



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 02:33 PM EDT

PDB ID : 1B4Z / pdb_00001b4z
Title : OLIGO-PEPTIDE BINDING PROTEIN (OPPA) COMPLEXED WITH KDK
Authors : Tame, J.R.H.; Sleigh, S.H.; Wilkinson, A.J.
Deposited on : 1999-01-04
Resolution : 1.75 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)	:	2.50

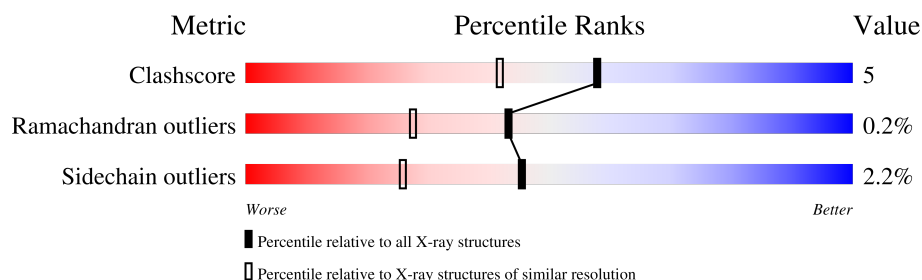
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)

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	517	
2	B	3	

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
4	ACT	A	528	-	-	X	-
4	ACT	A	529	-	-	X	-
4	ACT	A	530	-	-	X	-
4	ACT	A	531	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4690 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.

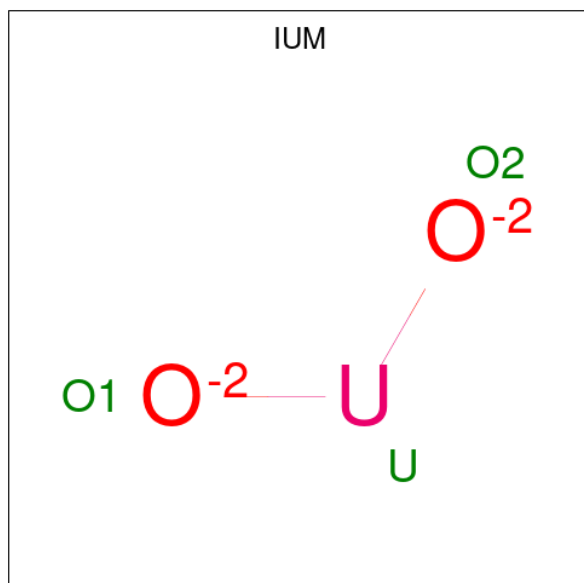
- Molecule 1 is a protein called PROTEIN (OLIGO-PEPTIDE BINDING PROTEIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	517	Total	C	N	O	S	0	3	0
			4175	2675	700	794	6			

- Molecule 2 is a protein called PROTEIN (PEPTIDE LYS-ASP-LYS).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	3	Total	C	N	O	0	0	0
			27	16	5	6			

- Molecule 3 is URANYL (VI) ION (CCD ID: IUM) (formula: O₂U).



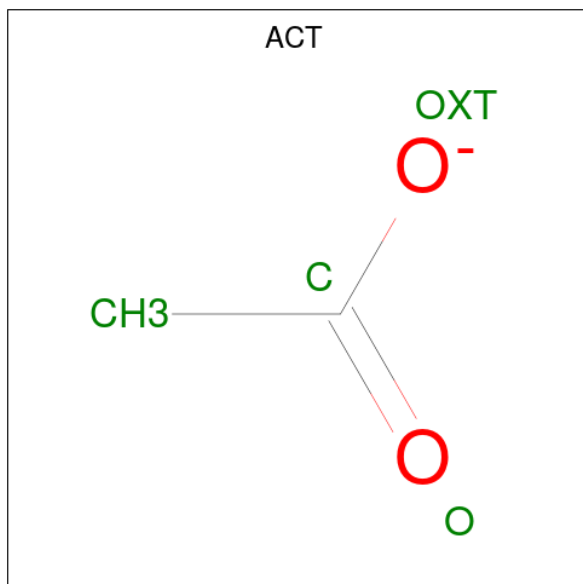
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	U	0	0
			1	1		
3	A	1	Total	U	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0
3	A	1	Total U 1 1	0	0

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 3 1 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0

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

- Molecule 5 is water.

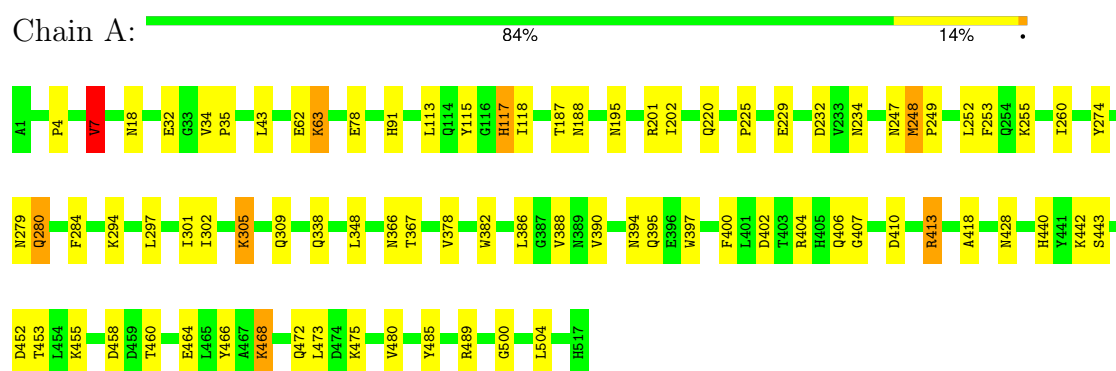
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	452	Total	O	0	0
			452	452		
5	B	4	Total	O	0	1
			5	5		

3 Residue-property plots

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. 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. 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: PROTEIN (OLIGO-PEPTIDE BINDING PROTEIN)



• Molecule 2: PROTEIN (PEPTIDE LYS-ASP-LYS)



There are no outlier residues recorded for this chain.

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	110.01Å 75.55Å 70.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 1.75	Depositor
% Data completeness (in resolution range)	92.2 (15.00-1.75)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.179 , 0.214	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4690	wwPDB-VP
Average B, all atoms (Å ²)	21.0	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: IUM, ACT

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	0.82	0/4300	1.53	45/5863 (0.8%)
2	B	0.87	0/26	1.33	0/30
All	All	0.82	0/4326	1.53	45/5893 (0.8%)

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	ASN	CA-CB-CG	11.65	124.25	112.60
1	A	280	GLN	OE1-CD-NE2	8.33	130.93	122.60
1	A	413	ARG	NE-CZ-NH2	-7.85	112.14	119.20
1	A	443	SER	CA-C-O	6.95	126.65	119.49
1	A	18	ASN	CA-CB-CG	-6.83	105.77	112.60
1	A	284	PHE	CA-CB-CG	-6.80	107.00	113.80
1	A	32	GLU	CA-C-N	-6.29	115.90	122.69
1	A	32	GLU	C-N-CA	-6.29	115.90	122.69
1	A	407	GLY	N-CA-C	-6.11	104.38	114.76
1	A	485	TYR	CA-CB-CG	6.03	124.75	113.90
1	A	394	ASN	CB-CG-ND2	-5.93	107.50	116.40
1	A	252	LEU	N-CA-CB	-5.89	103.07	110.90
1	A	201	ARG	CD-NE-CZ	5.80	132.51	124.40
1	A	188	ASN	CA-C-N	-5.79	116.18	122.55
1	A	188	ASN	C-N-CA	-5.79	116.18	122.55
1	A	91	HIS	CA-CB-CG	-5.76	108.04	113.80
1	A	280	GLN	CG-CD-NE2	-5.63	107.95	116.40
1	A	274	TYR	CA-C-N	-5.56	115.37	122.77
1	A	274	TYR	C-N-CA	-5.56	115.37	122.77
1	A	500	GLY	N-CA-C	-5.55	107.97	115.36
1	A	202	ILE	CA-C-N	-5.53	115.75	123.10
1	A	202	ILE	C-N-CA	-5.53	115.75	123.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	279	ASN	OD1-CG-ND2	-5.48	117.12	122.60
1	A	260	ILE	CA-C-N	5.44	125.09	119.05
1	A	260	ILE	C-N-CA	5.44	125.09	119.05
1	A	407	GLY	CA-C-N	-5.39	114.37	122.24
1	A	407	GLY	C-N-CA	-5.39	114.37	122.24
1	A	7	VAL	N-CA-CB	5.38	117.63	110.26
1	A	117	HIS	CA-CB-CG	-5.29	108.51	113.80
1	A	367	THR	CA-CB-OG1	-5.27	101.70	109.60
1	A	473	LEU	O-C-N	-5.26	116.54	122.12
1	A	248[A]	MET	CA-C-N	5.22	124.98	119.76
1	A	248[A]	MET	C-N-CA	5.22	124.98	119.76
1	A	248[B]	MET	CA-C-N	5.22	124.98	119.76
1	A	248[B]	MET	C-N-CA	5.22	124.98	119.76
1	A	232	ASP	CA-CB-CG	5.21	117.81	112.60
1	A	400	PHE	CA-CB-CG	5.17	118.97	113.80
1	A	410	ASP	CA-CB-CG	-5.16	107.44	112.60
1	A	458	ASP	CA-CB-CG	5.15	117.75	112.60
1	A	458	ASP	N-CA-CB	-5.15	102.37	110.77
1	A	397	TRP	N-CA-C	5.08	119.70	111.37
1	A	386	LEU	N-CA-CB	-5.06	104.30	110.98
1	A	220	GLN	CG-CD-NE2	5.05	123.98	116.40
1	A	234	ASN	N-CA-CB	5.01	117.27	110.01
1	A	195	ASN	CA-C-O	-5.00	114.60	120.60

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	4175	0	4095	31	0
2	B	27	0	32	0	0
3	A	8	0	0	0	0
4	A	23	0	15	7	0
5	A	452	0	0	3	0
5	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4690	0	4142	38	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 5.

All (38) 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:A:229:GLU:HB3	1:A:249:PRO:HD3	1.74	0.68
1:A:248[B]:MET:HE3	1:A:253:PHE:HD1	1.59	0.68
4:A:528:ACT:H3	4:A:529:ACT:H2	1.78	0.65
1:A:452:ASP:HA	1:A:455:LYS:HE2	1.78	0.65
4:A:528:ACT:H3	4:A:529:ACT:CH3	2.28	0.64
1:A:113:LEU:HD22	1:A:118:ILE:HD12	1.85	0.58
1:A:402:ASP:OD2	1:A:406:GLN:NE2	2.36	0.56
1:A:418:ALA:HB3	1:A:504:LEU:HD22	1.87	0.56
4:A:530:ACT:H2	4:A:531:ACT:CH3	2.37	0.55
4:A:528:ACT:CH3	4:A:529:ACT:H3	2.37	0.55
1:A:382:TRP:HB3	1:A:388:VAL:CG2	2.38	0.54
1:A:117:HIS:HE1	5:A:744:HOH:O	1.91	0.53
1:A:382:TRP:HB3	1:A:388:VAL:HG22	1.91	0.53
1:A:455:LYS:HD3	5:A:926:HOH:O	2.09	0.52
1:A:460:THR:O	1:A:464:GLU:HG3	2.09	0.52
4:A:530:ACT:H2	4:A:531:ACT:H3	1.92	0.52
1:A:297[B]:LEU:HD11	1:A:378:VAL:HG22	1.92	0.51
1:A:253:PHE:CD2	1:A:309:GLN:HG2	2.45	0.51
4:A:528:ACT:H2	4:A:529:ACT:H3	1.93	0.51
1:A:43:LEU:O	1:A:187:THR:HB	2.12	0.49
4:A:528:ACT:CH3	4:A:529:ACT:CH3	2.92	0.48
1:A:404:ARG:O	1:A:440:HIS:HE1	1.97	0.47
1:A:4:PRO:O	1:A:7:VAL:HG13	2.14	0.47
1:A:297[B]:LEU:HD11	1:A:302:ILE:HD11	1.96	0.47
1:A:294:LYS:HA	1:A:480:VAL:HG13	1.98	0.45
1:A:453:THR:HG21	1:A:466:TYR:CE1	2.51	0.45
1:A:115:TYR:CE1	1:A:428:ASN:HB3	2.51	0.45
1:A:248[B]:MET:HE3	1:A:253:PHE:CD1	2.46	0.44
1:A:280:GLN:HE21	1:A:442:LYS:HE2	1.83	0.43
1:A:301:ILE:HA	1:A:305:LYS:HG3	1.99	0.43
1:A:489:ARG:HG2	5:A:982:HOH:O	2.17	0.43
1:A:472:GLN:HA	1:A:475:LYS:HG2	2.00	0.43
1:A:464:GLU:O	1:A:468:LYS:HG2	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297[B]:LEU:HD11	1:A:302:ILE:CD1	2.50	0.42
1:A:62:GLU:HG3	1:A:63:LYS:HD3	2.02	0.41
1:A:34:VAL:HB	1:A:35:PRO:HD3	2.02	0.41
1:A:113:LEU:CD2	1:A:118:ILE:HD12	2.50	0.41
1:A:366:ASN:HA	1:A:395:GLN:O	2.21	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	518/517 (100%)	503 (97%)	14 (3%)	1 (0%)	43	27
2	B	1/3 (33%)	1 (100%)	0	0	100	100
All	All	519/520 (100%)	504 (97%)	14 (3%)	1 (0%)	43	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	225	PRO

5.3.2 Protein sidechains [i](#)

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	458/455 (101%)	448 (98%)	10 (2%)	45	26
2	B	3/3 (100%)	3 (100%)	0	100	100
All	All	461/458 (101%)	451 (98%)	10 (2%)	45	26

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	63	LYS
1	A	78	GLU
1	A	255	LYS
1	A	305	LYS
1	A	338	GLN
1	A	348	LEU
1	A	390	VAL
1	A	413	ARG
1	A	468	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	117	HIS
1	A	184	ASN
1	A	199	ASN
1	A	279	ASN
1	A	338	GLN
1	A	391	ASN
1	A	395	GLN
1	A	440	HIS

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](#)

Of 14 ligands modelled in this entry, 8 are modelled with single atom - leaving 6 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
4	ACT	A	531	3	3,3,3	0.86	0	3,3,3	1.57	1 (33%)
4	ACT	A	526	3	2,2,3	0.57	0	1,1,3	0.73	0
4	ACT	A	527	3	3,3,3	0.83	0	3,3,3	2.11	2 (66%)
4	ACT	A	530	3	3,3,3	0.83	0	3,3,3	0.95	0
4	ACT	A	528	3	3,3,3	0.79	0	3,3,3	0.49	0
4	ACT	A	529	3	3,3,3	1.17	0	3,3,3	2.17	1 (33%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	529	ACT	OXT-C-O	3.15	133.72	122.03
4	A	527	ACT	OXT-C-CH3	2.62	126.05	115.05
4	A	527	ACT	O-C-CH3	-2.54	112.13	122.53
4	A	531	ACT	O-C-CH3	-2.04	114.17	122.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	531	ACT	2	0
4	A	530	ACT	2	0
4	A	528	ACT	5	0
4	A	529	ACT	5	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.