Full wwPDB X-ray Structure Validation Report (i)

Oct 25, 2023 – 09:20 PM EDT

PDB ID	:	3DIT
Title	:	Crystal structure of MAD MH2 domain
Authors	:	Hao, R.; Wu, J.W.; Wang, Z.X.
Deposited on	:	2008-06-20
Resolution	:	3.20 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:

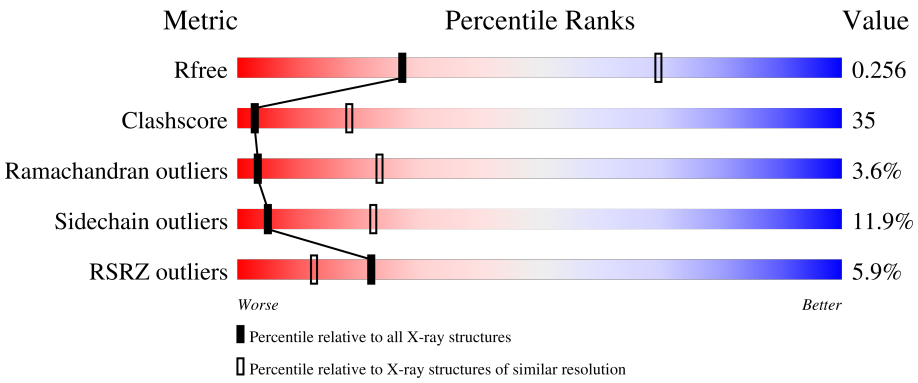
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	188	% 45% 44% 9% ••
1	B	188	% 43% 47% 9% •
1	C	188	16% 37% 56% 7% •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4548 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 Protein mothers against dpp.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	186	1482	931	263	275	13	0	0	0
1	B	186	1481	932	262	274	13	0	0	0
1	C	188	1495	939	265	278	13	0	0	0

- Molecule 2 is water.

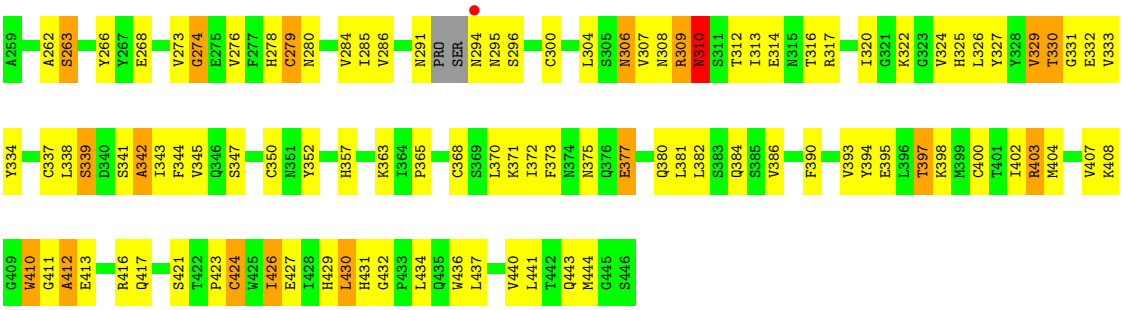
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	33	33	33	0	0
2	B	35	35	35	0	0
2	C	22	22	22	0	0

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.

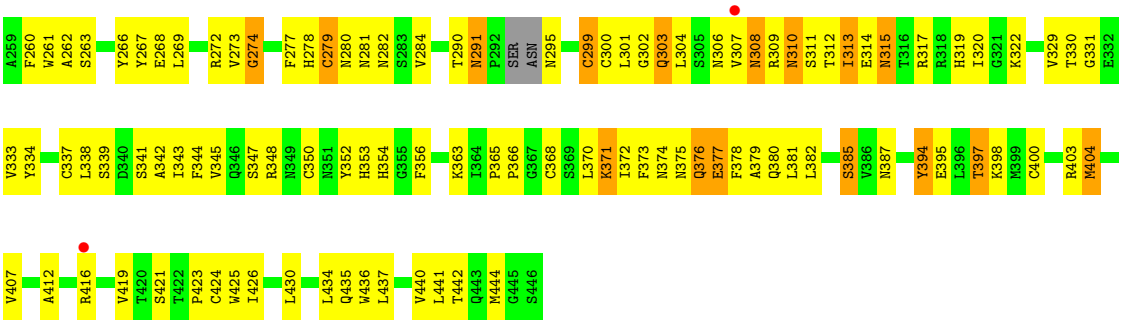
• Molecule 1: Protein mothers against dpp

Chain A:



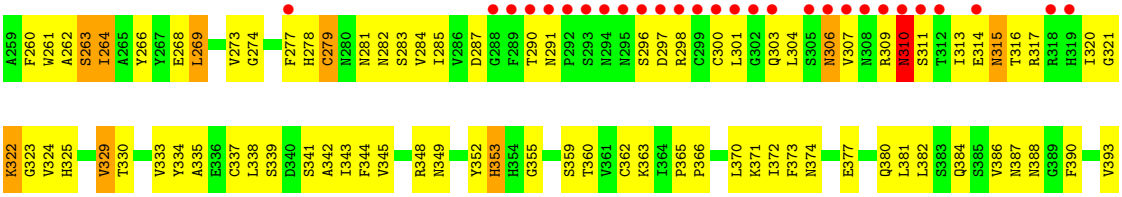
• Molecule 1: Protein mothers against dpp

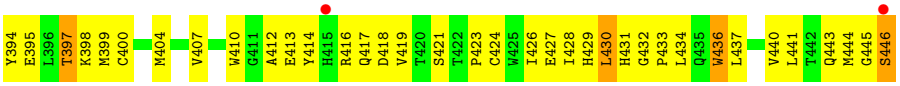
Chain B:



• Molecule 1: Protein mothers against dpp

Chain C:




Y394
E395
L396
T397
K398
M399
C400
M404
V407
W410
G411
A412
E413
Y414
H415
R416
Q417
D418
V419
T420
S421
T422
P423
C424
W425
I426
E427
I428
H429
L430
H431
G432
P433
L434
Q435
W436
L437
V440
L441
T442
Q443
M444
G445
S446

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α, β, γ	95.86Å 95.86Å 66.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 3.20 35.25 – 3.20	Depositor EDS
% Data completeness (in resolution range)	85.0 (50.00-3.20) 89.9 (35.25-3.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.40 (at 3.18Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.244 , 0.310 0.249 , 0.256	Depositor DCC
R_{free} test set	490 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å^2)	59.9	Xtriage
Anisotropy	0.187	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 66.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l 0.037 for h,-h-k,-l 0.025 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.87	EDS
Total number of atoms	4548	wwPDB-VP
Average B, all atoms (Å^2)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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)

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	0.49	0/1519	0.70	0/2059
1	B	0.49	0/1519	0.73	0/2060
1	C	0.45	0/1534	0.72	1/2082 (0.0%)
All	All	0.48	0/4572	0.72	1/6201 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	269	LEU	N-CA-C	-7.05	91.97	111.00

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	1482	0	1406	91	0
1	B	1481	0	1407	99	0
1	C	1495	0	1419	118	0
2	A	33	0	0	0	0
2	B	35	0	0	0	0
2	C	22	0	0	0	0
All	All	4548	0	4232	305	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 35.

All (305) 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:322:LYS:HB3	1:C:338:LEU:HB3	1.36	1.05
1:B:333:VAL:HG23	1:B:372:ILE:HG13	1.50	0.94
1:C:333:VAL:HG23	1:C:372:ILE:HG13	1.50	0.93
1:A:382:LEU:HD21	1:A:440:VAL:HG21	1.50	0.93
1:C:333:VAL:HG22	1:C:373:PHE:HB3	1.53	0.91
1:B:306:ASN:ND2	1:B:307:VAL:H	1.72	0.87
1:C:348:ARG:HD2	1:C:399:MET:HG2	1.55	0.87
1:B:352:TYR:CZ	1:B:381:LEU:HD21	2.12	0.84
1:A:268:GLU:HG3	1:A:273:VAL:HG21	1.57	0.83
1:B:310:ASN:ND2	1:B:310:ASN:H	1.74	0.83
1:B:308:ASN:N	1:B:308:ASN:HD22	1.78	0.81
1:A:310:ASN:OD1	1:A:312:THR:HG22	1.80	0.81
1:C:374:ASN:HD22	1:C:377:GLU:HB2	1.46	0.81
1:B:333:VAL:HG22	1:B:373:PHE:HB3	1.62	0.80
1:A:268:GLU:OE1	1:A:313:ILE:HD13	1.83	0.79
1:B:302:GLY:HA2	1:B:317:ARG:NH2	1.98	0.79
1:C:337:CYS:SG	1:C:366:PRO:HA	2.24	0.77
1:A:394:TYR:O	1:A:397:THR:HB	1.83	0.77
1:B:310:ASN:H	1:B:310:ASN:HD22	1.28	0.77
1:C:394:TYR:O	1:C:397:THR:HB	1.86	0.76
1:B:374:ASN:OD1	1:B:377:GLU:HB2	1.86	0.75
1:C:309:ARG:HG2	1:C:310:ASN:H	1.52	0.74
1:C:380:GLN:HB3	1:C:384:GLN:HE21	1.51	0.74
1:C:268:GLU:HA	1:C:423:PRO:O	1.88	0.73
1:B:311:SER:O	1:B:315:ASN:HB2	1.89	0.73
1:B:333:VAL:HG22	1:B:373:PHE:CB	2.19	0.73
1:B:333:VAL:CG2	1:B:372:ILE:HG13	2.19	0.73
1:C:333:VAL:CG2	1:C:372:ILE:HG13	2.19	0.73
1:C:268:GLU:HG3	1:C:273:VAL:HG21	1.69	0.72
1:A:363:LYS:O	1:A:365:PRO:HD3	1.88	0.72
1:A:333:VAL:CG2	1:A:372:ILE:HG13	2.19	0.72
1:B:310:ASN:HD22	1:B:310:ASN:N	1.87	0.72
1:C:374:ASN:HD22	1:C:377:GLU:CB	2.03	0.72
1:C:352:TYR:CZ	1:C:381:LEU:HD21	2.24	0.72
1:C:382:LEU:HD21	1:C:440:VAL:HG21	1.70	0.72
1:C:390:PHE:CZ	1:C:441:LEU:HD22	2.25	0.72
1:B:307:VAL:HG23	1:B:308:ASN:ND2	2.05	0.71
1:B:416:ARG:HB3	1:B:421:SER:CB	2.20	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:ILE:O	1:A:339:SER:HB2	1.91	0.71
1:C:352:TYR:CE1	1:C:381:LEU:HD21	2.25	0.71
1:A:390:PHE:CZ	1:A:441:LEU:HD22	2.27	0.70
1:A:370:LEU:HD12	1:A:371:LYS:H	1.55	0.70
1:C:287:ASP:O	1:C:301:LEU:HD13	1.92	0.70
1:C:263:SER:OG	1:C:429:HIS:HB2	1.92	0.69
1:A:331:GLY:O	1:A:375:ASN:HB2	1.93	0.69
1:C:333:VAL:HG22	1:C:373:PHE:CB	2.21	0.69
1:B:345:VAL:HG22	1:B:404:MET:HB3	1.74	0.69
1:B:307:VAL:HG23	1:B:308:ASN:HD22	1.57	0.69
1:A:333:VAL:HG23	1:A:372:ILE:HG13	1.75	0.68
1:B:394:TYR:O	1:B:397:THR:HB	1.92	0.68
1:C:279:CYS:HB3	1:C:284:VAL:HG11	1.76	0.68
1:A:310:ASN:O	1:A:314:GLU:HB2	1.94	0.68
1:B:310:ASN:ND2	1:B:310:ASN:N	2.42	0.68
1:A:352:TYR:CE1	1:A:381:LEU:HD21	2.30	0.67
1:A:370:LEU:HD12	1:A:371:LYS:N	2.10	0.67
1:C:287:ASP:OD1	1:C:300:CYS:HA	1.93	0.67
1:B:347:SER:HB3	1:B:350:CYS:HB3	1.76	0.66
1:B:333:VAL:HG23	1:B:372:ILE:CG1	2.25	0.66
1:B:416:ARG:HB3	1:B:421:SER:HB3	1.76	0.66
1:C:374:ASN:HD22	1:C:377:GLU:CG	2.09	0.66
1:B:370:LEU:HD12	1:B:371:LYS:N	2.11	0.66
1:A:408:LYS:HD2	1:A:416:ARG:NH2	2.12	0.65
1:C:363:LYS:O	1:C:365:PRO:HD3	1.97	0.65
1:A:408:LYS:HB3	1:A:416:ARG:CZ	2.26	0.65
1:C:264:ILE:HG23	1:C:428:ILE:HG12	1.78	0.65
1:A:310:ASN:ND2	1:A:313:ILE:HB	2.12	0.64
1:B:313:ILE:HG22	1:B:314:GLU:N	2.11	0.64
1:B:290:THR:HG22	1:B:290:THR:O	1.98	0.64
1:A:322:LYS:HB3	1:A:338:LEU:HB3	1.78	0.64
1:A:326:LEU:HD21	1:A:404:MET:CE	2.28	0.63
1:B:331:GLY:O	1:B:375:ASN:HB2	1.98	0.63
1:A:345:VAL:HG22	1:A:404:MET:HB3	1.79	0.63
1:C:266:TYR:CD2	1:C:273:VAL:HB	2.32	0.63
1:C:329:VAL:O	1:C:330:THR:HG22	1.99	0.63
1:B:260:PHE:HD1	1:B:278:HIS:HB3	1.63	0.63
1:B:404:MET:HG3	1:B:426:ILE:HG13	1.78	0.63
1:B:279:CYS:HB3	1:B:284:VAL:HG11	1.80	0.62
1:A:337:CYS:SG	1:A:343:ILE:HG13	2.40	0.62
1:A:333:VAL:HG22	1:A:373:PHE:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:337:CYS:HA	1:A:343:ILE:HD12	1.81	0.61
1:B:352:TYR:CE1	1:B:381:LEU:HD21	2.35	0.61
1:B:308:ASN:N	1:B:308:ASN:ND2	2.46	0.61
1:A:263:SER:HB3	1:A:278:HIS:HD2	1.66	0.61
1:A:400:CYS:CB	1:A:434:LEU:HG	2.30	0.61
1:A:427:GLU:HG2	1:A:429:HIS:CE1	2.37	0.60
1:C:380:GLN:HB3	1:C:384:GLN:NE2	2.15	0.60
1:C:345:VAL:HG11	1:C:372:ILE:CG2	2.32	0.60
1:B:268:GLU:HG3	1:B:273:VAL:HG21	1.82	0.60
1:B:370:LEU:HD12	1:B:371:LYS:H	1.66	0.59
1:B:437:LEU:O	1:B:441:LEU:HD12	2.01	0.59
1:A:395:GLU:C	1:A:397:THR:H	2.04	0.59
1:C:300:CYS:SG	1:C:303:GLN:HG2	2.43	0.59
1:C:334:TYR:OH	1:C:371:LYS:HD3	2.03	0.59
1:C:430:LEU:HB3	1:C:433:PRO:HD2	1.85	0.59
1:A:333:VAL:HG23	1:A:372:ILE:CG1	2.33	0.58
1:A:390:PHE:CE1	1:A:441:LEU:HD22	2.39	0.58
1:C:291:ASN:HB2	1:C:298:ARG:NH2	2.18	0.58
1:A:395:GLU:C	1:A:397:THR:N	2.58	0.57
1:C:310:ASN:HD21	1:C:313:ILE:H	1.50	0.57
1:A:310:ASN:ND2	1:A:313:ILE:H	2.01	0.57
1:C:390:PHE:CE1	1:C:441:LEU:HD22	2.40	0.57
1:C:374:ASN:ND2	1:C:377:GLU:OE2	2.38	0.57
1:C:309:ARG:NH1	1:C:317:ARG:NH1	2.53	0.57
1:A:279:CYS:HB3	1:A:284:VAL:HG11	1.87	0.57
1:B:290:THR:HG23	1:B:303:GLN:HA	1.86	0.57
1:C:290:THR:HG23	1:C:303:GLN:HB3	1.86	0.57
1:A:403:ARG:HB3	1:A:403:ARG:NH1	2.20	0.56
1:B:267:TYR:CE1	1:B:272:ARG:HB2	2.40	0.56
1:C:337:CYS:SG	1:C:343:ILE:HG13	2.44	0.56
1:C:404:MET:O	1:C:404:MET:HG3	2.06	0.56
1:C:431:HIS:O	1:C:432:GLY:C	2.43	0.56
1:A:306:ASN:HD22	1:A:307:VAL:H	1.54	0.56
1:C:344:PHE:O	1:C:404:MET:HB2	2.06	0.56
1:C:374:ASN:ND2	1:C:377:GLU:CG	2.69	0.56
1:B:291:ASN:HB3	1:B:300:CYS:SG	2.46	0.56
1:C:310:ASN:ND2	1:C:313:ILE:H	2.04	0.55
1:A:284:VAL:HG23	1:A:284:VAL:O	2.06	0.55
1:B:268:GLU:HA	1:B:423:PRO:O	2.07	0.55
1:C:393:VAL:HG13	1:C:441:LEU:HD21	1.87	0.55
1:A:326:LEU:HD21	1:A:404:MET:HE1	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:353:HIS:HB2	1:C:377:GLU:OE1	2.07	0.55
1:A:307:VAL:C	1:A:309:ARG:H	2.10	0.55
1:A:341:SER:O	1:A:342:ALA:O	2.25	0.55
1:A:404:MET:CG	1:A:426:ILE:HD11	2.36	0.55
1:B:273:VAL:HG12	1:B:304:LEU:HD13	1.89	0.55
1:C:353:HIS:C	1:C:353:HIS:CD2	2.80	0.55
1:A:333:VAL:HG22	1:A:373:PHE:CB	2.36	0.55
1:A:344:PHE:CE2	1:A:363:LYS:HB2	2.42	0.55
1:C:268:GLU:O	1:C:269:LEU:HB2	2.06	0.55
1:C:334:TYR:CZ	1:C:371:LYS:HD3	2.42	0.55
1:A:309:ARG:O	1:A:310:ASN:HB3	2.05	0.54
1:A:352:TYR:CZ	1:A:381:LEU:HD21	2.43	0.54
1:B:306:ASN:CG	1:B:307:VAL:H	2.09	0.54
1:B:310:ASN:O	1:B:314:GLU:HB2	2.08	0.54
1:B:436:TRP:O	1:B:440:VAL:HG23	2.07	0.54
1:B:320:ILE:O	1:B:339:SER:HB2	2.08	0.54
1:B:400:CYS:CB	1:B:434:LEU:HG	2.38	0.53
1:C:345:VAL:HG11	1:C:372:ILE:HG22	1.90	0.53
1:A:306:ASN:HD22	1:A:307:VAL:N	2.06	0.53
1:B:416:ARG:HB3	1:B:421:SER:HB2	1.89	0.53
1:C:262:ALA:HA	1:C:429:HIS:O	2.08	0.53
1:C:333:VAL:HG23	1:C:372:ILE:CG1	2.33	0.53
1:C:436:TRP:O	1:C:440:VAL:HG23	2.08	0.53
1:A:274:GLY:HA3	1:A:304:LEU:HD13	1.90	0.53
1:C:260:PHE:CD1	1:C:278:HIS:HB3	2.43	0.53
1:C:374:ASN:ND2	1:C:377:GLU:HB2	2.21	0.53
1:A:263:SER:HG	1:A:431:HIS:HE2	1.57	0.53
1:A:324:VAL:HG22	1:A:325:HIS:N	2.24	0.52
1:C:440:VAL:O	1:C:444:MET:HG3	2.09	0.52
1:B:322:LYS:HB3	1:B:338:LEU:HB3	1.90	0.52
1:C:374:ASN:ND2	1:C:377:GLU:H	2.08	0.52
1:C:324:VAL:HB	1:C:343:ILE:HD13	1.90	0.52
1:A:330:THR:CG2	1:A:330:THR:O	2.58	0.52
1:B:306:ASN:ND2	1:B:307:VAL:N	2.50	0.52
1:A:310:ASN:HD22	1:A:313:ILE:HB	1.74	0.52
1:A:382:LEU:HD21	1:A:440:VAL:CG2	2.32	0.51
1:C:337:CYS:HA	1:C:343:ILE:HD12	1.91	0.51
1:A:377:GLU:O	1:A:380:GLN:HB2	2.10	0.51
1:C:400:CYS:SG	1:C:433:PRO:HB2	2.51	0.51
1:A:416:ARG:HB3	1:A:421:SER:HB2	1.92	0.51
1:A:395:GLU:HG2	1:A:398:LYS:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:404:MET:O	1:A:426:ILE:HG12	2.10	0.50
1:C:281:ASN:OD1	1:C:282:ASN:N	2.44	0.50
1:B:365:PRO:O	1:B:368:CYS:HB2	2.12	0.50
1:A:410:TRP:CD1	1:A:411:GLY:N	2.80	0.50
1:C:445:GLY:O	1:C:446:SER:O	2.29	0.50
1:A:347:SER:HB3	1:A:350:CYS:HB3	1.93	0.50
1:A:412:ALA:O	1:A:413:GLU:HG2	2.12	0.49
1:C:341:SER:CB	1:C:407:VAL:HG22	2.42	0.49
1:B:404:MET:SD	1:B:426:ILE:HD11	2.52	0.49
1:B:313:ILE:CG2	1:B:314:GLU:N	2.76	0.49
1:A:306:ASN:HD21	1:A:308:ASN:HB2	1.78	0.48
1:A:400:CYS:HB3	1:A:434:LEU:HG	1.95	0.48
1:B:260:PHE:HD2	1:B:260:PHE:N	2.11	0.48
1:C:341:SER:HB2	1:C:407:VAL:HG22	1.95	0.48
1:A:339:SER:OG	1:A:341:SER:N	2.42	0.48
1:B:441:LEU:O	1:B:444:MET:HB3	2.14	0.48
1:C:397:THR:HA	1:C:437:LEU:CD2	2.43	0.48
1:B:339:SER:O	1:B:366:PRO:HB3	2.13	0.48
1:B:281:ASN:OD1	1:B:282:ASN:N	2.47	0.48
1:C:348:ARG:NH2	1:C:398:LYS:HD2	2.29	0.48
1:C:374:ASN:ND2	1:C:377:GLU:CB	2.75	0.48
1:B:260:PHE:N	1:B:260:PHE:CD2	2.81	0.48
1:B:291:ASN:ND2	1:B:300:CYS:SG	2.83	0.48
1:A:397:THR:HA	1:A:437:LEU:CD2	2.44	0.48
1:B:268:GLU:O	1:B:269:LEU:C	2.51	0.48
1:C:359:SER:O	1:C:360:THR:C	2.50	0.47
1:C:283:SER:OG	1:C:325:HIS:NE2	2.39	0.47
1:B:363:LYS:O	1:B:365:PRO:HD3	2.15	0.47
1:B:380:GLN:O	1:B:381:LEU:C	2.53	0.47
1:C:344:PHE:HA	1:C:362:CYS:O	2.14	0.47
1:B:376:GLN:NE2	1:B:380:GLN:HE21	2.12	0.47
1:A:310:ASN:HD22	1:A:310:ASN:N	2.11	0.47
1:B:353:HIS:C	1:B:353:HIS:CD2	2.87	0.47
1:B:397:THR:HA	1:B:437:LEU:CD2	2.45	0.47
1:C:353:HIS:C	1:C:355:GLY:H	2.18	0.47
1:C:416:ARG:HB3	1:C:421:SER:HB3	1.97	0.47
1:C:335:ALA:HB2	1:C:372:ILE:HD13	1.97	0.47
1:A:310:ASN:CG	1:A:312:THR:HG22	2.34	0.46
1:B:269:LEU:HD23	1:B:269:LEU:HA	1.77	0.46
1:B:337:CYS:HA	1:B:343:ILE:HD12	1.97	0.46
1:B:395:GLU:HG2	1:B:398:LYS:HE3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:431:HIS:O	1:A:432:GLY:C	2.52	0.46
1:C:374:ASN:ND2	1:C:377:GLU:HG3	2.31	0.46
1:A:334:TYR:CZ	1:A:371:LYS:HD3	2.50	0.46
1:C:274:GLY:HA3	1:C:304:LEU:HD22	1.98	0.46
1:A:327:TYR:HB2	1:A:334:TYR:HB2	1.98	0.46
1:C:291:ASN:HB2	1:C:298:ARG:CZ	2.46	0.46
1:B:334:TYR:CE1	1:B:371:LYS:HB2	2.51	0.46
1:C:285:ILE:N	1:C:285:ILE:HD12	2.31	0.46
1:C:301:LEU:O	1:C:317:ARG:NH2	2.49	0.46
1:C:345:VAL:HG11	1:C:372:ILE:HG21	1.97	0.45
1:B:394:TYR:CD1	1:B:394:TYR:C	2.88	0.45
1:B:376:GLN:HE22	1:B:380:GLN:NE2	2.13	0.45
1:A:416:ARG:HG3	1:A:416:ARG:NH1	2.30	0.45
1:A:357:HIS:CE1	1:B:322:LYS:NZ	2.84	0.45
1:B:381:LEU:O	1:B:385:SER:HB2	2.16	0.45
1:C:266:TYR:HD2	1:C:273:VAL:HB	1.79	0.45
1:B:404:MET:HG3	1:B:426:ILE:CG1	2.46	0.45
1:A:365:PRO:O	1:A:368:CYS:HB2	2.17	0.45
1:B:400:CYS:HB3	1:B:434:LEU:HG	1.99	0.44
1:B:300:CYS:HB2	1:B:303:GLN:HG2	1.99	0.44
1:B:356:PHE:CE2	1:C:310:ASN:OD1	2.70	0.44
1:C:281:ASN:OD1	1:C:283:SER:N	2.49	0.44
1:C:414:TYR:O	1:C:417:GLN:NE2	2.50	0.44
1:B:306:ASN:CG	1:B:307:VAL:N	2.70	0.44
1:B:320:ILE:O	1:B:339:SER:CB	2.65	0.44
1:C:285:ILE:HD12	1:C:285:ILE:H	1.83	0.44
1:C:418:ASP:OD2	1:C:419:VAL:N	2.51	0.44
1:B:266:TYR:CD2	1:B:301:LEU:HD22	2.53	0.44
1:A:307:VAL:C	1:A:309:ARG:N	2.70	0.44
1:B:309:ARG:HB3	1:B:314:GLU:OE1	2.18	0.44
1:B:352:TYR:OH	1:B:381:LEU:HD21	2.16	0.44
1:B:382:LEU:HD13	1:B:437:LEU:HD12	2.00	0.44
1:C:410:TRP:O	1:C:414:TYR:HD2	2.01	0.44
1:A:266:TYR:OH	1:A:424:CYS:HB3	2.18	0.43
1:C:337:CYS:HA	1:C:343:ILE:CD1	2.48	0.43
1:A:331:GLY:C	1:A:332:GLU:HG2	2.38	0.43
1:A:386:VAL:CG2	1:A:440:VAL:HG11	2.48	0.43
1:C:309:ARG:CG	1:C:310:ASN:H	2.24	0.43
1:C:315:ASN:HD22	1:C:315:ASN:HA	1.57	0.43
1:B:378:PHE:O	1:B:379:ALA:C	2.57	0.43
1:B:350:CYS:O	1:B:350:CYS:SG	2.77	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:ASN:HD22	1:B:315:ASN:HA	1.50	0.43
1:C:311:SER:O	1:C:315:ASN:HB2	2.18	0.43
1:C:400:CYS:CB	1:C:434:LEU:HG	2.48	0.43
1:C:441:LEU:O	1:C:444:MET:HB2	2.18	0.43
1:B:268:GLU:HG3	1:B:273:VAL:CG2	2.47	0.43
1:B:268:GLU:O	1:B:269:LEU:HB2	2.19	0.43
1:B:277:PHE:CD1	1:B:299:CYS:HB2	2.53	0.43
1:B:403:ARG:HG2	1:B:403:ARG:HH11	1.84	0.43
1:C:320:ILE:O	1:C:339:SER:HB2	2.17	0.43
1:A:436:TRP:O	1:A:440:VAL:HG23	2.18	0.43
1:C:310:ASN:HD21	1:C:313:ILE:CB	2.32	0.43
1:A:316:THR:HG21	1:A:423:PRO:HG2	2.01	0.43
1:C:310:ASN:ND2	1:C:310:ASN:N	2.67	0.43
1:B:261:TRP:CE3	1:B:262:ALA:HB2	2.54	0.43
1:A:402:ILE:HD12	1:A:430:LEU:CD2	2.49	0.42
1:A:306:ASN:ND2	1:A:307:VAL:N	2.68	0.42
1:C:348:ARG:CD	1:C:399:MET:HG2	2.39	0.42
1:B:344:PHE:CE2	1:B:363:LYS:HB2	2.55	0.42
1:C:370:LEU:HD12	1:C:371:LYS:N	2.34	0.42
1:C:300:CYS:C	1:C:301:LEU:HD12	2.40	0.42
1:C:310:ASN:HD21	1:C:313:ILE:HG13	1.84	0.42
1:C:345:VAL:HG22	1:C:404:MET:HB3	2.01	0.42
1:C:437:LEU:HD12	1:C:437:LEU:HA	1.88	0.42
1:A:380:GLN:O	1:A:384:GLN:HB2	2.20	0.42
1:C:310:ASN:C	1:C:310:ASN:HD22	2.23	0.42
1:B:376:GLN:NE2	1:B:380:GLN:NE2	2.68	0.41
1:B:309:ARG:HG2	1:B:313:ILE:HG21	2.03	0.41
1:C:314:GLU:OE1	1:C:314:GLU:HA	2.21	0.41
1:B:302:GLY:HA2	1:B:317:ARG:HH22	1.81	0.41
1:B:319:HIS:O	1:B:341:SER:HB2	2.21	0.41
1:C:263:SER:HA	1:C:277:PHE:O	2.21	0.41
1:C:281:ASN:ND2	1:C:297:ASP:OD2	2.53	0.41
1:C:386:VAL:HG13	1:C:387:ASN:N	2.36	0.41
1:A:262:ALA:HB2	1:A:430:LEU:HD12	2.02	0.41
1:A:309:ARG:HG3	1:A:313:ILE:HG21	2.03	0.41
1:B:263:SER:HB2	1:B:278:HIS:HD2	1.86	0.41
1:C:321:GLY:C	1:C:323:GLY:H	2.24	0.41
1:A:268:GLU:OE2	1:A:309:ARG:NH2	2.53	0.41
1:C:349:ASN:HD22	1:C:399:MET:CE	2.33	0.41
1:A:285:ILE:CG2	1:A:286:VAL:N	2.84	0.41
1:A:397:THR:HA	1:A:437:LEU:HD21	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:274:GLY:CA	1:C:304:LEU:HD22	2.50	0.41
1:C:316:THR:HG21	1:C:423:PRO:HG2	2.02	0.41
1:C:395:GLU:C	1:C:397:THR:N	2.71	0.41
1:B:273:VAL:O	1:B:274:GLY:O	2.38	0.41
1:B:442:THR:C	1:B:444:MET:N	2.73	0.41
1:C:261:TRP:CE3	1:C:262:ALA:HB2	2.55	0.41
1:C:306:ASN:ND2	1:C:307:VAL:H	2.19	0.41
1:C:353:HIS:CB	1:C:377:GLU:OE1	2.68	0.41
1:A:276:VAL:O	1:A:276:VAL:HG13	2.20	0.41
1:B:354:HIS:C	1:C:311:SER:HB2	2.42	0.41
1:A:268:GLU:OE2	1:A:317:ARG:HD2	2.21	0.40
1:A:337:CYS:HA	1:A:343:ILE:CD1	2.48	0.40
1:A:393:VAL:HG11	1:A:444:MET:SD	2.61	0.40
1:C:321:GLY:O	1:C:323:GLY:N	2.54	0.40
1:B:419:VAL:HG22	1:B:425:TRP:CD2	2.56	0.40
1:C:264:ILE:HA	1:C:427:GLU:O	2.21	0.40
1:A:310:ASN:HD21	1:A:313:ILE:H	1.70	0.40
1:A:386:VAL:O	1:A:444:MET:CE	2.69	0.40
1:B:277:PHE:CG	1:B:299:CYS:HB2	2.56	0.40
1:A:329:VAL:C	1:A:331:GLY:H	2.25	0.40
1:C:310:ASN:ND2	1:C:313:ILE:HB	2.36	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	182/188 (97%)	153 (84%)	20 (11%)	9 (5%)	2 17
1	B	182/188 (97%)	148 (81%)	30 (16%)	4 (2%)	6 35
1	C	186/188 (99%)	151 (81%)	28 (15%)	7 (4%)	3 22
All	All	550/564 (98%)	452 (82%)	78 (14%)	20 (4%)	3 23

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	SER
1	A	342	ALA
1	B	342	ALA
1	C	263	SER
1	C	296	SER
1	C	310	ASN
1	C	342	ALA
1	A	330	THR
1	A	412	ALA
1	B	274	GLY
1	A	295	ASN
1	C	388	ASN
1	C	412	ALA
1	A	274	GLY
1	A	309	ARG
1	A	310	ASN
1	B	412	ALA
1	C	322	LYS
1	B	407	VAL
1	A	407	VAL

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	165/167 (99%)	146 (88%)	19 (12%)	5 24
1	B	165/167 (99%)	140 (85%)	25 (15%)	3 13
1	C	167/167 (100%)	152 (91%)	15 (9%)	9 34
All	All	497/501 (99%)	438 (88%)	59 (12%)	5 22

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	263	SER

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Mol	Chain	Res	Type
1	A	279	CYS
1	A	280	ASN
1	A	291	ASN
1	A	294	ASN
1	A	300	CYS
1	A	306	ASN
1	A	310	ASN
1	A	329	VAL
1	A	339	SER
1	A	377	GLU
1	A	397	THR
1	A	403	ARG
1	A	410	TRP
1	A	417	GLN
1	A	424	CYS
1	A	426	ILE
1	A	430	LEU
1	A	443	GLN
1	B	279	CYS
1	B	280	ASN
1	B	291	ASN
1	B	295	ASN
1	B	299	CYS
1	B	303	GLN
1	B	308	ASN
1	B	310	ASN
1	B	312	THR
1	B	313	ILE
1	B	315	ASN
1	B	329	VAL
1	B	330	THR
1	B	348	ARG
1	B	371	LYS
1	B	376	GLN
1	B	377	GLU
1	B	385	SER
1	B	387	ASN
1	B	394	TYR
1	B	397	THR
1	B	404	MET
1	B	424	CYS
1	B	430	LEU

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Mol	Chain	Res	Type
1	B	435	GLN
1	C	264	ILE
1	C	279	CYS
1	C	306	ASN
1	C	310	ASN
1	C	315	ASN
1	C	329	VAL
1	C	353	HIS
1	C	397	THR
1	C	413	GLU
1	C	424	CYS
1	C	426	ILE
1	C	430	LEU
1	C	436	TRP
1	C	443	GLN
1	C	446	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	278	HIS
1	A	294	ASN
1	A	306	ASN
1	A	310	ASN
1	A	315	ASN
1	A	357	HIS
1	A	376	GLN
1	A	417	GLN
1	A	429	HIS
1	B	278	HIS
1	B	291	ASN
1	B	306	ASN
1	B	308	ASN
1	B	310	ASN
1	B	315	ASN
1	B	351	ASN
1	B	376	GLN
1	B	380	GLN
1	B	417	GLN
1	C	294	ASN
1	C	306	ASN
1	C	308	ASN

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Mol	Chain	Res	Type
1	C	310	ASN
1	C	315	ASN
1	C	349	ASN
1	C	351	ASN
1	C	374	ASN
1	C	384	GLN
1	C	387	ASN
1	C	435	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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	186/188 (98%)	-0.24	1 (0%)	91	86	11, 42, 100, 166	0
1	B	186/188 (98%)	-0.04	2 (1%)	80	69	9, 53, 109, 148	0
1	C	188/188 (100%)	0.58	30 (15%)	1	1	22, 66, 181, 198	0
All	All	560/564 (99%)	0.10	33 (5%)	22	13	9, 54, 148, 198	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	308	ASN	9.0
1	C	306	ASN	9.0
1	C	294	ASN	8.8
1	C	297	ASP	7.6
1	C	295	ASN	7.1
1	C	296	SER	5.3
1	C	299	CYS	5.0
1	C	303	GLN	4.1
1	C	291	ASN	3.9
1	C	298	ARG	3.9
1	C	309	ARG	3.8
1	C	310	ASN	3.7
1	C	293	SER	3.4
1	C	290	THR	3.4
1	C	302	GLY	3.2
1	C	314	GLU	3.1
1	A	294	ASN	3.1
1	C	446	SER	3.0
1	C	312	THR	2.9
1	C	319	HIS	2.8
1	B	416	ARG	2.8
1	C	292	PRO	2.4
1	C	301	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	289	PHE	2.3
1	C	300	CYS	2.3
1	B	307	VAL	2.3
1	C	305	SER	2.2
1	C	311	SER	2.2
1	C	288	GLY	2.2
1	C	318	ARG	2.1
1	C	415	HIS	2.1
1	C	307	VAL	2.1
1	C	277	PHE	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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.