



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 07:19 PM GMT

PDB ID : 2BEL  
Title : STRUCTURE OF HUMAN 11-BETA-HYDROXYSTEROIDDEHYDROGENASE IN COMPLEX WITH NADP AND CARBENOXOLONE  
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Deposited on : 2004-11-25  
Resolution : 2.11 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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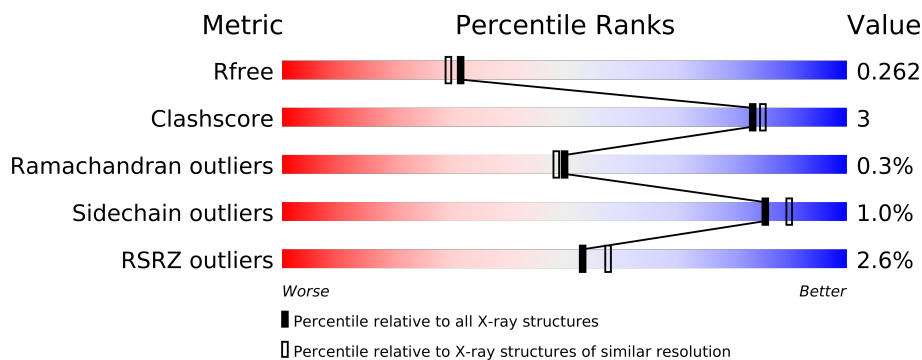
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.11 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))
$R_{free}$	66092	3409 (2.14-2.10)
Clashscore	79885	4090 (2.14-2.10)
Ramachandran outliers	78287	4048 (2.14-2.10)
Sidechain outliers	78261	4049 (2.14-2.10)
RSRZ outliers	66119	3410 (2.14-2.10)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	283	
1	B	283	
1	C	283	
1	D	283	

## 2 Entry composition i

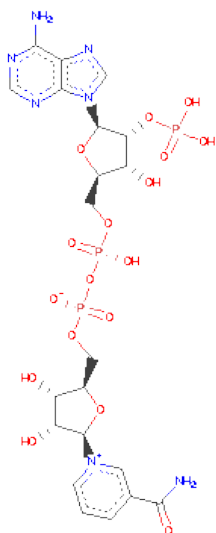
There are 5 unique types of molecules in this entry. The entry contains 8071 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CORTICOSTEROID 11-BETA-DEHYDROGENASE ISOZYME 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1809	1158	302	334	15			
1	B	251	Total	C	N	O	S	0	3	0
			1874	1202	312	345	15			
1	C	254	Total	C	N	O	S	0	0	0
			1868	1197	312	345	14			
1	D	253	Total	C	N	O	S	0	2	0
			1904	1215	320	352	17			

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDEPHOSPHATE (three-letter code: NAP) (formula:  $C_{21}H_{28}N_7O_{17}P_3$ ).



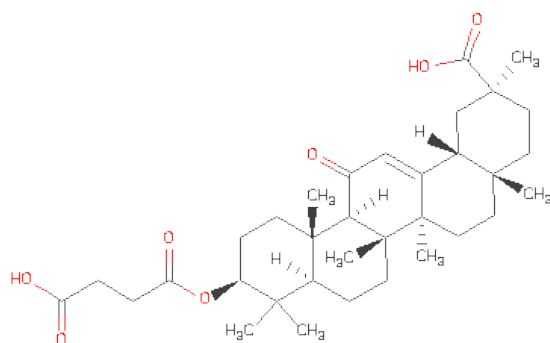
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

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

- Molecule 3 is CARBENOXOLONE (three-letter code: CBO) (formula:  $C_{34}H_{50}O_7$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			41	34	7		
3	B	1	Total	C	O	0	0
			37	32	5		
3	C	1	Total	C	O	0	0
			37	32	5		
3	D	1	Total	C	O	0	0
			37	32	5		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	A	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	1	Total 1	Cl 1	0	0
4	C	1	Total 1	Cl 1	0	0

- Molecule 5 is water.

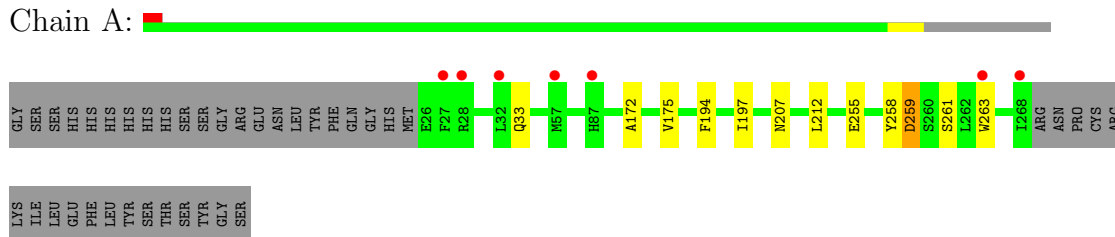
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	52	Total 52	O 52	0	0
5	B	80	Total 80	O 80	0	0
5	C	54	Total 54	O 54	0	0
5	D	82	Total 82	O 82	0	0

### 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.

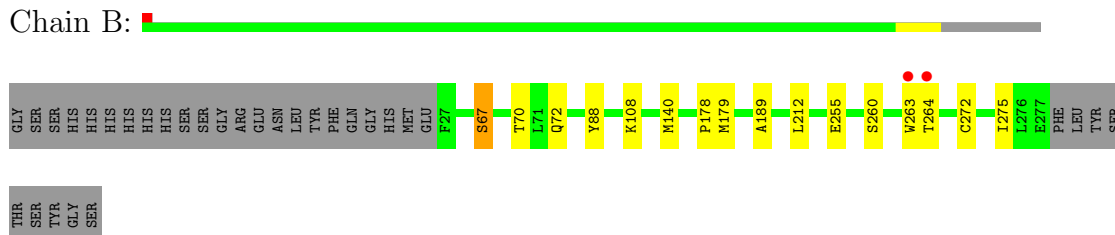
#### • Molecule 1: CORTICOSTEROID 11-BETA-DEHYDROGENASEISOZYME 1

Chain A:



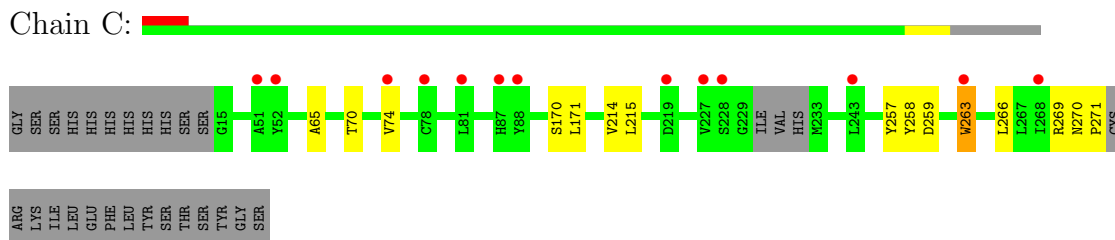
#### • Molecule 1: CORTICOSTEROID 11-BETA-DEHYDROGENASEISOZYME 1

Chain B:



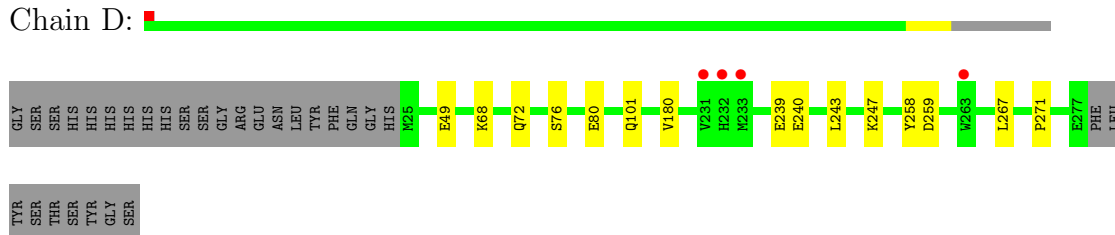
#### • Molecule 1: CORTICOSTEROID 11-BETA-DEHYDROGENASEISOZYME 1

Chain C:



#### • Molecule 1: CORTICOSTEROID 11-BETA-DEHYDROGENASEISOZYME 1

Chain D:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	160.16Å 112.98Å 66.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.52 – 2.11 53.87 – 2.11	Depositor EDS
% Data completeness (in resolution range)	98.0 (56.52-2.11) 98.0 (53.87-2.11)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.41 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.188 , 0.246 0.201 , 0.262	Depositor DCC
$R_{free}$ test set	1953 reflections (2.92%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.9	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 68774 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8071	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.15% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.66	0/1839	0.70	0/2489
1	B	0.74	0/1915	0.72	0/2592
1	C	0.68	0/1899	0.70	0/2572
1	D	0.78	0/1943	0.76	0/2625
All	All	0.72	0/7596	0.72	0/10278

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1809	0	1813	5	0
1	B	1874	0	1891	13	0
1	C	1868	0	1849	8	0
1	D	1904	0	1920	7	0
2	A	48	0	25	0	0
2	B	48	0	25	0	0
2	C	48	0	25	0	0
2	D	48	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	41	0	48	3	0
3	B	37	0	44	3	0
3	C	37	0	44	6	0
3	D	37	0	44	4	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	52	0	0	0	0
5	B	80	0	0	1	0
5	C	54	0	0	0	0
5	D	82	0	0	0	0
All	All	8071	0	7753	47	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (47) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:B:1279:CBO:H263	3:B:1279:CBO:H191	1.65	0.79
3:D:1279:CBO:H192	3:D:1279:CBO:H243	1.72	0.69
1:B:67[B]:SER:HB3	1:B:70[B]:THR:HG22	1.73	0.69
3:D:1279:CBO:H263	3:D:1279:CBO:H191	1.75	0.68
1:C:258:TYR:HA	1:C:259:ASP:C	2.14	0.67
1:C:214:VAL:HG11	1:C:266:LEU:HD11	1.81	0.63
1:C:70:THR:O	1:C:74:VAL:HG23	2.02	0.59
3:C:1273:CBO:H191	3:C:1273:CBO:H263	1.85	0.57
1:B:140:MET:HE1	1:B:189:ALA:CB	2.35	0.57
1:B:140:MET:HE1	1:B:189:ALA:HB2	1.87	0.56
3:C:1273:CBO:H243	3:C:1273:CBO:H192	1.88	0.55
3:A:1270:CBO:H263	3:A:1270:CBO:H191	1.89	0.54
3:A:1270:CBO:H243	3:A:1270:CBO:H192	1.91	0.53
3:B:1279:CBO:H192	3:B:1279:CBO:H243	1.92	0.52
1:B:67[B]:SER:HB3	1:B:70[B]:THR:CG2	2.37	0.52
3:D:1279:CBO:H271	3:D:1279:CBO:H202	1.91	0.52
3:B:1279:CBO:H271	3:B:1279:CBO:H202	1.94	0.49
1:B:272:CYS:HA	1:B:275:ILE:HG22	1.94	0.48
1:B:260:SER:O	1:B:263:TRP:HD1	1.97	0.47
1:D:68:LYS:O	1:D:72:GLN:HG3	2.15	0.47
1:B:67[B]:SER:CB	1:B:70[B]:THR:HG22	2.43	0.47
1:B:140:MET:CE	1:B:189:ALA:CB	2.93	0.47
3:A:1270:CBO:H202	3:A:1270:CBO:H271	1.96	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:171:LEU:HD23	3:C:1273:CBO:H22	1.97	0.46
1:A:212:LEU:O	1:A:255:GLU:HA	2.16	0.46
1:A:258:TYR:HA	1:A:259:ASP:C	2.37	0.45
1:D:243:LEU:O	1:D:247:LYS:HG3	2.17	0.45
1:D:76:SER:O	1:D:80:GLU:HG3	2.17	0.45
1:B:70[B]:THR:HG21	5:B:2080:HOH:O	2.16	0.45
1:A:194:PHE:HA	1:A:197:ILE:HG12	1.99	0.45
1:C:170:SER:OG	3:C:1273:CBO:O11	2.35	0.44
1:C:215:LEU:HA	1:C:258:TYR:O	2.18	0.44
1:B:72:GLN:HG2	1:B:88:TYR:CE2	2.52	0.44
3:C:1273:CBO:H202	3:C:1273:CBO:H271	2.01	0.42
1:B:212:LEU:O	1:B:255:GLU:HA	2.19	0.42
3:C:1273:CBO:H261	3:C:1273:CBO:C12	2.49	0.42
1:A:172:ALA:HA	1:A:175:VAL:O	2.19	0.42
1:A:259:ASP:OD2	1:A:261:SER:N	2.52	0.42
1:C:270:ASN:N	1:C:271:PRO:HD2	2.34	0.42
1:D:49:GLU:OE1	1:D:239:GLU:HG3	2.20	0.41
3:D:1279:CBO:H11	3:D:1279:CBO:O11	2.21	0.41
1:B:263:TRP:CZ2	1:B:264:THR:HG22	2.55	0.41
1:D:267:LEU:O	1:D:271:PRO:HG2	2.21	0.41
1:C:257:TYR:CD1	1:C:263:TRP:HB2	2.56	0.41
1:B:178:PRO:C	1:B:179:MET:HG2	2.42	0.40
1:D:258:TYR:HA	1:D:259:ASP:C	2.42	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	241/283 (85%)	230 (95%)	11 (5%)	0	100	100
1	B	252/283 (89%)	241 (96%)	11 (4%)	0	100	100
1	C	250/283 (88%)	234 (94%)	14 (6%)	2 (1%)	27	19
1	D	253/283 (89%)	239 (94%)	13 (5%)	1 (0%)	43	39

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	996/1132 (88%)	944 (95%)	49 (5%)	3 (0%)	50 48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	269	ARG
1	C	65	ALA
1	D	180	VAL

### 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	187/236 (79%)	183 (98%)	4 (2%)	66 69
1	B	196/236 (83%)	193 (98%)	3 (2%)	76 81
1	C	189/236 (80%)	188 (100%)	1 (0%)	94 96
1	D	201/236 (85%)	200 (100%)	1 (0%)	94 96
All	All	773/944 (82%)	764 (99%)	9 (1%)	85 87

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	207	ASN
1	A	259	ASP
1	A	263	TRP
1	B	67[A]	SER
1	B	67[B]	SER
1	B	108	LYS
1	C	263	TRP
1	D	101	GLN

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	234	GLN
1	B	127	ASN
1	C	123	ASN
1	C	160	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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, 4 are monoatomic - leaving 8 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	NAP	A	1269	-	52,52,52	1.35	3 (5%)	80,80,80	1.62	6 (7%)
3	CBO	A	1270	-	45,45,45	0.94	2 (4%)	76,76,76	1.15	6 (7%)
2	NAP	B	1278	-	52,52,52	1.31	4 (7%)	80,80,80	1.74	11 (13%)
3	CBO	B	1279	-	40,41,45	1.12	1 (2%)	69,71,76	1.34	7 (10%)
2	NAP	C	1272	-	52,52,52	1.21	3 (5%)	80,80,80	1.77	9 (11%)
3	CBO	C	1273	-	40,41,45	1.08	1 (2%)	69,71,76	1.15	4 (5%)
2	NAP	D	1278	-	52,52,52	1.22	3 (5%)	80,80,80	1.75	10 (12%)
3	CBO	D	1279	-	40,41,45	1.04	1 (2%)	69,71,76	1.16	5 (7%)

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	NAP	A	1269	-	-	0/35/67/67	0/3/5/5
3	CBO	A	1270	-	-	0/15/109/109	0/0/5/5
2	NAP	B	1278	-	-	0/35/67/67	0/3/5/5
3	CBO	B	1279	-	-	0/10/104/109	0/0/5/5
2	NAP	C	1272	-	-	0/35/67/67	0/3/5/5
3	CBO	C	1273	-	-	0/10/104/109	0/0/5/5
2	NAP	D	1278	-	-	0/35/67/67	0/3/5/5
3	CBO	D	1279	-	-	0/10/104/109	0/0/5/5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1269	NAP	O7N-C7N	7.03	1.40	1.24
2	B	1278	NAP	O7N-C7N	6.36	1.39	1.24
2	C	1272	NAP	O7N-C7N	6.10	1.38	1.24
2	D	1278	NAP	O7N-C7N	5.90	1.38	1.24
3	C	1273	CBO	O3-C29	5.54	1.46	1.34
3	B	1279	CBO	O3-C29	5.20	1.46	1.34
3	D	1279	CBO	O3-C29	4.95	1.45	1.34
3	A	1270	CBO	O3-C29	3.87	1.46	1.34
2	D	1278	NAP	C2A-N3A	3.12	1.38	1.32
2	A	1269	NAP	C2A-N3A	2.98	1.38	1.32
2	B	1278	NAP	C2A-N3A	2.72	1.37	1.32
2	B	1278	NAP	C2N-N1N	2.53	1.38	1.35
2	C	1272	NAP	C2A-N3A	2.50	1.37	1.32
2	A	1269	NAP	C2A-N1A	2.46	1.38	1.33
2	D	1278	NAP	C2A-N1A	2.41	1.38	1.33
3	A	1270	CBO	O3-C3	-2.26	1.42	1.46
2	C	1272	NAP	C2A-N1A	2.19	1.38	1.33
2	B	1278	NAP	PA-O3	2.10	1.63	1.59

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1278	NAP	N3A-C2A-N1A	-10.58	119.86	128.71
2	C	1272	NAP	N3A-C2A-N1A	-10.47	119.95	128.71
2	D	1278	NAP	N3A-C2A-N1A	-10.46	119.96	128.71
2	A	1269	NAP	N3A-C2A-N1A	-8.76	121.38	128.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1269	NAP	O4D-C1D-N1N	5.73	113.81	107.95
2	C	1272	NAP	O4D-C1D-N1N	5.17	113.24	107.95
2	C	1272	NAP	O4B-C1B-N9A	4.78	112.88	108.44
2	A	1269	NAP	O4B-C1B-N9A	4.58	112.70	108.44
2	D	1278	NAP	C3N-C7N-N7N	4.53	122.93	117.77
3	B	1279	CBO	C8-C9-C10	-4.03	115.51	118.02
2	C	1272	NAP	C3N-C7N-N7N	3.79	122.09	117.77
2	B	1278	NAP	C3N-C7N-N7N	3.52	121.78	117.77
3	B	1279	CBO	C20-C18-C17	-3.22	110.26	112.94
3	A	1270	CBO	C23-C17-C18	3.14	113.41	108.95
3	D	1279	CBO	C27-C14-C13	-3.13	103.76	106.97
2	B	1278	NAP	C4A-C5A-N7A	-3.12	106.85	109.52
2	C	1272	NAP	N3A-C4A-N9A	3.12	131.07	125.43
2	D	1278	NAP	N3A-C4A-N9A	3.08	130.99	125.43
3	C	1273	CBO	C7-C8-C9	3.02	113.16	109.61
2	A	1269	NAP	N3A-C4A-N9A	3.01	130.86	125.43
2	D	1278	NAP	O4D-C1D-N1N	2.88	110.90	107.95
2	C	1272	NAP	O4B-C1B-C2B	-2.84	104.30	106.95
2	B	1278	NAP	C2A-N1A-C6A	2.76	123.76	118.77
2	D	1278	NAP	P2B-O2B-C2B	2.68	127.59	121.96
2	A	1269	NAP	C2N-C3N-C4N	2.64	121.31	118.31
3	A	1270	CBO	C28-C17-C23	-2.64	104.37	108.86
2	B	1278	NAP	C1B-N9A-C4A	-2.61	122.12	126.64
2	C	1272	NAP	O7N-C7N-C3N	-2.56	116.69	119.58
2	B	1278	NAP	O2A-PA-O3	2.55	117.21	105.14
3	D	1279	CBO	C20-C18-C13	-2.53	107.13	111.53
3	B	1279	CBO	O3-C3-C4	2.53	112.20	107.62
3	B	1279	CBO	C26-C8-C7	-2.50	103.50	108.02
3	C	1273	CBO	C28-C17-C16	2.50	113.11	108.86
2	D	1278	NAP	O4B-C1B-N9A	2.49	110.75	108.44
2	A	1269	NAP	C2D-C1D-N1N	-2.48	109.65	113.86
2	D	1278	NAP	C4D-O4D-C1D	-2.44	107.10	109.75
3	A	1270	CBO	C28-C17-C16	2.43	113.00	108.86
2	B	1278	NAP	N7A-C8A-N9A	-2.37	107.64	114.36
2	B	1278	NAP	O4B-C1B-N9A	2.28	110.56	108.44
3	B	1279	CBO	C9-C11-C12	2.28	119.26	116.75
2	B	1278	NAP	N3A-C4A-N9A	2.26	129.52	125.43
3	A	1270	CBO	C26-C8-C14	-2.25	107.30	109.96
3	D	1279	CBO	C34-C21-C20	2.25	113.89	109.70
2	B	1278	NAP	C8A-N9A-C4A	2.19	108.57	106.90
2	D	1278	NAP	O7N-C7N-N7N	-2.19	119.42	122.59
2	B	1278	NAP	C4D-O4D-C1D	-2.19	107.38	109.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1278	NAP	C2D-C1D-N1N	-2.17	110.18	113.86
2	C	1272	NAP	C5A-C4A-N3A	-2.11	121.11	125.70
3	C	1273	CBO	C27-C14-C15	-2.09	104.18	107.92
3	C	1273	CBO	C20-C18-C13	-2.08	107.91	111.53
2	D	1278	NAP	O2A-PA-O3	2.06	114.91	105.14
3	B	1279	CBO	C26-C8-C14	2.04	112.38	109.96
3	B	1279	CBO	C2-C3-C4	-2.04	111.02	114.51
2	C	1272	NAP	C2N-C3N-C4N	2.04	120.62	118.31
3	D	1279	CBO	C25-C4-C3	-2.04	104.62	109.39
3	D	1279	CBO	C16-C15-C14	2.03	117.04	113.80
3	A	1270	CBO	C7-C8-C14	2.00	112.65	110.29
3	A	1270	CBO	O32-C32-C31	-2.00	116.15	123.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	243/283 (85%)	0.30	7 (2%) 49 53	33, 41, 54, 69	0
1	B	251/283 (88%)	0.01	2 (0%) 83 87	33, 40, 62, 74	0
1	C	254/283 (89%)	0.39	13 (5%) 27 29	29, 39, 57, 73	0
1	D	253/283 (89%)	0.31	4 (1%) 68 72	28, 38, 58, 74	0
All	All	1001/1132 (88%)	0.25	26 (2%) 53 58	28, 40, 59, 74	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	74	VAL	5.2
1	C	263	TRP	5.1
1	B	263	TRP	4.5
1	A	263	TRP	4.4
1	C	227	VAL	3.4
1	A	28	ARG	3.4
1	A	32	LEU	3.3
1	C	78	CYS	3.1
1	D	231	VAL	3.1
1	D	263	TRP	2.9
1	A	27	PHE	2.7
1	C	243	LEU	2.6
1	C	87	HIS	2.5
1	C	219	ASP	2.5
1	A	87	HIS	2.4
1	C	52	TYR	2.3
1	C	88	TYR	2.2
1	D	232	HIS	2.2
1	C	228	SER	2.2
1	A	268	ILE	2.2
1	B	264	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	268	ILE	2.1
1	C	81	LEU	2.1
1	D	233	MET	2.1
1	A	57	MET	2.0
1	C	51	ALA	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	CBO	B	1279	37/41	0.13	0.82	35,42,61,67	0
3	CBO	C	1273	37/41	0.16	0.30	30,38,64,68	0
2	NAP	B	1278	48/48	0.11	-0.19	34,39,45,49	0
2	NAP	D	1278	48/48	0.13	-0.28	26,36,40,42	0
3	CBO	D	1279	37/41	0.12	-0.36	30,36,50,56	0
2	NAP	A	1269	48/48	0.11	-0.38	29,36,45,47	0
3	CBO	A	1270	41/41	0.12	-0.65	34,39,56,67	0
4	CL	C	1274	1/1	0.17	-0.88	55,55,55,55	0
2	NAP	C	1272	48/48	0.09	-1.26	26,36,41,44	0
4	CL	A	1271	1/1	0.08	-1.59	44,44,44,44	0
4	CL	D	1280	1/1	0.07	-2.63	34,34,34,34	0
4	CL	B	1280	1/1	0.05	-6.39	30,30,30,30	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.