



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:59 AM GMT

PDB ID : 2Z2M
Title : Cefditoren-Acylated Penicillin-Binding Protein 2X (PBP2X) from *Streptococcus pneumoniae*
Authors : Yamada, M.; Watanabe, T.; Takeuchi, Y.
Deposited on : 2007-05-23
Resolution : 2.60 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

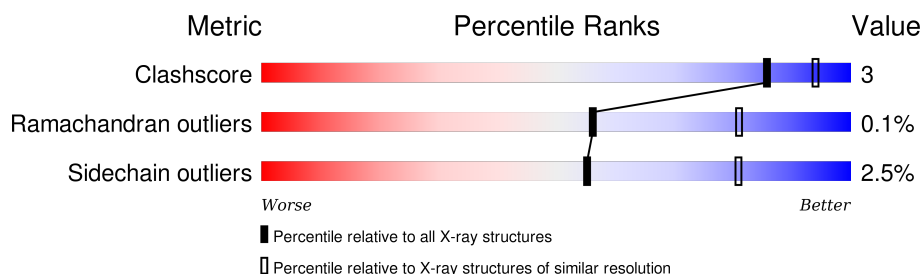
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	168	
1	D	168	
2	B	385	
2	E	385	
3	C	125	
3	F	125	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9617 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 Penicillin-binding protein 2X.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	159	Total	C	N	O	S	0	0	0
			1224	766	204	250	4			
1	D	100	Total	C	N	O	S	0	0	0
			743	459	126	155	3			

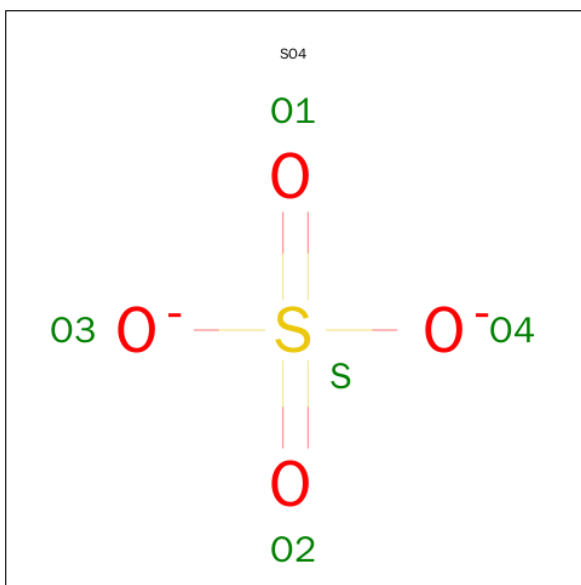
- Molecule 2 is a protein called Penicillin-binding protein 2X.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	363	Total	C	N	O	S	0	0	0
			2784	1752	459	556	17			
2	E	362	Total	C	N	O	S	0	0	0
			2778	1748	459	554	17			

- Molecule 3 is a protein called Penicillin-binding protein 2X.

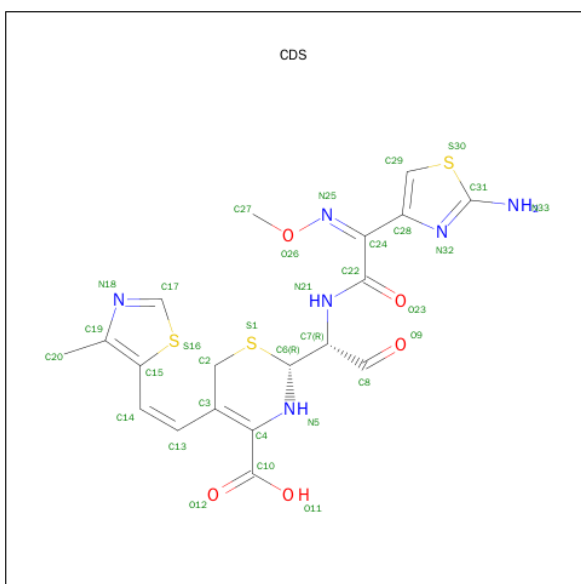
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	125	Total	C	N	O	S	0	0	0
			962	601	162	197	2			
3	F	125	Total	C	N	O	S	0	0	0
			962	601	162	197	2			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is (2R)-2-[(1R)-1-{|(2Z)-2-(2-AMINO-1,3-THIAZOL-4-YL)-2-(METHOXYIMINO)ACETYL}AMINO}-2-OXOETHYL]-5-[(Z)-2-(4-METHYL-1,3-THIAZOL-5-YL)VINYLYL]-3,6-DIHYDRO-2H-1,3-THIAZINE-4-CARBOXYLIC ACID (three-letter code: CDS) (formula: C₁₉H₂₀N₆O₅S₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			33	19	6	5	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	E	1	Total	C	N	O	S	0	0
			33	19	6	5	3		

- Molecule 6 is water.


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total	O	0	0
			3	3		
6	B	30	Total	O	0	0
			30	30		
6	C	8	Total	O	0	0
			8	8		
6	D	13	Total	O	0	0
			13	13		
6	E	37	Total	O	0	0
			37	37		
6	F	2	Total	O	0	0
			2	2		

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.

Note EDS was not executed.

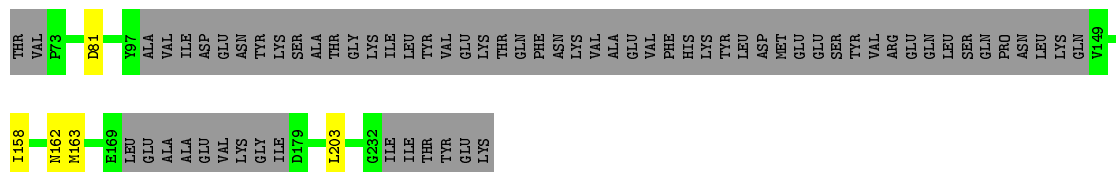
- Molecule 1: Penicillin-binding protein 2X

Chain A: 




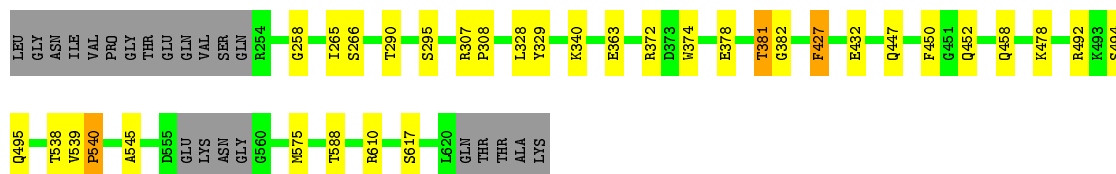
- Molecule 1: Penicillin-binding protein 2X

Chain D: 




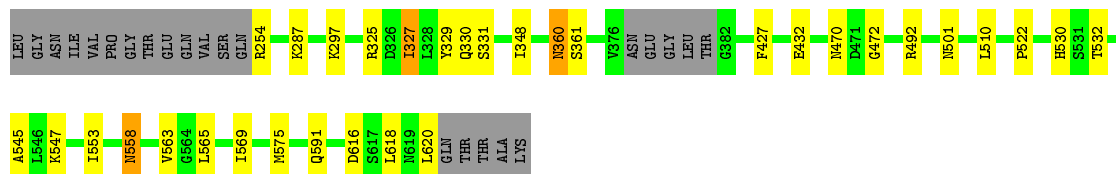
- Molecule 2: Penicillin-binding protein 2X

Chain B: 



- Molecule 2: Penicillin-binding protein 2X

Chain E: 



- Molecule 3: Penicillin-binding protein 2X

Chain C:

86%

13%

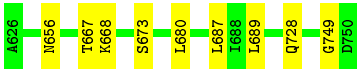


● Molecule 3: Penicillin-binding protein 2X

Chain F:

93%

7%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.73Å 171.21Å 89.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.42 – 2.60	Depositor
% Data completeness (in resolution range)	89.8 (29.42-2.60)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.221 , 0.267	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9617	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

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: SO4, CDS

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	0.35	0/1244	0.49	0/1676
1	D	0.36	0/754	0.51	0/1013
2	B	0.38	0/2837	0.51	0/3848
2	E	0.38	0/2831	0.52	0/3838
3	C	0.33	0/975	0.47	0/1320
3	F	0.34	0/975	0.47	0/1320
All	All	0.37	0/9616	0.50	0/13015

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	1224	0	1191	7	0
1	D	743	0	708	1	0
2	B	2784	0	2717	18	0
2	E	2778	0	2712	17	0
3	C	962	0	975	11	0
3	F	962	0	975	5	0
4	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	33	0	18	0	0
5	E	33	0	18	0	0
6	A	3	0	0	0	0
6	B	30	0	0	0	0
6	C	8	0	0	0	0
6	D	13	0	0	0	0
6	E	37	0	0	0	0
6	F	2	0	0	0	0
All	All	9617	0	9314	52	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 3.

All (52) 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)
2:E:327:ILE:HG22	2:E:331:SER:HB2	1.59	0.84
2:E:492:ARG:NH2	3:F:656:ASN:O	2.11	0.83
2:B:492:ARG:NH1	3:C:680:LEU:O	2.28	0.66
3:F:728:GLN:HG3	3:F:749:GLY:HA3	1.77	0.66
2:B:492:ARG:NH2	3:C:656:ASN:O	2.28	0.65
3:F:673:SER:HB3	3:F:687:LEU:HB2	1.78	0.64
2:E:558:ASN:H	2:E:558:ASN:HD22	1.46	0.64
3:C:728:GLN:HG3	3:C:749:GLY:HA3	1.82	0.62
2:E:297:LYS:HA	2:E:470:ASN:HD21	1.64	0.61
2:E:492:ARG:NH1	3:F:680:LEU:O	2.34	0.61
3:C:671:ASN:HD22	3:C:672:SER:H	1.48	0.60
2:B:329:TYR:HB2	2:B:432:GLU:HG3	1.83	0.60
2:B:290:THR:HG1	2:B:588:THR:HG1	1.51	0.59
2:E:297:LYS:HA	2:E:470:ASN:ND2	2.20	0.55
1:A:135:SER:O	1:A:139:GLU:HG2	2.06	0.55
2:E:360:ASN:HD22	2:E:361:SER:N	2.05	0.55
3:F:667:THR:HG23	3:F:668:LYS:HG3	1.89	0.54
1:A:206:ASN:HD21	1:A:210:SER:HB3	1.72	0.53
2:B:328:LEU:O	2:B:458:GLN:HB2	2.09	0.52
2:E:329:TYR:HB2	2:E:432:GLU:HG3	1.92	0.52
1:A:227:LEU:O	2:B:258:GLY:HA3	2.10	0.52
2:E:553:ILE:CD1	2:E:569:ILE:HG13	2.40	0.51
3:C:658:VAL:O	3:C:660:PRO:HD3	2.11	0.50
2:E:348:ILE:HD13	2:E:510:LEU:HB3	1.92	0.50
2:B:478:LYS:HE3	2:B:495:GLN:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:522:PRO:O	2:E:530:HIS:HE1	1.96	0.49
3:C:661:ILE:HD12	3:C:706:THR:HG21	1.94	0.49
2:E:472:GLY:O	2:E:501:ASN:ND2	2.45	0.48
2:B:494:SER:HB3	3:C:657:LEU:HD22	1.95	0.48
2:E:618:LEU:O	2:E:620:LEU:HG	2.14	0.47
1:D:158:ILE:HG23	1:D:162:ASN:HB2	1.94	0.47
1:A:81:ASP:HB2	2:B:265:ILE:O	2.15	0.47
3:C:729:LYS:HB2	3:C:747:THR:HB	1.97	0.46
2:B:427:PHE:H	2:B:427:PHE:HD1	1.63	0.46
2:B:374:TRP:CE2	2:B:378:GLU:HG3	2.51	0.46
3:C:667:THR:HG23	3:C:668:LYS:HG3	1.97	0.46
3:C:671:ASN:HD22	3:C:672:SER:N	2.14	0.45
2:E:287:LYS:HD3	2:E:591:GLN:OE1	2.17	0.44
2:B:447:GLN:HB3	2:B:452:GLN:HB3	1.97	0.44
2:B:363:GLU:HG2	2:B:372:ARG:HG2	1.99	0.44
2:B:539:VAL:HA	2:B:540:PRO:HD2	1.80	0.43
2:B:545:ALA:HB3	2:B:575:MET:HB2	1.99	0.43
2:E:325:ARG:HD3	2:E:330:GLN:O	2.19	0.43
2:E:327:ILE:HG22	2:E:331:SER:CB	2.42	0.43
2:B:381:THR:HG22	2:B:382:GLY:H	1.81	0.43
3:C:696:VAL:HG22	3:C:738:ILE:HD12	2.01	0.43
2:B:340:LYS:HE2	2:B:450:PHE:HB2	2.01	0.42
1:A:94:TYR:CD2	1:A:182:THR:HG22	2.56	0.41
2:E:545:ALA:HB3	2:E:575:MET:HB2	2.02	0.41
1:A:206:ASN:ND2	1:A:210:SER:HB3	2.35	0.41
1:A:79:ILE:HB	1:A:88:ALA:HB3	2.03	0.41
2:B:307:ARG:HA	2:B:308:PRO:C	2.41	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	157/168 (94%)	151 (96%)	6 (4%)	0	100	100
1	D	94/168 (56%)	90 (96%)	4 (4%)	0	100	100
2	B	359/385 (93%)	347 (97%)	11 (3%)	1 (0%)	46	72
2	E	358/385 (93%)	344 (96%)	14 (4%)	0	100	100
3	C	123/125 (98%)	122 (99%)	1 (1%)	0	100	100
3	F	123/125 (98%)	118 (96%)	5 (4%)	0	100	100
All	All	1214/1356 (90%)	1172 (96%)	41 (3%)	1 (0%)	56	81

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	540	PRO

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	132/141 (94%)	130 (98%)	2 (2%)	72	90
1	D	80/141 (57%)	77 (96%)	3 (4%)	40	68
2	B	304/322 (94%)	297 (98%)	7 (2%)	58	83
2	E	303/322 (94%)	293 (97%)	10 (3%)	45	73
3	C	108/108 (100%)	105 (97%)	3 (3%)	51	78
3	F	108/108 (100%)	107 (99%)	1 (1%)	84	95
All	All	1035/1142 (91%)	1009 (98%)	26 (2%)	55	81

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	168	LYS
1	A	204	HIS
2	B	266	SER
2	B	295	SER

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Mol	Chain	Res	Type
2	B	381	THR
2	B	427	PHE
2	B	538	THR
2	B	610	ARG
2	B	617	SER
3	C	671	ASN
3	C	738	ILE
3	C	750	ASP
1	D	81	ASP
1	D	163	MET
1	D	203	LEU
2	E	254	ARG
2	E	327	ILE
2	E	360	ASN
2	E	427	PHE
2	E	532	THR
2	E	547	LYS
2	E	558	ASN
2	E	563	VAL
2	E	565	LEU
2	E	616	ASP
3	F	689	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	ASN
1	A	206	ASN
2	B	458	GLN
2	B	487	ASN
2	B	514	ASN
3	C	671	ASN
1	D	162	ASN
2	E	360	ASN
2	E	458	GLN
2	E	470	ASN
2	E	530	HIS
2	E	558	ASN
3	F	715	ASN
3	F	721	GLN

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 ⓘ

3 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$
4	SO4	B	801	-	4,4,4	0.14	0	6,6,6	0.15	0
5	CDS	B	901	2	22,35,35	1.50	2 (9%)	14,48,48	2.45	5 (35%)
5	CDS	E	901	2	22,35,35	1.36	2 (9%)	14,48,48	2.50	4 (28%)

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
4	SO4	B	801	-	-	0/0/0/0	0/0/0/0
5	CDS	B	901	2	-	1/13/43/43	0/2/3/3
5	CDS	E	901	2	-	1/13/43/43	0/2/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	901	CDS	O26-N25	2.04	1.44	1.40
5	E	901	CDS	C29-S30	2.06	1.73	1.70
5	E	901	CDS	C10-C4	4.49	1.59	1.52
5	B	901	CDS	C10-C4	5.29	1.61	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	901	CDS	O23-C22-N21	-3.01	117.60	123.01
5	B	901	CDS	O23-C22-N21	-2.66	118.23	123.01
5	B	901	CDS	C7-N21-C22	2.14	125.53	122.01
5	E	901	CDS	C2-S1-C6	3.21	100.55	94.36
5	B	901	CDS	C2-S1-C6	4.10	102.27	94.36
5	B	901	CDS	C6-C7-N21	4.21	119.40	109.57
5	E	901	CDS	C6-C7-N21	4.60	120.31	109.57
5	B	901	CDS	O23-C22-C24	4.80	125.88	120.53
5	E	901	CDS	O23-C22-C24	5.50	126.66	120.53

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	901	CDS	C14-C13-C3-C2
5	B	901	CDS	C14-C13-C3-C2

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.