



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:49 PM GMT

PDB ID : 4M7V
Title : Dihydrofolate reductase from Enterococcus faecalis complexed with NADP(H) and RAB-propyl
Authors : Bourne, C.R.
Deposited on : 2013-08-12
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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

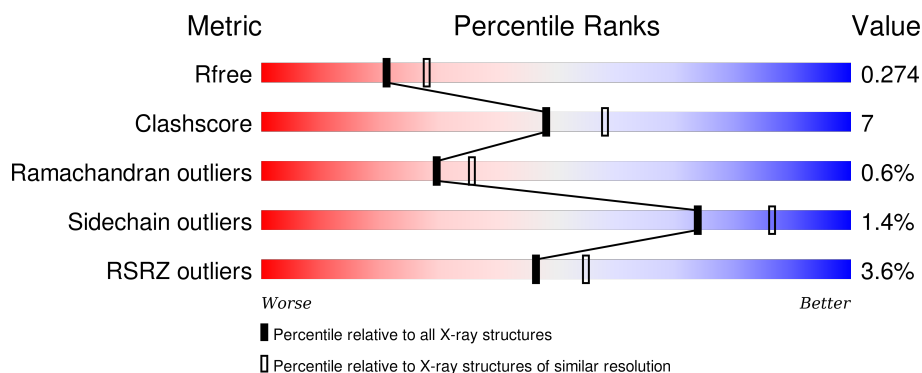
1 Overall quality at a glance

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))
R_{free}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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.

Mol	Chain	Length	Quality of chain
1	A	175	<div> <div>3%</div> <div>83%</div> <div>12%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	RAR	A	200[A]	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAP	A	201[A]	-	-	-	X
3	NAP	A	201[B]	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 1565 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.

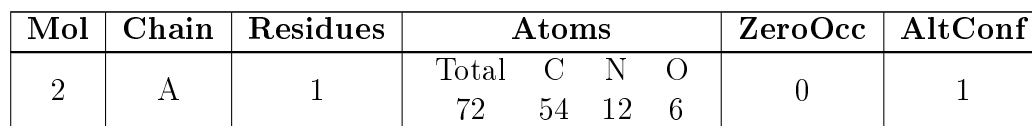
- Molecule 1 is a protein called Dihydrofolate reductase.

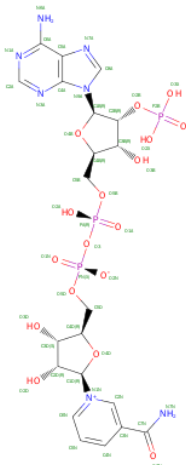
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1334	860	217	251	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	EXPRESSION TAG	UNP Q834R2
A	-6	GLU	-	EXPRESSION TAG	UNP Q834R2
A	-5	ASN	-	EXPRESSION TAG	UNP Q834R2
A	-4	LEU	-	EXPRESSION TAG	UNP Q834R2
A	-3	TYR	-	EXPRESSION TAG	UNP Q834R2
A	-2	PHE	-	EXPRESSION TAG	UNP Q834R2
A	-1	GLN	-	EXPRESSION TAG	UNP Q834R2
A	0	GLY	-	EXPRESSION TAG	UNP Q834R2
A	164	LEU	-	EXPRESSION TAG	UNP Q834R2
A	165	VAL	-	EXPRESSION TAG	UNP Q834R2
A	166	PRO	-	EXPRESSION TAG	UNP Q834R2
A	167	ARG	-	EXPRESSION TAG	UNP Q834R2

- Molecule 2 is 5-(3,4-DIMETHOXY-5-{(1E)-3-OXO-3-[(1S)-1-PROPYLPHTHALAZIN-2(1H)-YL]PROP-1-EN-1-YL}BENZYL)PYRIMIDINE-2,4-DIAMINE (three-letter code: RAR) (formula: C₂₇H₃₀N₆O₃).



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- The chemical structure of Naproxen is shown, featuring a naphthalene ring system. It has an amide group (-CONH₂) at position 1 and a propionic acid side chain (-CH₂-CH₂-COOH) at position 2. The structure is labeled with atom numbers and partial charges.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total 96	C 42	N 14	O 34	P 6	0	1

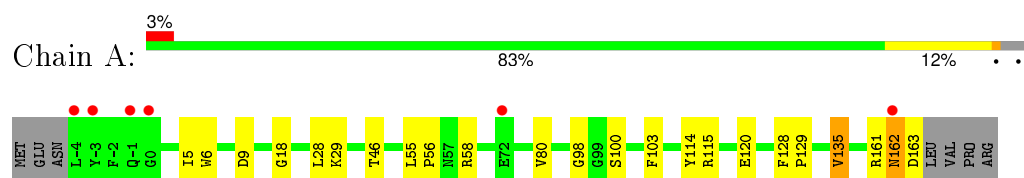
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	63	Total	O	0	0
			63	63		

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 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 ($\text{RSRZ} > 2$). 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: Dihydrofolate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	63.25Å 63.25Å 97.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.63 – 2.30 53.01 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.1 (40.63-2.30) 97.5 (53.01-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.69 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.208 , 0.266 0.222 , 0.274	Depositor DCC
R_{free} test set	460 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 9197 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	1565	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.44	0/1370	0.56	0/1866

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 [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1334	0	1285	20	0
2	A	72	0	59	4	0
3	A	96	0	44	5	0
4	A	63	0	0	1	0
All	All	1565	0	1388	21	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 7.

All (21) 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:28:LEU:HD22	2:A:200[A]:RAR:H15	1.75	0.69
1:A:18:GLY:HA2	3:A:201[A]:NAP:H3D	1.76	0.66
1:A:56:PRO:O	1:A:58:ARG:NH1	2.31	0.64
1:A:162:ASN:HD22	1:A:162:ASN:H	1.51	0.57
1:A:58:ARG:HH22	2:A:200[A]:RAR:C26	2.19	0.56
1:A:46:THR:OG1	3:A:201[A]:NAP:H6N	2.07	0.54
1:A:98:GLY:HA3	3:A:201[A]:NAP:C5N	2.39	0.53
1:A:80:VAL:HG23	3:A:201[A]:NAP:N6A	2.27	0.50
1:A:9:ASP:HB2	1:A:120:GLU:O	2.12	0.49
1:A:100:SER:N	3:A:201[B]:NAP:O2N	2.41	0.48
1:A:55:LEU:HB3	1:A:58:ARG:NH2	2.29	0.47
1:A:162:ASN:N	1:A:162:ASN:HD22	2.12	0.46
1:A:128:PHE:CD1	1:A:129:PRO:HD2	2.50	0.46
1:A:55:LEU:HD13	1:A:58:ARG:HH21	1.80	0.45
1:A:6:TRP:CZ2	1:A:115:ARG:HD3	2.52	0.45
1:A:29:LYS:HE2	4:A:336:HOH:O	2.15	0.45
1:A:135:VAL:HG11	1:A:161:ARG:NH2	2.32	0.44
1:A:58:ARG:NH2	2:A:200[A]:RAR:C26	2.81	0.43
1:A:5:ILE:HA	1:A:114:TYR:O	2.19	0.43
2:A:200[A]:RAR:O30	2:A:200[A]:RAR:H21	2.22	0.40
1:A:6:TRP:HB3	1:A:103:PHE:CZ	2.56	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	166/175 (95%)	161 (97%)	4 (2%)	1 (1%)	30 36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	ASN

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	142/151 (94%)	140 (99%)	2 (1%)	74 86

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	VAL
1	A	163	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	RAR	A	200[A]	-	38,39,39	3.79	15 (39%)	46,54,54	2.18	11 (23%)
2	RAR	A	200[B]	-	38,39,39	3.79	15 (39%)	46,54,54	2.56	15 (32%)
3	NAP	A	201[A]	-	42,52,52	3.33	14 (33%)	54,80,80	3.06	20 (37%)
3	NAP	A	201[B]	-	42,52,52	3.37	17 (40%)	54,80,80	2.89	17 (31%)

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	RAR	A	200[A]	-	-	0/19/33/33	0/3/4/4
2	RAR	A	200[B]	-	-	0/19/33/33	0/3/4/4
3	NAP	A	201[A]	-	-	0/27/67/67	0/5/5/5
3	NAP	A	201[B]	-	-	0/27/67/67	0/5/5/5

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	201[A]	NAP	C3B-C2B	-4.92	1.41	1.53
2	A	200[B]	RAR	C25-C24	-3.73	1.36	1.40
2	A	200[B]	RAR	C03-C02	-3.69	1.34	1.42
2	A	200[A]	RAR	C25-C24	-3.69	1.36	1.40
2	A	200[A]	RAR	C21-C20	-3.58	1.49	1.53
2	A	200[A]	RAR	C03-C02	-3.48	1.34	1.42
2	A	200[B]	RAR	C21-C20	-3.11	1.50	1.53
3	A	201[B]	NAP	O3B-C3B	-2.43	1.37	1.43
3	A	201[B]	NAP	C3B-C2B	-2.42	1.47	1.53
2	A	200[A]	RAR	C13-C10	-2.39	1.37	1.41
3	A	201[A]	NAP	O3B-C3B	-2.32	1.37	1.43
3	A	201[B]	NAP	O2B-C2B	-2.21	1.37	1.44
3	A	201[B]	NAP	C5A-C4A	-2.18	1.35	1.40
3	A	201[A]	NAP	C5A-C4A	-2.10	1.35	1.40
3	A	201[B]	NAP	O2D-C2D	-2.07	1.38	1.43
2	A	200[A]	RAR	C27-C26	2.10	1.43	1.38
3	A	201[B]	NAP	O4B-C4B	2.15	1.50	1.45
2	A	200[B]	RAR	C13-C14	2.16	1.50	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	200[B]	RAR	C27-C26	2.24	1.43	1.38
2	A	200[B]	RAR	C31-C05	2.59	1.43	1.39
2	A	200[A]	RAR	C31-C05	2.75	1.44	1.39
3	A	201[B]	NAP	C6A-N6A	2.93	1.44	1.34
3	A	201[A]	NAP	C6A-N6A	2.94	1.44	1.34
3	A	201[A]	NAP	C6N-N1N	3.12	1.43	1.35
3	A	201[B]	NAP	C6N-N1N	3.50	1.44	1.35
2	A	200[B]	RAR	C02-N01	3.70	1.43	1.34
2	A	200[A]	RAR	C02-N01	3.71	1.43	1.34
2	A	200[A]	RAR	C16-N17	3.78	1.43	1.35
2	A	200[A]	RAR	C07-C10	3.86	1.49	1.40
3	A	201[A]	NAP	O4D-C1D	4.02	1.46	1.41
3	A	201[B]	NAP	C2A-N3A	4.05	1.39	1.32
2	A	200[B]	RAR	C07-C10	4.24	1.49	1.40
3	A	201[A]	NAP	C2A-N3A	4.25	1.39	1.32
3	A	201[B]	NAP	O4D-C1D	4.27	1.46	1.41
2	A	200[A]	RAR	C34-N35	4.28	1.42	1.34
3	A	201[B]	NAP	C8A-N7A	4.29	1.42	1.34
2	A	200[B]	RAR	C34-N35	4.31	1.42	1.34
2	A	200[B]	RAR	C16-N17	4.32	1.44	1.35
3	A	201[A]	NAP	C8A-N7A	4.40	1.43	1.34
2	A	200[B]	RAR	C25-C19	4.47	1.51	1.44
3	A	201[B]	NAP	C7N-N7N	4.65	1.42	1.33
2	A	200[A]	RAR	C25-C19	4.77	1.52	1.44
2	A	200[B]	RAR	C02-N36	5.02	1.42	1.35
2	A	200[A]	RAR	C02-N36	5.22	1.43	1.35
3	A	201[A]	NAP	C7N-N7N	5.38	1.43	1.33
3	A	201[A]	NAP	O4B-C1B	5.83	1.48	1.41
3	A	201[B]	NAP	O4B-C1B	5.83	1.48	1.41
3	A	201[A]	NAP	C6N-C5N	6.33	1.52	1.38
2	A	200[B]	RAR	N17-N18	6.57	1.47	1.37
2	A	200[A]	RAR	N17-N18	6.59	1.47	1.37
2	A	200[B]	RAR	C29-C24	6.69	1.48	1.39
2	A	200[A]	RAR	C29-C24	6.88	1.49	1.39
3	A	201[B]	NAP	C6N-C5N	6.94	1.54	1.38
3	A	201[A]	NAP	C5N-C4N	6.99	1.53	1.38
3	A	201[B]	NAP	C5N-C4N	8.01	1.55	1.38
3	A	201[A]	NAP	C4N-C3N	8.60	1.54	1.39
3	A	201[B]	NAP	C4N-C3N	9.21	1.55	1.39
3	A	201[B]	NAP	C2N-C3N	9.50	1.53	1.39
3	A	201[A]	NAP	C2N-C3N	10.20	1.54	1.39
2	A	200[B]	RAR	C19-N18	16.14	1.47	1.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	200[A]	RAR	C19-N18	16.26	1.47	1.29

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	201[B]	NAP	N3A-C2A-N1A	-11.68	119.95	128.89
3	A	201[A]	NAP	N3A-C2A-N1A	-11.37	120.19	128.89
2	A	200[B]	RAR	C25-C19-N18	-11.30	111.67	125.69
2	A	200[A]	RAR	C25-C19-N18	-9.33	114.12	125.69
3	A	201[A]	NAP	C5N-C4N-C3N	-8.44	109.72	120.33
3	A	201[A]	NAP	C4B-O4B-C1B	-8.03	100.89	109.72
3	A	201[B]	NAP	C4B-O4B-C1B	-7.19	101.82	109.72
2	A	200[A]	RAR	N33-C34-N36	-5.50	120.04	125.78
2	A	200[B]	RAR	N33-C34-N36	-5.47	120.07	125.78
3	A	201[B]	NAP	C5N-C4N-C3N	-5.12	113.90	120.33
3	A	201[A]	NAP	PN-O3-PA	-3.58	122.67	132.73
2	A	200[B]	RAR	C14-C15-C16	-3.36	115.84	120.79
2	A	200[B]	RAR	C04-C03-C02	-3.12	118.22	122.21
2	A	200[B]	RAR	O08-C07-C06	-3.11	118.89	124.21
2	A	200[A]	RAR	C13-C14-C15	-3.08	120.73	126.87
3	A	201[B]	NAP	PN-O3-PA	-2.74	125.04	132.73
2	A	200[A]	RAR	C04-C03-C02	-2.55	118.96	122.21
3	A	201[B]	NAP	C3B-C2B-C1B	-2.47	97.96	102.73
2	A	200[B]	RAR	O30-C16-C15	-2.44	117.40	122.15
3	A	201[A]	NAP	O7N-C7N-C3N	-2.37	117.00	119.59
2	A	200[B]	RAR	C03-C32-N33	-2.23	119.99	123.86
2	A	200[A]	RAR	O08-C07-C06	-2.23	120.41	124.21
3	A	201[A]	NAP	O4B-C1B-C2B	-2.13	102.75	106.60
3	A	201[B]	NAP	O4B-C1B-C2B	-2.13	102.75	106.60
3	A	201[B]	NAP	C4D-O4D-C1D	-2.09	107.43	109.72
3	A	201[B]	NAP	O5D-C5D-C4D	2.01	116.54	109.12
2	A	200[B]	RAR	C12-O11-C10	2.12	120.41	114.82
3	A	201[A]	NAP	O2B-C2B-C1B	2.13	118.31	110.02
3	A	201[A]	NAP	O4B-C4B-C3B	2.16	109.50	105.15
3	A	201[B]	NAP	O2B-C2B-C1B	2.20	118.61	110.02
3	A	201[A]	NAP	O5D-C5D-C4D	2.21	117.26	109.12
2	A	200[B]	RAR	N01-C02-N36	2.29	120.27	116.95
3	A	201[A]	NAP	O4D-C4D-C5D	2.32	117.61	109.32
3	A	201[B]	NAP	O2D-C2D-C3D	2.33	119.40	111.83
3	A	201[A]	NAP	O2B-C2B-C3B	2.44	120.99	111.51
3	A	201[A]	NAP	O3-PN-O5D	2.46	109.45	102.94
2	A	200[A]	RAR	C32-C03-C02	2.46	119.25	115.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	200[A]	RAR	O08-C07-C10	2.49	119.75	115.26
3	A	201[B]	NAP	O3B-C3B-C2B	2.54	118.49	111.16
3	A	201[B]	NAP	O4D-C4D-C5D	2.56	118.46	109.32
2	A	200[B]	RAR	C34-N36-C02	2.66	120.06	117.04
3	A	201[B]	NAP	O3D-C3D-C4D	2.78	119.40	111.05
2	A	200[B]	RAR	C32-C03-C02	2.80	119.74	115.72
2	A	200[A]	RAR	N35-C34-N33	2.81	119.92	117.39
2	A	200[A]	RAR	C34-N36-C02	2.84	120.28	117.04
3	A	201[A]	NAP	C2N-C3N-C4N	2.85	121.47	118.29
2	A	200[A]	RAR	C32-N33-C34	2.86	119.99	116.05
2	A	200[A]	RAR	C24-C25-C19	2.97	120.45	117.68
2	A	200[B]	RAR	C32-N33-C34	2.98	120.17	116.05
3	A	201[B]	NAP	C2N-C3N-C4N	3.02	121.65	118.29
2	A	200[B]	RAR	N35-C34-N33	3.15	120.23	117.39
3	A	201[A]	NAP	O3D-C3D-C4D	3.23	120.75	111.05
2	A	200[B]	RAR	O08-C07-C10	3.30	121.21	115.26
3	A	201[A]	NAP	O3B-C3B-C2B	3.32	120.74	111.16
3	A	201[A]	NAP	O2D-C2D-C3D	3.32	122.63	111.83
3	A	201[B]	NAP	O3B-C3B-C4B	3.33	121.03	111.05
3	A	201[A]	NAP	O3B-C3B-C4B	3.33	121.05	111.05
3	A	201[A]	NAP	O3D-C3D-C2D	3.42	122.96	111.83
2	A	200[B]	RAR	C15-C16-N17	4.73	122.77	117.67
3	A	201[A]	NAP	O4D-C1D-N1N	5.92	114.63	108.13
3	A	201[B]	NAP	O4B-C1B-N9A	6.03	120.71	108.10
3	A	201[A]	NAP	O4B-C1B-N9A	6.77	122.28	108.10
3	A	201[B]	NAP	O4D-C1D-N1N	8.90	117.91	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	200[A]	RAR	4	0
3	A	201[A]	NAP	4	0
3	A	201[B]	NAP	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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, 95th 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 ²)	Q<0.9
1	A	168/175 (96%)	0.14	6 (3%) 46 55	27, 35, 55, 70	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	-3	TYR	9.0
1	A	-1	GLN	2.5
1	A	0	GLY	2.4
1	A	72	GLU	2.2
1	A	162	ASN	2.1
1	A	-4	LEU	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, 95th 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(\AA^2)	Q<0.9
2	RAR	A	200[A]	36/36	0.88	0.23	2.90	27,34,45,47	36
3	NAP	A	201[A]	48/48	0.92	0.21	2.90	26,29,31,33	48
3	NAP	A	201[B]	48/48	0.92	0.21	2.51	27,30,38,42	48
2	RAR	A	200[B]	36/36	0.88	0.23	1.61	27,34,43,44	36

6.5 Other polymers [i](#)

There are no such residues in this entry.