



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:00 PM GMT

PDB ID : 1DLS
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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

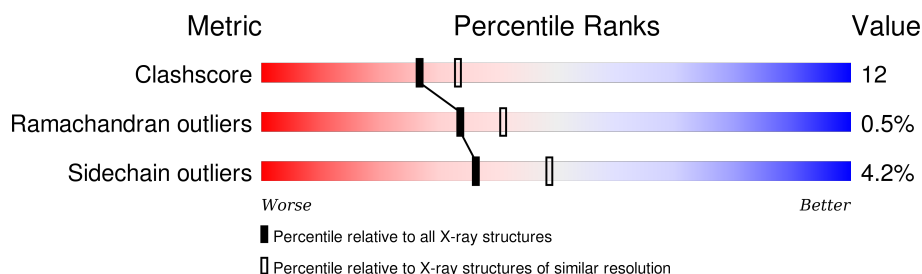
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (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	<div>49% 39% 11% •</div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1698 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
			1506	966	253	280	7			

There is a discrepancy between the modelled and reference sequences:

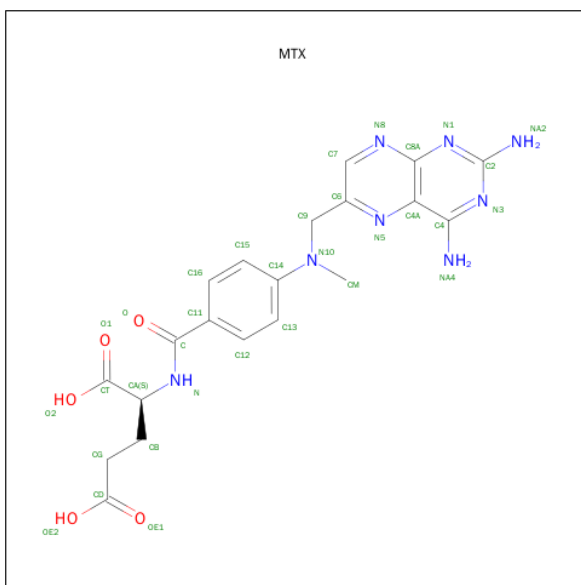
Chain	Residue	Modelled	Actual	Comment	Reference
A	22	TYR	LEU	CONFLICT	UNP P00374

- 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 METHOTREXATE (three-letter code: MTX) (formula: $C_{20}H_{22}N_8O_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			33	20	8	5		

- Molecule 4 is water.

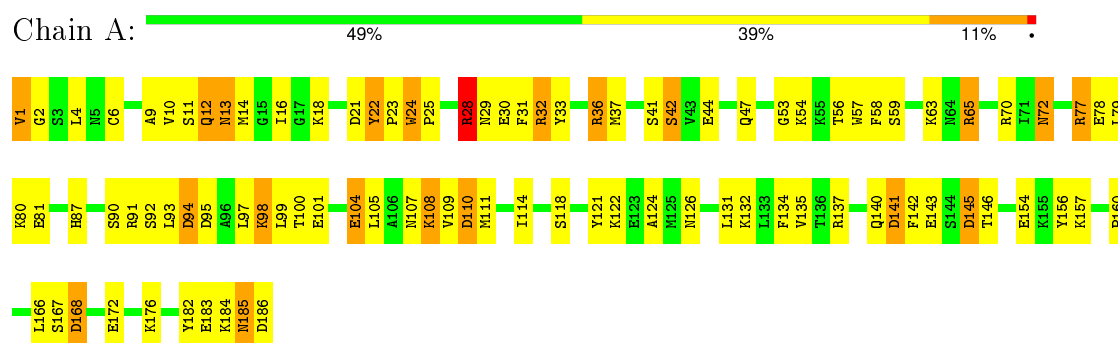
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	111	Total O 111 111	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: DIHYDROFOLATE REDUCTASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	87.50 Å 87.50 Å 76.73 Å 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}	0.156 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1698	wwPDB-VP
Average B, all atoms (Å ²)	34.0	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, MTX

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	1.39	5/1542 (0.3%)	2.73	95/2080 (4.6%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	172	GLU	CD-OE1	-5.86	1.19	1.25
1	A	78	GLU	CD-OE2	5.68	1.31	1.25
1	A	183	GLU	CD-OE1	-5.19	1.20	1.25
1	A	185	ASN	C-N	5.13	1.45	1.34
1	A	53	GLY	N-CA	5.10	1.53	1.46

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	ARG	NE-CZ-NH2	34.17	137.39	120.30
1	A	141	ASP	CB-CG-OD1	24.00	139.90	118.30
1	A	65	ARG	CD-NE-CZ	20.93	152.90	123.60
1	A	28	ARG	NE-CZ-NH1	-17.88	111.36	120.30
1	A	33	TYR	CB-CG-CD1	-14.32	112.41	121.00
1	A	154	GLU	OE1-CD-OE2	13.97	140.06	123.30
1	A	137	ARG	NE-CZ-NH1	-12.99	113.80	120.30
1	A	21	ASP	CB-CG-OD2	-12.63	106.93	118.30
1	A	65	ARG	NH1-CZ-NH2	-12.51	105.64	119.40
1	A	91	ARG	CD-NE-CZ	11.62	139.87	123.60
1	A	145	ASP	CB-CG-OD1	11.33	128.50	118.30
1	A	33	TYR	CB-CG-CD2	11.32	127.79	121.00
1	A	141	ASP	CB-CG-OD2	-11.07	108.34	118.30
1	A	77	ARG	NE-CZ-NH1	-10.88	114.86	120.30
1	A	182	TYR	CB-CG-CD2	-10.81	114.51	121.00
1	A	65	ARG	CG-CD-NE	10.35	133.53	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	ARG	NE-CZ-NH1	10.34	125.47	120.30
1	A	90	SER	N-CA-CB	10.09	125.63	110.50
1	A	137	ARG	NE-CZ-NH2	10.01	125.30	120.30
1	A	104	GLU	CA-CB-CG	9.16	133.54	113.40
1	A	81	GLU	OE1-CD-OE2	-8.99	112.51	123.30
1	A	28	ARG	CD-NE-CZ	-8.66	111.48	123.60
1	A	134	PHE	CB-CG-CD2	-8.58	114.79	120.80
1	A	121	TYR	CB-CG-CD2	-8.40	115.96	121.00
1	A	140	GLN	N-CA-CB	8.39	125.70	110.60
1	A	32	ARG	NE-CZ-NH1	8.23	124.41	120.30
1	A	32	ARG	NE-CZ-NH2	-8.06	116.27	120.30
1	A	94	ASP	CB-CG-OD1	8.05	125.55	118.30
1	A	154	GLU	CG-CD-OE2	-7.93	102.44	118.30
1	A	30	GLU	OE1-CD-OE2	-7.91	113.81	123.30
1	A	111	MET	CG-SD-CE	7.91	112.85	100.20
1	A	4	LEU	CB-CG-CD2	-7.89	97.59	111.00
1	A	37	MET	CG-SD-CE	-7.72	87.84	100.20
1	A	78	GLU	CG-CD-OE1	7.71	133.72	118.30
1	A	172	GLU	OE1-CD-OE2	7.63	132.46	123.30
1	A	78	GLU	OE1-CD-OE2	-7.56	114.22	123.30
1	A	18	LYS	CB-CG-CD	-7.50	92.10	111.60
1	A	58	PHE	CB-CG-CD1	-7.47	115.57	120.80
1	A	44	GLU	OE1-CD-OE2	7.24	131.98	123.30
1	A	172	GLU	CB-CA-C	7.22	124.84	110.40
1	A	143	GLU	OE1-CD-OE2	-7.07	114.82	123.30
1	A	59	SER	CB-CA-C	6.99	123.39	110.10
1	A	78	GLU	CA-CB-CG	6.96	128.71	113.40
1	A	63	LYS	CB-CA-C	-6.86	96.69	110.40
1	A	65	ARG	NE-CZ-NH1	-6.85	116.88	120.30
1	A	79	LEU	O-C-N	6.85	133.66	122.70
1	A	36	ARG	NE-CZ-NH2	6.81	123.71	120.30
1	A	168	ASP	CB-CG-OD1	6.79	124.41	118.30
1	A	156	TYR	CB-CG-CD2	-6.78	116.94	121.00
1	A	31	PHE	CB-CG-CD1	-6.76	116.07	120.80
1	A	12	GLN	CA-CB-CG	-6.72	98.61	113.40
1	A	36	ARG	CD-NE-CZ	-6.63	114.32	123.60
1	A	44	GLU	CB-CA-C	-6.59	97.22	110.40
1	A	95	ASP	CB-CG-OD2	-6.58	112.38	118.30
1	A	28	ARG	NH1-CZ-NH2	6.49	126.54	119.40
1	A	81	GLU	CG-CD-OE1	6.40	131.11	118.30
1	A	132	LYS	CD-CE-NZ	-6.39	96.99	111.70
1	A	29	ASN	CA-CB-CG	-6.37	99.40	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	135	VAL	O-C-N	6.36	132.87	122.70
1	A	24	TRP	CH2-CZ2-CE2	-6.32	111.08	117.40
1	A	118	SER	CB-CA-C	-6.27	98.19	110.10
1	A	41	SER	C-N-CA	6.25	137.31	121.70
1	A	22	TYR	CB-CG-CD1	-6.23	117.26	121.00
1	A	141	ASP	OD1-CG-OD2	-6.10	111.70	123.30
1	A	54	LYS	N-CA-CB	-6.08	99.65	110.60
1	A	91	ARG	O-C-N	6.08	132.43	122.70
1	A	110	ASP	CB-CA-C	6.07	122.54	110.40
1	A	36	ARG	CA-CB-CG	6.04	126.68	113.40
1	A	1	VAL	CA-C-N	-5.92	104.35	116.20
1	A	176	LYS	CB-CG-CD	-5.86	96.37	111.60
1	A	2	GLY	N-CA-C	-5.78	98.66	113.10
1	A	11	SER	O-C-N	-5.70	113.58	122.70
1	A	121	TYR	CZ-CE2-CD2	-5.68	114.69	119.80
1	A	98	LYS	O-C-N	-5.67	113.63	122.70
1	A	126	ASN	CB-CA-C	5.63	121.66	110.40
1	A	172	GLU	CG-CD-OE2	-5.61	107.08	118.30
1	A	131	LEU	CB-CG-CD2	5.60	120.52	111.00
1	A	1	VAL	CA-C-O	5.53	131.71	120.10
1	A	182	TYR	CD1-CE1-CZ	-5.46	114.89	119.80
1	A	53	GLY	CA-C-O	5.46	130.42	120.60
1	A	9	ALA	N-CA-CB	5.41	117.68	110.10
1	A	122	LYS	CA-CB-CG	-5.38	101.56	113.40
1	A	118	SER	O-C-N	5.31	131.19	122.70
1	A	6	CYS	N-CA-CB	5.31	120.15	110.60
1	A	53	GLY	N-CA-C	-5.30	99.85	113.10
1	A	185	ASN	N-CA-CB	5.26	120.08	110.60
1	A	108	LYS	CB-CA-C	-5.25	99.91	110.40
1	A	160	PRO	O-C-N	5.19	131.01	122.70
1	A	118	SER	CA-C-O	-5.13	109.33	120.10
1	A	56	THR	CA-CB-OG1	-5.12	98.26	109.00
1	A	2	GLY	O-C-N	5.08	130.83	122.70
1	A	157	LYS	CB-CA-C	-5.08	100.25	110.40
1	A	166	LEU	C-N-CA	5.06	134.34	121.70
1	A	92	SER	O-C-N	-5.01	114.69	122.70
1	A	12	GLN	CG-CD-OE1	-5.00	111.59	121.60

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	1506	0	1509	37	0
2	A	48	0	26	3	0
3	A	33	0	20	0	0
4	A	111	0	0	3	0
All	All	1698	0	1555	38	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 12.

All (38) 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:98:LYS:O	1:A:101:GLU:HG2	1.73	0.87
1:A:13:ASN:HD22	1:A:13:ASN:H	1.24	0.83
1:A:72:ASN:H	1:A:87:HIS:HD2	1.31	0.76
1:A:99:LEU:CD2	1:A:105:LEU:HD12	2.15	0.76
1:A:10:VAL:HG22	1:A:14:MET:HA	1.70	0.74
1:A:99:LEU:HD23	1:A:105:LEU:HD12	1.70	0.73
1:A:94:ASP:OD1	4:A:293:HOH:O	2.11	0.68
2:A:187:NDP:H52A	2:A:187:NDP:H8A	1.79	0.63
1:A:146:THR:HG21	2:A:187:NDP:H4D	1.79	0.63
1:A:72:ASN:H	1:A:87:HIS:CD2	2.15	0.62
1:A:114:ILE:HD13	1:A:124:ALA:HB2	1.81	0.62
1:A:28:ARG:O	1:A:32:ARG:HG3	2.02	0.60
1:A:13:ASN:ND2	1:A:13:ASN:H	1.93	0.59
1:A:184:LYS:NZ	1:A:186:ASP:OD1	2.36	0.58
1:A:13:ASN:HD21	1:A:142:PHE:H	1.53	0.57
1:A:93:LEU:O	1:A:97:LEU:HG	2.07	0.55
1:A:10:VAL:CG2	1:A:14:MET:HA	2.36	0.53
1:A:12:GLN:HB3	1:A:141:ASP:OD1	2.10	0.52
1:A:32:ARG:O	1:A:36:ARG:HG2	2.12	0.50
1:A:47:GLN:O	1:A:109:VAL:HA	2.12	0.49
1:A:23:PRO:HD2	1:A:24:TRP:CZ3	2.47	0.49
1:A:107:ASN:OD1	1:A:108:LYS:HG2	2.13	0.49
1:A:99:LEU:HD21	1:A:105:LEU:HD12	1.93	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:TRP:HB2	1:A:25:PRO:HD2	1.98	0.46
1:A:1:VAL:N	1:A:100:THR:CG2	2.78	0.46
1:A:145:ASP:OD1	1:A:146:THR:HG22	2.16	0.45
1:A:57:TRP:CE2	1:A:65:ARG:HD2	2.53	0.44
1:A:16:ILE:O	2:A:187:NDP:H2N	2.16	0.44
1:A:99:LEU:HD21	1:A:105:LEU:CD1	2.48	0.44
1:A:80:LYS:HA	1:A:80:LYS:HD3	1.58	0.43
1:A:98:LYS:HG3	4:A:292:HOH:O	2.17	0.43
1:A:1:VAL:H1	1:A:100:THR:HG21	1.83	0.43
1:A:114:ILE:HD13	1:A:124:ALA:CB	2.49	0.41
1:A:42:SER:HB2	1:A:110:ASP:OD2	2.21	0.41
1:A:28:ARG:HD3	1:A:28:ARG:HA	1.62	0.41
1:A:168:ASP:N	1:A:168:ASP:OD1	2.48	0.41
1:A:13:ASN:ND2	1:A:141:ASP:HA	2.36	0.40
1:A:77:ARG:NH2	4:A:206:HOH:O	2.54	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%)	179 (97%)	4 (2%)	1 (0%)	34	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	GLU

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	168/168 (100%)	161 (96%)	7 (4%)	36	49

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ASN
1	A	22	TYR
1	A	28	ARG
1	A	42	SER
1	A	72	ASN
1	A	167	SER
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	13	ASN
1	A	47	GLN
1	A	87	HIS
1	A	185	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

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	-	42,52,52	2.28	16 (38%)	55,80,80	2.35	20 (36%)
3	MTX	A	188	-	27,35,35	1.28	2 (7%)	30,49,49	1.75	8 (26%)

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
3	MTX	A	188	-	-	0/19/25/25	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	187	NDP	O4B-C4B	-6.53	1.29	1.45
2	A	187	NDP	C5A-C4A	-4.11	1.31	1.40
3	A	188	MTX	C8A-N8	-3.97	1.30	1.37
2	A	187	NDP	P2B-O3X	-3.15	1.43	1.54
2	A	187	NDP	PN-O2N	-3.06	1.41	1.54
2	A	187	NDP	C6N-C5N	-2.96	1.27	1.33
2	A	187	NDP	O7N-C7N	-2.60	1.17	1.24
2	A	187	NDP	PA-O1A	-2.31	1.42	1.51
2	A	187	NDP	O5D-C5D	-2.03	1.36	1.44
3	A	188	MTX	C2-NA2	-2.02	1.29	1.34
2	A	187	NDP	C4N-C5N	2.33	1.54	1.49
2	A	187	NDP	C1D-N1N	2.36	1.53	1.46
2	A	187	NDP	C5B-C4B	2.52	1.59	1.51
2	A	187	NDP	C3B-C4B	2.56	1.59	1.53
2	A	187	NDP	O3B-C3B	3.39	1.51	1.43
2	A	187	NDP	O4D-C1D	3.63	1.51	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	187	NDP	C3D-C4D	3.82	1.63	1.53
2	A	187	NDP	O4B-C1B	4.96	1.47	1.41

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	187	NDP	C3N-C2N-N1N	-5.19	115.71	123.14
3	A	188	MTX	N1-C2-N3	-4.87	120.03	127.44
2	A	187	NDP	O4D-C1D-N1N	-4.65	98.25	108.07
2	A	187	NDP	O3D-C3D-C4D	-4.50	97.55	111.05
2	A	187	NDP	O5B-C5B-C4B	-4.07	94.11	109.12
2	A	187	NDP	C5D-C4D-C3D	-3.87	99.84	115.21
2	A	187	NDP	C1D-N1N-C2N	-3.75	114.38	120.91
2	A	187	NDP	O3-PA-O5B	-3.38	93.96	102.94
2	A	187	NDP	C2A-N1A-C6A	-3.26	112.94	118.77
2	A	187	NDP	C5B-C4B-C3B	-3.06	103.06	115.21
3	A	188	MTX	C13-C12-C11	-2.77	117.55	120.76
2	A	187	NDP	O2B-C2B-C1B	-2.33	100.95	110.02
3	A	188	MTX	C4A-C8A-N1	-2.31	118.36	122.11
2	A	187	NDP	O3B-C3B-C4B	-2.28	104.23	111.05
2	A	187	NDP	C3B-C2B-C1B	-2.25	98.38	102.73
3	A	188	MTX	O-C-C11	-2.00	117.55	120.97
2	A	187	NDP	N6A-C6A-N1A	2.03	123.55	119.20
2	A	187	NDP	O4B-C4B-C3B	2.14	109.45	105.15
2	A	187	NDP	O3D-C3D-C2D	2.44	119.77	111.83
3	A	188	MTX	NA2-C2-N3	2.44	121.25	117.20
3	A	188	MTX	N8-C8A-N1	2.60	119.86	116.14
3	A	188	MTX	C11-C-N	2.64	121.63	116.93
2	A	187	NDP	C2D-C1D-N1N	2.81	120.92	113.34
3	A	188	MTX	C6-C9-N10	3.10	119.16	113.78
2	A	187	NDP	O7N-C7N-N7N	3.22	130.76	122.76
2	A	187	NDP	O2A-PA-O1A	3.31	130.45	112.53
2	A	187	NDP	N3A-C2A-N1A	4.38	132.25	128.89
2	A	187	NDP	PN-O3-PA	6.80	151.83	132.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	187	NDP	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.