



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:45 AM GMT

PDB ID : 2Y8V  
Title : Structure of chitinase, ChiC, from *Aspergillus fumigatus*.  
Authors : Rush, C.L.; Schuettelkopf, A.W.; Gay, L.M.; van Aalten, D.M.F.  
Deposited on : 2011-02-10  
Resolution : 1.99 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

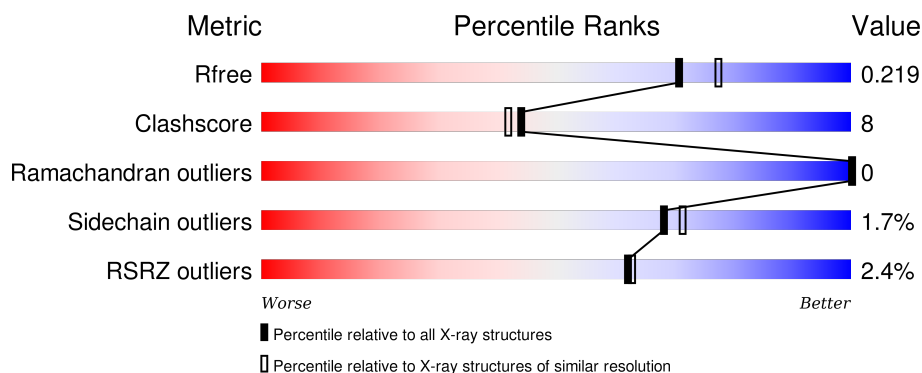
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.99 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	290	<div> <div>2%</div> <div>79%</div> <div>16%</div> <div>..</div> </div>
1	B	290	<div> <div>2%</div> <div>82%</div> <div>13%</div> <div>..</div> </div>
1	C	290	<div> <div>3%</div> <div>82%</div> <div>13%</div> <div>..</div> </div>
1	D	290	<div> <div>%</div> <div>83%</div> <div>10%</div> <div>..</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9395 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 CLASS III CHITINASE, PUTATIVE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	0	0
			2205	1408	385	400	12			
1	B	279	Total	C	N	O	S	0	0	0
			2204	1407	385	400	12			
1	C	281	Total	C	N	O	S	0	0	0
			2212	1411	383	405	13			
1	D	278	Total	C	N	O	S	0	0	0
			2196	1401	382	401	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q4WEQ6
A	0	PRO	-	EXPRESSION TAG	UNP Q4WEQ6
B	-1	GLY	-	EXPRESSION TAG	UNP Q4WEQ6
B	0	PRO	-	EXPRESSION TAG	UNP Q4WEQ6
C	-1	GLY	-	EXPRESSION TAG	UNP Q4WEQ6
C	0	PRO	-	EXPRESSION TAG	UNP Q4WEQ6
D	-1	GLY	-	EXPRESSION TAG	UNP Q4WEQ6
D	0	PRO	-	EXPRESSION TAG	UNP Q4WEQ6

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Na	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	127	Total	O	0	0
			127	127		

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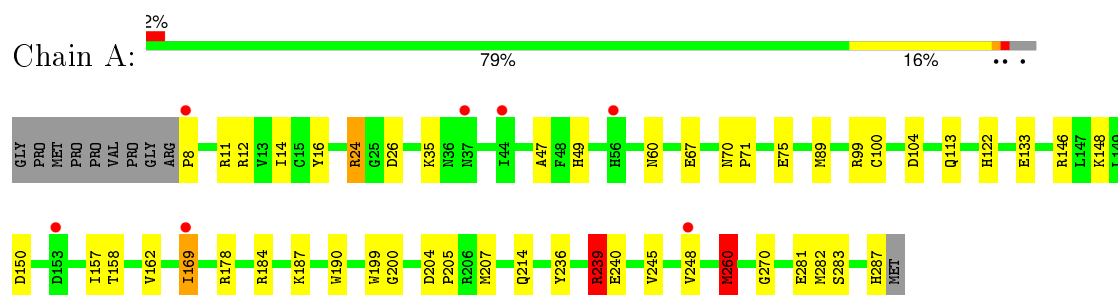
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	162	Total 162	O 162	0	0
3	C	131	Total 131	O 131	0	0
3	D	157	Total 157	O 157	0	0

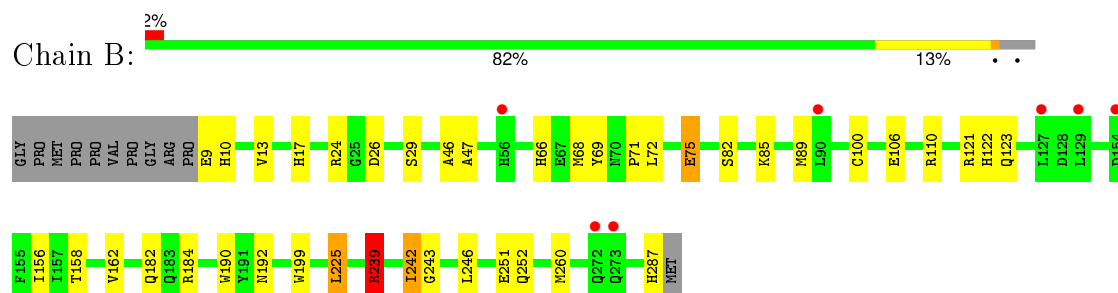
### 3 Residue-property plots [i](#)

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.

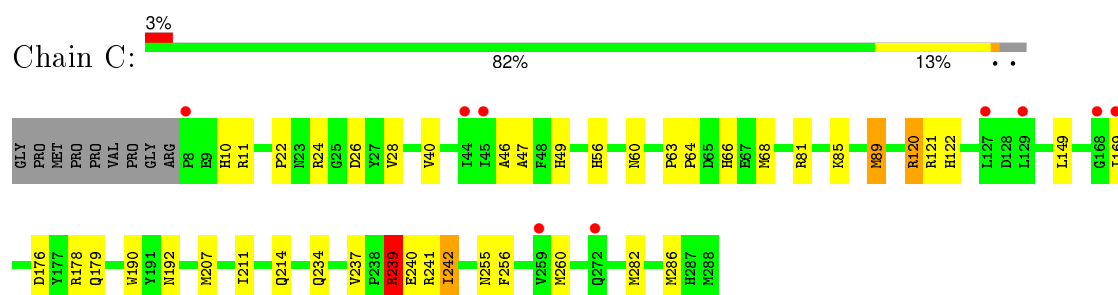
- Molecule 1: CLASS III CHITINASE, PUTATIVE



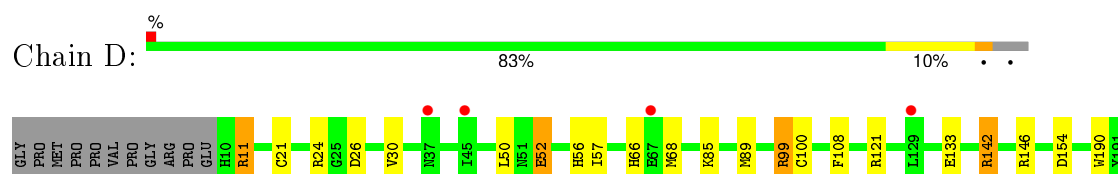
- Molecule 1: CLASS III CHITINASE, PUTATIVE



- Molecule 1: CLASS III CHITINASE, PUTATIVE



- Molecule 1: CLASS III CHITINASE, PUTATIVE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.14Å 145.75Å 82.90Å 90.00° 93.29° 90.00°	Depositor
Resolution (Å)	82.76 – 1.99 24.85 – 1.99	Depositor EDS
% Data completeness (in resolution range)	89.6 (82.76-1.99) 89.7 (24.85-1.99)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.59 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.169 , 0.213 0.172 , 0.219	Depositor DCC
$R_{free}$ test set	1507 reflections (2.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	19.4	Xtriage
Anisotropy	0.453	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 45.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 76641 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9395	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: NA

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	1.11	3/2265 (0.1%)	0.99	8/3082 (0.3%)
1	B	1.19	2/2264 (0.1%)	0.98	10/3080 (0.3%)
1	C	1.10	1/2273 (0.0%)	1.11	8/3093 (0.3%)
1	D	1.18	2/2255 (0.1%)	1.08	9/3068 (0.3%)
All	All	1.15	8/9057 (0.1%)	1.04	35/12323 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	9	GLU	CD-OE2	13.17	1.40	1.25
1	A	75	GLU	CB-CG	-8.55	1.35	1.52
1	D	52	GLU	CG-CD	8.41	1.64	1.51
1	B	75	GLU	CB-CG	-8.02	1.36	1.52
1	A	100	CYS	CB-SG	-5.31	1.73	1.81
1	C	240	GLU	CD-OE1	5.25	1.31	1.25
1	A	236	TYR	CE1-CZ	5.20	1.45	1.38
1	D	276	TRP	CB-CG	5.10	1.59	1.50

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	142	ARG	NE-CZ-NH1	18.14	129.37	120.30
1	C	89	MET	CG-SD-CE	-16.89	73.18	100.20
1	D	142	ARG	NE-CZ-NH2	-14.96	112.82	120.30
1	C	239	ARG	NE-CZ-NH2	-13.84	113.38	120.30
1	C	120	ARG	NE-CZ-NH1	13.19	126.90	120.30
1	C	120	ARG	NE-CZ-NH2	-10.36	115.12	120.30
1	C	242	ILE	CG1-CB-CG2	-9.96	89.50	111.40
1	B	239	ARG	NE-CZ-NH2	-9.21	115.69	120.30
1	D	89	MET	CG-SD-CE	-9.13	85.58	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	239	ARG	NE-CZ-NH1	8.76	124.68	120.30
1	D	100	CYS	CA-CB-SG	-7.74	100.06	114.00
1	C	239	ARG	NE-CZ-NH1	7.58	124.09	120.30
1	A	24	ARG	NE-CZ-NH1	7.18	123.89	120.30
1	D	99	ARG	NE-CZ-NH1	7.11	123.86	120.30
1	A	89	MET	CG-SD-CE	-7.08	88.88	100.20
1	D	259	VAL	CG1-CB-CG2	-7.05	99.62	110.90
1	B	121	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	B	242	ILE	CG1-CB-CG2	-6.38	97.36	111.40
1	A	8	PRO	N-CA-CB	6.36	110.94	103.30
1	D	241	ARG	NE-CZ-NH1	6.17	123.39	120.30
1	B	89	MET	CG-SD-CE	-6.10	90.44	100.20
1	B	121	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	A	239	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	D	246	LEU	CB-CG-CD1	-5.76	101.20	111.00
1	B	184	ARG	NE-CZ-NH2	-5.75	117.42	120.30
1	C	85	LYS	CD-CE-NZ	-5.61	98.79	111.70
1	D	99	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	C	241	ARG	NE-CZ-NH1	5.39	122.99	120.30
1	A	260	MET	CG-SD-CE	5.28	108.65	100.20
1	A	169	ILE	CB-CA-C	-5.22	101.15	111.60
1	B	100	CYS	CA-CB-SG	-5.15	104.73	114.00
1	B	75	GLU	CB-CA-C	-5.12	100.16	110.40
1	A	150	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	225	LEU	CB-CG-CD2	-5.10	102.33	111.00
1	A	12	ARG	NE-CZ-NH2	-5.05	117.77	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-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 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	2205	0	2136	46	0
1	B	2204	0	2140	36	0
1	C	2212	0	2135	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2196	0	2133	35	0
2	C	1	0	0	0	0
3	A	127	0	0	9	1
3	B	162	0	0	10	0
3	C	131	0	0	7	0
3	D	157	0	0	12	1
All	All	9395	0	8544	139	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (139) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:121:ARG:HD3	3:C:2064:HOH:O	1.42	1.15
1:A:67:GLU:HG3	3:A:2039:HOH:O	1.58	1.03
1:B:182:GLN:HB3	3:B:2103:HOH:O	1.60	1.00
1:C:178:ARG:HD3	1:C:214:GLN:NE2	1.84	0.92
1:B:252:GLN:HG2	3:B:2149:HOH:O	1.72	0.89
1:A:239:ARG:HG3	1:A:239:ARG:HH11	1.40	0.86
1:B:66:HIS:HD2	1:B:68:MET:H	1.25	0.80
1:D:52:GLU:H	1:D:52:GLU:CD	1.87	0.78
1:D:99:ARG:HD3	3:D:2070:HOH:O	1.83	0.77
1:A:169:ILE:O	1:A:169:ILE:HG22	1.86	0.76
1:C:169:ILE:CG2	1:C:169:ILE:O	2.33	0.75
1:A:99:ARG:HD3	3:A:2067:HOH:O	1.86	0.75
1:A:239:ARG:HG3	1:A:239:ARG:NH1	1.99	0.74
1:C:169:ILE:HG22	1:C:169:ILE:O	1.88	0.74
1:C:22:PRO:HD2	1:C:28:VAL:HG13	1.69	0.73
1:D:68:MET:HE1	3:D:2012:HOH:O	1.88	0.72
1:A:282:MET:HE2	1:B:225:LEU:HD12	1.71	0.72
1:C:178:ARG:HD3	1:C:214:GLN:HE21	1.54	0.72
1:D:85:LYS:CE	3:D:2031:HOH:O	2.38	0.71
1:A:16:TYR:HE1	1:A:260:MET:HG3	1.55	0.71
1:C:239:ARG:HD2	1:D:281:GLU:OE1	1.91	0.70
1:C:178:ARG:CD	1:C:214:GLN:NE2	2.54	0.69
1:A:104:ASP:OD2	3:A:2049:HOH:O	2.11	0.68
1:C:190:TRP:HE1	1:C:192:ASN:HD21	1.42	0.66
1:B:10:HIS:HD2	3:B:2002:HOH:O	1.78	0.66
1:D:56:HIS:HD2	3:D:2025:HOH:O	1.76	0.66
1:C:120:ARG:HD3	3:C:2061:HOH:O	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:ARG:HD2	3:D:2050:HOH:O	1.95	0.64
1:A:282:MET:CE	1:B:225:LEU:HD12	2.27	0.64
1:C:286:MET:O	1:D:11:ARG:HD2	1.98	0.64
1:A:184:ARG:HD2	1:A:187:LYS:HD2	1.80	0.63
1:B:17:HIS:HD2	3:B:2155:HOH:O	1.83	0.61
1:A:239:ARG:HH11	1:A:239:ARG:CG	2.12	0.61
1:A:287:HIS:C	3:A:2127:HOH:O	2.38	0.60
1:C:207:MET:O	1:C:211:ILE:HD13	2.01	0.60
3:A:2003:HOH:O	1:B:287:HIS:HE1	1.85	0.60
1:D:85:LYS:HE2	3:D:2031:HOH:O	2.02	0.60
1:C:47:ALA:H	1:C:60:ASN:ND2	2.00	0.59
1:A:16:TYR:HE1	1:A:260:MET:CG	2.15	0.59
1:D:190:TRP:HE1	1:D:192:ASN:ND2	2.01	0.58
1:C:176:ASP:HB3	1:C:179:GLN:HG3	1.84	0.58
1:B:106:GLU:HG2	1:B:110:ARG:NH1	2.19	0.58
1:C:22:PRO:HG2	1:C:28:VAL:HG11	1.84	0.58
1:C:239:ARG:NH2	1:D:285:SER:OG	2.37	0.58
1:C:122:HIS:HD2	3:C:2065:HOH:O	1.86	0.57
1:A:113:GLN:NE2	1:A:146:ARG:NH2	2.52	0.57
1:C:190:TRP:HE1	1:C:192:ASN:ND2	2.02	0.57
1:A:169:ILE:O	1:A:169:ILE:CG2	2.52	0.57
1:C:47:ALA:H	1:C:60:ASN:HD21	1.51	0.57
1:C:66:HIS:ND1	1:C:68:MET:HG2	2.19	0.57
1:A:113:GLN:HE22	1:A:146:ARG:NH2	2.03	0.57
1:D:190:TRP:HE1	1:D:192:ASN:HD21	1.54	0.56
1:B:252:GLN:CG	3:B:2149:HOH:O	2.43	0.56
1:A:283:SER:HB3	1:B:13:VAL:HG21	1.87	0.56
1:D:108:PHE:CD1	1:D:142:ARG:HD3	2.40	0.56
1:B:24:ARG:HD3	1:C:26:ASP:OD2	2.05	0.56
1:A:24:ARG:HD3	1:D:26:ASP:OD2	2.06	0.56
1:A:47:ALA:H	1:A:60:ASN:ND2	2.06	0.54
1:C:286:MET:HB3	1:D:11:ARG:HB3	1.90	0.54
1:A:200:GLY:HA2	1:A:207:MET:SD	2.48	0.54
1:D:52:GLU:N	1:D:52:GLU:CD	2.60	0.53
1:D:154:ASP:HB2	3:D:2082:HOH:O	2.09	0.53
1:C:237:VAL:CG1	1:C:242:ILE:HG12	2.38	0.53
1:B:190:TRP:HE1	1:B:192:ASN:ND2	2.06	0.52
1:A:158:THR:HG21	1:A:190:TRP:CE2	2.44	0.52
1:D:68:MET:CE	3:D:2012:HOH:O	2.49	0.52
1:D:50:LEU:HD23	1:D:57:ILE:HD12	1.92	0.52
1:C:178:ARG:HH11	1:C:214:GLN:NE2	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:GLU:OE1	1:B:239:ARG:HD2	2.10	0.51
1:B:251:GLU:OE2	3:B:2148:HOH:O	2.18	0.51
1:C:56:HIS:HD2	3:C:2024:HOH:O	1.92	0.51
1:D:121:ARG:HG2	1:D:121:ARG:HH11	1.75	0.51
1:C:56:HIS:HE1	3:C:2026:HOH:O	1.93	0.51
1:B:66:HIS:CD2	1:B:68:MET:H	2.17	0.50
1:D:50:LEU:CD2	1:D:57:ILE:HD12	2.41	0.50
1:A:35:LYS:HG3	3:A:2018:HOH:O	2.11	0.50
1:B:182:GLN:CB	3:B:2103:HOH:O	2.38	0.50
1:A:282:MET:CE	1:B:225:LEU:CD1	2.90	0.49
1:A:205:PRO:HD3	1:A:245:VAL:HG13	1.94	0.49
1:A:113:GLN:HE22	1:A:146:ARG:HH21	1.60	0.49
1:D:99:ARG:HG3	1:D:133:GLU:OE2	2.13	0.49
1:B:26:ASP:OD2	1:C:24:ARG:HD3	2.12	0.49
1:D:56:HIS:HE1	3:D:2027:HOH:O	1.95	0.49
1:A:282:MET:HE2	1:B:225:LEU:CD1	2.41	0.49
1:A:47:ALA:H	1:A:60:ASN:HD21	1.59	0.48
1:D:146:ARG:HG2	3:D:2077:HOH:O	2.11	0.48
1:C:178:ARG:CD	1:C:214:GLN:HE21	2.19	0.48
1:B:122:HIS:HD2	3:B:2075:HOH:O	1.95	0.48
1:C:22:PRO:CD	1:C:28:VAL:HG13	2.41	0.48
1:A:239:ARG:NH1	1:A:270:GLY:O	2.47	0.47
1:C:10:HIS:CD2	1:C:255:ASN:HA	2.49	0.47
1:A:204:ASP:HB2	1:C:234:GLN:NE2	2.29	0.47
3:B:2015:HOH:O	1:C:28:VAL:HG12	2.14	0.47
1:C:81:ARG:HD3	3:C:2043:HOH:O	2.14	0.47
1:C:282:MET:HB3	1:D:259:VAL:HG11	1.97	0.47
1:B:162:VAL:HB	1:B:199:TRP:HB3	1.96	0.47
1:B:199:TRP:HD1	3:B:2108:HOH:O	1.97	0.46
1:C:11:ARG:HD2	1:D:286:MET:HB3	1.97	0.46
1:B:85:LYS:HE2	1:B:156:ILE:HD11	1.98	0.46
1:D:108:PHE:CE1	1:D:142:ARG:HD3	2.51	0.46
1:A:204:ASP:HB2	1:C:234:GLN:HE22	1.81	0.46
1:B:190:TRP:HE1	1:B:192:ASN:HD21	1.64	0.46
1:A:16:TYR:CE1	1:A:260:MET:CG	2.98	0.45
1:A:148:LYS:HE2	1:A:157:ILE:HG12	1.99	0.45
1:C:178:ARG:HD3	1:C:214:GLN:HE22	1.73	0.45
1:B:242:ILE:HD13	1:B:242:ILE:HA	1.88	0.45
1:C:40:VAL:HG13	1:C:40:VAL:O	2.15	0.45
1:A:282:MET:HE1	1:B:246:LEU:HD11	1.98	0.45
1:B:71:PRO:O	1:B:75:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:HIS:HE1	3:A:2043:HOH:O	2.01	0.44
1:D:85:LYS:HE3	3:D:2031:HOH:O	2.12	0.44
1:D:21:CYS:SG	1:D:68:MET:CE	3.05	0.44
1:C:10:HIS:HB2	1:C:256:PHE:O	2.17	0.43
1:D:224:GLY:HA2	1:D:260:MET:HG3	2.01	0.43
1:A:99:ARG:HG3	1:A:133:GLU:OE2	2.19	0.42
1:B:162:VAL:HB	1:B:199:TRP:CB	2.48	0.42
1:A:214:GLN:HE21	1:A:214:GLN:HB3	1.73	0.42
1:D:66:HIS:ND1	1:D:68:MET:HG2	2.34	0.42
1:B:46:ALA:HA	1:B:47:ALA:HA	1.88	0.42
1:C:49:HIS:HE1	3:C:2044:HOH:O	2.02	0.42
1:A:14:ILE:HB	1:A:260:MET:HB2	2.01	0.41
1:A:169:ILE:HG21	1:A:169:ILE:HD13	1.72	0.41
1:A:11:ARG:NH1	1:B:287:HIS:CD2	2.89	0.41
1:D:241:ARG:O	1:D:244:PRO:HD2	2.21	0.41
1:B:242:ILE:HD12	1:B:242:ILE:HG23	1.46	0.41
1:A:199:TRP:HD1	3:A:2095:HOH:O	2.03	0.41
1:A:162:VAL:HB	1:A:199:TRP:CB	2.50	0.41
1:B:239:ARG:NH2	1:B:243:GLY:HA3	2.36	0.41
1:A:26:ASP:OD2	1:D:24:ARG:HD3	2.20	0.41
1:A:240:GLU:H	1:A:240:GLU:CD	2.24	0.41
1:D:200:GLY:HA2	1:D:207:MET:SD	2.61	0.41
1:B:29:SER:OG	1:B:75:GLU:OE2	2.26	0.41
1:A:70:ASN:HB2	1:A:71:PRO:HD3	2.03	0.41
1:D:242:ILE:HG21	3:D:2124:HOH:O	2.20	0.40
1:A:122:HIS:HD2	3:A:2062:HOH:O	2.04	0.40
1:C:63:PRO:HA	1:C:64:PRO:HD3	1.99	0.40
1:C:46:ALA:HA	1:C:47:ALA:HA	1.91	0.40
1:B:158:THR:HG21	1:B:190:TRP:CE2	2.56	0.40
1:B:69:TYR:O	1:B:72:LEU:HB3	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:2090:HOH:O	3:D:2072:HOH:O[2_656]	2.07	0.13

## 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	278/290 (96%)	272 (98%)	6 (2%)	0	100	100
1	B	277/290 (96%)	271 (98%)	6 (2%)	0	100	100
1	C	279/290 (96%)	272 (98%)	7 (2%)	0	100	100
1	D	276/290 (95%)	273 (99%)	3 (1%)	0	100	100
All	All	1110/1160 (96%)	1088 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	230/242 (95%)	226 (98%)	4 (2%)	68	71
1	B	231/242 (96%)	227 (98%)	4 (2%)	68	71
1	C	232/242 (96%)	228 (98%)	4 (2%)	68	71
1	D	231/242 (96%)	227 (98%)	4 (2%)	68	71
All	All	924/968 (96%)	908 (98%)	16 (2%)	68	71

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

Mol	Chain	Res	Type
1	A	178	ARG
1	A	239	ARG

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Mol	Chain	Res	Type
1	A	248	VAL
1	A	260	MET
1	B	82	SER
1	B	123	GLN
1	B	239	ARG
1	B	260	MET
1	C	89	MET
1	C	149	LEU
1	C	239	ARG
1	C	260	MET
1	D	11	ARG
1	D	30	VAL
1	D	239	ARG
1	D	260	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	49	HIS
1	A	60	ASN
1	A	105	GLN
1	A	113	GLN
1	A	122	HIS
1	A	214	GLN
1	B	17	HIS
1	B	49	HIS
1	B	66	HIS
1	B	122	HIS
1	B	192	ASN
1	B	287	HIS
1	C	36	ASN
1	C	49	HIS
1	C	56	HIS
1	C	60	ASN
1	C	122	HIS
1	C	179	GLN
1	C	192	ASN
1	C	214	GLN
1	C	234	GLN
1	C	252	GLN
1	C	287	HIS

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Mol	Chain	Res	Type
1	D	18	GLN
1	D	56	HIS
1	D	192	ASN

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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	280/290 (96%)	0.23	7 (2%) 61 61	13, 21, 33, 39	0
1	B	279/290 (96%)	0.11	7 (2%) 61 61	10, 18, 31, 36	0
1	C	281/290 (96%)	0.12	9 (3%) 51 52	10, 18, 31, 42	0
1	D	278/290 (95%)	0.00	4 (1%) 78 78	8, 16, 28, 35	0
All	All	1118/1160 (96%)	0.12	27 (2%) 62 63	8, 19, 31, 42	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	8	PRO	3.9
1	A	169	ILE	3.6
1	C	169	ILE	3.2
1	A	8	PRO	3.2
1	A	37	ASN	3.0
1	C	272	GLN	3.0
1	C	168	GLY	2.7
1	D	37	ASN	2.7
1	B	272	GLN	2.6
1	A	248	VAL	2.4
1	D	67	GLU	2.4
1	C	44	ILE	2.3
1	A	44	ILE	2.3
1	C	129	LEU	2.2
1	A	56	HIS	2.2
1	B	56	HIS	2.2
1	B	273	GLN	2.2
1	D	129	LEU	2.2
1	D	45	ILE	2.1
1	B	129	LEU	2.1
1	C	127	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	153	ASP	2.1
1	C	45	ILE	2.0
1	B	90	LEU	2.0
1	B	127	LEU	2.0
1	C	259	VAL	2.0
1	B	154	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 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	NA	C	1291	1/1	0.95	0.19	0.96	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.