



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 10:48 pm BST

PDB ID : 3TG0  
Title : E. coli alkaline phosphatase with bound inorganic phosphate  
Authors : Bobyr, E.; Lassila, J.K.; Wiersma-Koch, H.I.; Fenn, T.D.; Lee, J.J.; Nikolic-Hughes, I.; Hodgson, K.O.; Rees, D.C.; Hedman, B.; Herschlag, D.  
Deposited on : 2011-08-16  
Resolution : 1.20 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](mailto: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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

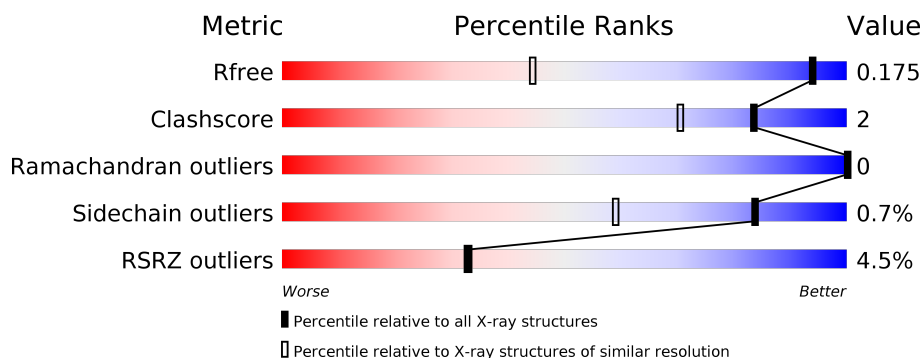
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.20 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))
$R_{free}$	130704	1223 (1.22-1.18)
Clashscore	141614	1286 (1.22-1.18)
Ramachandran outliers	138981	1240 (1.22-1.18)
Sidechain outliers	138945	1239 (1.22-1.18)
RSRZ outliers	127900	1200 (1.22-1.18)

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 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	449	<div> <div>3%</div> <div>94%</div> <div>• •</div> </div>
1	B	449	<div> <div>2%</div> <div>94%</div> <div>• •</div> </div>
1	C	449	<div> <div>8%</div> <div>95%</div> <div>• •</div> </div>
1	D	449	<div> <div>4%</div> <div>96%</div> <div>• •</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15183 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 Alkaline phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	13	0
			3315	2049	585	669	12			
1	B	440	Total	C	N	O	S	0	15	0
			3334	2059	594	669	12			
1	C	440	Total	C	N	O	S	0	12	0
			3300	2042	579	667	12			
1	D	443	Total	C	N	O	S	0	10	0
			3324	2058	584	670	12			

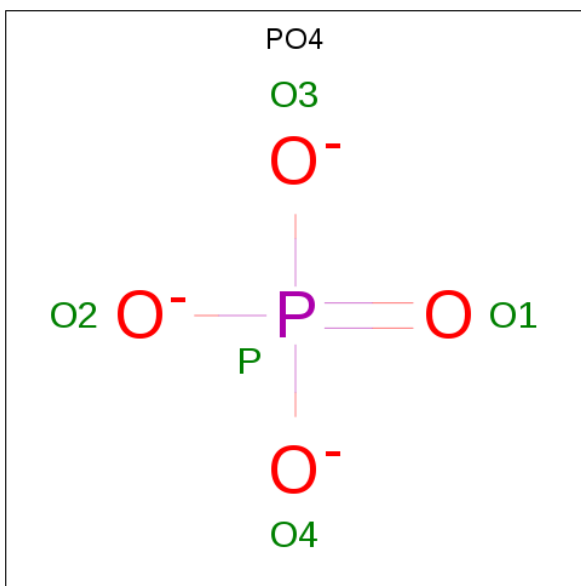
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula:  $\text{O}_4\text{P}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	D	1	Total	O	P	0	0
			5	4	1		

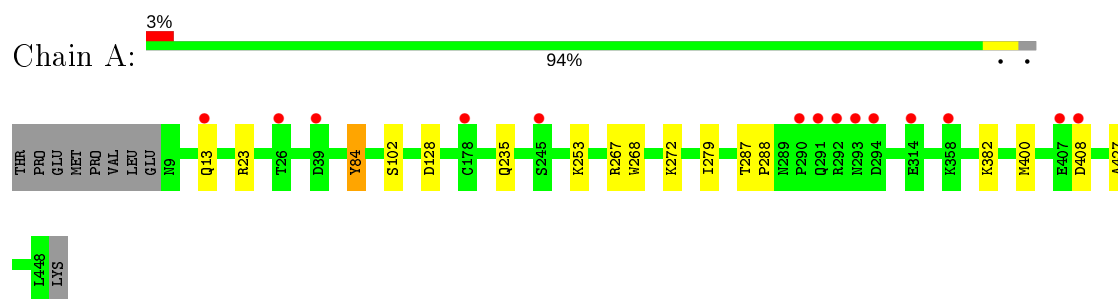
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	485	Total	O	0	0
			485	485		
5	B	489	Total	O	0	0
			489	489		
5	C	415	Total	O	0	0
			415	415		
5	D	489	Total	O	0	0
			489	489		

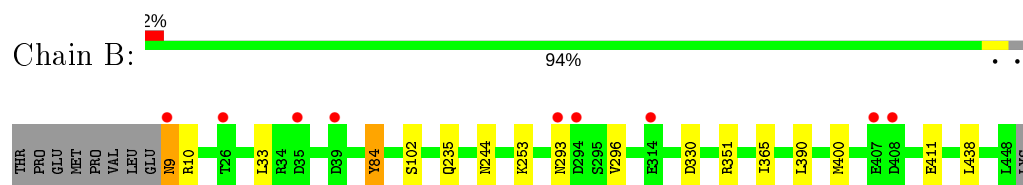
### 3 Residue-property plots [i](#)

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

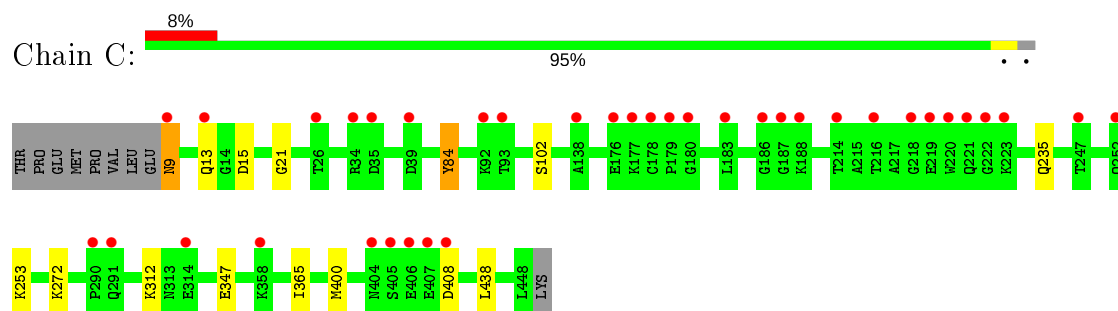
- Molecule 1: Alkaline phosphatase



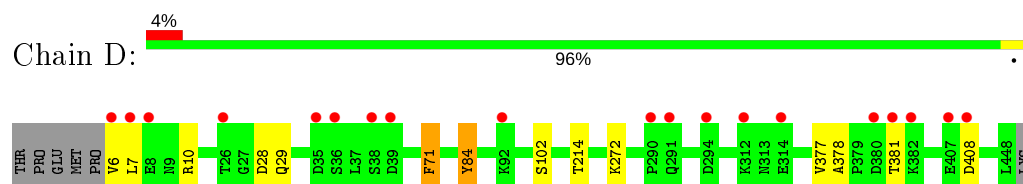
- Molecule 1: Alkaline phosphatase



- Molecule 1: Alkaline phosphatase



- Molecule 1: Alkaline phosphatase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.96 Å 98.46 Å 152.15 Å 90.00° 94.61° 90.00°	Depositor
Resolution (Å)	50.00 – 1.20 24.62 – 1.20	Depositor EDS
% Data completeness (in resolution range)	87.4 (50.00-1.20) 87.5 (24.62-1.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.36 (at 1.20 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.148 , 0.169 0.157 , 0.175	Depositor DCC
$R_{free}$ test set	26546 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	13.9	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 55.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	15183	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, MG, ZN

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	0.60	0/3389	0.70	2/4599 (0.0%)
1	B	0.61	0/3411	0.71	3/4626 (0.1%)
1	C	0.57	0/3380	0.69	1/4587 (0.0%)
1	D	0.58	0/3390	0.69	1/4602 (0.0%)
All	All	0.59	0/13570	0.70	7/18414 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	84	TYR	CB-CG-CD1	6.37	124.82	121.00
1	B	84	TYR	CB-CG-CD1	6.04	124.62	121.00
1	D	84	TYR	CB-CG-CD1	5.93	124.56	121.00
1	A	84	TYR	CB-CG-CD1	5.82	124.49	121.00
1	B	330	ASP	CB-CG-OD2	-5.47	113.37	118.30
1	A	128	ASP	CB-CG-OD1	5.34	123.10	118.30
1	B	330	ASP	CB-CG-OD1	5.22	123.00	118.30

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	3315	0	3270	16	0
1	B	3334	0	3295	15	0
1	C	3300	0	3259	12	0
1	D	3324	0	3273	22	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	5	0	0	1	0
4	B	5	0	0	1	0
4	C	5	0	0	1	0
4	D	5	0	0	1	0
5	A	485	0	0	4	0
5	B	489	0	0	5	0
5	C	415	0	0	5	0
5	D	489	0	0	1	0
All	All	15183	0	13097	64	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 2.

All (64) 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:D:378:ALA:O	1:D:381[B]:THR:HG23	1.51	1.07
1:D:6:VAL:HG21	1:D:29[A]:GLN:NE2	1.71	1.06
1:D:6:VAL:HG21	1:D:29[A]:GLN:HE22	1.19	1.03
1:D:6:VAL:CG2	1:D:29[A]:GLN:HE22	1.77	0.97
1:A:102:SER:OG	4:A:504:PO4:P	2.24	0.89
1:D:6:VAL:CG2	1:D:29[A]:GLN:NE2	2.34	0.84
1:D:102:SER:OG	4:D:516:PO4:P	2.36	0.81
1:B:235[A]:GLN:CG	1:B:253:LYS:HG2	2.17	0.75
1:C:102:SER:OG	4:C:512:PO4:P	2.36	0.74
1:B:9:ASN:HD22	1:B:9:ASN:C	1.92	0.73
1:A:287[B]:THR:HG23	5:A:5994:HOH:O	1.87	0.73
1:A:279[A]:ILE:HD11	1:A:382:LYS:HG2	1.70	0.72
1:B:235[B]:GLN:NE2	5:B:6247:HOH:O	2.22	0.69
1:B:102:SER:OG	4:B:508:PO4:P	2.50	0.69

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:29[A]:GLN:HA	1:D:29[A]:GLN:NE2	2.09	0.68
1:D:377:VAL:HB	1:D:381[B]:THR:HG21	1.77	0.67
1:A:235:GLN:HG3	1:A:253:LYS:HG2	1.77	0.66
1:B:244[A]:ASN:OD1	5:B:6956:HOH:O	2.15	0.64
1:C:235[A]:GLN:CG	1:C:253:LYS:HG2	2.30	0.62
1:A:279[A]:ILE:CD1	1:A:382:LYS:HG2	2.31	0.59
1:A:400[B]:MET:CE	5:A:5571:HOH:O	2.50	0.59
1:C:347[B]:GLU:HG3	5:C:6049:HOH:O	2.01	0.58
1:A:267[A]:ARG:HG2	1:A:268:TRP:CD1	2.39	0.58
1:C:235[A]:GLN:HG3	1:C:253:LYS:HG2	1.86	0.58
1:A:287[B]:THR:HG22	1:A:288:PRO:HD2	1.86	0.57
1:A:400[B]:MET:HE1	5:A:5571:HOH:O	2.03	0.57
1:D:6:VAL:HG22	1:D:29[A]:GLN:OE1	2.05	0.57
1:B:293:ASN:OD1	1:B:296:VAL:HG23	2.05	0.57
1:B:400[B]:MET:CE	5:B:6042:HOH:O	2.53	0.56
1:D:29[A]:GLN:CA	1:D:29[A]:GLN:NE2	2.68	0.56
1:B:411:GLU:OE1	5:B:6028:HOH:O	2.17	0.56
1:C:347[B]:GLU:CG	5:C:6049:HOH:O	2.54	0.55
1:D:28:ASP:C	1:D:29[A]:GLN:HE21	2.10	0.55
1:D:6:VAL:CG1	1:D:29[A]:GLN:HE22	2.20	0.54
1:B:400[B]:MET:HE1	5:B:6042:HOH:O	2.06	0.54
1:D:272:LYS:NZ	1:D:408:ASP:OD1	2.37	0.54
1:D:7:LEU:HD23	1:D:29[A]:GLN:OE1	2.08	0.53
1:B:9:ASN:ND2	1:B:9:ASN:C	2.60	0.53
1:A:235:GLN:CG	1:A:253:LYS:HG2	2.38	0.53
1:D:377:VAL:CG2	1:D:381[B]:THR:HG21	2.38	0.53
1:D:6:VAL:HG22	1:D:29[A]:GLN:CD	2.31	0.51
1:C:9:ASN:ND2	1:C:9:ASN:C	2.64	0.51
1:C:400[B]:MET:CE	5:C:5881:HOH:O	2.60	0.49
1:C:400[B]:MET:HE1	5:C:5881:HOH:O	2.12	0.49
1:D:377:VAL:CB	1:D:381[B]:THR:HG21	2.43	0.49
1:B:235[A]:GLN:HG3	1:B:253:LYS:HG2	1.92	0.48
1:C:272:LYS:NZ	1:C:408:ASP:OD2	2.36	0.48
1:D:6:VAL:HG11	1:D:29[A]:GLN:HE22	1.79	0.47
1:D:6:VAL:HG22	1:D:29[A]:GLN:NE2	2.26	0.47
1:C:312:LYS:HG3	5:C:6189:HOH:O	2.15	0.45
1:A:427:ALA:HB1	1:B:33[A]:LEU:HD12	1.98	0.44
1:D:214[A]:THR:HG23	5:D:6785:HOH:O	2.17	0.44
1:A:279[A]:ILE:HD13	1:A:382:LYS:HE3	1.99	0.44
1:A:13:GLN:OE1	1:A:23[B]:ARG:NH2	2.51	0.43
1:A:272:LYS:NZ	1:A:408[B]:ASP:OD1	2.32	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287[B]:THR:HG22	1:A:288:PRO:CD	2.48	0.43
1:A:400[B]:MET:HE3	5:A:5571:HOH:O	2.16	0.43
1:C:365:ILE:HD13	1:C:438:LEU:HD11	2.00	0.43
1:D:10:ARG:HH22	1:D:29[B]:GLN:NE2	2.17	0.43
1:B:365:ILE:HD13	1:B:438:LEU:HD11	2.02	0.42
1:B:9:ASN:HD22	1:B:10:ARG:N	2.18	0.42
1:B:351[B]:ARG:HA	1:B:351[B]:ARG:HD2	1.91	0.41
1:C:15:ASP:O	1:C:21:GLY:HA3	2.21	0.40
1:D:10:ARG:HB2	1:D:71[A]:PHE:CD1	2.56	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	451/449 (100%)	443 (98%)	8 (2%)	0	100	100
1	B	453/449 (101%)	446 (98%)	7 (2%)	0	100	100
1	C	450/449 (100%)	442 (98%)	8 (2%)	0	100	100
1	D	451/449 (100%)	445 (99%)	6 (1%)	0	100	100
All	All	1805/1796 (100%)	1776 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	344/340 (101%)	343 (100%)	1 (0%)	92	79
1	B	346/340 (102%)	343 (99%)	3 (1%)	78	50
1	C	343/340 (101%)	340 (99%)	3 (1%)	78	50
1	D	344/340 (101%)	341 (99%)	3 (1%)	78	50
All	All	1377/1360 (101%)	1367 (99%)	10 (1%)	84	59

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

Mol	Chain	Res	Type
1	A	84	TYR
1	B	9	ASN
1	B	84	TYR
1	B	390	LEU
1	C	9	ASN
1	C	13	GLN
1	C	84	TYR
1	D	71[A]	PHE
1	D	71[B]	PHE
1	D	84	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	235	GLN
1	B	9	ASN
1	B	13	GLN
1	C	13	GLN
1	C	83	GLN
1	C	410	GLN
1	D	291	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 16 ligands modelled in this entry, 12 are 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	$\# Z  > 2$
4	PO4	A	504	2	4,4,4	1.15	0	6,6,6	0.86	0
4	PO4	B	508	2	4,4,4	0.67	0	6,6,6	1.12	1 (16%)
4	PO4	D	516	2	4,4,4	1.05	0	6,6,6	1.23	1 (16%)
4	PO4	C	512	2	4,4,4	0.95	0	6,6,6	0.65	0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
4	D	516	PO4	O3-P-O1	-2.43	102.00	110.89
4	B	508	PO4	O2-P-O1	-2.25	102.66	110.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	504	PO4	1	0
4	B	508	PO4	1	0
4	D	516	PO4	1	0
4	C	512	PO4	1	0

## 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	440/449 (97%)	-0.02	14 (3%) 47 46	8, 13, 22, 30	0
1	B	440/449 (97%)	-0.01	9 (2%) 65 66	8, 13, 22, 30	0
1	C	440/449 (97%)	0.15	37 (8%) 11 11	10, 15, 24, 31	0
1	D	443/449 (98%)	0.02	19 (4%) 35 35	10, 14, 24, 30	0
All	All	1763/1796 (98%)	0.04	79 (4%) 33 33	8, 14, 23, 31	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	408	ASP	6.0
1	D	408	ASP	5.8
1	D	39	ASP	5.6
1	A	408[A]	ASP	4.7
1	B	293	ASN	4.6
1	D	6	VAL	4.4
1	D	290	PRO	4.4
1	C	220	TRP	4.2
1	C	26	THR	4.1
1	D	38	SER	4.1
1	C	407	GLU	4.1
1	C	39	ASP	3.9
1	D	381[A]	THR	3.8
1	A	314	GLU	3.8
1	C	92	LYS	3.7
1	A	26	THR	3.7
1	D	26	THR	3.7
1	D	407	GLU	3.7
1	D	314	GLU	3.7
1	C	222	GLY	3.6
1	C	35	ASP	3.6

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	314	GLU	3.6
1	D	36	SER	3.5
1	A	292	ARG	3.5
1	A	294[A]	ASP	3.5
1	C	358	LYS	3.4
1	D	294	ASP	3.4
1	B	26	THR	3.3
1	B	408	ASP	3.2
1	D	35	ASP	3.2
1	C	218	GLY	3.1
1	C	291	GLN	3.1
1	A	39	ASP	3.1
1	C	183	LEU	3.1
1	C	179	PRO	3.0
1	C	405	SER	3.0
1	C	176	GLU	2.9
1	C	178	CYS	2.9
1	C	219	GLU	2.9
1	C	216	THR	2.9
1	B	39	ASP	2.9
1	C	404	ASN	2.9
1	A	290	PRO	2.9
1	C	180	GLY	2.8
1	C	187	GLY	2.8
1	D	291	GLN	2.8
1	A	407	GLU	2.8
1	C	188	LYS	2.8
1	C	9	ASN	2.7
1	D	8	GLU	2.7
1	B	314	GLU	2.6
1	B	294	ASP	2.6
1	C	138	ALA	2.6
1	C	252[A]	GLN	2.6
1	C	223	LYS	2.5
1	B	35	ASP	2.5
1	C	290	PRO	2.5
1	D	380	ASP	2.5
1	D	92	LYS	2.5
1	C	214[A]	THR	2.4
1	D	382	LYS	2.4
1	A	358	LYS	2.4
1	A	178	CYS	2.3

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	407	GLU	2.3
1	C	177	LYS	2.2
1	A	293	ASN	2.2
1	C	93	THR	2.2
1	D	7	LEU	2.2
1	C	34	ARG	2.2
1	D	312	LYS	2.2
1	A	291	GLN	2.1
1	C	247	THR	2.1
1	C	406	GLU	2.1
1	C	221	GLN	2.1
1	C	186	GLY	2.1
1	A	13	GLN	2.1
1	B	9	ASN	2.0
1	A	245[A]	SER	2.0
1	C	13	GLN	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates 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, 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	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	PO4	A	504	5/5	0.96	0.09	12,13,13,18	5
4	PO4	B	508	5/5	0.97	0.08	11,11,12,16	5
4	PO4	D	516	5/5	0.98	0.08	11,11,14,17	5
4	PO4	C	512	5/5	0.98	0.07	15,15,16,19	5
2	ZN	D	514	1/1	1.00	0.03	12,12,12,12	0
2	ZN	A	502	1/1	1.00	0.03	10,10,10,10	0
3	MG	B	507	1/1	1.00	0.07	8,8,8,8	0

*Continued on next page...*



*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	C	509	1/1	1.00	0.03	14,14,14,14	0
2	ZN	B	506	1/1	1.00	0.04	10,10,10,10	0
3	MG	D	515	1/1	1.00	0.06	9,9,9,9	0
3	MG	C	511	1/1	1.00	0.07	12,12,12,12	0
3	MG	A	503	1/1	1.00	0.07	8,8,8,8	0
2	ZN	B	505	1/1	1.00	0.03	10,10,10,10	0
2	ZN	C	510	1/1	1.00	0.03	15,15,15,15	0
2	ZN	A	501	1/1	1.00	0.04	11,11,11,11	0
2	ZN	D	513	1/1	1.00	0.03	12,12,12,12	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.