



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 12:09 am BST

PDB ID : 1IXO
Title : Enzyme-analogue substrate complex of Pyridoxine 5'-Phosphate Synthase
Authors : Garrido-Franco, M.; Laber, B.; Huber, R.; Clausen, T.
Deposited on : 2002-06-28
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

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

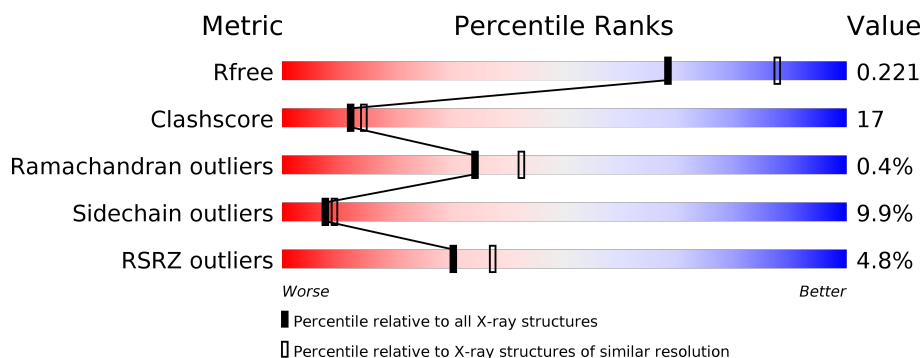
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 ≥ 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	242	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>• •</div> </div> </div>
1	B	242	<div> <div>6%</div> <div> <div></div> <div>68%</div> <div>25%</div> <div>5%</div> <div>•</div> </div> </div>
1	C	242	<div> <div>5%</div> <div> <div></div> <div>71%</div> <div>21%</div> <div>• •</div> </div> </div>
1	D	242	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>22%</div> <div>• •</div> </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	G3P	C	1003	-	-	X	-
2	G3P	C	1005	-	-	-	X

2 Entry composition [i](#)

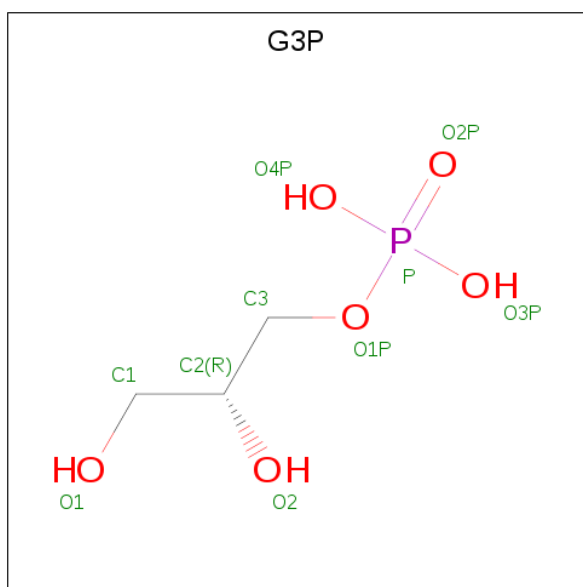
There are 3 unique types of molecules in this entry. The entry contains 7474 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	235	Total	C	N	O	S	0	0	0
			1774	1101	325	336	12			
1	B	235	Total	C	N	O	S	0	0	0
			1778	1103	325	338	12			
1	C	235	Total	C	N	O	S	0	0	0
			1770	1099	325	334	12			
1	D	235	Total	C	N	O	S	0	0	0
			1774	1100	324	338	12			

- Molecule 2 is SN-GLYCEROL-3-PHOSPHATE (three-letter code: G3P) (formula: $C_3H_9O_6P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			10	3	6	1		
2	B	1	Total	C	O	P	0	0
			10	3	6	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	O	P	0	0
			10	3	6	1		
2	C	1	Total	C	O	P	0	0
			10	3	6	1		
2	D	1	Total	C	O	P	0	0
			10	3	6	1		

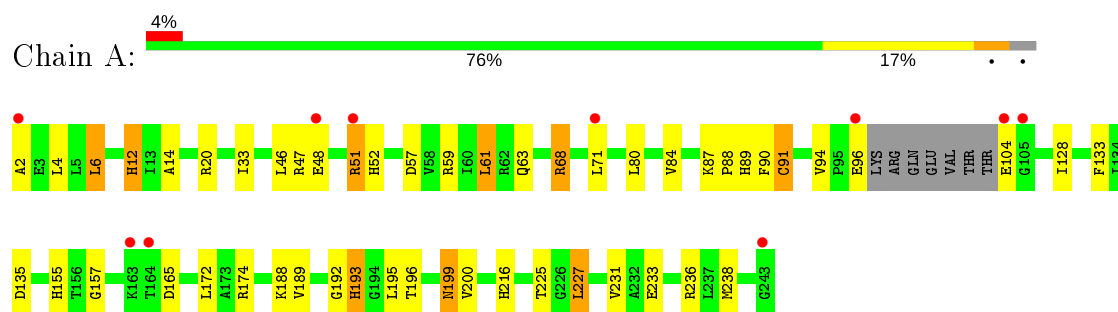
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	85	Total	O	0	0
			85	85		
3	B	75	Total	O	0	0
			75	75		
3	C	90	Total	O	0	0
			90	90		
3	D	78	Total	O	0	0
			78	78		

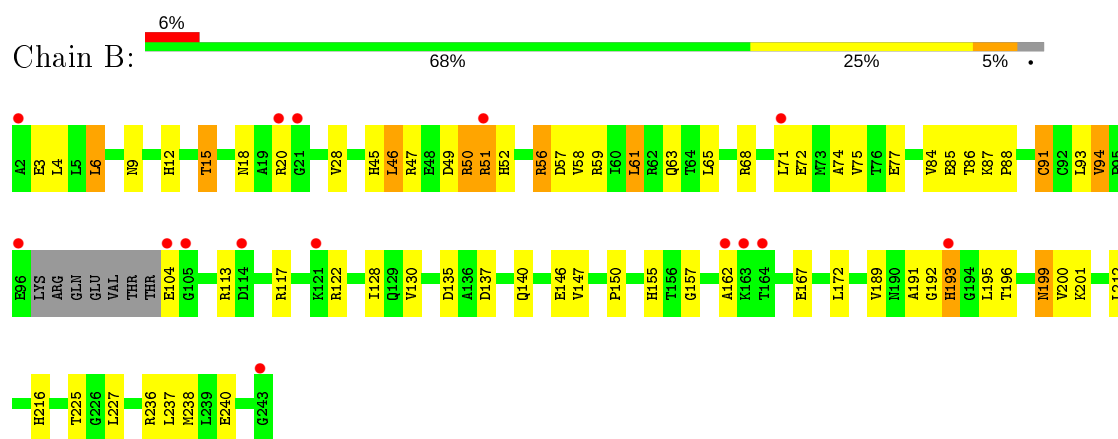
3 Residue-property plots [i](#)

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.

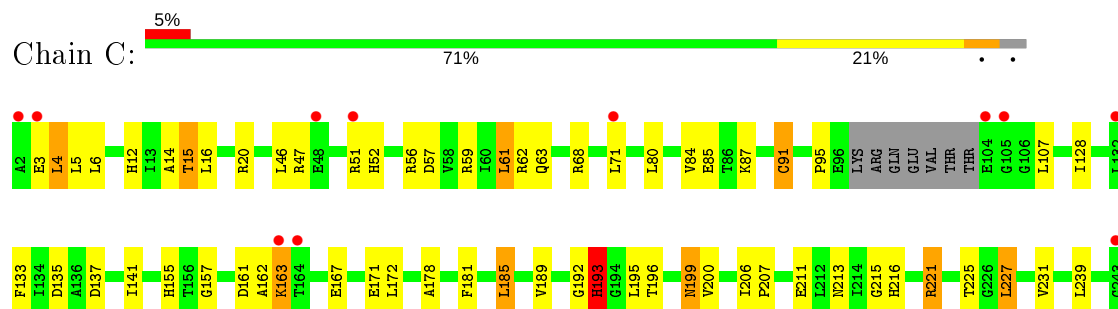
• Molecule 1: Pyridoxine 5'-Phosphate synthase



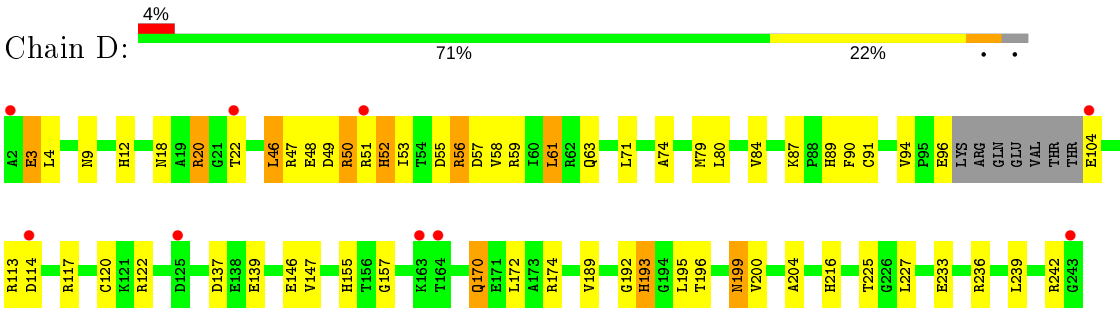
• Molecule 1: Pyridoxine 5'-Phosphate synthase



• 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, α , β , γ	129.80 Å 155.90 Å 127.40 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.96 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.30) 99.4 (19.96-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.50	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.30 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.220 , 0.260 0.213 , 0.221	Depositor DCC
R_{free} test set	2892 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	37.4	Xtriage
Anisotropy	0.924	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7474	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: G3P

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	0.47	0/1795	0.71	0/2422
1	B	0.48	0/1799	0.71	0/2427
1	C	0.48	0/1791	0.72	0/2417
1	D	0.47	0/1795	0.71	0/2423
All	All	0.48	0/7180	0.71	0/9689

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	1774	0	1786	46	0
1	B	1778	0	1789	83	1
1	C	1770	0	1782	51	0
1	D	1774	0	1779	67	0
2	A	10	0	6	0	0
2	B	10	0	6	3	0
2	C	20	0	11	6	1
2	D	10	0	6	1	0
3	A	85	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	75	0	0	2	0
3	C	90	0	0	4	0
3	D	78	0	0	5	0
All	All	7474	0	7165	244	1

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 17.

All (244) 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:B:15:THR:HG22	1:B:51:ARG:NH1	1.53	1.22
1:B:15:THR:HG22	1:B:51:ARG:HH12	0.99	1.10
1:B:51:ARG:HD3	1:B:52:HIS:CD2	1.89	1.05
1:B:12:HIS:HA	1:B:15:THR:HG23	1.34	1.04
1:B:51:ARG:NE	1:B:52:HIS:N	2.07	1.00
1:C:12:HIS:HA	1:C:15:THR:HG23	1.46	0.98
1:B:15:THR:CG2	1:B:51:ARG:HH12	1.78	0.97
1:A:63:GLN:HE22	1:C:59:ARG:HH12	1.06	0.94
1:C:192:GLY:O	1:C:193:HIS:HB2	1.66	0.93
1:B:15:THR:HG22	1:B:51:ARG:CZ	2.00	0.91
1:B:15:THR:HA	1:B:51:ARG:NH2	1.85	0.90
1:A:12:HIS:HD2	1:A:216:HIS:HD2	1.20	0.88
1:B:49:ASP:O	1:B:50:ARG:HB2	1.71	0.88
1:B:59:ARG:HH12	1:D:63:GLN:HE22	1.19	0.86
1:B:46:LEU:HD11	1:B:50:ARG:NH1	1.91	0.86
1:D:52:HIS:CD2	1:D:53:ILE:H	1.94	0.85
1:B:155:HIS:HD2	1:B:157:GLY:H	1.23	0.84
1:C:161:ASP:O	1:C:163:LYS:HE2	1.78	0.83
1:B:15:THR:CG2	1:B:51:ARG:NH1	2.40	0.82
1:B:51:ARG:HE	1:B:52:HIS:N	1.72	0.82
1:B:155:HIS:CD2	1:B:157:GLY:H	1.97	0.82
2:C:1003:G3P:O2P	3:C:1036:HOH:O	1.97	0.81
1:D:52:HIS:HD2	1:D:53:ILE:H	1.29	0.80
1:B:15:THR:CA	1:B:51:ARG:HH22	1.95	0.80
1:B:15:THR:CA	1:B:51:ARG:NH2	2.44	0.80
1:A:199:ASN:C	1:A:199:ASN:HD22	1.85	0.79
1:A:51:ARG:C	1:A:51:ARG:HD2	2.01	0.79
1:B:75:VAL:HG22	1:B:93:LEU:HD13	1.63	0.79
1:D:117:ARG:HA	1:D:147:VAL:HG13	1.65	0.79
1:B:51:ARG:CD	1:B:52:HIS:CD2	2.67	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:199:ASN:C	1:C:199:ASN:HD22	1.86	0.78
1:D:51:ARG:HG2	1:D:51:ARG:HH11	1.49	0.77
1:D:12:HIS:HD2	1:D:216:HIS:HD2	1.33	0.75
1:A:155:HIS:CD2	1:A:157:GLY:H	2.05	0.74
1:D:20:ARG:NH1	1:D:20:ARG:HB2	2.01	0.74
1:B:225:THR:HG22	3:B:1010:HOH:O	1.85	0.74
1:D:20:ARG:HG3	1:D:22:THR:HG23	1.70	0.73
1:B:15:THR:HG22	1:B:51:ARG:NH2	2.04	0.72
1:D:50:ARG:N	1:D:50:ARG:HD3	2.03	0.72
1:D:199:ASN:HD22	1:D:199:ASN:C	1.93	0.72
1:A:59:ARG:HH12	1:C:63:GLN:HE22	1.38	0.71
1:A:196:THR:H	1:A:199:ASN:ND2	1.89	0.71
1:A:84:VAL:O	1:A:87:LYS:HD2	1.91	0.70
1:A:14:ALA:HB3	1:A:52:HIS:HB2	1.73	0.70
1:B:195:LEU:HA	1:B:199:ASN:HD21	1.57	0.70
1:D:47:ARG:NH1	1:D:51:ARG:HD2	2.07	0.69
1:B:15:THR:N	1:B:51:ARG:HH22	1.89	0.69
1:B:199:ASN:HD22	1:B:199:ASN:C	1.95	0.69
1:C:5:LEU:HB2	1:C:211:GLU:HG3	1.74	0.69
1:B:51:ARG:HE	1:B:51:ARG:C	1.94	0.69
1:C:196:THR:H	1:C:199:ASN:ND2	1.91	0.69
1:D:46:LEU:HB3	1:D:79:MET:HE1	1.74	0.69
1:D:12:HIS:CD2	1:D:216:HIS:HD2	2.11	0.68
1:B:196:THR:H	1:B:199:ASN:ND2	1.91	0.68
1:B:63:GLN:HE22	1:D:59:ARG:HH12	1.40	0.68
1:D:155:HIS:CD2	1:D:157:GLY:H	2.11	0.68
1:C:12:HIS:HD2	1:C:216:HIS:ND1	1.92	0.68
1:A:96:GLU:H	1:A:96:GLU:CD	1.96	0.67
1:D:204:ALA:O	1:D:242:ARG:HD3	1.94	0.67
1:A:225:THR:HG22	3:A:1011:HOH:O	1.96	0.66
1:A:48:GLU:HG3	3:A:1035:HOH:O	1.95	0.66
1:B:47:ARG:CZ	1:B:51:ARG:HG2	2.26	0.66
1:D:57:ASP:O	1:D:61:LEU:HB2	1.97	0.65
1:D:196:THR:H	1:D:199:ASN:ND2	1.95	0.65
1:A:68:ARG:NE	1:A:90:PHE:HE2	1.95	0.64
1:A:12:HIS:CD2	1:A:216:HIS:HD2	2.10	0.64
1:C:12:HIS:HE1	2:C:1005:G3P:O4P	1.80	0.64
1:D:12:HIS:HD2	1:D:216:HIS:CD2	2.15	0.64
1:A:155:HIS:HD2	1:A:157:GLY:H	1.44	0.64
1:A:192:GLY:O	1:A:193:HIS:HB2	1.97	0.64
1:C:84:VAL:O	1:C:87:LYS:HD2	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:ARG:HD2	3:D:1010:HOH:O	1.97	0.63
1:D:104:GLU:OE2	1:D:174:ARG:NH2	2.32	0.63
1:D:50:ARG:HG3	1:D:50:ARG:HH11	1.64	0.63
1:C:199:ASN:HD22	1:C:200:VAL:N	1.97	0.62
1:B:12:HIS:ND1	1:B:15:THR:HG21	2.14	0.62
1:D:47:ARG:O	1:D:50:ARG:HD2	2.00	0.62
1:D:20:ARG:HH11	1:D:20:ARG:HB2	1.63	0.61
1:B:61:LEU:O	1:B:65:LEU:HG	2.00	0.61
1:B:196:THR:H	1:B:199:ASN:HD21	1.48	0.61
1:D:170:GLN:NE2	1:D:170:GLN:HA	2.15	0.61
1:C:12:HIS:HA	1:C:15:THR:CG2	2.25	0.61
1:A:196:THR:H	1:A:199:ASN:HD21	1.46	0.61
1:B:46:LEU:CD1	1:B:50:ARG:NH1	2.63	0.61
1:D:225:THR:HG22	3:D:1012:HOH:O	2.01	0.60
1:C:12:HIS:CD2	1:C:216:HIS:ND1	2.69	0.60
1:D:12:HIS:CD2	1:D:216:HIS:CD2	2.90	0.60
1:B:12:HIS:HA	1:B:15:THR:CG2	2.21	0.59
1:B:15:THR:HG22	1:B:51:ARG:HH22	1.67	0.59
1:B:18:ASN:ND2	1:B:51:ARG:O	2.34	0.59
1:D:117:ARG:CA	1:D:147:VAL:HG13	2.32	0.59
1:A:12:HIS:HD2	1:A:216:HIS:CD2	2.10	0.59
1:A:233:GLU:CD	1:A:236:ARG:HH21	2.05	0.59
1:A:104:GLU:HB2	1:A:135:ASP:OD1	2.03	0.59
1:C:215:GLY:N	2:C:1003:G3P:O4P	2.30	0.59
1:D:74:ALA:HB2	1:D:96:GLU:HB2	1.84	0.58
1:C:196:THR:H	1:C:199:ASN:HD21	1.52	0.58
1:B:117:ARG:HA	1:B:147:VAL:HG13	1.85	0.58
1:D:50:ARG:CD	1:D:50:ARG:N	2.66	0.58
1:C:80:LEU:O	1:C:84:VAL:HG23	2.02	0.58
1:C:12:HIS:CE1	2:C:1005:G3P:O4P	2.56	0.57
1:C:47:ARG:NH2	1:C:51:ARG:HH21	2.02	0.57
1:C:221:ARG:HD3	3:C:1073:HOH:O	2.05	0.57
1:D:89:HIS:HD2	1:D:90:PHE:CE1	2.22	0.57
1:D:84:VAL:O	1:D:87:LYS:HD2	2.05	0.57
1:B:12:HIS:CD2	1:B:216:HIS:HD2	2.23	0.57
1:B:130:VAL:O	1:B:150:PRO:HD2	2.05	0.56
1:B:199:ASN:HD22	1:B:200:VAL:N	2.02	0.56
1:A:51:ARG:HE	1:A:52:HIS:CD2	2.23	0.56
1:B:104:GLU:O	1:B:104:GLU:HG2	2.06	0.56
1:D:120:CYS:HB3	3:D:1049:HOH:O	2.05	0.56
1:B:45:HIS:HA	1:B:72:GLU:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:181:PHE:CZ	1:C:185:LEU:HD21	2.42	0.55
1:B:74:ALA:HA	1:B:94:VAL:O	2.07	0.55
1:A:199:ASN:C	1:A:199:ASN:ND2	2.58	0.54
1:D:50:ARG:HH12	1:D:55:ASP:CG	2.10	0.54
1:A:51:ARG:HD2	1:A:51:ARG:O	2.07	0.54
1:D:233:GLU:CD	1:D:236:ARG:HH21	2.11	0.54
1:C:47:ARG:HH21	1:C:51:ARG:HH21	1.54	0.54
1:B:137:ASP:OD2	1:B:140:GLN:HG2	2.09	0.53
1:B:51:ARG:HD3	1:B:52:HIS:CG	2.41	0.53
1:D:196:THR:H	1:D:199:ASN:HD21	1.55	0.53
1:A:80:LEU:O	1:A:84:VAL:HG23	2.07	0.53
1:C:15:THR:HG22	1:C:52:HIS:HB3	1.90	0.53
1:C:195:LEU:HA	1:C:199:ASN:HD21	1.73	0.53
1:A:6:LEU:HG	1:A:238:MET:SD	2.49	0.53
1:A:87:LYS:N	1:A:88:PRO:HD3	2.24	0.53
1:B:15:THR:CG2	1:B:51:ARG:HH22	2.22	0.52
1:A:199:ASN:HD22	1:A:200:VAL:N	2.06	0.52
1:C:199:ASN:C	1:C:199:ASN:ND2	2.57	0.52
1:C:5:LEU:CB	1:C:211:GLU:HG3	2.38	0.52
1:B:9:ASN:HD22	2:B:1002:G3P:H2	1.74	0.51
1:C:133:PHE:CZ	1:C:193:HIS:CD2	2.98	0.51
1:B:201:LYS:HE3	1:B:237:LEU:HD22	1.93	0.51
1:C:221:ARG:NH1	3:C:1037:HOH:O	2.42	0.51
1:D:113:ARG:HD3	1:D:146:GLU:OE1	2.11	0.51
1:D:3:GLU:OE2	1:D:239:LEU:HD21	2.11	0.51
1:B:236:ARG:NH1	1:B:240:GLU:OE2	2.43	0.51
1:C:3:GLU:CG	1:C:239:LEU:HD22	2.41	0.51
1:D:56:ARG:C	1:D:56:ARG:HD3	2.30	0.51
1:B:15:THR:CG2	1:B:51:ARG:NH2	2.74	0.51
1:B:9:ASN:ND2	2:B:1002:G3P:H2	2.26	0.50
1:D:20:ARG:HG3	1:D:20:ARG:O	2.11	0.50
1:B:50:ARG:O	1:B:51:ARG:C	2.49	0.50
1:D:74:ALA:HA	1:D:94:VAL:O	2.12	0.49
1:B:236:ARG:O	1:B:240:GLU:HG3	2.13	0.49
1:B:9:ASN:HD22	2:B:1002:G3P:H11	1.77	0.49
1:C:56:ARG:HD3	1:C:56:ARG:C	2.33	0.49
1:C:137:ASP:O	1:C:141:ILE:HG13	2.11	0.49
1:D:20:ARG:HH11	1:D:20:ARG:CB	2.26	0.49
1:B:49:ASP:O	1:B:50:ARG:CB	2.45	0.49
1:D:199:ASN:ND2	1:D:199:ASN:C	2.65	0.49
1:D:195:LEU:HA	1:D:199:ASN:HD21	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:LEU:HA	1:A:199:ASN:HD21	1.78	0.48
1:A:57:ASP:O	1:A:61:LEU:HB2	2.14	0.48
1:B:28:VAL:HG22	1:B:61:LEU:HD13	1.94	0.48
1:D:137:ASP:OD2	1:D:139:GLU:HB3	2.14	0.48
1:B:51:ARG:CD	1:B:52:HIS:CG	2.95	0.48
1:B:75:VAL:CG2	1:B:93:LEU:HD13	2.39	0.48
1:D:114:ASP:HB2	3:D:1041:HOH:O	2.13	0.48
1:C:52:HIS:HB3	3:C:1093:HOH:O	2.12	0.48
1:C:213:ASN:HB3	2:C:1003:G3P:H11	1.95	0.47
1:C:14:ALA:HB3	1:C:52:HIS:HB2	1.96	0.47
1:C:3:GLU:HG3	1:C:4:LEU:N	2.30	0.47
1:A:51:ARG:HE	1:A:52:HIS:HD2	1.60	0.47
1:B:51:ARG:HG3	1:B:52:HIS:H	1.79	0.47
1:D:89:HIS:CD2	1:D:90:PHE:CE1	3.02	0.47
1:D:49:ASP:OD1	1:D:51:ARG:HB2	2.15	0.47
1:B:84:VAL:O	1:B:87:LYS:HD2	2.14	0.47
1:A:104:GLU:OE2	1:A:174:ARG:NH2	2.42	0.47
1:D:50:ARG:H	1:D:50:ARG:HD3	1.77	0.47
1:B:225:THR:CG2	3:B:1015:HOH:O	2.63	0.46
1:D:192:GLY:O	1:D:193:HIS:CB	2.62	0.46
1:A:133:PHE:CE1	1:A:193:HIS:CE1	3.02	0.46
1:C:62:ARG:HG3	1:C:62:ARG:HH11	1.81	0.46
1:B:104:GLU:HB3	1:B:135:ASP:CG	2.36	0.46
1:B:51:ARG:CD	1:B:52:HIS:N	2.77	0.46
1:C:95:PRO:HG3	1:C:107:LEU:HA	1.98	0.46
1:D:104:GLU:CD	1:D:174:ARG:HH22	2.19	0.46
1:D:50:ARG:CG	1:D:50:ARG:HH11	2.26	0.46
1:A:12:HIS:CD2	1:A:216:HIS:CD2	2.96	0.46
1:B:51:ARG:CG	1:B:52:HIS:N	2.79	0.46
1:D:170:GLN:CA	1:D:170:GLN:NE2	2.77	0.46
1:A:33:ILE:HG21	1:A:227:LEU:HB3	1.97	0.45
1:B:57:ASP:O	1:B:61:LEU:HB2	2.16	0.45
1:C:161:ASP:O	1:C:162:ALA:C	2.55	0.45
1:C:62:ARG:HG3	1:C:62:ARG:NH1	2.31	0.45
1:A:233:GLU:HG3	3:A:1036:HOH:O	2.15	0.45
1:A:236:ARG:HH11	1:A:236:ARG:HG2	1.82	0.45
1:A:47:ARG:HH11	1:A:51:ARG:CB	2.29	0.45
1:C:133:PHE:CZ	1:C:193:HIS:HD2	2.34	0.45
1:A:227:LEU:HD22	1:A:231:VAL:HG23	1.99	0.44
1:B:199:ASN:ND2	1:B:199:ASN:C	2.67	0.44
1:B:113:ARG:O	1:B:117:ARG:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:HIS:CD2	1:C:157:GLY:H	2.35	0.44
1:D:199:ASN:HD22	1:D:200:VAL:N	2.15	0.44
1:C:91:CYS:HB2	1:C:128:ILE:HG21	1.98	0.44
1:B:51:ARG:CG	1:B:52:HIS:H	2.31	0.44
1:C:227:LEU:HD22	1:C:231:VAL:HG23	1.99	0.44
1:C:141:ILE:HD11	1:C:178:ALA:HB1	1.99	0.44
1:D:18:ASN:ND2	1:D:51:ARG:O	2.49	0.44
1:B:147:VAL:O	1:B:147:VAL:HG12	2.16	0.43
1:D:80:LEU:HD13	1:D:122:ARG:NH1	2.34	0.43
1:D:9:ASN:HD22	2:D:1004:G3P:H2	1.83	0.43
1:B:104:GLU:HB3	1:B:135:ASP:OD1	2.18	0.43
1:B:15:THR:CB	1:B:51:ARG:NH2	2.80	0.43
1:B:6:LEU:HG	1:B:238:MET:SD	2.58	0.43
1:D:170:GLN:HE21	1:D:170:GLN:HA	1.80	0.43
1:D:46:LEU:HD13	1:D:79:MET:HE1	2.01	0.42
1:D:51:ARG:HG2	1:D:51:ARG:NH1	2.23	0.42
1:C:163:LYS:HD2	1:C:167:GLU:OE2	2.19	0.42
1:A:2:ALA:HB1	3:A:1068:HOH:O	2.20	0.42
1:A:51:ARG:NE	1:A:52:HIS:HD2	2.17	0.42
1:A:89:HIS:HD2	1:A:90:PHE:CE2	2.37	0.42
1:A:94:VAL:HG12	1:A:133:PHE:CB	2.49	0.42
1:B:147:VAL:CG1	1:B:147:VAL:O	2.67	0.42
1:A:233:GLU:OE2	1:A:236:ARG:NH2	2.34	0.42
1:B:46:LEU:HD23	1:B:58:VAL:HG21	2.02	0.42
1:B:47:ARG:NE	1:B:51:ARG:HG2	2.34	0.42
1:B:91:CYS:HB2	1:B:128:ILE:CG2	2.50	0.42
1:B:15:THR:HA	1:B:51:ARG:HH21	1.76	0.42
1:B:77:GLU:CD	1:B:122:ARG:HH12	2.23	0.42
1:D:74:ALA:CB	1:D:96:GLU:HB2	2.48	0.42
1:C:15:THR:HG22	1:C:52:HIS:CB	2.49	0.42
1:A:91:CYS:HB2	1:A:128:ILE:CG2	2.50	0.41
1:A:192:GLY:O	1:A:193:HIS:CB	2.65	0.41
1:A:14:ALA:CB	1:A:52:HIS:HB2	2.46	0.41
1:B:162:ALA:HA	1:B:167:GLU:OE1	2.20	0.41
1:C:193:HIS:HA	2:C:1003:G3P:O2P	2.20	0.41
1:D:147:VAL:HG12	1:D:147:VAL:O	2.20	0.41
1:B:192:GLY:O	1:B:193:HIS:CB	2.68	0.41
1:C:57:ASP:O	1:C:61:LEU:HB2	2.21	0.41
1:D:117:ARG:NH2	1:D:146:GLU:OE2	2.52	0.41
1:D:52:HIS:CD2	1:D:53:ILE:N	2.76	0.41
1:D:80:LEU:O	1:D:84:VAL:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:ASP:O	1:C:137:ASP:N	2.53	0.41
1:D:46:LEU:HG	1:D:58:VAL:HG21	2.03	0.41
1:D:12:HIS:HE1	3:D:1078:HOH:O	2.02	0.41
1:B:86:THR:HB	1:B:88:PRO:HD3	2.04	0.40
1:C:171:GLU:OE2	1:C:171:GLU:HA	2.21	0.40
1:D:51:ARG:CG	1:D:51:ARG:NH1	2.85	0.40
1:B:191:ALA:O	1:B:212:LEU:HA	2.22	0.40
1:B:113:ARG:NH1	1:B:146:GLU:OE1	2.52	0.40
1:C:206:ILE:HA	1:C:207:PRO:HD3	1.78	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:ARG:NH1	2:C:1005:G3P:O2P[4_555]	2.19	0.01

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	231/242 (96%)	223 (96%)	7 (3%)	1 (0%)	34	42
1	B	231/242 (96%)	221 (96%)	9 (4%)	1 (0%)	34	42
1	C	231/242 (96%)	221 (96%)	9 (4%)	1 (0%)	34	42
1	D	231/242 (96%)	222 (96%)	8 (4%)	1 (0%)	34	42
All	All	924/968 (96%)	887 (96%)	33 (4%)	4 (0%)	34	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	193	HIS

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Mol	Chain	Res	Type
1	A	193	HIS
1	D	193	HIS
1	B	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	177/185 (96%)	161 (91%)	16 (9%)	9	11
1	B	178/185 (96%)	160 (90%)	18 (10%)	7	9
1	C	176/185 (95%)	156 (89%)	20 (11%)	5	6
1	D	177/185 (96%)	161 (91%)	16 (9%)	9	11
All	All	708/740 (96%)	638 (90%)	70 (10%)	8	9

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	6	LEU
1	A	12	HIS
1	A	20	ARG
1	A	46	LEU
1	A	51	ARG
1	A	61	LEU
1	A	68	ARG
1	A	71	LEU
1	A	91	CYS
1	A	165	ASP
1	A	172	LEU
1	A	188	LYS
1	A	189	VAL
1	A	199	ASN
1	A	227	LEU
1	B	3	GLU
1	B	4	LEU

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Mol	Chain	Res	Type
1	B	6	LEU
1	B	15	THR
1	B	46	LEU
1	B	50	ARG
1	B	51	ARG
1	B	56	ARG
1	B	61	LEU
1	B	68	ARG
1	B	71	LEU
1	B	85	GLU
1	B	91	CYS
1	B	94	VAL
1	B	172	LEU
1	B	189	VAL
1	B	199	ASN
1	B	227	LEU
1	C	4	LEU
1	C	6	LEU
1	C	15	THR
1	C	16	LEU
1	C	20	ARG
1	C	46	LEU
1	C	61	LEU
1	C	68	ARG
1	C	71	LEU
1	C	85	GLU
1	C	91	CYS
1	C	163	LYS
1	C	172	LEU
1	C	185	LEU
1	C	189	VAL
1	C	193	HIS
1	C	199	ASN
1	C	221	ARG
1	C	225	THR
1	C	227	LEU
1	D	3	GLU
1	D	4	LEU
1	D	20	ARG
1	D	46	LEU
1	D	48	GLU
1	D	50	ARG

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Mol	Chain	Res	Type
1	D	52	HIS
1	D	56	ARG
1	D	61	LEU
1	D	71	LEU
1	D	91	CYS
1	D	170	GLN
1	D	172	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	45	HIS
1	A	63	GLN
1	A	89	HIS
1	A	155	HIS
1	A	199	ASN
1	A	216	HIS
1	B	45	HIS
1	B	63	GLN
1	B	89	HIS
1	B	155	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	193	HIS
1	C	199	ASN
1	D	12	HIS
1	D	52	HIS
1	D	63	GLN
1	D	89	HIS
1	D	155	HIS
1	D	170	GLN
1	D	199	ASN
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 ⓘ

5 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$
2	G3P	A	1001	-	9,9,9	2.49	3 (33%)	11,12,12	1.62	2 (18%)
2	G3P	B	1002	-	9,9,9	2.46	3 (33%)	11,12,12	1.88	4 (36%)
2	G3P	C	1005	2	9,9,9	2.52	3 (33%)	11,12,12	2.04	3 (27%)
2	G3P	C	1003	1,2	9,9,9	2.46	3 (33%)	11,12,12	1.77	3 (27%)
2	G3P	D	1004	-	9,9,9	2.49	3 (33%)	11,12,12	1.81	3 (27%)

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	G3P	A	1001	-	-	5/8/8/8	-
2	G3P	B	1002	-	-	4/8/8/8	-
2	G3P	C	1005	2	-	3/8/8/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	G3P	C	1003	1,2	-	3/8/8/8	-
2	G3P	D	1004	-	-	5/8/8/8	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1002	G3P	P-O4P	-5.39	1.34	1.54
2	D	1004	G3P	P-O4P	-5.32	1.34	1.54
2	C	1005	G3P	P-O4P	-5.29	1.34	1.54
2	A	1001	G3P	P-O4P	-4.88	1.36	1.54
2	C	1003	G3P	P-O4P	-4.60	1.37	1.54
2	C	1003	G3P	O1-C1	-4.55	1.23	1.42
2	A	1001	G3P	O1-C1	-4.34	1.24	1.42
2	D	1004	G3P	O1-C1	-4.19	1.24	1.42
2	C	1005	G3P	O1-C1	-3.99	1.25	1.42
2	B	1002	G3P	O1-C1	-3.98	1.25	1.42
2	A	1001	G3P	P-O1P	-3.27	1.49	1.60
2	C	1005	G3P	P-O1P	-3.22	1.49	1.60
2	C	1003	G3P	P-O1P	-3.20	1.49	1.60
2	D	1004	G3P	P-O1P	-2.79	1.51	1.60
2	B	1002	G3P	P-O1P	-2.63	1.51	1.60

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1005	G3P	O1P-P-O2P	4.77	119.84	106.47
2	D	1004	G3P	O1P-P-O2P	4.07	117.89	106.47
2	A	1001	G3P	O1P-P-O2P	3.76	117.02	106.47
2	B	1002	G3P	O4P-P-O1P	-3.57	97.22	106.73
2	C	1003	G3P	P-O1P-C3	-3.54	108.56	118.30
2	C	1003	G3P	O1P-P-O2P	3.45	116.14	106.47
2	B	1002	G3P	O1P-P-O2P	3.03	114.97	106.47
2	A	1001	G3P	O4P-P-O1P	-2.65	99.67	106.73
2	C	1005	G3P	O1-C1-C2	2.65	122.92	110.20
2	B	1002	G3P	O1-C1-C2	2.64	122.86	110.20
2	D	1004	G3P	O1-C1-C2	2.60	122.67	110.20
2	B	1002	G3P	O3P-P-O1P	2.58	113.60	106.73
2	D	1004	G3P	O4P-P-O1P	-2.58	99.87	106.73
2	C	1003	G3P	O4P-P-O1P	-2.56	99.92	106.73
2	C	1005	G3P	O4P-P-O1P	-2.49	100.10	106.73

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	G3P	O1-C1-C2-C3
2	A	1001	G3P	C3-O1P-P-O4P
2	A	1001	G3P	C3-O1P-P-O2P
2	A	1001	G3P	C3-O1P-P-O3P
2	B	1002	G3P	C3-O1P-P-O4P
2	B	1002	G3P	C3-O1P-P-O2P
2	B	1002	G3P	C3-O1P-P-O3P
2	C	1005	G3P	C3-O1P-P-O4P
2	C	1005	G3P	C3-O1P-P-O2P
2	C	1005	G3P	C3-O1P-P-O3P
2	C	1003	G3P	C3-O1P-P-O4P
2	C	1003	G3P	C3-O1P-P-O2P
2	C	1003	G3P	C3-O1P-P-O3P
2	D	1004	G3P	C3-O1P-P-O4P
2	D	1004	G3P	C3-O1P-P-O2P
2	D	1004	G3P	C3-O1P-P-O3P
2	A	1001	G3P	O1-C1-C2-O2
2	B	1002	G3P	O1-C1-C2-C3
2	D	1004	G3P	O1-C1-C2-C3
2	D	1004	G3P	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1002	G3P	3	0
2	C	1005	G3P	2	1
2	C	1003	G3P	4	0
2	D	1004	G3P	1	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, 95th 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 ²)	Q<0.9
1	A	235/242 (97%)	0.01	10 (4%) 35 42	28, 41, 60, 72	0
1	B	235/242 (97%)	0.22	15 (6%) 19 25	29, 45, 66, 84	0
1	C	235/242 (97%)	0.06	11 (4%) 31 38	29, 41, 60, 68	0
1	D	235/242 (97%)	0.10	9 (3%) 40 47	28, 41, 63, 74	0
All	All	940/968 (97%)	0.10	45 (4%) 30 37	28, 42, 63, 84	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	ALA	7.0
1	B	104	GLU	5.7
1	B	51	ARG	5.0
1	D	104	GLU	4.6
1	D	51	ARG	4.3
1	B	105	GLY	4.2
1	B	163	LYS	4.2
1	D	2	ALA	4.2
1	A	104	GLU	4.0
1	C	104	GLU	3.7
1	A	243	GLY	3.5
1	A	96	GLU	3.5
1	C	48	GLU	3.5
1	A	163	LYS	3.4
1	B	162	ALA	3.4
1	B	114	ASP	3.3
1	D	163	LYS	3.2
1	D	243	GLY	3.2
1	C	163	LYS	3.0
1	C	105	GLY	3.0
1	B	121	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	164	THR	2.8
1	C	243	GLY	2.8
1	A	2	ALA	2.7
1	A	51	ARG	2.6
1	B	193	HIS	2.5
1	A	48	GLU	2.5
1	A	71	LEU	2.5
1	C	164	THR	2.4
1	D	125	ASP	2.4
1	B	164	THR	2.4
1	D	164	THR	2.4
1	B	20	ARG	2.3
1	C	3	GLU	2.2
1	C	2	ALA	2.2
1	D	22	THR	2.2
1	C	132	LEU	2.2
1	D	114	ASP	2.2
1	C	51	ARG	2.1
1	C	71	LEU	2.1
1	B	96	GLU	2.1
1	B	21	GLY	2.1
1	A	105	GLY	2.0
1	B	71	LEU	2.0
1	B	243	GLY	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, 95th 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(\AA^2)	Q<0.9
2	G3P	C	1005	10/10	0.75	0.44	87,88,91,92	0
2	G3P	C	1003	10/10	0.88	0.21	51,52,55,57	0
2	G3P	D	1004	10/10	0.93	0.18	69,71,73,75	0
2	G3P	B	1002	10/10	0.96	0.12	51,55,61,63	0
2	G3P	A	1001	10/10	0.96	0.16	49,52,57,58	0

6.5 Other polymers ⓘ

There are no such residues in this entry.