



# Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:16 AM GMT

PDB ID : 1O12  
Title : Crystal structure of N-acetylglucosamine-6-phosphatedeacetylase (TM0814)  
from *Thermotoga maritima* at 2.5 Å resolution  
Authors : Joint Center for Structural Genomics (JCSG)  
Deposited on : 2002-10-15  
Resolution : 2.50 Å (reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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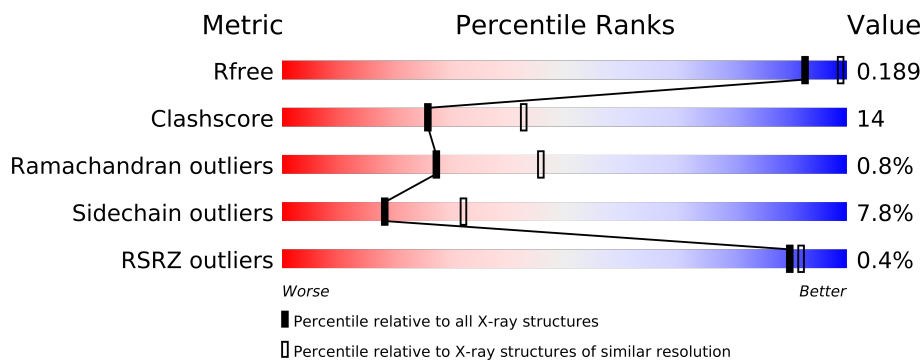
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

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))
$R_{free}$	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	376	
1	B	376	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5828 atoms, of which 0 are hydrogen and 0 are deuterium.

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 N-acetylglucosamine-6-phosphatedeacetylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	Se	0	0	1
			2810	1784	473	539	5	9			
1	B	364	Total	C	N	O	S	Se	0	0	1
			2810	1784	473	539	5	9			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MSE	-	LEADER SEQUENCE	UNP Q9WZS1
A	-10	GLY	-	LEADER SEQUENCE	UNP Q9WZS1
A	-9	SER	-	LEADER SEQUENCE	UNP Q9WZS1
A	-8	ASP	-	LEADER SEQUENCE	UNP Q9WZS1
A	-7	LYS	-	LEADER SEQUENCE	UNP Q9WZS1
A	-6	ILE	-	LEADER SEQUENCE	UNP Q9WZS1
A	-5	HIS	-	EXPRESSION TAG	UNP Q9WZS1
A	-4	HIS	-	EXPRESSION TAG	UNP Q9WZS1
A	-3	HIS	-	EXPRESSION TAG	UNP Q9WZS1
A	-2	HIS	-	EXPRESSION TAG	UNP Q9WZS1
A	-1	HIS	-	EXPRESSION TAG	UNP Q9WZS1
A	0	HIS	-	EXPRESSION TAG	UNP Q9WZS1
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	42	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	59	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	66	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	90	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	149	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	185	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	237	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
A	352	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	-11	MSE	-	LEADER SEQUENCE	UNP Q9WZS1
B	-10	GLY	-	LEADER SEQUENCE	UNP Q9WZS1
B	-9	SER	-	LEADER SEQUENCE	UNP Q9WZS1
B	-8	ASP	-	LEADER SEQUENCE	UNP Q9WZS1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	LYS	-	LEADER SEQUENCE	UNP Q9WZS1
B	-6	ILE	-	LEADER SEQUENCE	UNP Q9WZS1
B	-5	HIS	-	EXPRESSION TAG	UNP Q9WZS1
B	-4	HIS	-	EXPRESSION TAG	UNP Q9WZS1
B	-3	HIS	-	EXPRESSION TAG	UNP Q9WZS1
B	-2	HIS	-	EXPRESSION TAG	UNP Q9WZS1
B	-1	HIS	-	EXPRESSION TAG	UNP Q9WZS1
B	0	HIS	-	EXPRESSION TAG	UNP Q9WZS1
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	42	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	59	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	66	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	90	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	149	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	185	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	237	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1
B	352	MSE	MET	MODIFIED RESIDUE	UNP Q9WZS1

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Fe 1 1	0	0
2	A	1	Total Fe 1 1	0	0

- Molecule 3 is water.

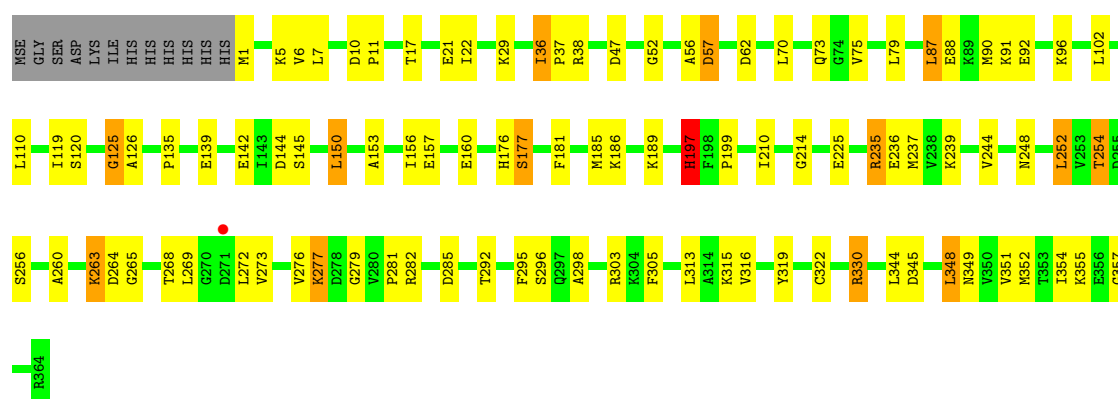
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	91	Total O 91 91	0	0
3	B	115	Total O 115 115	0	0

### 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.

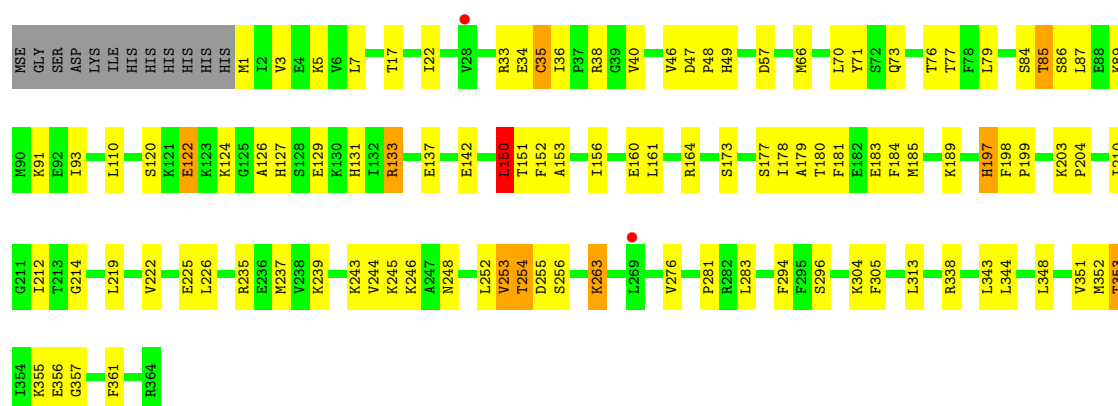
- Molecule 1: N-acetylglucosamine-6-phosphatideacetylase

Chain A:



- Molecule 1: N-acetylglucosamine-6-phosphatideacetylase

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.07Å 85.07Å 206.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.53 – 2.50 42.54 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.53-2.50) 99.9 (42.54-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.78 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.195 , 0.251 0.199 , 0.189	Depositor DCC
$R_{free}$ test set	1516 reflections (5.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.8	Xtriage
Anisotropy	0.420	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 31.9	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30653 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5828	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE

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	0.60	0/2845	0.76	1/3822 (0.0%)
1	B	0.60	0/2845	0.77	1/3822 (0.0%)
All	All	0.60	0/5690	0.76	2/7644 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	150	LEU	CA-CB-CG	6.39	129.99	115.30
1	A	150	LEU	CA-CB-CG	5.93	128.93	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2810	0	2885	76	0
1	B	2810	0	2885	79	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	91	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	115	0	0	3	0
All	All	5828	0	5770	154	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (154) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:263:LYS:H	1:B:263:LYS:HE3	1.05	1.07
1:B:185:MSE:HE1	1:B:219:LEU:HD11	1.14	1.06
1:B:254:THR:HG23	1:B:256:SER:H	1.22	1.02
1:A:17:THR:HG22	1:A:37:PRO:HG3	1.46	0.94
1:B:71:TYR:O	1:B:353:THR:HG21	1.70	0.91
1:A:352:MSE:HE3	1:A:354:ILE:HD11	1.52	0.90
1:A:254:THR:HG23	1:A:256:SER:H	1.34	0.90
1:B:263:LYS:N	1:B:263:LYS:HE3	1.89	0.88
1:B:263:LYS:CE	1:B:263:LYS:H	1.90	0.85
1:B:122:GLU:HG3	1:B:178:ILE:HD13	1.60	0.82
1:B:254:THR:HG23	1:B:256:SER:N	1.98	0.78
1:A:17:THR:CG2	1:A:37:PRO:HG3	2.15	0.76
1:B:353:THR:HG23	1:B:361:PHE:HB3	1.69	0.74
1:A:352:MSE:CE	1:A:354:ILE:HD11	2.18	0.73
1:A:276:VAL:HG22	1:A:281:PRO:HB3	1.71	0.73
1:B:225:GLU:O	1:B:226:LEU:HD23	1.90	0.71
1:B:344:LEU:N	1:B:344:LEU:HD12	2.05	0.71
1:B:185:MSE:HE1	1:B:219:LEU:CD1	2.08	0.71
1:B:57:ASP:OD1	3:B:411:HOH:O	2.08	0.71
1:A:197:HIS:H	1:A:225:GLU:CG	2.04	0.70
1:A:303:ARG:HB2	1:A:313:LEU:HD11	1.73	0.70
1:B:5:LYS:HG3	1:B:33:ARG:HH22	1.58	0.68
1:A:254:THR:HG23	1:A:256:SER:N	2.07	0.68
1:B:254:THR:CG2	1:B:256:SER:H	2.04	0.68
1:A:252:LEU:HD22	1:A:316:VAL:HG23	1.75	0.67
1:B:239:LYS:HE2	1:B:305:PHE:O	1.94	0.66
1:B:34:GLU:O	1:B:35:CYS:HB3	1.95	0.66
1:B:161:LEU:HD22	1:B:164:ARG:NH2	2.09	0.66
1:B:179:ALA:HB3	1:B:212:ILE:HD11	1.76	0.66
1:A:254:THR:HG23	1:A:256:SER:CB	2.26	0.66
1:B:126:ALA:HA	1:B:283:LEU:HD11	1.78	0.65
1:B:338:ARG:HD3	1:B:356:GLU:OE2	1.97	0.65
1:A:345:ASP:OD2	1:A:349:ASN:HB2	1.98	0.64
1:A:47:ASP:OD2	1:A:254:THR:HB	1.96	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:73:GLN:OE1	1:B:296:SER:HB2	1.99	0.63
1:A:254:THR:CG2	1:A:256:SER:H	2.10	0.63
1:B:33:ARG:HH11	1:B:33:ARG:HG2	1.66	0.61
1:A:239:LYS:HE2	1:A:305:PHE:O	2.00	0.61
1:A:277:LYS:HA	1:A:277:LYS:HE3	1.83	0.61
1:A:5:LYS:NZ	1:A:5:LYS:HB2	2.16	0.60
1:B:49:HIS:HB3	1:B:253:VAL:HG21	1.84	0.60
1:A:125:GLY:O	1:A:126:ALA:HB3	2.02	0.59
1:B:124:LYS:HE2	1:B:129:GLU:HB2	1.85	0.59
1:A:197:HIS:N	1:A:225:GLU:HG2	2.19	0.58
1:A:263:LYS:N	1:A:263:LYS:HD2	2.19	0.57
1:B:76:THR:O	1:B:355:LYS:HE2	2.05	0.57
1:B:36:ILE:HD12	1:B:36:ILE:O	2.06	0.56
1:B:47:ASP:OD2	1:B:254:THR:HB	2.06	0.56
1:A:17:THR:HG22	1:A:37:PRO:CG	2.28	0.55
1:B:49:HIS:HB3	1:B:253:VAL:CG2	2.37	0.55
1:A:153:ALA:HB1	1:A:177:SER:HB2	1.89	0.55
1:A:268:THR:HG22	1:A:273:VAL:HA	1.87	0.55
1:A:57:ASP:HB3	3:A:443:HOH:O	2.06	0.55
1:B:254:THR:HG23	1:B:256:SER:CB	2.37	0.54
1:A:73:GLN:OE1	1:A:296:SER:CB	2.55	0.54
1:B:160:GLU:H	1:B:160:GLU:CD	2.10	0.54
1:A:197:HIS:H	1:A:225:GLU:HG2	1.71	0.54
1:A:87:LEU:CD1	1:A:139:GLU:HG2	2.37	0.54
1:A:102:LEU:O	1:A:102:LEU:HD13	2.06	0.54
1:B:34:GLU:O	1:B:35:CYS:CB	2.56	0.53
1:B:203:LYS:HD3	1:B:204:PRO:HD2	1.91	0.53
1:A:153:ALA:CB	1:A:177:SER:HB2	2.39	0.52
1:A:197:HIS:N	1:A:225:GLU:CG	2.72	0.51
1:A:185:MSE:O	1:A:189:LYS:HG3	2.09	0.51
1:B:84:SER:H	1:B:127:HIS:CE1	2.29	0.51
1:B:91:LYS:HD3	1:B:142:GLU:CD	2.31	0.51
1:A:315:LYS:HA	1:A:319:TYR:HB3	1.92	0.51
1:B:71:TYR:O	1:B:353:THR:CG2	2.50	0.51
1:A:263:LYS:O	1:A:265:GLY:N	2.45	0.50
1:A:87:LEU:HG	1:A:139:GLU:HG2	1.93	0.50
1:B:48:PRO:HD2	1:B:253:VAL:HG23	1.93	0.50
1:B:33:ARG:NH1	1:B:33:ARG:HG2	2.26	0.50
1:B:181:PHE:CE1	1:B:210:ILE:HD11	2.46	0.50
1:B:181:PHE:CD1	1:B:210:ILE:HD11	2.47	0.49
1:B:17:THR:HG22	1:B:33:ARG:O	2.11	0.49
1:A:322:CYS:SG	1:A:330:ARG:HB3	2.53	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:199:PRO:HG3	1:B:237:MSE:HE2	1.95	0.49
1:A:244:VAL:HG22	1:B:244:VAL:HG22	1.95	0.48
1:A:91:LYS:HD2	1:A:142:GLU:OE2	2.13	0.48
1:B:40:VAL:O	1:B:343:LEU:HD12	2.13	0.48
1:B:73:GLN:OE1	1:B:296:SER:CB	2.62	0.48
1:A:260:ALA:HB2	1:A:292:THR:HA	1.96	0.48
1:B:338:ARG:NH1	3:B:479:HOH:O	2.47	0.48
1:B:343:LEU:O	1:B:351:VAL:HG22	2.14	0.48
1:B:184:PHE:HD2	1:B:185:MSE:HE2	1.79	0.47
1:B:254:THR:CG2	1:B:256:SER:HB3	2.45	0.47
1:B:153:ALA:CB	1:B:177:SER:HB2	2.43	0.47
1:B:343:LEU:HD23	1:B:352:MSE:HE2	1.95	0.47
1:A:252:LEU:HB2	1:A:316:VAL:O	2.15	0.47
1:B:137:GLU:OE2	1:B:164:ARG:NH1	2.47	0.47
1:B:89:LYS:HB2	1:B:89:LYS:HE3	1.65	0.46
1:B:180:THR:OG1	1:B:183:GLU:HG3	2.15	0.46
1:B:254:THR:CG2	1:B:255:ASP:N	2.79	0.46
1:A:92:GLU:OE1	1:A:96:LYS:HE3	2.15	0.46
1:A:120:SER:HB3	1:A:176:HIS:O	2.15	0.46
1:A:254:THR:HG23	1:A:256:SER:HB2	1.98	0.46
1:B:203:LYS:HD3	1:B:204:PRO:CD	2.45	0.46
1:A:268:THR:HB	1:A:272:LEU:O	2.16	0.46
1:A:236:GLU:H	1:A:236:GLU:CD	2.19	0.46
1:A:73:GLN:OE1	1:A:296:SER:HB3	2.16	0.46
1:A:254:THR:HG23	1:A:256:SER:HB3	1.99	0.45
1:B:66:MSE:O	1:B:70:LEU:HG	2.15	0.45
1:A:254:THR:CG2	1:A:256:SER:HB3	2.46	0.45
1:A:252:LEU:HD12	1:A:252:LEU:HA	1.74	0.45
1:A:56:ALA:HB1	1:A:62:ASP:HB2	1.99	0.45
1:B:248:ASN:H	1:B:248:ASN:ND2	2.16	0.44
1:A:36:ILE:HB	3:A:428:HOH:O	2.16	0.44
1:B:254:THR:HG23	1:B:256:SER:HB3	1.99	0.44
1:A:73:GLN:OE1	1:A:296:SER:HB2	2.15	0.44
1:B:210:ILE:HB	1:B:214:GLY:HA3	1.99	0.44
1:A:119:ILE:O	1:A:156:ILE:HD12	2.18	0.43
1:A:210:ILE:HB	1:A:214:GLY:HA3	1.98	0.43
1:A:1:MSE:HG3	1:A:22:ILE:O	2.17	0.43
1:B:222:VAL:O	1:B:245:LYS:NZ	2.51	0.43
1:A:181:PHE:O	1:A:185:MSE:HG2	2.19	0.43
1:B:73:GLN:OE1	1:B:294:PHE:HB3	2.19	0.43
1:A:295:PHE:O	1:A:298:ALA:HB3	2.19	0.43
1:A:144:ASP:O	1:A:145:SER:C	2.55	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:5:LYS:NZ	1:B:33:ARG:HH12	2.16	0.43
1:A:10:ASP:HA	1:A:11:PRO:HD3	1.90	0.43
1:B:22:ILE:HG21	1:B:357:GLY:HA2	2.01	0.42
1:A:348:LEU:HA	1:A:348:LEU:HD12	1.84	0.42
1:A:345:ASP:OD1	1:A:345:ASP:C	2.58	0.42
1:B:243:LYS:O	1:B:246:LYS:NZ	2.40	0.42
1:B:151:THR:HA	1:B:173:SER:O	2.20	0.42
1:A:279:GLY:O	1:A:281:PRO:HD3	2.19	0.42
1:B:1:MSE:HB3	1:B:22:ILE:O	2.20	0.42
1:B:131:HIS:O	1:B:133:ARG:HD3	2.20	0.42
1:B:276:VAL:HG22	1:B:281:PRO:HB3	2.02	0.42
1:B:150:LEU:CD1	1:B:152:PHE:HB3	2.50	0.42
1:B:91:LYS:HD3	1:B:142:GLU:OE2	2.20	0.41
1:A:17:THR:CG2	1:A:37:PRO:CG	2.93	0.41
1:B:5:LYS:O	1:B:38:ARG:N	2.49	0.41
1:A:6:VAL:O	1:A:17:THR:HA	2.20	0.41
1:B:46:VAL:HG22	1:B:77:THR:HB	2.02	0.41
1:A:235:ARG:HG3	1:A:236:GLU:N	2.35	0.41
1:A:125:GLY:O	1:A:126:ALA:CB	2.67	0.41
1:A:70:LEU:HB3	1:A:75:VAL:HB	2.03	0.41
1:A:254:THR:CG2	1:A:256:SER:CB	2.96	0.41
1:A:102:LEU:C	1:A:102:LEU:HD13	2.41	0.41
1:B:85:THR:HG23	1:B:86:SER:O	2.21	0.41
1:B:89:LYS:O	1:B:93:ILE:HG13	2.21	0.41
1:A:344:LEU:HD12	1:A:344:LEU:N	2.36	0.41
1:B:263:LYS:CE	1:B:263:LYS:N	2.66	0.41
1:B:197:HIS:N	1:B:225:GLU:HB2	2.35	0.41
1:A:235:ARG:HD2	1:A:239:LYS:NZ	2.36	0.40
1:B:198:PHE:HA	1:B:199:PRO:HA	1.80	0.40
1:A:52:GLY:HA3	1:A:56:ALA:O	2.21	0.40
1:B:120:SER:HB2	3:B:453:HOH:O	2.20	0.40
1:A:21:GLU:CD	1:A:29:LYS:HE3	2.41	0.40
1:A:313:LEU:HA	1:A:316:VAL:HG22	2.03	0.40
1:A:5:LYS:HB2	1:A:5:LYS:HZ3	1.85	0.40
1:A:87:LEU:CG	1:A:139:GLU:HG2	2.51	0.40
1:A:135:PRO:HD2	1:A:157:GLU:O	2.22	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	362/376 (96%)	341 (94%)	17 (5%)	4 (1%)	21	34
1	B	362/376 (96%)	341 (94%)	19 (5%)	2 (1%)	33	55
All	All	724/752 (96%)	682 (94%)	36 (5%)	6 (1%)	27	46

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	35	CYS
1	A	125	GLY
1	A	264	ASP
1	A	357	GLY
1	A	197	HIS
1	B	197	HIS

### 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	316/318 (99%)	287 (91%)	29 (9%)	13	24
1	B	316/318 (99%)	296 (94%)	20 (6%)	25	44
All	All	632/636 (99%)	583 (92%)	49 (8%)	18	32

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	36	ILE
1	A	38	ARG

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Mol	Chain	Res	Type
1	A	57	ASP
1	A	79	LEU
1	A	87	LEU
1	A	88	GLU
1	A	90	MSE
1	A	110	LEU
1	A	150	LEU
1	A	160	GLU
1	A	177	SER
1	A	186	LYS
1	A	197	HIS
1	A	199	PRO
1	A	235	ARG
1	A	237	MSE
1	A	248	ASN
1	A	252	LEU
1	A	254	THR
1	A	263	LYS
1	A	269	LEU
1	A	277	LYS
1	A	282	ARG
1	A	285	ASP
1	A	330	ARG
1	A	348	LEU
1	A	351	VAL
1	A	355	LYS
1	B	3	VAL
1	B	7	LEU
1	B	79	LEU
1	B	85	THR
1	B	87	LEU
1	B	110	LEU
1	B	122	GLU
1	B	133	ARG
1	B	150	LEU
1	B	156	ILE
1	B	189	LYS
1	B	235	ARG
1	B	252	LEU
1	B	253	VAL
1	B	254	THR
1	B	263	LYS

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Mol	Chain	Res	Type
1	B	304	LYS
1	B	313	LEU
1	B	348	LEU
1	B	353	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	HIS
1	B	248	ASN
1	B	297	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 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.

## 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 ⓘ

### 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	364/376 (96%)	-0.21	1 (0%) 91 93	27, 39, 62, 75	0
1	B	364/376 (96%)	-0.15	2 (0%) 88 90	25, 38, 58, 71	0
All	All	728/752 (96%)	-0.18	3 (0%) 90 92	25, 38, 61, 75	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	28	VAL	3.3
1	B	269	LEU	2.6
1	A	271	ASP	2.6

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FE	B	401	1/1	0.09	-1.62	52,52,52,52	0
2	FE	A	401	1/1	0.04	-2.82	67,67,67,67	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.