



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:47 AM UTC

PDB ID : 3GWB / pdb_00003gwb
Title : Crystal structure of peptidase M16 inactive domain from *Pseudomonas fluorescens*. Northeast Structural Genomics target PIR293L
Authors : Seetharaman, J.; Chen, Y.; Fang, F.; Xiao, R.; Everett, J.K.; Acton, T.B.; Rost, B.; Montelione, G.T.; Tong, L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2009-03-31
Resolution : 1.90 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

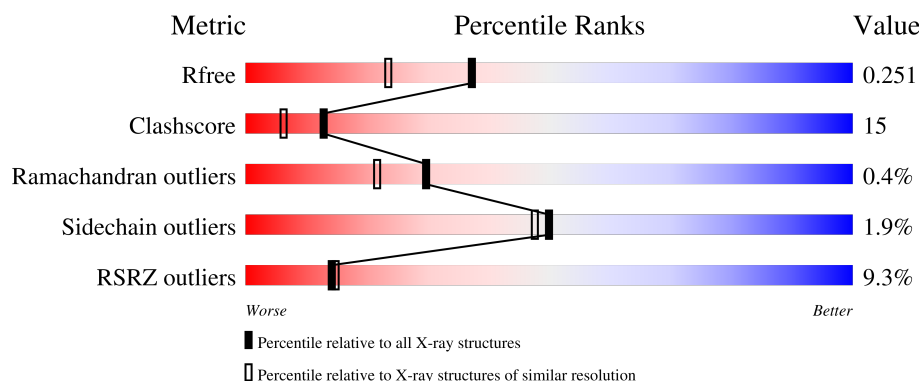
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	434	9% 68% 25% • 5%
1	B	434	8% 67% 24% • 5%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6636 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 Peptidase M16 inactive domain family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	412	Total	C	N	O	Se	0	0	0
			3074	1941	529	590	14			
1	B	412	Total	C	N	O	Se	0	0	0
			3078	1945	529	590	14			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	428	GLU	-	expression tag	UNP Q4K4B6
A	429	HIS	-	expression tag	UNP Q4K4B6
A	430	HIS	-	expression tag	UNP Q4K4B6
A	431	HIS	-	expression tag	UNP Q4K4B6
A	432	HIS	-	expression tag	UNP Q4K4B6
A	433	HIS	-	expression tag	UNP Q4K4B6
A	434	HIS	-	expression tag	UNP Q4K4B6
B	428	GLU	-	expression tag	UNP Q4K4B6
B	429	HIS	-	expression tag	UNP Q4K4B6
B	430	HIS	-	expression tag	UNP Q4K4B6
B	431	HIS	-	expression tag	UNP Q4K4B6
B	432	HIS	-	expression tag	UNP Q4K4B6
B	433	HIS	-	expression tag	UNP Q4K4B6
B	434	HIS	-	expression tag	UNP Q4K4B6

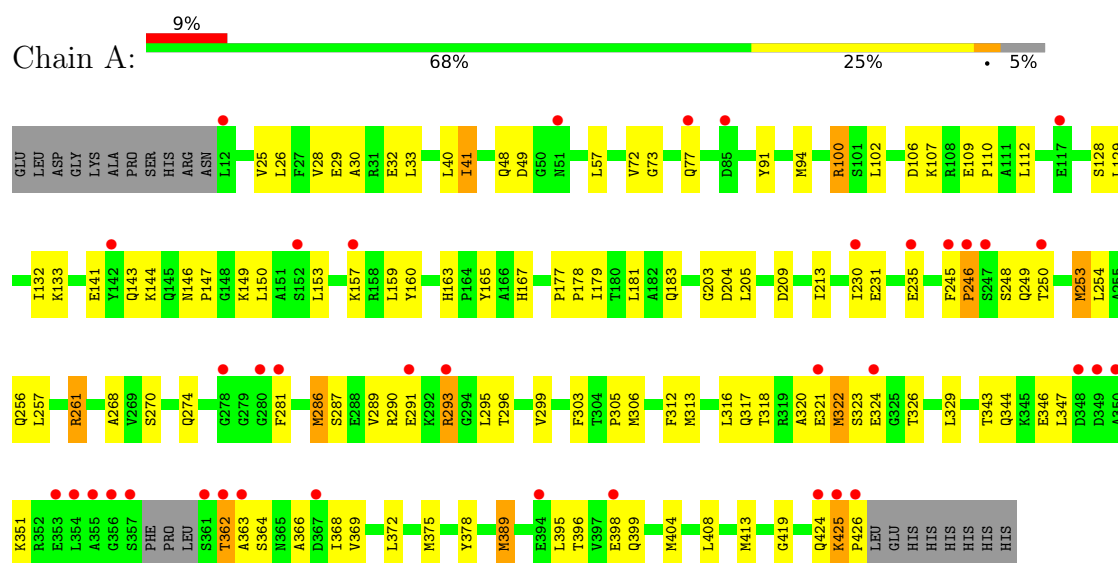
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	228	Total	O	0	0
			228	228		
2	B	256	Total	O	0	0
			256	256		

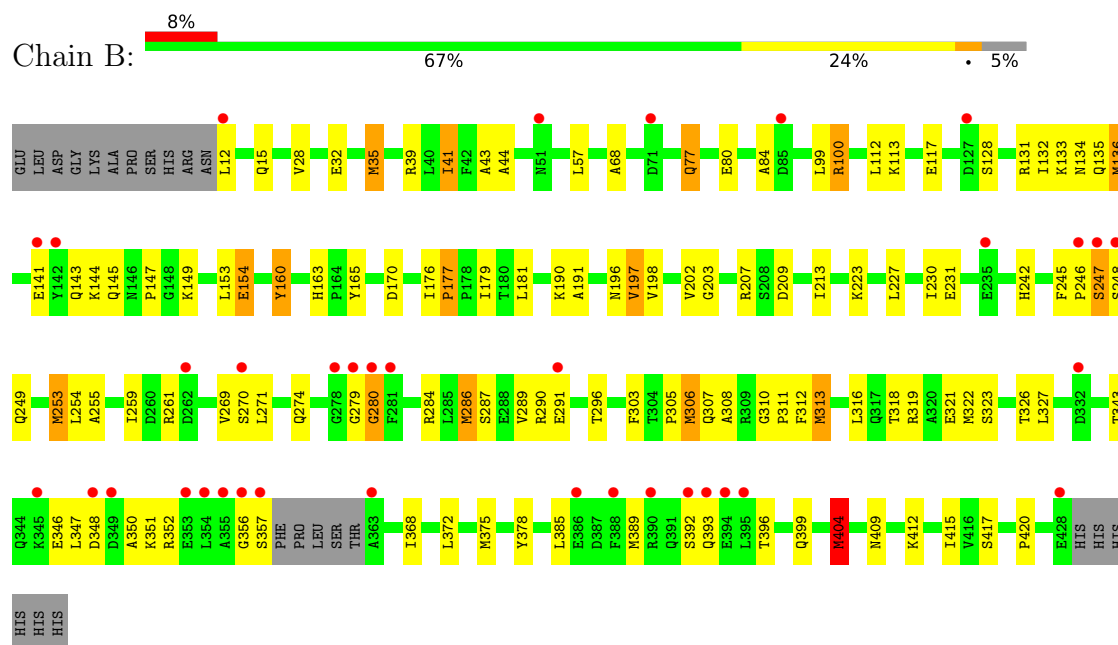
3 Residue-property plots [i](#)

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: Peptidase M16 inactive domain family protein



- Molecule 1: Peptidase M16 inactive domain family protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.67Å 64.84Å 114.97Å 90.00° 96.39° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90 50.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	77.5 (50.00-1.90) 90.8 (50.00-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 1.89Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.204 , 0.240 (Not available) , 0.251	Depositor DCC
R_{free} test set	5077 reflections (3.96%)	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6636	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% 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 [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	0.62	11/3113 (0.4%)	0.89	6/4184 (0.1%)
1	B	0.62	9/3117 (0.3%)	0.96	13/4189 (0.3%)
All	All	0.62	20/6230 (0.3%)	0.93	19/8373 (0.2%)

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	404	MSE	SE-CE	-9.08	1.68	1.95
1	B	306	MSE	SE-CE	-8.77	1.69	1.95
1	B	35	MSE	SE-CE	-7.20	1.73	1.95
1	A	404	MSE	SE-CE	-6.91	1.74	1.95
1	B	136	MSE	SE-CE	-6.79	1.75	1.95
1	A	306	MSE	SE-CE	-6.67	1.75	1.95
1	A	322	MSE	SE-CE	-6.54	1.75	1.95
1	A	253	MSE	SE-CE	-6.21	1.76	1.95
1	A	375	MSE	SE-CE	-6.00	1.77	1.95
1	A	286	MSE	SE-CE	-5.97	1.77	1.95
1	B	375	MSE	SE-CE	-5.62	1.78	1.95
1	A	94	MSE	SE-CE	-5.57	1.78	1.95
1	B	286	MSE	SE-CE	-5.52	1.78	1.95
1	A	389	MSE	CG-SE	-5.50	1.78	1.95
1	B	253	MSE	SE-CE	-5.41	1.79	1.95
1	A	253	MSE	CG-SE	-5.19	1.79	1.95
1	A	413	MSE	SE-CE	-5.18	1.79	1.95
1	B	313	MSE	SE-CE	-5.07	1.80	1.95
1	B	322	MSE	SE-CE	-5.02	1.80	1.95
1	A	286	MSE	CG-SE	-5.01	1.80	1.95

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	280	GLY	N-CA-C	13.65	145.54	113.18
1	B	279	GLY	N-CA-C	9.54	125.85	110.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	SER	N-CA-C	8.18	123.37	113.23
1	B	392	SER	N-CA-C	7.94	122.23	112.54
1	A	312	PHE	N-CA-C	-6.85	98.03	109.07
1	B	177	PRO	N-CA-C	6.50	118.63	110.70
1	A	203	GLY	N-CA-C	6.45	118.64	110.38
1	A	261	ARG	N-CA-C	6.25	118.89	111.33
1	B	312	PHE	N-CA-C	-6.06	99.32	109.07
1	B	323	SER	N-CA-C	5.75	117.23	111.07
1	A	102	LEU	N-CA-C	-5.69	102.61	110.35
1	B	43	ALA	N-CA-C	-5.64	101.16	109.62
1	B	100	ARG	N-CA-C	-5.49	99.95	108.90
1	A	100	ARG	N-CA-C	-5.36	100.44	109.07
1	B	248	SER	N-CA-C	-5.36	106.74	113.28
1	B	203	GLY	N-CA-C	5.17	117.72	110.38
1	B	197	VAL	N-CA-C	5.08	117.34	108.95
1	B	160	TYR	N-CA-C	5.05	119.59	113.38
1	B	308	ALA	N-CA-C	-5.03	103.02	110.46

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	3074	0	3091	91	0
1	B	3078	0	3096	100	0
2	A	228	0	0	10	0
2	B	256	0	0	7	0
All	All	6636	0	6187	186	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 15.

All (186) 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:B:280:GLY:O	1:B:286:MSE:HE3	1.64	0.96
1:A:141:GLU:HA	1:A:144:LYS:HE2	1.50	0.90
1:B:404:MSE:HE2	1:B:404:MSE:HA	1.51	0.90
1:A:324:GLU:OE2	1:A:424:GLN:HG3	1.74	0.88
1:A:246:PRO:HG2	1:B:246:PRO:HB3	1.55	0.88
1:B:230:ILE:H	1:B:307:GLN:NE2	1.74	0.84
1:A:143:GLN:HG2	1:A:149:LYS:HD2	1.62	0.81
1:B:190:LYS:HG2	1:B:227:LEU:HD21	1.65	0.78
1:A:296:THR:HG21	1:A:299:VAL:CG2	2.13	0.77
1:B:35:MSE:SE	1:B:100:ARG:HH21	2.18	0.77
1:B:112:LEU:HD13	1:B:213:ILE:HD12	1.67	0.76
1:B:141:GLU:O	1:B:144:LYS:HG2	1.87	0.75
1:A:112:LEU:HD13	1:A:213:ILE:HD12	1.68	0.75
1:A:296:THR:HG21	1:A:299:VAL:HG23	1.68	0.74
1:B:269:VAL:HG12	1:B:404:MSE:HE3	1.72	0.72
1:B:15:GLN:NE2	1:B:207:ARG:HG3	2.05	0.71
1:B:113:LYS:O	1:B:117:GLU:HG3	1.92	0.70
1:B:270:SER:O	1:B:274:GLN:HG3	1.94	0.68
1:B:149:LYS:O	1:B:153:LEU:HD13	1.94	0.68
1:A:321:GLU:O	1:A:322:MSE:HE2	1.94	0.67
1:B:41:ILE:HD12	1:B:41:ILE:N	2.10	0.66
1:B:147:PRO:HG3	1:B:245:PHE:CD2	2.30	0.66
1:A:318:THR:HG21	1:A:326:THR:HG21	1.76	0.66
1:B:348:ASP:O	1:B:352:ARG:HG3	1.95	0.66
1:A:177:PRO:HB2	1:A:178:PRO:HD3	1.79	0.65
1:A:425:LYS:N	1:A:426:PRO:HD3	2.12	0.65
1:A:249:GLN:HE21	1:A:317:GLN:HG2	1.62	0.64
1:A:41:ILE:N	1:A:41:ILE:HD12	2.13	0.64
1:B:128:SER:O	1:B:132:ILE:HG12	1.97	0.64
1:B:254:LEU:HD22	1:B:415:ILE:HG12	1.79	0.63
1:A:396:THR:OG1	1:A:398:GLU:HG2	1.99	0.63
1:B:249:GLN:NE2	1:B:319:ARG:HD3	2.15	0.62
1:A:287:SER:O	1:A:291:GLU:HB3	1.99	0.62
1:B:163:HIS:HE1	1:B:231:GLU:O	1.82	0.62
1:B:39:ARG:NH1	2:B:436:HOH:O	2.32	0.61
1:A:246:PRO:HG2	1:B:246:PRO:CB	2.29	0.59
1:B:404:MSE:HA	1:B:404:MSE:CE	2.27	0.59
1:A:112:LEU:CD1	1:A:213:ILE:HD12	2.33	0.59
1:B:261:ARG:HG2	1:B:378:TYR:CD2	2.39	0.58
1:B:15:GLN:HE22	1:B:207:ARG:HG3	1.67	0.57
1:B:247:SER:O	1:B:420:PRO:HG3	2.04	0.57
1:A:28:VAL:HG23	1:A:372:LEU:HD11	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:ILE:H	1:B:307:GLN:HE21	1.51	0.57
1:A:250:THR:OG1	1:A:323:SER:HB3	2.04	0.57
1:B:287:SER:HA	1:B:291:GLU:HB3	1.87	0.57
1:A:324:GLU:CD	1:A:424:GLN:HG3	2.29	0.56
1:A:293:ARG:NH2	1:A:329:LEU:HD11	2.20	0.56
1:A:347:LEU:HG	1:A:351:LYS:HD2	1.86	0.56
1:A:303:PHE:C	1:A:305:PRO:HD3	2.30	0.56
1:A:425:LYS:N	1:A:426:PRO:CD	2.68	0.56
1:B:230:ILE:H	1:B:307:GLN:HE22	1.50	0.56
1:A:235:GLU:OE1	1:A:257:LEU:HD12	2.06	0.55
1:A:250:THR:HG23	1:A:320:ALA:HA	1.88	0.55
1:A:163:HIS:HE1	1:A:231:GLU:O	1.89	0.54
1:A:289:VAL:HG12	1:A:296:THR:HG22	1.89	0.54
1:B:271:LEU:HD12	1:B:274:GLN:OE1	2.07	0.54
1:A:73:GLY:O	1:A:77:GLN:HG3	2.08	0.54
1:A:362:THR:OG1	1:B:77:GLN:NE2	2.41	0.54
1:A:49:ASP:OD2	1:A:167:HIS:HE1	1.91	0.54
1:A:163:HIS:CD2	1:A:165:TYR:H	2.26	0.53
1:B:181:LEU:H	1:B:181:LEU:HD12	1.74	0.53
1:A:48:GLN:NE2	1:A:230:ILE:HD11	2.24	0.53
1:A:100:ARG:HH11	1:A:100:ARG:HB2	1.73	0.53
1:A:146:ASN:CG	1:A:149:LYS:HG3	2.33	0.53
1:A:209:ASP:O	1:A:213:ILE:HG12	2.08	0.53
1:B:409:ASN:HB3	1:B:412:LYS:CG	2.39	0.53
1:A:424:GLN:HG2	2:A:662:HOH:O	2.09	0.53
1:B:112:LEU:CD1	1:B:213:ILE:HD12	2.39	0.53
1:B:253:MSE:C	1:B:254:LEU:HD23	2.34	0.53
1:A:295:LEU:O	1:A:318:THR:HG22	2.07	0.53
1:A:425:LYS:O	1:A:425:LYS:HG2	2.09	0.53
1:B:153:LEU:HD12	2:B:655:HOH:O	2.08	0.52
1:B:57:LEU:CD2	1:B:136:MSE:HE2	2.40	0.51
1:A:153:LEU:O	1:A:157:LYS:HG3	2.09	0.51
1:B:318:THR:HG21	1:B:326:THR:HG21	1.92	0.51
1:B:176:ILE:HB	1:B:177:PRO:HD3	1.92	0.51
1:B:196:ASN:ND2	1:B:227:LEU:H	2.09	0.51
1:B:163:HIS:CD2	1:B:165:TYR:H	2.27	0.51
1:B:255:ALA:HB2	1:B:313:MSE:HG2	1.92	0.51
1:A:321:GLU:CD	1:A:425:LYS:HE3	2.36	0.51
1:B:246:PRO:O	1:B:247:SER:HB2	2.11	0.51
1:A:29:GLU:HG3	1:A:205:LEU:O	2.10	0.50
1:A:128:SER:O	1:A:132:ILE:HG12	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:TYR:HB3	1:A:163:HIS:HB3	1.93	0.50
1:B:249:GLN:HE22	1:B:319:ARG:HD3	1.76	0.50
1:B:132:ILE:HG22	1:B:136:MSE:HE3	1.92	0.49
1:A:163:HIS:HD2	1:A:165:TYR:H	1.60	0.49
1:B:209:ASP:O	1:B:213:ILE:HG12	2.12	0.49
1:B:396:THR:HG23	1:B:399:GLN:HE21	1.78	0.49
1:A:30:ALA:HB1	1:A:32:GLU:OE1	2.12	0.49
1:B:409:ASN:HB3	1:B:412:LYS:HG3	1.94	0.49
1:A:167:HIS:HD2	2:A:534:HOH:O	1.95	0.49
1:A:100:ARG:HB2	1:A:100:ARG:NH1	2.29	0.48
1:B:80:GLU:HG3	2:B:634:HOH:O	2.13	0.48
1:B:385:LEU:O	1:B:389:MSE:HG2	2.13	0.48
1:B:28:VAL:HG23	1:B:372:LEU:HD11	1.96	0.48
1:B:254:LEU:HD21	1:B:327:LEU:CD1	2.43	0.48
1:B:44:ALA:HB2	1:B:191:ALA:HB1	1.95	0.48
1:B:284:ARG:HD3	1:B:346:GLU:OE2	2.14	0.48
1:A:343:THR:OG1	1:A:346:GLU:HG3	2.14	0.47
1:A:261:ARG:HG2	1:A:378:TYR:CD2	2.49	0.47
1:B:181:LEU:HD12	1:B:181:LEU:N	2.30	0.47
1:A:363:ALA:HB3	1:A:368:ILE:HD11	1.96	0.47
1:B:154:GLU:HA	1:B:154:GLU:OE1	2.15	0.47
1:B:197:VAL:HG22	1:B:198:VAL:N	2.29	0.47
1:B:254:LEU:CD2	1:B:415:ILE:HG23	2.44	0.47
1:A:408:LEU:HD22	2:A:660:HOH:O	2.14	0.47
1:B:286:MSE:O	1:B:290:ARG:HB3	2.14	0.47
1:A:290:ARG:HE	1:B:134:ASN:ND2	2.13	0.47
1:B:351:LYS:HB3	1:B:393:GLN:OE1	2.14	0.47
1:A:147:PRO:HG3	1:A:245:PHE:CD2	2.50	0.46
1:A:40:LEU:C	1:A:41:ILE:HD12	2.39	0.46
1:A:129:LEU:O	1:A:133:LYS:HG3	2.15	0.46
1:A:270:SER:O	1:A:274:GLN:HG3	2.15	0.46
1:B:32:GLU:H	1:B:32:GLU:CD	2.24	0.46
1:B:133:LYS:NZ	1:B:179:ILE:HB	2.30	0.46
1:B:223:LYS:HD2	2:B:522:HOH:O	2.16	0.46
1:B:289:VAL:HG12	1:B:296:THR:HG22	1.97	0.46
1:A:322:MSE:CE	1:A:425:LYS:HB3	2.46	0.46
1:B:163:HIS:HD2	1:B:165:TYR:H	1.64	0.46
1:B:242:HIS:HD2	1:B:417:SER:OG	1.99	0.46
1:A:109:GLU:HB2	1:A:110:PRO:HD3	1.97	0.46
1:B:343:THR:OG1	1:B:346:GLU:HG3	2.16	0.46
1:B:254:LEU:HD21	1:B:327:LEU:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:GLN:NE2	1:B:145:GLN:HE22	2.15	0.45
1:B:259:ILE:HD13	1:B:269:VAL:HG21	1.99	0.45
1:A:396:THR:OG1	1:A:399:GLN:HG3	2.16	0.45
1:A:268:ALA:HA	1:A:395:LEU:HD13	1.99	0.45
1:A:129:LEU:HD22	1:A:181:LEU:HD23	1.99	0.45
1:B:41:ILE:N	1:B:41:ILE:CD1	2.80	0.45
1:B:280:GLY:O	1:B:286:MSE:CE	2.52	0.45
1:B:68:ALA:HA	2:B:551:HOH:O	2.17	0.45
1:B:147:PRO:HG3	1:B:245:PHE:CE2	2.52	0.45
1:A:321:GLU:OE2	1:A:425:LYS:HE3	2.16	0.44
1:B:132:ILE:CG2	1:B:136:MSE:HE3	2.47	0.44
1:B:287:SER:O	1:B:291:GLU:HB3	2.17	0.44
1:B:303:PHE:C	1:B:305:PRO:HD3	2.43	0.44
1:A:41:ILE:N	1:A:41:ILE:CD1	2.81	0.44
1:A:254:LEU:HD12	1:A:254:LEU:N	2.32	0.44
1:A:256:GLN:NE2	2:A:660:HOH:O	2.50	0.44
1:B:213:ILE:HD11	2:B:495:HOH:O	2.17	0.44
1:A:253:MSE:HE2	1:A:313:MSE:SE	2.68	0.44
1:B:356:GLY:O	1:B:357:SER:HB3	2.17	0.44
1:A:91:TYR:HB2	2:A:530:HOH:O	2.17	0.44
1:A:153:LEU:HD23	2:A:533:HOH:O	2.17	0.43
1:A:250:THR:HG22	1:A:419:GLY:HA3	1.99	0.43
1:B:269:VAL:HG12	1:B:404:MSE:CE	2.46	0.43
1:B:149:LYS:HE3	1:B:153:LEU:HD11	2.00	0.43
1:A:250:THR:CG2	1:A:320:ALA:HA	2.49	0.43
1:B:356:GLY:O	1:B:357:SER:CB	2.66	0.43
1:B:352:ARG:HH11	1:B:352:ARG:HG2	1.84	0.43
1:B:160:TYR:HB3	1:B:163:HIS:HB3	2.02	0.42
1:B:321:GLU:HG2	2:B:534:HOH:O	2.18	0.42
1:A:144:LYS:HA	1:A:150:LEU:HD21	2.00	0.42
1:A:179:ILE:HA	1:A:183:GLN:OE1	2.20	0.42
1:B:15:GLN:HE21	1:B:207:ARG:HH11	1.67	0.42
1:B:57:LEU:HD13	1:B:170:ASP:HA	2.01	0.42
1:A:72:VAL:HG13	1:A:73:GLY:N	2.33	0.42
1:A:321:GLU:OE1	1:A:425:LYS:HE3	2.20	0.42
1:A:426:PRO:HG2	2:A:448:HOH:O	2.20	0.42
1:B:35:MSE:HE3	1:B:35:MSE:HB2	1.90	0.42
1:B:143:GLN:C	1:B:145:GLN:H	2.28	0.42
1:A:100:ARG:HD3	2:A:592:HOH:O	2.18	0.42
1:A:177:PRO:CB	1:A:178:PRO:HD3	2.44	0.42
1:A:299:VAL:HG22	1:A:316:LEU:CB	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:245:PHE:CD1	1:B:246:PRO:HD2	2.54	0.42
1:A:296:THR:CG2	1:A:299:VAL:HG23	2.46	0.41
1:B:133:LYS:HZ2	1:B:179:ILE:HB	1.85	0.41
1:A:366:ALA:O	1:A:369:VAL:HG12	2.20	0.41
1:B:306:MSE:HE3	1:B:310:GLY:HA2	2.02	0.41
1:A:106:ASP:OD2	1:A:107:LYS:HE3	2.21	0.41
1:A:25:VAL:C	1:A:26:LEU:HD12	2.45	0.41
1:B:84:ALA:HB1	1:B:99:LEU:HD11	2.02	0.41
1:A:268:ALA:HA	1:A:395:LEU:CD1	2.51	0.41
1:A:363:ALA:CB	1:A:368:ILE:HD11	2.50	0.41
1:B:347:LEU:O	1:B:350:ALA:HB3	2.20	0.41
1:B:131:ARG:O	1:B:135:GLN:HG3	2.21	0.41
1:A:33:LEU:O	1:A:204:ASP:HB2	2.21	0.40
1:B:132:ILE:O	1:B:136:MSE:HG3	2.21	0.40
1:B:202:VAL:HG21	1:B:368:ILE:HG22	2.02	0.40
1:A:344:GLN:NE2	2:A:446:HOH:O	2.40	0.40
1:B:306:MSE:HE2	1:B:311:PRO:HD2	2.03	0.40
1:A:281:PHE:HD1	1:A:286:MSE:HE3	1.86	0.40
1:A:299:VAL:HG22	1:A:316:LEU:HB2	2.03	0.40
1:A:364:SER:HB2	2:A:498:HOH:O	2.21	0.40
1:B:254:LEU:HD22	1:B:415:ILE:HG23	2.03	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	408/434 (94%)	398 (98%)	8 (2%)	2 (0%)	24	16
1	B	408/434 (94%)	397 (97%)	10 (2%)	1 (0%)	43	36
All	All	816/868 (94%)	795 (97%)	18 (2%)	3 (0%)	30	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	247	SER
1	A	246	PRO
1	A	425	LYS

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	316/322 (98%)	310 (98%)	6 (2%)	50	47
1	B	316/322 (98%)	310 (98%)	6 (2%)	50	47
All	All	632/644 (98%)	620 (98%)	12 (2%)	50	47

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	ILE
1	A	57	LEU
1	A	159	LEU
1	A	293	ARG
1	A	362	THR
1	A	389	MSE
1	B	12	LEU
1	B	41	ILE
1	B	77	GLN
1	B	154	GLU
1	B	316	LEU
1	B	404	MSE

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

Mol	Chain	Res	Type
1	A	48	GLN
1	A	88	ASN

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Mol	Chain	Res	Type
1	A	163	HIS
1	A	167	HIS
1	A	249	GLN
1	A	273	ASN
1	A	274	GLN
1	A	317	GLN
1	A	331	GLN
1	A	344	GLN
1	A	391	GLN
1	A	405	ASN
1	B	13	ASN
1	B	15	GLN
1	B	77	GLN
1	B	145	GLN
1	B	163	HIS
1	B	183	GLN
1	B	196	ASN
1	B	242	HIS
1	B	249	GLN
1	B	273	ASN
1	B	307	GLN
1	B	317	GLN
1	B	371	GLN
1	B	399	GLN
1	B	405	ASN

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

There are no ligands in this entry.

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	398/434 (91%)	0.66	38 (9%) 14 14	7, 19, 35, 49	0
1	B	398/434 (91%)	0.66	36 (9%) 15 16	7, 18, 34, 46	0
All	All	796/868 (91%)	0.66	74 (9%) 14 15	7, 19, 34, 49	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	280	GLY	8.8
1	B	281	PHE	8.1
1	A	361	SER	6.1
1	A	362	THR	4.6
1	B	356	GLY	4.5
1	B	355	ALA	4.4
1	A	426	PRO	4.1
1	B	279	GLY	4.0
1	B	428	GLU	4.0
1	B	357	SER	4.0
1	B	12	LEU	4.0
1	A	425	LYS	3.9
1	A	246	PRO	3.9
1	B	390	ARG	3.6
1	A	247	SER	3.5
1	B	393	GLN	3.5
1	B	142	TYR	3.5
1	B	348	ASP	3.5
1	B	392	SER	3.4
1	A	356	GLY	3.3
1	A	142	TYR	3.3
1	B	262	ASP	3.2
1	B	278	GLY	3.2
1	B	349	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	354	LEU	3.1
1	B	395	LEU	3.0
1	B	246	PRO	3.0
1	A	354	LEU	3.0
1	A	394	GLU	3.0
1	A	424	GLN	3.0
1	A	357	SER	2.9
1	B	127	ASP	2.9
1	A	398	GLU	2.9
1	A	245	PHE	2.9
1	A	291	GLU	2.9
1	A	12	LEU	2.8
1	B	141	GLU	2.8
1	A	355	ALA	2.8
1	A	353	GLU	2.7
1	B	353	GLU	2.7
1	B	394	GLU	2.7
1	B	388	PHE	2.7
1	B	247	SER	2.6
1	B	332	ASP	2.6
1	A	51	ASN	2.6
1	A	367	ASP	2.5
1	A	293	ARG	2.5
1	B	235	GLU	2.5
1	A	281	PHE	2.4
1	A	230	ILE	2.4
1	A	250	THR	2.4
1	A	321	GLU	2.4
1	B	248	SER	2.4
1	B	51	ASN	2.4
1	A	348	ASP	2.3
1	A	85	ASP	2.3
1	A	235	GLU	2.3
1	A	363	ALA	2.3
1	A	157	LYS	2.3
1	B	71	ASP	2.3
1	A	117	GLU	2.3
1	B	345	LYS	2.2
1	A	77	GLN	2.2
1	A	350	ALA	2.2
1	A	280	GLY	2.2
1	B	270	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	363	ALA	2.1
1	A	349	ASP	2.1
1	A	324	GLU	2.1
1	B	85	ASP	2.1
1	A	152	SER	2.1
1	B	386	GLU	2.1
1	B	291	GLU	2.0
1	A	278	GLY	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.