



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 04:32 AM GMT

PDB ID : 3E6Y  
Title : Structure of 14-3-3 in complex with the differentiation-inducing agent  
Cotylenin A  
Authors : Ottmann, C.; Weyand, M.; Wittinghofer, A.; Oecking, C.  
Deposited on : 2008-08-17  
Resolution : 2.50 Å (reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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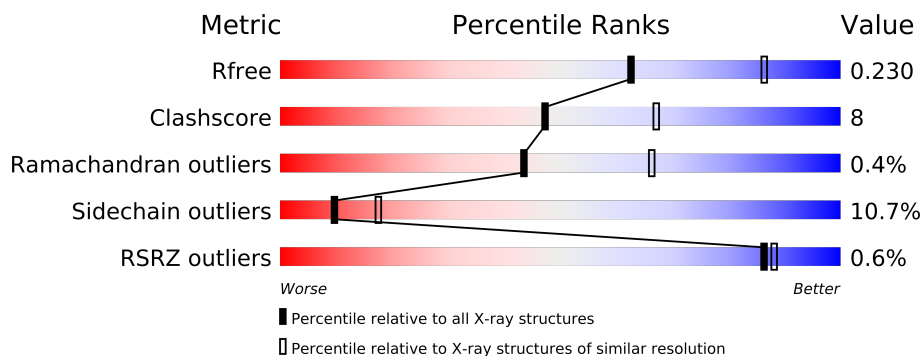
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.50 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	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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  $\geq 3$ , 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	260	
1	B	260	
2	C	5	
2	D	5	

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	CL	A	262	-	X

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 3926 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.

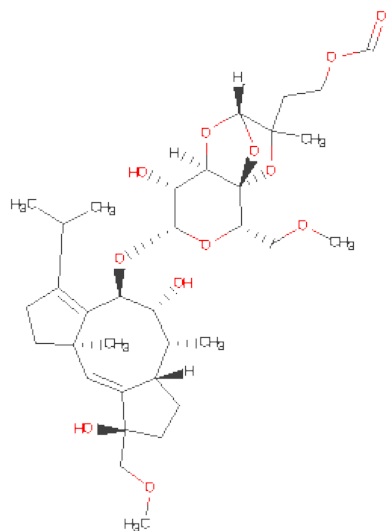
- Molecule 1 is a protein called 14-3-3-like protein C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1840	1156	309	368	7			
1	B	232	Total	C	N	O	S	0	0	0
			1801	1130	302	362	7			

- Molecule 2 is a protein called H<sup>+</sup>-ATPase phosphopeptide QSYpTV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	5	Total	C	N	O	P	0	0	0
			46	26	6	13	1			
2	D	5	Total	C	N	O	P	0	0	0
			46	26	6	13	1			

- Molecule 3 is COTYLENIN A (three-letter code: CW1) (formula: C<sub>34</sub>H<sub>52</sub>O<sub>12</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			46	34	12		
3	B	1	Total	C	O	0	0
			46	34	12		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	55	Total	O	0	0
			55	55		
5	B	42	Total	O	0	0
			42	42		
5	C	3	Total	O	0	0
			3	3		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	41.80Å 141.90Å 51.40Å 90.00° 114.10° 90.00°	Depositor
Resolution (Å)	14.95 – 2.50 14.95 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (14.95-2.50) 98.4 (14.95-2.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.62 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.5.0035	Depositor
R, $R_{free}$	0.195 , 0.285 0.192 , 0.230	Depositor DCC
$R_{free}$ test set	930 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.4	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 14.8	EDS
Estimated twinning fraction	0.177 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 18588 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3926	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.29% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities 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: TPO, CW1, 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	0.86	0/1868	0.97	5/2524 (0.2%)
1	B	0.83	0/1828	0.82	3/2471 (0.1%)
2	C	1.09	0/34	0.73	0/42
2	D	0.87	0/34	0.69	0/42
All	All	0.85	0/3764	0.90	8/5079 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	ARG	NE-CZ-NH1	12.30	126.45	120.30
1	A	63	ARG	NE-CZ-NH2	-9.47	115.56	120.30
1	A	110	LEU	CA-CB-CG	6.78	130.89	115.30
1	A	38	LEU	CA-CB-CG	6.10	129.34	115.30
1	B	110	LEU	CA-CB-CG	6.02	129.15	115.30
1	B	89	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	92	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	A	232	ASP	CB-CG-OD2	-5.05	113.75	118.30

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	1840	0	1787	26	0
1	B	1801	0	1720	30	0
2	C	46	0	38	4	0
2	D	46	0	38	1	0
3	A	46	0	52	3	0
3	B	46	0	52	5	0
4	A	1	0	0	1	0
5	A	55	0	0	2	0
5	B	42	0	0	3	0
5	C	3	0	0	1	0
All	All	3926	0	3687	60	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 8.

All (60) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:233:ASN:HD21	2:C:3:TYR:H	1.27	0.79
1:B:128:LEU:HB3	1:B:161:ALA:HB2	1.70	0.74
1:B:173:HIS:HD2	1:B:175:ILE:H	1.33	0.74
1:A:96:GLU:HG2	1:A:141:PHE:CD2	2.30	0.67
1:B:73:GLU:HB3	1:B:88:ILE:HG21	1.77	0.66
1:B:162:GLN:O	1:B:166:THR:HG22	1.96	0.65
1:A:181:LEU:CD1	2:C:4:TPO:HA	2.26	0.65
1:A:69:ILE:HD13	1:B:21:ALA:HB2	1.80	0.63
1:B:156:THR:HG23	1:B:157:ALA:N	2.16	0.61
1:B:52:SER:OG	5:B:275:HOH:O	2.15	0.61
1:B:76:GLU:HG3	1:B:81:ASN:HB2	1.82	0.60
1:A:53:VAL:HG21	3:A:261:CW1:HAA	1.83	0.60
1:B:187:TYR:HA	1:B:191:LEU:HB2	1.84	0.60
1:A:76:GLU:OE1	1:A:81:ASN:ND2	2.28	0.60
1:B:173:HIS:CD2	1:B:175:ILE:H	2.19	0.58
1:B:118:ALA:HB1	1:B:123:SER:HB2	1.86	0.57
1:A:35:SER:O	1:A:38:LEU:HG	2.05	0.57
1:A:199:ASN:HB3	5:A:264:HOH:O	2.07	0.55
1:B:156:THR:CG2	1:B:157:ALA:N	2.71	0.54
3:B:261:CW1:HBP	3:B:261:CW1:CBF	2.38	0.53
1:B:156:THR:CG2	1:B:157:ALA:H	2.22	0.53
1:A:173:HIS:HE1	1:A:175:ILE:HD12	1.74	0.53
1:A:20:GLN:HG2	1:B:72:ILE:HD11	1.91	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:209:ILE:HG22	5:A:276:HOH:O	2.08	0.53
1:B:110:LEU:HD21	1:B:130:MET:HB3	1.92	0.52
1:B:12:VAL:HA	1:B:31:MET:CE	2.41	0.51
1:A:69:ILE:CD1	1:B:21:ALA:HB2	2.40	0.50
1:B:164:ILE:O	1:B:168:GLU:HB2	2.12	0.50
1:B:205:PHE:O	1:B:209:ILE:HG12	2.12	0.49
1:B:153:SER:O	1:B:156:THR:HG22	2.13	0.49
3:B:261:CW1:CAA	5:B:264:HOH:O	2.61	0.48
3:B:261:CW1:HAAB	5:B:264:HOH:O	2.12	0.48
1:A:132:GLY:HA3	1:A:158:TYR:CE1	2.48	0.48
1:B:73:GLU:HB3	1:B:88:ILE:CG2	2.43	0.48
1:A:63:ARG:HG3	4:A:262:CL:CL	2.50	0.48
1:A:173:HIS:CE1	1:A:175:ILE:HD12	2.49	0.48
3:B:261:CW1:HBK	3:B:261:CW1:HAAA	1.63	0.47
1:B:25:GLU:O	1:B:29:GLU:HG3	2.13	0.47
1:A:236:LEU:O	1:A:236:LEU:HD12	2.15	0.46
1:B:120:SER:OG	1:B:122:ASP:HB2	2.15	0.46
1:A:228:GLN:HA	1:A:231:ARG:HD2	1.97	0.46
1:B:142:LYS:O	1:B:147:ARG:HG3	2.15	0.46
1:A:5:PRO:O	1:A:9:GLU:OE1	2.34	0.45
1:A:31:MET:O	1:A:35:SER:HB2	2.16	0.45
2:C:1:GLN:HG2	5:C:27:HOH:O	2.15	0.45
1:B:164:ILE:O	1:B:168:GLU:CB	2.64	0.45
1:A:115:ILE:HB	1:A:116:PRO:HD3	1.98	0.45
3:A:261:CW1:CBF	3:A:261:CW1:HBP	2.47	0.45
3:B:261:CW1:CBF	3:B:261:CW1:CBP	2.95	0.44
1:B:172:THR:O	1:B:177:LEU:HD12	2.18	0.44
1:A:203:GLN:O	1:A:207:GLU:HG3	2.18	0.43
1:A:63:ARG:NH2	2:C:4:TPO:O1P	2.52	0.42
1:B:81:ASN:O	1:B:85:VAL:HG23	2.19	0.42
1:A:21:ALA:HB2	1:B:69:ILE:CD1	2.49	0.42
1:B:63:ARG:HH22	2:D:4:TPO:P	2.43	0.41
3:A:261:CW1:CBF	3:A:261:CW1:CBP	2.98	0.41
1:B:187:TYR:CD1	1:B:191:LEU:HD12	2.55	0.41
1:A:187:TYR:CD1	1:A:191:LEU:HD12	2.55	0.41
1:A:155:LEU:HD13	1:A:186:PHE:HE2	1.86	0.41
1:A:220:TYR:O	1:A:224:THR:HG23	2.22	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	232/260 (89%)	223 (96%)	8 (3%)	1 (0%)	43	66
1	B	228/260 (88%)	210 (92%)	17 (8%)	1 (0%)	43	66
2	C	2/5 (40%)	2 (100%)	0	0	100	100
2	D	2/5 (40%)	2 (100%)	0	0	100	100
All	All	464/530 (88%)	437 (94%)	25 (5%)	2 (0%)	43	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	B	39	GLY

### 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	191/221 (86%)	169 (88%)	22 (12%)	8	15
1	B	183/221 (83%)	166 (91%)	17 (9%)	13	23
2	C	4/4 (100%)	3 (75%)	1 (25%)	1	1
2	D	4/4 (100%)	3 (75%)	1 (25%)	1	1
All	All	382/450 (85%)	341 (89%)	41 (11%)	10	17

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU

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Mol	Chain	Res	Type
1	A	20	GLN
1	A	32	GLU
1	A	35	SER
1	A	36	ASN
1	A	37	SER
1	A	43	LEU
1	A	56	LYS
1	A	110	LEU
1	A	123	SER
1	A	149	GLU
1	A	156	THR
1	A	166	THR
1	A	177	LEU
1	A	181	LEU
1	A	183	PHE
1	A	196	ARG
1	A	209	ILE
1	A	211	GLU
1	A	215	LEU
1	A	218	GLU
1	A	225	LEU
1	B	32	GLU
1	B	52	SER
1	B	70	SER
1	B	76	GLU
1	B	78	SER
1	B	82	GLU
1	B	93	SER
1	B	110	LEU
1	B	122	ASP
1	B	123	SER
1	B	149	GLU
1	B	166	THR
1	B	168	GLU
1	B	183	PHE
1	B	219	SER
1	B	225	LEU
1	B	232	ASP
2	C	1	GLN
2	D	1	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	199	ASN
1	A	233	ASN
1	B	49	ASN
1	B	74	GLN
1	B	173	HIS
1	B	233	ASN
2	D	1	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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	$\# Z  > 2$
2	TPO	C	4	2	10,10,11	5.14	1 (10%)	12,14,16	1.22	0
2	TPO	D	4	2	10,10,11	5.08	1 (10%)	12,14,16	1.61	3 (25%)

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	TPO	C	4	2	-	0/9/11/13	0/0/0/0
2	TPO	D	4	2	-	0/9/11/13	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	TPO	O-C	16.06	1.22	1.11
2	D	4	TPO	O-C	15.82	1.22	1.11

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	TPO	CG2-CB-CA	-2.76	107.54	113.20
2	D	4	TPO	OG1-P-O1P	-2.68	99.32	106.79
2	D	4	TPO	O3P-P-O1P	2.06	117.19	110.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z  > 2$
3	CW1	A	261	-	51,51,51	2.77	12 (23%)	81,81,81	4.28	34 (41%)
3	CW1	B	261	-	51,51,51	2.64	14 (27%)	81,81,81	3.87	23 (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
3	CW1	A	261	-	1/1/17/19	0/21/119/119	0/0/6/6
3	CW1	B	261	-	1/1/17/19	0/21/119/119	0/0/6/6

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	261	CW1	CAM-CBE	11.17	1.42	1.33
3	A	261	CW1	CAM-CBE	10.29	1.41	1.33
3	A	261	CW1	CBR-CBM	9.35	1.67	1.54
3	B	261	CW1	CBR-CBM	6.74	1.63	1.54
3	B	261	CW1	CBD-CBF	6.38	1.41	1.33
3	A	261	CW1	CBD-CBF	6.36	1.41	1.33
3	A	261	CW1	CBO-CAM	-4.92	1.42	1.51
3	B	261	CW1	OBC-CBR	4.81	1.55	1.45
3	A	261	CW1	OBB-CBM	-4.60	1.34	1.41
3	B	261	CW1	CBG-CBD	4.03	1.55	1.51
3	A	261	CW1	CBG-CBD	4.01	1.55	1.51
3	A	261	CW1	CBT-CBN	-3.42	1.49	1.54
3	B	261	CW1	CBT-CBN	-3.32	1.49	1.54
3	B	261	CW1	CBO-CAM	-3.25	1.45	1.51
3	A	261	CW1	OBC-CBR	2.99	1.52	1.45
3	A	261	CW1	CBT-CBK	2.99	1.58	1.53
3	A	261	CW1	OAX-CAN	-2.74	1.40	1.44
3	B	261	CW1	OBB-CBT	-2.66	1.39	1.42
3	B	261	CW1	CBP-CBH	2.63	1.58	1.55
3	B	261	CW1	CAU-CBS	-2.60	1.48	1.53
3	B	261	CW1	CBQ-CBF	2.58	1.55	1.50
3	B	261	CW1	OAY-CBL	2.57	1.48	1.41
3	A	261	CW1	OAZ-CBM	2.56	1.46	1.41
3	B	261	CW1	OAY-CBK	-2.49	1.40	1.44
3	B	261	CW1	OAX-CAN	2.32	1.47	1.44
3	A	261	CW1	CAS-CBS	-2.31	1.49	1.53

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	261	CW1	OBC-CBR-CAQ	21.05	126.67	107.69
3	B	261	CW1	OBC-CBR-CAQ	18.96	124.79	107.69
3	A	261	CW1	OAZ-CBM-OBB	-14.42	90.97	107.96
3	B	261	CW1	CAG-CBR-CBM	-13.15	85.88	114.45
3	A	261	CW1	CAQ-CBR-CBM	-11.68	81.46	113.06
3	B	261	CW1	CAQ-CBR-CBM	-10.92	83.53	113.06
3	A	261	CW1	CAG-CBR-CBM	-10.64	91.33	114.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	261	CW1	OBC-CBR-CBM	-10.40	89.92	99.74
3	B	261	CW1	OAZ-CBM-OB	-8.67	97.75	107.96
3	A	261	CW1	OBC-CBR-CBM	-7.98	92.20	99.74
3	A	261	CW1	OAZ-CBM-CBR	7.74	125.20	108.40
3	B	261	CW1	CAS-CBS-CBE	6.37	108.87	102.27
3	B	261	CW1	OBC-CBT-CBK	6.19	122.29	108.51
3	A	261	CW1	CBT-CBN-CBI	-6.16	106.20	114.31
3	A	261	CW1	OBC-CBT-CBK	5.97	121.81	108.51
3	B	261	CW1	CBS-CBE-CAM	5.35	129.77	121.28
3	B	261	CW1	CAT-CBK-CBT	5.31	122.67	113.49
3	A	261	CW1	OAZ-CBN-CBI	5.19	118.79	109.43
3	B	261	CW1	CAB-OAW-CAU	-5.12	102.08	112.26
3	A	261	CW1	OAY-CBK-CAT	4.89	118.74	107.33
3	B	261	CW1	OAZ-CBM-CBR	4.69	118.59	108.40
3	B	261	CW1	CBT-CBN-CBI	-4.56	108.29	114.31
3	A	261	CW1	CAS-CBS-CBE	4.50	106.93	102.27
3	A	261	CW1	CAD-CBG-CBD	4.48	116.87	111.57
3	A	261	CW1	OAZ-CBN-CBT	-4.41	96.94	103.81
3	B	261	CW1	OBC-CBR-CAG	4.22	117.86	108.71
3	A	261	CW1	CBS-CBE-CAM	4.04	127.69	121.28
3	A	261	CW1	OAW-CAU-CBS	-3.98	103.97	109.67
3	A	261	CW1	OBA-CBL-OAY	-3.90	100.97	110.69
3	A	261	CW1	CAP-CBO-CAM	-3.88	106.09	109.51
3	A	261	CW1	CAC-CBG-CBD	-3.73	107.15	111.57
3	A	261	CW1	CBG-CBD-CBF	-3.69	120.72	129.68
3	A	261	CW1	CAO-CBD-CBG	3.65	127.22	119.72
3	B	261	CW1	CBT-OB	-3.63	96.17	107.62
3	A	261	CW1	CBT-OB	-3.56	96.41	107.62
3	A	261	CW1	OBA-CBQ-CBF	-3.45	104.47	113.32
3	A	261	CW1	OBC-CBR-CAG	3.38	116.04	108.71
3	A	261	CW1	OB	-3.33	100.87	105.19
3	B	261	CW1	OB	-3.19	100.12	105.07
3	B	261	CW1	CBO-CBF-CBD	-3.18	105.75	111.23
3	A	261	CW1	CAN-CAQ-CBR	-3.16	108.03	116.84
3	B	261	CW1	CBP-CBE-CAM	-3.15	124.83	130.11
3	B	261	CW1	CAN-CAQ-CBR	-3.06	108.33	116.84
3	A	261	CW1	CAG-CBR-CAQ	3.04	117.00	111.64
3	A	261	CW1	CAR-CAS-CBS	2.96	107.98	103.54
3	A	261	CW1	CAR-CBP-CBE	2.83	107.05	102.85
3	B	261	CW1	OAZ-CBN-CBT	-2.81	99.43	103.81
3	A	261	CW1	CAS-CBS-CAU	-2.75	109.92	113.95
3	B	261	CW1	CAS-CBS-CAU	-2.75	109.92	113.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	261	CW1	CBP-CBE-CAM	-2.73	125.52	130.11
3	A	261	CW1	OBB-CBT-CBN	-2.62	101.00	105.07
3	B	261	CW1	CAG-CBR-CAQ	2.60	116.22	111.64
3	A	261	CW1	CBL-OBA-CBQ	-2.58	110.88	115.24
3	A	261	CW1	CAA-OAV-CAT	-2.51	102.27	112.35
3	A	261	CW1	CAE-CBH-CBJ	2.27	115.29	110.42
3	B	261	CW1	CAD-CBG-CBD	-2.04	109.15	111.57
3	B	261	CW1	CAR-CBP-CBE	2.01	105.83	102.85

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	261	CW1	CBR
3	B	261	CW1	CBR

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	234/260 (90%)	-0.33	0 100 100	13, 21, 33, 39	0
1	B	232/260 (89%)	-0.17	3 (1%) 74 76	12, 21, 36, 46	0
2	C	5/5 (100%)	0.31	0 100 100	12, 15, 17, 19	0
2	D	5/5 (100%)	0.46	0 100 100	15, 16, 20, 24	0
All	All	476/530 (89%)	-0.24	3 (0%) 86 88	12, 21, 34, 46	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	220	TYR	2.5
1	B	240	ASP	2.2
1	B	145	ALA	2.0

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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	TPO	D	4	11/12	0.14	0.07	14,16,17,17	0
2	TPO	C	4	11/12	0.13	-0.12	8,11,14,16	0

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	CL	A	262	1/1	0.24	8.31	43,43,43,43	0
3	CW1	B	261	46/46	0.13	-0.05	15,20,34,42	0
3	CW1	A	261	46/46	0.11	-0.78	10,17,26,29	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.