Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 – 08:57 PM EDT

PDB ID	:	2OAM
Title	:	Apo RebH from Lechevalieria aerocolonigenes
Authors	:	Blasiak, L.C.; Drennan, C.L.
Deposited on	:	2006-12-16
Resolution	:	2.30 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

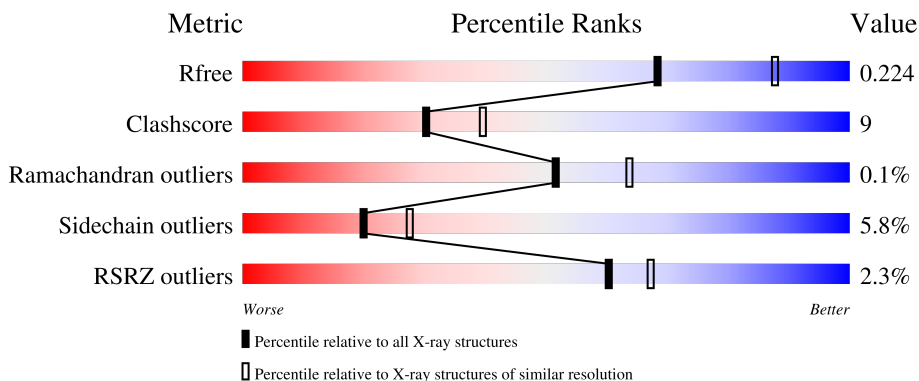|		
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.30 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	Percentile Ranks	Value
Rfree		0.224
Clashscore		9
Ramachandran outliers		0.1%
Sidechain outliers		5.8%
RSRZ outliers		2.3%

Worse ────────────────── *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	550	2% · 74% · 19% · 5%
1	B	550	2% · 75% · 19% · 5%

2 Entry composition ⓘ

There are 2 unique types of molecules in this entry. The entry contains 8973 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 Tryptophan halogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	523	Total 4200	C 2668	N 732	O 781	S 19	0	0	0
1	B	524	Total 4199	C 2667	N 731	O 782	S 19	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q8KHZ8
A	-18	GLY	-	cloning artifact	UNP Q8KHZ8
A	-17	SER	-	cloning artifact	UNP Q8KHZ8
A	-16	SER	-	cloning artifact	UNP Q8KHZ8
A	-15	HIS	-	expression tag	UNP Q8KHZ8
A	-14	HIS	-	expression tag	UNP Q8KHZ8
A	-13	HIS	-	expression tag	UNP Q8KHZ8
A	-12	HIS	-	expression tag	UNP Q8KHZ8
A	-11	HIS	-	expression tag	UNP Q8KHZ8
A	-10	HIS	-	expression tag	UNP Q8KHZ8
A	-9	SER	-	cloning artifact	UNP Q8KHZ8
A	-8	SER	-	cloning artifact	UNP Q8KHZ8
A	-7	GLY	-	cloning artifact	UNP Q8KHZ8
A	-6	LEU	-	cloning artifact	UNP Q8KHZ8
A	-5	VAL	-	cloning artifact	UNP Q8KHZ8
A	-4	PRO	-	cloning artifact	UNP Q8KHZ8
A	-3	ARG	-	cloning artifact	UNP Q8KHZ8
A	-2	GLY	-	cloning artifact	UNP Q8KHZ8
A	-1	SER	-	cloning artifact	UNP Q8KHZ8
A	0	HIS	-	cloning artifact	UNP Q8KHZ8
B	-19	MET	-	initiating methionine	UNP Q8KHZ8
B	-18	GLY	-	cloning artifact	UNP Q8KHZ8
B	-17	SER	-	cloning artifact	UNP Q8KHZ8
B	-16	SER	-	cloning artifact	UNP Q8KHZ8
B	-15	HIS	-	expression tag	UNP Q8KHZ8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q8KHZ8
B	-13	HIS	-	expression tag	UNP Q8KHZ8
B	-12	HIS	-	expression tag	UNP Q8KHZ8
B	-11	HIS	-	expression tag	UNP Q8KHZ8
B	-10	HIS	-	expression tag	UNP Q8KHZ8
B	-9	SER	-	cloning artifact	UNP Q8KHZ8
B	-8	SER	-	cloning artifact	UNP Q8KHZ8
B	-7	GLY	-	cloning artifact	UNP Q8KHZ8
B	-6	LEU	-	cloning artifact	UNP Q8KHZ8
B	-5	VAL	-	cloning artifact	UNP Q8KHZ8
B	-4	PRO	-	cloning artifact	UNP Q8KHZ8
B	-3	ARG	-	cloning artifact	UNP Q8KHZ8
B	-2	GLY	-	cloning artifact	UNP Q8KHZ8
B	-1	SER	-	cloning artifact	UNP Q8KHZ8
B	0	HIS	-	cloning artifact	UNP Q8KHZ8

- Molecule 2 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	289	Total 289	O 289	0	0
2	B	285	Total 285	O 285	0	0

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: Tryptophan halogenase
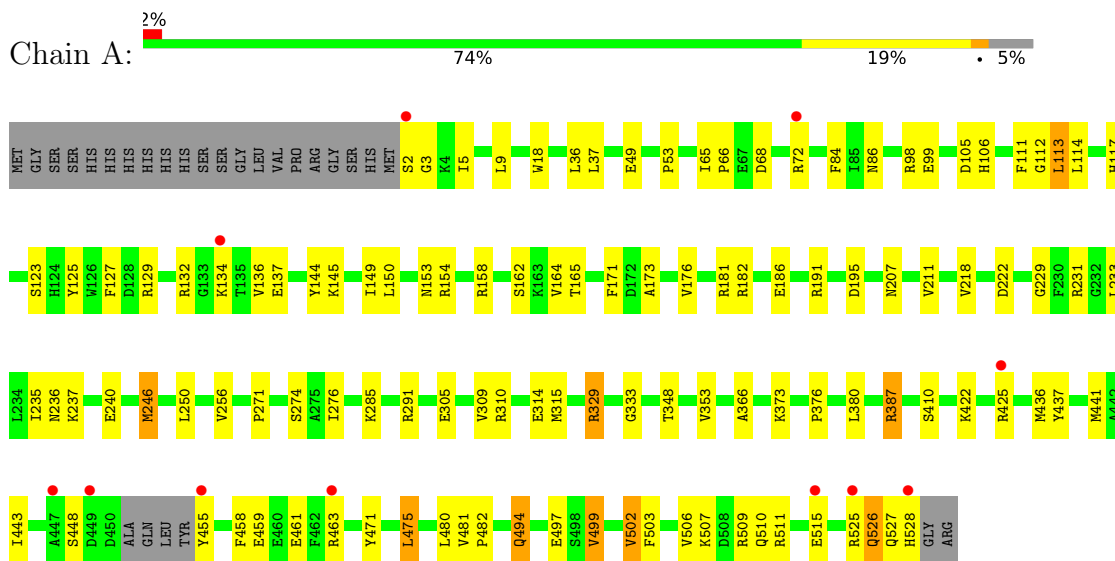
- Molecule 1: Tryptophan halogenase
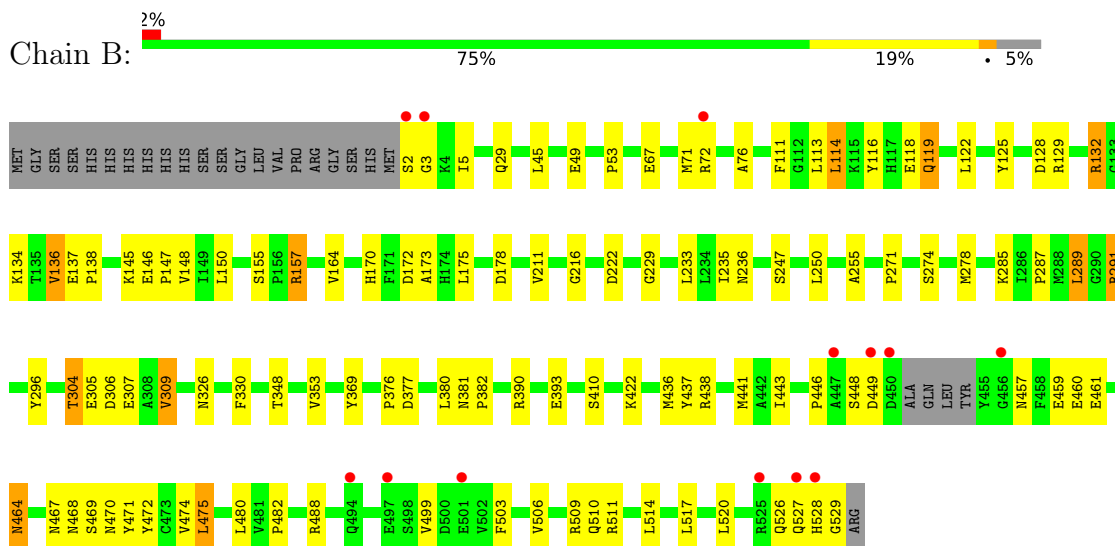
4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α, β, γ	114.82Å　114.82Å　230.63Å 90.00°　90.00°　120.00°	Depositor
Resolution (Å)	41.70　–　2.30 41.75　–　2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (41.70-2.30) 99.8 (41.75-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	2.55 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.199　,　0.224 0.198　,　0.224	Depositor DCC
R_{free} test set	3797 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	36.4	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 42.5	EDS
L-test for twinning[2]	$< |L| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.039 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	8973	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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 ⓘ

5.1 Standard geometry ⓘ

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	0.41	0/4315	0.61	1/5854 (0.0%)
1	B	0.37	0/4313	0.61	1/5851 (0.0%)
All	All	0.39	0/8628	0.61	2/11705 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	3	GLY	N-CA-C	-7.19	95.13	113.10
1	A	3	GLY	N-CA-C	-5.36	99.71	113.10

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	4200	0	4004	79	0
1	B	4199	0	4002	74	0
2	A	289	0	0	6	0
2	B	285	0	0	4	0
All	All	8973	0	8006	152	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 9.

All (152) 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:113:LEU:HD22	1:A:113:LEU:H	1.08	1.13
1:B:45:LEU:H	1:B:326:ASN:HD21	1.11	0.97
1:A:113:LEU:H	1:A:113:LEU:CD2	1.81	0.93
1:A:436:MET:HG2	1:A:441:MET:HE3	1.53	0.89
1:A:113:LEU:HD22	1:A:113:LEU:N	1.89	0.86
1:B:436:MET:HG2	1:B:441:MET:HE3	1.61	0.83
1:A:49:GLU:HG3	1:A:173:ALA:HB2	1.62	0.79
1:A:145:LYS:NZ	1:A:165:THR:HG22	2.01	0.76
1:B:172:ASP:HB3	1:B:175:LEU:HD13	1.67	0.74
1:B:506:VAL:O	1:B:510:GLN:HG3	1.90	0.72
1:A:329:ARG:HH11	1:A:329:ARG:HB2	1.56	0.71
1:B:467:ASN:H	1:B:470:ASN:HD22	1.38	0.70
1:B:471:TYR:O	1:B:475:LEU:HB2	1.94	0.68
1:B:503:PHE:O	1:B:506:VAL:HG22	1.94	0.66
1:B:467:ASN:HD22	1:B:469:SER:H	1.43	0.66
1:B:369:TYR:OH	1:B:459:GLU:HG2	1.96	0.66
1:A:231:ARG:HG3	2:A:776:HOH:O	1.95	0.66
1:B:509:ARG:HH11	1:B:509:ARG:HG3	1.60	0.65
1:A:526:GLN:HG3	1:A:527:GLN:N	2.11	0.64
1:B:148:VAL:HG21	1:B:510:GLN:HB3	1.79	0.63
1:B:528:HIS:O	1:B:529:GLY:C	2.38	0.62
1:A:310:ARG:O	1:A:314:GLU:HG3	2.00	0.61
1:A:499:VAL:O	1:A:502:VAL:HG13	2.00	0.61
1:B:304:THR:HB	1:B:307:GLU:HG3	1.81	0.61
1:A:117:HIS:CD2	1:A:499:VAL:HG13	2.35	0.61
1:A:145:LYS:HZ1	1:A:165:THR:HG22	1.64	0.61
1:A:129:ARG:HD3	2:A:780:HOH:O	2.01	0.60
1:A:471:TYR:O	1:A:475:LEU:HB2	2.01	0.60
1:B:437:TYR:HB2	1:B:443:ILE:HD11	1.83	0.59
1:A:164:VAL:HG23	1:A:165:THR:HG23	1.84	0.58
1:B:45:LEU:N	1:B:326:ASN:HD21	1.91	0.58
1:A:117:HIS:HB3	1:A:125:TYR:CE2	2.39	0.57
1:A:499:VAL:HA	1:A:502:VAL:CG1	2.34	0.57
1:A:233:LEU:O	1:A:237:LYS:HB3	2.05	0.56
1:A:511:ARG:O	1:A:515:GLU:HG2	2.05	0.56
1:A:182:ARG:O	1:A:186:GLU:HB2	2.05	0.56
1:B:116:TYR:HD2	1:B:449:ASP:HB3	1.71	0.56
1:A:246:MET:HG3	1:A:333:GLY:HA2	1.88	0.55
1:A:459:GLU:HG3	1:A:463:ARG:HH11	1.72	0.55
1:B:125:TYR:CG	1:B:499:VAL:HG21	2.43	0.54
1:B:49:GLU:HG3	1:B:173:ALA:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:72:ARG:HH11	1:B:72:ARG:HG3	1.72	0.53
1:A:154:ARG:HD2	1:A:154:ARG:N	2.21	0.53
1:A:410:SER:O	1:A:422:LYS:NZ	2.42	0.53
1:B:306:ASP:HA	1:B:309:VAL:HG13	1.91	0.53
1:A:475:LEU:HG	1:A:480:LEU:HD23	1.90	0.53
1:B:129:ARG:HB3	1:B:134:LYS:HB3	1.90	0.53
1:A:459:GLU:O	1:A:463:ARG:HD3	2.08	0.53
1:B:526:GLN:C	1:B:528:HIS:H	2.12	0.53
1:B:305:GLU:O	1:B:309:VAL:HG12	2.10	0.52
1:B:304:THR:HG22	1:B:306:ASP:N	2.23	0.52
1:A:236:ASN:O	1:A:240:GLU:HA	2.09	0.52
1:B:72:ARG:HG3	1:B:72:ARG:NH1	2.24	0.52
1:B:128:ASP:OD2	1:B:132:ARG:NH2	2.42	0.52
1:A:114:LEU:HD21	1:A:123:SER:HB3	1.92	0.52
1:B:113:LEU:HD13	1:B:449:ASP:O	2.10	0.52
1:B:509:ARG:HG3	1:B:509:ARG:NH1	2.25	0.51
1:A:84:PHE:O	1:A:106:HIS:HA	2.10	0.51
1:B:250:LEU:HD13	1:B:353:VAL:HG23	1.93	0.51
1:A:455:TYR:HA	1:A:461:GLU:CD	2.31	0.51
1:B:255:ALA:HA	1:B:296:TYR:O	2.12	0.50
1:B:475:LEU:HG	1:B:480:LEU:HD23	1.93	0.50
1:B:438:ARG:HA	1:B:482:PRO:HA	1.93	0.50
1:A:171:PHE:CE1	1:A:176:VAL:HG21	2.47	0.50
1:B:271:PRO:HB3	1:B:520:LEU:HD22	1.93	0.50
1:A:527:GLN:O	1:A:528:HIS:O	2.29	0.49
1:A:18:TRP:CE2	1:A:181:ARG:HG3	2.47	0.49
1:B:29:GLN:NE2	2:B:709:HOH:O	2.44	0.49
1:A:125:TYR:CG	1:A:499:VAL:HG11	2.47	0.49
1:B:467:ASN:H	1:B:470:ASN:ND2	2.08	0.49
1:B:216:GLY:HA2	2:B:698:HOH:O	2.12	0.49
1:A:18:TRP:CZ2	1:A:181:ARG:HG3	2.48	0.49
1:A:310:ARG:HD2	1:A:314:GLU:OE2	2.13	0.48
1:B:53:PRO:HD3	1:B:111:PHE:CE1	2.48	0.47
1:A:276:ILE:HG21	1:A:315:MET:CE	2.45	0.47
1:A:276:ILE:HG21	1:A:315:MET:HE1	1.97	0.47
1:B:304:THR:HG22	1:B:307:GLU:H	1.78	0.47
1:B:467:ASN:ND2	1:B:469:SER:H	2.09	0.47
1:B:128:ASP:O	1:B:132:ARG:HG3	2.13	0.47
1:B:136:VAL:HG13	1:B:136:VAL:O	2.15	0.47
1:B:150:LEU:HD22	1:B:271:PRO:HG2	1.97	0.47
1:B:274:SER:HB2	1:B:285:LYS:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:ARG:HA	1:A:387:ARG:HE	1.79	0.47
1:A:113:LEU:CD2	1:A:113:LEU:N	2.55	0.46
1:A:207:ASN:ND2	2:A:669:HOH:O	2.47	0.46
1:A:507:LYS:HZ2	1:A:510:GLN:HE22	1.63	0.46
1:A:9:LEU:HD11	1:A:37:LEU:HD13	1.98	0.45
1:A:112:GLY:O	1:A:144:TYR:HE2	1.99	0.45
1:B:526:GLN:O	1:B:528:HIS:N	2.48	0.45
1:B:145:LYS:HA	1:B:510:GLN:HG2	1.99	0.45
1:A:229:GLY:HA2	1:A:348:THR:OG1	2.16	0.45
1:B:164:VAL:O	1:B:164:VAL:HG23	2.16	0.45
1:B:71:MET:HB2	1:B:76:ALA:HB3	1.99	0.45
1:A:503:PHE:O	1:A:506:VAL:CG1	2.64	0.45
1:A:503:PHE:O	1:A:506:VAL:HG13	2.17	0.45
1:B:305:GLU:O	1:B:309:VAL:CG1	2.65	0.45
1:A:158:ARG:HD2	1:A:162:SER:OG	2.17	0.45
1:B:164:VAL:HG11	1:B:517:LEU:HD21	1.99	0.45
1:B:304:THR:HG22	1:B:306:ASP:H	1.83	0.45
1:A:218:VAL:O	1:A:218:VAL:HG23	2.17	0.44
1:B:377:ASP:OD1	1:B:377:ASP:C	2.56	0.44
1:A:125:TYR:O	1:A:129:ARG:HG3	2.17	0.44
1:B:526:GLN:C	1:B:528:HIS:N	2.71	0.44
1:A:68:ASP:O	1:A:72:ARG:HG2	2.18	0.44
1:A:237:LYS:HD3	2:A:677:HOH:O	2.18	0.44
1:A:329:ARG:HH11	1:A:329:ARG:CB	2.28	0.44
1:A:437:TYR:HB2	1:A:443:ILE:HD11	2.00	0.44
1:B:235:ILE:HG23	1:B:236:ASN:N	2.32	0.44
1:A:153:ASN:ND2	2:A:591:HOH:O	2.51	0.44
1:A:49:GLU:CG	1:A:173:ALA:HB2	2.41	0.43
1:A:191:ARG:HD3	2:A:657:HOH:O	2.17	0.43
1:A:274:SER:HB2	1:A:285:LYS:HB3	1.99	0.43
1:B:53:PRO:HD3	1:B:111:PHE:CZ	2.53	0.43
1:A:5:ILE:HG23	1:A:222:ASP:HB2	2.01	0.43
1:B:157:ARG:NH1	2:B:741:HOH:O	2.52	0.43
1:B:446:PRO:HB2	1:B:448:SER:O	2.18	0.43
1:B:155:SER:HB2	1:B:520:LEU:HA	2.01	0.43
1:A:305:GLU:O	1:A:309:VAL:HG23	2.19	0.43
1:A:366:ALA:HA	1:A:458:PHE:CZ	2.54	0.43
1:B:5:ILE:HG23	1:B:222:ASP:HB2	2.00	0.42
1:A:53:PRO:HD3	1:A:111:PHE:CE1	2.54	0.42
1:B:114:LEU:HD22	1:B:122:LEU:HB3	2.02	0.42
1:B:118:GLU:O	1:B:119:GLN:HB2	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:LEU:HD11	1:B:503:PHE:CE1	2.55	0.42
1:B:467:ASN:N	1:B:470:ASN:HD22	2.12	0.42
1:A:145:LYS:CD	1:A:509:ARG:HG2	2.50	0.42
1:A:494:GLN:H	1:A:494:GLN:HG2	1.45	0.42
1:A:150:LEU:HD22	1:A:271:PRO:HG2	2.00	0.42
1:A:235:ILE:HG23	1:A:236:ASN:N	2.35	0.42
1:B:229:GLY:HA2	1:B:348:THR:OG1	2.19	0.42
1:A:132:ARG:HH12	1:A:134:LYS:HG3	1.85	0.42
1:B:170:HIS:CE1	1:B:287:PRO:HD2	2.54	0.42
1:A:250:LEU:HD13	1:A:353:VAL:HG23	2.02	0.42
1:A:86:ASN:HB3	1:A:98:ARG:NH2	2.35	0.42
1:B:67:GLU:O	1:B:71:MET:HG2	2.20	0.41
1:A:86:ASN:HA	1:A:105:ASP:OD2	2.21	0.41
1:B:247:SER:HA	1:B:250:LEU:O	2.19	0.41
1:B:468:ASN:HB3	2:B:633:HOH:O	2.19	0.41
1:A:463:ARG:NH2	1:B:464:ASN:OD1	2.53	0.41
1:B:146:GLU:HB2	1:B:147:PRO:HD3	2.03	0.41
1:A:65:ILE:HA	1:A:66:PRO:HD3	1.91	0.41
1:A:127:PHE:CE1	1:A:481:VAL:HB	2.55	0.41
1:A:149:ILE:O	1:A:154:ARG:HG2	2.21	0.41
1:A:36:LEU:HD23	1:A:36:LEU:C	2.41	0.41
1:A:127:PHE:CE2	1:A:482:PRO:HB2	2.56	0.41
1:A:329:ARG:HB2	1:A:329:ARG:NH1	2.32	0.41
1:B:137:GLU:HA	1:B:138:PRO:HD3	1.92	0.41
1:B:278:MET:HG3	1:B:296:TYR:CE1	2.56	0.41
1:B:289:LEU:O	1:B:291:ARG:HD2	2.21	0.40
1:A:137:GLU:OE2	1:A:507:LYS:NZ	2.54	0.40
1:B:410:SER:O	1:B:422:LYS:NZ	2.50	0.40
1:B:381:ASN:HA	1:B:382:PRO:HD3	1.97	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	519/550 (94%)	506 (98%)	13 (2%)	0	100	100
1	B	520/550 (94%)	500 (96%)	19 (4%)	1 (0%)	47	58
All	All	1039/1100 (94%)	1006 (97%)	32 (3%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	527	GLN

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	441/463 (95%)	418 (95%)	23 (5%)	23	32
1	B	440/463 (95%)	412 (94%)	28 (6%)	17	23
All	All	881/926 (95%)	830 (94%)	51 (6%)	20	27

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	99	GLU
1	A	113	LEU
1	A	136	VAL
1	A	195	ASP
1	A	211	VAL
1	A	246	MET
1	A	256	VAL
1	A	291	ARG
1	A	329	ARG
1	A	373	LYS
1	A	376	PRO
1	A	380	LEU
1	A	387	ARG
1	A	425	ARG

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Mol	Chain	Res	Type
1	A	448	SER
1	A	475	LEU
1	A	494	GLN
1	A	497	GLU
1	A	499	VAL
1	A	502	VAL
1	A	525	ARG
1	A	526	GLN
1	B	2	SER
1	B	114	LEU
1	B	119	GLN
1	B	132	ARG
1	B	136	VAL
1	B	157	ARG
1	B	178	ASP
1	B	211	VAL
1	B	233	LEU
1	B	289	LEU
1	B	291	ARG
1	B	304	THR
1	B	309	VAL
1	B	330	PHE
1	B	376	PRO
1	B	380	LEU
1	B	390	ARG
1	B	393	GLU
1	B	457	ASN
1	B	460	GLU
1	B	461	GLU
1	B	464	ASN
1	B	472	TYR
1	B	474	VAL
1	B	475	LEU
1	B	488	ARG
1	B	511	ARG
1	B	514	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	153	ASN

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Mol	Chain	Res	Type
1	A	166	ASN
1	A	205	ASN
1	A	207	ASN
1	A	252	ASN
1	A	510	GLN
1	A	526	GLN
1	B	29	GLN
1	B	38	GLN
1	B	124	HIS
1	B	153	ASN
1	B	166	ASN
1	B	205	ASN
1	B	207	ASN
1	B	252	ASN
1	B	326	ASN
1	B	457	ASN
1	B	467	ASN
1	B	470	ASN
1	B	494	GLN

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 monosaccharides 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(Å²)	Q<0.9
1	A	523/550 (95%)	-0.19	11 (2%)	63	70	19, 34, 58, 96	0
1	B	524/550 (95%)	-0.14	13 (2%)	57	64	21, 37, 66, 96	0
All	All	1047/1100 (95%)	-0.17	24 (2%)	60	67	19, 35, 64, 96	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	528	HIS	5.3
1	A	2	SER	4.2
1	B	449	ASP	4.2
1	B	501	GLU	3.4
1	A	447	ALA	2.9
1	B	527	GLN	2.9
1	A	449	ASP	2.9
1	B	525	ARG	2.9
1	B	447	ALA	2.9
1	B	497	GLU	2.8
1	A	425	ARG	2.7
1	B	494	GLN	2.7
1	A	455	TYR	2.6
1	B	528	HIS	2.6
1	B	450	ASP	2.5
1	B	3	GLY	2.4
1	A	515	GLU	2.3
1	B	2	SER	2.3
1	A	525	ARG	2.2
1	B	456	GLY	2.1
1	A	72	ARG	2.1
1	B	72	ARG	2.0
1	A	134	LYS	2.0
1	A	463	ARG	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 ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.