



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:52 PM GMT

PDB ID : 1HO4  
Title : CRYSTAL STRUCTURE OF PYRIDOXINE 5'-PHOSPHATE SYNTHASE  
IN COMPLEX WITH PYRIDOXINE 5'-PHOSPHATE AND INORGANIC  
PHOSPHATE  
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Deposited on : 2000-12-08  
Resolution : 2.30 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

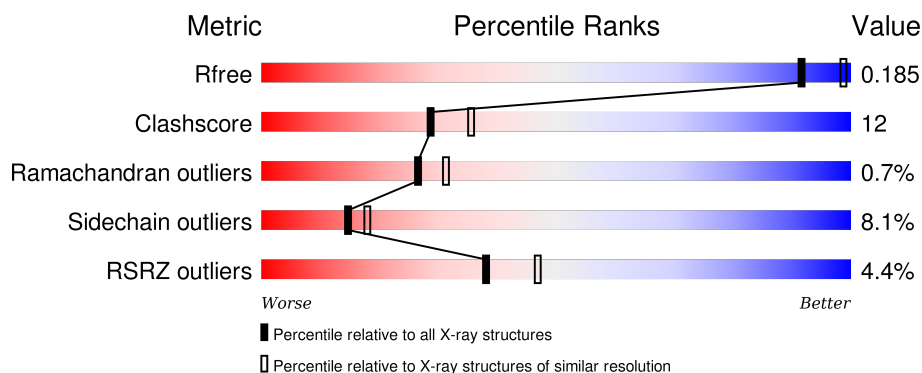
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	242	<div> <div>4%</div> <div>77%</div> <div>20%</div> <div>.</div> </div>
1	B	242	<div> <div>6%</div> <div>79%</div> <div>18%</div> <div>.</div> </div>
1	C	242	<div> <div>5%</div> <div>81%</div> <div>17%</div> <div>.</div> </div>
1	D	242	<div> <div>3%</div> <div>72%</div> <div>24%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8148 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 PYRIDOXINE 5'-PHOSPHATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1833	1136	337	348	12			
1	B	242	Total	C	N	O	S	0	0	0
			1837	1138	337	350	12			
1	C	242	Total	C	N	O	S	0	0	0
			1837	1138	337	350	12			
1	D	242	Total	C	N	O	S	0	0	0
			1827	1132	333	350	12			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



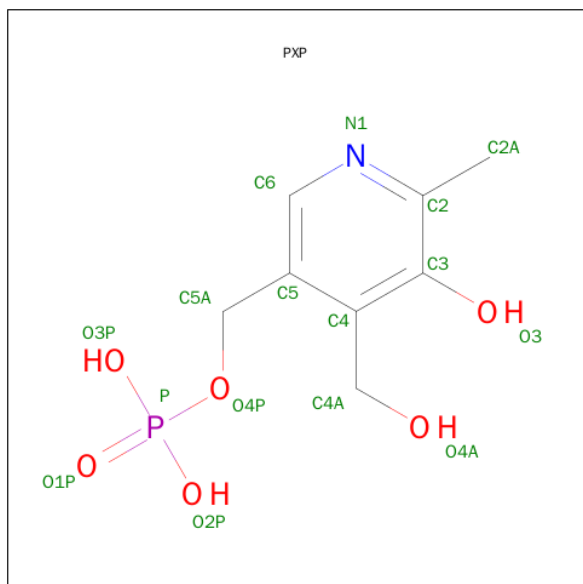
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is PYRIDOXINE-5'-PHOSPHATE (three-letter code: PXP) (formula: C<sub>8</sub>H<sub>12</sub>NO<sub>6</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	A	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	C	1	Total	C	N	O	P	0	0
			16	8	1	6	1		
3	B	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

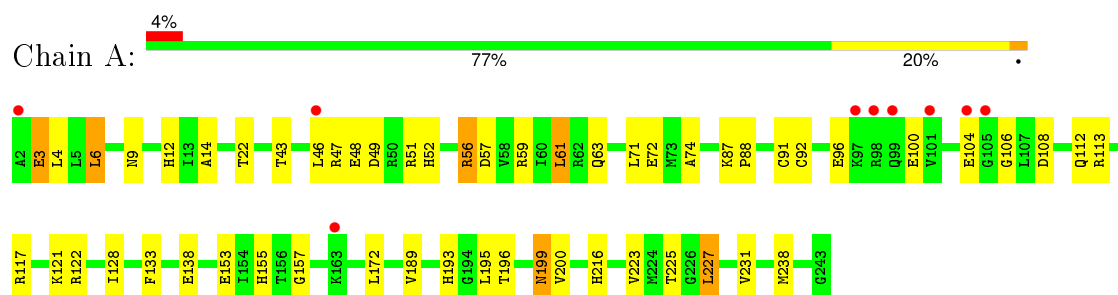
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	194	Total	O	0	0
			194	194		
4	B	168	Total	O	0	0
			168	168		
4	C	182	Total	O	0	0
			182	182		
4	D	191	Total	O	0	0
			191	191		

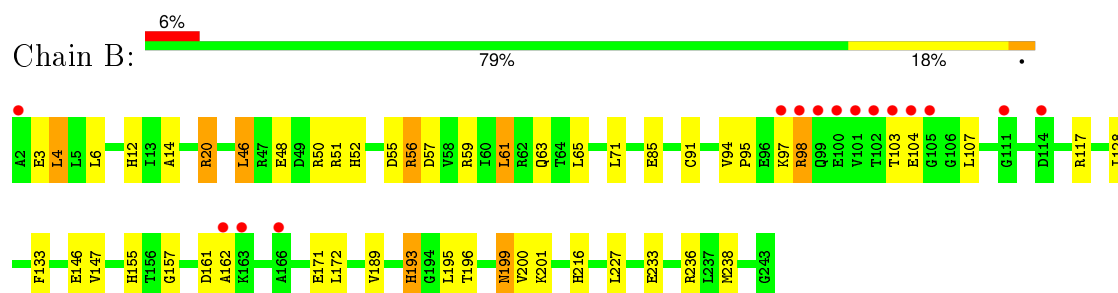
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

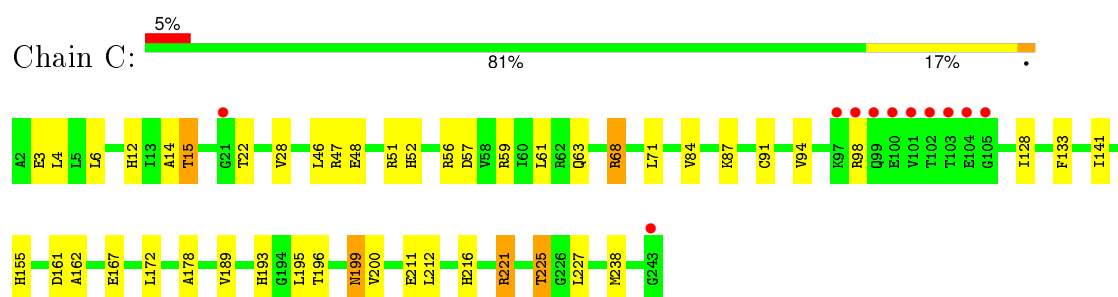
#### • Molecule 1: PYRIDOXINE 5'-PHOSPHATE SYNTHASE



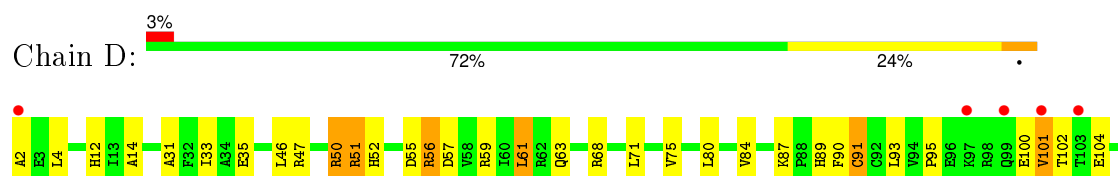
#### • Molecule 1: PYRIDOXINE 5'-PHOSPHATE SYNTHASE

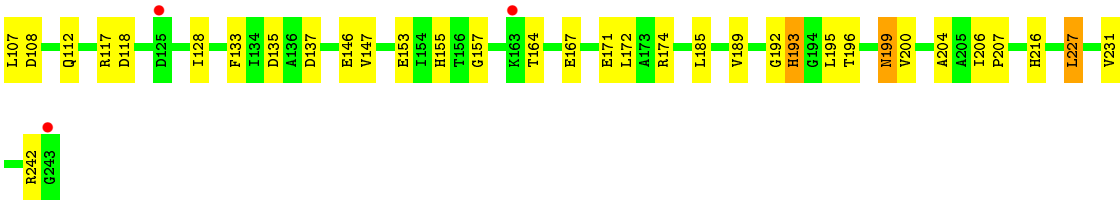


#### • Molecule 1: PYRIDOXINE 5'-PHOSPHATE SYNTHASE



#### • Molecule 1: PYRIDOXINE 5'-PHOSPHATE SYNTHASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	128.80 Å   156.30 Å   127.50 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.88 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.30) 95.9 (19.88-2.30)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.56 (at 2.30 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.187   ,   0.245 0.191   ,   0.185	Depositor DCC
$R_{free}$ test set	2782 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.1	Xtriage
Anisotropy	0.815	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 56.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 54834 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8148	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.92% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: PO4, PXP

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	0.49	0/1855	0.74	0/2504
1	B	0.52	0/1859	0.72	0/2509
1	C	0.51	0/1859	0.74	0/2509
1	D	0.52	0/1849	0.75	0/2498
All	All	0.51	0/7422	0.74	0/10020

There are no bond length outliers.

There are no bond angle outliers.

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	1833	0	1850	38	0
1	B	1837	0	1854	46	0
1	C	1837	0	1854	37	0
1	D	1827	0	1832	56	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	A	16	0	10	2	0
3	B	16	0	10	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	16	0	10	2	0
3	D	16	0	10	2	0
4	A	194	0	0	6	0
4	B	168	0	0	4	0
4	C	182	0	0	3	0
4	D	191	0	0	11	0
All	All	8148	0	7430	171	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 12.

All (171) 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:C:12:HIS:HA	1:C:15:THR:HG23	1.20	1.15
1:B:63:GLN:HE22	1:D:59:ARG:HH12	1.02	0.95
1:D:102:THR:HG23	4:D:870:HOH:O	1.69	0.90
1:A:63:GLN:HE22	1:C:59:ARG:HH12	0.93	0.89
1:B:59:ARG:HH12	1:D:63:GLN:HE22	1.18	0.88
1:D:155:HIS:HD2	1:D:157:GLY:H	1.17	0.88
1:C:12:HIS:HA	1:C:15:THR:CG2	2.02	0.86
1:A:199:ASN:C	1:A:199:ASN:HD22	1.79	0.85
1:D:155:HIS:CD2	1:D:157:GLY:H	1.98	0.81
1:A:59:ARG:HH12	1:C:63:GLN:HE22	1.28	0.81
1:C:221:ARG:HB2	1:C:221:ARG:HH11	1.48	0.79
1:B:199:ASN:HD22	1:B:199:ASN:C	1.86	0.78
1:A:63:GLN:HE22	1:C:59:ARG:NH1	1.77	0.77
1:A:63:GLN:NE2	1:C:59:ARG:HH12	1.79	0.76
1:B:56:ARG:HD2	4:B:906:HOH:O	1.85	0.75
1:A:196:THR:H	1:A:199:ASN:ND2	1.85	0.75
1:C:199:ASN:C	1:C:199:ASN:HD22	1.90	0.75
1:A:12:HIS:CD2	1:A:216:HIS:HD2	2.04	0.74
1:B:97:LYS:O	1:B:98:ARG:HB2	1.87	0.74
1:D:196:THR:H	1:D:199:ASN:ND2	1.88	0.71
1:C:162:ALA:HA	1:C:167:GLU:OE2	1.91	0.71
1:B:57:ASP:O	1:B:61:LEU:HB2	1.91	0.70
1:B:155:HIS:CE1	1:B:157:GLY:HA3	2.26	0.69
1:A:225:THR:HG22	4:A:921:HOH:O	1.93	0.68
1:D:50:ARG:N	1:D:50:ARG:HD2	2.09	0.68
1:D:199:ASN:HD22	1:D:199:ASN:C	1.96	0.68
1:A:14:ALA:HB3	1:A:52:HIS:HB2	1.73	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:PRO:HB2	1:D:101:VAL:HG22	1.76	0.67
1:B:117:ARG:HA	1:B:147:VAL:HG13	1.78	0.66
1:A:12:HIS:HD2	1:A:216:HIS:HD2	1.43	0.65
1:D:100:GLU:HG2	4:D:893:HOH:O	1.97	0.63
1:C:195:LEU:HA	1:C:199:ASN:HD21	1.63	0.63
1:D:89:HIS:HD2	1:D:90:PHE:CE1	2.17	0.63
1:A:196:THR:H	1:A:199:ASN:HD21	1.45	0.62
1:A:12:HIS:CD2	1:A:216:HIS:CD2	2.87	0.62
1:A:74:ALA:HB2	1:A:96:GLU:HB2	1.82	0.62
1:D:107:LEU:HD12	4:D:832:HOH:O	1.98	0.61
1:D:102:THR:HG23	4:D:842:HOH:O	1.99	0.61
1:A:12:HIS:HD2	1:A:216:HIS:CD2	2.19	0.61
1:B:95:PRO:HG3	1:B:107:LEU:HA	1.81	0.61
1:B:94:VAL:HG11	3:B:804:PXP:H2A1	1.82	0.61
1:C:48:GLU:OE1	1:C:98:ARG:HB2	2.00	0.61
1:A:47:ARG:HD3	1:A:51:ARG:HG2	1.84	0.60
1:C:12:HIS:CA	1:C:15:THR:HG23	2.14	0.60
1:B:14:ALA:HB3	1:B:52:HIS:HB2	1.83	0.60
1:B:51:ARG:NH2	1:B:98:ARG:O	2.35	0.59
1:C:14:ALA:HB3	1:C:52:HIS:HB2	1.85	0.59
1:B:195:LEU:HA	1:B:199:ASN:HD21	1.67	0.58
1:A:199:ASN:C	1:A:199:ASN:ND2	2.52	0.58
1:D:112:GLN:NE2	4:D:832:HOH:O	2.36	0.58
1:C:57:ASP:O	1:C:61:LEU:HD13	2.03	0.58
1:D:50:ARG:NH1	1:D:55:ASP:OD1	2.37	0.58
1:C:221:ARG:NH1	4:C:934:HOH:O	2.37	0.57
1:A:91:CYS:HB2	1:A:128:ILE:HG21	1.86	0.57
1:D:195:LEU:HA	1:D:199:ASN:HD21	1.70	0.57
1:B:199:ASN:HD22	1:B:200:VAL:N	2.03	0.56
1:B:97:LYS:O	1:B:98:ARG:CB	2.53	0.56
1:C:28:VAL:HG22	1:C:61:LEU:HD12	1.87	0.56
1:D:196:THR:H	1:D:199:ASN:HD21	1.50	0.56
1:C:84:VAL:O	1:C:87:LYS:HD2	2.06	0.56
1:D:12:HIS:CD2	1:D:216:HIS:HD2	2.23	0.56
1:D:102:THR:N	4:D:870:HOH:O	2.38	0.56
1:C:68:ARG:NH2	1:C:211:GLU:OE1	2.39	0.55
1:A:199:ASN:HD22	1:A:200:VAL:N	2.04	0.55
1:C:196:THR:H	1:C:199:ASN:ND2	2.05	0.55
1:D:199:ASN:HD22	1:D:200:VAL:N	2.04	0.54
1:A:108:ASP:O	1:A:112:GLN:HG3	2.07	0.54
1:C:199:ASN:HD22	1:C:200:VAL:N	2.04	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:33:ILE:HG21	1:D:227:LEU:HB3	1.90	0.53
1:C:221:ARG:HB2	1:C:221:ARG:NH1	2.19	0.53
1:A:122:ARG:NE	4:A:984:HOH:O	2.36	0.53
1:A:133:PHE:CD2	3:A:802:PXP:H2A2	2.44	0.53
1:B:63:GLN:HE22	1:D:59:ARG:NH1	1.87	0.53
1:D:133:PHE:CD2	3:D:801:PXP:H2A2	2.43	0.53
1:C:12:HIS:HD2	1:C:216:HIS:ND1	2.06	0.53
1:C:56:ARG:HD3	1:C:56:ARG:C	2.29	0.53
1:D:56:ARG:HD3	1:D:56:ARG:C	2.28	0.53
1:B:103:THR:CB	3:B:804:PXP:H2A3	2.39	0.52
1:A:227:LEU:HD22	1:A:231:VAL:HG23	1.92	0.52
1:A:56:ARG:HD2	4:C:947:HOH:O	2.10	0.52
1:B:20:ARG:HG2	1:B:20:ARG:HH11	1.75	0.52
1:D:108:ASP:HB2	4:D:870:HOH:O	2.10	0.52
1:B:46:LEU:HD21	1:B:50:ARG:CD	2.40	0.52
1:A:195:LEU:HA	1:A:199:ASN:HD21	1.74	0.51
1:B:196:THR:H	1:B:199:ASN:ND2	2.07	0.51
1:C:225:THR:HG21	4:C:919:HOH:O	2.10	0.51
1:D:100:GLU:O	4:D:832:HOH:O	2.19	0.51
1:A:57:ASP:O	1:A:61:LEU:HB2	2.10	0.51
1:B:51:ARG:HG2	4:B:1024:HOH:O	2.12	0.50
1:D:164:THR:OG1	1:D:167:GLU:HB2	2.11	0.50
1:B:199:ASN:C	1:B:199:ASN:ND2	2.60	0.50
1:A:87:LYS:N	1:A:88:PRO:HD3	2.25	0.50
1:B:95:PRO:HG3	1:B:107:LEU:HD23	1.93	0.50
1:B:12:HIS:CD2	1:B:216:HIS:HD2	2.30	0.50
1:B:91:CYS:HB2	1:B:128:ILE:HG21	1.94	0.50
1:D:57:ASP:O	1:D:61:LEU:HB2	2.12	0.49
1:A:3:GLU:HB3	4:A:1032:HOH:O	2.11	0.49
1:D:75:VAL:HG22	1:D:93:LEU:HD13	1.94	0.49
1:C:12:HIS:CD2	1:C:216:HIS:ND1	2.81	0.49
1:B:133:PHE:HZ	1:B:155:HIS:HD2	1.61	0.48
1:D:50:ARG:HH11	1:D:50:ARG:HG2	1.78	0.48
1:D:47:ARG:NE	1:D:51:ARG:NH1	2.60	0.48
1:C:91:CYS:HB2	1:C:128:ILE:HG21	1.95	0.48
1:C:28:VAL:HG22	1:C:61:LEU:CD1	2.43	0.48
1:B:20:ARG:HG2	1:B:20:ARG:NH1	2.28	0.48
1:D:117:ARG:NH1	1:D:146:GLU:OE2	2.44	0.48
1:B:61:LEU:O	1:B:65:LEU:HG	2.14	0.47
1:D:84:VAL:O	1:D:87:LYS:HD2	2.15	0.47
1:B:133:PHE:CZ	1:B:155:HIS:HD2	2.32	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:LEU:HG	1:A:238:MET:SD	2.55	0.47
1:D:91:CYS:HB2	1:D:128:ILE:HG21	1.95	0.47
1:C:199:ASN:C	1:C:199:ASN:ND2	2.63	0.46
1:B:147:VAL:CG1	1:B:147:VAL:O	2.63	0.46
1:D:12:HIS:CD2	1:D:216:HIS:CD2	3.03	0.46
1:D:227:LEU:HD22	1:D:231:VAL:HG23	1.96	0.46
1:D:242:ARG:NH2	4:D:913:HOH:O	2.46	0.46
1:D:14:ALA:HB3	1:D:52:HIS:HB2	1.97	0.46
1:D:153:GLU:OE1	3:D:801:PXP:N1	2.49	0.45
1:C:141:ILE:HD11	1:C:178:ALA:HB1	1.99	0.45
1:B:48:GLU:OE1	1:B:98:ARG:HD2	2.16	0.45
1:C:133:PHE:HZ	1:C:155:HIS:HD2	1.63	0.45
1:D:117:ARG:NH2	1:D:146:GLU:OE2	2.49	0.45
1:B:4:LEU:HD13	1:B:238:MET:HE3	1.99	0.45
1:D:95:PRO:HD3	1:D:107:LEU:HB2	2.00	0.44
1:D:104:GLU:OE2	1:D:174:ARG:NH2	2.41	0.44
1:D:2:ALA:HB2	4:D:901:HOH:O	2.18	0.44
1:B:193:HIS:CD2	4:B:972:HOH:O	2.70	0.44
1:A:9:ASN:HA	1:A:43:THR:O	2.18	0.43
1:A:113:ARG:HD3	4:A:992:HOH:O	2.18	0.43
1:C:133:PHE:CD2	3:C:803:PXP:H2A2	2.53	0.43
1:D:171:GLU:OE2	1:D:174:ARG:NH1	2.50	0.43
1:D:199:ASN:ND2	1:D:199:ASN:C	2.66	0.43
1:D:80:LEU:O	1:D:84:VAL:HG23	2.18	0.43
1:A:72:GLU:HG2	1:A:92:CYS:HB3	2.00	0.43
1:B:46:LEU:CD2	1:B:50:ARG:HD3	2.49	0.43
1:B:46:LEU:HD21	1:B:50:ARG:NE	2.34	0.43
1:D:135:ASP:O	1:D:137:ASP:N	2.47	0.42
1:C:161:ASP:O	1:C:162:ALA:C	2.55	0.42
1:B:146:GLU:HG2	4:B:1044:HOH:O	2.19	0.42
1:C:221:ARG:CB	1:C:221:ARG:HH11	2.25	0.42
1:C:47:ARG:HD3	1:C:51:ARG:HB3	2.01	0.42
1:A:121:LYS:HE3	4:A:1086:HOH:O	2.19	0.42
1:B:20:ARG:HA	1:B:20:ARG:HD3	1.87	0.42
1:D:91:CYS:HB2	1:D:128:ILE:CG2	2.49	0.42
1:C:12:HIS:ND1	1:C:15:THR:HG21	2.34	0.42
1:D:155:HIS:CD2	1:D:157:GLY:N	2.79	0.42
1:D:31:ALA:O	1:D:35:GLU:HG3	2.19	0.41
1:D:204:ALA:O	1:D:242:ARG:HD3	2.19	0.41
1:A:113:ARG:O	1:A:117:ARG:HB2	2.20	0.41
1:B:233:GLU:CD	1:B:236:ARG:HH21	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:VAL:HG13	1:B:201:LYS:N	2.35	0.41
1:A:153:GLU:OE1	3:A:802:PXP:N1	2.53	0.41
1:B:20:ARG:HH11	1:B:20:ARG:CG	2.32	0.41
1:C:212:LEU:HD12	1:C:238:MET:HE3	2.01	0.41
1:D:206:ILE:HA	1:D:207:PRO:HD3	1.75	0.41
1:B:12:HIS:CD2	1:B:216:HIS:CD2	3.08	0.41
1:D:50:ARG:HH12	1:D:55:ASP:CG	2.24	0.41
1:B:171:GLU:OE2	1:B:171:GLU:HA	2.20	0.41
1:B:196:THR:H	1:B:199:ASN:HD21	1.69	0.41
1:A:47:ARG:NH2	1:A:100:GLU:OE1	2.54	0.41
1:B:161:ASP:O	1:B:162:ALA:C	2.58	0.41
1:A:49:ASP:C	1:A:49:ASP:OD1	2.60	0.41
1:D:192:GLY:O	1:D:193:HIS:CB	2.68	0.41
1:A:155:HIS:CE1	1:A:157:GLY:HA3	2.56	0.41
1:D:68:ARG:HG2	4:D:963:HOH:O	2.20	0.41
1:B:133:PHE:CZ	1:B:155:HIS:CD2	3.10	0.40
1:B:103:THR:OG1	3:B:804:PXP:H2A3	2.21	0.40
1:C:94:VAL:HG11	3:C:803:PXP:H2A1	2.04	0.40
1:A:225:THR:CG2	4:A:942:HOH:O	2.70	0.40
3:B:804:PXP:H4A1	3:B:804:PXP:H5A2	1.88	0.40
1:D:89:HIS:CD2	1:D:90:PHE:CE1	3.04	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	240/242 (99%)	230 (96%)	8 (3%)	2 (1%)	24	27
1	B	240/242 (99%)	229 (95%)	8 (3%)	3 (1%)	15	15
1	C	240/242 (99%)	231 (96%)	8 (3%)	1 (0%)	39	48
1	D	240/242 (99%)	228 (95%)	11 (5%)	1 (0%)	39	48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	960/968 (99%)	918 (96%)	35 (4%)	7 (1%)	26	31

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	98	ARG
1	A	106	GLY
1	B	104	GLU
1	C	193	HIS
1	D	193	HIS
1	B	193	HIS
1	A	193	HIS

### 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	184/185 (100%)	168 (91%)	16 (9%)	13	15
1	B	185/185 (100%)	171 (92%)	14 (8%)	16	20
1	C	185/185 (100%)	171 (92%)	14 (8%)	16	20
1	D	183/185 (99%)	167 (91%)	16 (9%)	13	15
All	All	737/740 (100%)	677 (92%)	60 (8%)	15	18

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	4	LEU
1	A	6	LEU
1	A	22	THR
1	A	46	LEU
1	A	48	GLU
1	A	56	ARG
1	A	61	LEU

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Mol	Chain	Res	Type
1	A	71	LEU
1	A	104	GLU
1	A	138	GLU
1	A	172	LEU
1	A	189	VAL
1	A	199	ASN
1	A	223	VAL
1	A	227	LEU
1	B	3	GLU
1	B	4	LEU
1	B	6	LEU
1	B	20	ARG
1	B	46	LEU
1	B	55	ASP
1	B	56	ARG
1	B	61	LEU
1	B	71	LEU
1	B	85	GLU
1	B	172	LEU
1	B	189	VAL
1	B	199	ASN
1	B	227	LEU
1	C	3	GLU
1	C	4	LEU
1	C	6	LEU
1	C	15	THR
1	C	22	THR
1	C	46	LEU
1	C	68	ARG
1	C	71	LEU
1	C	172	LEU
1	C	189	VAL
1	C	199	ASN
1	C	221	ARG
1	C	225	THR
1	C	227	LEU
1	D	4	LEU
1	D	46	LEU
1	D	50	ARG
1	D	51	ARG
1	D	56	ARG
1	D	61	LEU

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Mol	Chain	Res	Type
1	D	71	LEU
1	D	91	CYS
1	D	101	VAL
1	D	118	ASP
1	D	147	VAL
1	D	172	LEU
1	D	185	LEU
1	D	189	VAL
1	D	199	ASN
1	D	227	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	HIS
1	A	63	GLN
1	A	89	HIS
1	A	155	HIS
1	A	199	ASN
1	A	210	HIS
1	A	216	HIS
1	B	63	GLN
1	B	89	HIS
1	B	129	GLN
1	B	155	HIS
1	B	193	HIS
1	B	199	ASN
1	B	210	HIS
1	B	216	HIS
1	C	12	HIS
1	C	63	GLN
1	C	89	HIS
1	C	155	HIS
1	C	199	ASN
1	C	210	HIS
1	D	12	HIS
1	D	63	GLN
1	D	89	HIS
1	D	155	HIS
1	D	199	ASN
1	D	210	HIS
1	D	216	HIS



### 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 ⓘ

7 ligands 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	PXP	A	802	-	16,16,16	1.55	4 (25%)	20,23,23	1.39	3 (15%)
2	PO4	A	901	-	4,4,4	0.35	0	6,6,6	0.29	0
3	PXP	B	804	-	16,16,16	1.66	5 (31%)	20,23,23	1.97	6 (30%)
2	PO4	B	903	-	4,4,4	0.24	0	6,6,6	0.30	0
3	PXP	C	803	-	16,16,16	1.63	4 (25%)	20,23,23	1.50	5 (25%)
2	PO4	C	902	-	4,4,4	0.21	0	6,6,6	0.30	0
3	PXP	D	801	-	16,16,16	1.62	6 (37%)	20,23,23	1.51	3 (15%)

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	PXP	A	802	-	-	0/8/8/8	0/1/1/1
2	PO4	A	901	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PXP	B	804	-	-	0/8/8/8	0/1/1/1
2	PO4	B	903	-	-	0/0/0/0	0/0/0/0
3	PXP	C	803	-	-	0/8/8/8	0/1/1/1
2	PO4	C	902	-	-	0/0/0/0	0/0/0/0
3	PXP	D	801	-	-	0/8/8/8	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	804	PXP	P-O3P	-2.84	1.44	1.54
3	C	803	PXP	P-O3P	-2.32	1.46	1.54
3	A	802	PXP	P-O3P	-2.26	1.46	1.54
3	D	801	PXP	P-O3P	-2.20	1.46	1.54
3	A	802	PXP	C3-C2	-2.18	1.39	1.40
3	D	801	PXP	C3-C2	-2.18	1.39	1.40
3	C	803	PXP	C3-C2	-2.11	1.39	1.40
3	B	804	PXP	P-O4P	-2.08	1.53	1.60
3	B	804	PXP	C4A-C4	-2.08	1.48	1.51
3	D	801	PXP	C5-C4	2.01	1.43	1.40
3	B	804	PXP	C6-N1	2.11	1.39	1.34
3	A	802	PXP	C2-N1	2.32	1.39	1.34
3	D	801	PXP	C6-N1	2.37	1.39	1.34
3	D	801	PXP	C2-N1	2.57	1.39	1.34
3	A	802	PXP	C6-N1	2.64	1.40	1.34
3	C	803	PXP	C2A-C2	2.68	1.55	1.50
3	C	803	PXP	C2-N1	2.82	1.40	1.34
3	B	804	PXP	C2-N1	2.85	1.40	1.34
3	D	801	PXP	C2A-C2	2.94	1.56	1.50

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	804	PXP	O4P-P-O1P	-4.12	96.65	107.14
3	B	804	PXP	O4P-C5A-C5	-3.65	102.96	108.99
3	A	802	PXP	O4P-P-O1P	-2.59	100.56	107.14
3	C	803	PXP	O4P-P-O1P	-2.40	101.03	107.14
3	B	804	PXP	C5-C6-N1	-2.34	119.79	123.86
3	D	801	PXP	C5-C6-N1	-2.31	119.84	123.86
3	C	803	PXP	C5-C6-N1	-2.25	119.95	123.86
3	D	801	PXP	O4P-P-O1P	-2.15	101.68	107.14
3	A	802	PXP	C5-C6-N1	-2.14	120.15	123.86
3	C	803	PXP	O2P-P-O4P	-2.08	100.58	106.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	804	PXP	O4A-C4A-C4	-2.06	105.16	111.64
3	B	804	PXP	O3-C3-C2	2.00	121.14	117.66
3	C	803	PXP	O3P-P-O1P	2.01	117.06	110.58
3	C	803	PXP	O3P-P-O4P	2.75	114.47	106.56
3	A	802	PXP	O3P-P-O4P	3.06	115.36	106.56
3	D	801	PXP	O3P-P-O4P	3.27	115.99	106.56
3	B	804	PXP	O3P-P-O4P	3.93	117.88	106.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	802	PXP	2	0
3	B	804	PXP	4	0
3	C	803	PXP	2	0
3	D	801	PXP	2	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 ⓘ

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	242/242 (100%)	-0.19	9 (3%) 45 54	21, 34, 53, 80	0
1	B	242/242 (100%)	-0.09	15 (6%) 24 32	22, 34, 58, 93	0
1	C	242/242 (100%)	-0.11	11 (4%) 37 46	20, 32, 53, 81	0
1	D	242/242 (100%)	-0.16	8 (3%) 50 59	21, 32, 50, 82	0
All	All	968/968 (100%)	-0.14	43 (4%) 38 47	20, 33, 54, 93	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	98	ARG	6.0
1	B	99	GLN	5.6
1	D	99	GLN	5.5
1	B	97	LYS	5.4
1	A	98	ARG	5.3
1	C	97	LYS	4.8
1	B	102	THR	4.6
1	C	99	GLN	4.5
1	C	98	ARG	4.3
1	C	101	VAL	4.3
1	D	103	THR	4.3
1	B	104	GLU	4.2
1	A	99	GLN	4.0
1	B	114	ASP	3.9
1	B	2	ALA	3.5
1	C	104	GLU	3.4
1	C	102	THR	3.1
1	D	2	ALA	3.1
1	D	101	VAL	3.0
1	A	104	GLU	2.9
1	B	105	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	243	GLY	2.8
1	B	163	LYS	2.8
1	B	103	THR	2.7
1	B	101	VAL	2.7
1	C	100	GLU	2.6
1	A	101	VAL	2.5
1	C	103	THR	2.4
1	A	46	LEU	2.4
1	C	105	GLY	2.4
1	D	163	LYS	2.4
1	A	97	LYS	2.4
1	B	162	ALA	2.4
1	A	163	LYS	2.3
1	D	125	ASP	2.3
1	C	243	GLY	2.3
1	A	2	ALA	2.2
1	C	21	GLY	2.2
1	A	105	GLY	2.2
1	B	100	GLU	2.1
1	D	97	LYS	2.0
1	B	111	GLY	2.0
1	B	166	ALA	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. 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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	PXP	B	804	16/16	0.96	0.17	0.30	30,42,47,49	0
2	PO4	B	903	5/5	0.89	0.15	-0.25	67,70,72,73	0
2	PO4	C	902	5/5	0.91	0.16	-0.28	60,64,67,69	0
3	PXP	C	803	16/16	0.98	0.09	-0.68	26,32,38,45	0
2	PO4	A	901	5/5	0.97	0.10	-0.92	51,54,55,56	0
3	PXP	D	801	16/16	0.98	0.08	-1.12	24,32,38,46	0
3	PXP	A	802	16/16	0.99	0.08	-1.35	28,32,37,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.