



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:34 PM EDT

PDB ID : 3PO5 / pdb_00003po5
Title : Structure of a mutant of the large fragment of DNA polymerase I from *Thermus*
Auquaticus in complex with an abasic site and ddATP
Authors : Marx, A.; Diederichs, K.; Obeid, S.
Deposited on : 2010-11-22
Resolution : 2.39 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

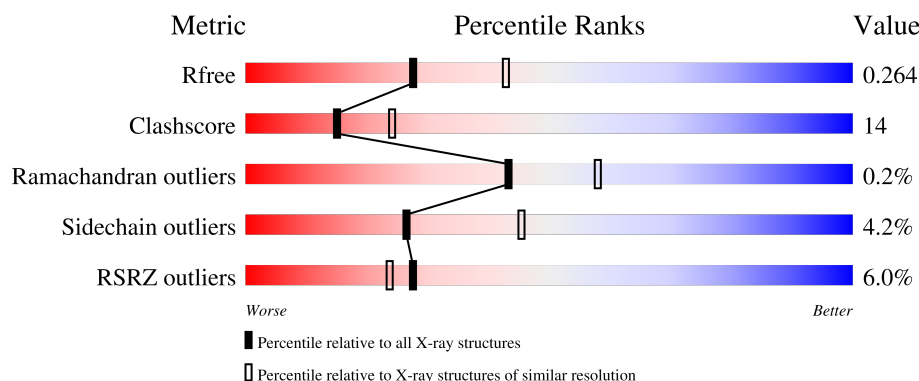
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 $\leq 5\%$. 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	540	
2	B	12	
3	C	13	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 4871 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 DNA polymerase I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	528	Total	C	N	O	S	0	0	0
			4197	2666	756	764	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	614	LYS	ILE	engineered mutation	UNP P19821
A	747	LYS	MET	engineered mutation	UNP P19821

- Molecule 2 is a DNA chain called DNA (5'-D(*GP*AP*CP*CP*AP*CP*GP*GP*CP*GP*CP*(2DA))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	12	Total	C	N	O	P	0	0	0
			242	115	50	66	11			

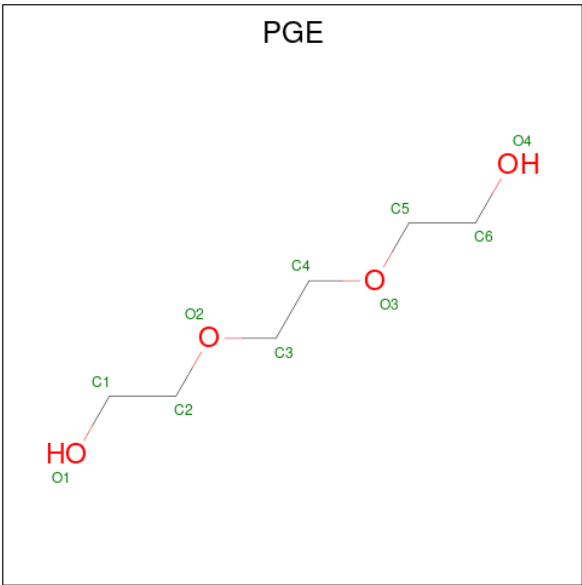
- Molecule 3 is a DNA chain called DNA (5'-D(P*(3DR)P*TP*GP*CP*GP*CP*CP*GP*TP*GP*GP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	P	0	0	0
			257	121	43	80	13			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

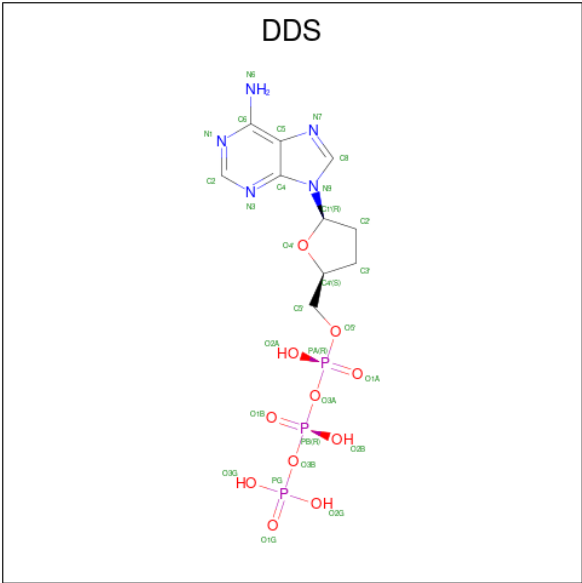
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



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

- Molecule 6 is 2',3'-dideoxyadenosine triphosphate (CCD ID: DDS) (formula: C₁₀H₁₆N₅O₁₁P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			29	10	5	11	3		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			6	3	3		

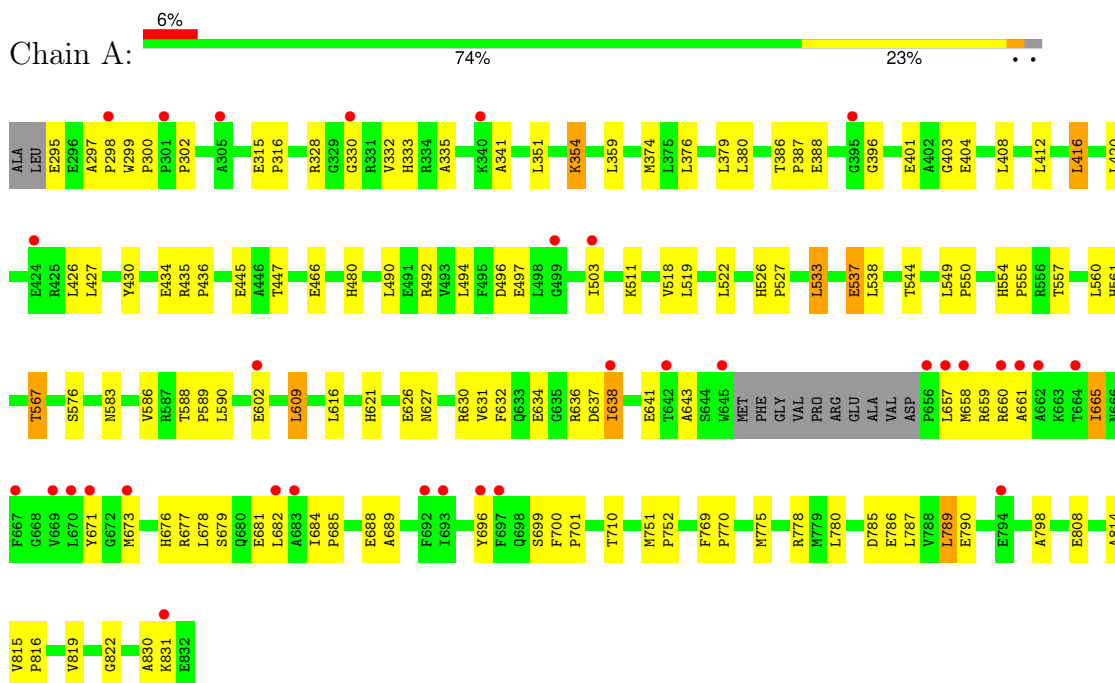
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	96	Total	O	0	0
			96	96		
8	B	12	Total	O	0	0
			12	12		
8	C	23	Total	O	0	0
			23	23		

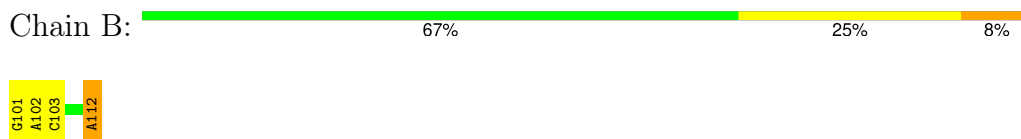
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

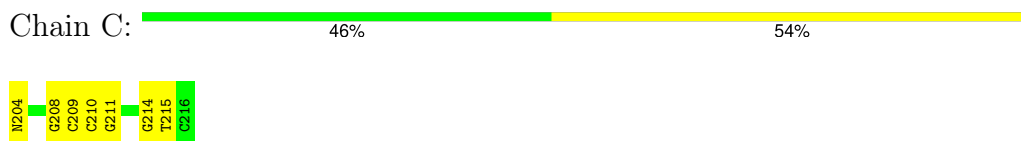
• Molecule 1: DNA polymerase I



• Molecule 2: DNA (5'-D(*GP*AP*CP*CP*AP*CP*GP*GP*CP*GP*CP*(2DA))-3')



• Molecule 3: DNA (5'-D(P*(3DR)P*TP*GP*CP*GP*CP*CP*GP*TP*GP*GP*TP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.30Å 109.30Å 90.32Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.33 – 2.39 47.33 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.33-2.39) 99.7 (47.33-2.39)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_572)	Depositor
R, R_{free}	0.210 , 0.266 0.207 , 0.264	Depositor DCC
R_{free} test set	1287 reflections (3.56%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4871	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DDS, PGE, MG, GOL, 3DR, 2DA

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	0.40	0/4284	0.71	2/5799 (0.0%)
2	B	0.24	0/249	0.83	0/382
3	C	0.26	0/274	0.86	0/421
All	All	0.39	0/4807	0.73	2/6602 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	700	PHE	CA-C-N	5.49	124.94	119.24
1	A	700	PHE	C-N-CA	5.49	124.94	119.24

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	4197	0	4246	117	0
2	B	242	0	134	5	0
3	C	257	0	144	4	0
4	A	1	0	0	0	0
5	A	8	0	8	2	0
6	A	29	0	12	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	6	0	8	0	0
8	A	96	0	0	0	0
8	B	12	0	0	0	0
8	C	23	0	0	0	0
All	All	4871	0	4552	127	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 14.

All (127) 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:B:102:DA:H2''	2:B:103:DC:H5'	1.33	1.05
6:A:835:DDS:H8	6:A:835:DDS:H5'	1.43	0.96
6:A:835:DDS:H5'	6:A:835:DDS:C8	2.02	0.89
2:B:102:DA:C2'	2:B:103:DC:H5'	2.11	0.80
1:A:374:MET:CE	1:A:387:PRO:HA	2.18	0.74
1:A:567:THR:HG22	1:A:567:THR:O	1.87	0.72
1:A:626:GLU:O	1:A:630:ARG:HD3	1.88	0.71
1:A:588:THR:HG23	1:A:589:PRO:HD2	1.73	0.70
1:A:785:ASP:OD1	1:A:786:GLU:HG3	1.93	0.68
1:A:354:LYS:HE2	1:A:445:GLU:OE1	1.94	0.66
1:A:685:PRO:HB2	1:A:688:GLU:HG3	1.79	0.65
1:A:632:PHE:CE1	1:A:638:ILE:HG12	2.32	0.65
1:A:567:THR:O	1:A:567:THR:CG2	2.45	0.64
1:A:302:PRO:HB2	1:A:328:ARG:HD3	1.81	0.62
1:A:376:LEU:HD22	1:A:420:LEU:HD22	1.82	0.62
1:A:386:THR:HG23	1:A:387:PRO:HD2	1.83	0.60
1:A:374:MET:HE1	1:A:387:PRO:HA	1.84	0.60
1:A:638:ILE:N	1:A:638:ILE:HD12	2.17	0.60
2:B:101:DG:H2''	2:B:102:DA:O5'	2.03	0.59
1:A:503:ILE:CD1	1:A:522:LEU:HD11	2.32	0.58
1:A:335:ALA:HB1	1:A:341:ALA:HB2	1.86	0.58
1:A:657:LEU:C	1:A:659:ARG:H	2.12	0.58
1:A:519:LEU:CB	1:A:533:LEU:HD11	2.33	0.57
1:A:815:VAL:HG23	1:A:816:PRO:HD2	1.86	0.56
2:B:102:DA:H2''	2:B:103:DC:C5'	2.24	0.56
1:A:299:TRP:CZ2	1:A:341:ALA:HB1	2.39	0.56
1:A:609:LEU:HD11	1:A:789:LEU:HD22	1.88	0.56
1:A:636:ARG:NE	1:A:641:GLU:HG3	2.21	0.56
1:A:557:THR:OG1	1:A:561:HIS:HE1	1.90	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:661:ALA:O	1:A:665:ILE:HG21	2.07	0.55
1:A:657:LEU:C	1:A:659:ARG:N	2.64	0.55
1:A:447:THR:HG21	1:A:778:ARG:HD3	1.88	0.54
1:A:518:VAL:O	1:A:522:LEU:HD13	2.08	0.54
1:A:576:SER:O	3:C:208:DG:H4'	2.07	0.54
1:A:401:GLU:HG3	1:A:404:GLU:H	1.72	0.54
1:A:374:MET:CE	1:A:387:PRO:CA	2.84	0.53
1:A:627:ASN:O	1:A:631:VAL:HG23	2.09	0.53
1:A:466:GLU:HG2	1:A:538:LEU:HD21	1.90	0.53
1:A:519:LEU:HB2	1:A:533:LEU:HD11	1.89	0.53
1:A:632:PHE:CB	1:A:815:VAL:CG2	2.87	0.52
1:A:780:LEU:HD11	1:A:790:GLU:HB2	1.91	0.52
1:A:299:TRP:CE2	1:A:341:ALA:HB1	2.44	0.51
1:A:609:LEU:CD1	1:A:789:LEU:HD22	2.41	0.51
1:A:588:THR:HG23	1:A:589:PRO:CD	2.41	0.50
1:A:643:ALA:HB1	1:A:658:MET:O	2.11	0.50
1:A:699:SER:C	1:A:701:PRO:HD3	2.37	0.50
1:A:673:MET:HG3	1:A:677:ARG:HB2	1.94	0.50
1:A:661:ALA:O	1:A:665:ILE:CG2	2.59	0.49
1:A:526:HIS:ND1	1:A:527:PRO:HD2	2.27	0.49
1:A:751:MET:HB3	1:A:752:PRO:HD3	1.93	0.49
1:A:299:TRP:CG	1:A:300:PRO:HA	2.46	0.49
1:A:374:MET:HE1	1:A:386:THR:C	2.37	0.49
1:A:549:LEU:HD22	1:A:560:LEU:HD21	1.93	0.49
1:A:678:LEU:O	1:A:682:LEU:HB2	2.13	0.49
1:A:632:PHE:CB	1:A:815:VAL:HG23	2.43	0.49
1:A:480:HIS:HE1	1:A:497:GLU:OE2	1.96	0.48
1:A:630:ARG:N	1:A:630:ARG:HD2	2.27	0.48
1:A:335:ALA:HB1	1:A:341:ALA:CB	2.42	0.48
1:A:295:GLU:HG2	1:A:332:VAL:HG12	1.95	0.48
1:A:632:PHE:HB3	1:A:815:VAL:CG2	2.44	0.48
1:A:386:THR:HG22	1:A:388:GLU:N	2.28	0.48
1:A:537:GLU:OE1	1:A:588:THR:HG21	2.14	0.48
1:A:503:ILE:CD1	1:A:522:LEU:CD1	2.92	0.47
1:A:658:MET:O	1:A:658:MET:SD	2.72	0.47
1:A:671:TYR:HE2	6:A:835:DDS:H2	1.79	0.47
1:A:351:LEU:HD11	1:A:412:LEU:HD12	1.95	0.47
1:A:374:MET:HE1	1:A:386:THR:O	2.14	0.47
1:A:380:LEU:HG	1:A:420:LEU:HD11	1.95	0.47
1:A:643:ALA:HB2	1:A:659:ARG:HA	1.96	0.47
1:A:302:PRO:HD3	1:A:333:HIS:NE2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:643:ALA:CB	1:A:659:ARG:HA	2.45	0.47
1:A:808:GLU:HG2	1:A:819:VAL:HG23	1.96	0.47
1:A:679:SER:HB3	1:A:689:ALA:HB2	1.97	0.47
1:A:549:LEU:HB2	1:A:550:PRO:HD3	1.97	0.46
1:A:616:LEU:HD23	1:A:638:ILE:HG21	1.98	0.46
1:A:632:PHE:HB2	1:A:815:VAL:HG23	1.98	0.46
3:C:210:DC:H2'	3:C:211:DG:C8	2.51	0.46
1:A:297:ALA:HB1	1:A:298:PRO:HD2	1.98	0.46
1:A:676:HIS:O	1:A:677:ARG:C	2.58	0.46
1:A:396:GLY:HA3	1:A:408:LEU:CD1	2.46	0.45
1:A:380:LEU:HD12	1:A:416:LEU:HD12	1.98	0.45
1:A:526:HIS:CG	1:A:527:PRO:HD2	2.52	0.45
1:A:602:GLU:H	1:A:602:GLU:CD	2.25	0.45
1:A:630:ARG:N	1:A:630:ARG:CD	2.80	0.45
5:A:834:PGE:H32	3:C:209:DC:H4'	1.97	0.45
1:A:490:LEU:O	1:A:494:LEU:HD13	2.16	0.45
1:A:632:PHE:HE1	1:A:638:ILE:HG12	1.80	0.45
1:A:822:GLY:HA3	1:A:830:ALA:O	2.17	0.45
1:A:445:GLU:O	1:A:561:HIS:HD2	2.00	0.45
1:A:682:LEU:HD23	1:A:684:ILE:CD1	2.47	0.45
1:A:544:THR:O	5:A:834:PGE:H4	2.18	0.44
6:A:835:DDS:O4'	2:B:112:2DA:H2''	2.17	0.44
1:A:386:THR:CG2	1:A:388:GLU:OE1	2.65	0.44
1:A:434:GLU:OE1	1:A:434:GLU:HA	2.17	0.44
1:A:519:LEU:HB3	1:A:533:LEU:HD11	2.00	0.44
1:A:379:LEU:CD1	1:A:430:TYR:HB2	2.48	0.43
1:A:684:ILE:HB	1:A:685:PRO:HD2	2.00	0.43
1:A:815:VAL:CG2	1:A:816:PRO:HD2	2.46	0.43
1:A:435:ARG:N	1:A:436:PRO:HD2	2.33	0.43
1:A:769:PHE:N	1:A:770:PRO:HD2	2.33	0.43
1:A:637:ASP:O	1:A:641:GLU:CB	2.67	0.43
1:A:374:MET:HE1	1:A:387:PRO:CA	2.48	0.43
1:A:315:GLU:HA	1:A:316:PRO:HD2	1.94	0.43
1:A:359:LEU:C	1:A:359:LEU:HD23	2.43	0.43
1:A:519:LEU:HD13	1:A:533:LEU:CD1	2.50	0.42
1:A:676:HIS:C	1:A:678:LEU:N	2.77	0.42
1:A:665:ILE:HD11	1:A:696:TYR:CD1	2.54	0.42
1:A:374:MET:HE3	1:A:387:PRO:CA	2.49	0.42
1:A:665:ILE:HD11	1:A:696:TYR:CG	2.54	0.42
1:A:684:ILE:HB	1:A:685:PRO:CD	2.50	0.42
1:A:775:MET:HE1	1:A:798:ALA:HB1	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:LYS:NZ	1:A:354:LYS:HB3	2.34	0.41
1:A:621:HIS:CD2	1:A:814:ALA:H	2.38	0.41
1:A:685:PRO:CB	1:A:688:GLU:HG3	2.47	0.41
1:A:492:ARG:O	1:A:496:ASP:HB2	2.21	0.41
1:A:634:GLU:OE1	1:A:636:ARG:HD2	2.20	0.41
3:C:214:DG:C8	3:C:215:DT:H72	2.55	0.41
1:A:637:ASP:O	1:A:641:GLU:HB3	2.21	0.41
1:A:537:GLU:HG3	1:A:538:LEU:N	2.32	0.40
1:A:588:THR:HG22	1:A:590:LEU:H	1.86	0.40
1:A:295:GLU:HG3	1:A:330:GLY:O	2.21	0.40
1:A:588:THR:CG2	1:A:589:PRO:N	2.84	0.40
1:A:490:LEU:HD11	1:A:494:LEU:HD11	2.03	0.40
1:A:554:HIS:HA	1:A:555:PRO:HD3	1.96	0.40
1:A:638:ILE:N	1:A:638:ILE:CD1	2.83	0.40
1:A:778:ARG:HD2	1:A:790:GLU:OE1	2.22	0.40
1:A:401:GLU:CD	1:A:403:GLY:H	2.29	0.40

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	524/540 (97%)	494 (94%)	29 (6%)	1 (0%)	43 58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	586	VAL

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	431/441 (98%)	413 (96%)	18 (4%)	26	45

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	354	LYS
1	A	416	LEU
1	A	426	LEU
1	A	427	LEU
1	A	511	LYS
1	A	533	LEU
1	A	537	GLU
1	A	567	THR
1	A	583	ASN
1	A	609	LEU
1	A	638	ILE
1	A	660	ARG
1	A	665	ILE
1	A	681	GLU
1	A	710	THR
1	A	787	LEU
1	A	789	LEU
1	A	831	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	480	HIS
1	A	534	GLN
1	A	561	HIS
1	A	582	GLN
1	A	583	ASN
1	A	621	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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
3	3DR	C	204	3	8,11,12	0.42	0	7,14,17	1.37	1 (14%)
2	2DA	B	112	3,2	19,22,23	1.63	4 (21%)	25,31,34	2.20	10 (40%)

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
3	3DR	C	204	3	-	0/3/15/16	0/1/1/1
2	2DA	B	112	3,2	-	0/7/18/19	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	112	2DA	C5-C4	4.72	1.47	1.39
2	B	112	2DA	C5-C6	2.64	1.48	1.41
2	B	112	2DA	C5-N7	-2.34	1.34	1.39
2	B	112	2DA	C8-N7	2.21	1.35	1.31

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	112	2DA	C5-C4-N3	-5.30	119.41	126.72
2	B	112	2DA	N3-C4-N9	4.38	134.61	127.17
2	B	112	2DA	C2-N3-C4	3.52	120.42	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	112	2DA	N3-C2-N1	-3.36	123.50	128.58
2	B	112	2DA	C4-C5-N7	-2.92	107.24	110.58
3	C	204	3DR	C1'-C2'-C3'	2.59	106.04	103.26
2	B	112	2DA	C4-N9-C8	2.56	108.42	105.74
2	B	112	2DA	C2'-C1'-N9	-2.47	107.83	112.48
2	B	112	2DA	C5-N7-C8	2.27	107.01	103.45
2	B	112	2DA	C3'-C2'-C1'	2.03	105.21	102.87
2	B	112	2DA	C2-N1-C6	2.03	122.06	118.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	112	2DA	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	$\# Z > 2$
6	DDS	A	835	4	30,31,31	1.29	4 (13%)	42,48,48	2.04	12 (28%)
7	GOL	B	1	-	5,5,5	0.26	0	5,5,5	0.31	0
5	PGE	A	833	-	3,3,9	0.40	0	2,2,8	0.38	0
5	PGE	A	834	-	3,3,9	0.49	0	2,2,8	0.29	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
6	DDS	A	835	4	-	4/22/31/31	0/3/3/3
7	GOL	B	1	-	-	1/4/4/4	-
5	PGE	A	833	-	-	0/1/1/7	-
5	PGE	A	834	-	-	0/1/1/7	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	835	DDS	C6-N6	3.57	1.43	1.34
6	A	835	DDS	C5'-C4'	-2.54	1.42	1.50
6	A	835	DDS	C3'-C2'	-2.34	1.47	1.54
6	A	835	DDS	C5-N7	-2.08	1.35	1.39

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	835	DDS	C5-C4-N3	-5.83	118.69	126.72
6	A	835	DDS	N3-C2-N1	-4.60	121.62	128.58
6	A	835	DDS	N3-C4-N9	4.02	134.01	127.17
6	A	835	DDS	C2-N3-C4	3.72	120.92	111.83
6	A	835	DDS	C3'-C2'-C1'	3.47	106.87	102.87
6	A	835	DDS	C4-C5-N7	-3.02	107.13	110.58
6	A	835	DDS	C5-N7-C8	2.98	108.13	103.45
6	A	835	DDS	N9-C8-N7	-2.88	109.85	113.94
6	A	835	DDS	C4'-O4'-C1'	-2.46	107.49	109.81
6	A	835	DDS	O2B-PB-O3A	2.38	113.71	107.27
6	A	835	DDS	O3A-PB-O1B	-2.18	104.16	110.70
6	A	835	DDS	O3G-PG-O3B	2.13	111.76	104.64

There are no chirality outliers.

All (5) torsion outliers are listed below:

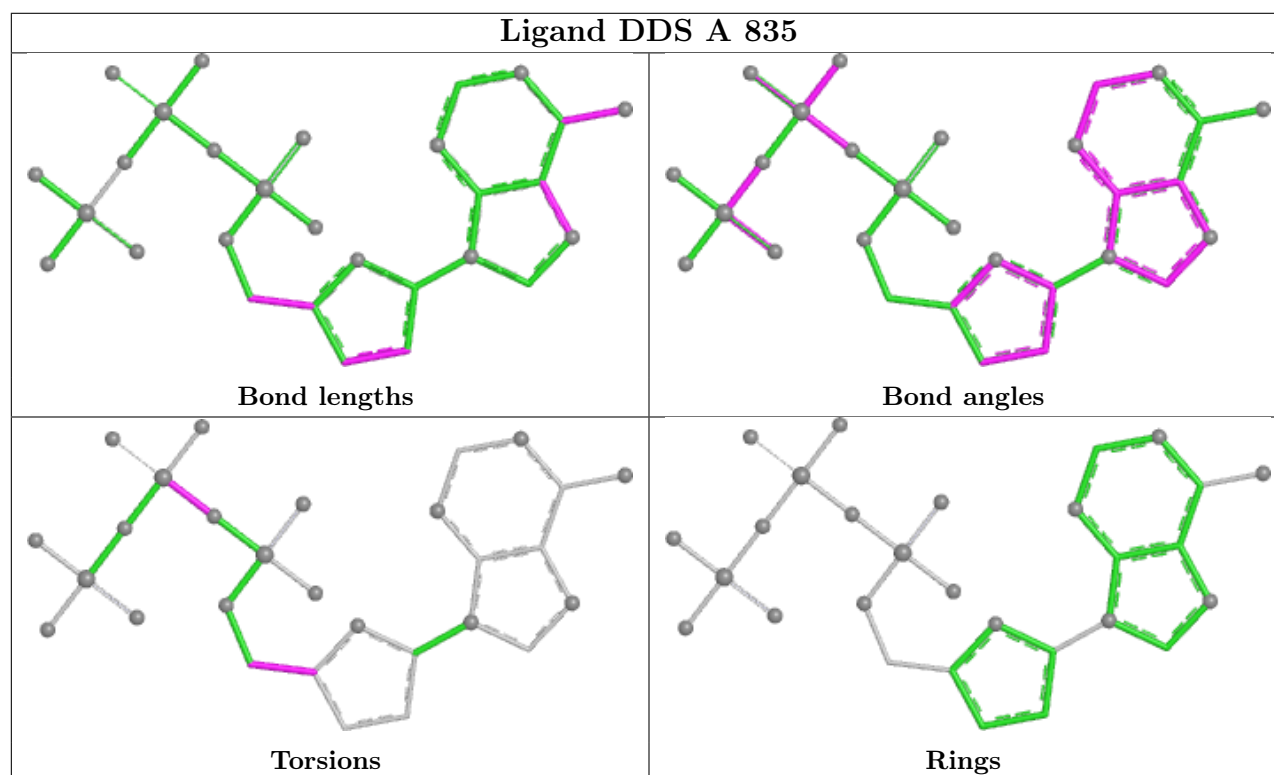
Mol	Chain	Res	Type	Atoms
6	A	835	DDS	C3'-C4'-C5'-O5'
6	A	835	DDS	O4'-C4'-C5'-O5'
7	B	1	GOL	O1-C1-C2-O2
6	A	835	DDS	PA-O3A-PB-O2B
6	A	835	DDS	PA-O3A-PB-O1B

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	835	DDS	4	0
5	A	834	PGE	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

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	528/540 (97%)	0.45	33 (6%) 26 22	26, 49, 105, 132	0
2	B	11/12 (91%)	-0.65	0 100 100	28, 32, 62, 71	0
3	C	12/13 (92%)	-0.93	0 100 100	24, 32, 51, 54	0
All	All	551/565 (97%)	0.40	33 (5%) 27 24	24, 48, 105, 132	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	658	MET	4.5
1	A	656	PRO	3.9
1	A	657	LEU	3.9
1	A	330	GLY	3.9
1	A	667	PHE	3.8
1	A	499	GLY	3.5
1	A	682	LEU	3.2
1	A	673	MET	3.1
1	A	645	TRP	3.0
1	A	661	ALA	2.9
1	A	298	PRO	2.7
1	A	697	PHE	2.6
1	A	395	GLY	2.6
1	A	670	LEU	2.5
1	A	660	ARG	2.5
1	A	602	GLU	2.4
1	A	693	ILE	2.4
1	A	696	TYR	2.4
1	A	642	THR	2.3
1	A	305	ALA	2.3
1	A	664	THR	2.3
1	A	503	ILE	2.2
1	A	340	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	671	TYR	2.2
1	A	662	ALA	2.2
1	A	831	LYS	2.1
1	A	669	VAL	2.1
1	A	424	GLU	2.1
1	A	794	GLU	2.1
1	A	692	PHE	2.1
1	A	301	PRO	2.0
1	A	683	ALA	2.0
1	A	638	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors(Å ²)	Q<0.9
3	3DR	C	204	11/12	0.84	0.13	66,67,85,86	0
2	2DA	B	112	20/21	0.97	0.05	27,30,40,42	0

6.3 Carbohydrates [i](#)

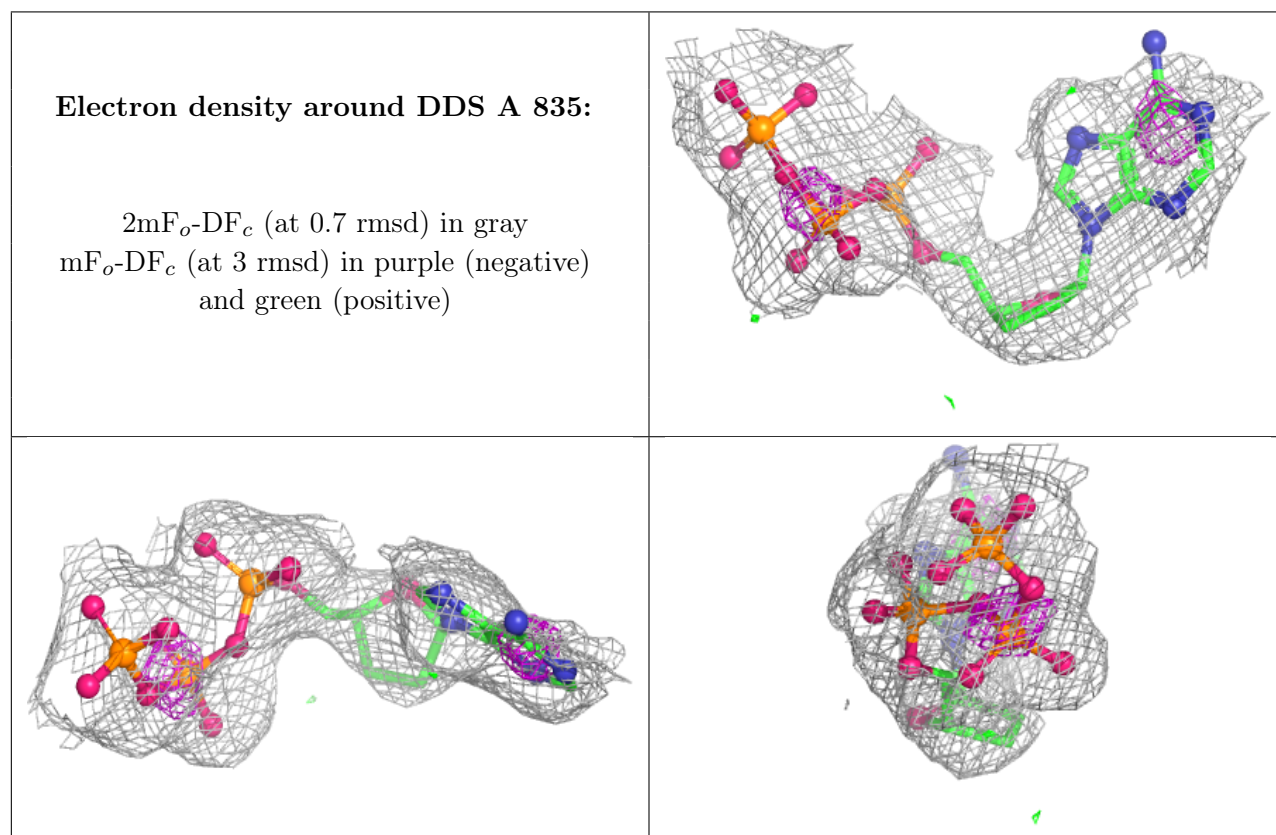
There are no oligosaccharides in this entry.

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. 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	B-factors(Å ²)	Q<0.9
4	MG	A	1	1/1	0.78	0.12	64,64,64,64	0
7	GOL	B	1	6/6	0.79	0.13	43,44,44,44	0
6	DDS	A	835	29/29	0.86	0.12	50,72,94,271	0
5	PGE	A	833	4/10	0.90	0.10	46,47,47,47	0
5	PGE	A	834	4/10	0.94	0.08	29,29,30,30	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.