Full wwPDB X-ray Structure Validation Report (i)

Nov 1, 2023 – 03:22 PM JST

PDB ID	:	5JGF
Title	:	Crystal structure of mApe1
Authors	:	Noda, N.N.; Adachi, W.; Inagaki, F.
Deposited on	:	2016-04-20
Resolution	:	1.83 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 (i) 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 (1)) were used in the production of this report:

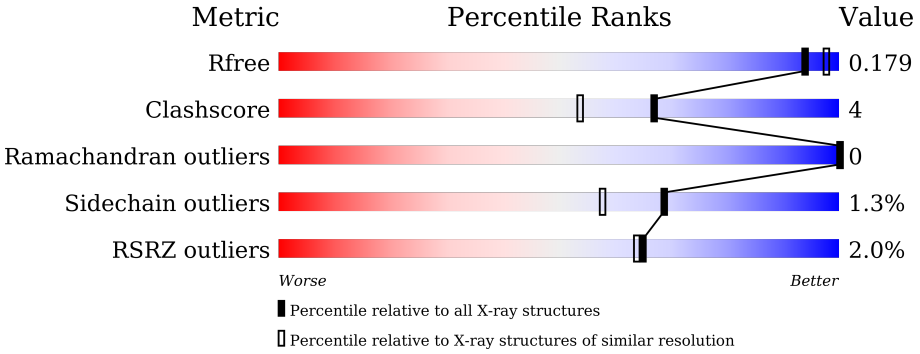
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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}	130704	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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 <=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	473	.% 88% 10% •
1	B	473	3% 88% 8% • •
1	C	473	2% 88% 9% •
1	D	473	2% 88% 9% •

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 16586 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 Vacuolar aminopeptidase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	460	3529	2260	595	666	8	0	0	0
1	B	460	3540	2266	597	669	8	0	0	0
1	C	460	3544	2268	597	671	8	0	0	0
1	D	462	3564	2280	602	674	8	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	GLY	-	expression tag	UNP P14904
A	43	PRO	-	expression tag	UNP P14904
A	44	HIS	-	expression tag	UNP P14904
A	45	MET	-	expression tag	UNP P14904
B	42	GLY	-	expression tag	UNP P14904
B	43	PRO	-	expression tag	UNP P14904
B	44	HIS	-	expression tag	UNP P14904
B	45	MET	-	expression tag	UNP P14904
C	42	GLY	-	expression tag	UNP P14904
C	43	PRO	-	expression tag	UNP P14904
C	44	HIS	-	expression tag	UNP P14904
C	45	MET	-	expression tag	UNP P14904
D	42	GLY	-	expression tag	UNP P14904
D	43	PRO	-	expression tag	UNP P14904
D	44	HIS	-	expression tag	UNP P14904
D	45	MET	-	expression tag	UNP P14904

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is water.

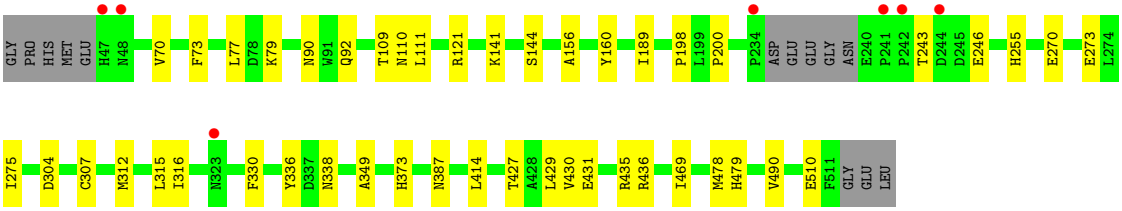
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	594	594	594	0	0
			Total	O		
3	B	569	569	569	0	0
			Total	O		
3	C	614	614	614	0	0
			Total	O		
3	D	624	624	624	0	0

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.

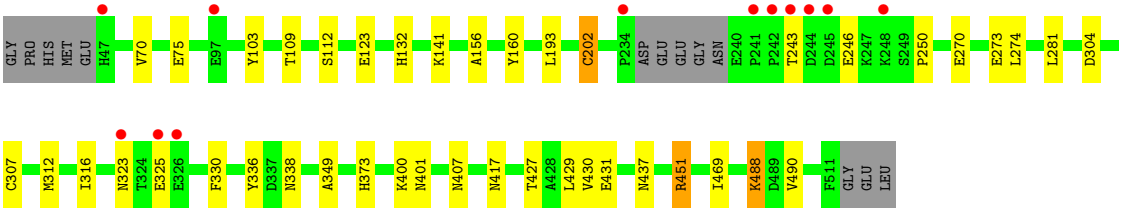
• Molecule 1: Vacuolar aminopeptidase 1

Chain A:



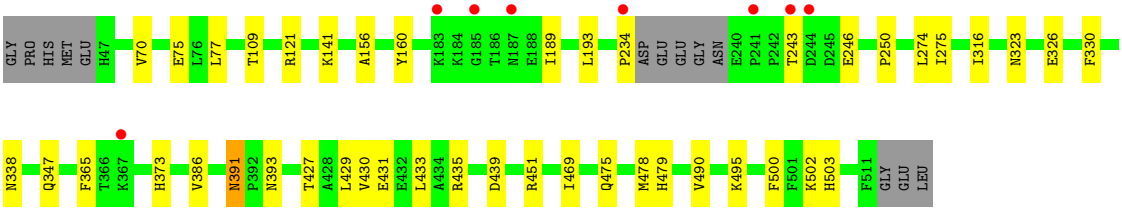
• Molecule 1: Vacuolar aminopeptidase 1

Chain B:



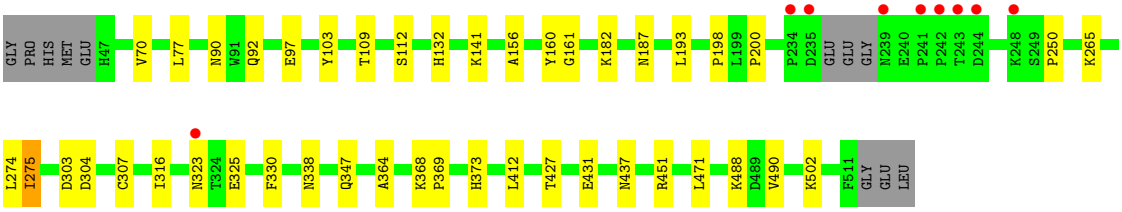
• Molecule 1: Vacuolar aminopeptidase 1

Chain C:



• Molecule 1: Vacuolar aminopeptidase 1

Chain D:


GLY PRO HIS MET GLU H47 V70 L77 N90 W91 Q92 E97 Y103 T109 S112 H132 K141 A156 Y160 G161 K182 N187 L193 P198 L199 P200 P234 D235 GLU GLU GLY N239 E240 P241 P242 T243 D244 K248 S249 P250 K265
L274 I275 D303 D304 C307 I316 N323 T324 E325 F330 N338 Q347 A364 K368 P369 H373 L412 T427 E431 M437 R451 L471 K488 D489 V490 K502 F511 GLY GLU LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	140.97Å 140.97Å 348.92Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.49 – 1.83 38.49 – 1.83	Depositor EDS
% Data completeness (in resolution range)	96.5 (38.49-1.83) 96.7 (38.49-1.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.23 (at 1.83Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.162 , 0.179 0.162 , 0.179	Depositor DCC
R_{free} test set	22019 reflections (9.67%)	wwPDB-VP
Wilson B-factor (Å^2)	17.1	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 67.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.005 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/3*k+1/3*l 0.007 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+1/3*l,-4/3*h+4/3*k+1/3*l 0.006 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k+1/3*l 0.006 for -h,2/3*h+1/3*k+1/3*l,4/3*h+8/3*k-1/3*l 0.007 for 1/3*h+2/3*k-1/3*l,-k,-8/3*h-4/3*k-1/3*l 0.007 for -1/3*h-2/3*k+1/3*l,-2/3*h-1/3*k-1/3*l,4/3*h-4/3*k-1/3*l 0.019 for -h-k,k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	16586	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.16% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN

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	0/3613	0.64	0/4908
1	B	0.33	1/3624 (0.0%)	0.64	0/4921
1	C	0.32	0/3628	0.64	0/4926
1	D	0.32	0/3648	0.64	0/4951
All	All	0.32	1/14513 (0.0%)	0.64	0/19706

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	202	CYS	CB-SG	-6.40	1.71	1.82

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	3529	0	3441	26	0
1	B	3540	0	3456	26	0
1	C	3544	0	3460	28	0
1	D	3564	0	3488	31	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	594	0	0	2	0
3	B	569	0	0	6	0
3	C	614	0	0	4	0
3	D	624	0	0	7	0
All	All	16586	0	13845	109	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 4.

All (109) 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:502:LYS:HG3	1:C:503:HIS:CE1	2.31	0.65
1:D:265:LYS:HD3	3:D:1100:HOH:O	1.96	0.65
1:B:417:ASN:HD21	1:D:161:GLY:H	1.45	0.65
1:D:437:ASN:HB2	3:D:1054:HOH:O	1.97	0.64
1:C:323:ASN:ND2	1:C:326:GLU:HG3	2.13	0.63
1:B:123:GLU:HG3	3:B:1057:HOH:O	1.99	0.62
1:B:243:THR:OG1	1:B:246:GLU:HG3	2.02	0.59
1:A:243:THR:OG1	1:A:246:GLU:HG3	2.01	0.59
1:B:437:ASN:HB3	3:B:1112:HOH:O	2.05	0.56
1:A:144:SER:OG	1:A:255:HIS:HD2	1.87	0.56
1:C:490:VAL:HG23	3:C:837:HOH:O	2.07	0.55
1:B:75:GLU:HG3	3:B:990:HOH:O	2.07	0.54
1:B:427:THR:O	1:B:431:GLU:HG3	2.07	0.54
1:A:73:PHE:CD2	1:A:312:MET:HG2	2.43	0.53
1:B:373:HIS:HD2	3:B:711:HOH:O	1.93	0.52
1:D:427:THR:O	1:D:431:GLU:HG3	2.09	0.52
1:D:70:VAL:HG21	1:D:109:THR:HA	1.90	0.52
1:A:189:ILE:HD12	1:A:275:ILE:HD13	1.92	0.52
1:D:92:GLN:NE2	1:D:364:ALA:HA	2.26	0.51
1:C:391:ASN:C	1:C:391:ASN:HD22	2.12	0.51
1:D:437:ASN:HB2	3:D:713:HOH:O	2.10	0.51
1:A:70:VAL:HG21	1:A:109:THR:HA	1.91	0.51
1:A:90:ASN:OD1	1:A:92:GLN:HB2	2.10	0.51
1:C:391:ASN:HD22	1:C:393:ASN:H	1.58	0.51
1:D:490:VAL:HG23	3:D:886:HOH:O	2.10	0.50
1:D:182:LYS:HD2	1:D:187:ASN:HA	1.93	0.50
1:B:323:ASN:OD1	1:B:325:GLU:HB2	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:LYS:HD2	1:A:156:ALA:HB1	1.94	0.49
1:A:77:LEU:HD21	1:A:312:MET:HE2	1.94	0.49
1:B:407:ASN:OD1	1:B:488:LYS:HB2	2.12	0.49
1:C:386:VAL:HG21	1:C:475:GLN:NE2	2.28	0.49
1:A:430:VAL:HG23	1:A:469:ILE:HD13	1.95	0.48
1:A:490:VAL:HG23	3:A:916:HOH:O	2.13	0.48
1:C:234:PRO:HD3	3:C:1207:HOH:O	2.14	0.48
1:B:417:ASN:ND2	1:D:161:GLY:H	2.10	0.47
1:C:386:VAL:HG21	1:C:475:GLN:HE21	1.79	0.47
1:D:90:ASN:OD1	1:D:92:GLN:HB2	2.14	0.47
1:C:70:VAL:HG21	1:C:109:THR:HA	1.96	0.47
1:A:427:THR:O	1:A:431:GLU:HG3	2.14	0.47
1:C:427:THR:O	1:C:431:GLU:HG3	2.14	0.47
1:C:323:ASN:ND2	1:C:326:GLU:OE2	2.48	0.47
1:B:70:VAL:HG21	1:B:109:THR:HA	1.98	0.46
1:C:250:PRO:HD2	1:C:274:LEU:O	2.15	0.46
1:C:373:HIS:HD2	3:C:1073:HOH:O	1.97	0.46
1:C:189:ILE:HD12	1:C:275:ILE:HD13	1.96	0.46
1:D:97:GLU:HG2	3:D:1102:HOH:O	2.16	0.46
1:D:502:LYS:HE3	3:D:1116:HOH:O	2.15	0.46
1:A:77:LEU:HD23	1:A:316:ILE:HD11	1.97	0.46
1:D:77:LEU:HD23	1:D:316:ILE:HD11	1.98	0.46
1:B:490:VAL:HG23	3:B:905:HOH:O	2.16	0.46
1:B:304:ASP:OD1	1:B:307:CYS:HB2	2.16	0.45
1:B:312:MET:O	1:B:316:ILE:HG12	2.17	0.45
1:D:412:LEU:HD12	1:D:471:LEU:HB3	1.98	0.45
1:C:243:THR:OG1	1:C:246:GLU:HG3	2.17	0.45
1:A:111:LEU:C	1:A:111:LEU:HD23	2.37	0.45
1:D:323:ASN:OD1	1:D:325:GLU:HB2	2.17	0.45
1:C:439:ASP:OD1	1:C:495:LYS:HE3	2.17	0.44
1:B:132:HIS:NE2	1:B:304:ASP:HB2	2.32	0.44
1:C:75:GLU:HG3	3:C:1067:HOH:O	2.15	0.44
1:D:451:ARG:HG2	1:D:451:ARG:HH11	1.82	0.44
1:D:141:LYS:HD2	1:D:156:ALA:HB1	1.99	0.44
1:D:250:PRO:HD2	1:D:274:LEU:O	2.17	0.44
1:A:160:TYR:OH	1:A:338:ASN:HA	2.18	0.44
1:A:304:ASP:OD1	1:A:307:CYS:HB2	2.17	0.44
1:C:160:TYR:OH	1:C:338:ASN:HA	2.18	0.44
1:C:429:LEU:C	1:C:429:LEU:HD23	2.38	0.44
1:D:132:HIS:NE2	1:D:304:ASP:HB2	2.32	0.44
1:C:141:LYS:HD2	1:C:156:ALA:HB1	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:304:ASP:OD1	1:D:307:CYS:HB2	2.17	0.44
1:B:430:VAL:HG23	1:B:469:ILE:HD13	2.00	0.43
1:C:478:MET:O	1:C:479:HIS:HB2	2.18	0.43
1:B:202:CYS:SG	1:B:281:LEU:HD13	2.57	0.43
1:D:373:HIS:HD2	3:D:729:HOH:O	2.01	0.43
1:B:141:LYS:HD2	1:B:156:ALA:HB1	2.00	0.43
1:A:336:TYR:CD2	1:A:349:ALA:HA	2.53	0.43
1:B:160:TYR:OH	1:B:338:ASN:HA	2.17	0.43
1:B:429:LEU:HD23	1:B:429:LEU:C	2.38	0.43
1:C:77:LEU:HD23	1:C:316:ILE:HD11	2.00	0.43
1:C:430:VAL:HG23	1:C:469:ILE:HD13	2.01	0.43
1:C:391:ASN:ND2	1:C:393:ASN:H	2.16	0.43
1:D:303:ASP:HA	1:D:304:ASP:HA	1.84	0.43
1:D:488:LYS:HE2	1:D:488:LYS:HB3	1.94	0.43
1:A:429:LEU:C	1:A:429:LEU:HD23	2.39	0.42
1:C:431:GLU:O	1:C:435:ARG:HG3	2.19	0.42
1:D:160:TYR:OH	1:D:338:ASN:HA	2.19	0.42
1:D:451:ARG:HG2	1:D:451:ARG:NH1	2.35	0.42
1:A:436:ARG:NH2	1:A:510:GLU:CD	2.73	0.41
1:B:336:TYR:CD2	1:B:349:ALA:HA	2.55	0.41
1:D:471:LEU:HD12	1:D:471:LEU:C	2.41	0.41
1:A:312:MET:CE	1:A:315:LEU:HD23	2.51	0.41
1:D:182:LYS:HE2	1:D:275:ILE:HD11	2.03	0.41
1:A:336:TYR:CE2	1:A:349:ALA:HA	2.56	0.41
1:B:451:ARG:NH2	3:B:714:HOH:O	2.53	0.41
1:B:250:PRO:HD2	1:B:274:LEU:O	2.21	0.41
1:A:270:GLU:HB2	1:A:273:GLU:HG3	2.02	0.41
1:B:400:LYS:O	1:B:401:ASN:HB2	2.21	0.41
1:B:103:TYR:HA	1:B:112:SER:O	2.21	0.41
1:B:270:GLU:HB2	1:B:273:GLU:HG3	2.03	0.41
1:C:323:ASN:HD22	1:C:326:GLU:CG	2.34	0.41
1:D:198:PRO:O	1:D:200:PRO:HD3	2.21	0.41
1:A:90:ASN:OD1	1:A:92:GLN:CB	2.69	0.40
1:D:103:TYR:HA	1:D:112:SER:O	2.20	0.40
1:A:373:HIS:HD2	3:A:719:HOH:O	2.05	0.40
1:A:431:GLU:O	1:A:435:ARG:HG3	2.20	0.40
1:D:368:LYS:HA	1:D:369:PRO:HD3	1.95	0.40
1:A:198:PRO:O	1:A:200:PRO:HD3	2.21	0.40
1:A:478:MET:O	1:A:479:HIS:HB2	2.21	0.40
1:C:433:LEU:HD21	1:C:500:PHE:HA	2.03	0.40
1:C:451:ARG:HH11	1:C:451:ARG:HG2	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	456/473 (96%)	448 (98%)	8 (2%)	0	100	100
1	B	456/473 (96%)	449 (98%)	7 (2%)	0	100	100
1	C	456/473 (96%)	449 (98%)	7 (2%)	0	100	100
1	D	458/473 (97%)	452 (99%)	6 (1%)	0	100	100
All	All	1826/1892 (96%)	1798 (98%)	28 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	375/398 (94%)	369 (98%)	6 (2%)	62	49
1	B	377/398 (95%)	373 (99%)	4 (1%)	73	64
1	C	378/398 (95%)	372 (98%)	6 (2%)	62	49
1	D	381/398 (96%)	377 (99%)	4 (1%)	76	68
All	All	1511/1592 (95%)	1491 (99%)	20 (1%)	69	58

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	LYS
1	A	110	ASN
1	A	121	ARG
1	A	330	PHE
1	A	387	ASN
1	A	414	LEU
1	B	193	LEU
1	B	330	PHE
1	B	451	ARG
1	B	488	LYS
1	C	121	ARG
1	C	193	LEU
1	C	330	PHE
1	C	347	GLN
1	C	365	PHE
1	C	391	ASN
1	D	193	LEU
1	D	275	ILE
1	D	330	PHE
1	D	347	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	255	HIS
1	B	80	HIS
1	B	92	GLN
1	B	417	ASN
1	B	503	HIS
1	C	81	ASN
1	C	323	ASN
1	C	338	ASN
1	C	347	GLN
1	C	391	ASN
1	C	437	ASN
1	C	475	GLN
1	D	92	GLN
1	D	338	ASN
1	D	347	GLN
1	D	475	GLN
1	D	498	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 monosaccharides 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 (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	460/473 (97%)	-0.42	7 (1%)	73	73	12, 18, 35, 48	0
1	B	460/473 (97%)	-0.43	12 (2%)	56	53	12, 18, 35, 50	0
1	C	460/473 (97%)	-0.47	8 (1%)	70	69	11, 16, 32, 47	0
1	D	462/473 (97%)	-0.42	9 (1%)	66	65	11, 17, 34, 46	0
All	All	1842/1892 (97%)	-0.44	36 (1%)	65	64	11, 17, 34, 50	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	239	ASN	4.7
1	A	48	ASN	4.0
1	D	235	ASP	4.0
1	A	234	PRO	3.6
1	D	241	PRO	3.1
1	C	244	ASP	3.1
1	D	243	THR	3.0
1	B	326	GLU	2.9
1	C	234	PRO	2.9
1	D	244	ASP	2.9
1	A	244	ASP	2.8
1	C	185	GLY	2.8
1	B	325	GLU	2.7
1	C	241	PRO	2.7
1	D	323	ASN	2.7
1	C	367	LYS	2.6
1	D	248	LYS	2.6
1	B	244	ASP	2.6
1	B	234	PRO	2.6
1	A	241	PRO	2.6
1	A	47	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	242	PRO	2.6
1	D	234	PRO	2.5
1	B	323	ASN	2.4
1	B	248	LYS	2.4
1	A	323	ASN	2.4
1	B	243	THR	2.4
1	C	243	THR	2.2
1	B	241	PRO	2.2
1	C	187	ASN	2.2
1	B	245	ASP	2.1
1	B	242	PRO	2.1
1	B	47	HIS	2.0
1	C	183	LYS	2.0
1	D	242	PRO	2.0
1	B	97	GLU	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 monosaccharides 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	ZN	A	601	1/1	1.00	0.02	17,17,17,17	0
2	ZN	A	602	1/1	1.00	0.02	15,15,15,15	0
2	ZN	B	601	1/1	1.00	0.04	19,19,19,19	0
2	ZN	B	602	1/1	1.00	0.04	18,18,18,18	0
2	ZN	C	601	1/1	1.00	0.01	14,14,14,14	0
2	ZN	C	602	1/1	1.00	0.01	15,15,15,15	0
2	ZN	D	601	1/1	1.00	0.02	15,15,15,15	0
2	ZN	D	602	1/1	1.00	0.03	16,16,16,16	0

6.5 Other polymers (i)

There are no such residues in this entry.