



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:59 PM UTC

PDB ID : 3HLH / pdb_00003hllh
Title : Diisopropyl fluorophosphatase (DFPase), active site mutants
Authors : Chen, J.C.-H.; Blum, M.-M.
Deposited on : 2009-05-27
Resolution : 1.80 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
Xtrriage (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

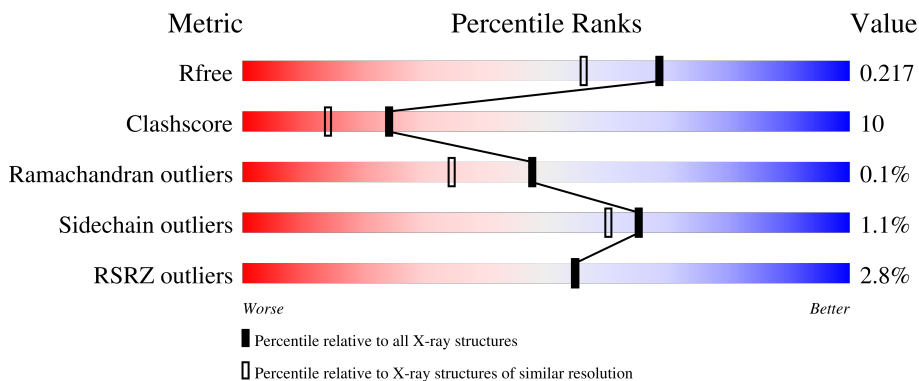
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

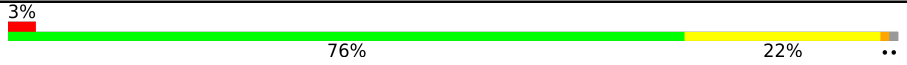



The reported resolution of this entry is 1.80 Å.

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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	314	 3% 76% 22% ..
1	B	314	 4% 74% 24% ..
1	C	314	 % 79% 19% ..
1	D	314	 4% 83% 14% ..

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 10655 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 Diisopropyl-fluorophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	312	2435	1547	413	459	16	0	0	0
1	B	312	2435	1547	413	459	16	0	0	0
1	C	312	2435	1547	413	459	16	0	0	0
1	D	312	2435	1547	413	459	16	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	37	ALA	GLU	engineered mutation	UNP Q7SIG4
A	144	ALA	TYR	engineered mutation	UNP Q7SIG4
A	146	ALA	ARG	engineered mutation	UNP Q7SIG4
A	195	MET	THR	engineered mutation	UNP Q7SIG4
B	37	ALA	GLU	engineered mutation	UNP Q7SIG4
B	144	ALA	TYR	engineered mutation	UNP Q7SIG4
B	146	ALA	ARG	engineered mutation	UNP Q7SIG4
B	195	MET	THR	engineered mutation	UNP Q7SIG4
C	37	ALA	GLU	engineered mutation	UNP Q7SIG4
C	144	ALA	TYR	engineered mutation	UNP Q7SIG4
C	146	ALA	ARG	engineered mutation	UNP Q7SIG4
C	195	MET	THR	engineered mutation	UNP Q7SIG4
D	37	ALA	GLU	engineered mutation	UNP Q7SIG4
D	144	ALA	TYR	engineered mutation	UNP Q7SIG4
D	146	ALA	ARG	engineered mutation	UNP Q7SIG4
D	195	MET	THR	engineered mutation	UNP Q7SIG4

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Ca 2 2	0	0
2	B	2	Total Ca 2 2	0	0
2	C	2	Total Ca 2 2	0	0
2	D	2	Total Ca 2 2	0	0

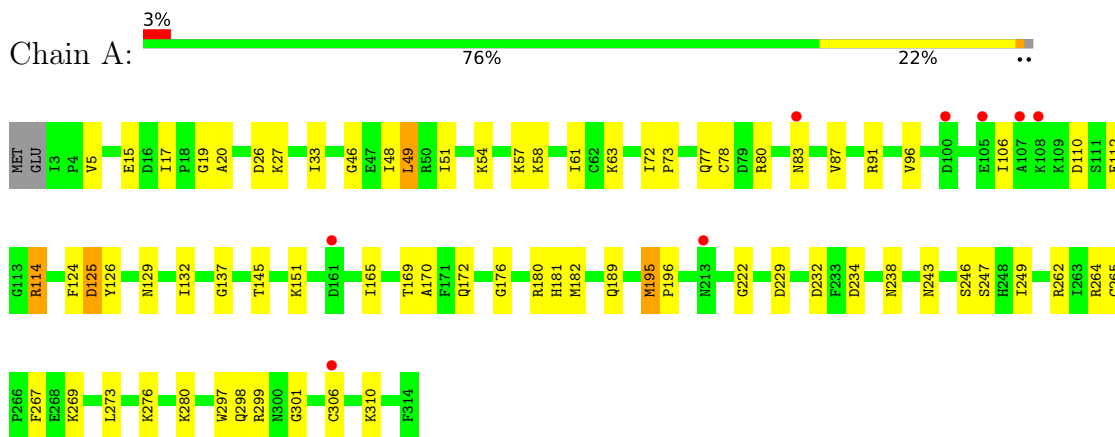
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	208	Total O 208 208	0	0
3	B	244	Total O 244 244	0	0
3	C	231	Total O 231 231	0	0
3	D	224	Total O 224 224	0	0

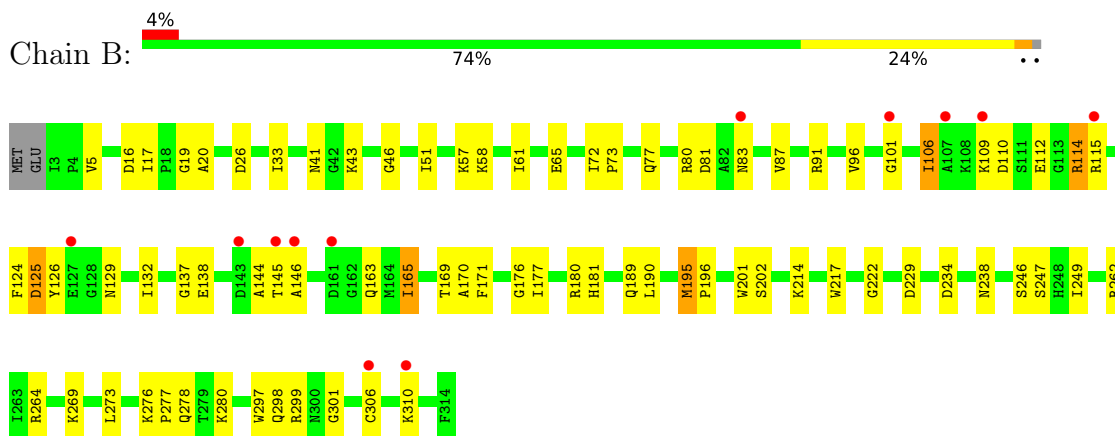
3 Residue-property plots [i](#)

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.

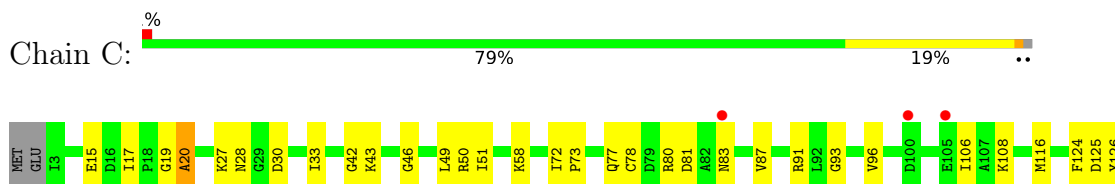
- Molecule 1: Diisopropyl-fluorophosphatase

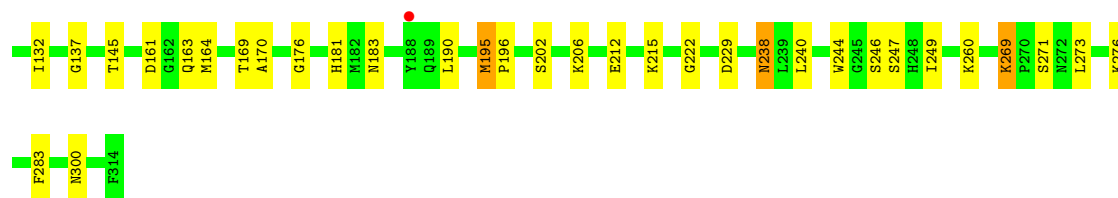


- Molecule 1: Diisopropyl-fluorophosphatase

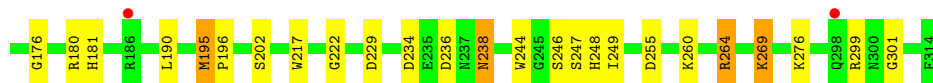
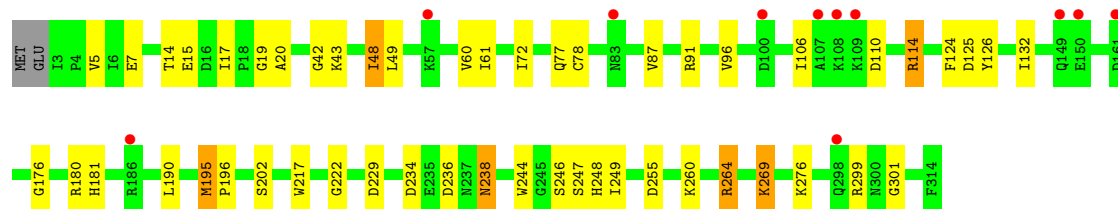
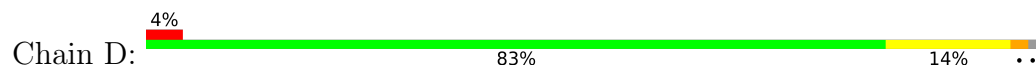


- Molecule 1: Diisopropyl-fluorophosphatase





● Molecule 1: Diisopropyl-fluorophosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.11Å 74.39Å 118.64Å 90.00° 99.90° 90.00°	Depositor
Resolution (Å)	47.77 – 1.80 47.77 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.77-1.80) 99.2 (47.77-1.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 1.79Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.195 , 0.218 0.193 , 0.217	Depositor DCC
R_{free} test set	5332 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	8.7	Xtrriage
Anisotropy	0.362	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10655	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.84 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.8044e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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:
CA

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	0.33	0/2497	0.96	13/3379 (0.4%)
1	B	0.33	0/2497	0.95	11/3379 (0.3%)
1	C	0.33	0/2497	0.95	10/3379 (0.3%)
1	D	0.33	0/2497	0.96	10/3379 (0.3%)
All	All	0.33	0/9988	0.95	44/13516 (0.3%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	249	ILE	N-CA-C	-7.03	98.33	108.17
1	D	238	ASN	N-CA-C	-7.01	99.12	110.20
1	A	222	GLY	N-CA-C	6.93	120.05	112.29
1	A	238	ASN	N-CA-C	-6.55	99.85	110.20
1	D	269	LYS	N-CA-C	6.51	120.83	108.97
1	B	238	ASN	N-CA-C	-6.43	100.04	110.20
1	A	276	LYS	N-CA-C	-6.43	101.68	109.72
1	D	276	LYS	N-CA-C	-6.40	101.72	109.72
1	A	269	LYS	N-CA-C	6.36	120.54	108.97
1	A	125	ASP	N-CA-C	-6.31	101.44	110.59
1	B	276	LYS	N-CA-C	-6.30	101.84	109.72
1	C	238	ASN	N-CA-C	-6.26	100.31	110.20
1	C	269	LYS	N-CA-C	6.22	121.10	110.02
1	B	269	LYS	N-CA-C	6.19	120.24	108.97
1	C	276	LYS	N-CA-C	-6.04	102.17	109.72
1	B	132	ILE	N-CA-C	6.01	117.14	108.48
1	C	132	ILE	N-CA-C	5.98	117.37	108.46
1	B	106	ILE	N-CA-C	5.93	116.71	110.72
1	B	125	ASP	N-CA-C	-5.90	101.73	110.46
1	D	249	ILE	N-CA-C	-5.87	99.29	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	222	GLY	N-CA-C	5.84	120.04	112.68
1	A	132	ILE	N-CA-C	5.84	116.89	108.48
1	D	132	ILE	N-CA-C	5.83	117.15	108.46
1	A	182	MET	N-CA-C	-5.83	103.01	110.53
1	A	195	MET	N-CA-C	5.80	122.62	109.81
1	C	222	GLY	N-CA-C	5.64	119.78	112.68
1	A	165	ILE	N-CA-C	5.59	116.36	108.36
1	D	195	MET	N-CA-C	5.57	122.11	109.81
1	C	125	ASP	N-CA-C	-5.56	102.53	110.59
1	C	195	MET	N-CA-C	5.51	121.98	109.81
1	B	195	MET	N-CA-C	5.35	121.63	109.81
1	D	87	VAL	N-CA-C	5.29	115.52	108.11
1	C	249	ILE	N-CA-C	-5.29	100.14	107.80
1	C	87	VAL	N-CA-C	5.28	115.50	108.11
1	B	165	ILE	N-CA-C	5.25	115.51	108.17
1	B	87	VAL	N-CA-C	5.22	115.42	108.11
1	C	81	ASP	N-CA-C	5.22	119.34	112.92
1	D	125	ASP	N-CA-C	-5.10	103.20	110.59
1	B	222	GLY	N-CA-C	5.09	119.10	112.68
1	A	87	VAL	N-CA-C	5.08	115.23	108.11
1	D	48	ILE	N-CA-C	-5.06	101.09	108.17
1	A	63	LYS	CA-C-N	5.05	124.97	119.76
1	A	63	LYS	C-N-CA	5.05	124.97	119.76
1	A	249	ILE	N-CA-C	-5.05	100.47	107.80

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	2435	0	2359	53	0
1	B	2435	0	2359	61	0
1	C	2435	0	2359	43	0
1	D	2435	0	2358	40	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	208	0	0	1	0
3	B	244	0	0	5	0
3	C	231	0	0	5	0
3	D	224	0	0	2	0
All	All	10655	0	9435	191	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 10.

All (191) 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:17:ILE:HD13	1:B:33:ILE:HD12	1.47	0.97
1:D:264:ARG:HB2	1:D:264:ARG:HH11	1.37	0.89
1:B:17:ILE:HD13	1:B:33:ILE:CD1	2.04	0.86
1:B:126:TYR:H	1:B:181:HIS:CE1	1.92	0.86
1:A:180:ARG:HE	1:A:189:GLN:NE2	1.75	0.84
1:A:126:TYR:H	1:A:181:HIS:CE1	1.99	0.81
1:D:126:TYR:H	1:D:181:HIS:HE1	1.29	0.80
1:A:77:GLN:HE21	1:A:124:PHE:H	1.29	0.79
1:D:77:GLN:HE21	1:D:124:PHE:H	1.31	0.79
1:D:126:TYR:H	1:D:181:HIS:CE1	1.99	0.78
1:C:126:TYR:H	1:C:181:HIS:CE1	2.01	0.78
1:B:41:ASN:HD21	1:D:7:GLU:H	1.31	0.77
1:C:77:GLN:HE21	1:C:124:PHE:H	1.31	0.75
1:C:126:TYR:H	1:C:181:HIS:HE1	1.35	0.75
1:B:77:GLN:HE21	1:B:124:PHE:H	1.33	0.74
1:D:96:VAL:CG2	1:D:106:ILE:HD11	2.18	0.74
1:B:126:TYR:H	1:B:181:HIS:HE1	1.35	0.72
1:A:310:LYS:NZ	1:B:144:ALA:HB3	2.07	0.70
1:A:126:TYR:H	1:A:181:HIS:HE1	1.37	0.70
1:B:96:VAL:CG2	1:B:106:ILE:HD11	2.21	0.69
1:A:151:LYS:N	1:A:151:LYS:HD2	2.08	0.69
1:C:116:MET:HE1	1:C:164:MET:SD	2.35	0.67
1:B:110:ASP:OD2	1:B:114:ARG:HD3	1.94	0.67
1:D:5:VAL:HG11	1:D:264:ARG:HH12	1.60	0.65
1:A:96:VAL:CG2	1:A:106:ILE:HD11	2.27	0.65
1:D:96:VAL:HG21	1:D:106:ILE:HD11	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:VAL:HG21	1:A:106:ILE:HD11	1.80	0.64
1:B:201:TRP:CE3	1:B:214:LYS:HE2	2.33	0.63
1:C:206:LYS:HE3	1:C:212:GLU:OE1	1.98	0.62
1:B:41:ASN:O	1:B:43:LYS:HD2	2.00	0.62
1:A:15:GLU:HB2	3:A:822:HOH:O	2.00	0.62
1:B:33:ILE:HD11	1:B:51:ILE:HD11	1.81	0.61
1:C:30:ASP:OD2	1:C:50:ARG:HD2	2.01	0.61
1:C:96:VAL:HG21	1:C:106:ILE:HD11	1.82	0.61
1:B:96:VAL:HG21	1:B:106:ILE:HD11	1.83	0.60
1:B:80:ARG:CZ	1:B:310:LYS:HB3	2.31	0.60
1:C:51:ILE:HD12	1:C:58:LYS:HG2	1.81	0.60
1:D:96:VAL:HG23	1:D:106:ILE:HD11	1.84	0.59
1:B:41:ASN:ND2	1:D:7:GLU:H	2.01	0.59
1:B:65:GLU:HG2	3:B:670:HOH:O	2.03	0.59
1:A:80:ARG:HB2	1:A:306:CYS:HA	1.86	0.57
1:D:264:ARG:HB2	1:D:264:ARG:NH1	2.14	0.57
1:B:80:ARG:HB2	1:B:306:CYS:HA	1.86	0.57
1:A:151:LYS:HE3	1:A:172:GLN:OE1	2.04	0.57
1:D:15:GLU:C	1:D:17:ILE:HD12	2.30	0.57
1:D:234:ASP:HA	1:D:301:GLY:HA2	1.86	0.57
1:A:17:ILE:HD11	1:A:58:LYS:HD3	1.87	0.56
1:A:48:ILE:C	1:A:49:LEU:HD12	2.30	0.56
1:B:112:GLU:HB2	1:B:114:ARG:HD2	1.86	0.56
1:B:96:VAL:HG23	1:B:106:ILE:HD11	1.87	0.55
1:A:49:LEU:HD12	1:A:49:LEU:N	2.21	0.55
1:A:15:GLU:C	1:A:17:ILE:HD12	2.33	0.54
1:C:42:GLY:C	1:C:43:LYS:HD2	2.32	0.54
1:B:163:GLN:NE2	1:B:165:ILE:HD11	2.22	0.54
1:D:299:ARG:HG2	1:D:299:ARG:HH11	1.72	0.54
1:D:202:SER:HB3	1:D:217:TRP:HB2	1.89	0.54
1:A:72:ILE:HD12	1:A:91:ARG:HG2	1.88	0.54
1:C:15:GLU:C	1:C:17:ILE:HD12	2.34	0.53
1:D:180:ARG:HD2	3:D:716:HOH:O	2.08	0.53
1:A:280:LYS:NZ	1:A:299:ARG:HA	2.23	0.53
1:C:161:ASP:OD1	1:C:163:GLN:HG3	2.09	0.52
1:C:96:VAL:CG2	1:C:106:ILE:HD11	2.39	0.51
1:B:57:LYS:HE3	3:B:633:HOH:O	2.10	0.51
1:D:238:ASN:HD21	1:D:260:LYS:NZ	2.08	0.51
1:B:115:ARG:HE	1:B:138:GLU:CD	2.19	0.51
1:C:72:ILE:HD12	1:C:91:ARG:HG2	1.92	0.51
1:D:244:TRP:CH2	1:D:269:LYS:HE3	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:ARG:HA	1:A:306:CYS:HB2	1.92	0.51
1:C:51:ILE:CD1	1:C:58:LYS:HG2	2.40	0.51
1:A:54:LYS:NZ	1:D:180:ARG:HH22	2.09	0.51
1:C:238:ASN:HD21	1:C:260:LYS:NZ	2.09	0.50
1:D:48:ILE:HG22	1:D:61:ILE:HD11	1.93	0.50
1:A:48:ILE:HG22	1:A:61:ILE:HD11	1.93	0.50
1:B:177:ILE:HD11	1:B:190:LEU:HD11	1.92	0.50
1:D:110:ASP:OD2	1:D:114:ARG:HD3	2.11	0.50
1:C:96:VAL:HG23	1:C:106:ILE:HG12	1.93	0.50
1:C:19:GLY:O	1:C:20:ALA:C	2.55	0.49
1:B:19:GLY:O	1:B:20:ALA:C	2.55	0.49
1:B:72:ILE:HD12	1:B:91:ARG:HG2	1.95	0.49
1:B:81:ASP:OD1	1:B:310:LYS:HD3	2.12	0.49
1:B:234:ASP:HA	1:B:301:GLY:HA2	1.93	0.49
1:D:195:MET:HB3	1:D:196:PRO:CD	2.43	0.49
1:C:176:GLY:HA3	1:C:229:ASP:O	2.12	0.49
1:D:5:VAL:HG11	1:D:264:ARG:NH1	2.26	0.48
1:D:246:SER:O	1:D:247:SER:HB2	2.13	0.48
1:A:112:GLU:HB2	1:A:114:ARG:HD2	1.94	0.48
1:A:232:ASP:HB2	1:A:273:LEU:HD11	1.95	0.48
1:C:17:ILE:HG12	1:C:33:ILE:HD11	1.96	0.47
1:A:246:SER:O	1:A:247:SER:HB2	2.15	0.47
1:D:42:GLY:C	1:D:43:LYS:HD2	2.40	0.47
1:D:49:LEU:CD2	1:D:60:VAL:HG22	2.45	0.47
1:A:5:VAL:HG22	1:A:262:ARG:HB2	1.96	0.47
1:A:151:LYS:N	1:A:151:LYS:CD	2.77	0.47
1:A:51:ILE:CD1	1:A:58:LYS:HG2	2.44	0.47
1:D:49:LEU:HD23	1:D:60:VAL:HA	1.95	0.47
1:A:19:GLY:O	1:A:20:ALA:C	2.57	0.46
1:A:110:ASP:OD2	1:A:114:ARG:HD3	2.16	0.46
1:B:273:LEU:HD12	1:B:273:LEU:C	2.40	0.46
1:C:17:ILE:HG12	1:C:33:ILE:CD1	2.44	0.46
1:C:137:GLY:HA3	1:C:145:THR:CG2	2.45	0.46
1:D:19:GLY:O	1:D:20:ALA:C	2.58	0.46
1:A:46:GLY:HA3	1:A:73:PRO:HD2	1.97	0.46
1:A:96:VAL:HG23	1:A:106:ILE:HD11	1.97	0.46
1:A:280:LYS:HZ2	1:A:299:ARG:HA	1.80	0.46
1:C:28:ASN:ND2	1:C:83:ASN:ND2	2.63	0.46
1:C:43:LYS:HD2	1:C:43:LYS:N	2.31	0.46
1:D:238:ASN:ND2	1:D:260:LYS:NZ	2.64	0.46
1:C:195:MET:HB3	1:C:196:PRO:CD	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:LYS:HD2	1:A:80:ARG:O	2.16	0.46
1:C:96:VAL:HG23	1:C:106:ILE:CG1	2.46	0.46
1:D:176:GLY:HA3	1:D:229:ASP:O	2.14	0.46
1:A:137:GLY:HA3	1:A:145:THR:CG2	2.46	0.46
1:B:246:SER:O	1:B:247:SER:HB2	2.16	0.46
1:D:14:THR:HG23	1:D:17:ILE:HD11	1.98	0.45
1:C:27:LYS:HD2	1:C:80:ARG:O	2.17	0.45
1:B:17:ILE:HD13	1:B:33:ILE:HD11	1.95	0.45
1:B:145:THR:HG22	3:B:504:HOH:O	2.17	0.45
1:B:146:ALA:HB2	3:B:436:HOH:O	2.15	0.45
1:C:240:LEU:HD12	1:C:240:LEU:N	2.32	0.45
1:C:246:SER:O	1:C:247:SER:HB2	2.17	0.45
1:B:137:GLY:HA3	1:B:145:THR:CG2	2.46	0.45
1:C:58:LYS:HB2	1:C:58:LYS:NZ	2.32	0.45
1:A:54:LYS:HD3	1:D:255:ASP:OD2	2.17	0.44
1:C:46:GLY:HA3	1:C:73:PRO:HD2	1.99	0.44
1:C:181:HIS:HD2	3:C:337:HOH:O	2.01	0.44
1:B:46:GLY:HA3	1:B:73:PRO:HD2	1.99	0.44
1:A:51:ILE:HD12	1:A:58:LYS:HG2	2.00	0.44
1:B:51:ILE:HD12	1:B:58:LYS:HG2	2.00	0.44
1:B:77:GLN:NE2	1:B:124:PHE:H	2.08	0.44
1:B:202:SER:HB3	1:B:217:TRP:HB2	1.99	0.44
1:C:300:ASN:HB3	3:C:379:HOH:O	2.18	0.44
1:A:273:LEU:C	1:A:273:LEU:HD12	2.42	0.44
1:B:176:GLY:HA3	1:B:229:ASP:O	2.17	0.44
1:D:72:ILE:HD12	1:D:91:ARG:HG2	2.00	0.44
1:A:176:GLY:HA3	1:A:229:ASP:O	2.17	0.44
1:B:278:GLN:NE2	1:B:278:GLN:HA	2.33	0.44
1:A:72:ILE:HG22	1:A:72:ILE:O	2.17	0.44
1:B:80:ARG:CB	1:B:306:CYS:HA	2.48	0.44
1:A:17:ILE:HG12	1:A:33:ILE:HD11	2.00	0.43
1:B:280:LYS:NZ	1:B:299:ARG:HA	2.33	0.43
1:D:195:MET:HB3	1:D:196:PRO:HD3	2.00	0.43
1:A:234:ASP:HA	1:A:301:GLY:HA2	2.01	0.43
1:B:61:ILE:O	1:B:101:GLY:HA2	2.18	0.43
1:D:181:HIS:HD2	3:D:349:HOH:O	2.02	0.43
1:A:264:ARG:HG2	1:A:264:ARG:HH11	1.84	0.43
1:A:169:THR:O	1:A:170:ALA:HB3	2.19	0.43
1:A:195:MET:HB3	1:A:196:PRO:CD	2.49	0.43
1:B:26:ASP:C	1:B:26:ASP:OD1	2.62	0.43
1:B:163:GLN:HE21	1:B:165:ILE:HD11	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:ARG:CB	1:A:306:CYS:HA	2.49	0.42
1:A:243:ASN:C	1:A:243:ASN:HD22	2.26	0.42
1:A:15:GLU:O	1:A:17:ILE:HD12	2.18	0.42
1:C:108:LYS:HA	3:C:780:HOH:O	2.19	0.42
1:C:169:THR:O	1:C:170:ALA:HB3	2.18	0.42
1:D:77:GLN:NE2	1:D:124:PHE:H	2.08	0.42
1:D:190:LEU:O	1:D:202:SER:HA	2.19	0.42
1:C:190:LEU:O	1:C:202:SER:HA	2.19	0.42
1:B:5:VAL:HG22	1:B:262:ARG:HB2	2.02	0.42
1:B:169:THR:O	1:B:170:ALA:HB3	2.19	0.42
1:C:238:ASN:ND2	1:C:260:LYS:NZ	2.68	0.42
1:B:195:MET:HB3	1:B:196:PRO:CD	2.49	0.42
1:B:264:ARG:HG2	1:B:264:ARG:HH11	1.84	0.42
1:A:96:VAL:HG23	1:A:106:ILE:CG1	2.50	0.42
1:B:125:ASP:OD2	1:B:129:ASN:HB2	2.20	0.42
1:C:215:LYS:HE2	3:C:339:HOH:O	2.20	0.42
1:A:48:ILE:CG2	1:A:61:ILE:HD11	2.50	0.41
1:B:109:LYS:HA	1:B:114:ARG:O	2.19	0.41
1:A:125:ASP:OD2	1:A:129:ASN:HB2	2.20	0.41
1:C:93:GLY:HA3	3:C:779:HOH:O	2.19	0.41
1:C:229:ASP:HB3	1:C:271:SER:O	2.20	0.41
1:B:171:PHE:CZ	1:B:214:LYS:HE3	2.55	0.41
1:C:49:LEU:N	1:C:49:LEU:HD12	2.35	0.41
1:B:96:VAL:HG23	1:B:106:ILE:CG1	2.51	0.41
1:C:273:LEU:HA	1:C:283:PHE:O	2.20	0.41
1:B:106:ILE:HD13	1:B:106:ILE:HA	1.89	0.41
1:C:244:TRP:CZ2	1:C:269:LYS:HD3	2.55	0.41
1:A:297:TRP:CG	1:A:298:GLN:N	2.88	0.41
1:B:72:ILE:HG22	1:B:72:ILE:O	2.18	0.41
1:B:80:ARG:HA	1:B:306:CYS:HB2	2.03	0.41
1:B:277:PRO:HG3	3:B:415:HOH:O	2.20	0.41
1:B:297:TRP:CG	1:B:298:GLN:N	2.89	0.41
1:A:57:LYS:NZ	1:D:236:ASP:OD1	2.49	0.41
1:B:180:ARG:HB3	1:B:189:GLN:HB3	2.02	0.41
1:B:125:ASP:C	1:B:125:ASP:OD1	2.65	0.41
1:D:96:VAL:HG23	1:D:106:ILE:CD1	2.49	0.41
1:D:248:HIS:NE2	1:D:264:ARG:NH1	2.69	0.41
1:C:195:MET:HB3	1:C:196:PRO:HD3	2.03	0.40
1:A:26:ASP:HA	1:A:78:CYS:SG	2.62	0.40
1:A:265:CYS:C	1:A:267:PHE:H	2.28	0.40
1:B:190:LEU:O	1:B:202:SER:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:278:GLN:HA	1:B:278:GLN:HE21	1.86	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	310/314 (99%)	297 (96%)	13 (4%)	0	100	100
1	B	310/314 (99%)	300 (97%)	10 (3%)	0	100	100
1	C	310/314 (99%)	300 (97%)	9 (3%)	1 (0%)	36	25
1	D	310/314 (99%)	298 (96%)	12 (4%)	0	100	100
All	All	1240/1256 (99%)	1195 (96%)	44 (4%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	20	ALA

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	258/260 (99%)	255 (99%)	3 (1%)	63	57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	258/260 (99%)	255 (99%)	3 (1%)	63	57
1	C	258/260 (99%)	256 (99%)	2 (1%)	73	70
1	D	258/260 (99%)	255 (99%)	3 (1%)	63	57
All	All	1032/1040 (99%)	1021 (99%)	11 (1%)	65	60

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LEU
1	A	83	ASN
1	A	114	ARG
1	B	16	ASP
1	B	83	ASN
1	B	114	ARG
1	C	78	CYS
1	C	183	ASN
1	D	78	CYS
1	D	114	ARG
1	D	264	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	77	GLN
1	A	181	HIS
1	A	189	GLN
1	A	237	ASN
1	A	243	ASN
1	A	258	GLN
1	A	278	GLN
1	A	298	GLN
1	A	304	GLN
1	B	28	ASN
1	B	41	ASN
1	B	77	GLN
1	B	163	GLN
1	B	181	HIS
1	B	237	ASN
1	B	238	ASN

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Mol	Chain	Res	Type
1	B	243	ASN
1	B	258	GLN
1	B	278	GLN
1	B	304	GLN
1	C	28	ASN
1	C	77	GLN
1	C	163	GLN
1	C	181	HIS
1	C	238	ASN
1	C	243	ASN
1	C	258	GLN
1	C	278	GLN
1	C	298	GLN
1	C	304	GLN
1	D	28	ASN
1	D	41	ASN
1	D	77	GLN
1	D	98	GLN
1	D	163	GLN
1	D	181	HIS
1	D	238	ASN
1	D	243	ASN
1	D	258	GLN
1	D	278	GLN
1	D	304	GLN

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

Of 8 ligands modelled in this entry, 8 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

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	312/314 (99%)	-0.18	8 (2%) 57 57	2, 9, 21, 30	0
1	B	312/314 (99%)	-0.09	12 (3%) 44 44	2, 9, 24, 33	0
1	C	312/314 (99%)	-0.22	4 (1%) 75 75	2, 8, 21, 27	0
1	D	312/314 (99%)	-0.11	11 (3%) 47 47	2, 9, 22, 30	0
All	All	1248/1256 (99%)	-0.15	35 (2%) 55 55	2, 8, 22, 33	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	101	GLY	5.1
1	C	83	ASN	3.9
1	D	83	ASN	3.8
1	C	100	ASP	3.5
1	B	115	ARG	3.5
1	B	83	ASN	3.4
1	C	188	TYR	3.3
1	D	161	ASP	3.3
1	A	108	LYS	3.2
1	A	100	ASP	3.2
1	B	107	ALA	3.1
1	B	310	LYS	2.9
1	D	100	ASP	2.9
1	B	145	THR	2.9
1	B	109	LYS	2.8
1	D	108	LYS	2.6
1	B	146	ALA	2.6
1	A	107	ALA	2.5
1	D	186	ARG	2.5
1	D	149	GLN	2.3
1	D	107	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	105	GLU	2.3
1	D	298	GLN	2.2
1	B	306	CYS	2.2
1	A	213	ASN	2.2
1	B	161	ASP	2.2
1	D	109	LYS	2.2
1	B	127	GLU	2.1
1	C	105	GLU	2.1
1	A	83	ASN	2.1
1	D	57	LYS	2.1
1	A	161	ASP	2.1
1	B	143	ASP	2.1
1	D	150	GLU	2.0
1	A	306	CYS	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	CA	A	316	1/1	0.99	0.02	2,2,2,2	0
2	CA	A	315	1/1	1.00	0.01	3,3,3,3	0
2	CA	B	315	1/1	1.00	0.01	3,3,3,3	0
2	CA	B	316	1/1	1.00	0.02	2,2,2,2	0
2	CA	C	315	1/1	1.00	0.01	2,2,2,2	0
2	CA	C	316	1/1	1.00	0.01	2,2,2,2	0
2	CA	D	315	1/1	1.00	0.01	2,2,2,2	0
2	CA	D	316	1/1	1.00	0.01	1,1,1,1	0

6.5 Other polymers [i](#)

There are no such residues in this entry.