



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:56 PM UTC

PDB ID : 7NHF / pdb_00007nhf
Title : Crystal structure of Arabidopsis thaliana Pdx1K166R
Authors : Rodrigues, M.J.; Zhang, Y.; Bolton, R.; Evans, G.; Giri, N.; Royant, A.; Begley, T.; Ealick, S.E.; Tews, I.
Deposited on : 2021-02-10
Resolution : 2.35 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

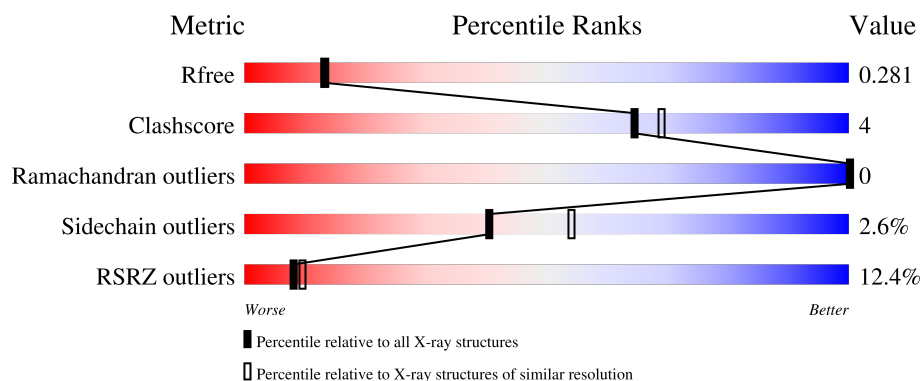
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	180053	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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 of 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	291	10% 84% 8% 8%
1	B	291	13% 84% 8% 7%
1	C	291	13% 80% 10% 9%
1	D	291	10% 82% 9% 8%

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	D	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7975 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 Pyridoxal 5'-phosphate synthase subunit PDX1.3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	2	0
			1998	1247	367	367	17			
1	B	270	Total	C	N	O	S	0	1	0
			1981	1239	362	364	16			
1	C	266	Total	C	N	O	S	0	1	0
			1965	1229	358	362	16			
1	D	267	Total	C	N	O	S	0	1	0
			1973	1233	361	362	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	166	ARG	LYS	engineered mutation	UNP Q8L940
B	166	ARG	LYS	engineered mutation	UNP Q8L940
C	166	ARG	LYS	engineered mutation	UNP Q8L940
D	166	ARG	LYS	engineered mutation	UNP Q8L940

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

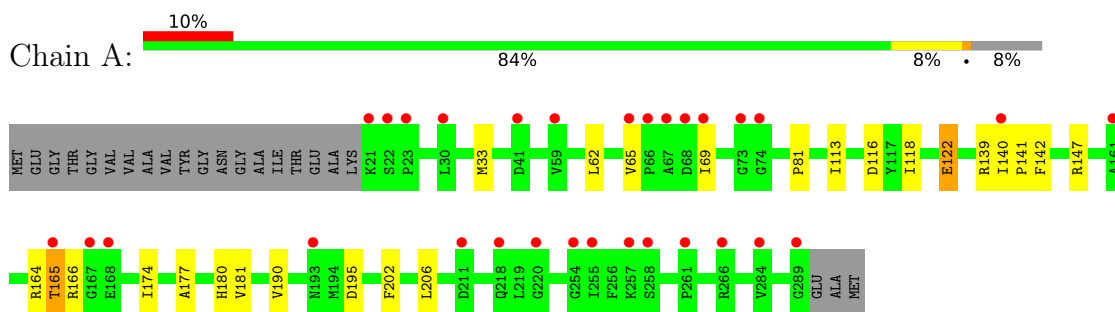
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	8	Total	O	0	0
			8	8		
3	C	9	Total	O	0	0
			9	9		
3	D	6	Total	O	0	0
			6	6		

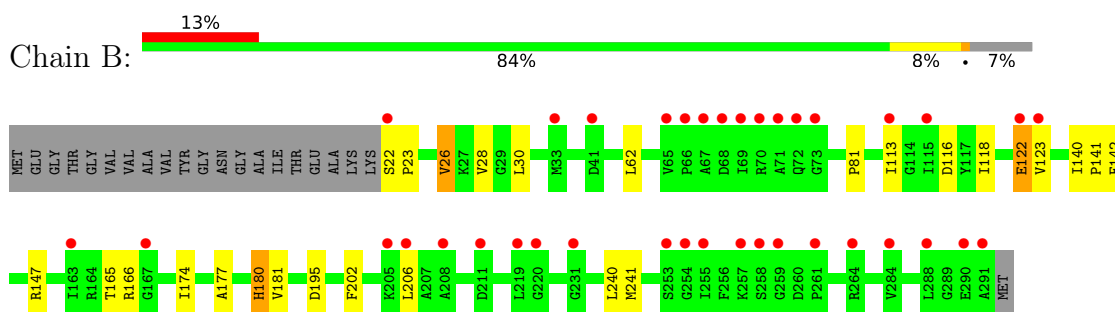
3 Residue-property plots [i](#)

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

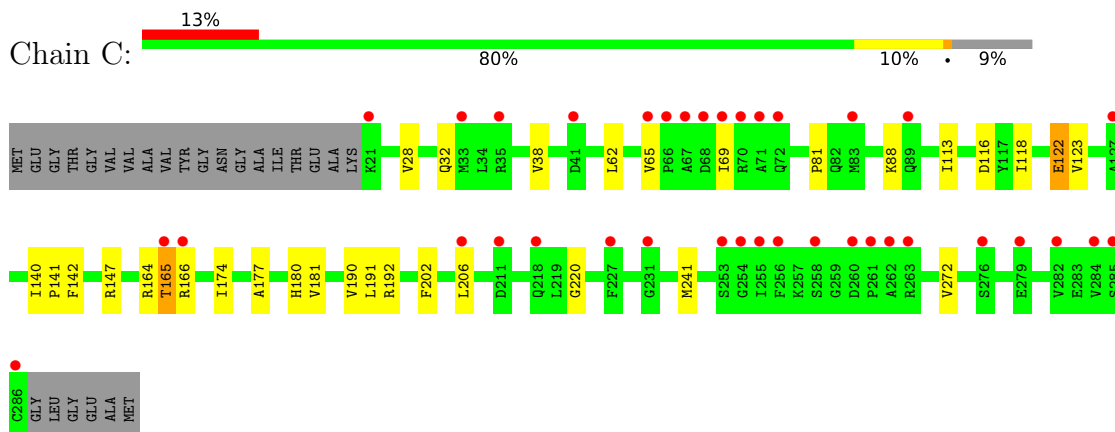
- Molecule 1: Pyridoxal 5'-phosphate synthase subunit PDX1.3



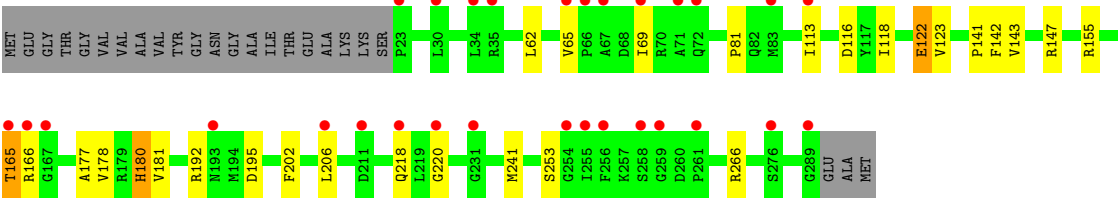
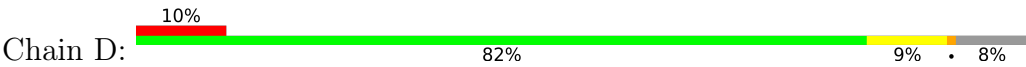
- Molecule 1: Pyridoxal 5'-phosphate synthase subunit PDX1.3



- Molecule 1: Pyridoxal 5'-phosphate synthase subunit PDX1.3



- Molecule 1: Pyridoxal 5'-phosphate synthase subunit PDX1.3



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	178.00Å 178.00Å 115.13Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.38 – 2.35 38.38 – 2.35	Depositor EDS
% Data completeness (in resolution range)	98.6 (38.38-2.35) 98.6 (38.38-2.35)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.220 , 0.249 0.258 , 0.281	Depositor DCC
R_{free} test set	2796 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.806	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 20.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for $-1/3^*h+1/3^*k+4/3^*l, -k, 2/3^*h+1/3^*k+1/3^*l$ 0.000 for $-2/3^*h-1/3^*k-4/3^*l, -1/3^*h-2/3^*k+4/3^*l, -1/3^*h+1/3^*k+1/3^*l$ 0.000 for $-h, 1/3^*h-1/3^*k-4/3^*l, -1/3^*h-2/3^*k+1/3^*l$ 0.000 for $-1/3^*h-2/3^*k+4/3^*l, -2/3^*h-1/3^*k-4/3^*l, 1/3^*h-1/3^*k-1/3^*l$ 0.000 for $-h, 2/3^*h+1/3^*k+4/3^*l, 1/3^*h+2/3^*k-1/3^*l$ 0.000 for $1/3^*h+2/3^*k-4/3^*l, -k, -2/3^*h-1/3^*k-1/3^*l$ 0.025 for h, -h-k, -l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7975	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% 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

5.1 Standard geometry

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

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	1.08	0/2025	1.49	3/2738 (0.1%)
1	B	1.10	1/2008 (0.0%)	1.50	4/2716 (0.1%)
1	C	1.08	0/1992	1.49	4/2696 (0.1%)
1	D	1.08	1/2000 (0.1%)	1.50	2/2705 (0.1%)
All	All	1.08	2/8025 (0.0%)	1.49	13/10855 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	180	HIS	CE1-NE2	5.17	1.37	1.32
1	B	180	HIS	CE1-NE2	5.04	1.37	1.32

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	165	THR	CA-CB-OG1	-5.90	100.75	109.60
1	A	165	THR	CA-CB-OG1	-5.89	100.77	109.60
1	B	165	THR	CA-CB-OG1	-5.58	101.24	109.60
1	D	165	THR	CA-CB-OG1	-5.48	101.38	109.60
1	A	164	ARG	CG-CD-NE	-5.16	100.64	112.00
1	B	28	VAL	CA-C-N	5.12	125.77	120.03
1	B	28	VAL	C-N-CA	5.12	125.77	120.03
1	A	195	ASP	CA-CB-CG	5.11	117.71	112.60
1	C	28	VAL	CA-C-N	5.07	125.71	120.03
1	C	28	VAL	C-N-CA	5.07	125.71	120.03
1	C	164	ARG	CG-CD-NE	-5.07	100.85	112.00
1	B	195	ASP	CA-CB-CG	5.03	117.63	112.60
1	D	195	ASP	CA-CB-CG	5.01	117.61	112.60

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	1998	0	1995	12	0
1	B	1981	0	1971	16	0
1	C	1965	0	1952	18	0
1	D	1973	0	1971	16	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	2	0
3	A	15	0	0	0	0
3	B	8	0	0	0	0
3	C	9	0	0	1	0
3	D	6	0	0	0	0
All	All	7975	0	7889	61	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 4.

All (61) 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:A:81:PRO:HB3	1:A:113:ILE:HD11	1.55	0.87
1:C:174:ILE:HD11	1:C:241:MET:HE3	1.68	0.76
1:A:81:PRO:CB	1:A:113:ILE:HD11	2.26	0.66
1:C:81:PRO:HB3	1:C:113:ILE:HD11	1.76	0.65
1:B:22:SER:CB	1:B:23:PRO:CD	2.76	0.63
1:C:202:PHE:CE2	1:C:206:LEU:HD11	2.35	0.61
1:C:174:ILE:CD1	1:C:241:MET:HE3	2.32	0.58
1:B:81:PRO:HB3	1:B:113:ILE:HD11	1.85	0.58
1:D:177:ALA:O	1:D:181:VAL:HG23	2.06	0.55
1:D:155:ARG:NH2	2:D:301:PO4:O2	2.38	0.54
1:D:81:PRO:HB3	1:D:113:ILE:HD11	1.89	0.54
1:B:177:ALA:O	1:B:181:VAL:HG23	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:ALA:O	1:A:181:VAL:HG23	2.09	0.52
1:C:177:ALA:O	1:C:181:VAL:HG23	2.10	0.51
1:D:122:GLU:HG3	1:D:166:ARG:HG3	1.93	0.51
1:C:122:GLU:HG3	1:C:166:ARG:HG3	1.93	0.51
1:D:155:ARG:NH1	2:D:301:PO4:O2	2.45	0.50
1:A:122:GLU:HG3	1:A:166:ARG:HG3	1.94	0.50
1:B:174:ILE:CG1	1:B:241:MET:HE3	2.42	0.50
1:B:174:ILE:HD11	1:B:241:MET:HE2	1.93	0.49
1:B:122:GLU:HG3	1:B:166:ARG:HG3	1.95	0.48
1:B:174:ILE:HD11	1:B:241:MET:CE	2.43	0.48
1:B:118:ILE:O	1:B:142:PHE:HA	2.15	0.47
1:B:174:ILE:HD13	1:B:240:LEU:HD23	1.96	0.47
1:D:65:VAL:HG11	1:D:253:SER:OG	2.15	0.47
1:C:165:THR:O	1:C:165:THR:OG1	2.34	0.46
1:D:178:VAL:HG22	1:D:241:MET:HE1	1.96	0.46
1:B:147:ARG:HG2	1:B:180:HIS:CE1	2.50	0.46
1:C:147:ARG:HG2	1:C:180:HIS:CE1	2.49	0.46
1:D:116:ASP:O	1:D:141:PRO:HD2	2.15	0.46
1:A:118:ILE:HG13	1:A:140:ILE:HD11	1.98	0.46
1:A:165:THR:O	1:A:165:THR:OG1	2.31	0.46
1:C:65:VAL:O	1:C:69:ILE:HG13	2.16	0.46
1:C:116:ASP:O	1:C:141:PRO:HD2	2.15	0.46
1:A:118:ILE:O	1:A:142:PHE:HA	2.16	0.46
1:D:147:ARG:HG2	1:D:180:HIS:CE1	2.50	0.46
1:A:65:VAL:O	1:A:69:ILE:HG13	2.15	0.45
1:C:32:GLN:NE2	3:C:401:HOH:O	2.49	0.45
1:A:116:ASP:O	1:A:141:PRO:HD2	2.17	0.45
1:D:202:PHE:CE2	1:D:206:LEU:HD11	2.52	0.45
1:A:202:PHE:CE2	1:A:206:LEU:HD11	2.52	0.45
1:A:147:ARG:HG2	1:A:180:HIS:CE1	2.51	0.45
1:D:118:ILE:O	1:D:142:PHE:HA	2.17	0.45
1:C:174:ILE:HG13	1:C:241:MET:CE	2.46	0.44
1:D:65:VAL:O	1:D:69:ILE:HG13	2.17	0.44
1:B:116:ASP:O	1:B:141:PRO:HD2	2.16	0.44
1:C:118:ILE:HG13	1:C:140:ILE:HD11	2.00	0.44
1:D:192:ARG:HD2	1:D:220:GLY:CA	2.47	0.44
1:B:202:PHE:CE2	1:B:206:LEU:HD11	2.52	0.44
1:C:118:ILE:O	1:C:142:PHE:HA	2.17	0.44
1:D:165:THR:O	1:D:165:THR:OG1	2.33	0.43
1:C:81:PRO:CB	1:C:113:ILE:HD11	2.45	0.42
1:C:192:ARG:HD3	1:C:220:GLY:CA	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:SER:O	1:B:26:VAL:HG13	2.19	0.42
1:B:118:ILE:HG13	1:B:140:ILE:HD11	2.01	0.41
1:B:22:SER:CB	1:B:23:PRO:HD3	2.49	0.41
1:B:81:PRO:CB	1:B:113:ILE:HD11	2.50	0.41
1:A:190:VAL:HG13	1:C:190:VAL:HG22	2.02	0.41
1:D:65:VAL:HG11	1:D:253:SER:CB	2.51	0.41
1:D:266:ARG:HA	1:D:266:ARG:HD2	1.82	0.41
1:C:38:VAL:HG22	1:C:272:VAL:HG21	2.03	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	269/291 (92%)	261 (97%)	8 (3%)	0	100	100
1	B	269/291 (92%)	259 (96%)	10 (4%)	0	100	100
1	C	265/291 (91%)	255 (96%)	10 (4%)	0	100	100
1	D	266/291 (91%)	259 (97%)	7 (3%)	0	100	100
All	All	1069/1164 (92%)	1034 (97%)	35 (3%)	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	200/223 (90%)	195 (98%)	5 (2%)	42	55
1	B	195/223 (87%)	190 (97%)	5 (3%)	40	54
1	C	195/223 (87%)	190 (97%)	5 (3%)	40	54
1	D	197/223 (88%)	192 (98%)	5 (2%)	42	55
All	All	787/892 (88%)	767 (98%)	20 (2%)	40	55

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	MET
1	A	62	LEU
1	A	122	GLU
1	A	139	ARG
1	A	174	ILE
1	B	26	VAL
1	B	30	LEU
1	B	62	LEU
1	B	122	GLU
1	B	123	VAL
1	C	62	LEU
1	C	88	LYS
1	C	122	GLU
1	C	123	VAL
1	C	191	LEU
1	D	62	LEU
1	D	122	GLU
1	D	123	VAL
1	D	143	VAL
1	D	218	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	132	HIS
1	A	148	ASN
1	A	180	HIS
1	A	243	GLN
1	B	108	GLN
1	B	132	HIS
1	B	148	ASN

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Mol	Chain	Res	Type
1	B	180	HIS
1	C	132	HIS
1	C	148	ASN
1	C	180	HIS
1	C	243	GLN
1	D	108	GLN
1	D	132	HIS
1	D	148	ASN
1	D	180	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	PO4	C	301	-	4,4,4	0.95	0	6,6,6	0.50	0
2	PO4	B	301	-	4,4,4	1.36	1 (25%)	6,6,6	0.29	0
2	PO4	A	301	-	4,4,4	0.97	0	6,6,6	0.39	0
2	PO4	D	301	-	4,4,4	0.66	0	6,6,6	0.58	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	PO4	P-O1	2.52	1.56	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	PO4	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	269/291 (92%)	1.03	30 (11%)	10 12	21, 50, 93, 123	2 (0%)
1	B	270/291 (92%)	1.13	37 (13%)	6 7	25, 52, 95, 117	1 (0%)
1	C	266/291 (91%)	1.18	37 (13%)	6 7	37, 56, 100, 127	1 (0%)
1	D	267/291 (91%)	1.04	29 (10%)	10 12	27, 55, 99, 136	1 (0%)
All	All	1072/1164 (92%)	1.09	133 (12%)	8 9	21, 54, 99, 136	5 (0%)

All (133) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	66	PRO	6.8
1	A	67	ALA	5.5
1	B	65	VAL	5.2
1	C	67	ALA	5.1
1	B	66	PRO	5.0
1	C	260	ASP	4.9
1	D	23	PRO	4.7
1	C	284	VAL	4.6
1	B	205[A]	LYS	4.3
1	B	22	SER	4.3
1	B	291	ALA	4.2
1	D	289	GLY	4.1
1	D	66	PRO	4.1
1	A	21	LYS	4.0
1	C	258	SER	3.9
1	D	67	ALA	3.9
1	A	66	PRO	3.8
1	B	254	GLY	3.8
1	C	68	ASP	3.8
1	D	218	GLN	3.8
1	B	70	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	254	GLY	3.6
1	B	290	GLU	3.6
1	A	254	GLY	3.6
1	C	65	VAL	3.6
1	C	254	GLY	3.5
1	B	253	SER	3.5
1	C	33	MET	3.4
1	A	289	GLY	3.4
1	B	67	ALA	3.4
1	B	167	GLY	3.4
1	A	261	PRO	3.3
1	D	167	GLY	3.3
1	D	35[A]	ARG	3.3
1	C	21	LYS	3.2
1	B	211	ASP	3.2
1	C	261	PRO	3.2
1	B	73	GLY	3.1
1	C	218	GLN	3.1
1	C	72	GLN	3.0
1	A	74	GLY	3.0
1	B	257	LYS	3.0
1	A	68	ASP	3.0
1	B	259	GLY	2.9
1	C	89[A]	GLN	2.9
1	C	70	ARG	2.8
1	C	256	PHE	2.8
1	C	165	THR	2.8
1	C	255	ILE	2.8
1	A	284	VAL	2.8
1	B	69	ILE	2.8
1	B	68	ASP	2.7
1	B	255	ILE	2.7
1	A	167	GLY	2.7
1	A	258	SER	2.7
1	A	165	THR	2.7
1	B	258	SER	2.7
1	B	261	PRO	2.7
1	A	257	LYS	2.6
1	D	259	GLY	2.6
1	D	71	ALA	2.6
1	D	193	ASN	2.6
1	A	22	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	255	ILE	2.6
1	D	72	GLN	2.5
1	B	71	ALA	2.5
1	B	231	GLY	2.5
1	C	231	GLY	2.5
1	C	211	ASP	2.5
1	C	69	ILE	2.5
1	B	219	LEU	2.5
1	D	231	GLY	2.5
1	C	285	SER	2.4
1	D	113	ILE	2.4
1	C	41	ASP	2.4
1	A	140	ILE	2.4
1	B	33	MET	2.4
1	B	206	LEU	2.4
1	C	71	ALA	2.4
1	A	69	ILE	2.4
1	A	30	LEU	2.4
1	B	264	ARG	2.4
1	C	263	ARG	2.4
1	C	279	GLU	2.4
1	D	83	MET	2.3
1	D	65	VAL	2.3
1	C	286	CYS	2.3
1	A	65	VAL	2.3
1	B	288	LEU	2.3
1	C	253	SER	2.3
1	B	284	VAL	2.3
1	D	34	LEU	2.2
1	D	256	PHE	2.2
1	A	218[A]	GLN	2.2
1	B	220	GLY	2.2
1	D	276	SER	2.2
1	D	206	LEU	2.2
1	A	220	GLY	2.2
1	C	206	LEU	2.2
1	D	255	ILE	2.2
1	C	282	VAL	2.2
1	C	262	ALA	2.2
1	A	266	ARG	2.1
1	B	113	ILE	2.1
1	D	30	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	73	GLY	2.1
1	D	258	SER	2.1
1	D	166	ARG	2.1
1	B	163	ILE	2.1
1	C	83	MET	2.1
1	D	69	ILE	2.1
1	B	122	GLU	2.1
1	A	23	PRO	2.1
1	A	59	VAL	2.1
1	D	165	THR	2.1
1	A	41	ASP	2.1
1	B	72	GLN	2.1
1	D	220	GLY	2.1
1	C	35	ARG	2.1
1	A	161	ALA	2.1
1	C	127	ALA	2.1
1	A	193	ASN	2.1
1	D	211	ASP	2.0
1	B	115	ILE	2.0
1	D	261	PRO	2.0
1	B	123	VAL	2.0
1	C	166	ARG	2.0
1	C	227	PHE	2.0
1	A	168	GLU	2.0
1	C	276	SER	2.0
1	B	208	ALA	2.0
1	A	211	ASP	2.0
1	B	41	ASP	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 oligosaccharides 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(Å ²)	Q<0.9
2	PO4	A	301	5/5	0.78	0.15	75,85,97,98	0
2	PO4	C	301	5/5	0.79	0.14	81,86,95,105	0
2	PO4	D	301	5/5	0.82	0.33	69,73,76,84	0
2	PO4	B	301	5/5	0.83	0.13	73,77,90,91	0

6.5 Other polymers [i](#)

There are no such residues in this entry.