



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:21 PM UTC

PDB ID : 7DOQ / pdb_00007doq
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 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

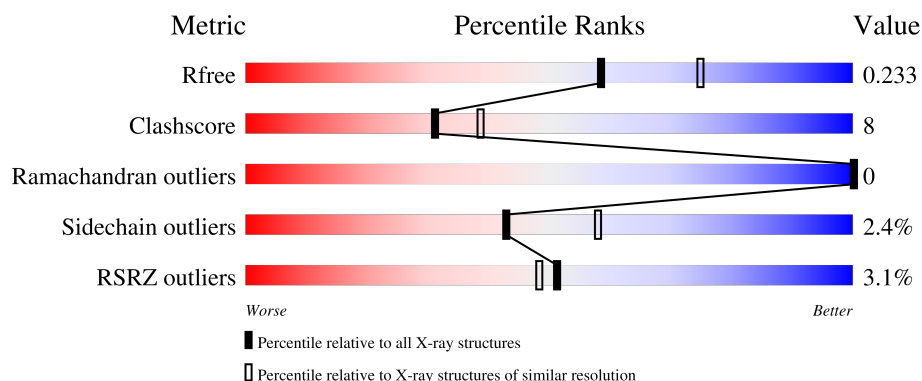
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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	2% 87% 12% .
1	B	335	3% 83% 15% ..
1	C	335	5% 84% 14% ..
1	D	335	2% 85% 13% ..

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 (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	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		

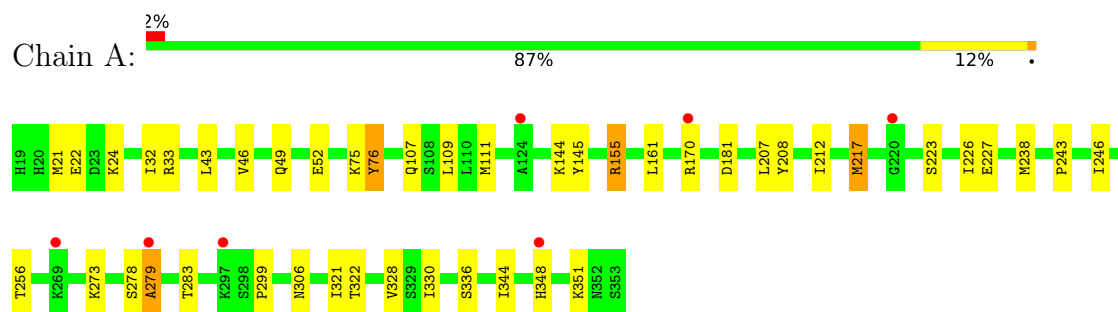
- Molecule 3 is water.

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		

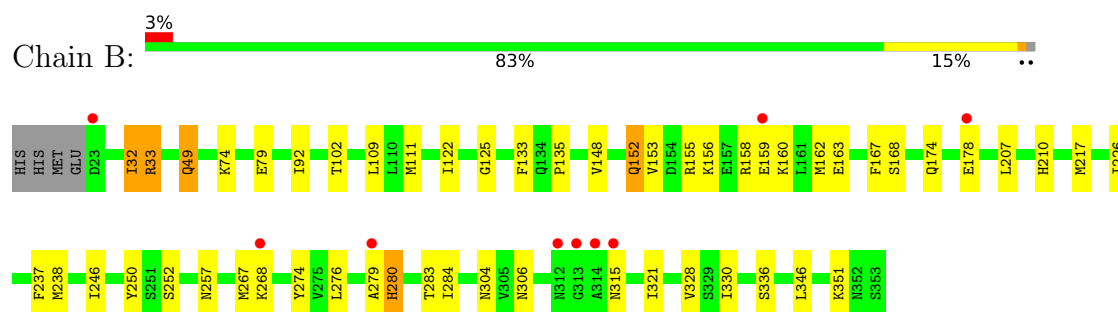
3 Residue-property plots [i](#)

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.

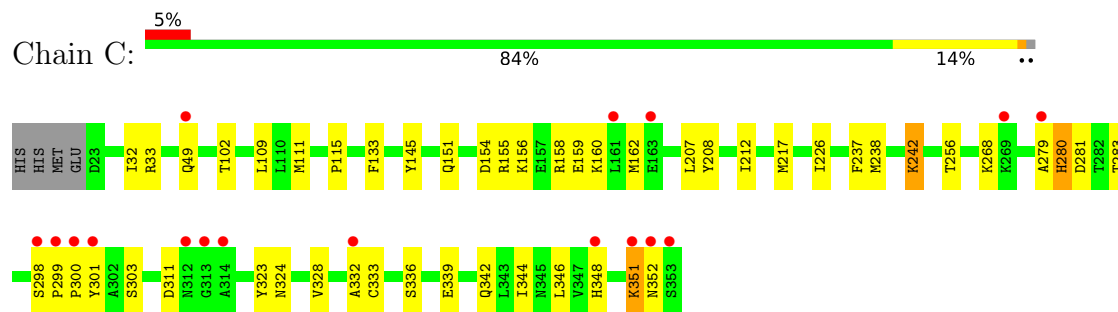
• Molecule 1: Acid phosphatase



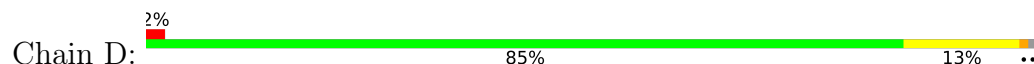
• Molecule 1: Acid phosphatase

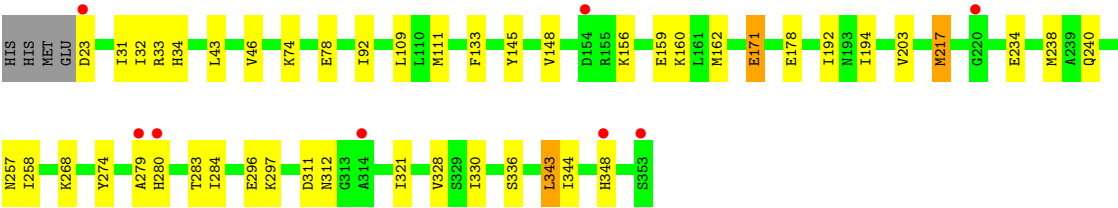


• Molecule 1: Acid phosphatase



• 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 35.62 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.4 (35.62-2.20) 96.5 (35.62-2.20)	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 (4.88%)	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)	0.39 , 38.1	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	0.95	0/2727	1.16	5/3708 (0.1%)
1	B	0.95	0/2700	1.18	0/3672
1	C	0.95	0/2696	1.14	1/3667 (0.0%)
1	D	0.95	0/2708	1.15	0/3683
All	All	0.95	0/10831	1.16	6/14730 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	339	GLU	CB-CA-C	5.27	119.23	109.54
1	A	279	ALA	O-C-N	5.12	128.69	122.14
1	A	279	ALA	CA-C-N	5.01	132.00	123.03
1	A	279	ALA	C-N-CA	5.01	132.00	123.03
1	A	76	TYR	CA-C-N	-5.00	118.07	122.97
1	A	76	TYR	C-N-CA	-5.00	118.07	122.97

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	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2629	0	2595	46	0
1	C	2626	0	2590	58	0
1	D	2634	0	2606	44	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
3	C	63	0	0	1	0
3	D	68	0	0	0	0
All	All	10841	0	10406	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 8.

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: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:D:162:MET:HE1	1:D:240:GLN:HG2	1.50	0.91
1:C:300:PRO:HD2	1:C:323:TYR:OH	1.72	0.88
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:B:207:LEU:HD22	1:B:217:MET:HE1	1.71	0.73
1:C:279:ALA:CB	1:C:283:THR:HG21	2.19	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:315:ASN:CG	1:C:256:THR:HG21	2.22	0.64
1:B:267:MET:HE2	1:C:151:GLN:NE2	2.13	0.63
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:267:MET:C	1:B:268:LYS:HG3	2.25	0.61
1:B:49:GLN:HA	1:B:49:GLN:OE1	1.99	0.61
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: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: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:A:238:MET:HA	1:A:238:MET:HE2	1.86	0.57
1:B:111:MET:HE2	1:B:133[A]:PHE:HE1	1.69	0.57
1:C:155:ARG:NH1	1:C:158:ARG:HD3	2.19	0.57
1:B:33:ARG:HB3	1:B:284:ILE:HD12	1.86	0.57
1:D:171:GLU:H	1:D:171:GLU:CD	2.12	0.56
1:B:162:MET:HE2	1:B:246:ILE:HD12	1.88	0.56
1:B:238:MET:HE2	1:B:238:MET:HA	1.87	0.56
1:B:163:GLU:HA	1:B:167:PHE:HB2	1.87	0.55
1:C:217:MET:HE2	1:C:226:ILE:HD12	1.89	0.55
1:A:279:ALA:HB1	1:A:283:THR:OG1	2.07	0.55
1:B:32[A]:ILE:HG13	1:B:276:LEU:HD11	1.89	0.55
1:B:217:MET:HE2	1:B:226:ILE:HD12	1.89	0.55
1:C:111:MET:HE1	1:D:133[A]:PHE:CD2	2.42	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:THR:HG21	1:B:280:HIS:CE1	2.44	0.53
1:C:300:PRO:O	1:C:303:SER:CB	2.58	0.52
1:D:238:MET:HA	1:D:238:MET:HE2	1.90	0.52
1:A:43:LEU:HB2	1:A:46:VAL:HG22	1.91	0.52
1:C:238:MET:HA	1:C:238:MET:HE2	1.90	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
1:D:344:ILE:CG2	1:D:348:HIS:NE2	2.75	0.50
1:D:217:MET:HE3	1:D:217:MET:HA	1.94	0.49
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:23:ASP:CG	1:D:312:ASN:HA	2.38	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: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:A:111:MET:HE1	1:B:133[A]:PHE:CD2	2.49	0.48
1:B:163:GLU:O	1:B:163:GLU:HG3	2.13	0.47
1:A:217:MET:HE2	1:A:226:ILE:HD12	1.96	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:207:LEU:HD22	1:C:217:MET:HE1	1.98	0.46
1:C:279:ALA:HB2	1:C:283:THR:HG21	1.95	0.46
1:A:207:LEU:HD22	1:A:217:MET:HE1	1.97	0.46
1:B:207:LEU:CD2	1:B:217:MET:HE1	2.44	0.46
1:D:234:GLU:OE2	1:D:297:LYS:CD	2.62	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:109:LEU:C	1:B:109:LEU:HD23	2.42	0.45
1:B:148:VAL:HA	1:B:257:ASN:ND2	2.32	0.45
1:C:111:MET:HE3	1:C:111:MET:HB2	1.80	0.45
1:D:156:LYS:O	1:D:159:GLU:HG2	2.16	0.45
1:C:346:LEU:C	1:C:346:LEU:HD23	2.42	0.45
1:D:279:ALA:HB1	1:D:283:THR:OG1	2.16	0.45
1:B:111:MET:HE3	1:B:111:MET:HB2	1.83	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:D:111:MET:HE2	1:D:133[A]:PHE:HE1	1.81	0.45
1:B:32[B]:ILE:HD13	1:B:109:LEU:HD13	1.99	0.44
1:D:343:LEU:O	1:D:343:LEU:HD23	2.18	0.44
1:A:321:ILE:HD12	1:A:330:ILE:CD1	2.45	0.44
1:C:102:THR:HG21	1:C:280:HIS:CE1	2.52	0.44
1:A:111:MET:HE3	1:A:111:MET:HB2	1.85	0.43
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:C:133:PHE:CD2	1:D:111:MET:HE1	2.53	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:D:343:LEU:HD23	1:D:343:LEU:C	2.43	0.43
1:B:163:GLU:O	1:B:168:SER:HB3	2.19	0.42
1:D:109:LEU:C	1:D:109:LEU:HD23	2.44	0.42
1:D:279:ALA:HB2	1:D:283:THR:HG21	2.00	0.42
1:C:111:MET:HE2	1:C:133:PHE:HE1	1.84	0.42
1:A:306:ASN:HB3	1:A:322:THR:OG1	2.19	0.42
1:C:217:MET:HE3	1:C:217:MET:HA	2.00	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:LEU:HB3	1:A:246:ILE:HD13	2.02	0.42
1:B:346:LEU:C	1:B:346:LEU:HD23	2.45	0.42
1:A:217:MET:HA	1:A:217:MET:HE3	2.00	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: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:D:33:ARG:HG2	1:D:34:HIS:N	2.34	0.41
1:D:111:MET:HE2	1:D:133[A]:PHE:CE1	2.55	0.41
1:B:111:MET:HE2	1:B:133[A]:PHE:CE1	2.51	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:237:PHE:CD2	1:B:238:MET:HE3	2.56	0.41
1:A:21:MET:HE2	1:A:24:LYS:NZ	2.36	0.40
1:B:158:ARG:NH1	1:B:250:TYR:CZ	2.89	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

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
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%)	286 (98%)	7 (2%)	43	58
1	B	290/292 (99%)	280 (97%)	10 (3%)	32	44
1	C	290/292 (99%)	285 (98%)	5 (2%)	53	69
1	D	291/292 (100%)	284 (98%)	7 (2%)	43	58
All	All	1164/1168 (100%)	1135 (98%)	29 (2%)	43	56

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	49	GLN
1	A	155	ARG
1	A	181	ASP
1	A	217	MET
1	A	273	LYS
1	A	278	SER
1	B	32[A]	ILE
1	B	32[B]	ILE
1	B	33	ARG
1	B	49	GLN
1	B	152	GLN
1	B	153	VAL

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Mol	Chain	Res	Type
1	B	156	LYS
1	B	160	LYS
1	B	210	HIS
1	B	280	HIS
1	C	160	LYS
1	C	162	MET
1	C	242	LYS
1	C	280	HIS
1	C	351	LYS
1	D	171	GLU
1	D	178	GLU
1	D	192	ILE
1	D	217	MET
1	D	280	HIS
1	D	296	GLU
1	D	343	LEU

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	81	HIS
1	A	107	GLN
1	A	131	HIS
1	A	248	ASN
1	A	257	ASN
1	B	81	HIS
1	B	131	HIS
1	B	306	ASN
1	B	352	ASN
1	C	107	GLN
1	C	151	GLN
1	C	257	ASN
1	C	304	ASN
1	C	306	ASN
1	C	352	ASN
1	D	306	ASN

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 oligosaccharides 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	0.73	0	6,6,6	0.46	0
2	PO4	C	403	-	4,4,4	0.63	0	6,6,6	0.51	0
2	PO4	A	401	-	4,4,4	0.58	0	6,6,6	0.49	0
2	PO4	C	401	-	4,4,4	0.70	0	6,6,6	0.53	0
2	PO4	B	403	-	4,4,4	0.83	0	6,6,6	0.44	0
2	PO4	B	402	-	4,4,4	0.90	0	6,6,6	0.60	0
2	PO4	C	402	-	4,4,4	0.78	0	6,6,6	0.49	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [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	335/335 (100%)	0.21	7 (2%) 63 60	17, 40, 69, 84	3 (0%)
1	B	331/335 (98%)	0.23	9 (2%) 56 53	17, 40, 70, 101	3 (0%)
1	C	331/335 (98%)	0.20	17 (5%) 33 30	17, 37, 73, 115	3 (0%)
1	D	331/335 (98%)	0.16	8 (2%) 59 56	16, 39, 65, 81	4 (1%)
All	All	1328/1340 (99%)	0.20	41 (3%) 51 48	16, 39, 69, 115	13 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	300	PRO	6.4
1	C	279	ALA	4.7
1	B	268	LYS	4.6
1	C	299	PRO	4.2
1	D	279	ALA	4.0
1	B	315	ASN	4.0
1	C	298	SER	3.9
1	D	23	ASP	3.9
1	C	348	HIS	3.6
1	A	348	HIS	3.6
1	B	279	ALA	3.5
1	A	279	ALA	3.3
1	C	314	ALA	3.2
1	B	312	ASN	3.1
1	A	220	GLY	3.0
1	B	314	ALA	2.9
1	B	313	GLY	2.9
1	C	351	LYS	2.8
1	A	124	ALA	2.7
1	C	332	ALA	2.6
1	C	312	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	352	ASN	2.5
1	D	353	SER	2.5
1	C	161	LEU	2.5
1	C	301	TYR	2.4
1	C	353	SER	2.4
1	D	280	HIS	2.3
1	A	269	LYS	2.3
1	D	220	GLY	2.3
1	B	23	ASP	2.2
1	C	269	LYS	2.2
1	C	313	GLY	2.2
1	D	314	ALA	2.2
1	B	178	GLU	2.2
1	C	49	GLN	2.2
1	D	348	HIS	2.2
1	C	163	GLU	2.1
1	A	170	ARG	2.1
1	B	159	GLU	2.1
1	D	154	ASP	2.0
1	A	297	LYS	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	B	403	5/5	0.54	0.33	230,231,239,258	0
2	PO4	C	402	5/5	0.76	0.18	88,102,106,106	1
2	PO4	C	403	5/5	0.77	0.18	91,93,97,101	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	A	401	5/5	0.83	0.14	82,88,97,101	0
2	PO4	B	401	5/5	0.86	0.09	91,92,102,106	0
2	PO4	B	402	5/5	0.89	0.15	53,66,70,71	5
2	PO4	C	401	5/5	0.90	0.11	67,69,74,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.