



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:55 PM UTC

PDB ID : 8GQG / pdb_00008gqg
Title : Crystal structure of Thioloase from Pseudomonas aeruginosa PAO1
Authors : Hong, J.; Son, H.F.; Kim, K.J.
Deposited on : 2022-08-30
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

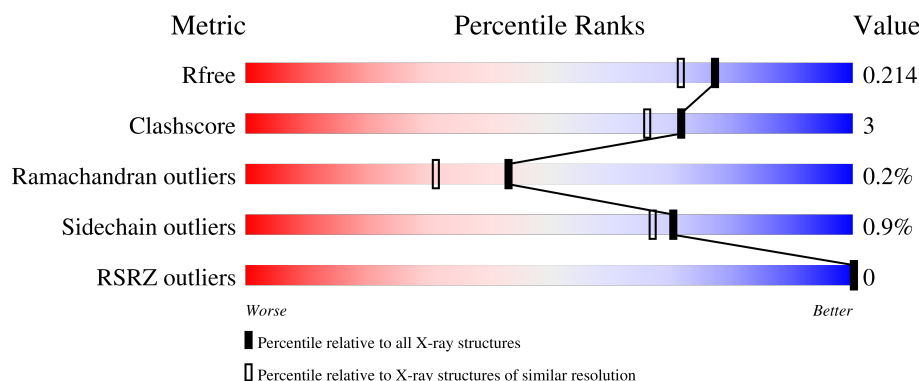
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	402	
1	B	402	
1	C	402	
1	D	402	
1	E	402	

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Mol	Chain	Length	Quality of chain
1	F	402	 90% 7% .

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 19363 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 Thiolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	1	0
			2903	1808	519	556	20			
1	B	394	Total	C	N	O	S	0	1	0
			2953	1836	527	570	20			
1	C	388	Total	C	N	O	S	0	1	0
			2912	1813	520	559	20			
1	D	394	Total	C	N	O	S	0	1	0
			2953	1836	527	570	20			
1	E	388	Total	C	N	O	S	0	1	0
			2912	1813	520	559	20			
1	F	394	Total	C	N	O	S	0	1	0
			2953	1836	527	570	20			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	395	LEU	-	expression tag	UNP Q9HYF2
A	396	GLU	-	expression tag	UNP Q9HYF2
A	397	HIS	-	expression tag	UNP Q9HYF2
A	398	HIS	-	expression tag	UNP Q9HYF2
A	399	HIS	-	expression tag	UNP Q9HYF2
A	400	HIS	-	expression tag	UNP Q9HYF2
A	401	HIS	-	expression tag	UNP Q9HYF2
A	402	HIS	-	expression tag	UNP Q9HYF2
B	395	LEU	-	expression tag	UNP Q9HYF2
B	396	GLU	-	expression tag	UNP Q9HYF2
B	397	HIS	-	expression tag	UNP Q9HYF2
B	398	HIS	-	expression tag	UNP Q9HYF2
B	399	HIS	-	expression tag	UNP Q9HYF2
B	400	HIS	-	expression tag	UNP Q9HYF2
B	401	HIS	-	expression tag	UNP Q9HYF2
B	402	HIS	-	expression tag	UNP Q9HYF2
C	395	LEU	-	expression tag	UNP Q9HYF2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	396	GLU	-	expression tag	UNP Q9HYF2
C	397	HIS	-	expression tag	UNP Q9HYF2
C	398	HIS	-	expression tag	UNP Q9HYF2
C	399	HIS	-	expression tag	UNP Q9HYF2
C	400	HIS	-	expression tag	UNP Q9HYF2
C	401	HIS	-	expression tag	UNP Q9HYF2
C	402	HIS	-	expression tag	UNP Q9HYF2
D	395	LEU	-	expression tag	UNP Q9HYF2
D	396	GLU	-	expression tag	UNP Q9HYF2
D	397	HIS	-	expression tag	UNP Q9HYF2
D	398	HIS	-	expression tag	UNP Q9HYF2
D	399	HIS	-	expression tag	UNP Q9HYF2
D	400	HIS	-	expression tag	UNP Q9HYF2
D	401	HIS	-	expression tag	UNP Q9HYF2
D	402	HIS	-	expression tag	UNP Q9HYF2
E	395	LEU	-	expression tag	UNP Q9HYF2
E	396	GLU	-	expression tag	UNP Q9HYF2
E	397	HIS	-	expression tag	UNP Q9HYF2
E	398	HIS	-	expression tag	UNP Q9HYF2
E	399	HIS	-	expression tag	UNP Q9HYF2
E	400	HIS	-	expression tag	UNP Q9HYF2
E	401	HIS	-	expression tag	UNP Q9HYF2
E	402	HIS	-	expression tag	UNP Q9HYF2
F	395	LEU	-	expression tag	UNP Q9HYF2
F	396	GLU	-	expression tag	UNP Q9HYF2
F	397	HIS	-	expression tag	UNP Q9HYF2
F	398	HIS	-	expression tag	UNP Q9HYF2
F	399	HIS	-	expression tag	UNP Q9HYF2
F	400	HIS	-	expression tag	UNP Q9HYF2
F	401	HIS	-	expression tag	UNP Q9HYF2
F	402	HIS	-	expression tag	UNP Q9HYF2

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	317	Total O 317 317	0	0
2	B	284	Total O 284 284	0	0
2	C	321	Total O 321 321	0	0
2	D	270	Total O 270 270	0	0

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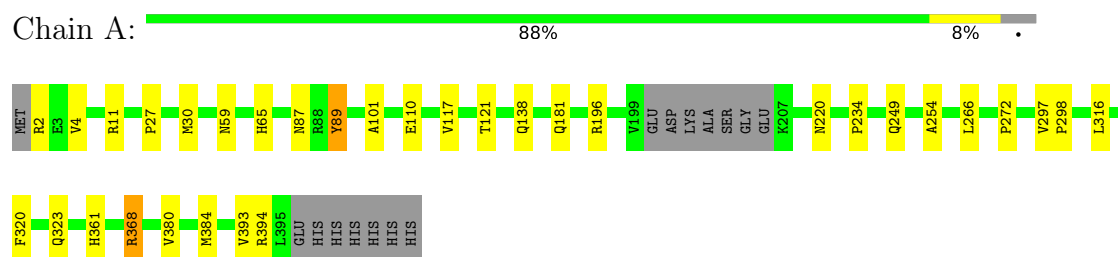
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	312	Total 312	O 312	0	0
2	F	273	Total 273	O 273	0	0

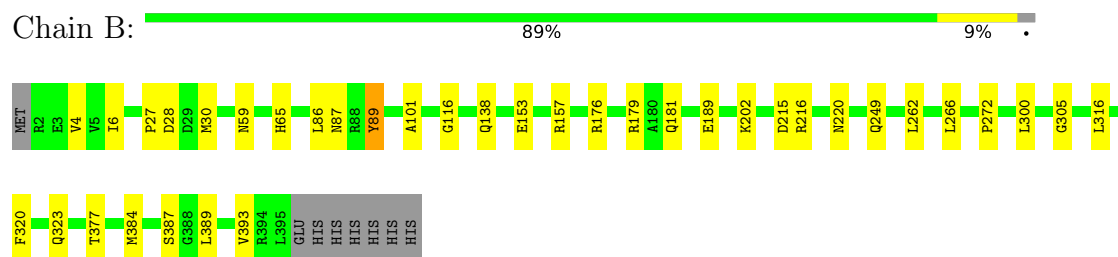
3 Residue-property plots

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.

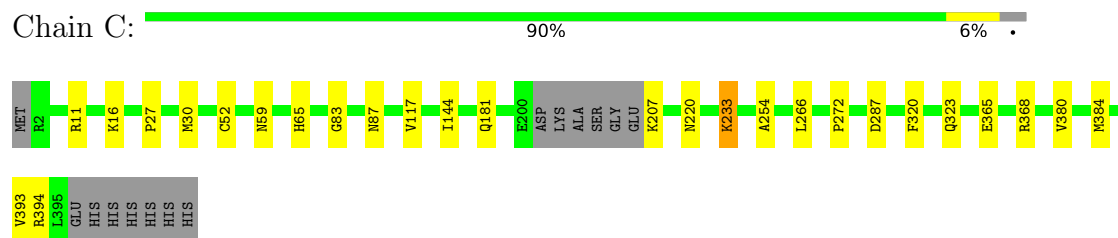
• Molecule 1: Thiolase



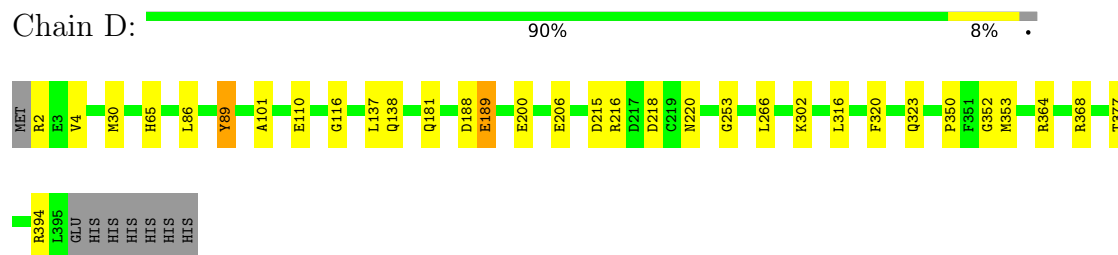
• Molecule 1: Thiolase



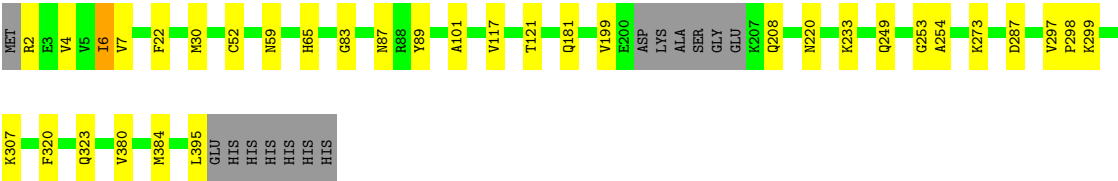
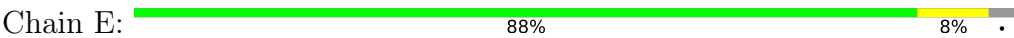
• Molecule 1: Thiolase



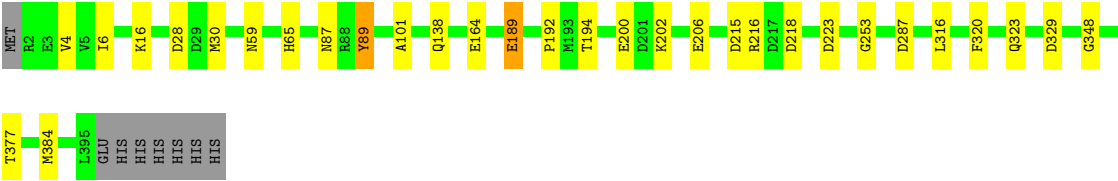
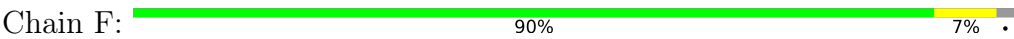
• Molecule 1: Thiolase



• Molecule 1: Thiolase



• Molecule 1: Thiolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.22Å 204.78Å 262.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.65 – 1.80 31.65 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.2 (31.65-1.80) 95.2 (31.65-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.172 , 0.207 0.182 , 0.214	Depositor DCC
R_{free} test set	13813 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.487 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.487 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19363	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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	1.09	3/2944 (0.1%)	1.33	2/3982 (0.1%)
1	B	1.10	1/2995 (0.0%)	1.34	6/4051 (0.1%)
1	C	1.09	1/2953 (0.0%)	1.31	2/3994 (0.1%)
1	D	1.12	0/2995	1.31	1/4051 (0.0%)
1	E	1.09	1/2953 (0.0%)	1.32	1/3994 (0.0%)
1	F	1.11	0/2995	1.35	7/4051 (0.2%)
All	All	1.10	6/17835 (0.0%)	1.33	19/24123 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	121	THR	C-O	6.27	1.31	1.24
1	A	121	THR	C-O	5.17	1.30	1.24
1	A	394	ARG	C-O	5.10	1.29	1.24
1	A	234	PRO	C-O	-5.03	1.17	1.23
1	B	116	GLY	C-O	5.02	1.28	1.23
1	C	394	ARG	C-O	5.01	1.30	1.24

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	384	MET	CA-C-N	6.28	126.92	122.33
1	F	384	MET	C-N-CA	6.28	126.92	122.33
1	F	218	ASP	CA-CB-CG	5.87	118.47	112.60
1	F	194	THR	CA-CB-OG1	-5.58	101.23	109.60
1	B	384	MET	CA-C-N	5.51	126.36	122.33
1	B	384	MET	C-N-CA	5.51	126.36	122.33
1	C	11	ARG	CG-CD-NE	5.34	123.76	112.00
1	F	287	ASP	CA-CB-CG	5.30	117.90	112.60
1	D	218	ASP	CA-CB-CG	5.29	117.89	112.60
1	B	387	SER	CA-C-N	5.26	125.23	121.65
1	B	387	SER	C-N-CA	5.26	125.23	121.65
1	E	287	ASP	CA-CB-CG	5.17	117.77	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	348	GLY	CA-C-O	-5.12	118.48	122.52
1	C	287	ASP	CA-CB-CG	5.11	117.71	112.60
1	A	11	ARG	CG-CD-NE	5.10	123.21	112.00
1	B	249	GLN	CB-CG-CD	-5.06	103.99	112.60
1	B	28	ASP	CA-CB-CG	5.06	117.66	112.60
1	A	361	HIS	CA-CB-CG	-5.00	108.80	113.80
1	F	28	ASP	CA-CB-CG	5.00	117.60	112.60

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	2903	0	2905	22	0
1	B	2953	0	2948	22	0
1	C	2912	0	2911	18	0
1	D	2953	0	2948	25	0
1	E	2912	0	2911	20	0
1	F	2953	0	2948	18	0
2	A	317	0	0	6	0
2	B	284	0	0	2	0
2	C	321	0	0	5	0
2	D	270	0	0	3	0
2	E	312	0	0	4	0
2	F	273	0	0	5	0
All	All	19363	0	17571	115	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 3.

All (115) 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:F:30:MET:HE1	1:F:253:GLY:HA2	1.46	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:30:MET:HE1	1:D:253:GLY:HA2	1.53	0.89
1:B:305:GLY:HA2	2:D:631:HOH:O	1.76	0.85
1:D:189:GLU:HB3	2:D:610:HOH:O	1.83	0.76
1:F:164:GLU:HG2	2:F:756:HOH:O	1.84	0.76
1:A:368:ARG:HD2	2:A:720:HOH:O	1.87	0.74
1:F:30:MET:HE1	1:F:253:GLY:CA	2.17	0.74
1:F:30:MET:CE	1:F:253:GLY:HA2	2.18	0.72
1:E:22:PHE:CE2	1:E:30:MET:HE3	2.25	0.72
1:D:30:MET:HE1	1:D:253:GLY:CA	2.19	0.72
1:D:138:GLN:NE2	2:D:501:HOH:O	2.22	0.71
1:B:215:ASP:OD1	1:B:216:ARG:NH2	2.23	0.70
1:E:181:GLN:HE22	1:E:220:ASN:HD21	1.40	0.69
1:D:30:MET:CE	1:D:253:GLY:HA2	2.23	0.68
1:A:27:PRO:HA	1:A:30:MET:HE3	1.77	0.67
1:A:138:GLN:CG	2:A:703:HOH:O	2.42	0.67
1:F:138:GLN:NE2	2:F:502:HOH:O	2.28	0.67
1:B:59:ASN:HD21	1:B:87:ASN:HB2	1.61	0.66
1:C:181:GLN:HE22	1:C:220:ASN:HD21	1.41	0.66
1:F:320:PHE:H	1:F:323:GLN:HE21	1.45	0.65
1:A:138:GLN:HG2	2:A:703:HOH:O	1.96	0.65
1:D:181:GLN:HE22	1:D:220:ASN:HD21	1.45	0.64
1:A:266:LEU:HD11	1:D:266:LEU:HD11	1.80	0.64
1:B:320:PHE:H	1:B:323:GLN:HE21	1.46	0.63
1:F:215:ASP:OD1	1:F:216:ARG:NH2	2.31	0.63
1:C:27:PRO:HA	1:C:30:MET:HE3	1.79	0.62
1:D:320:PHE:H	1:D:323:GLN:HE21	1.46	0.62
1:B:27:PRO:HA	1:B:30:MET:HE3	1.83	0.61
1:E:320:PHE:H	1:E:323:GLN:HE21	1.49	0.61
1:C:320:PHE:H	1:C:323:GLN:HE21	1.50	0.60
1:E:30:MET:HE1	1:E:253:GLY:HA2	1.84	0.59
1:D:215:ASP:OD1	1:D:216:ARG:NH2	2.35	0.58
1:A:320:PHE:H	1:A:323:GLN:HE21	1.49	0.57
1:B:181:GLN:HE22	1:B:220:ASN:HD21	1.51	0.57
1:A:196:ARG:HD3	2:A:686:HOH:O	2.05	0.57
1:B:138:GLN:NE2	2:B:501:HOH:O	2.22	0.57
1:E:7:VAL:O	1:E:273:LYS:HD2	2.05	0.56
1:E:30:MET:CE	1:E:253:GLY:HA2	2.35	0.56
1:C:117:VAL:HG23	1:C:254:ALA:HB2	1.87	0.55
1:B:266:LEU:HD13	1:C:266:LEU:HD11	1.88	0.55
1:A:59:ASN:HD21	1:A:87:ASN:HB2	1.72	0.54
1:D:188:ASP:OD2	1:D:364:ARG:NH2	2.35	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:272:PRO:HB2	1:C:393:VAL:HG21	1.89	0.54
1:A:272:PRO:HB2	1:A:393:VAL:HG21	1.89	0.54
1:A:181:GLN:HE22	1:A:220:ASN:HD21	1.56	0.54
1:E:117:VAL:HG23	1:E:254:ALA:HB2	1.91	0.53
1:A:266:LEU:CD1	1:D:266:LEU:HD11	2.38	0.53
1:F:59:ASN:HD21	1:F:87:ASN:HB2	1.73	0.53
1:A:2:ARG:NH2	1:A:110:GLU:OE1	2.41	0.53
1:E:299:LYS:NZ	2:E:505:HOH:O	2.43	0.52
1:B:65:HIS:HD2	2:B:743:HOH:O	1.93	0.51
1:F:192:PRO:HG3	2:F:750:HOH:O	2.09	0.51
2:E:768:HOH:O	1:F:65:HIS:HD2	1.94	0.51
1:B:6:ILE:HD12	1:B:6:ILE:N	2.25	0.50
1:A:138:GLN:HG3	2:A:703:HOH:O	2.09	0.50
1:B:262:LEU:HD11	1:C:266:LEU:HD12	1.94	0.50
1:C:365:GLU:OE2	1:C:368:ARG:NH1	2.43	0.50
1:C:16:LYS:NZ	2:C:505:HOH:O	2.45	0.49
1:F:189:GLU:HB3	2:F:572:HOH:O	2.12	0.49
1:E:249:GLN:NE2	2:E:509:HOH:O	2.46	0.49
1:A:117:VAL:HG23	1:A:254:ALA:HB2	1.95	0.48
1:F:200:GLU:HA	1:F:206:GLU:O	2.13	0.48
1:A:297:VAL:HB	1:A:298:PRO:HD3	1.95	0.47
1:B:153:GLU:HG3	1:B:157:ARG:HD2	1.95	0.47
1:B:272:PRO:HB2	1:B:393:VAL:HG21	1.97	0.47
1:D:181:GLN:NE2	1:D:220:ASN:HD21	2.11	0.47
2:C:777:HOH:O	1:D:65:HIS:HD2	1.98	0.47
1:C:233:LYS:HA	1:C:233:LYS:HD3	1.69	0.46
1:E:181:GLN:NE2	1:E:220:ASN:HD21	2.11	0.46
1:E:6:ILE:HD13	1:E:6:ILE:N	2.31	0.46
1:E:59:ASN:HD21	1:E:87:ASN:HB2	1.81	0.46
1:C:59:ASN:HD21	1:C:87:ASN:HB2	1.81	0.46
1:F:4:VAL:HG11	1:F:101:ALA:HA	1.97	0.46
1:F:329:ASP:OD2	2:F:501:HOH:O	2.21	0.45
1:A:380:VAL:HB	1:A:384:MET:HB2	1.98	0.45
1:B:181:GLN:NE2	1:B:220:ASN:HD21	2.12	0.45
1:B:316:LEU:HG	1:B:377:THR:HG23	1.97	0.45
1:C:16:LYS:HE2	2:C:519:HOH:O	2.16	0.45
1:A:249:GLN:NE2	2:A:510:HOH:O	2.51	0.44
1:A:4:VAL:HG11	1:A:101:ALA:HA	1.98	0.44
1:E:30:MET:HE1	1:E:253:GLY:CA	2.46	0.44
1:A:266:LEU:HD11	1:D:266:LEU:CD1	2.45	0.44
1:C:52:CYS:O	1:C:83:GLY:HA2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:368:ARG:CD	2:C:643:HOH:O	2.65	0.44
1:B:4:VAL:HG11	1:B:101:ALA:HA	2.00	0.43
1:E:4:VAL:HG11	1:E:101:ALA:HA	2.00	0.43
1:E:199:VAL:HG23	1:E:208:GLN:HG2	1.99	0.43
1:B:316:LEU:HA	1:B:377:THR:O	2.19	0.43
1:D:86:LEU:C	1:D:86:LEU:HD12	2.44	0.43
1:B:300:LEU:HD21	1:B:389:LEU:HB2	2.00	0.43
1:D:137:LEU:HD23	1:D:137:LEU:HA	1.88	0.43
1:B:176:ARG:HG2	1:B:179:ARG:HH22	1.84	0.43
1:C:368:ARG:HD2	2:C:643:HOH:O	2.19	0.43
1:D:316:LEU:HA	1:D:377:THR:O	2.19	0.43
1:F:316:LEU:HA	1:F:377:THR:O	2.19	0.43
1:C:65:HIS:HA	1:D:89[A]:TYR:OH	2.19	0.42
1:C:144:ILE:O	1:D:65:HIS:HE1	2.03	0.42
1:D:4:VAL:HG11	1:D:101:ALA:HA	2.01	0.42
1:E:395:LEU:O	2:E:501:HOH:O	2.22	0.42
1:E:52:CYS:O	1:E:83:GLY:HA2	2.19	0.42
1:A:65:HIS:HA	1:B:89[A]:TYR:OH	2.20	0.42
1:D:116:GLY:HA3	1:D:352:GLY:O	2.19	0.42
1:D:200:GLU:HA	1:D:206:GLU:O	2.19	0.42
1:E:65:HIS:HA	1:F:89[A]:TYR:OH	2.20	0.42
1:A:89[B]:TYR:OH	1:B:65:HIS:HA	2.19	0.42
1:E:380:VAL:HB	1:E:384:MET:HB2	2.01	0.42
1:A:316:LEU:C	1:A:316:LEU:HD23	2.45	0.41
1:D:353:MET:HE2	1:D:353:MET:HB2	1.99	0.41
1:F:316:LEU:HG	1:F:377:THR:HG23	2.02	0.41
1:B:86:LEU:C	1:B:86:LEU:HD12	2.46	0.41
1:D:2:ARG:HH12	1:D:110:GLU:CD	2.29	0.41
1:D:350:PRO:O	1:D:353:MET:HG3	2.21	0.41
1:F:223:ASP:OD1	1:F:223:ASP:N	2.52	0.40
1:C:380:VAL:HB	1:C:384:MET:HB2	2.02	0.40
1:E:297:VAL:HB	1:E:298:PRO:HD3	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	384/402 (96%)	370 (96%)	12 (3%)	2 (0%)	24	14
1	B	393/402 (98%)	380 (97%)	11 (3%)	2 (0%)	24	14
1	C	385/402 (96%)	374 (97%)	11 (3%)	0	100	100
1	D	393/402 (98%)	379 (96%)	12 (3%)	2 (0%)	24	14
1	E	385/402 (96%)	373 (97%)	10 (3%)	2 (0%)	24	14
1	F	393/402 (98%)	381 (97%)	10 (2%)	2 (0%)	24	14
All	All	2333/2412 (97%)	2257 (97%)	66 (3%)	10 (0%)	43	19

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	89[A]	TYR
1	B	89[B]	TYR
1	D	89[A]	TYR
1	D	89[B]	TYR
1	F	89[A]	TYR
1	F	89[B]	TYR
1	E	89[A]	TYR
1	E	89[B]	TYR
1	A	89[A]	TYR
1	A	89[B]	TYR

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	306/318 (96%)	305 (100%)	1 (0%)	86	86
1	B	311/318 (98%)	309 (99%)	2 (1%)	78	77
1	C	307/318 (96%)	305 (99%)	2 (1%)	76	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	311/318 (98%)	307 (99%)	4 (1%)	61	54
1	E	307/318 (96%)	303 (99%)	4 (1%)	61	54
1	F	311/318 (98%)	307 (99%)	4 (1%)	61	54
All	All	1853/1908 (97%)	1836 (99%)	17 (1%)	70	67

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	368	ARG
1	B	189	GLU
1	B	202	LYS
1	C	207	LYS
1	C	233	LYS
1	D	189	GLU
1	D	302	LYS
1	D	368	ARG
1	D	394	ARG
1	E	2	ARG
1	E	6	ILE
1	E	233	LYS
1	E	307	LYS
1	F	6	ILE
1	F	16	LYS
1	F	189	GLU
1	F	202	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	59	ASN
1	A	65	HIS
1	A	70	ASN
1	A	131	HIS
1	A	181	GLN
1	A	208	GLN
1	A	249	GLN
1	A	323	GLN
1	A	370	ASN
1	B	59	ASN

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Mol	Chain	Res	Type
1	B	65	HIS
1	B	181	GLN
1	B	323	GLN
1	B	341	ASN
1	C	42	ASN
1	C	59	ASN
1	C	65	HIS
1	C	131	HIS
1	C	166	GLN
1	C	181	GLN
1	C	323	GLN
1	C	367	HIS
1	C	370	ASN
1	D	65	HIS
1	D	80	GLN
1	D	131	HIS
1	D	138	GLN
1	D	181	GLN
1	D	323	GLN
1	D	341	ASN
1	D	367	HIS
1	E	42	ASN
1	E	59	ASN
1	E	65	HIS
1	E	181	GLN
1	E	323	GLN
1	E	367	HIS
1	E	370	ASN
1	F	59	ASN
1	F	65	HIS
1	F	66	ASN
1	F	80	GLN
1	F	103	GLN
1	F	323	GLN
1	F	341	ASN
1	F	367	HIS

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	387/402 (96%)	-1.19	0 100 100	13, 24, 42, 75	1 (0%)
1	B	394/402 (98%)	-1.15	0 100 100	13, 25, 47, 80	1 (0%)
1	C	388/402 (96%)	-1.18	0 100 100	13, 24, 41, 81	1 (0%)
1	D	394/402 (98%)	-1.12	0 100 100	13, 25, 48, 84	1 (0%)
1	E	388/402 (96%)	-1.17	0 100 100	14, 24, 41, 79	1 (0%)
1	F	394/402 (98%)	-1.08	0 100 100	13, 25, 47, 83	1 (0%)
All	All	2345/2412 (97%)	-1.15	0 100 100	13, 24, 46, 84	6 (0%)

There are no RSRZ outliers to report.

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.