



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:22 PM UTC

PDB ID : 2DBT / pdb_00002dbt
Title : Crystal structure of chitinase C from Streptomyces griseus HUT6037
Authors : Kezuka, Y.; Watanabe, T.; Nonaka, T.
Deposited on : 2005-12-16
Resolution : 3.14 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

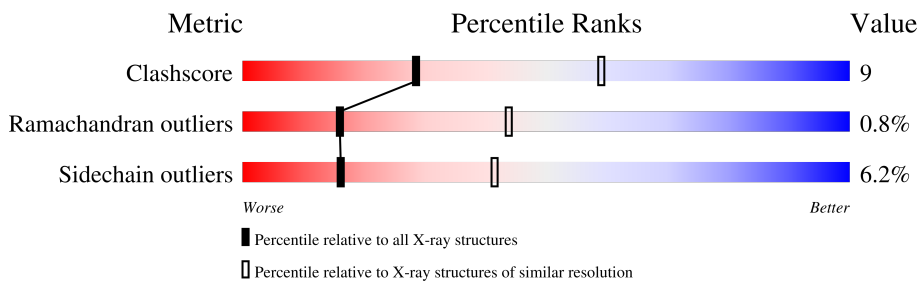
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	265	
1	B	265	
1	C	265	

2 Entry composition [i](#)

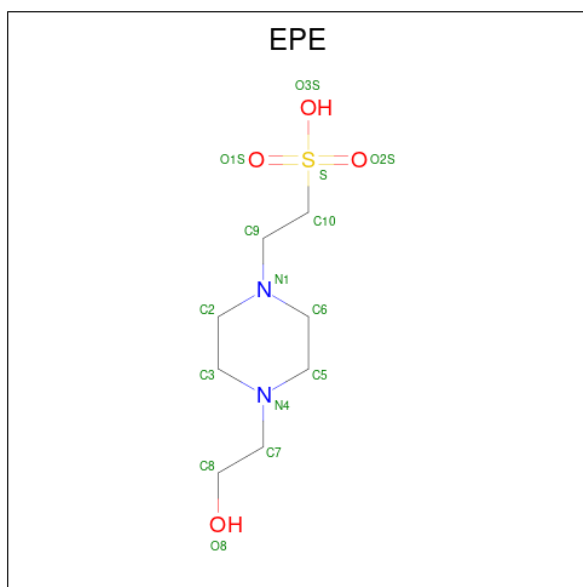
There are 3 unique types of molecules in this entry. The entry contains 4794 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 chitinase C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	0	0
			1576	999	269	302	6			
1	B	205	Total	C	N	O	S	0	0	0
			1576	999	269	302	6			
1	C	205	Total	C	N	O	S	0	0	0
			1576	999	269	302	6			

- Molecule 2 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
2	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total 8	O 8	0	0
3	B	6	Total 6	O 6	0	0
3	C	7	Total 7	O 7	0	0

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	153.12Å 153.12Å 90.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	58.32 – 3.14	Depositor
% Data completeness (in resolution range)	100.0 (58.32-3.14)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.174 , 0.225	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4794	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EPE

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	0.70	0/1620	1.00	3/2207 (0.1%)
1	B	0.74	0/1620	1.02	5/2207 (0.2%)
1	C	0.63	0/1620	0.94	2/2207 (0.1%)
All	All	0.69	0/4860	0.99	10/6621 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	VAL	N-CA-C	6.22	116.38	110.53
1	A	185	GLY	CA-C-N	6.08	125.97	119.28
1	A	185	GLY	C-N-CA	6.08	125.97	119.28
1	B	185	GLY	CA-C-N	5.98	125.66	119.56
1	B	185	GLY	C-N-CA	5.98	125.66	119.56
1	C	163	PRO	CA-C-N	-5.82	114.50	123.17
1	C	163	PRO	C-N-CA	-5.82	114.50	123.17
1	B	121	TYR	CA-C-N	5.21	125.51	119.47
1	B	121	TYR	C-N-CA	5.21	125.51	119.47
1	B	215	GLN	N-CA-C	5.11	117.63	111.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1576	0	1479	24	0
1	B	1576	0	1479	34	0
1	C	1576	0	1479	27	0
2	A	15	0	17	0	0
2	B	15	0	17	2	0
2	C	15	0	17	0	0
3	A	8	0	0	0	0
3	B	6	0	0	0	0
3	C	7	0	0	0	0
All	All	4794	0	4488	81	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 9.

All (81) 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:232:ASN:HB3	1:B:237:MET:O	1.72	0.90
1:A:164:HIS:ND1	1:C:164:HIS:HB3	1.95	0.81
1:A:126:LYS:HA	1:A:133:LYS:HD3	1.63	0.79
1:B:102:PHE:O	1:B:105:ARG:HG2	1.92	0.69
1:A:99:ASN:OD1	1:A:105:ARG:NH1	2.26	0.68
1:B:232:ASN:HB3	1:B:237:MET:C	2.19	0.67
1:B:248:GLY:HA3	1:B:291:ASN:HB2	1.78	0.65
1:A:272:SER:O	1:A:276:LYS:HD2	1.97	0.64
1:B:99:ASN:OD1	1:B:105:ARG:NH1	2.31	0.64
1:B:211:TYR:HB3	1:B:215:GLN:HE22	1.65	0.61
1:A:253:ILE:HG12	1:A:273:ARG:HD2	1.83	0.61
1:A:164:HIS:HA	1:C:164:HIS:HB3	1.84	0.60
1:A:190:SER:O	1:A:191:TRP:HB2	2.02	0.60
1:A:164:HIS:HA	1:C:164:HIS:CB	2.33	0.58
1:A:102:PHE:O	1:A:105:ARG:HG2	2.02	0.58
1:C:184:ARG:HB2	1:C:213:VAL:HG12	1.86	0.56
1:B:258:GLY:HA2	1:B:262:CYS:SG	2.47	0.54
1:C:257:ASN:HD22	1:C:261:GLU:HG3	1.73	0.53
1:B:190:SER:O	1:B:191:TRP:HB2	2.08	0.53
1:C:91:PHE:CE2	1:C:93:VAL:HB	2.44	0.53
1:A:241:ASN:O	1:A:242:ALA:C	2.53	0.52
1:B:232:ASN:CB	1:B:237:MET:O	2.54	0.51
1:C:165:TYR:CD2	1:C:182:TYR:HA	2.46	0.50
1:C:161:ASN:C	1:C:163:PRO:HD2	2.35	0.50
1:B:143:ASN:O	1:B:147:GLU:HG2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ASN:HD21	1:B:212:LEU:HG	1.77	0.50
1:C:204:ILE:HD13	1:C:219:VAL:HG22	1.93	0.50
1:B:114:LEU:HD13	1:B:221:TRP:CE3	2.48	0.49
1:B:146:HIS:CE1	1:B:269:GLN:HG3	2.46	0.49
1:A:264:GLY:O	1:A:265:GLY:C	2.55	0.49
1:B:134:LYS:HD2	1:B:284:LEU:HB3	1.94	0.48
1:B:109:TYR:CG	1:B:151:LEU:HD22	2.48	0.48
1:B:261:GLU:CD	1:B:269:GLN:HG2	2.38	0.48
1:C:288:THR:HB	1:C:292:LEU:HD21	1.95	0.48
1:A:136:GLU:OE2	1:A:240:HIS:ND1	2.44	0.48
1:A:164:HIS:CG	1:C:164:HIS:HB3	2.48	0.47
1:C:281:THR:HG22	1:C:286:THR:HG23	1.96	0.47
1:B:181:TYR:CE1	1:B:210:PRO:HG3	2.49	0.47
1:A:266:ASN:OD1	1:A:269:GLN:HG3	2.15	0.47
1:A:164:HIS:CE1	1:C:163:PRO:HG2	2.50	0.47
1:C:257:ASN:ND2	1:C:261:GLU:HG3	2.29	0.46
1:B:257:ASN:O	1:B:259:ALA:N	2.49	0.46
1:B:196:LYS:HB2	1:B:207:LEU:HD13	1.96	0.46
2:B:295:EPE:H101	2:B:295:EPE:H22	1.64	0.46
1:C:184:ARG:HB2	1:C:213:VAL:CG1	2.46	0.46
1:B:205:ASN:ND2	1:B:208:ALA:HB2	2.30	0.46
1:A:227:TYR:HA	1:A:231:GLN:HG2	1.97	0.46
1:C:206:LEU:HD22	1:C:213:VAL:HG23	1.97	0.45
1:C:280:PHE:CD1	1:C:280:PHE:N	2.83	0.45
1:B:91:PHE:CE2	1:B:93:VAL:HB	2.52	0.45
1:B:115:THR:HA	1:B:118:LEU:HD12	1.99	0.45
1:B:211:TYR:HB3	1:B:215:GLN:NE2	2.32	0.44
1:B:266:ASN:ND2	1:B:269:GLN:HB2	2.31	0.44
1:A:131:GLU:HG2	1:A:286:THR:OG1	2.18	0.43
1:C:174:CYS:HB3	1:C:181:TYR:CE2	2.52	0.43
1:B:109:TYR:CD2	1:B:151:LEU:HD22	2.54	0.43
1:A:143:ASN:OD1	1:A:273:ARG:HD3	2.19	0.43
1:B:198:ALA:HB2	1:B:226:TRP:CE3	2.54	0.42
1:A:155:LYS:H	1:A:155:LYS:HG2	1.59	0.42
1:A:258:GLY:HA2	1:A:262:CYS:SG	2.59	0.42
1:C:189:LEU:HD12	1:C:189:LEU:HA	1.86	0.42
1:A:109:TYR:CG	1:A:151:LEU:HD22	2.55	0.42
1:A:165:TYR:O	1:A:191:TRP:HA	2.20	0.42
1:B:229:ASN:OD1	1:B:240:HIS:HB2	2.19	0.42
1:C:253:ILE:HG12	1:C:273:ARG:HD2	2.01	0.42
1:B:168:THR:C	1:B:170:GLN:H	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:172:TYR:OH	1:C:193:PHE:HA	2.19	0.42
1:C:121:TYR:CZ	1:C:226:TRP:HA	2.54	0.42
1:B:167:ASP:O	1:B:192:ASN:ND2	2.50	0.41
1:A:211:TYR:HB3	1:A:215:GLN:NE2	2.36	0.41
1:C:162:TYR:N	1:C:163:PRO:HD2	2.36	0.41
1:B:126:LYS:HA	1:B:133:LYS:HD3	2.03	0.41
1:A:211:TYR:HB3	1:A:215:GLN:HE22	1.86	0.41
1:B:239:PRO:O	1:B:242:ALA:HB3	2.20	0.41
1:C:174:CYS:HB3	1:C:181:TYR:HE2	1.85	0.41
1:B:92:VAL:HG21	1:B:134:LYS:HA	2.03	0.40
1:B:188:GLN:O	2:B:295:EPE:H101	2.22	0.40
1:C:192:ASN:OD1	1:C:192:ASN:N	2.54	0.40
1:B:243:ILE:HD11	1:B:252:THR:HG21	2.03	0.40
1:C:159:GLU:HA	1:C:162:TYR:CD1	2.56	0.40
1:C:280:PHE:N	1:C:280:PHE:HD1	2.18	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	203/265 (77%)	185 (91%)	17 (8%)	1 (0%)	24	54
1	B	203/265 (77%)	182 (90%)	18 (9%)	3 (2%)	8	29
1	C	203/265 (77%)	188 (93%)	14 (7%)	1 (0%)	24	54
All	All	609/795 (77%)	555 (91%)	49 (8%)	5 (1%)	16	44

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	GLY
1	B	258	GLY

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Mol	Chain	Res	Type
1	B	129	SER
1	B	169	THR
1	C	191	TRP

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	160/203 (79%)	149 (93%)	11 (7%)	14	38
1	B	160/203 (79%)	151 (94%)	9 (6%)	19	45
1	C	160/203 (79%)	150 (94%)	10 (6%)	16	41
All	All	480/609 (79%)	450 (94%)	30 (6%)	16	41

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105	ARG
1	A	130	ASP
1	A	155	LYS
1	A	161	ASN
1	A	178	GLN
1	A	212	LEU
1	A	244	VAL
1	A	271	GLN
1	A	276	LYS
1	A	278	THR
1	A	279	GLN
1	B	130	ASP
1	B	204	ILE
1	B	212	LEU
1	B	215	GLN
1	B	236	THR
1	B	237	MET
1	B	243	ILE
1	B	246	ASN

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Mol	Chain	Res	Type
1	B	269	GLN
1	C	127	THR
1	C	133	LYS
1	C	243	ILE
1	C	269	GLN
1	C	271	GLN
1	C	272	SER
1	C	276	LYS
1	C	278	THR
1	C	281	THR
1	C	293	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	GLN
1	A	194	ASN
1	A	215	GLN
1	A	229	ASN
1	A	232	ASN
1	A	241	ASN
1	A	246	ASN
1	A	257	ASN
1	B	161	ASN
1	B	209	ASN
1	B	215	GLN
1	B	241	ASN
1	B	245	ASN
1	B	246	ASN
1	B	266	ASN
1	C	100	GLN
1	C	178	GLN
1	C	188	GLN
1	C	232	ASN
1	C	279	GLN
1	C	282	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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$
2	EPE	C	295	-	15,15,15	0.87	1 (6%)	19,20,20	1.93	8 (42%)
2	EPE	B	295	-	15,15,15	0.85	1 (6%)	19,20,20	2.16	7 (36%)
2	EPE	A	295	-	15,15,15	0.90	1 (6%)	19,20,20	1.77	6 (31%)

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	EPE	C	295	-	-	3/9/19/19	0/1/1/1
2	EPE	B	295	-	-	6/9/19/19	0/1/1/1
2	EPE	A	295	-	-	4/9/19/19	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	295	EPE	C10-S	3.03	1.81	1.77
2	A	295	EPE	C10-S	2.98	1.81	1.77
2	B	295	EPE	C10-S	2.85	1.81	1.77

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	295	EPE	C5-N4-C3	6.19	122.18	108.84
2	C	295	EPE	C5-N4-C3	4.08	117.62	108.84
2	A	295	EPE	C5-N4-C3	3.43	116.22	108.84
2	C	295	EPE	O2S-S-C10	3.42	111.89	106.73
2	C	295	EPE	O3S-S-C10	3.02	111.91	106.00
2	B	295	EPE	C6-C5-N4	2.91	116.51	110.65
2	B	295	EPE	C7-N4-C3	2.89	118.93	111.24
2	A	295	EPE	O1S-S-C10	2.85	111.03	106.73
2	B	295	EPE	O3S-S-C10	2.75	111.39	106.00
2	A	295	EPE	C7-N4-C5	2.70	118.43	111.24
2	C	295	EPE	C6-C5-N4	2.68	116.05	110.65
2	C	295	EPE	C7-N4-C3	2.67	118.36	111.24
2	B	295	EPE	C7-N4-C5	2.48	117.84	111.24
2	B	295	EPE	O1S-S-C10	2.26	110.15	106.73
2	A	295	EPE	C7-N4-C3	2.25	117.23	111.24
2	C	295	EPE	O3S-S-O2S	-2.18	105.94	111.40
2	A	295	EPE	O3S-S-C10	2.13	110.17	106.00
2	B	295	EPE	C9-N1-C6	-2.11	105.61	111.24
2	C	295	EPE	C9-N1-C6	-2.08	105.69	111.24
2	C	295	EPE	O2S-S-O1S	-2.05	107.15	113.82
2	A	295	EPE	C6-C5-N4	2.05	114.79	110.65

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	295	EPE	C10-C9-N1-C2
2	A	295	EPE	C10-C9-N1-C6
2	B	295	EPE	C10-C9-N1-C2
2	B	295	EPE	S-C10-C9-N1
2	B	295	EPE	C9-C10-S-O1S
2	C	295	EPE	N4-C7-C8-O8
2	B	295	EPE	C8-C7-N4-C3
2	B	295	EPE	C9-C10-S-O3S
2	B	295	EPE	C9-C10-S-O2S
2	A	295	EPE	C8-C7-N4-C3
2	C	295	EPE	C8-C7-N4-C3
2	C	295	EPE	C8-C7-N4-C5
2	A	295	EPE	N4-C7-C8-O8

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	295	EPE	2	0

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.