



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 04:41 PM GMT

PDB ID : 2HQY  
Title : Crystal Structure of Conserved Protein of Unknown Function from Bacteroides  
thetaiotaomicron VPI-5482  
Authors : Nocek, B.; Borovilos, M.; Abdullah, J.; Joachimiak, A.; Midwest Center for  
Structural Genomics (MCSG)  
Deposited on : 2006-07-19  
Resolution : 1.80 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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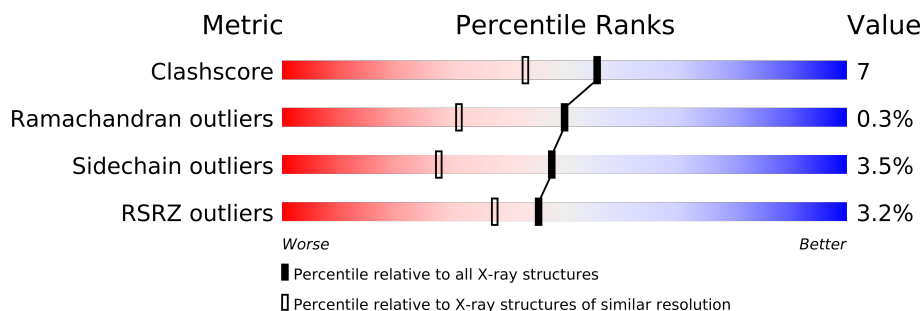
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.80 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))
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (1.80-1.80)

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 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	305	
1	B	305	

## 2 Entry composition i

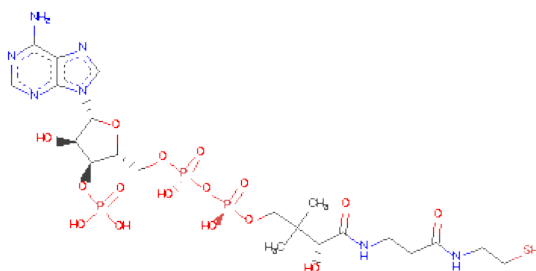
There are 3 unique types of molecules in this entry. The entry contains 5595 atoms, of which 0 are hydrogen and 0 are deuterium.

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 conserved hypothetical protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	28	9	0
			2483	1586	420	459	18			
1	B	296	Total	C	N	O	S	18	9	0
			2451	1571	410	452	18			

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: C<sub>21</sub>H<sub>36</sub>N<sub>7</sub>O<sub>16</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P S	19	0
			48	21	7	16	3 1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	309	Total	O	0	0
			309	309		

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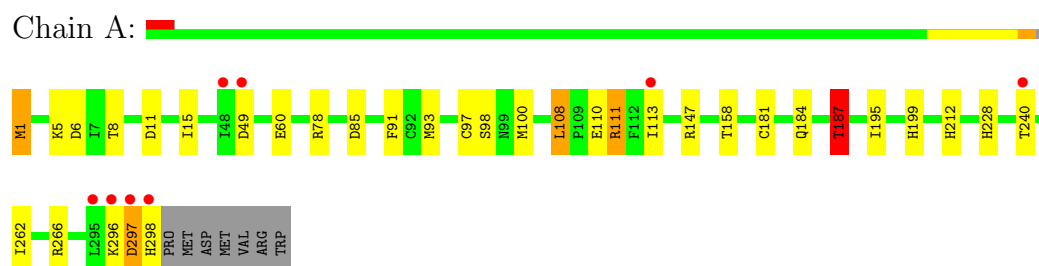
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	304	Total 304	O 304	0	0

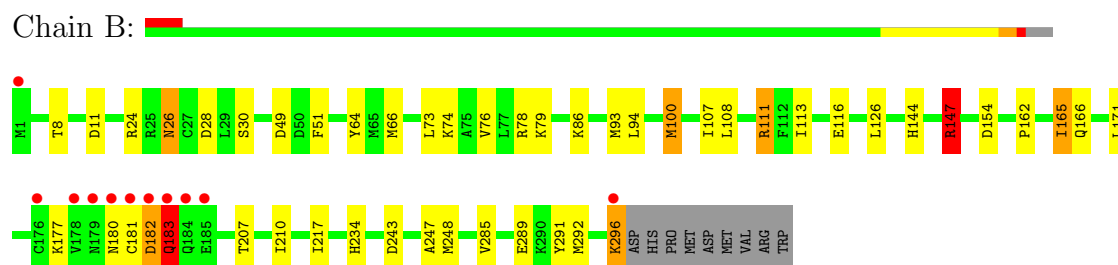
### 3 Residue-property plots

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: conserved hypothetical protein



- Molecule 1: conserved hypothetical protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.47Å 81.88Å 138.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.80 39.26 – 1.79	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.00-1.80) 98.9 (39.26-1.79)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.77 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.176 , 0.211 0.177 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	23.2	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 55.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 71634 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5595	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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: COA

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	1.53	5/2549 (0.2%)	0.91	13/3443 (0.4%)
1	B	0.85	2/2531 (0.1%)	0.83	5/3419 (0.1%)
All	All	1.24	7/5080 (0.1%)	0.87	18/6862 (0.3%)

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	1	1
1	B	0	1
All	All	1	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	111	ARG	CD-NE	44.91	2.22	1.46
1	A	110	GLU	CB-CG	33.36	2.15	1.52
1	A	147	ARG	CD-NE	-27.32	1.00	1.46
1	A	60	GLU	CB-CG	-26.01	1.02	1.52
1	B	147	ARG	NE-CZ	-19.68	1.07	1.33
1	B	177	LYS	CD-CE	12.38	1.82	1.51
1	A	108	LEU	CB-CG	-10.84	1.21	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	GLU	CA-CB-CG	15.89	148.35	113.40
1	B	147	ARG	NE-CZ-NH2	-13.56	113.52	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	ARG	NE-CZ-NH1	13.51	127.06	120.30
1	A	111	ARG	CD-NE-CZ	-12.99	105.42	123.60
1	A	110	GLU	CA-CB-CG	-12.98	84.84	113.40
1	B	147	ARG	CD-NE-CZ	12.22	140.71	123.60
1	A	111	ARG	CG-CD-NE	-9.94	90.92	111.80
1	A	147	ARG	CG-CD-NE	8.92	130.54	111.80
1	A	108	LEU	CA-CB-CG	8.31	134.41	115.30
1	A	248	MET	CG-SD-CE	-7.05	88.91	100.20
1	A	108	LEU	CB-CG-CD2	7.00	122.90	111.00
1	A	147	ARG	CD-NE-CZ	6.65	132.91	123.60
1	B	111	ARG	CD-NE-CZ	6.63	132.88	123.60
1	B	248	MET	CG-SD-CE	-6.32	90.09	100.20
1	A	78	ARG	CG-CD-NE	6.01	124.41	111.80
1	A	187	THR	OG1-CB-CG2	5.87	123.50	110.00
1	A	187	THR	N-CA-CB	-5.45	99.94	110.30
1	A	266	ARG	NE-CZ-NH2	-5.11	117.74	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	297	ASP	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	ARG	Sidechain
1	B	147	ARG	Sidechain

## 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	2483	0	2438	25	0
1	B	2451	0	2427	44	1
2	A	48	0	32	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	309	0	0	5	0
3	B	304	0	0	8	1
All	All	5595	0	4897	64	1

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 7.

All (64) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:66:MET:SD	1:B:93:MET:HE1	1.94	1.08
1:B:171:LEU:CD1	1:B:217:ILE:HG23	1.84	1.07
1:A:187:THR:HG21	3:A:1032:HOH:O	1.75	0.86
1:B:171:LEU:HD13	1:B:217:ILE:HG23	1.59	0.82
1:B:207[B]:THR:HG22	3:B:318:HOH:O	1.81	0.81
1:B:171:LEU:CD1	1:B:217:ILE:CG2	2.60	0.79
1:B:207[B]:THR:HG21	3:B:499:HOH:O	1.85	0.77
1:B:66:MET:HG2	1:B:100:MET:HG3	1.66	0.77
1:B:94:LEU:HD21	1:B:292[B]:MET:HE3	1.65	0.77
1:B:182:ASP:O	1:B:183:GLN:HB2	1.82	0.76
1:A:240:THR:HB	3:A:1236:HOH:O	1.89	0.72
1:A:100:MET:HG3	3:A:1047:HOH:O	1.89	0.72
1:A:85:ASP:OD1	1:A:298:HIS:CE1	2.43	0.71
1:B:66:MET:SD	1:B:93:MET:CE	2.78	0.70
1:A:8:THR:HG21	1:B:8[A]:THR:CG2	2.23	0.69
1:B:94:LEU:CD2	1:B:292[B]:MET:HE3	2.23	0.69
1:A:85:ASP:OD1	1:A:298:HIS:HE1	1.74	0.68
1:B:162:PRO:HA	1:B:165:ILE:HD13	1.76	0.68
1:A:8:THR:CG2	1:B:8[A]:THR:CG2	2.73	0.66
1:B:8[B]:THR:HG22	1:B:11:ASP:OD2	1.94	0.66
1:B:171:LEU:HD11	1:B:217:ILE:HG23	1.74	0.65
1:B:26:ASN:ND2	1:B:28:ASP:H	1.94	0.65
1:B:51:PHE:CD2	1:B:76[B]:VAL:HG21	2.32	0.64
1:A:260:GLN:HE21	1:A:260:GLN:H	1.45	0.62
1:A:97[A]:CYS:SG	1:A:100:MET:HG2	2.40	0.62
1:A:1:MET:O	1:A:1:MET:HG3	2.00	0.61
1:A:8:THR:CG2	1:B:8[A]:THR:HG21	2.32	0.60
1:A:184:GLN:HB2	1:A:187:THR:HG22	1.86	0.57
1:A:181:CYS:SG	1:A:187:THR:HG23	2.46	0.56
1:B:66:MET:HA	1:B:93:MET:HE3	1.86	0.56
1:A:8:THR:HG21	1:B:8[A]:THR:HG21	1.87	0.56
1:B:78:ARG:HD3	3:B:481:HOH:O	2.07	0.55
1:B:49:ASP:OD2	1:B:79:LYS:HE3	2.09	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:49:ASP:HB3	3:A:1115:HOH:O	2.10	0.51
1:A:228:HIS:HD2	3:A:1195:HOH:O	1.94	0.51
1:B:30:SER:OG	1:B:234:HIS:CE1	2.64	0.51
1:B:73:LEU:HD23	1:B:107:ILE:HD11	1.92	0.50
1:B:171:LEU:HD12	1:B:217:ILE:CG2	2.42	0.50
1:A:260:GLN:O	1:A:262:ILE:HD12	2.11	0.50
1:B:182:ASP:O	1:B:183:GLN:CB	2.57	0.50
1:B:182:ASP:HB2	3:B:574:HOH:O	2.12	0.48
1:A:113:ILE:HG13	1:A:296:LYS:HG2	1.94	0.48
1:B:210:ILE:HG23	1:B:217:ILE:HD12	1.96	0.48
1:A:195:ILE:CG2	1:A:199:HIS:CE1	2.97	0.47
1:B:116:GLU:OE2	1:B:289:GLU:OE1	2.33	0.46
1:B:126:LEU:HG	1:B:285:VAL:HG11	1.98	0.46
1:A:158[A]:THR:OG1	1:A:212:HIS:HE1	1.99	0.46
1:B:74:LYS:O	1:B:78:ARG:HG3	2.15	0.46
1:A:6:ASP:HB2	1:B:8[B]:THR:OG1	2.15	0.45
1:B:64:TYR:O	1:B:93:MET:HA	2.17	0.45
1:A:11:ASP:O	1:A:15[A]:ILE:HG12	2.18	0.44
1:B:291:TYR:HD1	3:B:370:HOH:O	2.02	0.43
1:A:113:ILE:HG13	1:A:296:LYS:CG	2.49	0.42
1:B:108:LEU:HB3	1:B:111:ARG:HB2	2.01	0.42
1:B:94:LEU:CD2	1:B:292[B]:MET:CE	2.96	0.42
1:B:180:ASN:O	1:B:182:ASP:O	2.37	0.42
1:B:51:PHE:HD2	1:B:76[B]:VAL:HG21	1.79	0.42
1:B:24:ARG:HD3	3:B:484:HOH:O	2.19	0.41
1:B:144:HIS:HE1	3:B:431:HOH:O	2.03	0.41
1:A:91:PHE:CZ	1:A:93:MET:HG3	2.55	0.41
1:B:247:ALA:HB1	3:B:554:HOH:O	2.20	0.41
1:B:113:ILE:HD11	1:B:296:LYS:HG3	2.03	0.40
1:B:165:ILE:HG22	1:B:166:GLN:N	2.35	0.40
1:A:195:ILE:HG23	1:A:199:HIS:CE1	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:180:ASN:ND2	3:B:370:HOH:O[4_455]	2.19	0.01

## 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	305/305 (100%)	302 (99%)	2 (1%)	1 (0%)	50	31
1	B	303/305 (99%)	299 (99%)	3 (1%)	1 (0%)	50	31
All	All	608/610 (100%)	601 (99%)	5 (1%)	2 (0%)	50	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	183	GLN
1	A	297	ASP

### 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	267/265 (101%)	260 (97%)	7 (3%)	59	41
1	B	265/265 (100%)	253 (96%)	12 (4%)	38	17
All	All	532/530 (100%)	513 (96%)	19 (4%)	48	27

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	LYS
1	A	98	SER
1	A	108	LEU
1	A	187	THR
1	A	260	GLN
1	A	297	ASP

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Mol	Chain	Res	Type
1	B	26	ASN
1	B	86	LYS
1	B	100	MET
1	B	147	ARG
1	B	154	ASP
1	B	165	ILE
1	B	181	CYS
1	B	182	ASP
1	B	183	GLN
1	B	243[A]	ASP
1	B	243[B]	ASP
1	B	296	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	212	HIS
1	A	260	GLN
1	A	298	HIS
1	B	26	ASN
1	B	33	ASN
1	B	144	HIS
1	B	228	HIS
1	B	234	HIS

### 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 ⓘ

1 ligand is 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	COA	A	1001	-	50,50,50	0.80	1 (2%)	75,75,75	2.31	22 (29%)

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	COA	A	1001	-	1/1/11/13	2/48/64/64	0/1/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	COA	O4B-C1B	2.31	1.44	1.41

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	COA	CBP-CAP-C9P	10.09	122.52	112.73
2	A	1001	COA	N3A-C2A-N1A	-8.91	121.26	128.71
2	A	1001	COA	CDP-CBP-CAP	4.12	115.97	108.82
2	A	1001	COA	N3A-C4A-N9A	3.79	132.27	125.43
2	A	1001	COA	C4A-C5A-N7A	-3.65	106.39	109.52
2	A	1001	COA	C6P-C7P-N8P	-3.64	104.03	111.87
2	A	1001	COA	C6P-C5P-N4P	-3.37	110.37	116.50
2	A	1001	COA	O5P-C5P-N4P	3.30	129.47	122.94
2	A	1001	COA	C2B-C1B-N9A	3.26	121.65	113.27
2	A	1001	COA	O3A-P2A-O6A	-2.75	91.09	103.41
2	A	1001	COA	C2P-C3P-N4P	-2.74	106.30	112.45
2	A	1001	COA	CDP-CBP-CCP	-2.74	104.81	108.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	COA	C5A-C4A-N3A	-2.72	119.78	125.70
2	A	1001	COA	C3P-N4P-C5P	2.70	128.33	122.84
2	A	1001	COA	O5A-P2A-O6A	2.57	121.45	108.51
2	A	1001	COA	OAP-CAP-C9P	-2.45	104.45	110.76
2	A	1001	COA	C8A-N9A-C4A	2.28	108.64	106.90
2	A	1001	COA	P2A-O3A-P1A	-2.24	125.12	131.68
2	A	1001	COA	C2A-N3A-C4A	2.20	120.26	114.01
2	A	1001	COA	C7P-N8P-C9P	-2.16	118.14	122.57
2	A	1001	COA	N7A-C8A-N9A	-2.08	108.47	114.36
2	A	1001	COA	CAP-C9P-N8P	2.05	120.97	116.57

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1001	COA	CAP

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	COA	O9P-C9P-CAP-CBP
2	A	1001	COA	CBP-CAP-C9P-N8P

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	298/305 (97%)	-0.02	8 (2%) 52 44	10, 20, 34, 43	8 (2%)
1	B	296/305 (97%)	0.00	11 (3%) 39 32	11, 19, 30, 38	6 (2%)
All	All	594/610 (97%)	-0.01	19 (3%) 45 37	10, 19, 32, 43	14 (2%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	297	ASP	6.6
1	B	183	GLN	5.8
1	A	298	HIS	5.7
1	B	184	GLN	5.5
1	B	178	VAL	4.7
1	B	181	CYS	4.4
1	A	240	THR	4.0
1	B	185	GLU	3.7
1	B	182	ASP	3.4
1	B	180	ASN	3.2
1	B	179	ASN	3.0
1	A	49	ASP	2.9
1	B	296	LYS	2.7
1	B	176	CYS	2.6
1	A	48	ILE	2.5
1	A	295	LEU	2.4
1	B	1	MET	2.4
1	A	296	LYS	2.3
1	A	113	ILE	2.2

### 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
2	COA	A	1001	48/48	0.11	0.30	29,59,84,86	19

### 6.5 Other polymers

There are no such residues in this entry.