



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:09 PM UTC

PDB ID : 1XEB / pdb_00001xeb
Title : Crystal Structure of an Acyl-CoA N-acyltransferase from *Pseudomonas aeruginosa*
Authors : Bertero, M.G.; Walker, J.R.; Skarina, T.; Gorodichtchenskaia, E.; Joachimiak, A.; Edwards, A.E.; Savchenko, A.; Strynadka, N.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2004-09-09
Resolution : 2.35 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

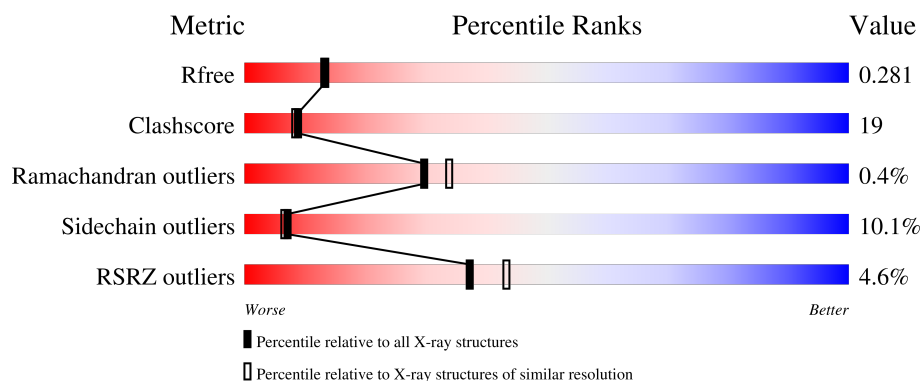
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	150	
1	B	150	
1	C	150	
1	D	150	

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Mol	Chain	Length	Quality of chain
1	E	150	7% 58% 33% 7%
1	F	150	5% 57% 34% 6%
1	G	150	7% 58% 35%
1	H	150	10% 57% 31% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9702 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 hypothetical protein PA0115.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	Se	0	0	0
			1196	760	214	217	2	3			
1	B	148	Total	C	N	O	S	Se	0	0	0
			1192	758	213	216	2	3			
1	C	149	Total	C	N	O	S	Se	0	0	0
			1196	760	214	217	2	3			
1	D	148	Total	C	N	O	S	Se	0	0	0
			1190	757	213	215	2	3			
1	E	146	Total	C	N	O	S	Se	0	0	0
			1172	746	211	210	2	3			
1	F	145	Total	C	N	O	S	Se	0	0	0
			1173	748	209	211	2	3			
1	G	147	Total	C	N	O	S	Se	0	0	0
			1184	754	212	213	2	3			
1	H	146	Total	C	N	O	S	Se	0	0	0
			1177	750	210	212	2	3			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	53	MSE	MET	modified residue	UNP Q9I717
A	97	MSE	MET	modified residue	UNP Q9I717
A	147	MSE	MET	modified residue	UNP Q9I717
B	53	MSE	MET	modified residue	UNP Q9I717
B	97	MSE	MET	modified residue	UNP Q9I717
B	147	MSE	MET	modified residue	UNP Q9I717
C	53	MSE	MET	modified residue	UNP Q9I717
C	97	MSE	MET	modified residue	UNP Q9I717
C	147	MSE	MET	modified residue	UNP Q9I717
D	53	MSE	MET	modified residue	UNP Q9I717
D	97	MSE	MET	modified residue	UNP Q9I717
D	147	MSE	MET	modified residue	UNP Q9I717
E	53	MSE	MET	modified residue	UNP Q9I717

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Chain	Residue	Modelled	Actual	Comment	Reference
E	97	MSE	MET	modified residue	UNP Q9I717
E	147	MSE	MET	modified residue	UNP Q9I717
F	53	MSE	MET	modified residue	UNP Q9I717
F	97	MSE	MET	modified residue	UNP Q9I717
F	147	MSE	MET	modified residue	UNP Q9I717
G	53	MSE	MET	modified residue	UNP Q9I717
G	97	MSE	MET	modified residue	UNP Q9I717
G	147	MSE	MET	modified residue	UNP Q9I717
H	53	MSE	MET	modified residue	UNP Q9I717
H	97	MSE	MET	modified residue	UNP Q9I717
H	147	MSE	MET	modified residue	UNP Q9I717

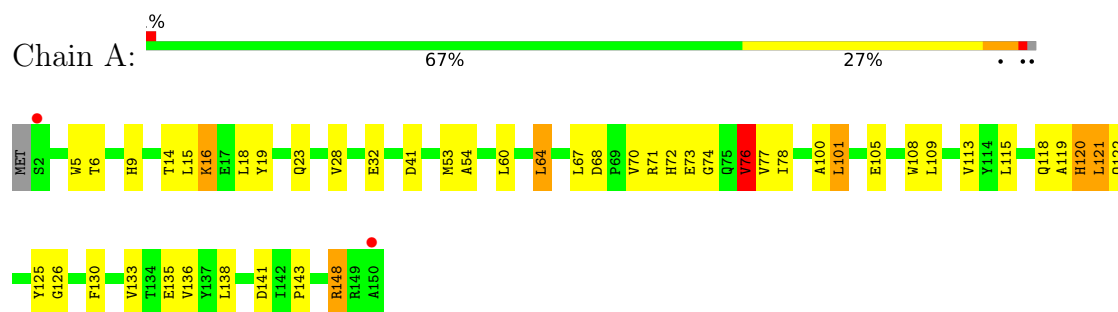
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	47	Total O 47 47	0	0
2	B	37	Total O 37 37	0	0
2	C	39	Total O 39 39	0	0
2	D	17	Total O 17 17	0	0
2	E	27	Total O 27 27	0	0
2	F	34	Total O 34 34	0	0
2	G	14	Total O 14 14	0	0
2	H	7	Total O 7 7	0	0

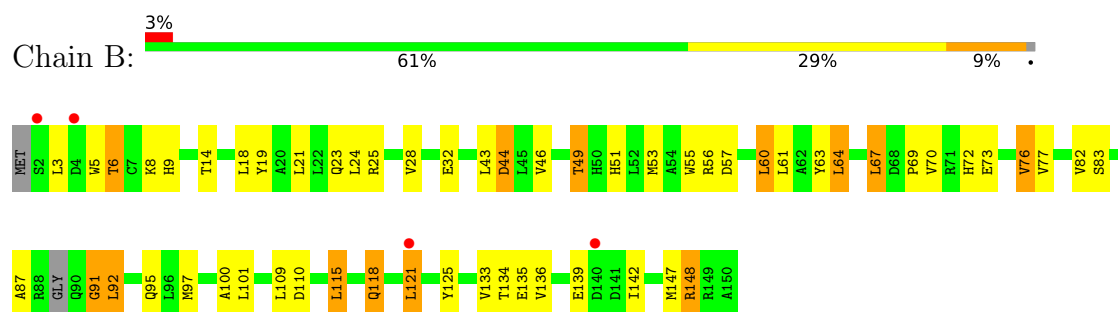
3 Residue-property plots

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.

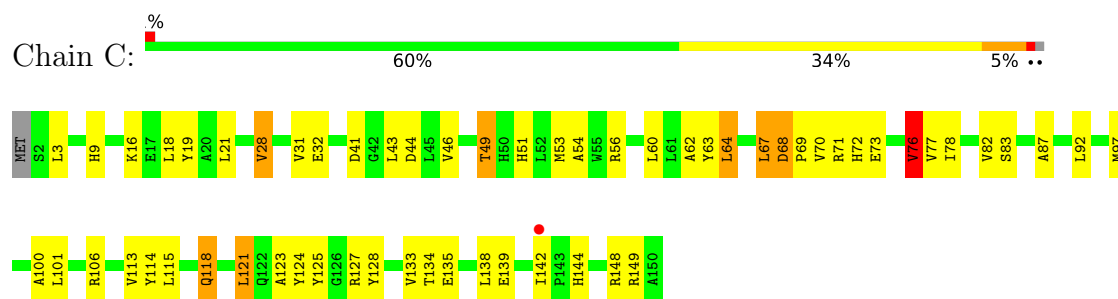
- Molecule 1: hypothetical protein PA0115



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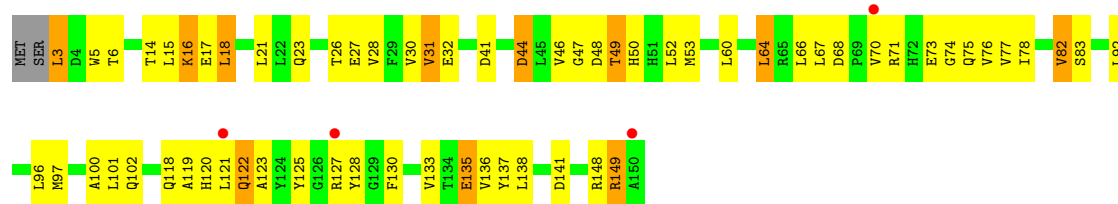


- Molecule 1: hypothetical protein PA0115

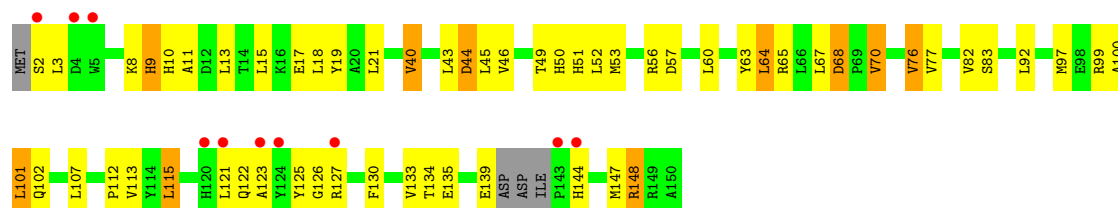


- Molecule 1: hypothetical protein PA0115

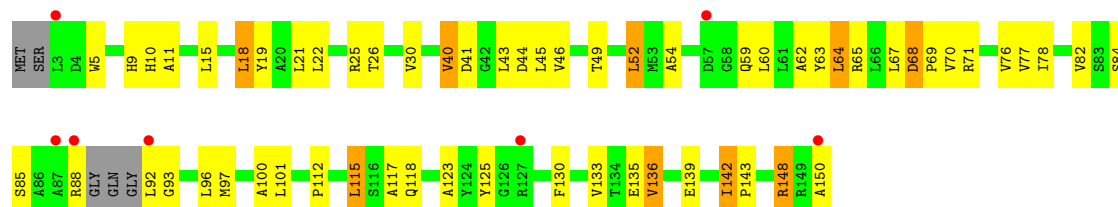




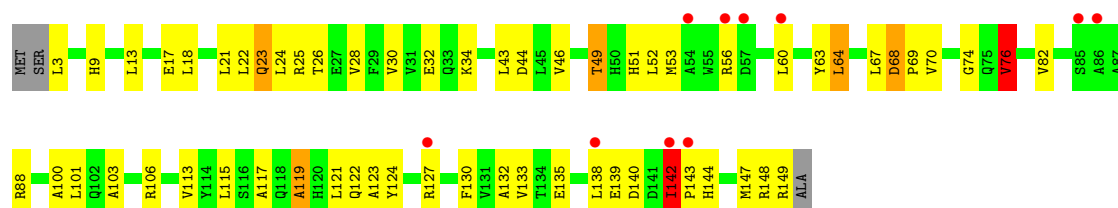
• Molecule 1: hypothetical protein PA0115



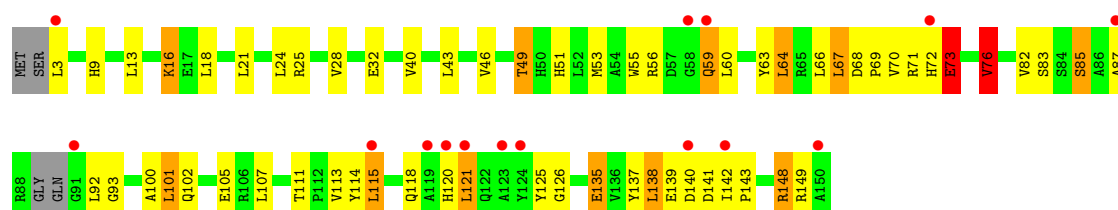
• Molecule 1: hypothetical protein PA0115



• Molecule 1: hypothetical protein PA0115



• Molecule 1: hypothetical protein PA0115



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	138.40Å 138.40Å 136.49Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.89 – 2.35 29.89 – 2.35	Depositor EDS
% Data completeness (in resolution range)	91.9 (29.89-2.35) 91.9 (29.89-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	93.19 (at 2.36Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.232 , 0.282 0.231 , 0.281	Depositor DCC
R_{free} test set	4765 reflections (8.20%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9702	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.14% 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	0.50	0/1220	0.98	7/1651 (0.4%)
1	B	0.52	0/1215	1.06	11/1643 (0.7%)
1	C	0.51	0/1220	1.03	10/1651 (0.6%)
1	D	0.46	0/1214	0.98	6/1643 (0.4%)
1	E	0.46	0/1195	1.00	8/1614 (0.5%)
1	F	0.49	0/1196	0.96	9/1618 (0.6%)
1	G	0.43	0/1208	0.96	7/1636 (0.4%)
1	H	0.43	0/1200	1.00	7/1623 (0.4%)
All	All	0.48	0/9668	1.00	65/13079 (0.5%)

There are no bond length outliers.

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	68	ASP	CA-C-N	8.74	128.33	119.24
1	G	68	ASP	C-N-CA	8.74	128.33	119.24
1	E	68	ASP	CA-C-N	8.25	128.72	119.32
1	E	68	ASP	C-N-CA	8.25	128.72	119.32
1	H	9	HIS	N-CA-C	-7.50	98.71	109.96
1	B	9	HIS	N-CA-C	-7.39	99.35	110.28
1	E	9	HIS	N-CA-C	-7.31	98.14	109.76
1	D	44	ASP	N-CA-C	7.31	120.68	111.69
1	C	31	VAL	N-CA-C	7.18	117.27	110.74
1	B	91	GLY	N-CA-C	-7.14	105.52	114.16
1	C	135	GLU	N-CA-C	-6.99	101.51	110.53
1	B	69	PRO	N-CA-C	6.90	122.61	113.57
1	H	107	LEU	N-CA-C	6.82	118.37	111.07
1	D	122	GLN	N-CA-C	-6.75	103.85	111.14
1	A	41	ASP	N-CA-C	6.62	121.03	113.15
1	A	135	GLU	N-CA-C	-6.60	101.02	110.59
1	E	68	ASP	N-CA-C	6.47	118.05	109.83
1	A	9	HIS	N-CA-C	-6.42	101.45	110.50
1	C	68	ASP	CA-C-N	6.35	126.83	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	68	ASP	C-N-CA	6.35	126.83	119.47
1	C	41	ASP	N-CA-C	6.34	120.87	113.20
1	H	135	GLU	N-CA-C	-6.33	101.41	110.59
1	A	120	HIS	N-CA-C	6.25	117.89	111.14
1	F	93	GLY	N-CA-C	-6.21	105.05	113.37
1	H	73	GLU	N-CA-C	6.16	120.14	111.52
1	C	9	HIS	N-CA-C	-6.14	101.19	110.28
1	B	118	GLN	N-CA-C	-6.13	100.76	109.96
1	F	76	VAL	N-CA-C	-6.06	100.97	109.21
1	F	68	ASP	CA-C-N	6.05	125.94	119.28
1	F	68	ASP	C-N-CA	6.05	125.94	119.28
1	F	135	GLU	N-CA-C	-5.97	102.83	110.53
1	D	82	VAL	N-CA-C	5.84	115.94	107.88
1	D	128	TYR	N-CA-C	-5.83	104.72	112.94
1	C	28	VAL	N-CA-C	5.73	117.12	111.67
1	D	76	VAL	N-CA-C	-5.68	101.40	108.89
1	B	73	GLU	N-CA-C	5.67	119.02	111.30
1	H	76	VAL	N-CA-C	-5.57	100.56	108.58
1	B	44	ASP	N-CA-C	5.55	118.52	111.69
1	B	76	VAL	N-CA-C	-5.52	101.60	108.89
1	F	41	ASP	N-CA-C	5.50	119.86	113.20
1	E	135	GLU	N-CA-C	-5.47	103.47	110.53
1	B	135	GLU	N-CA-C	-5.42	103.54	110.53
1	C	76	VAL	N-CA-C	-5.38	101.79	108.89
1	A	68	ASP	CA-C-N	5.37	125.01	119.05
1	A	68	ASP	C-N-CA	5.37	125.01	119.05
1	E	44	ASP	N-CA-C	5.36	118.98	112.23
1	C	118	GLN	N-CA-C	-5.34	102.14	110.10
1	G	142	ILE	CA-C-N	5.34	126.17	119.98
1	G	142	ILE	C-N-CA	5.34	126.17	119.98
1	G	9	HIS	N-CA-C	-5.33	102.98	110.50
1	E	101	LEU	N-CA-C	-5.32	105.48	111.28
1	F	82	VAL	N-CA-C	5.32	116.02	108.42
1	A	76	VAL	N-CA-C	-5.31	101.88	108.89
1	H	126	GLY	N-CA-C	-5.28	107.77	114.16
1	C	69	PRO	N-CA-C	5.26	121.44	113.81
1	F	118	GLN	N-CA-C	-5.22	102.56	110.28
1	G	44	ASP	N-CA-C	5.21	118.79	112.23
1	B	8	LYS	N-CA-C	5.20	117.97	109.59
1	E	8	LYS	N-CA-C	5.17	118.16	109.95
1	B	55	TRP	N-CA-C	5.13	117.86	109.59
1	F	69	PRO	N-CA-C	5.10	120.53	113.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	14	THR	N-CA-C	-5.08	102.94	110.46
1	H	72	HIS	N-CA-C	5.03	116.78	108.49
1	D	41	ASP	N-CA-C	5.00	119.36	113.16
1	G	76	VAL	N-CA-C	-5.00	102.28	108.89

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	1196	0	1178	32	0
1	B	1192	0	1174	34	0
1	C	1196	0	1178	47	0
1	D	1190	0	1173	46	0
1	E	1172	0	1159	43	0
1	F	1173	0	1158	48	0
1	G	1184	0	1168	59	0
1	H	1177	0	1161	59	0
2	A	47	0	0	2	0
2	B	37	0	0	0	0
2	C	39	0	0	3	0
2	D	17	0	0	1	0
2	E	27	0	0	0	0
2	F	34	0	0	4	0
2	G	14	0	0	2	0
2	H	7	0	0	0	0
All	All	9702	0	9349	351	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 19.

All (351) 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:H:16:LYS:H	1:H:16:LYS:HD2	1.00	1.13
1:F:64:LEU:HD13	1:F:78:ILE:HG23	1.47	0.97
1:D:64:LEU:HD21	1:D:100:ALA:HB2	1.49	0.94
1:A:64:LEU:HD21	1:A:100:ALA:HB2	1.50	0.92
1:E:64:LEU:HD21	1:E:100:ALA:HB2	1.49	0.91
1:H:16:LYS:H	1:H:16:LYS:CD	1.84	0.89
1:H:16:LYS:HD2	1:H:16:LYS:N	1.85	0.88
1:F:142:ILE:HD12	1:F:143:PRO:HD2	1.57	0.86
1:D:97:MSE:HE3	1:D:97:MSE:HA	1.59	0.84
1:H:28:VAL:HA	1:H:32:GLU:HG2	1.59	0.84
1:B:97:MSE:HE3	1:B:97:MSE:HA	1.60	0.83
1:H:59:GLN:HE21	1:H:59:GLN:HA	1.42	0.83
1:C:139:GLU:O	1:C:142:ILE:HG22	1.81	0.79
1:G:3:LEU:HD23	1:G:56:ARG:HD3	1.64	0.79
1:B:133:VAL:HG11	1:B:148:ARG:HG2	1.66	0.77
1:G:28:VAL:HA	1:G:32:GLU:HG2	1.66	0.76
1:C:106:ARG:HD3	2:C:172:HOH:O	1.86	0.76
1:D:28:VAL:HA	1:D:32:GLU:HG2	1.69	0.75
1:H:68:ASP:OD2	1:H:70:VAL:HG22	1.85	0.75
1:E:52:LEU:HD23	1:E:64:LEU:HG	1.69	0.75
1:G:133:VAL:HG11	1:G:148:ARG:HG2	1.69	0.75
1:H:142:ILE:HD12	1:H:143:PRO:HD2	1.68	0.74
1:B:64:LEU:HD21	1:B:100:ALA:HB2	1.70	0.73
1:G:130:PHE:HB3	1:G:147:MSE:HE2	1.70	0.73
1:G:140:ASP:HB2	2:G:156:HOH:O	1.87	0.73
1:H:59:GLN:HE21	1:H:59:GLN:CA	2.02	0.73
1:D:46:VAL:O	1:D:49:THR:HG23	1.89	0.72
1:C:121:LEU:HD13	1:C:125:TYR:HE2	1.54	0.72
1:C:142:ILE:HG23	1:C:144:HIS:HE1	1.54	0.72
1:G:64:LEU:HD21	1:G:100:ALA:HB2	1.70	0.71
1:C:118:GLN:HB2	1:C:121:LEU:HD11	1.73	0.70
1:A:118:GLN:HB2	1:A:121:LEU:HD11	1.73	0.70
1:C:28:VAL:HA	1:C:32:GLU:HG2	1.74	0.70
1:H:76:VAL:CG1	1:H:113:VAL:HG22	2.22	0.69
1:G:106:ARG:HD3	2:G:164:HOH:O	1.93	0.69
1:H:121:LEU:HD13	1:H:125:TYR:HE2	1.58	0.69
1:C:142:ILE:HG23	1:C:144:HIS:CE1	2.28	0.68
1:G:142:ILE:HD13	1:G:142:ILE:C	2.18	0.68
1:C:64:LEU:HD21	1:C:100:ALA:HB2	1.73	0.68
1:H:76:VAL:HG13	1:H:113:VAL:HA	1.76	0.68
1:E:133:VAL:HG11	1:E:148:ARG:HG2	1.76	0.67
1:F:84:SER:O	1:F:88:ARG:HG3	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:ASP:OD2	1:D:70:VAL:HG22	1.95	0.66
1:F:133:VAL:HG11	1:F:148:ARG:HG2	1.76	0.66
1:G:64:LEU:HD12	1:G:64:LEU:C	2.21	0.66
1:A:76:VAL:CG2	1:A:113:VAL:HG22	2.27	0.65
1:F:112:PRO:HB3	1:F:150:ALA:HB2	1.78	0.65
1:H:56:ARG:HG3	1:H:56:ARG:HH11	1.60	0.65
1:B:118:GLN:O	1:B:121:LEU:HD12	1.96	0.65
1:F:64:LEU:HD12	1:F:64:LEU:C	2.22	0.65
1:H:59:GLN:HA	1:H:59:GLN:NE2	2.12	0.65
1:A:28:VAL:HA	1:A:32:GLU:HG2	1.79	0.65
1:H:76:VAL:HG11	1:H:113:VAL:HG22	1.78	0.65
1:E:121:LEU:C	1:E:123:ALA:H	2.03	0.64
1:H:121:LEU:HD13	1:H:125:TYR:CE2	2.32	0.64
1:E:46:VAL:O	1:E:49:THR:HG23	1.98	0.64
1:C:133:VAL:HG11	1:C:148:ARG:HG2	1.80	0.64
1:G:25:ARG:HD2	1:G:63:TYR:HE2	1.61	0.64
1:F:97:MSE:O	1:F:101:LEU:HD13	1.98	0.64
1:H:101:LEU:HG	1:H:149:ARG:HH11	1.63	0.64
1:D:64:LEU:HD23	1:D:96:LEU:HD22	1.81	0.63
1:D:135:GLU:OE1	1:H:111:THR:HA	1.98	0.62
1:G:142:ILE:HD12	1:G:144:HIS:CE1	2.34	0.62
1:H:59:GLN:HE21	1:H:60:LEU:H	1.48	0.62
1:C:64:LEU:C	1:C:64:LEU:HD12	2.25	0.62
1:F:46:VAL:O	1:F:49:THR:HG23	2.00	0.62
1:F:52:LEU:HD13	1:F:96:LEU:HD23	1.82	0.62
1:D:137:TYR:CG	1:H:73:GLU:HA	2.36	0.61
1:B:24:LEU:HD22	1:B:60:LEU:HD13	1.82	0.61
1:A:78:ILE:HD12	1:A:113:VAL:HG11	1.82	0.60
1:C:76:VAL:HG13	1:C:113:VAL:HG22	1.83	0.60
1:B:64:LEU:HD12	1:B:64:LEU:C	2.26	0.60
1:E:76:VAL:CG1	1:E:113:VAL:HG22	2.32	0.60
1:D:21:LEU:HB3	1:D:53:MSE:HE3	1.82	0.60
1:G:76:VAL:CG1	1:G:113:VAL:HG22	2.31	0.60
1:G:46:VAL:O	1:G:49:THR:HG23	2.02	0.60
1:C:43:LEU:O	1:C:49:THR:HG21	2.01	0.59
1:A:64:LEU:C	1:A:64:LEU:HD12	2.27	0.59
1:D:97:MSE:HA	1:D:97:MSE:CE	2.32	0.59
1:D:133:VAL:HG11	1:D:148:ARG:HH11	1.68	0.59
1:G:76:VAL:HG13	1:G:113:VAL:HA	1.84	0.59
1:B:43:LEU:O	1:B:49:THR:HG21	2.02	0.59
1:H:24:LEU:HD22	1:H:60:LEU:HD12	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:MSE:O	1:B:101:LEU:HD13	2.03	0.59
1:H:139:GLU:O	1:H:142:ILE:HG22	2.03	0.59
1:G:76:VAL:HG13	1:G:113:VAL:HG22	1.84	0.58
1:H:56:ARG:HG3	1:H:56:ARG:NH1	2.18	0.58
1:F:68:ASP:OD2	1:F:70:VAL:HG12	2.02	0.58
1:F:115:LEU:HD21	1:F:117:ALA:HB2	1.86	0.58
1:A:133:VAL:HG11	1:A:148:ARG:HG2	1.85	0.58
1:D:67:LEU:HB2	1:D:77:VAL:HB	1.84	0.58
1:G:23:GLN:HB2	1:H:40:VAL:HG11	1.86	0.58
1:E:15:LEU:HD23	1:F:15:LEU:HD23	1.85	0.58
1:G:139:GLU:O	1:G:140:ASP:HB2	2.02	0.58
1:D:14:THR:OG1	1:D:16:LYS:HD2	2.04	0.58
1:C:118:GLN:O	1:C:121:LEU:HD12	2.04	0.58
1:G:25:ARG:HD2	1:G:63:TYR:CE2	2.38	0.58
1:H:101:LEU:HG	1:H:149:ARG:NH1	2.18	0.57
1:A:118:GLN:O	1:A:121:LEU:HD12	2.04	0.57
1:E:64:LEU:C	1:E:64:LEU:HD12	2.29	0.57
1:F:54:ALA:HB3	1:F:62:ALA:HB3	1.87	0.57
1:G:138:LEU:HD12	1:G:142:ILE:N	2.19	0.57
1:G:142:ILE:HD13	1:G:143:PRO:N	2.19	0.57
1:E:76:VAL:HG12	1:E:112:PRO:O	2.05	0.57
1:B:21:LEU:HB3	1:B:53:MSE:HE3	1.87	0.57
1:C:71:ARG:NH2	2:C:178:HOH:O	2.37	0.57
1:C:78:ILE:HD13	1:C:97:MSE:HE1	1.86	0.57
1:F:64:LEU:CD1	1:F:78:ILE:HG23	2.30	0.57
1:G:142:ILE:HG23	1:G:144:HIS:CE1	2.39	0.57
1:G:28:VAL:HG21	1:G:82:VAL:CG2	2.35	0.56
1:C:87:ALA:HB1	1:C:92:LEU:HD11	1.86	0.56
1:F:85:SER:HA	1:F:88:ARG:HE	1.69	0.56
1:G:64:LEU:HD21	1:G:100:ALA:CB	2.36	0.56
1:G:56:ARG:HG2	1:G:56:ARG:NH1	2.20	0.56
1:G:121:LEU:HD22	1:G:124:TYR:HD1	1.69	0.56
1:F:136:VAL:HG12	2:F:175:HOH:O	2.04	0.56
1:H:59:GLN:NE2	1:H:60:LEU:H	2.03	0.56
1:E:123:ALA:O	1:E:127:ARG:HG2	2.06	0.56
1:F:142:ILE:HD12	1:F:143:PRO:CD	2.33	0.56
1:G:56:ARG:HG2	1:G:56:ARG:HH11	1.70	0.56
1:E:76:VAL:HG13	1:E:113:VAL:HA	1.87	0.56
1:F:26:THR:O	1:F:30:VAL:HB	2.05	0.56
1:F:123:ALA:HB3	2:F:184:HOH:O	2.07	0.55
1:C:67:LEU:HB3	1:C:72:HIS:CD2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:121:LEU:HD22	1:G:124:TYR:CD1	2.42	0.55
1:A:76:VAL:HG22	1:A:113:VAL:HG22	1.87	0.55
1:C:128:TYR:O	1:C:149:ARG:HD2	2.07	0.55
1:D:3:LEU:HD11	1:D:92:LEU:HD21	1.90	0.54
1:D:66:LEU:CD2	1:D:78:ILE:HG12	2.37	0.54
1:D:5:TRP:O	1:D:6:THR:HG23	2.06	0.54
1:H:76:VAL:HG13	1:H:113:VAL:HG22	1.90	0.54
1:F:85:SER:OG	1:F:88:ARG:NH1	2.39	0.54
1:C:121:LEU:HD13	1:C:125:TYR:CE2	2.39	0.54
1:E:13:LEU:HD11	1:E:51:HIS:CD2	2.42	0.54
1:H:59:GLN:HE21	1:H:60:LEU:N	2.06	0.54
1:G:123:ALA:O	1:G:127:ARG:HG2	2.07	0.53
1:F:64:LEU:HD21	1:F:100:ALA:CB	2.38	0.53
1:A:76:VAL:HG22	1:A:113:VAL:HA	1.90	0.53
1:B:46:VAL:O	1:B:49:THR:HG23	2.09	0.53
1:B:82:VAL:HG22	1:B:83:SER:N	2.23	0.53
1:D:16:LYS:HD2	1:D:17:GLU:H	1.75	0.52
1:H:53:MSE:HB3	1:H:60:LEU:HD21	1.90	0.52
1:D:16:LYS:HD2	1:D:17:GLU:N	2.24	0.52
1:E:21:LEU:HB3	1:E:53:MSE:HE3	1.90	0.52
1:G:26:THR:O	1:G:30:VAL:HB	2.09	0.52
1:H:13:LEU:HD11	1:H:51:HIS:CD2	2.44	0.52
1:A:118:GLN:HB2	1:A:121:LEU:CD1	2.37	0.52
1:D:133:VAL:HG11	1:D:148:ARG:NH1	2.24	0.52
1:D:52:LEU:HD23	1:D:64:LEU:HG	1.92	0.52
1:D:135:GLU:OE1	1:H:111:THR:HG23	2.10	0.52
1:F:64:LEU:HD21	1:F:100:ALA:HB2	1.92	0.52
1:C:21:LEU:HB3	1:C:53:MSE:HE3	1.91	0.51
1:C:118:GLN:HB2	1:C:121:LEU:CD1	2.39	0.51
1:H:16:LYS:CD	1:H:16:LYS:N	2.60	0.51
1:E:9:HIS:O	1:E:11:ALA:N	2.44	0.51
1:G:142:ILE:HG23	1:G:144:HIS:HE1	1.74	0.51
1:C:64:LEU:HD21	1:C:100:ALA:CB	2.38	0.51
1:G:21:LEU:C	1:G:21:LEU:HD12	2.35	0.51
1:G:135:GLU:HA	1:G:135:GLU:OE1	2.09	0.51
1:E:9:HIS:C	1:E:11:ALA:N	2.67	0.51
1:H:64:LEU:C	1:H:64:LEU:HD12	2.35	0.51
1:E:121:LEU:C	1:E:123:ALA:N	2.68	0.51
1:D:137:TYR:CD1	1:H:73:GLU:HA	2.46	0.50
1:C:76:VAL:CG1	1:C:113:VAL:HG22	2.41	0.50
1:E:139:GLU:HB3	1:E:144:HIS:CE1	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:115:LEU:HD12	1:E:130:PHE:CE1	2.46	0.50
1:G:68:ASP:HB2	1:G:69:PRO:HD2	1.93	0.50
1:H:115:LEU:C	1:H:115:LEU:HD22	2.37	0.50
1:A:120:HIS:C	1:A:122:GLN:H	2.18	0.50
1:D:64:LEU:HD21	1:D:100:ALA:CB	2.32	0.50
1:F:21:LEU:C	1:F:21:LEU:HD12	2.36	0.50
1:H:64:LEU:HD21	1:H:100:ALA:HB2	1.94	0.50
1:A:67:LEU:HB3	1:A:72:HIS:CD2	2.47	0.50
1:G:119:ALA:O	1:G:122:GLN:HG3	2.12	0.49
1:H:24:LEU:CD2	1:H:60:LEU:HD12	2.42	0.49
1:E:76:VAL:HG11	1:E:113:VAL:HG22	1.95	0.49
1:H:82:VAL:HG22	1:H:83:SER:N	2.27	0.49
1:E:65:ARG:NE	1:E:67:LEU:HD11	2.26	0.49
1:E:97:MSE:HE3	1:E:100:ALA:HB3	1.94	0.49
1:A:76:VAL:HG21	1:A:113:VAL:HG22	1.93	0.49
1:G:142:ILE:C	1:G:142:ILE:CD1	2.86	0.49
1:C:124:TYR:HA	1:C:127:ARG:CD	2.42	0.49
1:F:21:LEU:HD12	1:F:22:LEU:HD23	1.95	0.49
1:C:78:ILE:CD1	1:C:97:MSE:HE1	2.41	0.49
1:D:127:ARG:HD3	1:D:127:ARG:HA	1.59	0.49
1:H:46:VAL:O	1:H:49:THR:HG23	2.13	0.49
1:B:133:VAL:HG23	1:B:134:THR:HG23	1.95	0.48
1:E:123:ALA:HB1	1:E:127:ARG:HH21	1.78	0.48
1:F:133:VAL:CG1	1:F:148:ARG:HG2	2.41	0.48
1:D:119:ALA:C	1:D:121:LEU:H	2.21	0.48
1:B:91:GLY:O	1:B:95:GLN:HG2	2.13	0.48
1:D:130:PHE:HA	1:D:149:ARG:HB3	1.96	0.48
1:C:68:ASP:OD2	1:C:70:VAL:HG22	2.13	0.48
1:H:25:ARG:HD2	1:H:63:TYR:CE2	2.48	0.48
1:D:82:VAL:HG22	1:D:83:SER:N	2.27	0.48
1:C:46:VAL:O	1:C:49:THR:HG23	2.14	0.48
1:D:119:ALA:O	1:D:121:LEU:N	2.45	0.48
1:E:40:VAL:HG12	1:F:19:TYR:CZ	2.49	0.48
1:G:138:LEU:HD22	1:G:138:LEU:N	2.29	0.48
1:C:124:TYR:O	1:C:127:ARG:HD3	2.13	0.48
1:A:121:LEU:HD13	1:A:125:TYR:HE2	1.79	0.48
1:E:76:VAL:HG13	1:E:113:VAL:HG22	1.96	0.47
1:A:67:LEU:HD13	2:A:176:HOH:O	2.13	0.47
1:E:64:LEU:HD21	1:E:100:ALA:CB	2.34	0.47
1:E:82:VAL:HG22	1:E:83:SER:N	2.29	0.47
1:B:67:LEU:HB2	1:B:77:VAL:HB	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:TYR:HA	1:C:127:ARG:HD3	1.97	0.47
1:D:118:GLN:O	1:D:121:LEU:HD12	2.14	0.47
1:A:67:LEU:HB2	1:A:77:VAL:HB	1.95	0.47
1:D:48:ASP:O	1:D:50:HIS:ND1	2.44	0.47
1:F:52:LEU:HD13	1:F:96:LEU:CD2	2.44	0.47
1:G:43:LEU:O	1:G:49:THR:HG21	2.14	0.47
1:G:88:ARG:HG3	1:G:88:ARG:HH11	1.80	0.47
1:G:101:LEU:HD21	1:G:149:ARG:CD	2.45	0.47
1:H:63:TYR:CD2	1:H:64:LEU:N	2.83	0.47
1:F:43:LEU:O	1:F:49:THR:HG21	2.15	0.47
1:F:139:GLU:O	1:F:142:ILE:HG23	2.15	0.47
1:F:67:LEU:HB2	1:F:77:VAL:HB	1.97	0.46
1:H:49:THR:HB	1:H:67:LEU:HD12	1.97	0.46
1:B:49:THR:HB	1:B:67:LEU:HD12	1.97	0.46
1:F:5:TRP:CZ3	1:F:54:ALA:HB2	2.50	0.46
1:F:59:GLN:HA	2:F:167:HOH:O	2.15	0.46
1:H:63:TYR:CG	1:H:64:LEU:N	2.84	0.46
1:B:3:LEU:HD11	1:B:61:LEU:HD12	1.95	0.46
1:B:28:VAL:HA	1:B:32:GLU:HG2	1.96	0.46
1:E:9:HIS:C	1:E:11:ALA:H	2.24	0.46
1:H:120:HIS:CE1	1:H:121:LEU:HG	2.51	0.46
1:F:143:PRO:HG2	2:F:164:HOH:O	2.15	0.46
1:C:70:VAL:HG13	2:C:169:HOH:O	2.16	0.46
2:A:182:HOH:O	1:C:70:VAL:HG12	2.15	0.46
1:C:64:LEU:C	1:C:64:LEU:CD1	2.88	0.46
1:C:123:ALA:O	1:C:127:ARG:HD2	2.14	0.46
1:B:44:ASP:HB2	1:C:19:TYR:CE2	2.50	0.46
1:E:13:LEU:HA	1:E:17:GLU:OE2	2.15	0.46
1:F:25:ARG:HD2	1:F:63:TYR:CE2	2.51	0.46
1:G:21:LEU:HD12	1:G:22:LEU:N	2.31	0.46
1:B:24:LEU:CD2	1:B:60:LEU:HD13	2.46	0.45
1:E:68:ASP:OD2	1:E:70:VAL:HG13	2.17	0.45
1:G:28:VAL:HA	1:G:32:GLU:CG	2.42	0.45
1:H:71:ARG:HG2	1:H:71:ARG:HH21	1.81	0.45
1:H:102:GLN:O	1:H:105:GLU:HB3	2.16	0.45
1:C:3:LEU:HG	1:C:56:ARG:NH2	2.31	0.45
1:A:19:TYR:CE2	1:D:44:ASP:HB2	2.51	0.45
1:B:121:LEU:HD13	1:B:125:TYR:CE2	2.52	0.45
1:C:82:VAL:HG22	1:C:83:SER:N	2.31	0.45
1:B:87:ALA:O	1:B:92:LEU:HD12	2.17	0.45
1:G:132:ALA:HA	1:G:147:MSE:HG2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:2:SER:O	1:E:56:ARG:HD3	2.16	0.45
1:H:3:LEU:HD11	1:H:92:LEU:HD21	1.98	0.45
1:B:109:LEU:O	1:B:110:ASP:HB2	2.17	0.45
1:A:115:LEU:HD12	1:A:115:LEU:C	2.41	0.45
1:D:18:LEU:O	1:D:21:LEU:HG	2.17	0.45
1:D:121:LEU:HD13	1:D:125:TYR:CE2	2.52	0.45
1:F:9:HIS:O	1:F:11:ALA:N	2.50	0.45
1:E:3:LEU:CD2	1:E:56:ARG:HB2	2.46	0.45
1:A:101:LEU:HD12	1:A:101:LEU:HA	1.73	0.45
1:B:63:TYR:CG	1:B:64:LEU:N	2.85	0.45
1:D:27:GLU:O	1:D:31:VAL:HB	2.16	0.45
1:D:47:GLY:HA3	1:D:71:ARG:HD2	1.99	0.45
1:F:64:LEU:HD12	1:F:65:ARG:N	2.31	0.44
1:F:21:LEU:HD12	1:F:22:LEU:N	2.32	0.44
1:D:14:THR:CB	1:D:16:LYS:HZ3	2.30	0.44
1:H:55:TRP:CE2	1:H:60:LEU:HD23	2.52	0.44
1:A:118:GLN:CB	1:A:121:LEU:HD11	2.46	0.44
1:A:138:LEU:N	1:A:138:LEU:HD22	2.32	0.44
1:F:44:ASP:O	1:F:45:LEU:HD23	2.17	0.44
1:E:134:THR:HG22	1:G:74:GLY:HA2	2.00	0.44
1:G:34:LYS:HE2	1:G:34:LYS:HB2	1.77	0.44
1:A:5:TRP:CZ3	1:A:54:ALA:HB2	2.53	0.44
1:C:72:HIS:C	1:C:73:GLU:OE1	2.60	0.44
1:E:99:ARG:O	1:E:102:GLN:HB3	2.18	0.44
1:E:122:GLN:HG3	1:E:147:MSE:HE3	1.98	0.44
1:G:130:PHE:CB	1:G:147:MSE:HE2	2.44	0.44
1:F:64:LEU:C	1:F:64:LEU:CD1	2.91	0.44
1:B:25:ARG:HD2	1:B:63:TYR:CE2	2.53	0.44
1:D:73:GLU:O	1:D:74:GLY:C	2.61	0.43
1:D:136:VAL:HG23	2:D:163:HOH:O	2.18	0.43
1:B:97:MSE:HA	1:B:97:MSE:CE	2.39	0.43
1:D:26:THR:O	1:D:30:VAL:HB	2.19	0.43
1:E:13:LEU:HD22	1:E:17:GLU:HB3	2.00	0.43
1:E:44:ASP:O	1:E:45:LEU:HD23	2.18	0.43
1:E:67:LEU:HB2	1:E:77:VAL:HB	2.00	0.43
1:C:67:LEU:HB2	1:C:77:VAL:HB	2.00	0.43
1:D:118:GLN:O	1:D:119:ALA:C	2.61	0.43
1:B:67:LEU:HB3	1:B:72:HIS:CD2	2.53	0.43
1:C:54:ALA:HB3	1:C:62:ALA:HB3	2.01	0.43
1:E:43:LEU:O	1:E:49:THR:HG21	2.19	0.43
1:H:68:ASP:HB2	1:H:69:PRO:HD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:3:LEU:HD23	1:E:56:ARG:HB2	2.01	0.43
1:E:125:TYR:O	1:E:126:GLY:C	2.62	0.43
1:B:115:LEU:CD1	1:B:147:MSE:HB2	2.49	0.43
1:G:76:VAL:HG11	1:G:113:VAL:HG22	2.00	0.43
1:A:108:TRP:O	1:A:109:LEU:C	2.62	0.42
1:G:3:LEU:CD2	1:G:56:ARG:HD3	2.43	0.42
1:H:118:GLN:O	1:H:121:LEU:HD12	2.19	0.42
1:B:43:LEU:HD13	1:B:67:LEU:HD21	2.01	0.42
1:H:49:THR:HA	1:H:66:LEU:O	2.20	0.42
1:B:19:TYR:CE2	1:C:44:ASP:HB2	2.54	0.42
1:C:115:LEU:HD12	1:C:115:LEU:O	2.19	0.42
1:G:21:LEU:HD23	1:G:53:MSE:HE1	2.01	0.42
1:G:23:GLN:HB2	1:H:40:VAL:CG1	2.49	0.42
1:G:51:HIS:HB3	1:G:53:MSE:CE	2.50	0.42
1:A:73:GLU:HA	1:C:114:TYR:CZ	2.55	0.42
1:C:142:ILE:CG2	1:C:144:HIS:HE1	2.27	0.42
1:D:26:THR:HA	1:D:30:VAL:HG23	2.02	0.42
1:A:126:GLY:HA2	1:A:130:PHE:O	2.19	0.42
1:B:139:GLU:O	1:B:142:ILE:HG22	2.19	0.42
1:G:53:MSE:HB2	1:G:60:LEU:CD2	2.50	0.42
1:G:24:LEU:HD22	1:G:60:LEU:CD1	2.50	0.42
1:H:138:LEU:N	1:H:138:LEU:CD1	2.83	0.42
1:B:5:TRP:C	1:B:6:THR:HG22	2.45	0.42
1:C:139:GLU:O	1:C:142:ILE:CG2	2.62	0.42
1:F:112:PRO:CB	1:F:150:ALA:HB2	2.48	0.42
1:H:43:LEU:O	1:H:49:THR:HG21	2.19	0.41
1:H:114:TYR:HD2	1:H:148:ARG:HB3	1.85	0.41
1:B:51:HIS:HB3	1:B:63:TYR:HE1	1.85	0.41
1:D:14:THR:HB	1:D:16:LYS:HZ3	1.85	0.41
1:H:82:VAL:CG2	1:H:83:SER:N	2.84	0.41
1:H:139:GLU:O	1:H:140:ASP:C	2.64	0.41
1:C:51:HIS:HB3	1:C:63:TYR:HE1	1.86	0.41
1:G:13:LEU:HA	1:G:17:GLU:OE2	2.19	0.41
1:A:74:GLY:HA2	1:C:134:THR:HG22	2.02	0.41
1:A:119:ALA:HB2	1:A:143:PRO:HB2	2.02	0.41
1:D:122:GLN:O	1:D:123:ALA:C	2.63	0.41
1:F:115:LEU:HD12	1:F:130:PHE:CD1	2.55	0.41
1:E:63:TYR:CG	1:E:64:LEU:N	2.88	0.41
1:F:115:LEU:HD22	1:F:117:ALA:N	2.35	0.41
1:G:121:LEU:O	1:G:124:TYR:HB3	2.20	0.41
1:A:6:THR:HG23	1:A:53:MSE:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:LEU:HD12	1:C:115:LEU:C	2.46	0.41
1:G:115:LEU:HD13	1:G:117:ALA:HB2	2.02	0.41
1:B:82:VAL:CG2	1:B:83:SER:N	2.83	0.41
1:D:73:GLU:HA	1:H:137:TYR:CG	2.56	0.41
1:F:18:LEU:O	1:F:21:LEU:HG	2.21	0.41
1:H:85:SER:C	1:H:87:ALA:H	2.29	0.41
1:H:92:LEU:O	1:H:93:GLY:C	2.64	0.41
1:D:78:ILE:HD13	1:D:97:MSE:CE	2.51	0.41
1:F:139:GLU:O	1:F:142:ILE:CG2	2.69	0.41
1:G:56:ARG:HH11	1:G:56:ARG:CG	2.34	0.41
1:G:138:LEU:HD12	1:G:142:ILE:H	1.83	0.41
1:A:15:LEU:HD23	1:D:15:LEU:HD23	2.04	0.40
1:A:14:THR:HB	1:A:16:LYS:HE2	2.02	0.40
1:A:67:LEU:N	1:A:67:LEU:CD1	2.85	0.40
1:G:103:ALA:HA	1:G:106:ARG:HG2	2.02	0.40
1:B:56:ARG:O	1:B:57:ASP:C	2.64	0.40
1:F:117:ALA:HB1	1:F:125:TYR:CE2	2.56	0.40
1:G:52:LEU:O	1:G:63:TYR:HA	2.21	0.40
1:E:50:HIS:CD2	1:E:107:LEU:HD11	2.56	0.40
1:E:19:TYR:CZ	1:F:40:VAL:HG12	2.57	0.40
1:F:63:TYR:CG	1:F:64:LEU:N	2.90	0.40
1:F:117:ALA:HB1	1:F:125:TYR:CD2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	147/150 (98%)	137 (93%)	10 (7%)	0	100	100
1	B	144/150 (96%)	135 (94%)	9 (6%)	0	100	100
1	C	147/150 (98%)	138 (94%)	9 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	146/150 (97%)	131 (90%)	13 (9%)	2 (1%)	9	7
1	E	142/150 (95%)	130 (92%)	11 (8%)	1 (1%)	18	20
1	F	141/150 (94%)	130 (92%)	10 (7%)	1 (1%)	18	20
1	G	145/150 (97%)	130 (90%)	14 (10%)	1 (1%)	18	20
1	H	142/150 (95%)	129 (91%)	13 (9%)	0	100	100
All	All	1154/1200 (96%)	1060 (92%)	89 (8%)	5 (0%)	30	34

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	120	HIS
1	D	31	VAL
1	E	10	HIS
1	F	10	HIS
1	G	119	ALA

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	124/122 (102%)	110 (89%)	14 (11%)	5	5
1	B	124/122 (102%)	110 (89%)	14 (11%)	5	5
1	C	124/122 (102%)	114 (92%)	10 (8%)	11	12
1	D	123/122 (101%)	109 (89%)	14 (11%)	5	5
1	E	121/122 (99%)	110 (91%)	11 (9%)	9	9
1	F	122/122 (100%)	111 (91%)	11 (9%)	9	9
1	G	123/122 (101%)	115 (94%)	8 (6%)	15	18
1	H	122/122 (100%)	105 (86%)	17 (14%)	3	3
All	All	983/976 (101%)	884 (90%)	99 (10%)	7	6

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	18	LEU
1	A	23	GLN
1	A	60	LEU
1	A	64	LEU
1	A	70	VAL
1	A	71	ARG
1	A	76	VAL
1	A	101	LEU
1	A	105	GLU
1	A	121	LEU
1	A	136	VAL
1	A	141	ASP
1	A	148	ARG
1	B	6	THR
1	B	18	LEU
1	B	23	GLN
1	B	49	THR
1	B	60	LEU
1	B	64	LEU
1	B	67	LEU
1	B	70	VAL
1	B	76	VAL
1	B	92	LEU
1	B	115	LEU
1	B	121	LEU
1	B	136	VAL
1	B	148	ARG
1	C	16	LYS
1	C	18	LEU
1	C	49	THR
1	C	60	LEU
1	C	64	LEU
1	C	67	LEU
1	C	76	VAL
1	C	101	LEU
1	C	121	LEU
1	C	138	LEU
1	D	3	LEU
1	D	16	LYS
1	D	18	LEU
1	D	23	GLN
1	D	49	THR

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Mol	Chain	Res	Type
1	D	60	LEU
1	D	64	LEU
1	D	75	GLN
1	D	101	LEU
1	D	102	GLN
1	D	135	GLU
1	D	138	LEU
1	D	141	ASP
1	D	149	ARG
1	E	18	LEU
1	E	40	VAL
1	E	57	ASP
1	E	60	LEU
1	E	64	LEU
1	E	70	VAL
1	E	76	VAL
1	E	92	LEU
1	E	101	LEU
1	E	115	LEU
1	E	148	ARG
1	F	18	LEU
1	F	40	VAL
1	F	52	LEU
1	F	60	LEU
1	F	64	LEU
1	F	71	ARG
1	F	92	LEU
1	F	115	LEU
1	F	136	VAL
1	F	142	ILE
1	F	148	ARG
1	G	18	LEU
1	G	23	GLN
1	G	49	THR
1	G	64	LEU
1	G	67	LEU
1	G	70	VAL
1	G	76	VAL
1	G	142	ILE
1	H	16	LYS
1	H	18	LEU
1	H	21	LEU

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Mol	Chain	Res	Type
1	H	49	THR
1	H	59	GLN
1	H	64	LEU
1	H	67	LEU
1	H	73	GLU
1	H	76	VAL
1	H	85	SER
1	H	101	LEU
1	H	115	LEU
1	H	121	LEU
1	H	135	GLU
1	H	138	LEU
1	H	141	ASP
1	H	148	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	122	GLN
1	B	23	GLN
1	B	75	GLN
1	B	122	GLN
1	C	95	GLN
1	C	122	GLN
1	D	90	GLN
1	D	102	GLN
1	F	94	HIS
1	G	23	GLN
1	G	59	GLN
1	G	144	HIS
1	H	33	GLN
1	H	59	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	146/150 (97%)	-0.15	2 (1%) 73 78	18, 29, 46, 52	0
1	B	145/150 (96%)	-0.06	4 (2%) 55 61	17, 30, 47, 56	0
1	C	146/150 (97%)	-0.11	1 (0%) 84 87	15, 29, 46, 61	0
1	D	145/150 (96%)	0.20	4 (2%) 55 61	22, 36, 50, 63	0
1	E	143/150 (95%)	0.44	10 (6%) 22 25	21, 39, 58, 72	0
1	F	142/150 (94%)	0.28	7 (4%) 35 41	21, 34, 56, 64	0
1	G	144/150 (96%)	0.71	10 (6%) 23 26	25, 48, 67, 79	0
1	H	143/150 (95%)	0.71	15 (10%) 11 13	30, 45, 65, 72	0
All	All	1154/1200 (96%)	0.25	53 (4%) 37 43	15, 36, 59, 79	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	150	ALA	4.5
1	E	143	PRO	4.2
1	B	140	ASP	3.9
1	F	92	LEU	3.6
1	E	2	SER	3.5
1	H	121	LEU	3.4
1	E	121	LEU	3.2
1	D	150	ALA	3.2
1	F	127	ARG	3.2
1	A	150	ALA	3.2
1	G	56	ARG	3.1
1	F	57	ASP	3.0
1	G	54	ALA	3.0
1	F	87	ALA	2.9
1	H	87	ALA	2.9
1	H	142	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	4	ASP	2.8
1	G	127	ARG	2.8
1	H	91	GLY	2.7
1	A	2	SER	2.7
1	H	119	ALA	2.7
1	H	140	ASP	2.6
1	F	3	LEU	2.6
1	H	123	ALA	2.5
1	B	121	LEU	2.5
1	D	70	VAL	2.5
1	E	123	ALA	2.5
1	G	85	SER	2.4
1	G	138	LEU	2.4
1	H	124	TYR	2.4
1	E	127	ARG	2.4
1	F	88	ARG	2.4
1	B	2	SER	2.4
1	G	57	ASP	2.4
1	H	150	ALA	2.3
1	H	59	GLN	2.3
1	G	142	ILE	2.3
1	G	60	LEU	2.3
1	E	144	HIS	2.2
1	H	72	HIS	2.2
1	H	115	LEU	2.2
1	E	124	TYR	2.1
1	D	121	LEU	2.1
1	C	142	ILE	2.1
1	D	127	ARG	2.1
1	G	86	ALA	2.1
1	G	143	PRO	2.1
1	E	5	TRP	2.1
1	H	58	GLY	2.1
1	H	120	HIS	2.0
1	H	3	LEU	2.0
1	E	120	HIS	2.0
1	E	4	ASP	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 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.