



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:23 PM UTC

PDB ID : 2PAJ / pdb_00002paj
Title : Crystal structure of an amidohydrolase from an environmental sample of Sargasso sea
Authors : Agarwal, R.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2007-03-27
Resolution : 2.70 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

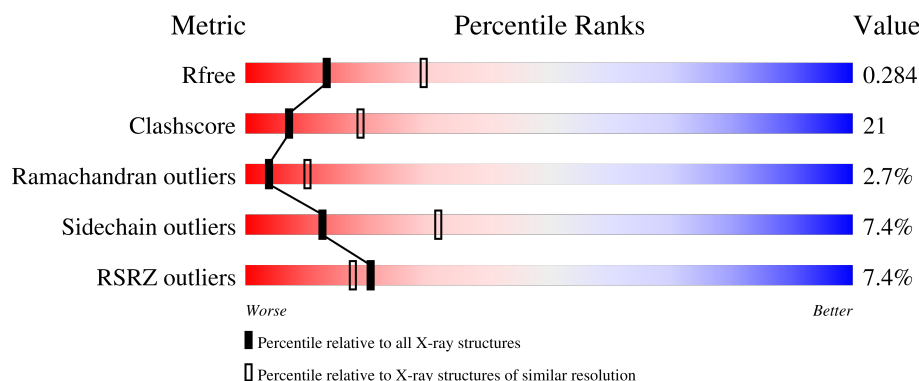
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 $\leq 5\%$. 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	492	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3280 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 putative cytosine/guanine deaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	421	Total	C	N	O	S	0	0	0
			3194	2006	584	589	15			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

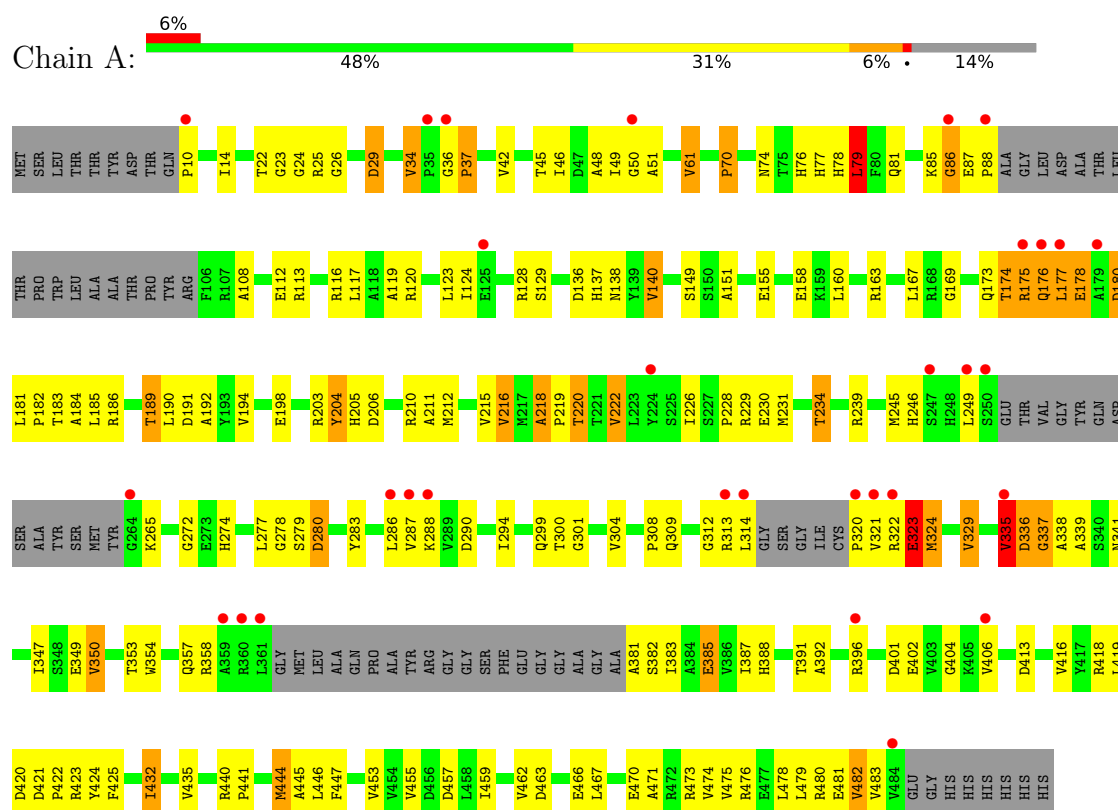
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	85	Total	O	0	0
			85	85		

3 Residue-property plots

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: putative cytosine/guanine deaminase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.87Å 88.87Å 161.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.44 – 2.70 44.44 – 2.70	Depositor EDS
% Data completeness (in resolution range)	91.6 (44.44-2.70) 91.6 (44.44-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.33 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.286 0.232 , 0.284	Depositor DCC
R_{free} test set	569 reflections (2.57%)	wwPDB-VP
Wilson B-factor (Å ²)	48.6	Xtriage
Anisotropy	0.539	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3280	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

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

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.54	0/3256	1.13	21/4419 (0.5%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	176	GLN	N-CA-C	-18.58	88.41	113.30
1	A	218	ALA	N-CA-C	8.20	121.52	109.30
1	A	79	LEU	N-CA-C	6.61	120.21	111.75
1	A	129	SER	N-CA-C	-6.58	104.40	112.88
1	A	184	ALA	N-CA-C	6.38	119.91	111.75
1	A	336	ASP	N-CA-C	6.23	119.09	109.07
1	A	401	ASP	N-CA-C	6.18	120.43	113.01
1	A	70	PRO	N-CA-C	-6.12	101.45	111.68
1	A	323	GLU	N-CA-C	-6.08	97.85	110.80
1	A	444	MET	N-CA-C	-5.94	104.93	111.82
1	A	229	ARG	N-CA-C	-5.92	104.91	111.36
1	A	335	VAL	N-CA-C	-5.84	99.07	107.78
1	A	204	TYR	N-CA-C	5.79	121.01	113.88
1	A	34	VAL	CA-C-N	-5.77	112.62	119.84
1	A	34	VAL	C-N-CA	-5.77	112.62	119.84
1	A	36	GLY	N-CA-C	-5.62	100.87	112.34
1	A	324	MET	N-CA-C	-5.50	105.20	111.14
1	A	481	GLU	N-CA-C	5.39	117.85	111.33
1	A	180	ASP	N-CA-C	-5.12	106.78	112.57
1	A	335	VAL	CA-C-N	-5.02	116.07	123.00
1	A	335	VAL	C-N-CA	-5.02	116.07	123.00

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	3194	0	3187	133	0
2	A	1	0	0	0	0
3	A	85	0	0	9	0
All	All	3280	0	3187	133	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 21.

All (133) 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:A:70:PRO:HG2	1:A:391:THR:HG21	1.52	0.91
1:A:347:ILE:HD13	1:A:441:PRO:HD2	1.65	0.77
1:A:335:VAL:HG12	1:A:349:GLU:HG3	1.67	0.76
1:A:228:PRO:HA	1:A:231:MET:HG2	1.68	0.75
1:A:176:GLN:HB2	1:A:181:LEU:HD11	1.67	0.74
1:A:137:HIS:HA	1:A:167:LEU:HB2	1.70	0.72
1:A:14:ILE:HA	1:A:61:VAL:HG13	1.73	0.70
1:A:181:LEU:HB3	3:A:525:HOH:O	1.90	0.69
1:A:383:ILE:HG12	1:A:432:ILE:HG12	1.75	0.69
1:A:37:PRO:HD2	1:A:50:GLY:HA2	1.75	0.67
1:A:124:ILE:O	1:A:128:ARG:HG2	1.94	0.66
1:A:413:ASP:HA	1:A:447:PHE:O	1.95	0.66
1:A:480:ARG:O	1:A:483:VAL:HG22	1.96	0.65
1:A:230:GLU:O	1:A:234:THR:HB	1.97	0.64
1:A:239:ARG:CZ	1:A:245:MET:HE1	2.28	0.63
1:A:76:HIS:CD2	1:A:336:ASP:HB3	2.33	0.63
1:A:354:TRP:CZ2	1:A:358:ARG:HD3	2.33	0.63
1:A:322:ARG:O	1:A:323:GLU:HG3	1.99	0.63
1:A:459:ILE:O	1:A:462:VAL:HG22	2.00	0.62
1:A:203:ARG:HD3	1:A:204:TYR:CE1	2.34	0.62
1:A:335:VAL:O	1:A:335:VAL:HG22	1.99	0.61
1:A:29:ASP:OD2	1:A:29:ASP:N	2.33	0.61
1:A:186:ARG:HG3	3:A:525:HOH:O	2.00	0.61
1:A:191:ASP:HB2	3:A:533:HOH:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:MET:O	1:A:329:VAL:HG13	1.99	0.61
1:A:205:HIS:HD2	1:A:215:VAL:H	1.47	0.60
1:A:335:VAL:CG1	1:A:349:GLU:HG3	2.31	0.60
1:A:37:PRO:HD2	1:A:51:ALA:H	1.67	0.60
1:A:120:ARG:O	1:A:124:ILE:HG12	2.01	0.60
1:A:34:VAL:HG13	1:A:34:VAL:O	2.01	0.59
1:A:140:VAL:O	1:A:140:VAL:CG2	2.51	0.59
1:A:312:GLY:C	1:A:314:LEU:H	2.11	0.59
1:A:278:GLY:C	1:A:280:ASP:H	2.11	0.59
1:A:119:ALA:O	1:A:123:LEU:HB2	2.04	0.58
1:A:322:ARG:HH21	1:A:385:GLU:CD	2.12	0.58
1:A:25:ARG:HE	1:A:396:ARG:NH2	2.02	0.58
1:A:163:ARG:HG2	1:A:212:MET:SD	2.44	0.58
1:A:182:PRO:HG2	1:A:185:LEU:HD23	1.85	0.58
1:A:421:ASP:OD2	1:A:423:ARG:HG2	2.03	0.58
1:A:169:GLY:HA2	1:A:218:ALA:O	2.04	0.57
1:A:335:VAL:O	1:A:335:VAL:CG2	2.52	0.57
1:A:116:ARG:HH22	1:A:155:GLU:CD	2.12	0.57
1:A:231:MET:O	1:A:234:THR:HG22	2.04	0.57
1:A:124:ILE:HD12	1:A:462:VAL:HG11	1.86	0.56
1:A:287:VAL:HG13	1:A:288:LYS:HG3	1.87	0.56
1:A:353:THR:O	1:A:357:GLN:HG2	2.05	0.56
1:A:81:GLN:OE1	1:A:337:GLY:HA3	2.06	0.56
1:A:70:PRO:HB3	1:A:406:VAL:HG23	1.88	0.56
1:A:286:LEU:O	1:A:286:LEU:HG	2.06	0.55
1:A:79:LEU:HD23	1:A:136:ASP:HB2	1.89	0.54
1:A:350:VAL:HG11	1:A:435:VAL:HG22	1.89	0.54
1:A:308:PRO:HD2	1:A:349:GLU:OE1	2.08	0.54
1:A:463:ASP:CG	1:A:466:GLU:HB2	2.31	0.54
1:A:108:ALA:HB2	1:A:185:LEU:HD21	1.88	0.54
1:A:418:ARG:HG2	1:A:420:ASP:OD2	2.08	0.53
1:A:190:LEU:O	1:A:194:VAL:HG23	2.09	0.52
1:A:87:GLU:CD	1:A:88:PRO:HD3	2.34	0.52
1:A:357:GLN:HB3	1:A:381:ALA:HB1	1.92	0.51
1:A:108:ALA:HB2	1:A:185:LEU:CD2	2.40	0.51
1:A:206:ASP:HB3	1:A:211:ALA:HB1	1.91	0.51
1:A:337:GLY:HA2	3:A:502:HOH:O	2.10	0.51
1:A:87:GLU:HB3	1:A:88:PRO:CD	2.41	0.50
1:A:203:ARG:HH11	1:A:203:ARG:HG2	1.76	0.50
1:A:286:LEU:HD21	1:A:304:VAL:HB	1.93	0.50
1:A:25:ARG:HH21	1:A:396:ARG:NH2	2.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:PRO:HD2	1:A:51:ALA:N	2.26	0.49
1:A:87:GLU:HA	1:A:87:GLU:OE1	2.12	0.49
1:A:336:ASP:O	1:A:337:GLY:O	2.30	0.49
1:A:478:LEU:O	1:A:482:VAL:HG22	2.13	0.49
1:A:210:ARG:HD2	1:A:402:GLU:OE1	2.12	0.48
1:A:470:GLU:O	1:A:474:VAL:HG23	2.13	0.48
1:A:419:LEU:HB3	1:A:424:TYR:CG	2.48	0.48
1:A:194:VAL:O	1:A:198:GLU:HG3	2.13	0.48
1:A:278:GLY:C	1:A:280:ASP:N	2.69	0.48
1:A:392:ALA:HB2	1:A:404:GLY:O	2.14	0.47
1:A:177:LEU:HB3	1:A:180:ASP:OD1	2.15	0.47
1:A:279:SER:HA	1:A:301:GLY:O	2.15	0.47
1:A:338:ALA:HA	1:A:341:ASN:O	2.15	0.47
1:A:382:SER:OG	1:A:385:GLU:HB2	2.15	0.46
1:A:88:PRO:HD3	1:A:313:ARG:HH12	1.81	0.46
1:A:416:VAL:HB	1:A:445:ALA:HB3	1.98	0.46
1:A:77:HIS:HA	1:A:336:ASP:HA	1.96	0.46
1:A:42:VAL:O	1:A:42:VAL:HG13	2.16	0.46
1:A:440:ARG:NH2	1:A:457:ASP:OD1	2.49	0.46
1:A:48:ALA:C	1:A:49:ILE:HG13	2.42	0.45
1:A:174:THR:HG23	1:A:175:ARG:N	2.31	0.45
1:A:462:VAL:O	1:A:462:VAL:HG23	2.16	0.45
1:A:113:ARG:NH2	1:A:117:LEU:HD13	2.31	0.45
1:A:140:VAL:O	1:A:140:VAL:HG23	2.16	0.45
1:A:312:GLY:C	1:A:314:LEU:N	2.74	0.45
1:A:151:ALA:O	1:A:155:GLU:HG3	2.17	0.45
1:A:455:VAL:O	1:A:455:VAL:HG13	2.17	0.45
1:A:70:PRO:HG2	1:A:391:THR:CG2	2.36	0.45
1:A:25:ARG:HH21	1:A:396:ARG:HH21	1.64	0.44
1:A:422:PRO:HA	1:A:425:PHE:CD1	2.52	0.44
1:A:10:PRO:HB2	3:A:539:HOH:O	2.17	0.44
1:A:167:LEU:HD23	1:A:216:VAL:HG13	2.00	0.44
1:A:76:HIS:HD1	1:A:246:HIS:HE1	1.66	0.44
1:A:312:GLY:HA3	3:A:575:HOH:O	2.18	0.44
1:A:350:VAL:CG1	1:A:435:VAL:HG22	2.48	0.44
1:A:228:PRO:HA	1:A:231:MET:CG	2.44	0.43
1:A:473:ARG:HG3	3:A:573:HOH:O	2.18	0.43
1:A:383:ILE:O	1:A:387:ILE:HG13	2.18	0.43
1:A:24:GLY:C	1:A:26:GLY:H	2.27	0.43
1:A:473:ARG:HA	1:A:476:ARG:HD2	1.99	0.43
1:A:158:GLU:O	1:A:158:GLU:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:471:ALA:O	1:A:475:VAL:HG23	2.19	0.42
1:A:45:THR:HG22	1:A:46:ILE:N	2.35	0.42
1:A:138:ASN:HD21	1:A:149:SER:HB3	1.85	0.42
1:A:190:LEU:C	1:A:190:LEU:HD13	2.44	0.42
1:A:85:LYS:O	1:A:86:GLY:C	2.61	0.42
1:A:160:LEU:O	1:A:453:VAL:HG13	2.20	0.42
1:A:290:ASP:O	1:A:294:ILE:HG13	2.20	0.42
1:A:473:ARG:HD3	3:A:558:HOH:O	2.20	0.42
1:A:77:HIS:ND1	1:A:78:HIS:N	2.63	0.41
1:A:219:PRO:O	1:A:220:THR:C	2.63	0.41
1:A:309:GLN:HB2	1:A:349:GLU:OE2	2.20	0.41
1:A:87:GLU:OE1	1:A:313:ARG:NH1	2.53	0.41
1:A:249:LEU:HB2	1:A:283:TYR:CD2	2.56	0.41
1:A:479:LEU:O	1:A:482:VAL:HG23	2.20	0.41
1:A:23:GLY:O	1:A:388:HIS:HD2	2.04	0.41
1:A:299:GLN:HG3	1:A:300:THR:N	2.35	0.41
1:A:272:GLY:HA2	1:A:277:LEU:HG	2.02	0.41
1:A:189:THR:HG23	1:A:192:ALA:HB2	2.02	0.41
1:A:117:LEU:HD22	1:A:474:VAL:HG21	2.03	0.40
1:A:182:PRO:O	1:A:183:THR:C	2.65	0.40
1:A:231:MET:HE2	1:A:274:HIS:HB2	2.03	0.40
1:A:74:ASN:HD21	1:A:77:HIS:HB2	1.87	0.40
1:A:249:LEU:C	1:A:249:LEU:HD23	2.46	0.40
1:A:354:TRP:CE2	1:A:358:ARG:HD3	2.56	0.40
1:A:22:THR:HG23	1:A:406:VAL:O	2.20	0.40
1:A:337:GLY:C	1:A:339:ALA:H	2.29	0.40
1:A:313:ARG:HD3	3:A:550:HOH:O	2.21	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	411/492 (84%)	368 (90%)	32 (8%)	11 (3%)	4	10

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	37	PRO
1	A	177	LEU
1	A	337	GLY
1	A	86	GLY
1	A	174	THR
1	A	178	GLU
1	A	220	THR
1	A	321	VAL
1	A	323	GLU
1	A	112	GLU
1	A	222	VAL

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	326/377 (86%)	302 (93%)	24 (7%)	13	32

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	61	VAL
1	A	79	LEU
1	A	140	VAL
1	A	173	GLN
1	A	175	ARG
1	A	178	GLU
1	A	189	THR
1	A	216	VAL
1	A	222	VAL

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Mol	Chain	Res	Type
1	A	226	ILE
1	A	234	THR
1	A	265	LYS
1	A	280	ASP
1	A	320	PRO
1	A	329	VAL
1	A	335	VAL
1	A	350	VAL
1	A	385	GLU
1	A	432	ILE
1	A	444	MET
1	A	446	LEU
1	A	467	LEU
1	A	482	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	138	ASN
1	A	173	GLN
1	A	205	HIS
1	A	246	HIS
1	A	274	HIS
1	A	309	GLN
1	A	341	ASN
1	A	351	HIS
1	A	388	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

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

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	421/492 (85%)	0.39	31 (7%) 20 18	42, 56, 77, 94	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	36	GLY	4.8
1	A	250	SER	4.7
1	A	288	LYS	4.7
1	A	484	VAL	4.7
1	A	361	LEU	4.2
1	A	35	PRO	3.9
1	A	249	LEU	3.7
1	A	88	PRO	3.7
1	A	322	ARG	3.5
1	A	179	ALA	3.4
1	A	359	ALA	3.4
1	A	314	LEU	3.2
1	A	86	GLY	3.1
1	A	320	PRO	3.1
1	A	224	TYR	3.0
1	A	287	VAL	3.0
1	A	264	GLY	2.9
1	A	286	LEU	2.9
1	A	313	ARG	2.8
1	A	50	GLY	2.5
1	A	335	VAL	2.4
1	A	406	VAL	2.4
1	A	360	ARG	2.4
1	A	175	ARG	2.3
1	A	396	ARG	2.3
1	A	177	LEU	2.3
1	A	10	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	176	GLN	2.3
1	A	125	GLU	2.2
1	A	247	SER	2.2
1	A	321	VAL	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	ZN	A	493	1/1	1.00	0.05	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.