



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:44 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 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)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

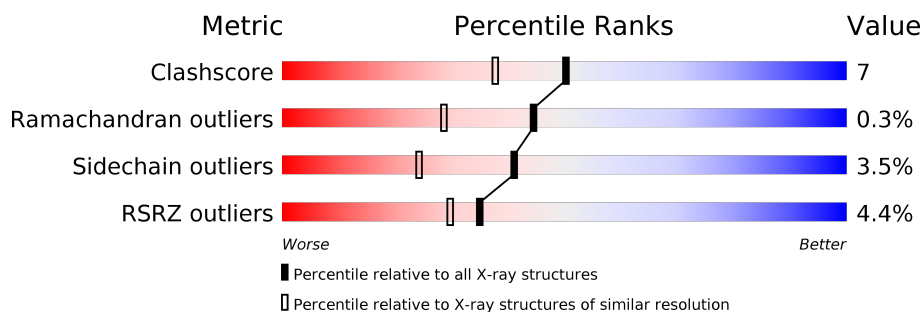
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	112137	5742 (1.80-1.80)
Ramachandran outliers	110173	5676 (1.80-1.80)
Sidechain outliers	110143	5675 (1.80-1.80)
RSRZ outliers	101464	4906 (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 ≥ 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	305	
1	B	305	

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	COA	A	1001	X	-	-	-

2 Entry composition [i](#)

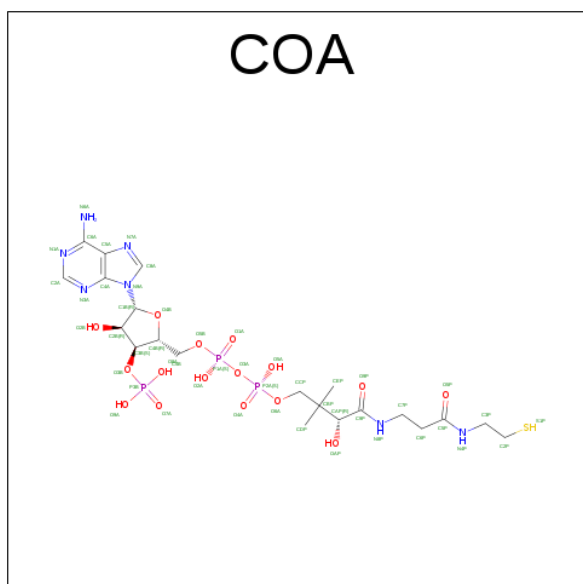
There are 3 unique types of molecules in this entry. The entry contains 5595 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 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: $C_{21}H_{36}N_7O_{16}P_3S$).



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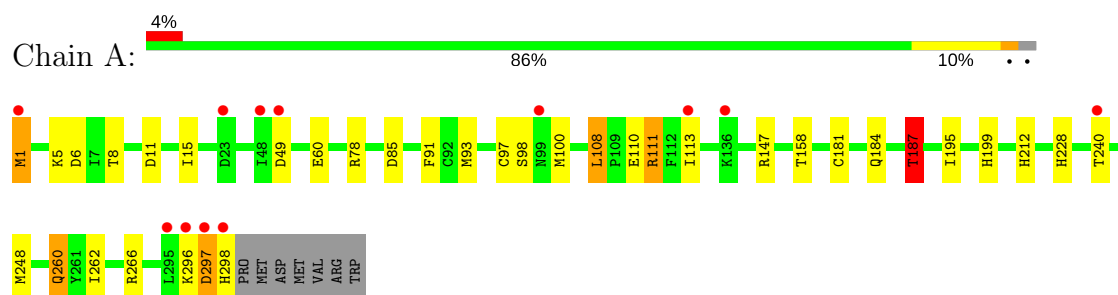
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	304	Total 304	O 304	0	0

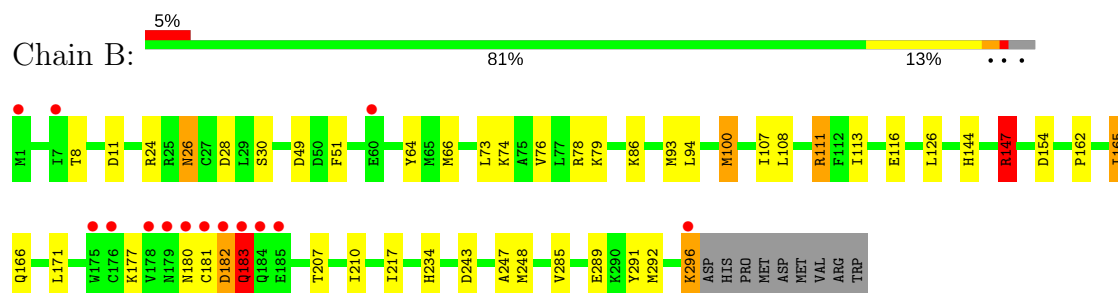
3 Residue-property plots [i](#)

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.

- 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, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	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 ²)	23.2	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5595	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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 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	2483	0	2438	25	0
1	B	2451	0	2427	44	1
2	A	48	0	32	0	0
3	A	309	0	0	5	0
3	B	304	0	0	8	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5595	0	4897	64	1

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 (64) 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: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	Interatomic distance (Å)	Clash overlap (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:A:158[A]:THR:OG1	1:A:212:HIS:HE1	1.99	0.46
1:B:126:LEU:HG	1:B:285:VAL:HG11	1.98	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	Interatomic distance (Å)	Clash overlap (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%)	44	29
1	B	303/305 (99%)	299 (99%)	3 (1%)	1 (0%)	44	29
All	All	608/610 (100%)	601 (99%)	5 (1%)	2 (0%)	44	29

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%)	51	36
1	B	265/265 (100%)	253 (96%)	12 (4%)	32	15
All	All	532/530 (100%)	513 (96%)	19 (4%)	41	23

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

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Mol	Chain	Res	Type
1	A	187	THR
1	A	260	GLN
1	A	297	ASP
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 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

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	-	43,50,50	0.79	1 (2%)	48,75,75	2.12	12 (25%)

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	2/44/64/64	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	COA	O4B-C1B	2.63	1.44	1.41

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	COA	N3A-C2A-N1A	-8.73	121.26	128.86
2	A	1001	COA	C6P-C7P-N8P	-3.79	104.03	111.87
2	A	1001	COA	C6P-C5P-N4P	-3.55	110.37	116.49
2	A	1001	COA	C4A-C5A-N7A	-3.12	106.39	109.41
2	A	1001	COA	C2P-C3P-N4P	-2.85	106.30	112.50
2	A	1001	COA	CDP-CBP-CCP	-2.42	104.81	108.37
2	A	1001	COA	C7P-N8P-C9P	-2.38	118.14	122.59
2	A	1001	COA	CAP-C9P-N8P	2.11	120.97	116.58
2	A	1001	COA	O5A-P2A-O6A	2.82	121.45	108.14
2	A	1001	COA	C3P-N4P-C5P	2.86	128.33	122.84
2	A	1001	COA	O5P-C5P-N4P	3.40	129.47	122.97
2	A	1001	COA	CDP-CBP-CAP	4.12	115.97	108.82

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	CBP-CAP-C9P-N8P
2	A	1001	COA	O9P-C9P-CAP-CBP

There are no ring outliers.

No monomer is involved in short contacts.

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

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	298/305 (97%)	0.23	12 (4%) 39 33	10, 20, 34, 43	8 (2%)
1	B	296/305 (97%)	0.31	14 (4%) 32 27	11, 19, 30, 38	6 (2%)
All	All	594/610 (97%)	0.27	26 (4%) 35 30	10, 19, 32, 43	14 (2%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	297	ASP	7.1
1	B	183	GLN	7.0
1	B	181	CYS	6.4
1	B	178	VAL	6.1
1	A	298	HIS	6.0
1	B	184	GLN	5.9
1	A	240	THR	4.6
1	B	185	GLU	4.3
1	B	182	ASP	4.0
1	B	180	ASN	3.9
1	B	179	ASN	3.8
1	B	176	CYS	3.7
1	A	49	ASP	3.0
1	B	7	ILE	2.9
1	B	175	TRP	2.9
1	A	296	LYS	2.9
1	A	1	MET	2.5
1	B	1	MET	2.4
1	A	48	ILE	2.4
1	A	295	LEU	2.4
1	A	99	ASN	2.3
1	A	136	LYS	2.2
1	B	60	GLU	2.1
1	B	296	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	23	ASP	2.0
1	A	113	ILE	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(Å ²)	Q<0.9
2	COA	A	1001	48/48	0.94	0.15	0.68	29,59,84,86	19

6.5 Other polymers [i](#)

There are no such residues in this entry.