



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:13 AM GMT

PDB ID : 3HLI  
Title : diisopropyl fluorophosphatase (DFPase), active site mutants  
Authors : Chen, J.C.-H.; Blum, M.-M.  
Deposited on : 2009-05-27  
Resolution : 1.40 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

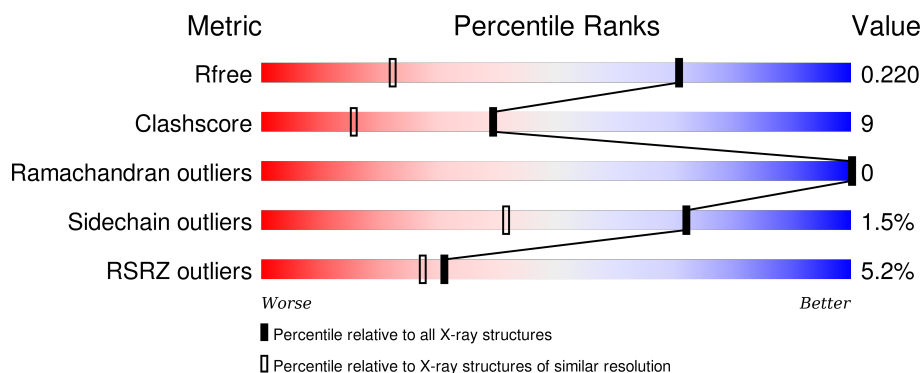
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.40 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))
$R_{free}$	91344	1199 (1.40-1.40)
Clashscore	102246	1295 (1.40-1.40)
Ramachandran outliers	100387	1259 (1.40-1.40)
Sidechain outliers	100360	1258 (1.40-1.40)
RSRZ outliers	91569	1198 (1.40-1.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>4%</div> <div>84%</div> <div>15%</div> </div>
1	B	314	<div> <div>6%</div> <div>80%</div> <div>20%</div> <div>•</div> </div>
1	C	314	<div> <div>4%</div> <div>83%</div> <div>16%</div> <div>•</div> </div>
1	D	314	<div> <div>6%</div> <div>80%</div> <div>19%</div> <div>•</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10803 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
1	A	313	Total	C	N	O	S	0	0	0
			2443	1551	414	462	16			
1	B	312	Total	C	N	O	S	0	0	0
			2438	1548	413	461	16			
1	C	312	Total	C	N	O	S	0	0	0
			2438	1548	413	461	16			
1	D	312	Total	C	N	O	S	0	0	0
			2438	1548	413	461	16			

There are 16 discrepancies between the modelled and reference sequences:

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

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

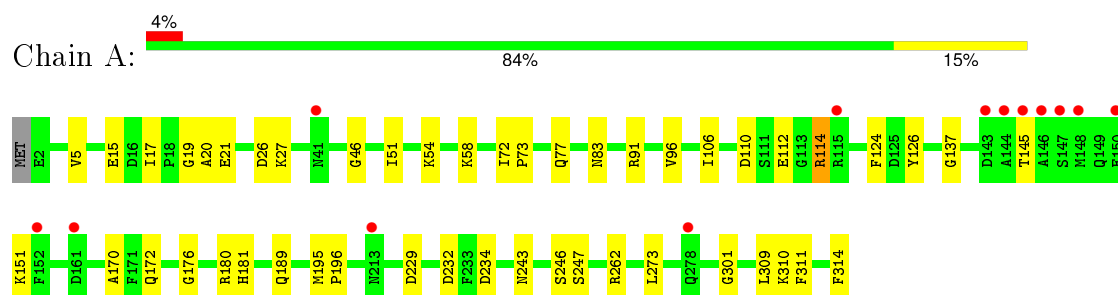
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	250	Total O 250 250	0	0
3	B	273	Total O 273 273	0	0
3	C	270	Total O 270 270	0	0
3	D	245	Total O 245 245	0	0

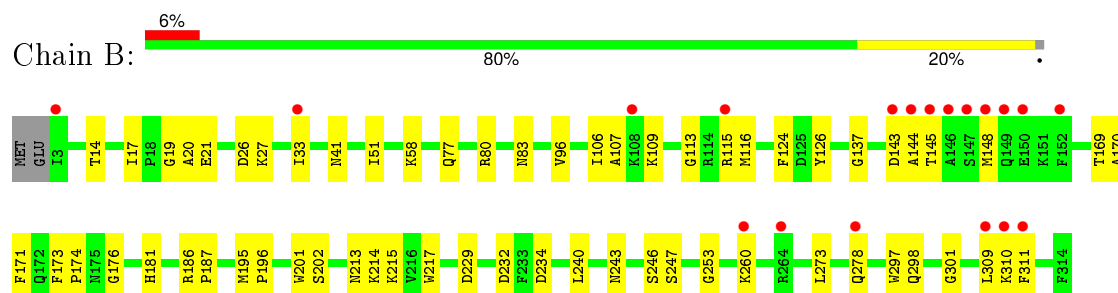
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

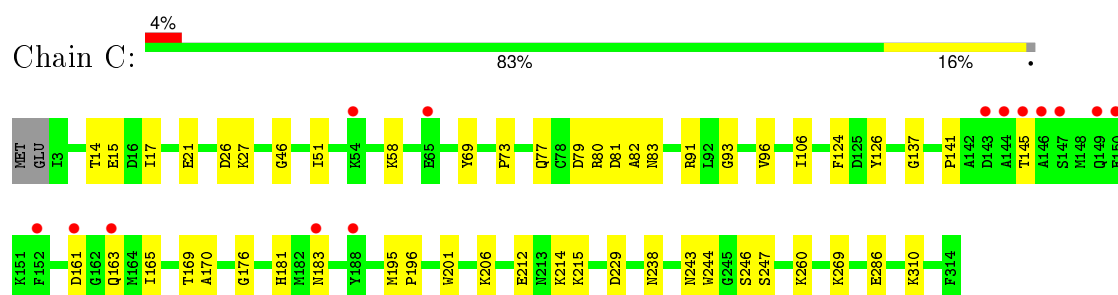
#### • Molecule 1: Diisopropyl-fluorophosphatase



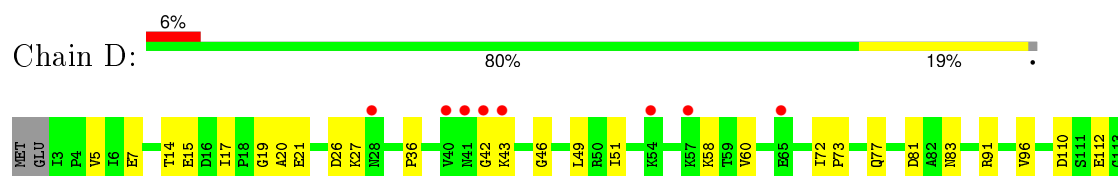
#### • Molecule 1: Diisopropyl-fluorophosphatase

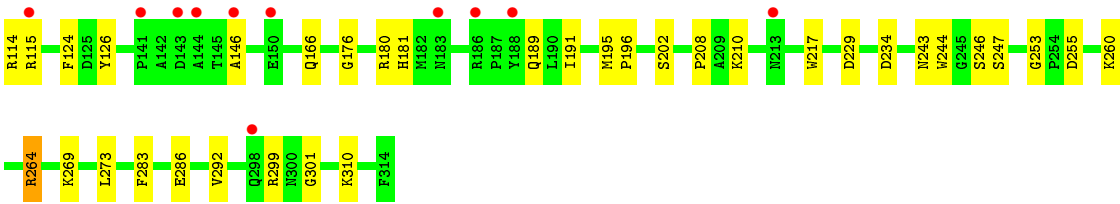


#### • Molecule 1: Diisopropyl-fluorophosphatase



#### • Molecule 1: Diisopropyl-fluorophosphatase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.95Å 74.49Å 118.95Å 90.00° 100.00° 90.00°	Depositor
Resolution (Å)	24.82 – 1.40 24.81 – 1.40	Depositor EDS
% Data completeness (in resolution range)	94.5 (24.82-1.40) 94.4 (24.81-1.40)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.85 (at 1.40Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.205 , 0.221 0.203 , 0.220	Depositor DCC
$R_{free}$ test set	10622 reflections (4.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	8.9	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.43 , 49.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 213602 reflections (0.000%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10803	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.49 % 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 5.6378e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.29	0/2505	0.66	0/3390
1	B	0.30	0/2500	0.66	0/3383
1	C	0.30	0/2500	0.65	0/3383
1	D	0.30	0/2500	0.66	0/3383
All	All	0.30	0/10005	0.66	0/13539

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	2443	0	2362	40	0
1	B	2438	0	2360	50	0
1	C	2438	0	2360	37	0
1	D	2438	0	2360	52	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	250	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	273	0	0	3	0
3	C	270	0	0	5	0
3	D	245	0	0	3	0
All	All	10803	0	9442	170	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 9.

All (170) 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:D:264:ARG:HB2	1:D:264:ARG:HH11	1.33	0.93
1:A:17:ILE:HD11	1:A:58:LYS:HD2	1.56	0.86
1:A:26:ASP:HB2	1:A:83:ASN:HD21	1.42	0.84
1:B:17:ILE:HD13	1:B:33:ILE:HD12	1.60	0.82
1:D:126:TYR:H	1:D:181:HIS:CE1	1.99	0.80
1:A:126:TYR:H	1:A:181:HIS:CE1	2.01	0.78
1:D:126:TYR:H	1:D:181:HIS:HE1	1.30	0.78
1:B:126:TYR:H	1:B:181:HIS:CE1	2.02	0.78
1:C:126:TYR:H	1:C:181:HIS:CE1	2.01	0.77
1:D:5:VAL:HG11	1:D:264:ARG:HH12	1.49	0.77
1:B:41:ASN:HD21	1:D:7:GLU:H	1.34	0.76
1:C:77:GLN:HE21	1:C:124:PHE:H	1.33	0.76
1:A:180:ARG:HE	1:A:189:GLN:NE2	1.82	0.76
1:C:126:TYR:H	1:C:181:HIS:HE1	1.33	0.75
1:A:77:GLN:HE21	1:A:124:PHE:H	1.33	0.75
1:A:310:LYS:NZ	1:B:144:ALA:HB3	2.03	0.74
1:D:26:ASP:HB2	1:D:83:ASN:HD21	1.53	0.74
1:B:17:ILE:HD13	1:B:33:ILE:CD1	2.17	0.74
1:B:26:ASP:HB2	1:B:83:ASN:HD21	1.54	0.72
1:A:17:ILE:HD11	1:A:58:LYS:CD	2.20	0.71
1:C:96:VAL:CG2	1:C:106:ILE:HD11	2.22	0.70
1:B:126:TYR:H	1:B:181:HIS:HE1	1.39	0.70
1:A:126:TYR:H	1:A:181:HIS:HE1	1.38	0.69
1:D:77:GLN:HE21	1:D:124:PHE:H	1.40	0.69
1:C:26:ASP:HB2	1:C:83:ASN:HD21	1.57	0.68
1:B:77:GLN:HE21	1:B:124:PHE:H	1.42	0.67
1:C:96:VAL:HG21	1:C:106:ILE:HD11	1.76	0.67
1:B:310:LYS:HZ2	1:B:311:PHE:HE2	1.41	0.66
1:D:112:GLU:HB2	1:D:114:ARG:HD2	1.79	0.64
1:D:43:LYS:HD2	1:D:43:LYS:N	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ILE:CD1	1:A:58:LYS:HD2	2.25	0.63
1:D:264:ARG:NH1	1:D:264:ARG:HB2	2.09	0.62
1:B:96:VAL:CG2	1:B:106:ILE:HD11	2.31	0.61
1:B:41:ASN:ND2	1:D:7:GLU:H	1.97	0.61
1:B:253:GLY:CA	1:B:260:LYS:HE3	2.32	0.59
1:B:311:PHE:CE1	1:C:91:ARG:HD3	2.37	0.59
1:A:112:GLU:HB2	1:A:114:ARG:HD2	1.85	0.58
1:A:151:LYS:HE3	1:A:172:GLN:OE1	2.04	0.58
1:A:110:ASP:OD2	1:A:114:ARG:HD3	2.04	0.58
1:B:96:VAL:HG23	1:B:106:ILE:HD11	1.87	0.57
1:A:151:LYS:HD2	1:A:151:LYS:N	2.19	0.57
1:A:54:LYS:NZ	1:D:180:ARG:HH22	2.03	0.56
1:D:202:SER:HB3	1:D:217:TRP:HB2	1.87	0.56
1:B:27:LYS:H	1:B:83:ASN:ND2	2.03	0.56
1:D:5:VAL:HG11	1:D:264:ARG:NH1	2.19	0.56
1:C:51:ILE:HD12	1:C:58:LYS:HG2	1.86	0.56
1:C:80:ARG:NH2	1:C:310:LYS:HG3	2.20	0.56
1:A:96:VAL:CG2	1:A:106:ILE:HD11	2.36	0.56
1:B:51:ILE:HD12	1:B:58:LYS:HG2	1.88	0.55
1:C:81:ASP:OD1	1:C:310:LYS:HD2	2.07	0.54
1:B:201:TRP:CE3	1:B:214:LYS:HE2	2.43	0.53
1:A:15:GLU:HB2	3:A:803:HOH:O	2.07	0.53
1:C:58:LYS:HB2	1:C:58:LYS:NZ	2.24	0.53
1:C:163:GLN:HE21	1:C:165:ILE:HD11	1.73	0.53
1:D:27:LYS:H	1:D:83:ASN:ND2	2.05	0.53
1:C:93:GLY:HA3	3:C:779:HOH:O	2.09	0.52
1:B:202:SER:HB3	1:B:217:TRP:HB2	1.92	0.52
1:C:215:LYS:HG2	3:C:340:HOH:O	2.09	0.52
1:C:15:GLU:C	1:C:17:ILE:HD12	2.30	0.52
1:C:201:TRP:CE3	1:C:214:LYS:HE2	2.45	0.52
1:A:26:ASP:HB2	1:A:83:ASN:ND2	2.20	0.52
1:C:96:VAL:HG23	1:C:106:ILE:HD11	1.91	0.52
1:B:213:ASN:OD1	1:B:215:LYS:HE3	2.10	0.51
1:C:51:ILE:CD1	1:C:58:LYS:HG2	2.40	0.51
1:A:5:VAL:HG22	1:A:262:ARG:HB2	1.91	0.51
1:A:96:VAL:HG21	1:A:106:ILE:HD11	1.94	0.50
1:B:27:LYS:H	1:B:83:ASN:HD21	1.60	0.50
1:D:114:ARG:NH1	1:D:166:GLN:OE1	2.45	0.49
1:D:189:GLN:HE21	1:D:191:ILE:HD11	1.77	0.49
1:A:180:ARG:HE	1:A:189:GLN:HE21	1.59	0.49
1:D:72:ILE:HD12	1:D:91:ARG:HG2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:246:SER:O	1:C:247:SER:HB2	2.13	0.49
1:B:33:ILE:HD11	1:B:51:ILE:HD11	1.93	0.49
1:A:54:LYS:HZ1	1:D:180:ARG:HH22	1.59	0.49
1:A:51:ILE:HD12	1:A:58:LYS:HG2	1.94	0.49
1:B:144:ALA:HA	3:B:504:HOH:O	2.12	0.49
1:C:215:LYS:HE2	3:C:340:HOH:O	2.11	0.49
1:D:15:GLU:C	1:D:17:ILE:HD12	2.33	0.49
1:B:310:LYS:NZ	1:B:311:PHE:HE2	2.09	0.49
1:B:253:GLY:HA3	1:B:260:LYS:HE3	1.94	0.48
1:D:27:LYS:H	1:D:83:ASN:HD21	1.61	0.48
1:C:27:LYS:H	1:C:83:ASN:ND2	2.11	0.48
1:A:46:GLY:HA3	1:A:73:PRO:HD2	1.95	0.48
1:A:54:LYS:HD3	1:D:255:ASP:OD2	2.14	0.48
1:B:26:ASP:HB2	1:B:83:ASN:ND2	2.27	0.48
1:D:180:ARG:HB3	1:D:189:GLN:HG2	1.95	0.48
1:C:137:GLY:HA3	1:C:145:THR:HG23	1.96	0.48
1:A:309:LEU:HA	3:A:678:HOH:O	2.14	0.47
1:D:110:ASP:OD2	1:D:114:ARG:HD3	2.15	0.47
1:B:148:MET:HG2	1:B:173:PHE:CD2	2.49	0.47
1:A:246:SER:O	1:A:247:SER:HB2	2.15	0.47
1:D:310:LYS:H	1:D:310:LYS:HG2	1.52	0.47
1:B:51:ILE:CD1	1:B:58:LYS:HG2	2.44	0.46
1:B:309:LEU:HB3	3:B:366:HOH:O	2.15	0.46
1:D:26:ASP:HB2	1:D:83:ASN:ND2	2.27	0.46
1:B:246:SER:O	1:B:247:SER:HB2	2.15	0.46
1:D:181:HIS:HD2	3:D:349:HOH:O	1.99	0.46
1:A:311:PHE:HE1	1:B:144:ALA:CB	2.29	0.46
1:A:310:LYS:HZ1	1:B:144:ALA:HB3	1.78	0.45
1:C:161:ASP:OD2	1:C:163:GLN:HG3	2.16	0.45
1:D:246:SER:O	1:D:247:SER:HB2	2.16	0.45
1:D:51:ILE:HD12	1:D:58:LYS:HG2	1.97	0.45
1:D:234:ASP:HA	1:D:301:GLY:HA2	1.98	0.45
1:D:180:ARG:HD2	3:D:716:HOH:O	2.17	0.45
1:C:69:TYR:CE2	1:C:141:PRO:HB3	2.52	0.45
1:B:169:THR:O	1:B:170:ALA:HB3	2.16	0.45
1:B:176:GLY:HA3	1:B:229:ASP:O	2.17	0.45
1:D:112:GLU:CB	1:D:114:ARG:HD2	2.47	0.44
1:D:46:GLY:HA3	1:D:73:PRO:HD2	2.00	0.44
1:B:14:THR:HG23	1:B:17:ILE:HD11	2.00	0.44
1:D:14:THR:HG23	1:D:17:ILE:HD11	2.00	0.44
1:A:176:GLY:HA3	1:A:229:ASP:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:238:ASN:HD21	1:C:260:LYS:NZ	2.15	0.44
1:C:79:ASP:HB3	1:C:82:ALA:O	2.18	0.44
1:D:195:MET:HB3	1:D:196:PRO:CD	2.48	0.44
1:B:234:ASP:HA	1:B:301:GLY:HA2	1.99	0.44
1:A:151:LYS:O	1:A:170:ALA:HA	2.18	0.43
1:B:96:VAL:HG21	1:B:106:ILE:HD11	2.00	0.43
1:C:195:MET:HB3	1:C:196:PRO:CD	2.48	0.43
1:A:72:ILE:HD12	1:A:91:ARG:HG2	1.99	0.43
1:B:171:PHE:HB3	1:B:174:PRO:HD3	1.99	0.43
1:D:176:GLY:HA3	1:D:229:ASP:O	2.17	0.43
1:B:240:LEU:N	1:B:240:LEU:HD12	2.33	0.43
1:D:42:GLY:C	1:D:43:LYS:HD2	2.39	0.43
1:A:27:LYS:H	1:A:83:ASN:ND2	2.16	0.43
1:C:181:HIS:HD2	3:C:338:HOH:O	2.02	0.43
1:C:244:TRP:CZ2	1:C:269:LYS:HD3	2.54	0.43
1:B:297:TRP:CG	1:B:298:GLN:N	2.87	0.43
1:A:96:VAL:HG23	1:A:106:ILE:HD11	2.00	0.43
1:A:314:PHE:HA	3:A:1055:HOH:O	2.19	0.42
1:C:46:GLY:HA3	1:C:73:PRO:HD2	2.01	0.42
1:D:244:TRP:CH2	1:D:269:LYS:HE3	2.54	0.42
1:D:253:GLY:CA	1:D:260:LYS:HE3	2.50	0.42
1:C:15:GLU:HB2	3:C:376:HOH:O	2.19	0.42
1:A:195:MET:HB3	1:A:196:PRO:CD	2.50	0.42
1:B:201:TRP:CZ3	1:B:214:LYS:HE2	2.55	0.42
1:D:81:ASP:OD1	1:D:310:LYS:HD2	2.19	0.42
1:D:299:ARG:HG2	1:D:299:ARG:HH11	1.84	0.42
1:B:181:HIS:HD2	3:B:392:HOH:O	2.03	0.42
1:D:180:ARG:HB3	1:D:189:GLN:CG	2.50	0.42
1:C:14:THR:HG23	1:C:17:ILE:HD11	2.02	0.42
1:A:234:ASP:HA	1:A:301:GLY:HA2	2.02	0.42
1:D:49:LEU:CD2	1:D:60:VAL:HG22	2.50	0.42
1:D:208:PRO:O	1:D:210:LYS:HG3	2.20	0.42
1:C:169:THR:O	1:C:170:ALA:HB3	2.20	0.42
1:D:96:VAL:O	1:D:96:VAL:HG13	2.20	0.42
1:D:43:LYS:CD	1:D:43:LYS:N	2.83	0.41
1:B:137:GLY:HA3	1:B:145:THR:HG22	2.01	0.41
1:B:186:ARG:HA	1:B:187:PRO:HD3	1.97	0.41
1:C:206:LYS:HE3	1:C:212:GLU:OE1	2.19	0.41
1:A:181:HIS:HD2	3:A:348:HOH:O	2.02	0.41
1:C:176:GLY:HA3	1:C:229:ASP:O	2.20	0.41
1:B:195:MET:HB3	1:B:196:PRO:CD	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:TRP:CZ3	1:C:214:LYS:HE2	2.56	0.41
1:A:19:GLY:O	1:A:20:ALA:C	2.59	0.41
1:B:107:ALA:O	1:B:116:MET:HG3	2.21	0.41
1:B:278:GLN:HA	1:B:278:GLN:HE21	1.86	0.41
1:C:77:GLN:NE2	1:C:124:PHE:H	2.11	0.40
1:B:80:ARG:CZ	1:B:310:LYS:HB3	2.51	0.40
1:D:36:PRO:HB3	1:D:73:PRO:O	2.21	0.40
1:B:109:LYS:HD3	1:B:113:GLY:O	2.21	0.40
1:D:19:GLY:O	1:D:20:ALA:C	2.59	0.40
1:D:110:ASP:OD1	1:D:114:ARG:HD3	2.21	0.40
1:A:137:GLY:HA3	1:A:145:THR:CG2	2.52	0.40
1:B:19:GLY:O	1:B:20:ALA:C	2.60	0.40
1:A:232:ASP:HB2	1:A:273:LEU:CD1	2.51	0.40
1:D:273:LEU:HA	1:D:283:PHE:O	2.22	0.40
1:D:146:ALA:HB3	3:D:1025:HOH:O	2.21	0.40
1:D:14:THR:HG22	1:D:292:VAL:HB	2.03	0.40
1:B:232:ASP:HB2	1:B:273:LEU:HD11	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	311/314 (99%)	301 (97%)	10 (3%)	0	100	100
1	B	310/314 (99%)	297 (96%)	13 (4%)	0	100	100
1	C	310/314 (99%)	299 (96%)	11 (4%)	0	100	100
1	D	310/314 (99%)	298 (96%)	12 (4%)	0	100	100
All	All	1241/1256 (99%)	1195 (96%)	46 (4%)	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	259/261 (99%)	256 (99%)	3 (1%)	78	52
1	B	259/261 (99%)	255 (98%)	4 (2%)	72	41
1	C	259/261 (99%)	255 (98%)	4 (2%)	72	41
1	D	259/261 (99%)	254 (98%)	5 (2%)	65	29
All	All	1036/1044 (99%)	1020 (98%)	16 (2%)	72	41

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	GLU
1	A	114	ARG
1	A	243	ASN
1	B	21	GLU
1	B	115	ARG
1	B	143	ASP
1	B	243	ASN
1	C	21	GLU
1	C	183	ASN
1	C	243	ASN
1	C	286	GLU
1	D	21	GLU
1	D	115	ARG
1	D	243	ASN
1	D	264	ARG
1	D	286	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	83	ASN
1	A	98	GLN

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Mol	Chain	Res	Type
1	A	181	HIS
1	A	189	GLN
1	A	238	ASN
1	A	243	ASN
1	A	258	GLN
1	A	278	GLN
1	A	304	GLN
1	B	41	ASN
1	B	77	GLN
1	B	83	ASN
1	B	98	GLN
1	B	181	HIS
1	B	238	ASN
1	B	243	ASN
1	B	258	GLN
1	B	278	GLN
1	B	304	GLN
1	C	77	GLN
1	C	83	ASN
1	C	98	GLN
1	C	163	GLN
1	C	181	HIS
1	C	238	ASN
1	C	243	ASN
1	C	258	GLN
1	C	289	ASN
1	C	304	GLN
1	D	41	ASN
1	D	77	GLN
1	D	83	ASN
1	D	98	GLN
1	D	181	HIS
1	D	189	GLN
1	D	238	ASN
1	D	243	ASN
1	D	258	GLN
1	D	278	GLN
1	D	304	GLN

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	313/314 (99%)	0.38	13 (4%) 40 38	4, 10, 21, 32	0
1	B	312/314 (99%)	0.51	19 (6%) 25 22	4, 10, 22, 33	0
1	C	312/314 (99%)	0.36	14 (4%) 37 34	4, 9, 20, 30	0
1	D	312/314 (99%)	0.43	19 (6%) 25 22	4, 9, 21, 30	0
All	All	1249/1256 (99%)	0.42	65 (5%) 31 27	4, 10, 21, 33	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	146	ALA	10.4
1	B	144	ALA	10.0
1	C	144	ALA	9.8
1	B	152	PHE	6.1
1	B	145	THR	5.4
1	B	148	MET	5.3
1	A	144	ALA	5.2
1	B	143	ASP	5.2
1	A	145	THR	5.1
1	D	144	ALA	5.0
1	D	41	ASN	4.7
1	B	147	SER	4.5
1	B	150	GLU	4.4
1	C	54	LYS	4.4
1	A	146	ALA	4.4
1	C	146	ALA	3.8
1	C	150	GLU	3.8
1	A	143	ASP	3.6
1	D	298	GLN	3.6
1	D	188	TYR	3.5
1	A	152	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	145	THR	3.3
1	B	115	ARG	3.3
1	D	54	LYS	3.2
1	B	311	PHE	3.1
1	C	143	ASP	3.1
1	B	310	LYS	3.1
1	D	115	ARG	3.0
1	A	150	GLU	3.0
1	D	40	VAL	3.0
1	D	42	GLY	2.9
1	D	57	LYS	2.9
1	D	186	ARG	2.9
1	C	161	ASP	2.8
1	C	152	PHE	2.8
1	A	148	MET	2.8
1	C	149	GLN	2.7
1	D	150	GLU	2.7
1	B	149	GLN	2.6
1	C	147	SER	2.6
1	C	183	ASN	2.5
1	A	147	SER	2.5
1	D	143	ASP	2.4
1	A	41	ASN	2.4
1	D	213	ASN	2.4
1	C	188	TYR	2.3
1	A	213	ASN	2.3
1	D	146	ALA	2.3
1	A	161	ASP	2.3
1	B	33	ILE	2.2
1	C	65	GLU	2.2
1	D	28	ASN	2.2
1	D	43	LYS	2.2
1	C	163	GLN	2.1
1	A	278	GLN	2.1
1	B	309	LEU	2.1
1	B	260	LYS	2.1
1	A	115	ARG	2.1
1	B	264	ARG	2.1
1	B	3	ILE	2.0
1	D	65	GLU	2.0
1	B	278	GLN	2.0
1	B	108	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	141	PRO	2.0
1	D	183	ASN	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CA	A	316	1/1	1.00	0.06	-2.38	4,4,4,4	0
2	CA	B	315	1/1	1.00	0.07	-2.95	3,3,3,3	0
2	CA	D	315	1/1	1.00	0.06	-3.08	3,3,3,3	0
2	CA	C	315	1/1	1.00	0.07	-3.33	4,4,4,4	0
2	CA	A	315	1/1	1.00	0.07	-3.39	4,4,4,4	0
2	CA	D	316	1/1	1.00	0.06	-3.45	4,4,4,4	0
2	CA	B	316	1/1	1.00	0.04	-4.43	3,3,3,3	0
2	CA	C	316	1/1	1.00	0.05	-4.88	4,4,4,4	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.