



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:52 PM UTC

PDB ID : 6S2Q / pdb_00006s2q
Title : Mycobacterial hydrolase 1
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Deposited on : 2019-06-21
Resolution : 2.50 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
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

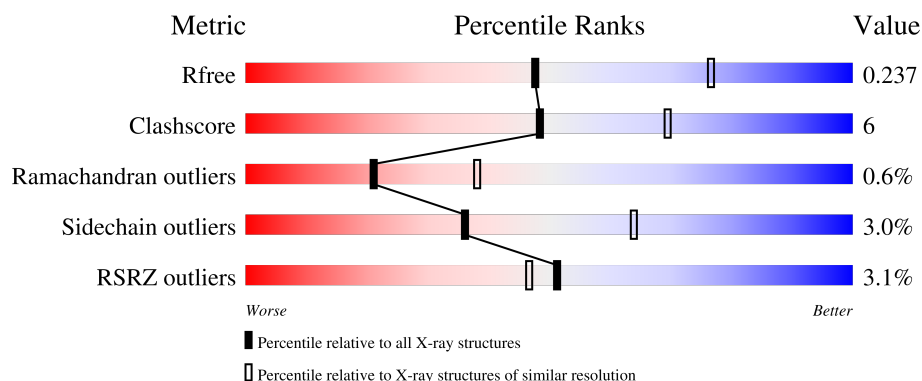
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	236	
1	B	236	
1	C	236	
1	D	236	

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	IOD	A	304	-	-	X	-
2	IOD	A	307	-	-	X	-
2	IOD	D	305	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6738 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 Glucosyl-3-phosphoglycerate phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	4	0
			1678	1048	318	309	3			
1	B	210	Total	C	N	O	S	0	2	0
			1631	1018	308	302	3			
1	C	210	Total	C	N	O	S	0	3	0
			1639	1019	314	303	3			
1	D	208	Total	C	N	O	S	0	0	0
			1595	993	307	292	3			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	224	LYS	-	expression tag	UNP P9WIC7
A	225	LEU	-	expression tag	UNP P9WIC7
A	226	ALA	-	expression tag	UNP P9WIC7
A	227	ALA	-	expression tag	UNP P9WIC7
A	228	ALA	-	expression tag	UNP P9WIC7
A	229	LEU	-	expression tag	UNP P9WIC7
A	230	GLU	-	expression tag	UNP P9WIC7
A	231	HIS	-	expression tag	UNP P9WIC7
A	232	HIS	-	expression tag	UNP P9WIC7
A	233	HIS	-	expression tag	UNP P9WIC7
A	234	HIS	-	expression tag	UNP P9WIC7
A	235	HIS	-	expression tag	UNP P9WIC7
A	236	HIS	-	expression tag	UNP P9WIC7
B	224	LYS	-	expression tag	UNP P9WIC7
B	225	LEU	-	expression tag	UNP P9WIC7
B	226	ALA	-	expression tag	UNP P9WIC7
B	227	ALA	-	expression tag	UNP P9WIC7
B	228	ALA	-	expression tag	UNP P9WIC7
B	229	LEU	-	expression tag	UNP P9WIC7
B	230	GLU	-	expression tag	UNP P9WIC7
B	231	HIS	-	expression tag	UNP P9WIC7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	232	HIS	-	expression tag	UNP P9WIC7
B	233	HIS	-	expression tag	UNP P9WIC7
B	234	HIS	-	expression tag	UNP P9WIC7
B	235	HIS	-	expression tag	UNP P9WIC7
B	236	HIS	-	expression tag	UNP P9WIC7
C	224	LYS	-	expression tag	UNP P9WIC7
C	225	LEU	-	expression tag	UNP P9WIC7
C	226	ALA	-	expression tag	UNP P9WIC7
C	227	ALA	-	expression tag	UNP P9WIC7
C	228	ALA	-	expression tag	UNP P9WIC7
C	229	LEU	-	expression tag	UNP P9WIC7
C	230	GLU	-	expression tag	UNP P9WIC7
C	231	HIS	-	expression tag	UNP P9WIC7
C	232	HIS	-	expression tag	UNP P9WIC7
C	233	HIS	-	expression tag	UNP P9WIC7
C	234	HIS	-	expression tag	UNP P9WIC7
C	235	HIS	-	expression tag	UNP P9WIC7
C	236	HIS	-	expression tag	UNP P9WIC7
D	224	LYS	-	expression tag	UNP P9WIC7
D	225	LEU	-	expression tag	UNP P9WIC7
D	226	ALA	-	expression tag	UNP P9WIC7
D	227	ALA	-	expression tag	UNP P9WIC7
D	228	ALA	-	expression tag	UNP P9WIC7
D	229	LEU	-	expression tag	UNP P9WIC7
D	230	GLU	-	expression tag	UNP P9WIC7
D	231	HIS	-	expression tag	UNP P9WIC7
D	232	HIS	-	expression tag	UNP P9WIC7
D	233	HIS	-	expression tag	UNP P9WIC7
D	234	HIS	-	expression tag	UNP P9WIC7
D	235	HIS	-	expression tag	UNP P9WIC7
D	236	HIS	-	expression tag	UNP P9WIC7

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	7	Total I 7 7	0	1
2	B	5	Total I 5 5	0	1
2	C	6	Total I 6 6	0	1
2	D	5	Total I 5 5	0	0

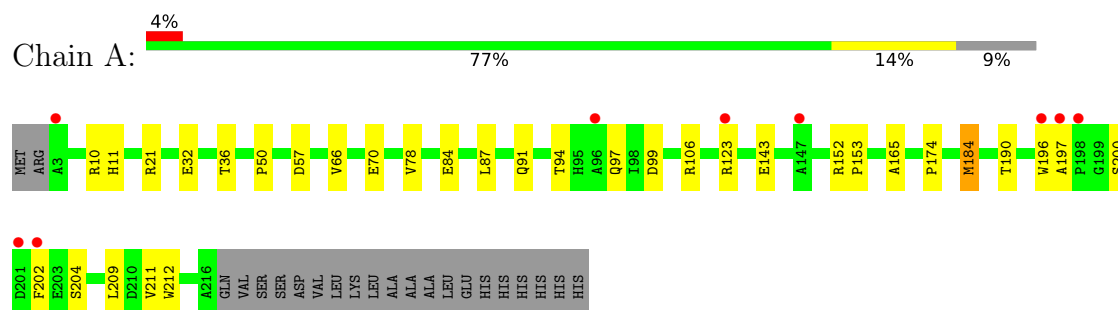
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total 48	O 48	0	0
3	B	17	Total 17	O 17	0	0
3	C	81	Total 81	O 81	0	0
3	D	26	Total 26	O 26	0	0

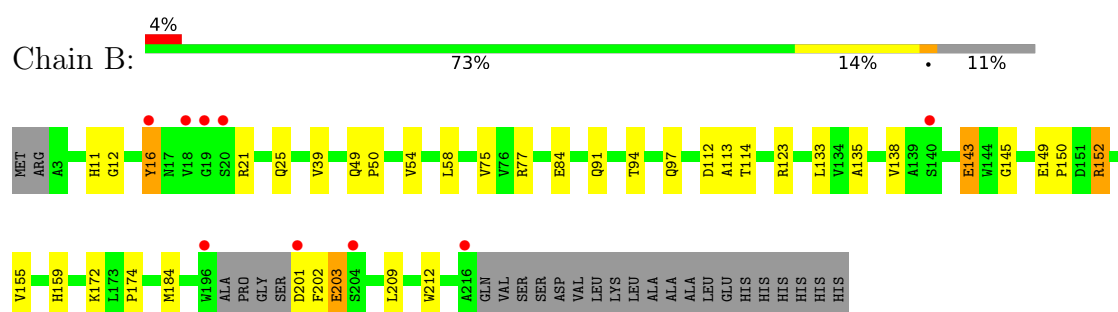
3 Residue-property plots [i](#)

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.

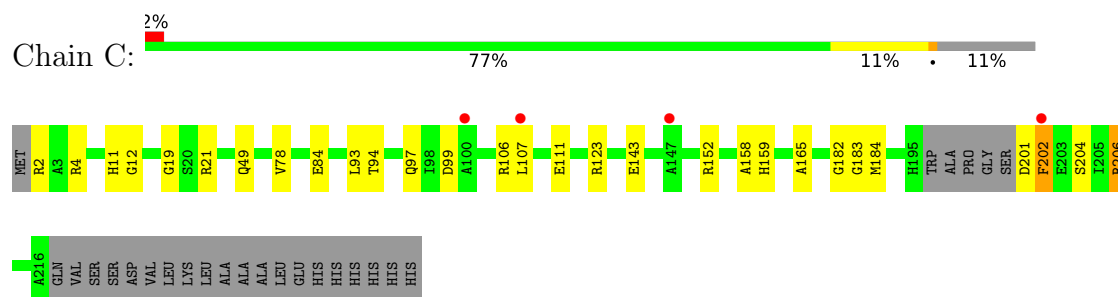
- Molecule 1: Glucosyl-3-phosphoglycerate phosphatase



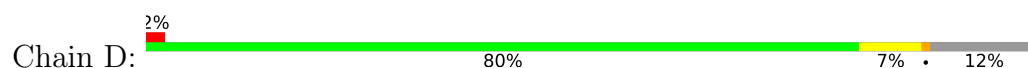
- Molecule 1: Glucosyl-3-phosphoglycerate phosphatase

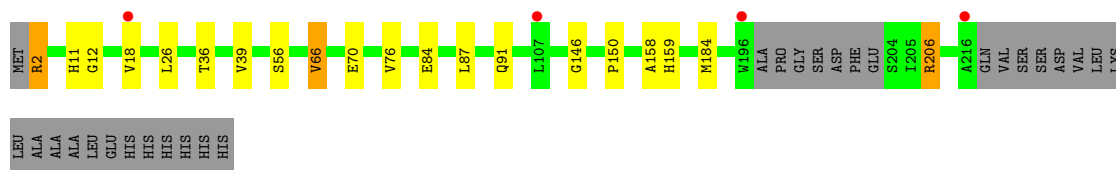


- Molecule 1: Glucosyl-3-phosphoglycerate phosphatase



- Molecule 1: Glucosyl-3-phosphoglycerate phosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	96.95Å 46.10Å 103.25Å 90.00° 104.29° 90.00°	Depositor
Resolution (Å)	50.03 – 2.50 50.03 – 2.50	Depositor EDS
% Data completeness (in resolution range)	88.0 (50.03-2.50) 88.2 (50.03-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???)	Depositor
R, R_{free}	0.187 , 0.233 0.191 , 0.237	Depositor DCC
R_{free} test set	1325 reflections (4.25%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6738	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.55% 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: IOD

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	0.41	0/1717	0.64	0/2340
1	B	0.34	0/1666	0.61	0/2268
1	C	0.48	0/1672	0.73	4/2273 (0.2%)
1	D	0.35	0/1628	0.60	0/2216
All	All	0.40	0/6683	0.65	4/9097 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	202[A]	PHE	CA-C-N	-5.09	112.21	120.72
1	C	202[A]	PHE	C-N-CA	-5.09	112.21	120.72
1	C	202[B]	PHE	CA-C-N	-5.09	112.21	120.72
1	C	202[B]	PHE	C-N-CA	-5.09	112.21	120.72

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	1678	0	1643	22	0
1	B	1631	0	1600	23	0
1	C	1639	0	1615	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1595	0	1583	11	0
2	A	7	0	0	7	0
2	B	5	0	0	1	0
2	C	6	0	0	1	0
2	D	5	0	0	3	0
3	A	48	0	0	0	0
3	B	17	0	0	0	0
3	C	81	0	0	3	0
3	D	26	0	0	0	0
All	All	6738	0	6441	72	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 6.

All (72) 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:11:HIS:NE2	2:A:307:IOD:I	2.77	0.88
1:C:11:HIS:NE2	2:C:306:IOD:I	2.82	0.82
1:C:201[B]:ASP:N	1:C:204:SER:HG	1.77	0.81
1:C:2:ARG:HH21	1:C:4:ARG:HH22	1.29	0.80
1:B:21:ARG:HG2	1:B:94:THR:HG22	1.71	0.71
1:C:111:GLU:OE1	1:D:206:ARG:NE	2.22	0.70
1:D:11:HIS:NE2	2:D:305:IOD:I	2.91	0.65
1:C:2:ARG:NH2	1:C:4:ARG:HH22	1.95	0.65
1:A:21:ARG:NH2	2:A:304:IOD:I	3.00	0.64
1:B:25:GLN:N	1:B:91:GLN:OE1	2.30	0.63
1:A:200:SER:HB3	1:A:204:SER:OG	1.99	0.63
1:B:94:THR:OG1	1:B:97:GLN:HG3	1.99	0.62
1:A:94:THR:HG22	1:A:97:GLN:HG3	1.80	0.61
1:D:87:LEU:HB2	1:D:91:GLN:HG2	1.84	0.59
1:C:99:ASP:OD2	1:C:106:ARG:NH1	2.38	0.56
2:A:304:IOD:I	1:C:21:ARG:NH1	3.06	0.56
1:C:183:GLY:HA2	3:C:440:HOH:O	2.04	0.56
1:B:149:GLU:OE1	1:B:152:ARG:NH2	2.38	0.56
1:A:32:GLU:O	1:A:36:THR:HG23	2.05	0.55
1:C:84:GLU:HB2	1:C:159:HIS:CG	2.41	0.55
1:B:50:PRO:HB3	1:B:155:VAL:HG23	1.88	0.55
1:B:201[A]:ASP:OD1	1:B:202[A]:PHE:N	2.41	0.54
1:C:49:GLN:O	1:C:152:ARG:HD2	2.07	0.54
1:C:94:THR:HG22	1:C:97:GLN:HG3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201[B]:ASP:OD1	1:B:202[B]:PHE:N	2.43	0.52
1:A:174:PRO:HG2	1:B:174:PRO:HG2	1.91	0.52
1:D:84:GLU:OE2	2:D:305:IOD:I	2.98	0.52
1:A:87:LEU:HB2	1:A:91:GLN:HG2	1.93	0.51
1:B:84:GLU:OE2	2:B:304:IOD:I	2.99	0.50
1:C:206:ARG:HD3	3:C:441:HOH:O	2.13	0.48
1:A:106:ARG:NE	2:A:303:IOD:I	3.12	0.48
1:B:112:ASP:OD1	1:B:114:THR:OG1	2.27	0.48
1:C:99:ASP:OD1	1:C:106:ARG:HD3	2.14	0.48
1:B:172:LYS:NZ	1:B:203:GLU:O	2.46	0.47
1:D:11:HIS:ND1	1:D:12:GLY:O	2.48	0.47
1:B:49:GLN:O	1:B:152:ARG:HG3	2.14	0.47
1:A:190:THR:HG23	1:A:211:VAL:O	2.15	0.47
1:C:11:HIS:ND1	1:C:12:GLY:O	2.48	0.47
1:A:99:ASP:OD1	1:A:106:ARG:NH1	2.48	0.46
2:A:304:IOD:I	1:C:19:GLY:HA3	2.86	0.46
1:B:16:TYR:CE2	1:B:21:ARG:HB2	2.51	0.46
1:A:84:GLU:OE2	2:A:307:IOD:I	3.03	0.46
1:C:123[A]:ARG:HD2	1:C:165:ALA:HB2	1.98	0.46
1:D:66:VAL:O	1:D:70:GLU:HG3	2.15	0.46
1:A:202[B]:PHE:N	1:A:202[B]:PHE:CD1	2.84	0.45
1:B:209:LEU:HD11	1:B:212:TRP:HB2	1.98	0.45
1:A:50:PRO:HA	1:A:153:PRO:HD2	1.98	0.45
1:A:123[A]:ARG:HD2	1:A:165:ALA:HB2	1.98	0.45
1:B:145:GLY:HA2	1:B:150:PRO:O	2.15	0.45
1:C:107:LEU:HG	3:C:418:HOH:O	2.17	0.45
1:B:77:ARG:HE	1:B:77:ARG:HB2	1.55	0.45
1:A:10:ARG:NH1	1:A:184:MET:O	2.50	0.44
1:A:50:PRO:O	1:A:152:ARG:NH2	2.28	0.44
1:A:209:LEU:HD21	1:A:212:TRP:HB2	2.00	0.44
1:A:66:VAL:O	1:A:70:GLU:HG3	2.17	0.43
1:B:58:LEU:HD12	1:B:159:HIS:CE1	2.53	0.43
1:B:135:ALA:O	1:B:138:VAL:HG12	2.18	0.43
1:C:99:ASP:CG	1:C:106:ARG:HH11	2.27	0.43
1:B:184:MET:HE2	1:B:184:MET:HB3	1.91	0.42
1:A:196:TRP:HD1	1:A:197:ALA:O	2.03	0.41
1:B:11:HIS:ND1	1:B:12:GLY:O	2.53	0.41
1:D:36:THR:O	1:D:39:VAL:HG12	2.20	0.41
1:A:123[A]:ARG:NH2	2:A:301:IOD:I	3.22	0.41
1:B:113:ALA:HB1	1:B:123:ARG:HB2	2.02	0.41
1:B:143:GLU:CD	1:B:143:GLU:H	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:HIS:HB3	2:D:305:IOD:I	2.91	0.41
1:B:54:VAL:HG11	1:B:133:LEU:HD21	2.02	0.40
1:D:146:GLY:O	1:D:150:PRO:HB3	2.21	0.40
1:D:2:ARG:HB3	1:D:2:ARG:NH1	2.36	0.40
1:D:184:MET:HB3	1:D:184:MET:HE3	1.75	0.40
1:A:57:ASP:HB3	1:A:78:VAL:HG13	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	216/236 (92%)	207 (96%)	9 (4%)	0	100	100
1	B	207/236 (88%)	199 (96%)	7 (3%)	1 (0%)	24	43
1	C	208/236 (88%)	200 (96%)	4 (2%)	4 (2%)	6	11
1	D	204/236 (86%)	199 (98%)	4 (2%)	1 (0%)	24	43
All	All	835/944 (88%)	805 (96%)	24 (3%)	6 (1%)	21	34

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	202[A]	PHE
1	C	202[B]	PHE
1	B	16	TYR
1	C	158	ALA
1	D	158	ALA
1	C	182	GLY

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	167/182 (92%)	165 (99%)	2 (1%)	63	83
1	B	163/182 (90%)	158 (97%)	5 (3%)	35	62
1	C	164/182 (90%)	159 (97%)	5 (3%)	36	64
1	D	159/182 (87%)	152 (96%)	7 (4%)	25	50
All	All	653/728 (90%)	634 (97%)	19 (3%)	36	65

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	GLU
1	A	184	MET
1	B	39	VAL
1	B	75	VAL
1	B	143	GLU
1	B	152	ARG
1	B	203	GLU
1	C	78	VAL
1	C	93	LEU
1	C	143	GLU
1	C	184	MET
1	C	206	ARG
1	D	2	ARG
1	D	18	VAL
1	D	26	LEU
1	D	56	SER
1	D	66	VAL
1	D	76	VAL
1	D	206	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	ASN
1	A	186	ASN
1	B	13	GLN
1	B	17	ASN
1	B	49	GLN
1	B	195	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](#)

Of 23 ligands modelled in this entry, 23 are monoatomic - leaving 0 for Mogul analysis.

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	214/236 (90%)	0.12	9 (4%)	40	36	16, 40, 73, 96	4 (1%)
1	B	210/236 (88%)	0.31	9 (4%)	40	35	35, 56, 82, 93	2 (0%)
1	C	210/236 (88%)	-0.16	4 (1%)	66	62	12, 32, 64, 104	3 (1%)
1	D	208/236 (88%)	0.26	4 (1%)	66	62	31, 56, 79, 87	0
All	All	842/944 (89%)	0.13	26 (3%)	51	47	12, 46, 77, 104	9 (1%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	202[A]	PHE	4.6
1	B	196	TRP	3.6
1	A	147	ALA	3.5
1	A	198	PRO	3.5
1	B	216	ALA	3.4
1	B	19	GLY	3.3
1	C	100	ALA	3.1
1	B	201[A]	ASP	2.9
1	B	18	VAL	2.9
1	A	202[A]	PHE	2.8
1	B	20	SER	2.8
1	D	18	VAL	2.8
1	C	147	ALA	2.6
1	A	196	TRP	2.5
1	D	196	TRP	2.5
1	A	3	ALA	2.5
1	A	96	ALA	2.5
1	D	216	ALA	2.4
1	C	107	LEU	2.4
1	A	197	ALA	2.4
1	D	107	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	140	SER	2.2
1	B	204	SER	2.2
1	B	16	TYR	2.1
1	A	123[A]	ARG	2.1
1	A	201[A]	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	IOD	A	304	1/1	0.92	0.08	67,67,67,67	0
2	IOD	D	304	1/1	0.92	0.08	94,94,94,94	1
2	IOD	D	303	1/1	0.93	0.09	79,79,79,79	1
2	IOD	B	305	1/1	0.93	0.09	86,86,86,86	1
2	IOD	A	303	1/1	0.94	0.09	80,80,80,80	1
2	IOD	B	302	1/1	0.96	0.11	90,90,90,90	0
2	IOD	C	303	1/1	0.96	0.12	77,77,77,77	0
2	IOD	C	305[B]	1/1	0.97	0.09	46,46,46,46	1
2	IOD	A	305	1/1	0.98	0.19	95,95,95,95	0
2	IOD	A	306[B]	1/1	0.98	0.06	54,54,54,54	1
2	IOD	C	304	1/1	0.98	0.06	61,61,61,61	1
2	IOD	B	301	1/1	0.98	0.09	63,63,63,63	0
2	IOD	D	301	1/1	0.98	0.11	69,69,69,69	0
2	IOD	D	302	1/1	0.98	0.10	70,70,70,70	0
2	IOD	A	302	1/1	0.98	0.19	82,82,82,82	0
2	IOD	B	303[B]	1/1	0.98	0.05	70,70,70,70	1
2	IOD	A	307	1/1	0.99	0.05	30,30,30,30	0
2	IOD	C	306	1/1	0.99	0.05	23,23,23,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	IOD	C	301	1/1	0.99	0.17	76,76,76,76	0
2	IOD	C	302	1/1	0.99	0.08	42,42,42,42	0
2	IOD	A	301	1/1	0.99	0.07	48,48,48,48	0
2	IOD	B	304	1/1	0.99	0.03	58,58,58,58	0
2	IOD	D	305	1/1	0.99	0.06	49,49,49,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.