



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 07:46 pm BST

PDB ID : 1GQ7
Title : PROCLAVAMINATE AMIDINO HYDROLASE FROM STREPTOMYCES CLAVULIGERUS
Authors : Elkins, J.M.; Clifton, I.J.; Hernandez, H.; Robinson, C.V.; Schofield, C.J.; Hewitson, K.S.
Deposited on : 2001-11-20
Resolution : 2.45 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

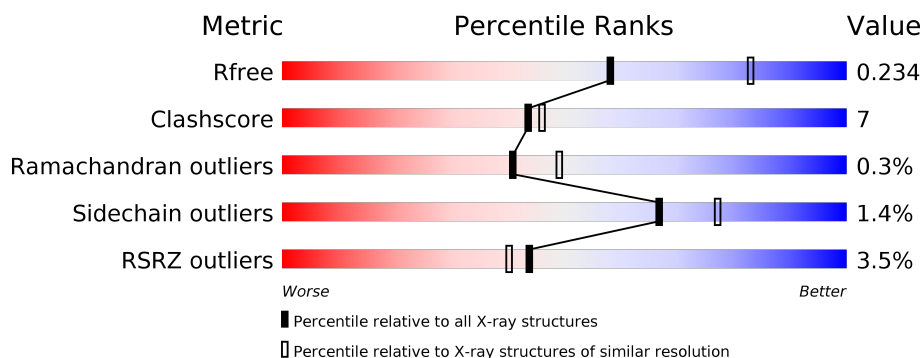
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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	313	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>12%</div> <div></div> </div> <div></div> </div>
1	B	313	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div></div> </div> <div></div> </div>
1	C	313	<div> <div>8%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div></div> </div> <div></div> </div>
1	D	313	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div></div> </div> <div></div> </div>
1	E	313	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div></div> </div> <div></div> </div>
1	F	313	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div></div> </div> <div></div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13457 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 PROCLAVAMINATE AMIDINO HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	S	0	0	0
			2185	1385	382	411	7			
1	B	301	Total	C	N	O	S	0	0	0
			2185	1385	382	411	7			
1	C	301	Total	C	N	O	S	0	0	0
			2185	1385	382	411	7			
1	D	301	Total	C	N	O	S	0	0	0
			2185	1385	382	411	7			
1	E	301	Total	C	N	O	S	0	0	0
			2188	1386	382	413	7			
1	F	301	Total	C	N	O	S	0	0	0
			2185	1385	382	411	7			

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total	Mn	0	0
			2	2		
2	E	2	Total	Mn	0	0
			2	2		
2	B	2	Total	Mn	0	0
			2	2		
2	C	2	Total	Mn	0	0
			2	2		
2	A	2	Total	Mn	0	0
			2	2		
2	F	2	Total	Mn	0	0
			2	2		

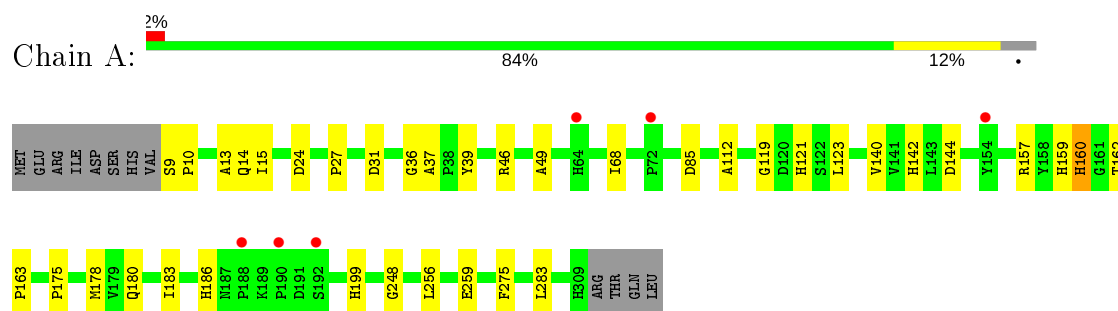
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	59	Total 59	O 59	0	0
3	B	41	Total 41	O 41	0	0
3	C	29	Total 29	O 29	0	0
3	D	61	Total 61	O 61	0	0
3	E	76	Total 76	O 76	0	0
3	F	66	Total 66	O 66	0	0

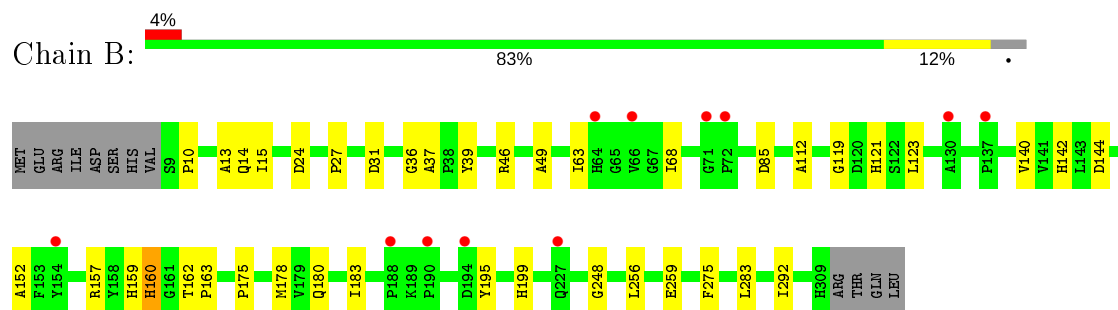
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

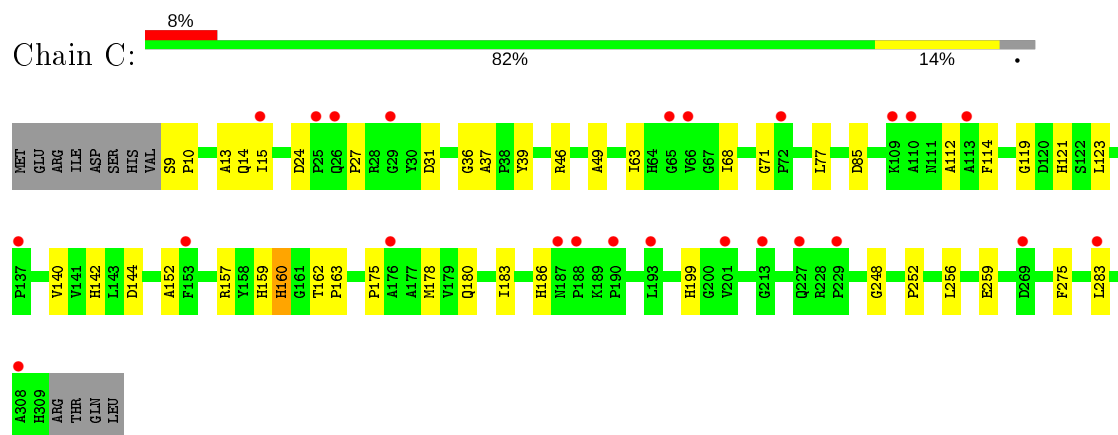
• Molecule 1: PROCLAVAMINATE AMIDINO HYDROLASE



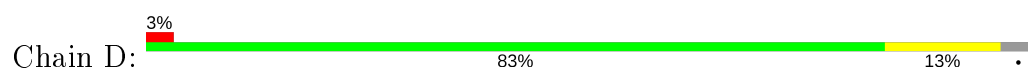
• Molecule 1: PROCLAVAMINATE AMIDINO HYDROLASE

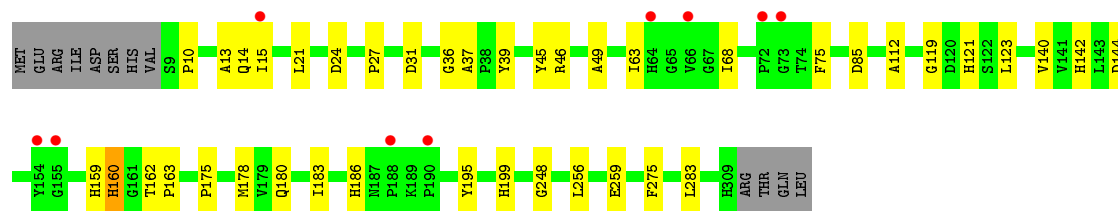


• Molecule 1: PROCLAVAMINATE AMIDINO HYDROLASE

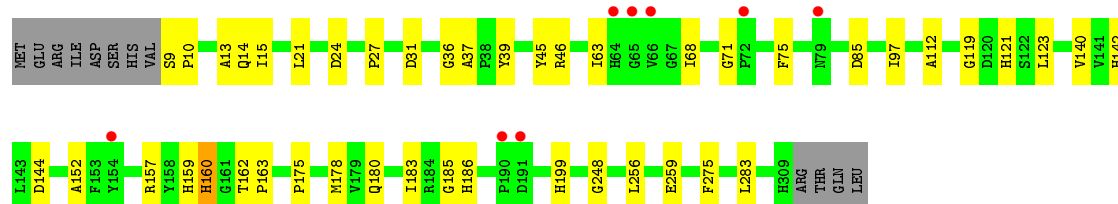
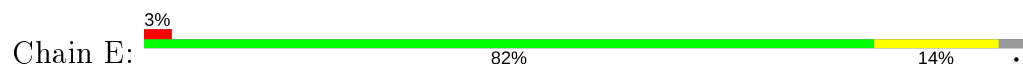


• Molecule 1: PROCLAVAMINATE AMIDINO HYDROLASE

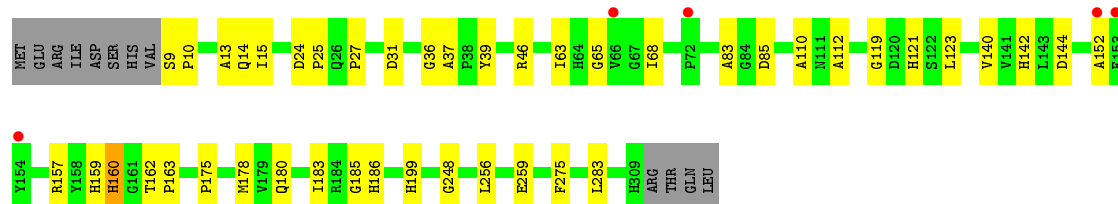
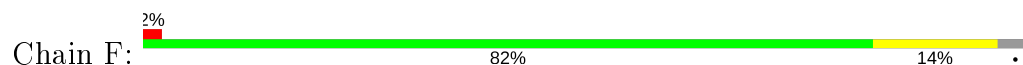




● Molecule 1: PROCLAVAMINATE AMIDINO HYDROLASE



● Molecule 1: PROCLAVAMINATE AMIDINO HYDROLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.92Å 81.35Å 120.39Å 90.00° 99.55° 90.00°	Depositor
Resolution (Å)	35.59 – 2.45 35.59 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.5 (35.59-2.45) 98.6 (35.59-2.45)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.45Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.222 , 0.236 0.221 , 0.234	Depositor DCC
R_{free} test set	3296 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13457	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.31 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7970e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: MN

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	0.39	0/2241	0.64	0/3068
1	B	0.39	0/2241	0.65	0/3068
1	C	0.39	0/2241	0.65	0/3068
1	D	0.39	0/2241	0.65	0/3068
1	E	0.39	0/2244	0.65	0/3072
1	F	0.39	0/2241	0.64	0/3068
All	All	0.39	0/13449	0.65	0/18412

There are no bond length outliers.

There are no bond angle outliers.

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	2185	0	2085	28	0
1	B	2185	0	2085	30	0
1	C	2185	0	2085	36	0
1	D	2185	0	2085	32	0
1	E	2188	0	2087	34	0
1	F	2185	0	2085	36	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	59	0	0	2	0
3	B	41	0	0	1	0
3	C	29	0	0	3	0
3	D	61	0	0	2	0
3	E	76	0	0	3	0
3	F	66	0	0	4	0
All	All	13457	0	12512	172	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 7.

All (172) 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:283:LEU:HD22	1:A:283:LEU:H	1.42	0.82
1:F:283:LEU:H	1:F:283:LEU:HD22	1.45	0.81
1:C:283:LEU:H	1:C:283:LEU:HD22	1.46	0.80
1:B:283:LEU:H	1:B:283:LEU:HD22	1.47	0.80
1:E:283:LEU:H	1:E:283:LEU:HD22	1.47	0.79
1:D:283:LEU:H	1:D:283:LEU:HD22	1.46	0.78
1:C:162:THR:N	1:C:163:PRO:HD2	2.08	0.69
1:F:162:THR:N	1:F:163:PRO:HD2	2.10	0.67
1:A:162:THR:N	1:A:163:PRO:HD2	2.10	0.66
1:E:162:THR:N	1:E:163:PRO:HD2	2.10	0.65
1:B:15:ILE:HG22	1:D:15:ILE:HG22	1.79	0.65
1:B:162:THR:N	1:B:163:PRO:HD2	2.11	0.65
1:D:162:THR:N	1:D:163:PRO:HD2	2.12	0.64
1:C:15:ILE:HB	3:C:2001:HOH:O	1.97	0.64
1:D:15:ILE:HG12	3:D:2003:HOH:O	1.98	0.62
1:D:39:TYR:O	1:D:119:GLY:HA3	2.01	0.61
1:D:283:LEU:N	1:D:283:LEU:HD22	2.17	0.60
1:C:39:TYR:O	1:C:119:GLY:HA3	2.01	0.60
1:C:15:ILE:CG2	1:F:15:ILE:HG22	2.32	0.60
1:E:68:ILE:HD13	1:F:183:ILE:HG22	1.83	0.60
1:B:68:ILE:HD13	1:C:183:ILE:HG22	1.82	0.59
1:B:283:LEU:N	1:B:283:LEU:HD22	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:TYR:O	1:B:119:GLY:HA3	2.03	0.58
1:A:15:ILE:HG22	1:E:15:ILE:HG22	1.85	0.58
1:C:15:ILE:HG22	1:F:15:ILE:CG2	2.33	0.58
1:C:15:ILE:HG22	1:F:15:ILE:HG22	1.86	0.58
1:B:63:ILE:HG23	1:C:186:HIS:NE2	2.19	0.58
1:F:39:TYR:O	1:F:119:GLY:HA3	2.02	0.58
1:A:39:TYR:O	1:A:119:GLY:HA3	2.03	0.58
1:E:63:ILE:HG23	1:F:186:HIS:NE2	2.18	0.58
1:E:283:LEU:HD22	1:E:283:LEU:N	2.18	0.57
1:E:39:TYR:O	1:E:119:GLY:HA3	2.05	0.56
1:A:283:LEU:HD22	1:A:283:LEU:N	2.14	0.56
1:C:283:LEU:N	1:C:283:LEU:HD22	2.18	0.56
1:F:283:LEU:N	1:F:283:LEU:HD22	2.16	0.56
1:E:71:GLY:HA3	3:E:2023:HOH:O	2.05	0.56
1:C:162:THR:N	1:C:163:PRO:CD	2.68	0.56
1:A:283:LEU:CD2	1:A:283:LEU:H	2.17	0.56
1:E:162:THR:N	1:E:163:PRO:CD	2.69	0.55
1:E:13:ALA:O	1:E:14:GLN:HB2	2.08	0.54
1:F:162:THR:N	1:F:163:PRO:CD	2.70	0.54
1:D:31:ASP:O	1:D:112:ALA:HB3	2.08	0.54
1:F:65:GLY:HA3	3:F:2014:HOH:O	2.06	0.54
1:A:256:LEU:HB2	1:A:259:GLU:HG3	1.88	0.54
1:A:162:THR:N	1:A:163:PRO:CD	2.69	0.54
1:B:31:ASP:O	1:B:112:ALA:HB3	2.08	0.54
1:D:162:THR:N	1:D:163:PRO:CD	2.71	0.54
1:A:183:ILE:HG22	1:C:68:ILE:HD13	1.89	0.53
1:B:15:ILE:CG2	1:D:15:ILE:HG22	2.38	0.53
1:B:162:THR:N	1:B:163:PRO:CD	2.70	0.53
1:D:283:LEU:H	1:D:283:LEU:CD2	2.19	0.53
1:F:13:ALA:O	1:F:14:GLN:HB2	2.08	0.53
1:B:15:ILE:HG22	1:D:15:ILE:CG2	2.38	0.53
1:B:63:ILE:HG23	1:C:186:HIS:CD2	2.44	0.53
1:B:283:LEU:CD2	1:B:283:LEU:H	2.21	0.53
1:D:13:ALA:O	1:D:14:GLN:HB2	2.09	0.53
1:F:37:ALA:HB1	1:F:123:LEU:HD13	1.91	0.53
1:A:13:ALA:O	1:A:14:GLN:HB2	2.08	0.52
1:C:31:ASP:O	1:C:112:ALA:HB3	2.10	0.52
1:F:283:LEU:H	1:F:283:LEU:CD2	2.19	0.52
1:C:13:ALA:O	1:C:14:GLN:HB2	2.08	0.52
1:B:13:ALA:O	1:B:14:GLN:HB2	2.10	0.51
1:B:140:VAL:HB	1:B:178:MET:HG3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:283:LEU:H	1:C:283:LEU:CD2	2.20	0.51
1:A:37:ALA:HB1	1:A:123:LEU:HD13	1.93	0.51
1:B:256:LEU:HB2	1:B:259:GLU:HG3	1.93	0.51
1:E:63:ILE:HG23	1:F:186:HIS:CD2	2.45	0.51
1:D:256:LEU:HB2	1:D:259:GLU:HG3	1.92	0.51
1:E:31:ASP:O	1:E:112:ALA:HB3	2.10	0.51
1:E:283:LEU:CD2	1:E:283:LEU:H	2.20	0.51
1:F:24:ASP:O	1:F:27:PRO:HD3	2.11	0.51
1:A:15:ILE:CG2	1:E:15:ILE:HG22	2.41	0.51
1:D:37:ALA:HB1	1:D:123:LEU:HD13	1.93	0.51
1:E:256:LEU:HB2	1:E:259:GLU:HG3	1.93	0.51
1:A:31:ASP:O	1:A:112:ALA:HB3	2.12	0.50
1:A:15:ILE:HG22	1:E:15:ILE:CG2	2.41	0.50
1:F:110:ALA:HB3	3:F:2024:HOH:O	2.11	0.50
1:D:10:PRO:HG2	1:D:13:ALA:HB2	1.93	0.50
1:E:37:ALA:HB1	1:E:123:LEU:HD13	1.93	0.50
1:B:10:PRO:HG2	1:B:13:ALA:HB2	1.93	0.50
1:F:31:ASP:O	1:F:112:ALA:HB3	2.12	0.49
1:A:159:HIS:CE1	1:A:162:THR:HG23	2.47	0.49
1:C:256:LEU:HB2	1:C:259:GLU:HG3	1.92	0.49
1:C:77:LEU:HD11	3:C:2009:HOH:O	2.12	0.49
1:F:256:LEU:HB2	1:F:259:GLU:HG3	1.93	0.49
1:D:142:HIS:O	1:D:180:GLN:HA	2.13	0.49
1:E:10:PRO:HG2	1:E:13:ALA:HB2	1.94	0.49
1:F:10:PRO:HG2	1:F:13:ALA:HB2	1.94	0.49
1:D:24:ASP:O	1:D:27:PRO:HD3	2.13	0.49
1:C:37:ALA:HB1	1:C:123:LEU:HD13	1.94	0.49
1:D:68:ILE:HD13	1:E:183:ILE:HG22	1.93	0.48
1:E:140:VAL:HB	1:E:178:MET:HG3	1.95	0.48
1:B:142:HIS:O	1:B:180:GLN:HA	2.13	0.48
1:A:10:PRO:HG2	1:A:13:ALA:HB2	1.95	0.48
1:A:142:HIS:O	1:A:180:GLN:HA	2.12	0.48
1:B:24:ASP:O	1:B:27:PRO:HD3	2.14	0.48
1:E:24:ASP:O	1:E:27:PRO:HD3	2.13	0.48
1:C:10:PRO:HG2	1:C:13:ALA:HB2	1.95	0.48
1:D:159:HIS:CE1	1:D:162:THR:HG23	2.49	0.48
1:F:142:HIS:O	1:F:180:GLN:HA	2.13	0.48
1:A:140:VAL:HB	1:A:178:MET:HG3	1.94	0.48
1:C:142:HIS:O	1:C:180:GLN:HA	2.13	0.48
1:D:186:HIS:NE2	1:F:63:ILE:HG23	2.29	0.48
1:B:37:ALA:HB1	1:B:123:LEU:HD13	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:ARG:NH1	3:A:2034:HOH:O	2.46	0.47
1:A:24:ASP:O	1:A:27:PRO:HD3	2.13	0.47
1:E:142:HIS:O	1:E:180:GLN:HA	2.14	0.47
1:C:46:ARG:HB2	1:C:248:GLY:HA2	1.96	0.47
1:A:142:HIS:NE2	1:A:144:ASP:HB2	2.30	0.47
1:D:140:VAL:HB	1:D:178:MET:HG3	1.96	0.47
1:E:46:ARG:HB2	1:E:248:GLY:HA2	1.96	0.47
1:F:140:VAL:HB	1:F:178:MET:HG3	1.96	0.47
3:A:2056:HOH:O	1:F:283:LEU:CD2	2.62	0.46
1:F:142:HIS:NE2	1:F:144:ASP:HB2	2.30	0.46
1:D:46:ARG:HB2	1:D:248:GLY:HA2	1.98	0.46
1:C:36:GLY:HA3	1:C:85:ASP:OD1	2.16	0.46
1:B:159:HIS:CE1	1:B:162:THR:HG23	2.51	0.46
1:C:24:ASP:O	1:C:27:PRO:HD3	2.15	0.46
1:A:186:HIS:NE2	1:C:63:ILE:HG23	2.31	0.45
1:C:159:HIS:CE1	1:C:162:THR:HG23	2.51	0.45
1:C:140:VAL:HB	1:C:178:MET:HG3	1.97	0.45
1:A:68:ILE:HD13	1:B:183:ILE:HG22	1.98	0.45
1:F:159:HIS:CE1	1:F:162:THR:HG23	2.51	0.45
1:E:159:HIS:CE1	1:E:162:THR:HG23	2.52	0.45
1:C:142:HIS:NE2	1:C:144:ASP:HB2	2.32	0.45
1:C:175:PRO:HB2	1:C:199:HIS:O	2.17	0.45
1:C:71:GLY:HA3	3:C:2009:HOH:O	2.17	0.45
1:A:36:GLY:HA3	1:A:85:ASP:OD1	2.17	0.45
1:D:142:HIS:NE2	1:D:144:ASP:HB2	2.32	0.45
1:A:175:PRO:HB2	1:A:199:HIS:O	2.18	0.44
1:A:46:ARG:HB2	1:A:248:GLY:HA2	2.00	0.44
1:F:36:GLY:HA3	1:F:85:ASP:OD1	2.18	0.44
1:B:142:HIS:NE2	1:B:144:ASP:HB2	2.32	0.44
1:E:36:GLY:HA3	1:E:85:ASP:OD1	2.17	0.44
1:F:175:PRO:HB2	1:F:199:HIS:O	2.17	0.44
1:F:46:ARG:HB2	1:F:248:GLY:HA2	1.99	0.44
1:B:292:ILE:HG22	1:C:252:PRO:CG	2.48	0.44
1:B:175:PRO:HB2	1:B:199:HIS:O	2.18	0.44
1:B:36:GLY:HA3	1:B:85:ASP:OD1	2.18	0.44
1:B:46:ARG:HB2	1:B:248:GLY:HA2	2.00	0.43
1:E:97:ILE:HG22	3:E:2038:HOH:O	2.17	0.43
1:E:142:HIS:NE2	1:E:144:ASP:HB2	2.32	0.43
1:E:185:GLY:HA3	3:E:2045:HOH:O	2.18	0.43
1:E:175:PRO:HB2	1:E:199:HIS:O	2.18	0.43
1:D:195:TYR:HB3	3:D:2029:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:183:ILE:HG22	1:F:68:ILE:HD13	2.00	0.43
1:F:9:SER:HA	1:F:10:PRO:HD2	1.91	0.42
1:F:152:ALA:HB1	1:F:157:ARG:HA	2.02	0.42
1:F:83:ALA:HB1	3:F:2022:HOH:O	2.17	0.42
1:E:9:SER:HA	1:E:10:PRO:HD2	1.91	0.42
1:A:9:SER:HA	1:A:10:PRO:HD2	1.91	0.42
1:D:175:PRO:HB2	1:D:199:HIS:O	2.20	0.42
1:D:21:LEU:HD21	1:D:75:PHE:CE1	2.54	0.41
1:D:63:ILE:HG23	1:E:186:HIS:NE2	2.36	0.41
1:E:45:TYR:HB3	1:E:46:ARG:H	1.75	0.41
1:B:195:TYR:HB3	3:B:2025:HOH:O	2.21	0.41
1:D:45:TYR:HB3	1:D:46:ARG:H	1.75	0.41
1:B:152:ALA:HB1	1:B:157:ARG:HA	2.02	0.41
1:D:49:ALA:HB1	1:D:119:GLY:HA2	2.03	0.41
1:D:36:GLY:HA3	1:D:85:ASP:OD1	2.20	0.41
1:A:49:ALA:HB1	1:A:119:GLY:HA2	2.03	0.41
1:C:152:ALA:HB1	1:C:157:ARG:HA	2.03	0.41
1:F:185:GLY:HA3	3:F:2038:HOH:O	2.20	0.41
1:C:114:PHE:C	1:C:114:PHE:CD1	2.95	0.41
1:C:49:ALA:HB1	1:C:119:GLY:HA2	2.03	0.41
1:D:10:PRO:HG2	1:D:13:ALA:CB	2.50	0.41
1:E:21:LEU:HD21	1:E:75:PHE:CE1	2.56	0.40
1:B:49:ALA:HB1	1:B:119:GLY:HA2	2.02	0.40
1:F:24:ASP:HA	1:F:25:PRO:HD2	1.96	0.40
1:E:152:ALA:HB1	1:E:157:ARG:HA	2.03	0.40
1:C:9:SER:HA	1:C:10:PRO:HD2	1.92	0.40
1:C:15:ILE:CG2	1:F:15:ILE:CG2	2.96	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	299/313 (96%)	285 (95%)	13 (4%)	1 (0%)	41	49
1	B	299/313 (96%)	284 (95%)	14 (5%)	1 (0%)	41	49
1	C	299/313 (96%)	286 (96%)	12 (4%)	1 (0%)	41	49
1	D	299/313 (96%)	285 (95%)	13 (4%)	1 (0%)	41	49
1	E	299/313 (96%)	285 (95%)	13 (4%)	1 (0%)	41	49
1	F	299/313 (96%)	284 (95%)	14 (5%)	1 (0%)	41	49
All	All	1794/1878 (96%)	1709 (95%)	79 (4%)	6 (0%)	41	49

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	160	HIS
1	B	160	HIS
1	C	160	HIS
1	D	160	HIS
1	E	160	HIS
1	F	160	HIS

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	218/245 (89%)	215 (99%)	3 (1%)	67	77
1	B	218/245 (89%)	215 (99%)	3 (1%)	67	77
1	C	218/245 (89%)	215 (99%)	3 (1%)	67	77
1	D	218/245 (89%)	215 (99%)	3 (1%)	67	77
1	E	219/245 (89%)	216 (99%)	3 (1%)	67	77
1	F	218/245 (89%)	215 (99%)	3 (1%)	67	77
All	All	1309/1470 (89%)	1291 (99%)	18 (1%)	67	77

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	HIS
1	A	160	HIS
1	A	275	PHE
1	B	121	HIS
1	B	160	HIS
1	B	275	PHE
1	C	121	HIS
1	C	160	HIS
1	C	275	PHE
1	D	121	HIS
1	D	160	HIS
1	D	275	PHE
1	E	121	HIS
1	E	160	HIS
1	E	275	PHE
1	F	121	HIS
1	F	160	HIS
1	F	275	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	HIS
1	A	187	ASN
1	B	64	HIS
1	B	187	ASN
1	C	64	HIS
1	C	187	ASN
1	D	64	HIS
1	D	187	ASN
1	D	199	HIS
1	D	227	GLN
1	E	64	HIS
1	E	187	ASN
1	F	64	HIS
1	F	187	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 12 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 [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 ⓘ

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	301/313 (96%)	0.02	6 (1%) 65 62	16, 30, 55, 68	0
1	B	301/313 (96%)	0.11	11 (3%) 41 38	18, 33, 55, 67	0
1	C	301/313 (96%)	0.47	24 (7%) 12 9	19, 34, 55, 67	0
1	D	301/313 (96%)	-0.06	9 (2%) 50 46	15, 31, 54, 67	0
1	E	301/313 (96%)	0.02	8 (2%) 54 50	15, 29, 54, 67	0
1	F	301/313 (96%)	0.05	5 (1%) 70 67	17, 31, 54, 66	0
All	All	1806/1878 (96%)	0.10	63 (3%) 44 40	15, 32, 55, 68	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	154	TYR	7.0
1	E	154	TYR	4.5
1	F	153	PHE	4.4
1	D	64	HIS	4.1
1	F	72	PRO	4.1
1	E	64	HIS	4.0
1	B	72	PRO	3.9
1	D	72	PRO	3.7
1	B	66	VAL	3.5
1	A	64	HIS	3.4
1	D	66	VAL	3.2
1	D	154	TYR	3.1
1	D	190	PRO	3.1
1	E	72	PRO	3.0
1	E	191	ASP	3.0
1	A	154	TYR	2.9
1	B	154	TYR	2.9
1	C	72	PRO	2.9
1	C	66	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	176	ALA	2.8
1	C	109	LYS	2.8
1	D	15	ILE	2.7
1	D	188	PRO	2.7
1	A	190	PRO	2.5
1	C	269	ASP	2.5
1	C	137	PRO	2.5
1	F	66	VAL	2.5
1	A	72	PRO	2.5
1	E	190	PRO	2.5
1	C	227	GLN	2.5
1	A	192	SER	2.5
1	C	113	ALA	2.4
1	C	193	LEU	2.4
1	C	15	ILE	2.4
1	C	25	PRO	2.4
1	B	190	PRO	2.3
1	C	201	VAL	2.3
1	C	153	PHE	2.3
1	B	64	HIS	2.3
1	B	130	ALA	2.3
1	D	155	GLY	2.2
1	E	79	ASN	2.2
1	D	73	GLY	2.2
1	B	71	GLY	2.2
1	C	110	ALA	2.1
1	C	283	LEU	2.1
1	C	213	GLY	2.1
1	C	308	ALA	2.1
1	C	229	PRO	2.1
1	C	29	GLY	2.1
1	C	188	PRO	2.1
1	B	194	ASP	2.1
1	B	227	GLN	2.1
1	B	137	PRO	2.1
1	C	187	ASN	2.1
1	B	188	PRO	2.0
1	F	152	ALA	2.0
1	C	65	GLY	2.0
1	E	65	GLY	2.0
1	E	66	VAL	2.0
1	A	188	PRO	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	26	GLN	2.0
1	C	190	PRO	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 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. 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(\AA^2)	Q<0.9
2	MN	C	350	1/1	0.95	0.10	34,34,34,34	0
2	MN	F	351	1/1	0.95	0.11	28,28,28,28	0
2	MN	A	351	1/1	0.96	0.08	34,34,34,34	0
2	MN	E	350	1/1	0.97	0.07	23,23,23,23	0
2	MN	A	350	1/1	0.97	0.07	29,29,29,29	0
2	MN	B	351	1/1	0.98	0.14	32,32,32,32	0
2	MN	D	350	1/1	0.98	0.08	31,31,31,31	0
2	MN	D	351	1/1	0.98	0.07	28,28,28,28	0
2	MN	C	351	1/1	0.98	0.12	35,35,35,35	0
2	MN	F	350	1/1	0.98	0.12	26,26,26,26	0
2	MN	B	350	1/1	0.98	0.09	37,37,37,37	0
2	MN	E	351	1/1	0.99	0.12	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.