



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:16 AM GMT

PDB ID : 3QZ7  
Title : T-3 ternary complex of Dpo4  
Authors : Pata, J.D.; Wu, Y.; Wilson, R.C.  
Deposited on : 2011-03-04  
Resolution : 2.00 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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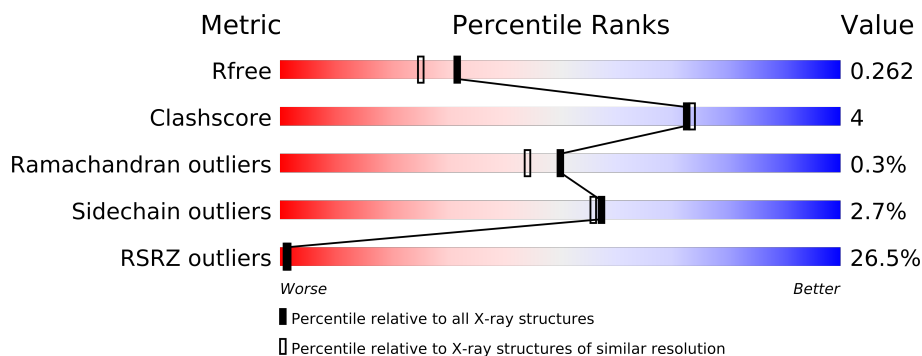
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.00 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	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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	360	
2	P	13	
3	T	19	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6635 atoms, of which 3219 are hydrogens 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 DNA polymerase IV.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	341	Total	C	H	N	O	S	0	0	0
			5634	1760	2891	472	504	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	353	GLY	-	EXPRESSION TAG	UNP Q97W02
A	354	GLY	-	EXPRESSION TAG	UNP Q97W02
A	355	HIS	-	EXPRESSION TAG	UNP Q97W02
A	356	HIS	-	EXPRESSION TAG	UNP Q97W02
A	357	HIS	-	EXPRESSION TAG	UNP Q97W02
A	358	HIS	-	EXPRESSION TAG	UNP Q97W02
A	359	HIS	-	EXPRESSION TAG	UNP Q97W02
A	360	HIS	-	EXPRESSION TAG	UNP Q97W02

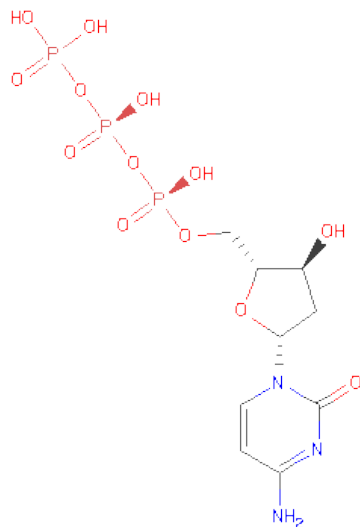
- Molecule 2 is a DNA chain called 5'-D(\*GP\*GP\*CP\*AP\*CP\*TP\*GP\*AP\*TP\*CP\*GP\*GP\*G)-3'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	P	13	Total	C	H	N	O	P	0	0	0
			415	127	147	53	76	12			

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	T	15	Total	C	H	N	O	P	0	0	0
			473	144	169	54	91	15			

- Molecule 4 is 2'-DEOXYCYTIDINE-5'-TRIPHOSPHATE (three-letter code: DCP) (formula: C<sub>9</sub>H<sub>16</sub>N<sub>3</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	
			40	9	12	3	13	3	
								0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total	Ca		
			3	3	0	0

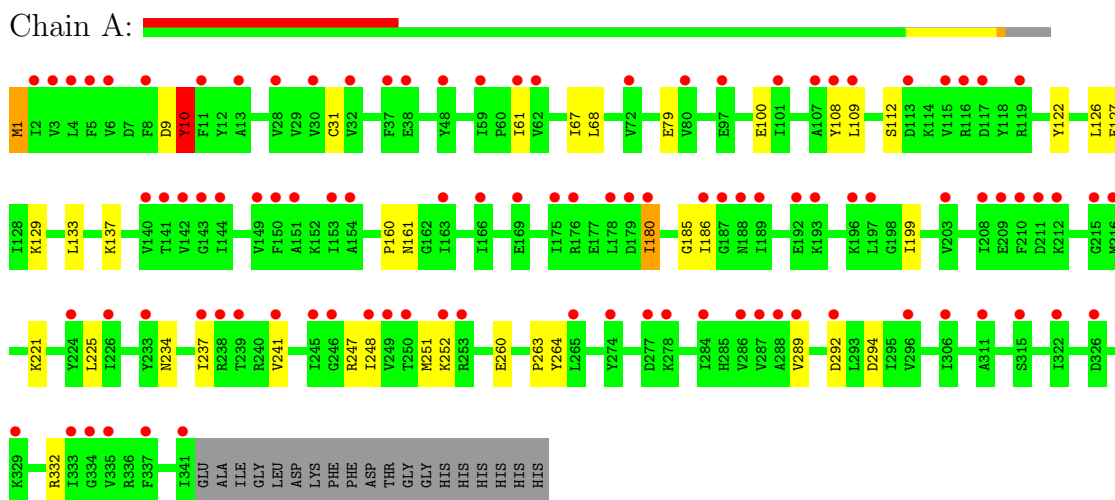
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	56	Total	O		
			56	56	0	0
6	P	6	Total	O		
			6	6	0	0
6	T	8	Total	O		
			8	8	0	0

### 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.

- Molecule 1: DNA polymerase IV



- Molecule 2: 5'-D(\*GP\*GP\*CP\*AP\*CP\*TP\*GP\*AP\*TP\*CP\*GP\*GP\*G)-3'



- Molecule 3: 5'-D(\*TP\*TP\*AP\*CP\*GP\*CP\*CP\*TP\*CP\*GP\*AP\*TP\*CP\*AP\*GP\*TP\*GP\*CP\*C)-3'



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.71Å 102.46Å 52.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.54 – 2.00 25.54 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.0 (25.54-2.00) 90.0 (25.54-2.00)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.71 (at 1.99Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.227 , 0.267 0.218 , 0.262	Depositor DCC
$R_{free}$ test set	1629 reflections (4.89%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.9	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.45 , 53.4	EDS
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 33331 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6635	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% 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: CA, DCP

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	0.37	0/2782	0.52	0/3736
2	P	0.69	0/301	1.43	3/464 (0.6%)
3	T	0.70	0/339	1.47	6/520 (1.2%)
All	All	0.45	0/3422	0.81	9/4720 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	16	DT	O4'-C4'-C3'	-8.59	100.84	106.00
3	T	16	DT	O4'-C1'-N1	7.09	112.96	108.00
3	T	9	DC	O4'-C1'-N1	7.08	112.96	108.00
2	P	6	DT	O4'-C1'-N1	-6.38	103.53	108.00
3	T	5	DG	O4'-C1'-N9	-6.04	103.78	108.00
2	P	12	DG	O4'-C1'-N9	5.61	111.93	108.00
2	P	12	DG	C1'-O4'-C4'	-5.23	104.87	110.10
3	T	10	DG	C5-C6-O6	-5.18	125.49	128.60
3	T	10	DG	N1-C6-O6	5.00	122.90	119.90

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	2743	2891	0	24	0
2	P	268	147	0	2	0
3	T	304	169	0	2	0
4	A	28	12	0	1	0
5	A	3	0	0	0	0
6	A	56	0	0	0	0
6	P	6	0	0	0	0
6	T	8	0	0	0	0
All	All	3416	3219	0	28	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 4.

All (28) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1:MET:HB2	1:A:112:SER:OG	1.82	0.79
1:A:247:ARG:C	1:A:248:ILE:HD12	2.22	0.59
3:T:16:DT:H2''	3:T:17:DG:O5'	2.04	0.58
1:A:122:TYR:CE2	1:A:126:LEU:HD11	2.39	0.57
1:A:186:ILE:HD11	1:A:225:LEU:HD21	1.85	0.57
1:A:133:LEU:O	1:A:137:LYS:HD3	2.04	0.57
1:A:108:TYR:C	1:A:109:LEU:HD12	2.26	0.56
2:P:5:DC:H2'	2:P:6:DT:H73	1.88	0.56
1:A:289:VAL:HB	1:A:332:ARG:HB2	1.87	0.55
3:T:16:DT:H2'	3:T:17:DG:C8	2.42	0.55
1:A:251:MET:HG2	1:A:264:TYR:CD2	2.46	0.50
1:A:185:GLY:O	1:A:221:LYS:CE	2.59	0.50
1:A:100:GLU:HB2	1:A:237:ILE:HG23	1.95	0.47
1:A:10:TYR:HA	4:A:361:DCP:PB	2.55	0.46
1:A:9:ASP:O	1:A:10:TYR:C	2.54	0.46
1:A:67:ILE:HG22	1:A:68:LEU:HG	1.97	0.46
1:A:260:GLU:O	1:A:263:PRO:HD2	2.15	0.46
1:A:122:TYR:HE2	1:A:126:LEU:HD11	1.79	0.45
1:A:248:ILE:HD12	1:A:248:ILE:N	2.32	0.45
1:A:241:VAL:HG23	1:A:241:VAL:O	2.17	0.45
2:P:6:DT:H2''	2:P:7:DG:C8	2.52	0.45
1:A:180:ILE:HD13	1:A:199:ILE:HG22	1.99	0.44
1:A:79:GLU:CD	1:A:79:GLU:H	2.21	0.44
1:A:31:CYS:HB3	1:A:61:ILE:HD11	1.99	0.43
1:A:109:LEU:HD12	1:A:109:LEU:N	2.34	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:129:LYS:HE3	1:A:129:LYS:HB3	1.89	0.42
1:A:160:PRO:O	1:A:161:ASN:C	2.58	0.42
1:A:31:CYS:HB3	1:A:61:ILE:CD1	2.51	0.41

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	339/360 (94%)	325 (96%)	13 (4%)	1 (0%)	50 44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	TYR

### 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	300/315 (95%)	292 (97%)	8 (3%)	57 56

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	10	TYR
1	A	127	GLU
1	A	180	ILE

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Mol	Chain	Res	Type
1	A	234	ASN
1	A	252	LYS
1	A	292	ASP
1	A	294	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	DCP	A	361	5	29,29,29	2.14	8 (27%)	42,45,45	2.62	14 (33%)

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	DCP	A	361	5	-	0/19/34/34	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	361	DCP	C2-N1	5.96	1.44	1.38
4	A	361	DCP	C6-N1	4.37	1.42	1.35
4	A	361	DCP	C6-C5	3.74	1.46	1.38
4	A	361	DCP	C4-N3	3.45	1.42	1.35
4	A	361	DCP	C2-N3	3.26	1.44	1.35
4	A	361	DCP	C4-N4	3.01	1.44	1.35
4	A	361	DCP	C5-C4	2.68	1.47	1.40
4	A	361	DCP	C3'-C4'	-2.58	1.45	1.53

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	361	DCP	O3B-PB-O3A	-7.49	86.43	101.66
4	A	361	DCP	O5'-PA-O1A	-7.38	80.47	109.37
4	A	361	DCP	C6-C5-C4	6.71	120.26	117.47
4	A	361	DCP	O2B-PB-O3A	-4.45	84.01	105.14
4	A	361	DCP	O2A-PA-O5'	-3.94	88.63	108.51
4	A	361	DCP	PB-O3B-PG	-3.81	120.50	131.68
4	A	361	DCP	O5'-C5'-C4'	3.22	120.76	108.94
4	A	361	DCP	O2B-PB-O3B	3.07	119.71	105.14
4	A	361	DCP	O2A-PA-O3A	2.85	118.65	105.14
4	A	361	DCP	C2-N3-C4	2.81	119.64	115.57
4	A	361	DCP	O3A-PB-O1B	-2.52	93.16	111.28
4	A	361	DCP	O3A-PA-O5'	-2.45	92.45	103.41
4	A	361	DCP	O4'-C1'-N1	2.29	111.97	107.68
4	A	361	DCP	O3G-PG-O3B	2.28	115.93	105.14

There are no chirality outliers.

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	341/360 (94%)	1.46	99 (29%) 1 1	36, 60, 83, 91	0
2	P	13/13 (100%)	0.53	0 100 100	54, 61, 72, 73	0
3	T	15/19 (78%)	0.16	0 100 100	51, 65, 69, 77	0
All	All	369/392 (94%)	1.37	99 (26%) 1 1	36, 61, 83, 91	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	37	PHE	6.3
1	A	116	ARG	6.3
1	A	241	VAL	5.5
1	A	62	VAL	5.1
1	A	334	GLY	4.8
1	A	189	ILE	4.7
1	A	341	ILE	4.6
1	A	209	GLU	4.5
1	A	274	TYR	4.5
1	A	48	TYR	4.4
1	A	239	THR	4.4
1	A	335	VAL	4.4
1	A	337	PHE	4.2
1	A	5	PHE	4.1
1	A	115	VAL	4.0
1	A	107	ALA	4.0
1	A	61	ILE	4.0
1	A	248	ILE	3.9
1	A	113	ASP	3.9
1	A	144	ILE	3.9
1	A	333	ILE	3.8
1	A	108	TYR	3.7
1	A	4	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	287	VAL	3.7
1	A	178	LEU	3.7
1	A	187	GLY	3.7
1	A	212	LYS	3.6
1	A	203	VAL	3.5
1	A	169	GLU	3.4
1	A	208	ILE	3.3
1	A	253	ARG	3.3
1	A	175	ILE	3.3
1	A	288	ALA	3.3
1	A	180	ILE	3.3
1	A	284	ILE	3.2
1	A	143	GLY	3.2
1	A	216	MET	3.1
1	A	186	ILE	3.1
1	A	286	VAL	3.0
1	A	197	LEU	3.0
1	A	142	VAL	3.0
1	A	153	ILE	2.9
1	A	249	VAL	2.9
1	A	289	VAL	2.9
1	A	296	VAL	2.9
1	A	278	LYS	2.9
1	A	238	ARG	2.8
1	A	210	PHE	2.8
1	A	6	VAL	2.8
1	A	97	GLU	2.8
1	A	252	LYS	2.8
1	A	179	ASP	2.8
1	A	306	ILE	2.8
1	A	151	ALA	2.8
1	A	188	ASN	2.7
1	A	192	GLU	2.7
1	A	246	GLY	2.7
1	A	193	LYS	2.7
1	A	245	ILE	2.7
1	A	196	LYS	2.6
1	A	176	ARG	2.6
1	A	215	GLY	2.6
1	A	117	ASP	2.6
1	A	237	ILE	2.6
1	A	2	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	149	VAL	2.6
1	A	150	PHE	2.5
1	A	329	LYS	2.5
1	A	3	VAL	2.5
1	A	322	ILE	2.5
1	A	38	GLU	2.5
1	A	119	ARG	2.5
1	A	211	ASP	2.4
1	A	166	ILE	2.4
1	A	226	ILE	2.4
1	A	109	LEU	2.4
1	A	154	ALA	2.4
1	A	163	ILE	2.4
1	A	265	LEU	2.3
1	A	233	TYR	2.3
1	A	224	TYR	2.3
1	A	250	THR	2.3
1	A	72	VAL	2.3
1	A	315	SER	2.3
1	A	11	PHE	2.2
1	A	80	VAL	2.2
1	A	326	ASP	2.2
1	A	28	VAL	2.2
1	A	101	ILE	2.2
1	A	13	ALA	2.2
1	A	311	ALA	2.2
1	A	8	PHE	2.2
1	A	32	VAL	2.2
1	A	277	ASP	2.2
1	A	140	VAL	2.1
1	A	30	VAL	2.1
1	A	292	ASP	2.1
1	A	141	THR	2.0
1	A	59	ILE	2.0

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

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

### 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	DCP	A	361	28/28	0.14	-0.55	34,43,52,56	0
5	CA	A	363	1/1	0.15	-0.76	71,71,71,71	0
5	CA	A	362	1/1	0.13	-1.19	38,38,38,38	0
5	CA	A	364	1/1	0.07	-2.30	94,94,94,94	0

### 6.5 Other polymers

There are no such residues in this entry.