



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 08:28 pm GMT

PDB ID : 1DLR
Title : METHOTREXATE-RESISTANT VARIANTS OF HUMAN DIHYDROFOLATE REDUCTASE WITH SUBSTITUTION OF LEUCINE 22: KINETICS, CRYSTALLOGRAPHY AND POTENTIAL AS SELECTABLE MARKERS
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Deposited on : 1995-01-25
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

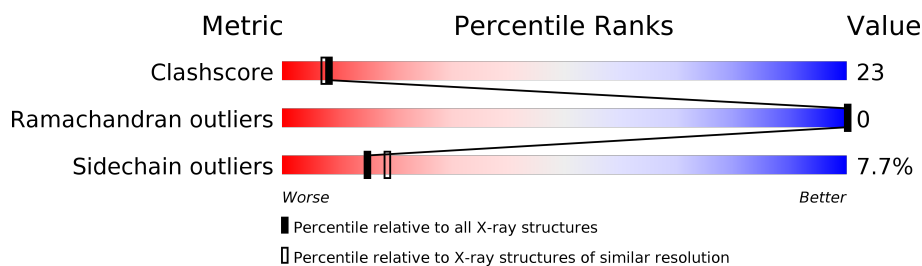
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	186	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1646 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.

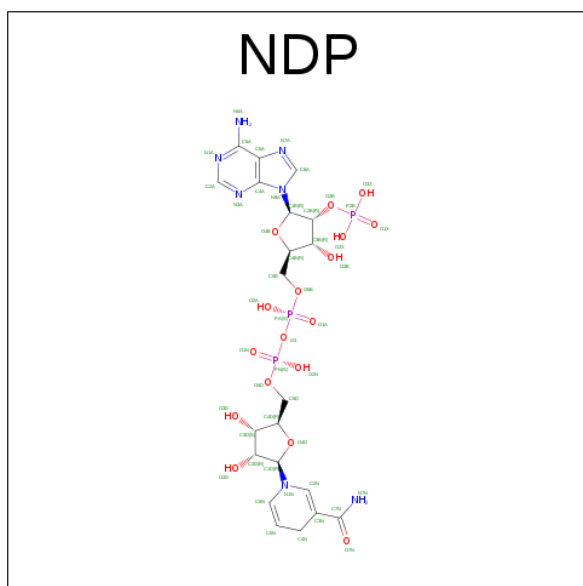
- Molecule 1 is a protein called DIHYDROFOLATE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	186	Total	C	N	O	S	0	0	0
			1505	966	253	279	7			

There is a discrepancy between the modelled and reference sequences:

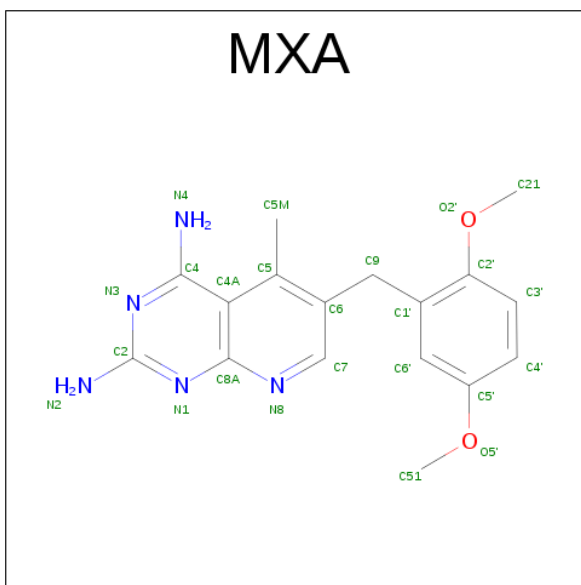
Chain	Residue	Modelled	Actual	Comment	Reference
A	22	PHE	LEU	CONFLICT	UNP P00374

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is 6-(2,5-DIMETHOXY-BENZYL)-5-METHYL-PYRIDO[2,3-D]PYRIMIDINE-2,4-DIAMINE (three-letter code: MXA) (formula: $C_{17}H_{19}N_5O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			24	17	5	2		

- Molecule 4 is water.

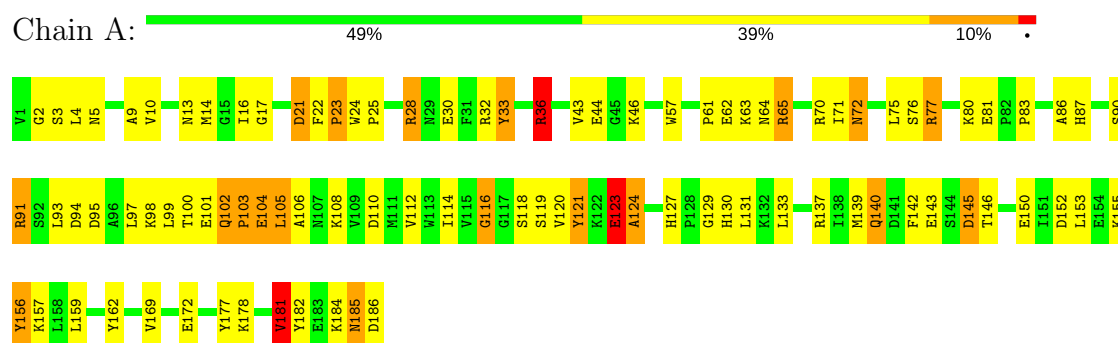
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	69	Total	O	0	0
			69	69		

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.

Note EDS was not executed.

• Molecule 1: DIHYDROFOLATE REDUCTASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	87.00Å 87.00Å 76.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.167 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1646	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.05	2/1541 (0.1%)	2.29	70/2078 (3.4%)

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	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	121	TYR	CZ-OH	-8.79	1.23	1.37
1	A	105	LEU	C-N	-6.00	1.20	1.34

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	ARG	NE-CZ-NH2	-16.01	112.30	120.30
1	A	77	ARG	NE-CZ-NH1	-14.23	113.19	120.30
1	A	110	ASP	CB-CG-OD2	12.89	129.90	118.30
1	A	110	ASP	CB-CG-OD1	-12.59	106.97	118.30
1	A	70	ARG	NE-CZ-NH1	12.27	126.44	120.30
1	A	36	ARG	NE-CZ-NH1	11.52	126.06	120.30
1	A	21	ASP	CB-CG-OD1	11.17	128.35	118.30
1	A	137	ARG	NE-CZ-NH1	10.73	125.66	120.30
1	A	162	TYR	CB-CG-CD1	-9.72	115.17	121.00
1	A	65	ARG	NE-CZ-NH2	8.53	124.56	120.30
1	A	116	GLY	CA-C-O	-8.48	105.33	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	TYR	CB-CG-CD2	8.35	126.01	121.00
1	A	90	SER	N-CA-CB	8.27	122.90	110.50
1	A	21	ASP	CB-CG-OD2	-8.18	110.94	118.30
1	A	145	ASP	CB-CG-OD1	-8.14	110.97	118.30
1	A	140	GLN	N-CA-CB	8.03	125.06	110.60
1	A	156	TYR	CB-CG-CD1	-7.93	116.24	121.00
1	A	23	PRO	O-C-N	-7.66	110.44	122.70
1	A	33	TYR	CB-CG-CD1	-7.35	116.59	121.00
1	A	33	TYR	CB-CG-CD2	7.15	125.29	121.00
1	A	28	ARG	CD-NE-CZ	7.03	133.44	123.60
1	A	81	GLU	CG-CD-OE1	7.01	132.32	118.30
1	A	28	ARG	NE-CZ-NH2	6.87	123.73	120.30
1	A	80	LYS	CA-CB-CG	6.64	128.01	113.40
1	A	121	TYR	CB-CG-CD2	-6.56	117.06	121.00
1	A	143	GLU	CB-CG-CD	6.53	131.82	114.20
1	A	80	LYS	N-CA-CB	6.44	122.19	110.60
1	A	152	ASP	CB-CG-OD2	6.37	124.03	118.30
1	A	145	ASP	CB-CG-OD2	6.33	124.00	118.30
1	A	91	ARG	CD-NE-CZ	6.29	132.41	123.60
1	A	181	VAL	CB-CA-C	6.23	123.23	111.40
1	A	123	GLU	O-C-N	-6.19	112.80	122.70
1	A	95	ASP	CB-CG-OD1	6.15	123.83	118.30
1	A	101	GLU	CA-CB-CG	6.14	126.92	113.40
1	A	5	ASN	CB-CG-OD1	6.10	133.79	121.60
1	A	9	ALA	N-CA-CB	6.07	118.60	110.10
1	A	157	LYS	CB-CA-C	-6.01	98.38	110.40
1	A	118	SER	CB-CA-C	-6.00	98.69	110.10
1	A	23	PRO	CA-C-N	5.90	130.17	117.20
1	A	80	LYS	O-C-N	5.89	132.13	122.70
1	A	156	TYR	CG-CD2-CE2	5.88	126.00	121.30
1	A	121	TYR	CD1-CE1-CZ	-5.75	114.63	119.80
1	A	150	GLU	CA-CB-CG	5.74	126.04	113.40
1	A	5	ASN	OD1-CG-ND2	-5.63	108.95	121.90
1	A	116	GLY	O-C-N	5.63	132.76	123.20
1	A	177	TYR	O-C-N	5.60	131.66	122.70
1	A	133	LEU	CA-CB-CG	5.60	128.18	115.30
1	A	91	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	A	3	SER	CB-CA-C	-5.51	99.64	110.10
1	A	182	TYR	O-C-N	5.50	131.49	122.70
1	A	81	GLU	CG-CD-OE2	-5.44	107.41	118.30
1	A	153	LEU	CB-CA-C	5.40	120.47	110.20
1	A	172	GLU	OE1-CD-OE2	5.36	129.73	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	TYR	CZ-CE2-CD2	-5.35	114.98	119.80
1	A	71	ILE	CB-CA-C	5.33	122.27	111.60
1	A	21	ASP	O-C-N	-5.32	114.19	122.70
1	A	145	ASP	CB-CA-C	5.31	121.02	110.40
1	A	44	GLU	CG-CD-OE1	5.29	128.88	118.30
1	A	77	ARG	CG-CD-NE	-5.28	100.71	111.80
1	A	13	ASN	C-N-CA	5.27	134.88	121.70
1	A	65	ARG	NH1-CZ-NH2	-5.20	113.68	119.40
1	A	77	ARG	NE-CZ-NH2	5.20	122.90	120.30
1	A	36	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	124	ALA	N-CA-CB	5.13	117.28	110.10
1	A	129	GLY	O-C-N	5.13	130.90	122.70
1	A	169	VAL	CA-CB-CG2	5.13	118.59	110.90
1	A	102	GLN	N-CA-CB	5.04	119.66	110.60
1	A	121	TYR	CZ-CE2-CD2	-5.03	115.27	119.80
1	A	44	GLU	CB-CA-C	-5.03	100.35	110.40
1	A	143	GLU	CA-CB-CG	5.02	124.44	113.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	GLU	Mainchain

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	1505	0	1507	68	0
2	A	48	0	26	6	0
3	A	24	0	19	6	0
4	A	69	0	0	3	0
All	All	1646	0	1552	73	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 23.

All (73) 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:139:MET:HE3	1:A:178:LYS:HE2	1.16	1.14
1:A:105:LEU:HD23	1:A:108:LYS:HD2	1.19	1.08
1:A:139:MET:CE	1:A:178:LYS:HE2	1.91	1.00
1:A:72:ASN:H	1:A:87:HIS:HD2	1.15	0.93
1:A:36:ARG:HG2	1:A:36:ARG:HH11	1.34	0.91
1:A:139:MET:HE3	1:A:178:LYS:CE	2.04	0.86
1:A:105:LEU:CD2	1:A:108:LYS:HD2	2.05	0.86
1:A:76:SER:O	1:A:91:ARG:HD3	1.76	0.84
1:A:114:ILE:HD13	1:A:124:ALA:HB2	1.59	0.83
1:A:36:ARG:HG2	1:A:36:ARG:NH1	1.91	0.81
1:A:61:PRO:HG2	1:A:64:ASN:HD22	1.49	0.77
1:A:114:ILE:HD13	1:A:124:ALA:CB	2.16	0.76
1:A:28:ARG:O	1:A:32:ARG:HG3	1.89	0.70
1:A:72:ASN:H	1:A:87:HIS:CD2	2.05	0.70
1:A:105:LEU:HD23	1:A:108:LYS:CD	2.10	0.70
1:A:159:LEU:HD12	1:A:181:VAL:HG22	1.75	0.69
1:A:94:ASP:O	1:A:98:LYS:HG3	1.93	0.68
1:A:94:ASP:HB3	1:A:98:LYS:HZ3	1.60	0.67
1:A:145:ASP:OD1	1:A:146:THR:HG22	1.96	0.65
1:A:99:LEU:HD23	1:A:99:LEU:O	1.96	0.65
1:A:22:PHE:HZ	3:A:188:MXA:C7	2.10	0.65
1:A:94:ASP:HB3	1:A:98:LYS:NZ	2.13	0.64
1:A:156:TYR:CZ	1:A:184:LYS:HD2	2.34	0.62
2:A:187:NDP:H8A	2:A:187:NDP:H52A	1.81	0.62
1:A:102:GLN:CB	1:A:104:GLU:HG3	2.29	0.61
1:A:4:LEU:HD12	1:A:112:VAL:HG22	1.82	0.61
1:A:103:PRO:HA	1:A:106:ALA:HB2	1.81	0.60
3:A:188:MXA:HN42	3:A:188:MXA:C5M	2.14	0.59
1:A:102:GLN:HG3	1:A:104:GLU:OE2	2.02	0.59
1:A:139:MET:HE1	1:A:178:LYS:HD3	1.85	0.59
1:A:155:LYS:HD2	1:A:184:LYS:HE3	1.84	0.58
1:A:102:GLN:HB3	1:A:104:GLU:HG3	1.85	0.58
1:A:139:MET:CE	1:A:178:LYS:CE	2.71	0.58
1:A:57:TRP:O	1:A:65:ARG:HD2	2.04	0.57
1:A:116:GLY:HA3	2:A:187:NDP:H5N	1.88	0.55
1:A:139:MET:HE1	1:A:178:LYS:CD	2.37	0.55
1:A:30:GLU:O	1:A:33:TYR:HB3	2.07	0.55
1:A:16:ILE:O	2:A:187:NDP:H2N	2.08	0.54
1:A:99:LEU:CD2	1:A:105:LEU:HD12	2.38	0.53
1:A:61:PRO:HG2	1:A:64:ASN:ND2	2.22	0.53
1:A:124:ALA:O	1:A:127:HIS:HB3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:TYR:OH	2:A:187:NDP:H41N	2.12	0.50
1:A:36:ARG:CG	1:A:36:ARG:HH11	2.16	0.49
1:A:93:LEU:O	1:A:93:LEU:HG	2.13	0.49
1:A:43:VAL:HG11	1:A:46:LYS:HD2	1.95	0.48
1:A:130:HIS:HB2	1:A:185:ASN:ND2	2.29	0.48
1:A:10:VAL:HG22	1:A:14:MET:HA	1.95	0.48
1:A:102:GLN:HB2	1:A:104:GLU:HG3	1.94	0.47
1:A:17:GLY:HA2	1:A:23:PRO:HD3	1.95	0.46
1:A:83:PRO:HD2	1:A:86:ALA:HB3	1.97	0.46
1:A:17:GLY:HA2	1:A:21:ASP:O	2.17	0.45
1:A:91:ARG:HD2	4:A:234:HOH:O	2.17	0.45
1:A:131:LEU:HD23	1:A:156:TYR:OH	2.18	0.44
1:A:114:ILE:HG23	1:A:120:VAL:HG12	2.00	0.44
1:A:2:GLY:HA3	4:A:189:HOH:O	2.18	0.43
1:A:24:TRP:HD1	1:A:25:PRO:O	2.02	0.43
1:A:184:LYS:HE2	1:A:186:ASP:CG	2.39	0.43
1:A:24:TRP:HB3	1:A:142:PHE:CE2	2.54	0.43
1:A:145:ASP:OD1	1:A:145:ASP:C	2.56	0.43
1:A:22:PHE:CZ	3:A:188:MXA:C7	2.96	0.42
1:A:75:LEU:O	2:A:187:NDP:H1B	2.19	0.42
1:A:155:LYS:HB3	1:A:184:LYS:HE3	2.02	0.42
3:A:188:MXA:HN42	3:A:188:MXA:H5M2	1.83	0.41
1:A:99:LEU:C	1:A:99:LEU:HD23	2.40	0.41
3:A:188:MXA:H91	3:A:188:MXA:H5M1	1.71	0.41
1:A:99:LEU:HA	1:A:102:GLN:HG2	2.02	0.41
3:A:188:MXA:C5M	3:A:188:MXA:N4	2.81	0.41
1:A:156:TYR:CE1	1:A:184:LYS:HD2	2.56	0.41
1:A:24:TRP:HB2	1:A:25:PRO:HD2	2.03	0.40
1:A:139:MET:CE	1:A:178:LYS:CD	2.99	0.40
1:A:17:GLY:HA3	2:A:187:NDP:O3D	2.21	0.40
1:A:93:LEU:O	1:A:97:LEU:HG	2.20	0.40
1:A:28:ARG:CG	4:A:245:HOH:O	2.70	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	184/186 (99%)	175 (95%)	9 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	168/168 (100%)	155 (92%)	13 (8%)	15	18

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ARG
1	A	62	GLU
1	A	63	LYS
1	A	72	ASN
1	A	77	ARG
1	A	100	THR
1	A	103	PRO
1	A	104	GLU
1	A	119	SER
1	A	123	GLU
1	A	140	GLN
1	A	181	VAL
1	A	185	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN

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Mol	Chain	Res	Type
1	A	64	ASN
1	A	87	HIS
1	A	185	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 ⓘ

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	$\# Z > 2$
2	NDP	A	187	-	43,52,52	2.56	14 (32%)	49,80,80	2.40	15 (30%)
3	MXA	A	188	-	26,26,26	0.98	1 (3%)	34,37,37	1.61	7 (20%)

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	NDP	A	187	-	-	0/30/77/77	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MXA	A	188	-	-	0/8/8/8	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	187	NDP	O4B-C4B	-9.52	1.23	1.45
2	A	187	NDP	C5A-C4A	-3.76	1.32	1.40
2	A	187	NDP	P2B-O2X	-2.39	1.45	1.54
2	A	187	NDP	PN-O2N	-2.15	1.44	1.55
2	A	187	NDP	O2B-C2B	-2.06	1.36	1.44
2	A	187	NDP	O7N-C7N	-2.01	1.19	1.24
2	A	187	NDP	PA-O5B	2.11	1.68	1.59
2	A	187	NDP	C4A-N3A	2.33	1.39	1.35
3	A	188	MXA	O5'-C5'	2.33	1.42	1.37
2	A	187	NDP	C6A-C5A	2.64	1.56	1.42
2	A	187	NDP	O4D-C1D	2.93	1.49	1.42
2	A	187	NDP	C1D-N1N	2.94	1.54	1.46
2	A	187	NDP	O3B-C3B	3.72	1.51	1.43
2	A	187	NDP	C3B-C4B	4.21	1.64	1.53
2	A	187	NDP	P2B-O2B	7.86	1.73	1.59

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	188	MXA	C51-O5'-C5'	-5.35	105.80	117.50
2	A	187	NDP	O3B-C3B-C4B	-4.93	96.69	111.09
2	A	187	NDP	O5D-C5D-C4D	-4.15	94.29	109.00
2	A	187	NDP	O3D-C3D-C4D	-4.12	99.06	111.09
2	A	187	NDP	C5B-C4B-C3B	-3.65	101.37	115.29
2	A	187	NDP	C1D-N1N-C2N	-3.43	115.26	121.09
2	A	187	NDP	C3N-C2N-N1N	-3.42	118.12	123.08
2	A	187	NDP	O4B-C1B-C2B	-2.50	102.22	106.59
3	A	188	MXA	N2-C2-N3	-2.30	113.56	117.24
2	A	187	NDP	O3X-P2B-O2B	-2.18	96.09	106.00
3	A	188	MXA	O2'-C2'-C1'	-2.16	113.12	115.86
2	A	187	NDP	O2B-P2B-O1X	-2.14	100.88	109.26
3	A	188	MXA	O5'-C5'-C6'	-2.04	110.20	119.97
2	A	187	NDP	O3B-C3B-C2B	2.04	116.98	111.18
3	A	188	MXA	C21-O2'-C2'	2.42	121.03	117.54
3	A	188	MXA	O2'-C2'-C3'	2.43	128.45	124.37
2	A	187	NDP	N6A-C6A-N1A	2.53	123.78	118.77
2	A	187	NDP	C2D-C1D-N1N	2.60	120.02	113.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	188	MXA	C6'-C1'-C2'	2.95	121.52	118.30
2	A	187	NDP	N3A-C2A-N1A	4.36	132.65	128.86
2	A	187	NDP	O3X-P2B-O2X	5.70	130.61	107.61
2	A	187	NDP	C4B-O4B-C1B	7.76	118.03	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	187	NDP	6	0
3	A	188	MXA	6	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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.