



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2021 – 10:08 AM JST

PDB ID : 7DOQ
Title : Lp major histidine acid phosphatase mutant D281A/5'-AMP
Authors : Guo, Y.; Teng, Y.
Deposited on : 2020-12-16
Resolution : 2.20 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

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 : 2.24
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) : 2.24

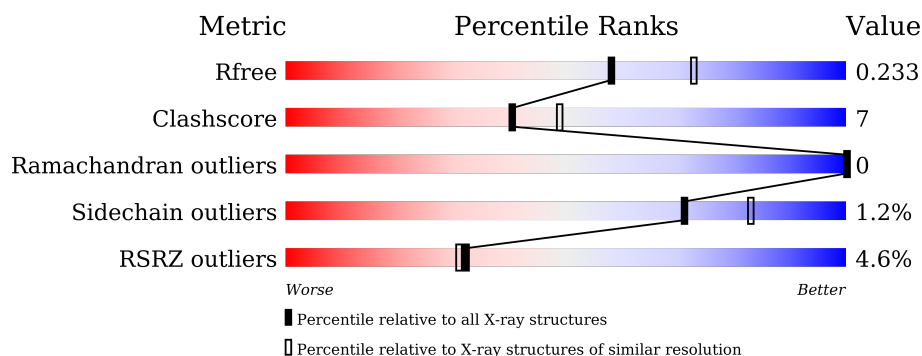
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	335	<div> <div>2%</div> <div>88% 12%</div> </div>
1	B	335	<div> <div>8%</div> <div>84% 13% ..</div> </div>
1	C	335	<div> <div>6%</div> <div>85% 14% .</div> </div>
1	D	335	<div> <div>2%</div> <div>86% 13% .</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	PO4	B	403	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10841 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.

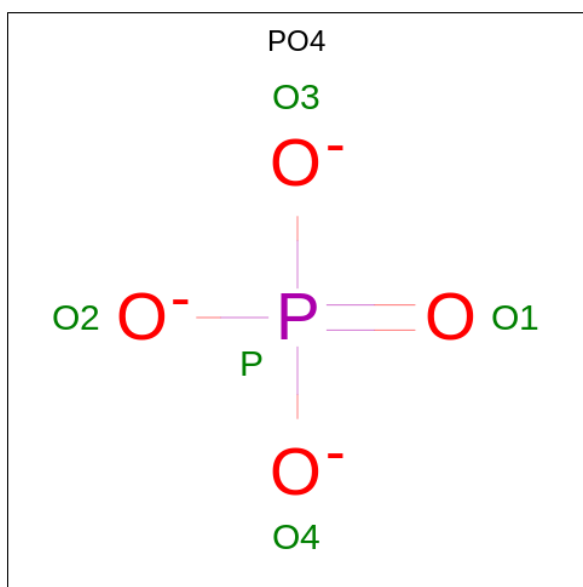
- Molecule 1 is a protein called Acid phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	335	Total	C	N	O	S	0	1	0
			2658	1695	444	507	12			
1	B	331	Total	C	N	O	S	0	2	0
			2629	1681	436	501	11			
1	C	331	Total	C	N	O	S	0	2	0
			2626	1676	436	503	11			
1	D	331	Total	C	N	O	S	0	3	0
			2634	1686	436	501	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	HIS	-	expression tag	UNP A0A2S6F805
A	20	HIS	-	expression tag	UNP A0A2S6F805
A	21	MET	-	expression tag	UNP A0A2S6F805
B	19	HIS	-	expression tag	UNP A0A2S6F805
B	20	HIS	-	expression tag	UNP A0A2S6F805
B	21	MET	-	expression tag	UNP A0A2S6F805
C	19	HIS	-	expression tag	UNP A0A2S6F805
C	20	HIS	-	expression tag	UNP A0A2S6F805
C	21	MET	-	expression tag	UNP A0A2S6F805
D	19	HIS	-	expression tag	UNP A0A2S6F805
D	20	HIS	-	expression tag	UNP A0A2S6F805
D	21	MET	-	expression tag	UNP A0A2S6F805

- Molecule 2 is PHOSPHATE ION (three-letter code: 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	B	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	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		

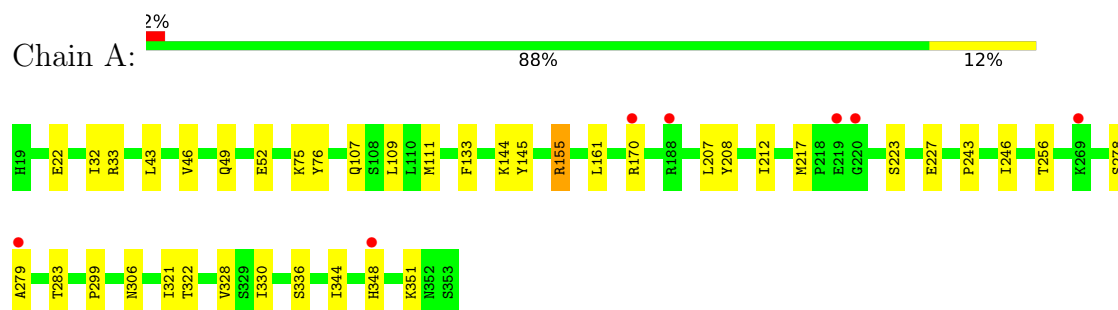
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	70	Total	O	0	0
			70	70		
3	B	58	Total	O	0	0
			58	58		
3	C	63	Total	O	0	0
			63	63		
3	D	68	Total	O	0	0
			68	68		

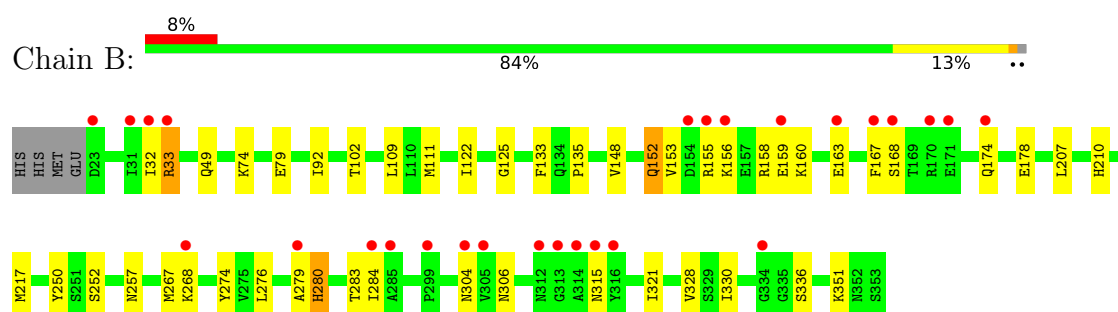
3 Residue-property plots [i](#)

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.

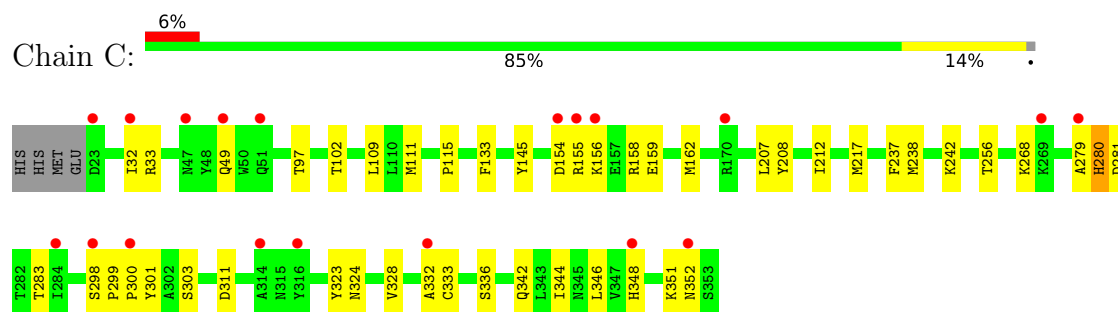
• Molecule 1: Acid phosphatase



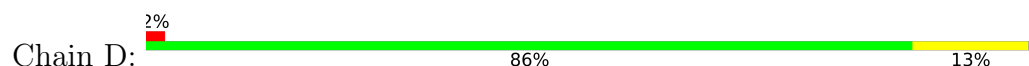
• Molecule 1: Acid phosphatase

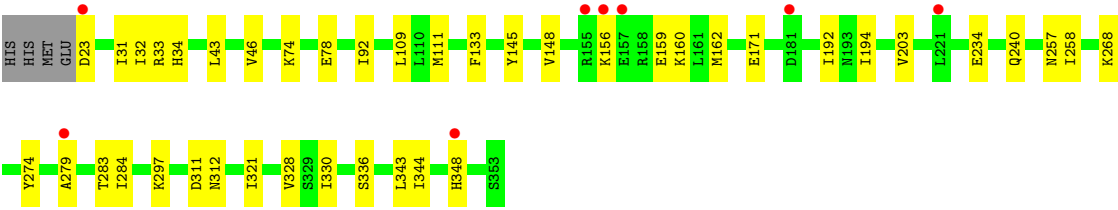


• Molecule 1: Acid phosphatase



• Molecule 1: Acid phosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.50Å 56.48Å 146.35Å 90.00° 110.01° 90.00°	Depositor
Resolution (Å)	35.62 – 2.20 35.59 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.4 (35.62-2.20) 96.4 (35.59-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.192 , 0.229 0.196 , 0.233	Depositor DCC
R_{free} test set	3517 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	32.8	Xtriage
Anisotropy	0.153	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10841	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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: 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	0.64	0/2727	0.68	0/3708
1	B	0.65	0/2700	0.70	0/3672
1	C	0.65	0/2696	0.69	0/3667
1	D	0.66	0/2708	0.69	0/3683
All	All	0.65	0/10831	0.69	0/14730

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	2658	0	2615	29	0
1	B	2629	0	2595	38	0
1	C	2626	0	2590	55	0
1	D	2634	0	2606	39	0
2	A	5	0	0	0	0
2	B	15	0	0	0	0
2	C	15	0	0	0	0
3	A	70	0	0	0	0
3	B	58	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	63	0	0	1	0
3	D	68	0	0	0	0
All	All	10841	0	10406	151	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 7.

All (151) 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:300:PRO:HG2	1:C:303:SER:OG	1.51	1.10
1:C:237:PHE:HD2	1:C:298:SER:HG	1.05	0.93
1:C:300:PRO:HD2	1:C:323:TYR:OH	1.72	0.88
1:D:162:MET:HE1	1:D:240:GLN:HG2	1.54	0.87
1:D:23:ASP:OD1	1:D:312:ASN:HA	1.77	0.85
1:B:315:ASN:OD1	1:C:256:THR:HG21	1.77	0.84
1:C:279:ALA:HB1	1:C:283:THR:CB	2.08	0.82
1:B:279:ALA:HB1	1:B:283:THR:OG1	1.80	0.82
1:C:237:PHE:HD2	1:C:298:SER:OG	1.65	0.79
1:C:300:PRO:CG	1:C:303:SER:OG	2.31	0.77
1:C:300:PRO:HG2	1:C:303:SER:HG	1.50	0.76
1:B:152:GLN:HA	1:B:152:GLN:OE1	1.88	0.74
1:C:279:ALA:CB	1:C:283:THR:HG21	2.19	0.73
1:B:207:LEU:HD22	1:B:217:MET:HE1	1.70	0.73
1:D:279:ALA:HB1	1:D:283:THR:CB	2.18	0.73
1:D:234:GLU:OE2	1:D:297:LYS:HD2	1.90	0.71
1:A:144:LYS:HE3	1:A:145:TYR:CZ	2.26	0.70
1:C:351:LYS:HE3	1:C:352:ASN:HD21	1.54	0.70
1:C:111:MET:HE1	1:D:133[A]:PHE:CD2	2.25	0.70
1:A:321:ILE:HD12	1:A:330:ILE:HD11	1.73	0.69
1:D:344:ILE:HG23	1:D:348:HIS:NE2	2.08	0.69
1:C:279:ALA:HB1	1:C:283:THR:OG1	1.92	0.68
1:D:279:ALA:CB	1:D:283:THR:HG21	2.24	0.67
1:C:279:ALA:HB1	1:C:283:THR:HG21	1.77	0.66
1:A:279:ALA:HB1	1:A:283:THR:CB	2.25	0.66
1:C:300:PRO:CD	1:C:323:TYR:OH	2.43	0.64
1:C:279:ALA:HB1	1:C:283:THR:CG2	2.28	0.63
1:B:207:LEU:HD22	1:B:217:MET:CE	2.28	0.62
1:B:279:ALA:HB1	1:B:283:THR:HG1	1.62	0.62
1:B:315:ASN:OD1	1:C:256:THR:CG2	2.48	0.62
1:B:49:GLN:HA	1:B:49:GLN:OE1	1.99	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:351:LYS:HE3	1:C:352:ASN:ND2	2.14	0.61
1:C:333:CYS:HA	1:C:342:GLN:NE2	2.16	0.61
1:B:279:ALA:HB1	1:B:283:THR:CB	2.30	0.61
1:B:304:ASN:HD21	1:B:306:ASN:HB2	1.65	0.61
1:C:155:ARG:HH12	1:C:158:ARG:HH11	1.46	0.60
1:A:133:PHE:CD2	1:B:111:MET:HE1	2.36	0.60
1:B:32[A]:ILE:HG22	1:B:33:ARG:O	2.00	0.60
1:D:23:ASP:OD1	1:D:311:ASP:O	2.20	0.60
1:C:300:PRO:HG2	1:C:324:ASN:HD21	1.66	0.60
1:C:155:ARG:HH11	1:C:158:ARG:HD3	1.67	0.60
1:C:281:ASP:HB2	1:C:299:PRO:HG2	1.83	0.60
1:B:315:ASN:CG	1:C:256:THR:HG21	2.22	0.59
1:D:92[A]:ILE:HG22	1:D:274:TYR:HB3	1.83	0.59
1:B:155:ARG:O	1:B:155:ARG:HG3	2.03	0.59
1:D:279:ALA:HB1	1:D:283:THR:HG21	1.83	0.59
1:B:328:VAL:O	1:B:336:SER:HA	2.04	0.58
1:B:279:ALA:CB	1:B:283:THR:HG21	2.34	0.58
1:C:155:ARG:NH1	1:C:158:ARG:HD3	2.19	0.57
1:A:111:MET:HE1	1:B:133[A]:PHE:CD2	2.39	0.57
1:B:267:MET:C	1:B:268:LYS:HG3	2.25	0.57
1:B:33:ARG:HB3	1:B:284:ILE:HD12	1.86	0.57
1:B:163:GLU:HA	1:B:167:PHE:HB2	1.87	0.55
1:A:279:ALA:HB1	1:A:283:THR:OG1	2.07	0.55
1:C:133:PHE:CD2	1:D:111:MET:HE1	2.42	0.55
1:B:32[A]:ILE:HG13	1:B:276:LEU:HD11	1.89	0.55
1:D:279:ALA:HB1	1:D:283:THR:CG2	2.37	0.54
1:C:154:ASP:OD2	1:C:156:LYS:HE2	2.08	0.53
1:D:156:LYS:O	1:D:160:LYS:HG2	2.08	0.53
1:A:207:LEU:HD22	1:A:217:MET:HE1	1.91	0.53
1:B:102:THR:HG21	1:B:280:HIS:CE1	2.44	0.53
1:D:171:GLU:H	1:D:171:GLU:CD	2.12	0.53
1:C:300:PRO:O	1:C:303:SER:CB	2.58	0.52
1:A:43:LEU:HB2	1:A:46:VAL:HG22	1.91	0.52
1:C:268:LYS:NZ	1:C:311:ASP:OD1	2.31	0.52
1:D:148:VAL:HA	1:D:257:ASN:ND2	2.25	0.51
1:C:281:ASP:O	1:C:299:PRO:HG3	2.11	0.51
1:B:174:GLN:O	1:B:178:GLU:HG2	2.10	0.51
1:D:268:LYS:NZ	1:D:311:ASP:OD2	2.43	0.51
1:C:300:PRO:O	1:C:303:SER:HB2	2.11	0.51
1:D:32[A]:ILE:HG22	1:D:33:ARG:O	2.11	0.50
1:C:344:ILE:O	1:C:348:HIS:ND1	2.37	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:344:ILE:CG2	1:D:348:HIS:NE2	2.75	0.50
1:D:321:ILE:HD12	1:D:330:ILE:HD11	1.93	0.49
1:A:207:LEU:HD22	1:A:217:MET:CE	2.42	0.49
1:C:300:PRO:CG	1:C:303:SER:HG	2.23	0.49
1:D:279:ALA:HB1	1:D:283:THR:HB	1.95	0.49
1:A:243:PRO:HB2	1:A:246:ILE:HG12	1.94	0.49
1:A:328:VAL:O	1:A:336:SER:HA	2.13	0.49
1:A:155:ARG:N	1:A:155:ARG:HD2	2.28	0.48
1:C:49:GLN:OE1	1:C:49:GLN:N	2.32	0.48
1:D:162:MET:HE1	1:D:240:GLN:CG	2.34	0.48
1:D:321:ILE:HD12	1:D:330:ILE:CD1	2.42	0.48
1:A:279:ALA:CB	1:A:283:THR:HG21	2.42	0.48
1:B:163:GLU:O	1:B:163:GLU:HG3	2.13	0.47
1:A:256:THR:OG1	1:A:351:LYS:HE2	2.14	0.47
1:A:33:ARG:NH1	1:A:299:PRO:O	2.48	0.47
1:B:321:ILE:HD12	1:B:330:ILE:CD1	2.45	0.47
1:C:32:ILE:HD13	1:C:109:LEU:HD13	1.96	0.47
1:C:279:ALA:CB	1:C:283:THR:CG2	2.91	0.47
1:A:208:TYR:CE2	1:A:212:ILE:HD11	2.50	0.47
1:D:74:LYS:HD2	1:D:78:GLU:HB2	1.97	0.47
1:C:279:ALA:HB2	1:C:283:THR:HG21	1.95	0.46
1:D:234:GLU:OE2	1:D:297:LYS:CD	2.62	0.46
1:B:74:LYS:HE3	1:B:79:GLU:OE1	2.16	0.46
1:B:155:ARG:HD2	1:B:159:GLU:OE1	2.16	0.46
1:C:207:LEU:HD22	1:C:217:MET:CE	2.46	0.46
1:C:238:MET:HA	1:C:238:MET:HE2	1.97	0.46
1:D:32[B]:ILE:HD13	1:D:109:LEU:HD13	1.98	0.46
1:C:155:ARG:CZ	1:C:159:GLU:OE2	2.64	0.46
1:A:52:GLU:OE1	1:A:52:GLU:N	2.46	0.45
1:C:328:VAL:O	1:C:336:SER:HA	2.17	0.45
1:C:344:ILE:HG23	1:C:348:HIS:CE1	2.52	0.45
1:B:148:VAL:HA	1:B:257:ASN:ND2	2.32	0.45
1:D:156:LYS:O	1:D:159:GLU:HG2	2.16	0.45
1:D:279:ALA:HB1	1:D:283:THR:OG1	2.16	0.45
1:C:111:MET:O	1:C:115:PRO:HA	2.17	0.45
1:D:43:LEU:HB2	1:D:46:VAL:HG22	1.99	0.45
1:C:207:LEU:HD22	1:C:217:MET:HE1	2.00	0.44
1:B:32[B]:ILE:HD13	1:B:109:LEU:HD13	1.99	0.44
1:C:97:THR:HG1	1:C:280:HIS:CE1	2.35	0.44
1:D:23:ASP:CG	1:D:312:ASN:HA	2.38	0.44
1:D:343:LEU:O	1:D:343:LEU:HD23	2.18	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ILE:HD12	1:A:330:ILE:CD1	2.45	0.44
1:B:217:MET:HE2	1:B:217:MET:HB3	1.86	0.44
1:C:102:THR:HG21	1:C:280:HIS:CE1	2.52	0.44
1:B:252:SER:OG	1:B:351:LYS:HA	2.17	0.43
1:D:33:ARG:HB2	1:D:284:ILE:HD12	2.01	0.43
1:B:92:ILE:HG22	1:B:274:TYR:HB3	2.01	0.43
1:C:333:CYS:HA	1:C:342:GLN:HE22	1.84	0.43
1:A:43:LEU:HD11	1:A:207:LEU:HD23	2.01	0.43
1:D:328:VAL:O	1:D:336:SER:HA	2.19	0.43
1:A:223:SER:O	1:A:227:GLU:HG2	2.19	0.43
1:C:208:TYR:CE2	1:C:212:ILE:HD11	2.53	0.43
1:C:217:MET:HE2	1:C:217:MET:HB3	1.79	0.42
1:B:163:GLU:O	1:B:168:SER:HB3	2.19	0.42
1:D:279:ALA:HB2	1:D:283:THR:HG21	2.00	0.42
1:A:306:ASN:HB3	1:A:322:THR:OG1	2.19	0.42
1:A:279:ALA:HB1	1:A:283:THR:HB	2.02	0.42
1:A:107:GLN:NE2	1:B:135:PRO:HA	2.35	0.42
1:A:161:LEU:HB3	1:A:246:ILE:HD13	2.02	0.42
1:A:75:LYS:HD2	1:A:76:TYR:CZ	2.55	0.42
1:C:145:TYR:CG	1:D:145:TYR:HB3	2.55	0.42
1:C:332:ALA:HB3	1:C:346:LEU:HD12	2.00	0.42
1:A:32:ILE:HD13	1:A:109:LEU:HD13	2.01	0.42
1:A:344:ILE:HG23	1:A:348:HIS:NE2	2.35	0.42
1:D:33:ARG:HG2	1:D:34:HIS:N	2.34	0.42
1:B:102:THR:CG2	1:B:280:HIS:CE1	3.03	0.41
1:A:161:LEU:HB3	1:A:246:ILE:CD1	2.50	0.41
1:A:170:ARG:HD2	1:A:170:ARG:HA	1.71	0.41
1:C:145:TYR:HB3	1:D:145:TYR:CD1	2.56	0.41
1:C:33:ARG:NH2	1:C:301:TYR:CE1	2.89	0.41
1:C:242:LYS:HG3	3:C:539:HOH:O	2.20	0.41
1:C:280:HIS:CD2	1:C:280:HIS:N	2.88	0.41
1:B:122:ILE:HG13	1:B:125:GLY:HA3	2.02	0.41
1:B:109:LEU:HD23	1:B:109:LEU:C	2.42	0.40
1:B:158:ARG:NH1	1:B:250:TYR:CZ	2.89	0.40
1:C:346:LEU:C	1:C:346:LEU:HD23	2.42	0.40
1:D:31:ILE:HD11	1:D:258:ILE:HD11	2.03	0.40
1:D:74:LYS:HD2	1:D:74:LYS:HA	1.75	0.40
1:D:194:ILE:HD11	1:D:203:VAL:HG21	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	334/335 (100%)	330 (99%)	4 (1%)	0	100	100
1	B	331/335 (99%)	324 (98%)	7 (2%)	0	100	100
1	C	331/335 (99%)	328 (99%)	3 (1%)	0	100	100
1	D	332/335 (99%)	326 (98%)	6 (2%)	0	100	100
All	All	1328/1340 (99%)	1308 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	293/292 (100%)	289 (99%)	4 (1%)	67	80
1	B	290/292 (99%)	283 (98%)	7 (2%)	49	62
1	C	290/292 (99%)	288 (99%)	2 (1%)	84	91
1	D	291/292 (100%)	290 (100%)	1 (0%)	92	97
All	All	1164/1168 (100%)	1150 (99%)	14 (1%)	71	83

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	49	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	155	ARG
1	A	278	SER
1	B	33	ARG
1	B	152	GLN
1	B	153	VAL
1	B	156	LYS
1	B	160	LYS
1	B	210	HIS
1	B	280	HIS
1	C	162	MET
1	C	280	HIS
1	D	192	ILE

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

Mol	Chain	Res	Type
1	A	107	GLN
1	B	210	HIS
1	C	151	GLN
1	C	304	ASN
1	C	342	GLN
1	C	352	ASN
1	D	63	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 monosaccharides 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	$\# Z > 2$
2	PO4	B	401	-	4,4,4	0.68	0	6,6,6	0.43	0
2	PO4	C	401	-	4,4,4	0.65	0	6,6,6	0.50	0
2	PO4	B	403	-	4,4,4	0.78	0	6,6,6	0.41	0
2	PO4	A	401	-	4,4,4	0.54	0	6,6,6	0.46	0
2	PO4	C	403	-	4,4,4	0.59	0	6,6,6	0.48	0
2	PO4	C	402	-	4,4,4	0.72	0	6,6,6	0.46	0
2	PO4	B	402	-	4,4,4	0.85	0	6,6,6	0.57	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	335/335 (100%)	-0.10	7 (2%) 63 61	24, 40, 69, 84	2 (0%)
1	B	331/335 (98%)	0.23	27 (8%) 11 10	23, 40, 70, 101	1 (0%)
1	C	331/335 (98%)	0.12	19 (5%) 23 22	23, 37, 73, 115	1 (0%)
1	D	331/335 (98%)	-0.11	8 (2%) 59 56	22, 39, 65, 86	1 (0%)
All	All	1328/1340 (99%)	0.04	61 (4%) 32 31	22, 39, 69, 115	5 (0%)

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	312	ASN	5.0
1	B	170	ARG	4.4
1	C	154	ASP	4.3
1	D	23	ASP	4.3
1	C	352	ASN	4.2
1	B	154	ASP	4.1
1	B	314	ALA	3.9
1	B	156	LYS	3.7
1	C	300	PRO	3.6
1	D	155	ARG	3.5
1	C	269	LYS	3.3
1	B	313	GLY	3.2
1	B	316	TYR	3.2
1	D	156	LYS	3.2
1	B	315	ASN	3.2
1	B	284	ILE	3.1
1	B	268	LYS	3.0
1	B	23	ASP	2.9
1	A	279	ALA	2.9
1	A	219	GLU	2.9
1	A	170	ARG	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	332	ALA	2.8
1	B	174	GLN	2.8
1	C	348	HIS	2.8
1	B	334	GLY	2.7
1	A	220	GLY	2.7
1	C	23	ASP	2.6
1	B	31	ILE	2.6
1	B	304	ASN	2.6
1	C	49	GLN	2.6
1	A	269	LYS	2.6
1	B	279	ALA	2.5
1	D	157	GLU	2.5
1	C	314	ALA	2.5
1	B	155	ARG	2.5
1	C	279	ALA	2.5
1	B	32[A]	ILE	2.5
1	C	155	ARG	2.4
1	B	167	PHE	2.4
1	D	279	ALA	2.4
1	B	299	PRO	2.4
1	B	168	SER	2.4
1	C	156	LYS	2.4
1	D	221	LEU	2.4
1	A	188	ARG	2.3
1	C	170	ARG	2.3
1	C	47	ASN	2.3
1	C	51	GLN	2.3
1	B	285	ALA	2.3
1	B	159	GLU	2.2
1	B	305	VAL	2.2
1	B	33	ARG	2.2
1	C	32	ILE	2.2
1	D	181	ASP	2.2
1	C	316	TYR	2.1
1	C	284	ILE	2.1
1	C	298	SER	2.1
1	B	171	GLU	2.1
1	D	348	HIS	2.0
1	B	163	GLU	2.0
1	A	348	HIS	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 monosaccharides 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	PO4	B	403	5/5	0.57	0.54	230,231,239,258	0
2	PO4	C	402	5/5	0.84	0.26	88,102,106,106	1
2	PO4	A	401	5/5	0.88	0.16	82,88,97,101	0
2	PO4	B	402	5/5	0.90	0.25	53,66,70,71	5
2	PO4	C	403	5/5	0.90	0.20	91,93,97,101	0
2	PO4	B	401	5/5	0.91	0.26	91,92,102,106	0
2	PO4	C	401	5/5	0.93	0.10	67,69,74,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.