



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:27 PM GMT

PDB ID : 4F0P
Title : MspJI Restriction Endonuclease - P31 Form
Authors : Horton, J.R.; Mabuchi, M.; Cohen-Karni, D.; Zhang, X.; Griggs, R.; Samaranyake, M.; Roberts, R.J.; Zheng, Y.; Cheng, X.
Deposited on : 2012-05-04
Resolution : 2.79 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

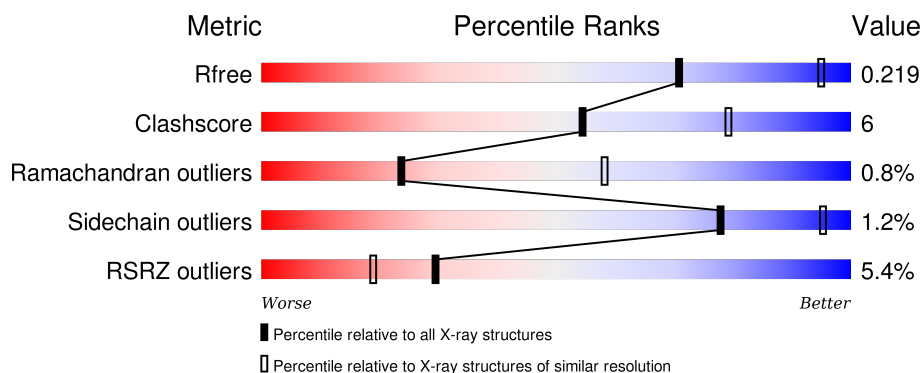
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 ≥ 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	456	<div> <div>3%</div> <div>82%</div> <div>16%</div> <div>.</div> </div>
1	B	456	<div> <div>2%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>
1	C	456	<div> <div>8%</div> <div>80%</div> <div>17%</div> <div>..</div> </div>
1	D	456	<div> <div>8%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13639 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 Restriction endonuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	0	0
			3402	2132	633	632	5			
1	B	452	Total	C	N	O	S	0	0	0
			3426	2154	637	630	5			
1	C	449	Total	C	N	O	S	0	0	0
			3257	2048	598	606	5			
1	D	452	Total	C	N	O	S	0	0	0
			3256	2052	600	600	4			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

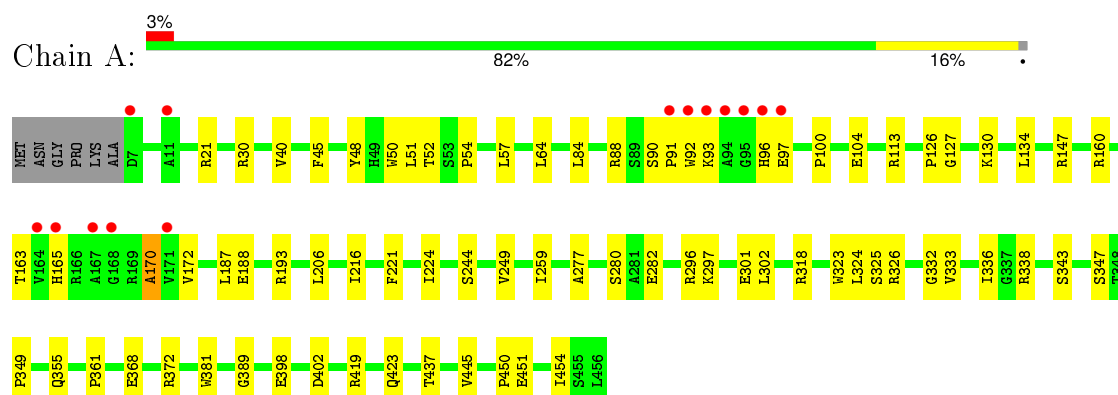
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	113	Total	O	0	0
			113	113		
3	B	111	Total	O	0	0
			111	111		
3	C	45	Total	O	0	0
			45	45		
3	D	27	Total	O	0	0
			27	27		

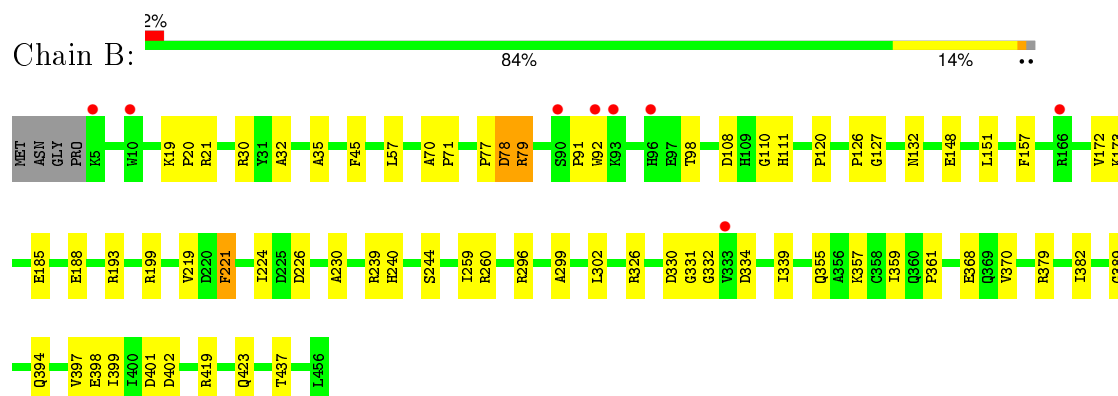
3 Residue-property plots [i](#)

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.

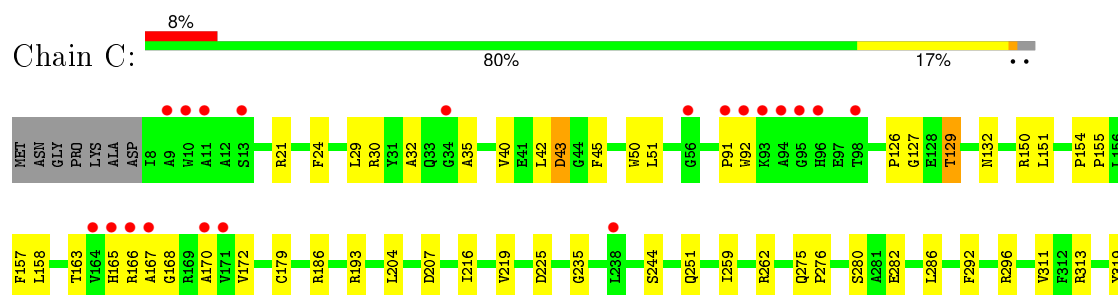
• Molecule 1: Restriction endonuclease

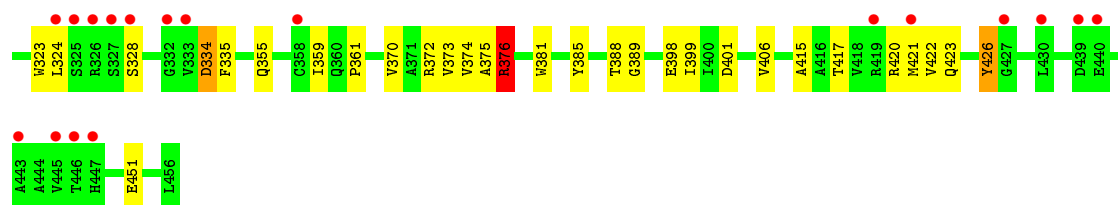


• Molecule 1: Restriction endonuclease

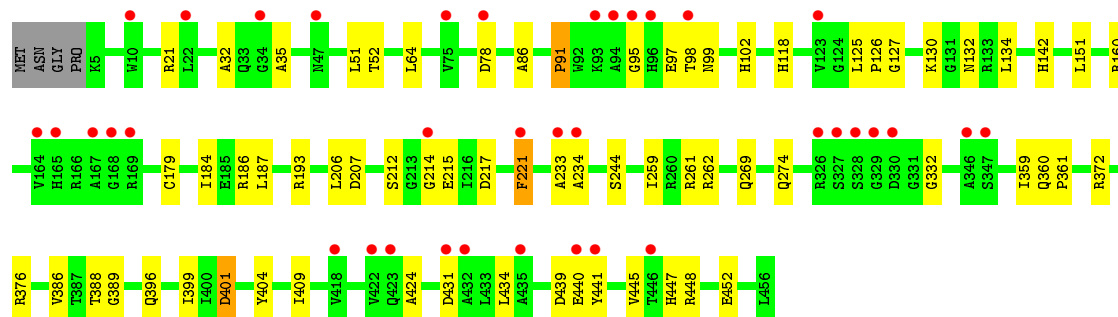
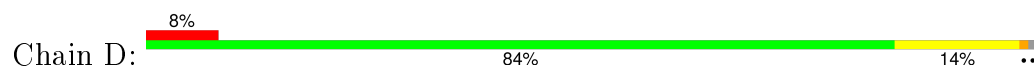


• Molecule 1: Restriction endonuclease





● Molecule 1: Restriction endonuclease



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	144.95Å 144.95Å 101.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.65 – 2.79 34.65 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.9 (34.65-2.79) 99.1 (34.65-2.79)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.98 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.164 , 0.221 0.163 , 0.219	Depositor DCC
R_{free} test set	2945 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 24.6	EDS
Estimated twinning fraction	0.059 for -h,-k,l 0.199 for h,-h-k,-l 0.060 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Outliers	0 of 59272 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13639	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: MG

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	0.21	0/3472	0.39	0/4721
1	B	0.22	0/3500	0.38	0/4761
1	C	0.21	0/3323	0.41	1/4537 (0.0%)
1	D	0.20	0/3325	0.38	1/4544 (0.0%)
All	All	0.21	0/13620	0.39	2/18563 (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	D	91	PRO	N-CA-CB	5.87	110.34	103.30
1	C	376	ARG	NE-CZ-NH1	5.61	123.11	120.30

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	3402	0	3328	43	0
1	B	3426	0	3351	41	0
1	C	3257	0	3081	55	0
1	D	3256	0	3056	41	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
3	A	113	0	0	0	0
3	B	111	0	0	1	0
3	C	45	0	0	1	0
3	D	27	0	0	1	0
All	All	13639	0	12816	167	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 6.

All (167) 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:186:ARG:HB3	1:C:207:ASP:HB2	1.67	0.77
1:D:186:ARG:HB3	1:D:207:ASP:HB2	1.68	0.76
1:B:379:ARG:NH2	1:C:375:ALA:O	2.20	0.74
1:B:185:GLU:OE1	1:B:260:ARG:NH2	2.21	0.71
1:C:165:HIS:HA	1:C:170:ALA:HA	1.73	0.71
1:C:21:ARG:NH2	1:C:225:ASP:OD2	2.26	0.67
1:B:368:GLU:HG3	1:D:372:ARG:HD2	1.77	0.67
1:B:326:ARG:NH2	1:B:334:ASP:OD1	2.28	0.67
1:D:269:GLN:H	1:D:396:GLN:HE22	1.43	0.65
1:A:163:THR:HG22	1:A:172:VAL:HG23	1.78	0.65
1:D:212:SER:HB3	1:D:261:ARG:HA	1.78	0.65
1:D:274:GLN:NE2	1:D:396:GLN:OE1	2.31	0.63
1:C:262:ARG:NH2	1:C:401:ASP:OD2	2.33	0.62
1:B:172:VAL:HG13	1:B:173:LYS:HG2	1.82	0.60
1:C:370:VAL:HG11	1:C:399:ILE:HD11	1.84	0.60
1:C:275:GLN:NE2	1:C:361:PRO:O	2.35	0.60
1:C:313:ARG:NH1	1:C:319:TYR:O	2.35	0.59
1:B:402:ASP:OD1	1:D:376:ARG:NH1	2.36	0.59
1:C:361:PRO:HA	1:C:389:GLY:HA2	1.83	0.59
1:A:332:GLY:O	1:A:355:GLN:NE2	2.37	0.58
1:B:361:PRO:HA	1:B:389:GLY:HA2	1.85	0.57
1:C:150:ARG:NH2	3:C:524:HOH:O	2.37	0.56
1:B:332:GLY:O	1:B:355:GLN:NE2	2.37	0.56
1:A:84:LEU:HD13	1:A:224:ILE:HD13	1.88	0.56
1:C:376:ARG:CG	1:C:376:ARG:HH11	2.18	0.55
1:B:32:ALA:HB3	1:B:35:ALA:HB2	1.87	0.55
1:D:332:GLY:HA3	1:D:376:ARG:HE	1.71	0.55
1:C:292:PHE:O	1:C:296:ARG:NH1	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:361:PRO:HA	1:D:389:GLY:HA2	1.88	0.54
1:A:244:SER:HB2	1:A:259:ILE:HB	1.90	0.54
1:C:311:VAL:HG22	1:C:417:THR:HG21	1.90	0.54
1:D:359:ILE:O	1:D:388:THR:OG1	2.21	0.54
1:C:334:ASP:HB3	1:C:355:GLN:HG2	1.90	0.54
1:A:402:ASP:OD2	1:C:376:ARG:NH1	2.42	0.53
1:A:398:GLU:OE1	1:C:372:ARG:NH2	2.40	0.53
1:C:359:ILE:O	1:C:388:THR:OG1	2.22	0.53
1:D:95:GLY:HA3	1:D:98:THR:HG23	1.90	0.53
1:B:120:PRO:O	1:B:199:ARG:NH1	2.41	0.53
1:D:52:THR:HG22	1:D:221:PHE:HE2	1.74	0.53
1:C:166:ARG:O	1:C:168:GLY:N	2.38	0.53
1:A:188:GLU:HA	1:B:188:GLU:HA	1.91	0.53
1:D:386:VAL:HG22	1:D:409:ILE:HD12	1.91	0.53
1:B:370:VAL:HG11	1:B:399:ILE:HD11	1.91	0.52
1:C:29:LEU:HD21	1:C:51:LEU:HD23	1.91	0.52
1:A:318:ARG:HH22	1:B:111:HIS:HB2	1.72	0.52
1:D:32:ALA:HB3	1:D:35:ALA:HB2	1.90	0.52
1:A:147:ARG:HG3	1:A:249:VAL:HG12	1.92	0.52
1:B:357:LYS:HG3	1:B:359:ILE:HG23	1.91	0.52
1:A:302:LEU:HD23	1:A:437:THR:HA	1.92	0.51
1:D:151:LEU:HB3	1:D:234:ALA:HB1	1.92	0.51
1:D:193:ARG:NH2	3:D:501:HOH:O	2.42	0.51
1:A:326:ARG:NH2	1:A:333:VAL:O	2.44	0.51
1:A:91:PRO:HG3	1:A:160:ARG:HD2	1.93	0.51
1:B:127:GLY:HA3	1:B:132:ASN:HB2	1.92	0.51
1:B:108:ASP:OD1	1:C:193:ARG:NH2	2.39	0.51
1:C:381:TRP:HZ2	1:C:451:GLU:HG3	1.76	0.51
1:A:48:TYR:CZ	1:A:52:THR:HG21	2.46	0.50
1:D:78:ASP:HB2	1:D:233:ALA:HA	1.93	0.50
1:A:40:VAL:HG12	1:A:50:TRP:CE2	2.47	0.50
1:D:244:SER:HB2	1:D:259:ILE:HB	1.94	0.50
1:B:368:GLU:OE2	1:D:372:ARG:NH1	2.45	0.50
1:D:99:ASN:O	1:D:102:HIS:ND1	2.26	0.50
1:C:376:ARG:HH11	1:C:376:ARG:HG2	1.76	0.49
1:A:30:ARG:HB2	1:A:45:PHE:HB2	1.95	0.49
1:D:262:ARG:NH2	1:D:401:ASP:OD2	2.46	0.49
1:C:24:PHE:CE2	1:C:216:ILE:HD11	2.48	0.48
1:D:187:LEU:HD12	1:D:206:LEU:HD11	1.95	0.48
1:C:30:ARG:HG2	1:C:42:LEU:HB3	1.95	0.48
1:C:32:ALA:HB3	1:C:35:ALA:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:THR:HG23	1:C:172:VAL:HA	1.94	0.48
1:A:91:PRO:O	1:A:93:LYS:N	2.42	0.48
1:B:77:PRO:HD3	1:B:230:ALA:O	2.14	0.48
1:C:324:LEU:HA	1:C:335:PHE:HA	1.96	0.48
1:B:110:GLY:O	1:B:260:ARG:NH1	2.46	0.48
1:B:126:PRO:HA	1:B:127:GLY:HA2	1.46	0.48
1:C:43:ASP:OD1	1:C:43:ASP:N	2.47	0.48
1:A:216:ILE:HG22	1:D:125:LEU:HD21	1.95	0.48
1:D:99:ASN:HB2	1:D:102:HIS:HB3	1.95	0.48
1:C:129:THR:HG23	1:C:132:ASN:HB2	1.96	0.48
1:C:376:ARG:HH11	1:C:376:ARG:HB3	1.79	0.47
1:C:417:THR:HA	1:C:420:ARG:HE	1.79	0.47
1:A:126:PRO:HA	1:A:127:GLY:HA2	1.45	0.47
1:D:448:ARG:NH2	1:D:452:GLU:HB3	2.28	0.47
1:C:244:SER:HB3	1:C:259:ILE:HB	1.95	0.47
1:D:399:ILE:HG23	1:D:404:TYR:HB2	1.96	0.47
1:C:376:ARG:CB	1:C:376:ARG:HH11	2.27	0.47
1:C:374:VAL:HG22	1:C:406:VAL:HG21	1.96	0.47
1:A:318:ARG:NH2	1:B:111:HIS:HB2	2.30	0.47
1:C:323:TRP:HD1	1:C:324:LEU:O	1.98	0.47
1:D:439:ASP:O	1:D:441:TYR:N	2.48	0.46
1:C:30:ARG:HB2	1:C:45:PHE:HB2	1.96	0.46
1:B:30:ARG:HB2	1:B:45:PHE:HB2	1.97	0.46
1:A:221:PHE:CE2	1:A:224:ILE:HD12	2.50	0.46
1:C:280:SER:HB3	1:C:282:GLU:HG2	1.96	0.46
1:B:21:ARG:NH2	1:B:221:PHE:HB3	2.30	0.46
1:B:226:ASP:OD2	1:B:239:ARG:NE	2.41	0.46
1:B:244:SER:HB2	1:B:259:ILE:HB	1.95	0.46
1:C:276:PRO:HB3	1:C:282:GLU:HG3	1.97	0.46
1:A:381:TRP:HZ2	1:A:451:GLU:HG3	1.81	0.46
1:A:372:ARG:NH2	1:C:398:GLU:OE1	2.40	0.46
1:D:142:HIS:CD2	1:D:184:ILE:H	2.34	0.46
1:B:78:ASP:O	1:B:79:ARG:HB3	2.15	0.46
1:A:324:LEU:HB3	1:A:445:VAL:HG22	1.97	0.46
1:B:239:ARG:NH2	3:B:672:HOH:O	2.36	0.46
1:B:239:ARG:HG3	1:B:240:HIS:ND1	2.31	0.46
1:D:97:GLU:N	1:D:98:THR:HA	2.32	0.45
1:A:64:LEU:HG	1:A:88:ARG:HB2	1.99	0.45
1:A:361:PRO:HA	1:A:389:GLY:HA2	1.99	0.45
1:C:417:THR:HG22	1:C:420:ARG:HH21	1.80	0.45
1:B:330:ASP:OD1	1:B:331:GLY:N	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:286:LEU:HB2	1:C:415:ALA:HB2	2.00	0.44
1:C:126:PRO:HB2	1:C:204:LEU:HD13	1.99	0.44
1:B:19:LYS:HA	1:B:20:PRO:HD3	1.83	0.44
1:A:277:ALA:HB3	1:A:280:SER:HB3	1.99	0.44
1:B:157:PHE:CE1	1:B:224:ILE:HD11	2.52	0.44
1:C:157:PHE:HE2	1:C:219:VAL:HG21	1.82	0.44
1:D:445:VAL:O	1:D:447:HIS:N	2.43	0.44
1:B:302:LEU:HD23	1:B:437:THR:HA	2.00	0.44
1:C:158:LEU:HB2	1:C:179:CYS:HB2	1.99	0.43
1:A:187:LEU:HG	1:A:206:LEU:HG	1.99	0.43
1:A:282:GLU:HG2	1:A:419:ARG:HD3	2.00	0.43
1:B:394:GLN:HA	1:B:397:VAL:HG22	1.99	0.43
1:D:126:PRO:HA	1:D:127:GLY:HA2	1.44	0.43
1:D:21:ARG:HG2	1:D:51:LEU:HD12	2.01	0.43
1:A:349:PRO:HG3	1:A:454:ILE:HG23	2.00	0.43
1:D:160:ARG:HB2	1:D:179:CYS:SG	2.59	0.43
1:D:431:ASP:HA	1:D:434:LEU:HB2	2.01	0.43
1:A:297:LYS:O	1:A:301:GLU:HG3	2.19	0.43
1:A:21:ARG:HG2	1:A:51:LEU:HD12	2.00	0.43
1:B:193:ARG:NH1	1:B:423:GLN:OE1	2.51	0.43
1:A:193:ARG:NH1	1:A:423:GLN:OE1	2.46	0.43
1:C:40:VAL:HG22	1:C:50:TRP:CE2	2.53	0.42
1:C:376:ARG:NH1	1:C:376:ARG:CG	2.83	0.42
1:A:338:ARG:HH11	1:A:454:ILE:HA	1.84	0.42
1:C:126:PRO:HA	1:C:127:GLY:HA2	1.56	0.42
1:A:54:PRO:HD2	1:A:57:LEU:HD12	2.01	0.42
1:B:398:GLU:HG2	1:D:376:ARG:HH22	1.85	0.42
1:D:214:GLY:HA2	1:D:215:GLU:HA	1.75	0.42
1:C:376:ARG:NH1	1:C:376:ARG:HG2	2.34	0.42
1:B:70:ALA:HA	1:B:71:PRO:HD3	1.94	0.42
1:C:91:PRO:HA	1:C:92:TRP:HA	1.77	0.42
1:D:269:GLN:H	1:D:396:GLN:NE2	2.14	0.42
1:A:324:LEU:O	1:A:326:ARG:N	2.51	0.42
1:A:323:TRP:NE1	1:A:336:ILE:HD12	2.34	0.42
1:A:165:HIS:HA	1:A:170:ALA:HA	2.01	0.41
1:D:130:LYS:O	1:D:134:LEU:HG	2.21	0.41
1:C:422:VAL:O	1:C:426:TYR:HB2	2.21	0.41
1:A:104:GLU:HG2	1:A:113:ARG:HB2	2.03	0.41
1:B:98:THR:O	1:B:419:ARG:NH2	2.53	0.41
1:D:360:GLN:HA	1:D:361:PRO:HD3	1.95	0.41
1:B:339:ILE:HD12	1:B:382:ILE:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:64:LEU:HD13	1:D:86:ALA:HB1	2.03	0.41
1:A:100:PRO:HG2	1:A:423:GLN:HG3	2.03	0.41
1:C:373:VAL:HG21	1:C:385:TYR:HD2	1.86	0.41
1:B:148:GLU:O	1:B:151:LEU:HB2	2.21	0.41
1:B:296:ARG:HB3	1:B:299:ALA:HB3	2.03	0.41
1:D:118:HIS:HB2	1:D:132:ASN:HD21	1.85	0.41
1:B:157:PHE:HE2	1:B:219:VAL:HG21	1.85	0.41
1:A:336:ILE:HD13	1:A:450:PRO:HA	2.03	0.41
1:D:359:ILE:HA	1:D:359:ILE:HD12	1.95	0.40
1:A:323:TRP:HE1	1:A:336:ILE:HD12	1.86	0.40
1:A:130:LYS:HE2	1:A:134:LEU:HD11	2.03	0.40
1:C:422:VAL:HG13	1:C:423:GLN:H	1.86	0.40
1:C:417:THR:O	1:C:421:MET:HB2	2.22	0.40
1:A:343:SER:O	1:A:347:SER:HB3	2.22	0.40
1:C:154:PRO:HA	1:C:155:PRO:HD3	1.93	0.40
1:C:151:LEU:HD11	1:C:235:GLY:HA2	2.02	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	448/456 (98%)	428 (96%)	15 (3%)	5 (1%)	17	50
1	B	450/456 (99%)	434 (96%)	14 (3%)	2 (0%)	39	74
1	C	447/456 (98%)	409 (92%)	35 (8%)	3 (1%)	26	62
1	D	450/456 (99%)	428 (95%)	18 (4%)	4 (1%)	21	55
All	All	1795/1824 (98%)	1699 (95%)	82 (5%)	14 (1%)	24	58

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	HIS
1	C	328	SER
1	D	91	PRO
1	B	79	ARG
1	A	92	TRP
1	C	167	ALA
1	D	440	GLU
1	A	97	GLU
1	A	325	SER
1	C	251	GLN
1	D	217	ASP
1	A	170	ALA
1	D	424	ALA
1	B	91	PRO

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	339/358 (95%)	336 (99%)	3 (1%)	84	96
1	B	339/358 (95%)	334 (98%)	5 (2%)	72	93
1	C	303/358 (85%)	298 (98%)	5 (2%)	68	92
1	D	297/358 (83%)	295 (99%)	2 (1%)	88	97
All	All	1278/1432 (89%)	1263 (99%)	15 (1%)	78	95

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	SER
1	A	296	ARG
1	A	368	GLU
1	B	57	LEU
1	B	78	ASP
1	B	92	TRP
1	B	221	PHE
1	B	401	ASP

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Mol	Chain	Res	Type
1	C	43	ASP
1	C	129	THR
1	C	334	ASP
1	C	376	ARG
1	C	426	TYR
1	D	221	PHE
1	D	401	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules 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.

No monomer is involved in short contacts.

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, 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	450/456 (98%)	0.13	14 (3%)	52 40	22, 45, 89, 201	0
1	B	452/456 (99%)	0.04	8 (1%)	71 61	24, 46, 86, 137	1 (0%)
1	C	449/456 (98%)	0.60	38 (8%)	13 6	34, 70, 123, 204	0
1	D	452/456 (99%)	0.61	37 (8%)	14 7	32, 79, 123, 157	0
All	All	1803/1824 (98%)	0.34	97 (5%)	29 19	22, 59, 114, 204	1 (0%)

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	95	GLY	32.8
1	C	94	ALA	16.8
1	A	95	GLY	14.9
1	D	167	ALA	11.1
1	C	96	HIS	10.7
1	D	96	HIS	10.1
1	A	94	ALA	9.8
1	D	328	SER	9.5
1	C	93	LYS	9.0
1	D	95	GLY	7.6
1	A	96	HIS	7.3
1	C	98	THR	6.6
1	C	9	ALA	6.5
1	A	93	LYS	6.0
1	D	446	THR	5.9
1	A	92	TRP	5.9
1	C	327	SER	5.7
1	C	171	VAL	5.7
1	A	91	PRO	5.5
1	D	94	ALA	5.2
1	D	329	GLY	5.2

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Mol	Chain	Res	Type	RSRZ
1	C	445	VAL	5.2
1	C	427	GLY	5.1
1	A	167	ALA	5.0
1	C	92	TRP	5.0
1	D	441	TYR	5.0
1	C	167	ALA	5.0
1	C	170	ALA	4.9
1	C	166	ARG	4.9
1	D	326	ARG	4.8
1	C	10	TRP	4.7
1	C	165	HIS	4.6
1	D	168	GLY	4.6
1	C	440	GLU	4.6
1	D	327	SER	4.5
1	D	440	GLU	4.3
1	D	432	ALA	4.1
1	C	11	ALA	3.8
1	D	435	ALA	3.7
1	C	443	ALA	3.7
1	C	439	ASP	3.5
1	D	123	VAL	3.4
1	C	326	ARG	3.3
1	C	430	LEU	3.3
1	C	332	GLY	3.3
1	D	10	TRP	3.3
1	D	347	SER	3.3
1	C	328	SER	3.2
1	C	333	VAL	3.1
1	B	96	HIS	3.1
1	C	91	PRO	3.1
1	A	168	GLY	3.1
1	C	13	SER	3.1
1	B	90	SER	3.1
1	D	431	ASP	3.0
1	B	166	ARG	3.0
1	D	75	VAL	3.0
1	D	165	HIS	2.9
1	C	325	SER	2.9
1	D	169	ARG	2.9
1	A	171	VAL	2.9
1	C	34	GLY	2.8
1	A	165	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	234	ALA	2.8
1	C	56	GLY	2.7
1	D	78	ASP	2.7
1	B	92	TRP	2.7
1	D	422	VAL	2.6
1	C	421	MET	2.6
1	D	214	GLY	2.5
1	C	447	HIS	2.5
1	C	419	ARG	2.5
1	B	333	VAL	2.5
1	B	93	LYS	2.4
1	D	34	GLY	2.4
1	D	423	GLN	2.4
1	C	164	VAL	2.3
1	A	11	ALA	2.3
1	D	221	PHE	2.3
1	D	233	ALA	2.3
1	B	5	LYS	2.2
1	D	164	VAL	2.2
1	D	346	ALA	2.2
1	D	98	THR	2.2
1	D	330	ASP	2.2
1	A	97	GLU	2.2
1	A	164	VAL	2.2
1	C	446	THR	2.1
1	D	93	LYS	2.1
1	A	7	ASP	2.1
1	D	418	VAL	2.1
1	D	47	ASN	2.1
1	C	358	CYS	2.1
1	C	238	LEU	2.1
1	B	10	TRP	2.0
1	C	324	LEU	2.0
1	D	22	LEU	2.0

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 [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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	MG	B	501	1/1	0.77	0.15	-0.39	65,65,65,65	0
2	MG	A	501	1/1	0.92	0.12	-2.02	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.