



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:54 PM GMT

PDB ID : 4B4U
Title : Crystal structure of Acinetobacter baumannii N5, N10- methylenetetrahydrofolate dehydrogenase-cyclohydrolase(FolD) complexed with NADP cofactor
Authors : Eadsforth, T.C.; Maluf, F.V.; Hunter, W.N.
Deposited on : 2012-08-01
Resolution : 1.45 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

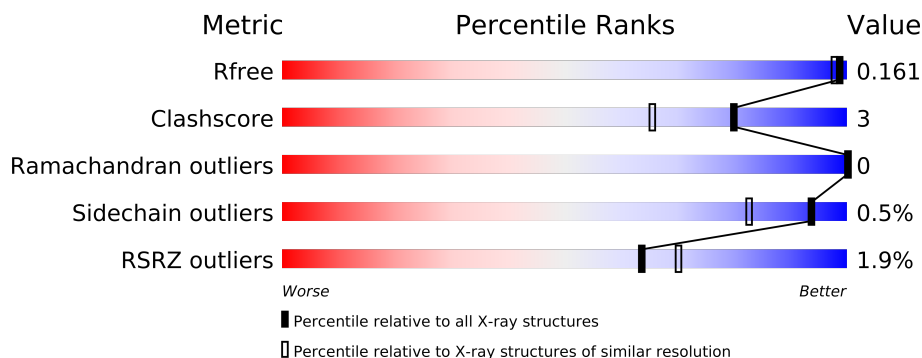
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.45 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	2476 (1.50-1.42)
Clashscore	79885	2834 (1.50-1.42)
Ramachandran outliers	78287	2769 (1.50-1.42)
Sidechain outliers	78261	2767 (1.50-1.42)
RSRZ outliers	66119	2477 (1.50-1.42)

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. 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	303	
1	B	303	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	EDO	A	1286	-	X
4	EDO	A	1287	-	X
4	EDO	B	1286	-	X
4	EDO	B	1287	-	X
5	PEG	A	1288	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5163 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 BIFUNCTIONAL PROTEIN FOLD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	282	Total	C	N	O	S	13	12	0
			2133	1342	369	404	18			
1	B	286	Total	C	N	O	S	11	16	0
			2196	1383	381	415	17			

There are 42 discrepancies between the modelled and reference sequences:

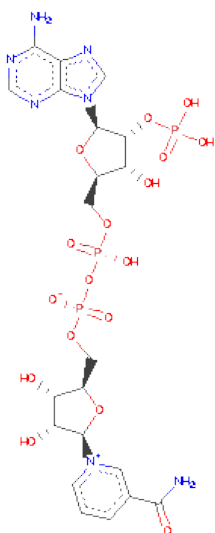
Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP D0CBC8
A	-19	GLY	-	EXPRESSION TAG	UNP D0CBC8
A	-18	SER	-	EXPRESSION TAG	UNP D0CBC8
A	-17	SER	-	EXPRESSION TAG	UNP D0CBC8
A	-16	HIS	-	EXPRESSION TAG	UNP D0CBC8
A	-15	HIS	-	EXPRESSION TAG	UNP D0CBC8
A	-14	HIS	-	EXPRESSION TAG	UNP D0CBC8
A	-13	HIS	-	EXPRESSION TAG	UNP D0CBC8
A	-12	HIS	-	EXPRESSION TAG	UNP D0CBC8
A	-11	HIS	-	EXPRESSION TAG	UNP D0CBC8
A	-10	SER	-	EXPRESSION TAG	UNP D0CBC8
A	-9	SER	-	EXPRESSION TAG	UNP D0CBC8
A	-8	GLY	-	EXPRESSION TAG	UNP D0CBC8
A	-7	GLU	-	EXPRESSION TAG	UNP D0CBC8
A	-6	ASN	-	EXPRESSION TAG	UNP D0CBC8
A	-5	LEU	-	EXPRESSION TAG	UNP D0CBC8
A	-4	TYR	-	EXPRESSION TAG	UNP D0CBC8
A	-3	PHE	-	EXPRESSION TAG	UNP D0CBC8
A	-2	GLN	-	EXPRESSION TAG	UNP D0CBC8
A	-1	GLY	-	EXPRESSION TAG	UNP D0CBC8
A	0	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-20	MET	-	EXPRESSION TAG	UNP D0CBC8
B	-19	GLY	-	EXPRESSION TAG	UNP D0CBC8
B	-18	SER	-	EXPRESSION TAG	UNP D0CBC8
B	-17	SER	-	EXPRESSION TAG	UNP D0CBC8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-15	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-14	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-13	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-12	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-11	HIS	-	EXPRESSION TAG	UNP D0CBC8
B	-10	SER	-	EXPRESSION TAG	UNP D0CBC8
B	-9	SER	-	EXPRESSION TAG	UNP D0CBC8
B	-8	GLY	-	EXPRESSION TAG	UNP D0CBC8
B	-7	GLU	-	EXPRESSION TAG	UNP D0CBC8
B	-6	ASN	-	EXPRESSION TAG	UNP D0CBC8
B	-5	LEU	-	EXPRESSION TAG	UNP D0CBC8
B	-4	TYR	-	EXPRESSION TAG	UNP D0CBC8
B	-3	PHE	-	EXPRESSION TAG	UNP D0CBC8
B	-2	GLN	-	EXPRESSION TAG	UNP D0CBC8
B	-1	GLY	-	EXPRESSION TAG	UNP D0CBC8
B	0	HIS	-	EXPRESSION TAG	UNP D0CBC8

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

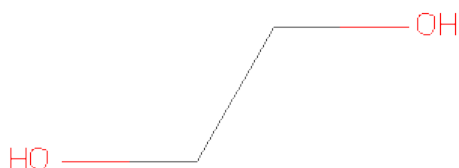


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

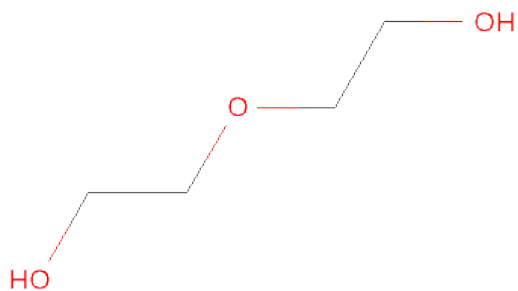
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		
3	A	3	Total	Cl	0	0
			3	3		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	388	Total	O	0	0
			388	388		
6	B	352	Total	O	0	0
			352	352		

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- Molecule 1: BIFUNCTIONAL PROTEIN FOLD

Amino Acid	Percentage (%)
MET	~1
GLY	~2
SER	~3
HIS	~4
HIS	~5
HIS	~6
HIS	~7
HIS	~8
SER	~9
SER	~10
GLY	~11
GLU	~12
ASN	~13
LEU	~14
TYR	~15
PHE	~16
GLN	~17
GLY	~18
HIS	~19
MI	~20
Q13	~1
K66	~2
R107	~3
C124	~4
Q180	~5
E197	~6
D285	~7
G236	~8
G237	~9
R271	~10
G242	~11

- Molecule 1: BIFUNCTIONAL PROTEIN FOLD

MET	GLY	SER	SER	HIS	HIS	HIS	HIS	HIS	HIS	SER	GLY	GLU	ASN	LEU	TYR	F-3	Q-2	H0	M1	I34	R51	R58	D63	K66	Q71	G128	R129	M132	G133	E134	R234	D235	G236	G237	R271	G284
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.04Å 80.13Å 68.62Å 90.00° 107.11° 90.00°	Depositor
Resolution (Å)	52.60 – 1.45 43.97 – 1.45	Depositor EDS
% Data completeness (in resolution range)	99.7 (52.60-1.45) 99.7 (43.97-1.45)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 1.45Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.122 , 0.152 0.134 , 0.161	Depositor DCC
R_{free} test set	5019 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	13.8	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 38.2	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	0 of 100364 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	5163	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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, PEG, EDO, 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	0.50	1/2193 (0.0%)	0.69	2/2964 (0.1%)
1	B	0.55	2/2270 (0.1%)	0.71	4/3067 (0.1%)
All	All	0.53	3/4463 (0.1%)	0.70	6/6031 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	66	LYS	CB-CG	-9.55	1.26	1.52
1	A	66	LYS	CD-CE	-7.26	1.33	1.51
1	B	235	ASP	CA-CB	6.12	1.67	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	LYS	CG-CD-CE	13.20	151.50	111.90
1	B	235	ASP	CB-CA-C	-7.46	95.49	110.40
1	B	63[A]	ASP	CB-CG-OD1	5.89	123.60	118.30
1	B	63[B]	ASP	CB-CG-OD1	5.89	123.60	118.30
1	B	271	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	A	271	ARG	NE-CZ-NH1	5.13	122.87	120.30

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	2133	0	7	9	0
1	B	2196	0	12	11	0
2	A	31	0	0	0	0
2	B	31	0	0	0	0
3	A	3	0	0	0	0
3	B	2	0	0	1	0
4	A	8	0	0	0	0
4	B	12	0	0	0	0
5	A	7	0	0	0	0
6	A	388	0	0	4	0
6	B	352	0	0	2	0
All	All	5163	0	19	15	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 (15) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:234[B]:ARG:CG	1:B:234[B]:ARG:NH1	2.36	0.87
1:B:71[B]:GLN:OE1	3:B:1284:CL:CL	2.42	0.74
1:B:34:ILE:CD1	1:B:63[B]:ASP:CB	2.72	0.67
1:B:51:ARG:NH1	6:B:2110:HOH:O	2.30	0.63
1:B:58:ARG:NH2	6:B:2107:HOH:O	2.37	0.57
1:A:13:GLN:NE2	6:A:2037:HOH:O	2.42	0.53
1:A:124[A]:CYS:SG	1:B:134[A]:GLU:OE2	2.69	0.51
1:A:124[B]:CYS:SG	1:B:129:ARG:N	2.84	0.50
1:A:124[B]:CYS:SG	1:B:128:GLY:C	2.93	0.47
1:A:107:ARG:CZ	1:B:132[B]:MET:SD	3.03	0.47
1:A:180[A]:GLN:NE2	6:A:2249:HOH:O	2.49	0.46
1:B:129:ARG:NH1	1:B:134[A]:GLU:OE1	2.50	0.45
1:A:124[B]:CYS:SG	1:B:129:ARG:CA	3.05	0.44
1:A:180[A]:GLN:NE2	6:A:2300:HOH:O	2.52	0.41
1:A:197[A]:GLU:CG	6:A:2316:HOH:O	2.69	0.41

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	292/303 (96%)	289 (99%)	3 (1%)	0	100	100
1	B	300/303 (99%)	297 (99%)	3 (1%)	0	100	100
All	All	592/606 (98%)	586 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	229/235 (97%)	228 (100%)	1 (0%)	95	85
1	B	236/235 (100%)	235 (100%)	1 (0%)	95	85
All	All	465/470 (99%)	463 (100%)	2 (0%)	94	85

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

Mol	Chain	Res	Type
1	A	235	ASP
1	B	-2	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 13 ligands modelled in this entry, 5 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	1001	-	33,33,52	1.63	3 (9%)	52,52,80	2.56	9 (17%)
4	EDO	A	1286	-	3,3,3	0.19	0	2,2,2	0.73	0
4	EDO	A	1287	-	3,3,3	0.67	0	2,2,2	0.41	0
5	PEG	A	1288	-	6,6,6	0.54	0	5,5,5	0.25	0
2	NAP	B	1001	-	33,33,52	1.67	3 (9%)	52,52,80	2.50	9 (17%)
4	EDO	B	1285	-	3,3,3	0.57	0	2,2,2	0.21	0
4	EDO	B	1286	-	3,3,3	0.34	0	2,2,2	0.33	0
4	EDO	B	1287	-	3,3,3	0.50	0	2,2,2	0.15	0

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	1001	-	-	0/21/37/67	0/1/3/5
4	EDO	A	1286	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1287	-	-	0/1/1/1	0/0/0/0
5	PEG	A	1288	-	-	0/4/4/4	0/0/0/0
2	NAP	B	1001	-	-	0/21/37/67	0/1/3/5
4	EDO	B	1285	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	1286	-	-	0/1/1/1	0/0/0/0
4	EDO	B	1287	-	-	0/1/1/1	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	NAP	C8A-N9A	5.82	1.45	1.36
2	A	1001	NAP	C8A-N7A	5.55	1.45	1.34
2	B	1001	NAP	C8A-N7A	5.39	1.45	1.34
2	A	1001	NAP	C8A-N9A	5.36	1.44	1.36
2	B	1001	NAP	C4A-N9A	-4.00	1.31	1.37
2	A	1001	NAP	C4A-N9A	-3.68	1.32	1.37

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	NAP	C8A-N9A-C4A	11.02	115.31	106.90
2	A	1001	NAP	C8A-N9A-C4A	11.01	115.30	106.90
2	B	1001	NAP	N3A-C2A-N1A	-8.56	121.55	128.71
2	A	1001	NAP	N3A-C2A-N1A	-7.97	122.04	128.71
2	A	1001	NAP	O4B-C1B-N9A	6.27	114.27	108.44
2	B	1001	NAP	N7A-C8A-N9A	-5.78	98.01	114.36
2	A	1001	NAP	N7A-C8A-N9A	-5.63	98.43	114.36
2	B	1001	NAP	O4B-C1B-N9A	4.64	112.75	108.44
2	A	1001	NAP	C8A-N9A-C1B	-4.30	117.90	126.38
2	B	1001	NAP	C8A-N9A-C1B	-3.92	118.66	126.38
2	B	1001	NAP	N3A-C4A-N9A	3.82	132.33	125.43
2	A	1001	NAP	N3A-C4A-N9A	3.50	131.74	125.43
2	A	1001	NAP	PN-O3-PA	-3.13	122.03	133.31
2	B	1001	NAP	PN-O3-PA	-2.91	122.81	133.31
2	B	1001	NAP	C8A-N7A-C5A	2.73	112.03	103.58
2	A	1001	NAP	C8A-N7A-C5A	2.67	111.86	103.58
2	A	1001	NAP	O3X-P2B-O2B	-2.44	100.06	107.09
2	B	1001	NAP	O3X-P2B-O2B	-2.02	101.29	107.09

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, 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	282/303 (93%)	-0.45	4 (1%) 72 78	9, 14, 28, 78	4 (1%)
1	B	286/303 (94%)	-0.36	7 (2%) 56 62	9, 14, 27, 75	3 (1%)
All	All	568/606 (93%)	-0.41	11 (1%) 64 70	9, 14, 28, 78	7 (1%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	GLY	6.2
1	B	0	HIS	4.5
1	A	237	GLY	3.9
1	B	235	ASP	3.5
1	B	237	GLY	3.0
1	B	-2	GLN	3.0
1	B	-3	PHE	2.5
1	B	282	GLY	2.4
1	A	235	ASP	2.4
1	A	1	MET	2.3
1	B	1	MET	2.1

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	EDO	B	1287	4/4	0.24	15.45	28,29,31,33	0
4	EDO	A	1287	4/4	0.18	10.15	28,31,32,36	0
4	EDO	B	1286	4/4	0.19	4.80	19,22,24,32	0
4	EDO	A	1286	4/4	0.18	4.25	19,24,25,31	0
5	PEG	A	1288	7/7	0.16	4.13	34,37,41,46	0
4	EDO	B	1285	4/4	0.31	1.97	40,46,51,51	0
2	NAP	A	1001	31/48	0.08	1.36	12,17,68,84	0
2	NAP	B	1001	31/48	0.07	0.32	12,17,68,81	0
3	CL	A	1284	1/1	0.06	0.25	15,15,15,15	0
3	CL	A	1285	1/1	0.06	-1.36	23,23,23,23	0
3	CL	A	1283	1/1	0.03	-2.15	23,23,23,23	0
3	CL	B	1284	1/1	0.03	-2.37	17,17,17,17	0
3	CL	B	1283	1/1	0.04	-4.53	24,24,24,24	0

6.5 Other polymers ⓘ

There are no such residues in this entry.