DGL	CA-CB-CG-CD

There are no ring outliers.

12 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	4	DTY	2	0
2	C	11	DLE	1	0
2	C	12	DAR	9	0
2	C	7	D0C	1	0
2	B	3	DTR	2	0
2	B	6	DGL	2	0
2	B	7	D0C	4	0
2	B	5	DVA	2	0
2	C	6	DGL	1	0
2	C	3	DTR	3	0
2	C	5	DVA	1	0
2	B	11	DLE	5	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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	85/85 (100%)	-0.38	0 <a href="#">100</a> <a href="#">100</a>	40, 55, 66, 94	0
2	B	0/12	-	-	-	-
2	C	0/12	-	-	-	-
All	All	85/109 (77%)	-0.38	0 <a href="#">100</a> <a href="#">100</a>	40, 55, 66, 94	0

There are no RSRZ outliers to report.

### 6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	DGL	C	1	9/10	0.74	0.31	106,115,121,125	0
2	DGL	B	1	9/10	0.85	0.24	71,85,90,92	0
2	DPN	C	2	11/12	0.87	0.21	83,92,103,104	0
2	DAR	C	12	12/12	0.89	0.30	65,74,87,87	0
2	DLY	C	9	9/10	0.91	0.21	52,53,70,71	0
2	DTY	B	4	12/13	0.91	0.20	52,55,58,59	0
2	DAR	B	12	12/12	0.92	0.15	64,74,79,80	0
2	DTY	C	4	12/13	0.92	0.17	49,53,59,61	0
2	DVA	C	5	7/8	0.93	0.20	52,53,58,70	0
2	DGL	C	6	9/10	0.93	0.22	48,57,90,99	0
2	D0C	C	7	12/13	0.93	0.20	48,52,59,77	0
2	DLY	B	9	9/10	0.94	0.15	49,53,58,66	0
2	D0C	B	7	12/13	0.94	0.22	48,62,69,84	0
2	DPN	B	2	11/12	0.94	0.18	52,64,67,67	0
2	DGL	C	8	9/10	0.94	0.23	55,62,90,97	0
2	DGL	B	6	9/10	0.95	0.19	57,63,78,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	DVA	B	5	7/8	0.95	0.23	52,54,57,59	0
2	DTR	C	3	14/15	0.95	0.20	58,62,70,77	0
2	DLE	B	11	8/9	0.96	0.22	63,66,74,75	0
2	DLE	C	10	8/9	0.97	0.23	49,50,51,51	0
2	DGL	B	8	9/10	0.97	0.22	49,54,57,58	0
2	DLE	C	11	8/9	0.97	0.24	53,59,61,61	0
2	DTR	B	3	14/15	0.97	0.21	46,50,54,58	0
2	DLE	B	10	8/9	0.97	0.18	59,61,67,68	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

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

There are no such residues in this entry.