



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:38 PM UTC

PDB ID : 3N24 / pdb_00003n24
Title : The crystal structure of a two-component sensor domain (3rd form) from *Pseudomonas aeruginosa* PA01
Authors : Tan, K.; Chhor, G.; Buck, K.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2010-05-17
Resolution : 2.06 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

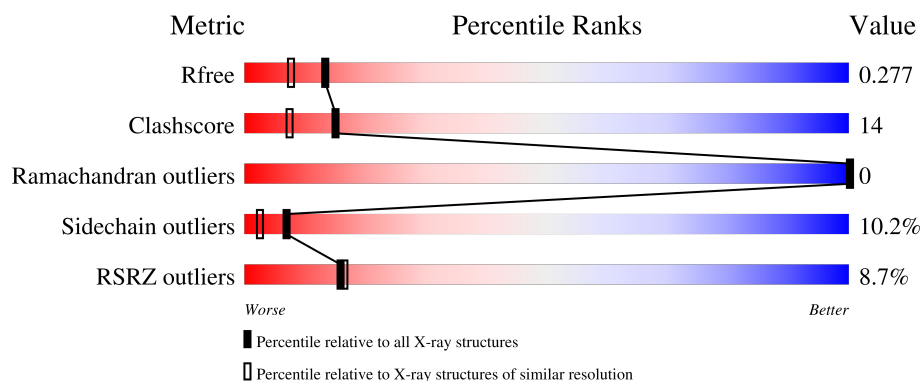
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	130	7% 62% 15% 7% 16%
1	B	130	10% 56% 19% • 22%
1	C	130	3% 68% 25% • 5%
1	D	130	7% 55% 23% • 19%
1	E	130	11% 42% 33% • 22%

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Mol	Chain	Length	Quality of chain
1	F	130	
1	G	130	
1	H	130	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 7213 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 Sensor protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	109	Total	C	N	O	Se	0	3	0
			883	542	170	170	1			
1	B	102	Total	C	N	O		0	1	0
			807	495	149	163				
1	C	123	Total	C	N	O	Se	0	7	0
			1005	614	189	200	2			
1	D	105	Total	C	N	O		0	1	0
			834	510	158	166				
1	E	102	Total	C	N	O	Se	0	0	0
			806	493	156	156	1			
1	F	114	Total	C	N	O	Se	0	3	0
			922	563	178	180	1			
1	G	110	Total	C	N	O	Se	0	0	0
			864	530	164	169	1			
1	H	111	Total	C	N	O	Se	0	2	0
			887	542	167	177	1			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	SER	-	expression tag	UNP Q9HT87
A	40	ASN	-	expression tag	UNP Q9HT87
B	39	SER	-	expression tag	UNP Q9HT87
B	40	ASN	-	expression tag	UNP Q9HT87
C	39	SER	-	expression tag	UNP Q9HT87
C	40	ASN	-	expression tag	UNP Q9HT87
D	39	SER	-	expression tag	UNP Q9HT87
D	40	ASN	-	expression tag	UNP Q9HT87
E	39	SER	-	expression tag	UNP Q9HT87
E	40	ASN	-	expression tag	UNP Q9HT87
F	39	SER	-	expression tag	UNP Q9HT87
F	40	ASN	-	expression tag	UNP Q9HT87
G	39	SER	-	expression tag	UNP Q9HT87

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Chain	Residue	Modelled	Actual	Comment	Reference
G	40	ASN	-	expression tag	UNP Q9HT87
H	39	SER	-	expression tag	UNP Q9HT87
H	40	ASN	-	expression tag	UNP Q9HT87

- Molecule 2 is water.

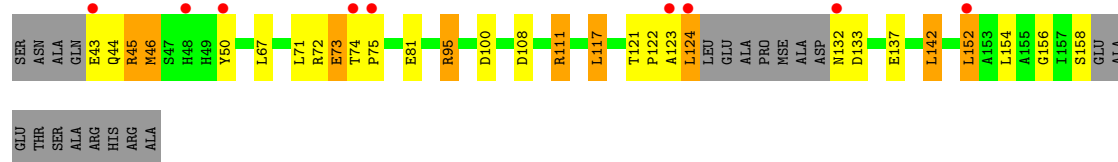
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	30	Total 30	O 30	0	0
2	B	22	Total 22	O 22	0	0
2	C	35	Total 35	O 35	0	0
2	D	21	Total 21	O 21	0	0
2	E	20	Total 20	O 20	0	0
2	F	22	Total 22	O 22	0	0
2	G	35	Total 35	O 35	0	0
2	H	20	Total 20	O 20	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

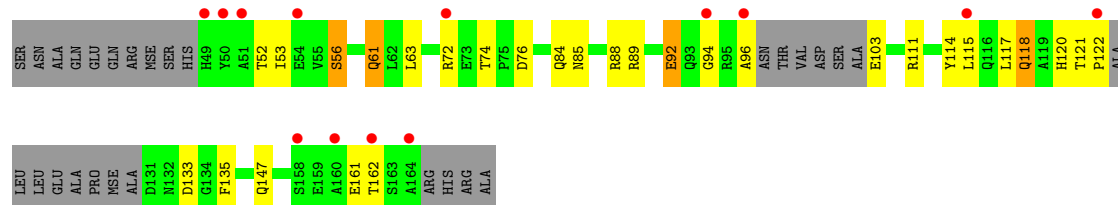
• Molecule 1: Sensor protein

Chain A: 



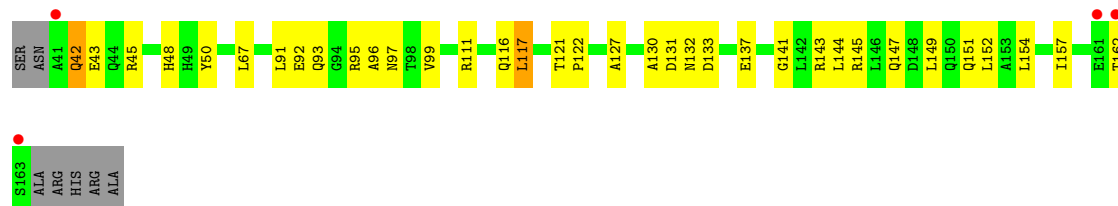
• Molecule 1: Sensor protein

Chain B: 



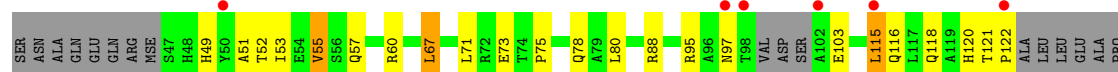
• Molecule 1: Sensor protein

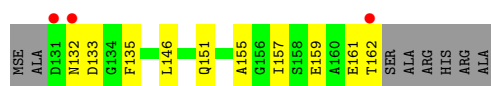
Chain C: 



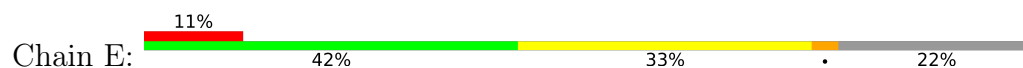
• Molecule 1: Sensor protein

Chain D: 

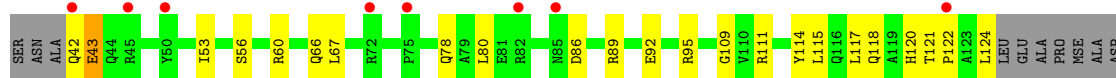




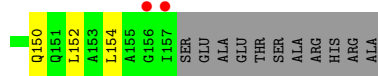
• Molecule 1: Sensor protein



• Molecule 1: Sensor protein



• Molecule 1: Sensor protein



• Molecule 1: Sensor protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.00Å 61.09Å 71.60Å 85.60° 70.47° 89.50°	Depositor
Resolution (Å)	36.63 – 2.06 36.63 – 2.06	Depositor EDS
% Data completeness (in resolution range)	90.3 (36.63-2.06) 96.8 (36.63-2.06)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.06Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.203 , 0.282 0.207 , 0.277	Depositor DCC
R_{free} test set	2317 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.449	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for h,-k,h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7213	wwPDB-VP
Average B, all atoms (Å ²)	44.0	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 61.40 % 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 1.3166e-05. 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 [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	0.52	0/895	0.74	0/1202
1	B	0.46	0/816	0.75	0/1099
1	C	0.59	0/1034	0.92	2/1392 (0.1%)
1	D	0.50	0/844	0.78	0/1137
1	E	0.51	0/810	0.80	1/1085 (0.1%)
1	F	0.54	0/938	0.78	1/1261 (0.1%)
1	G	0.51	0/871	0.76	0/1173
1	H	0.50	0/900	0.85	2/1211 (0.2%)
All	All	0.52	0/7108	0.80	6/9560 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	127	ALA	CA-C-N	9.59	130.18	119.92
1	C	127	ALA	C-N-CA	9.59	130.18	119.92
1	H	121	THR	CA-C-N	6.50	127.97	119.84
1	H	121	THR	C-N-CA	6.50	127.97	119.84
1	E	51	ALA	N-CA-C	-5.13	105.58	111.07
1	F	43	GLU	N-CA-C	-5.05	106.63	112.89

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	883	0	878	27	0
1	B	807	0	785	27	0
1	C	1005	0	995	23	0
1	D	834	0	812	22	0
1	E	806	0	789	41	0
1	F	922	0	913	27	0
1	G	864	0	846	25	0
1	H	887	0	866	33	0
2	A	30	0	0	1	0
2	B	22	0	0	2	0
2	C	35	0	0	3	0
2	D	21	0	0	1	0
2	E	20	0	0	1	0
2	F	22	0	0	0	0
2	G	35	0	0	1	0
2	H	20	0	0	2	0
All	All	7213	0	6884	190	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 14.

All (190) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:PRO:HA	2:B:177:HOH:O	1.69	0.92
1:A:123:ALA:O	1:A:124:LEU:HB2	1.73	0.88
1:A:133:ASP:O	1:A:137:GLU:HG3	1.72	0.88
1:A:45[A]:ARG:HH21	1:A:45[A]:ARG:CG	1.89	0.86
1:A:45[A]:ARG:HH21	1:A:45[A]:ARG:HG2	1.41	0.84
1:G:84:GLN:HE22	1:G:121:THR:HG21	1.42	0.84
1:H:76:ASP:CG	1:H:78:GLN:HG2	2.04	0.82
1:E:46:MSE:HG2	1:E:50:TYR:HE2	1.45	0.82
1:H:78:GLN:HG3	1:H:79:ALA:N	1.94	0.82
1:D:155:ALA:O	1:D:159:GLU:HG2	1.81	0.80
1:B:72:ARG:HH12	1:B:76:ASP:HB2	1.46	0.79
1:H:93:GLN:HG3	2:H:171:HOH:O	1.80	0.79
1:E:63:LEU:O	1:E:67:LEU:HD23	1.82	0.78
1:H:57:GLN:HE22	1:H:150:GLN:HE21	1.29	0.77
1:E:94:GLY:O	1:E:98:THR:HG23	1.86	0.76
1:H:147:GLN:O	1:H:151[B]:GLN:HG2	1.87	0.75
1:E:146:LEU:HD13	1:F:60:ARG:HG2	1.69	0.73
1:D:88:ARG:NH1	1:D:118:GLN:OE1	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:111:ARG:HD3	2:G:175:HOH:O	1.89	0.73
1:A:132:ASN:O	1:A:133:ASP:HB2	1.88	0.72
1:G:42:GLN:HE21	1:H:160:ALA:HB1	1.56	0.70
1:A:142:LEU:HD13	1:B:63:LEU:HD22	1.71	0.70
1:E:46:MSE:HG2	1:E:50:TYR:CE2	2.27	0.69
1:A:75:PRO:HG2	1:A:124:LEU:HD21	1.73	0.68
1:H:89:ARG:O	1:H:93:GLN:HG2	1.94	0.67
1:D:73:GLU:O	1:D:75:PRO:HD3	1.95	0.67
1:A:67:LEU:O	1:A:71[A]:LEU:HD13	1.95	0.66
1:A:73:GLU:HG2	1:A:74:THR:N	2.10	0.66
1:E:95:ARG:NH2	1:E:111:ARG:HD3	2.11	0.66
1:C:154:LEU:HD23	1:D:53:ILE:HD13	1.78	0.66
1:B:72:ARG:HH12	1:B:76:ASP:CB	2.08	0.65
1:E:150:GLN:HG2	1:F:56:SER:OG	1.98	0.64
1:A:95:ARG:HB3	1:A:95:ARG:NH2	2.14	0.63
1:D:95:ARG:HB3	1:D:95:ARG:NH2	2.14	0.63
1:B:85:ASN:O	1:B:89:ARG:HB2	1.99	0.63
1:C:133:ASP:O	1:C:137:GLU:HG3	1.99	0.63
1:C:117:LEU:HD21	1:D:135:PHE:CE1	2.33	0.63
1:D:51:ALA:O	1:D:55:VAL:HG13	1.98	0.63
1:E:142:LEU:HD22	1:F:117:LEU:HD22	1.81	0.62
1:E:56:SER:O	1:E:60:ARG:HG2	1.99	0.62
1:G:100:ASP:OD1	1:G:102:ALA:HB3	2.00	0.62
1:E:67:LEU:O	1:E:71:LEU:HG	2.00	0.61
1:G:63:LEU:HD22	1:H:142:LEU:HD13	1.80	0.61
1:E:111:ARG:NH1	1:E:112:ASP:OD1	2.33	0.61
1:E:150:GLN:HG3	1:F:60:ARG:HH11	1.66	0.61
1:G:84:GLN:NE2	1:G:121:THR:HG21	2.14	0.60
1:E:66:GLN:O	1:E:70:LEU:HD23	2.02	0.60
1:H:57:GLN:NE2	1:H:150:GLN:HE21	1.99	0.60
1:E:46:MSE:O	1:E:50:TYR:HD2	1.85	0.60
1:F:95:ARG:NH2	1:F:111:ARG:HD3	2.17	0.60
1:F:114:TYR:CE2	1:F:118:GLN:OE1	2.55	0.59
1:G:61:GLN:HG2	1:G:147:GLN:OE1	2.02	0.59
1:E:135:PHE:HB2	1:F:124:LEU:HD11	1.83	0.59
1:H:57:GLN:HE22	1:H:150:GLN:NE2	1.99	0.59
1:A:95:ARG:NH2	1:A:108:ASP:OD1	2.37	0.58
1:F:42:GLN:NE2	1:F:42:GLN:HA	2.18	0.58
1:G:117:LEU:HD22	1:H:142:LEU:HD12	1.84	0.58
1:B:85:ASN:OD1	1:B:88:ARG:NH1	2.36	0.58
1:H:89:ARG:HB3	1:H:89:ARG:NH1	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:LEU:HD12	1:B:117:LEU:HD22	1.85	0.57
1:C:132[B]:ASN:ND2	2:C:202:HOH:O	2.37	0.56
1:C:91:LEU:O	1:C:95:ARG:HG3	2.06	0.56
1:G:100:ASP:HB3	1:G:103:GLU:HB2	1.86	0.56
1:B:72:ARG:NH1	1:B:76:ASP:HB2	2.19	0.56
1:H:122:PRO:O	1:H:123:ALA:HB2	2.05	0.56
1:E:157:ILE:O	1:E:158:SER:C	2.49	0.55
1:B:88:ARG:NH2	1:B:118:GLN:OE1	2.40	0.55
1:C:121:THR:N	1:C:122:PRO:CD	2.70	0.55
1:A:45[A]:ARG:HG2	1:A:45[A]:ARG:NH2	2.16	0.54
1:G:91:LEU:HD21	1:G:111:ARG:HA	1.89	0.54
1:C:147[A]:GLN:HG3	1:C:151:GLN:NE2	2.23	0.54
1:A:72:ARG:HD3	2:A:183:HOH:O	2.07	0.54
1:B:111:ARG:O	1:B:115:LEU:HG	2.09	0.53
1:B:61:GLN:HE22	1:B:147:GLN:NE2	2.05	0.53
1:H:147:GLN:O	1:H:151[A]:GLN:HG3	2.09	0.53
1:C:147[B]:GLN:HG2	1:D:60:ARG:NH2	2.24	0.52
1:B:84:GLN:O	1:B:88:ARG:HG2	2.10	0.52
1:H:140:ASN:O	1:H:144:LEU:HD12	2.09	0.52
1:A:50:TYR:CZ	1:A:158:SER:HB2	2.45	0.52
1:G:80:LEU:HD21	1:G:121:THR:HG23	1.92	0.51
1:F:86:ASP:OD1	1:F:89[A]:ARG:NH2	2.36	0.51
1:A:46:MSE:HE1	1:B:161:GLU:OE1	2.10	0.51
1:G:123:ALA:C	1:G:124:LEU:HD23	2.34	0.51
1:A:45[A]:ARG:CG	1:A:45[A]:ARG:NH2	2.59	0.50
1:A:121:THR:N	1:A:122:PRO:CD	2.75	0.50
1:C:141:GLY:O	1:C:145[A]:ARG:HG3	2.12	0.50
1:A:95:ARG:HB3	1:A:95:ARG:HH21	1.75	0.49
1:G:121:THR:N	1:G:122:PRO:CD	2.75	0.49
1:H:133:ASP:O	1:H:137:GLU:HG2	2.12	0.49
1:G:46:MSE:HE3	1:H:160:ALA:HB3	1.94	0.49
1:B:94:GLY:C	1:B:96:ALA:H	2.21	0.48
1:D:95:ARG:HB3	1:D:95:ARG:HH21	1.77	0.48
1:H:78:GLN:CG	1:H:79:ALA:N	2.69	0.48
1:C:130:ALA:HB1	2:C:196:HOH:O	2.13	0.48
1:C:117:LEU:CD2	1:D:135:PHE:CE1	2.96	0.48
1:A:154:LEU:HD23	1:B:53:ILE:HD13	1.96	0.48
1:D:53:ILE:O	1:D:57:GLN:HG3	2.13	0.48
1:F:133:ASP:O	1:F:137:GLU:OE1	2.31	0.48
1:E:155:ALA:O	1:E:158:SER:HB3	2.14	0.48
1:E:151:GLN:HB3	2:E:174:HOH:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:48:HIS:CG	1:C:99:VAL:HG21	2.49	0.48
1:C:143:ARG:O	1:C:147[B]:GLN:HG3	2.13	0.48
1:F:121:THR:HB	1:F:122:PRO:HD3	1.96	0.48
1:H:76:ASP:OD2	1:H:78:GLN:HG2	2.14	0.47
1:H:51:ALA:O	1:H:54:GLU:HB3	2.14	0.47
1:C:97:ASN:HD22	1:C:97:ASN:N	2.13	0.47
1:G:121:THR:N	1:G:122:PRO:HD3	2.30	0.47
1:B:61:GLN:NE2	2:B:187:HOH:O	2.46	0.46
1:C:117:LEU:HD22	1:C:117:LEU:O	2.15	0.46
1:E:50:TYR:O	1:E:54:GLU:HG3	2.15	0.46
1:E:54:GLU:O	1:E:58:GLN:HG3	2.16	0.46
1:H:43:GLU:O	1:H:47:SER:HB2	2.15	0.46
1:H:67:LEU:O	1:H:71:LEU:HG	2.16	0.46
1:C:121:THR:N	1:C:122:PRO:HD3	2.30	0.46
1:C:42:GLN:HA	1:C:45:ARG:HG3	1.98	0.46
1:D:67:LEU:O	1:D:71:LEU:HG	2.16	0.46
1:E:49:HIS:O	1:E:52:THR:HB	2.16	0.45
1:E:111:ARG:CZ	1:E:112:ASP:OD1	2.65	0.45
1:B:61:GLN:NE2	1:B:147:GLN:NE2	2.64	0.45
1:H:159:GLU:HG3	1:H:160:ALA:N	2.28	0.45
1:G:53:ILE:HG12	1:H:153:ALA:HB1	1.98	0.45
1:H:137:GLU:HG2	1:H:137:GLU:H	1.57	0.45
1:C:92[A]:GLU:OE1	1:C:95:ARG:NH1	2.49	0.45
1:B:52:THR:O	1:B:56:SER:HB2	2.16	0.45
1:E:134:GLY:HA3	1:F:120:HIS:HD2	1.82	0.45
1:E:146:LEU:HD13	1:F:60:ARG:CG	2.44	0.45
1:B:120:HIS:O	1:B:122:PRO:HD3	2.17	0.44
1:C:137:GLU:HB3	2:C:194:HOH:O	2.17	0.44
1:E:142:LEU:CD2	1:F:117:LEU:HD22	2.47	0.44
1:E:145:ARG:NE	1:E:145:ARG:HA	2.32	0.44
1:G:42:GLN:NE2	1:H:160:ALA:HB1	2.29	0.44
1:G:122:PRO:C	1:G:124:LEU:H	2.25	0.44
1:E:53:ILE:HG13	1:F:153:ALA:HB3	1.99	0.44
1:B:61:GLN:HE22	1:B:147:GLN:HE22	1.64	0.44
1:D:97:ASN:C	1:G:41:ALA:HB2	2.42	0.44
1:A:121:THR:N	1:A:122:PRO:HD2	2.32	0.44
1:D:52:THR:OG1	1:D:103:GLU:OE2	2.36	0.44
1:A:45[B]:ARG:HH22	1:A:100:ASP:HB2	1.83	0.43
1:E:87:PHE:CE2	1:E:114:TYR:CD1	3.06	0.43
1:H:122:PRO:O	1:H:123:ALA:CB	2.66	0.43
1:C:50:TYR:CE2	1:D:157:ILE:HD13	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:147:GLN:O	1:E:151:GLN:HG3	2.18	0.43
1:H:157:ILE:O	1:H:161:GLU:HG2	2.18	0.43
1:A:43:GLU:CD	1:A:43:GLU:C	2.86	0.43
1:D:115:LEU:HD23	2:D:178:HOH:O	2.18	0.43
1:E:60:ARG:HD3	1:F:146:LEU:HB3	1.99	0.43
1:D:162:THR:HG21	1:F:133:ASP:HB3	2.00	0.43
1:E:57:GLN:O	1:E:61:GLN:HG3	2.19	0.43
1:F:53:ILE:HG21	1:F:154:LEU:CD2	2.49	0.43
1:G:150:GLN:OE1	1:H:150:GLN:NE2	2.52	0.43
1:G:46:MSE:HE1	1:H:161:GLU:HB3	2.00	0.43
1:H:145:ARG:HA	1:H:145:ARG:NE	2.34	0.43
1:E:78:GLN:H	1:E:78:GLN:HG2	1.49	0.43
1:E:145:ARG:HA	1:E:145:ARG:HE	1.83	0.43
1:B:92[B]:GLU:HA	1:B:92[B]:GLU:OE1	2.19	0.42
1:G:74:THR:HG22	1:G:74:THR:O	2.18	0.42
1:A:95:ARG:NH1	1:A:111:ARG:HH21	2.17	0.42
1:C:50:TYR:OH	1:C:157:ILE:HG22	2.18	0.42
1:G:91:LEU:HD21	1:G:111:ARG:CA	2.49	0.42
1:B:118:GLN:HE21	1:B:118:GLN:HB2	1.56	0.42
1:E:113:ALA:HB1	1:F:142:LEU:HA	2.02	0.42
1:E:153:ALA:HB1	1:F:53:ILE:HG12	2.01	0.42
1:F:42:GLN:HA	1:F:42:GLN:HE21	1.83	0.42
1:D:146:LEU:HD23	1:D:146:LEU:HA	1.92	0.42
1:A:152:LEU:O	1:A:156:GLY:N	2.52	0.42
1:E:114:TYR:O	1:E:117:LEU:HB3	2.19	0.42
1:H:61:GLN:NE2	2:H:21:HOH:O	2.51	0.42
1:B:121:THR:HA	1:B:122:PRO:HD3	1.89	0.42
1:D:116:GLN:O	1:D:120:HIS:HD2	2.02	0.42
1:F:53:ILE:HG21	1:F:154:LEU:HD22	2.02	0.42
1:E:56:SER:HB3	1:F:149:LEU:HG	2.02	0.41
1:E:145:ARG:HG2	1:F:109:GLY:O	2.20	0.41
1:A:67:LEU:HD13	1:A:67:LEU:HA	1.85	0.41
1:F:118:GLN:O	1:F:122:PRO:HD3	2.20	0.41
1:A:117:LEU:HD21	1:B:135:PHE:CE1	2.54	0.41
1:B:72:ARG:NH1	1:B:76:ASP:CB	2.81	0.41
1:B:85:ASN:CG	1:B:88:ARG:HH11	2.29	0.41
1:C:96:ALA:C	1:C:97:ASN:HD22	2.28	0.41
1:D:49:HIS:O	1:D:53:ILE:HG13	2.21	0.41
1:E:66:GLN:O	1:E:70:LEU:CD2	2.68	0.41
1:G:56:SER:O	1:G:60:ARG:HG3	2.21	0.41
1:H:121:THR:HA	1:H:122:PRO:HD3	1.58	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:LEU:HD21	1:C:111:ARG:HA	2.02	0.40
1:D:121:THR:HA	1:D:122:PRO:HD3	1.91	0.40
1:D:97:ASN:HA	1:G:41:ALA:HB2	2.02	0.40
1:E:55:VAL:O	1:E:59:LEU:HG	2.21	0.40
1:B:84:GLN:HA	1:B:114:TYR:OH	2.21	0.40
1:E:143:ARG:O	1:E:147:GLN:HG3	2.21	0.40
1:F:66:GLN:HG2	1:F:80:LEU:HD12	2.02	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	108/130 (83%)	105 (97%)	3 (3%)	0	100	100
1	B	97/130 (75%)	95 (98%)	2 (2%)	0	100	100
1	C	128/130 (98%)	127 (99%)	1 (1%)	0	100	100
1	D	100/130 (77%)	97 (97%)	3 (3%)	0	100	100
1	E	96/130 (74%)	95 (99%)	1 (1%)	0	100	100
1	F	113/130 (87%)	112 (99%)	1 (1%)	0	100	100
1	G	106/130 (82%)	105 (99%)	1 (1%)	0	100	100
1	H	109/130 (84%)	107 (98%)	2 (2%)	0	100	100
All	All	857/1040 (82%)	843 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	93/103 (90%)	81 (87%)	12 (13%)	4	1
1	B	84/103 (82%)	75 (89%)	9 (11%)	6	2
1	C	107/103 (104%)	95 (89%)	12 (11%)	6	2
1	D	87/103 (84%)	78 (90%)	9 (10%)	7	2
1	E	83/103 (81%)	74 (89%)	9 (11%)	6	2
1	F	97/103 (94%)	89 (92%)	8 (8%)	10	5
1	G	89/103 (86%)	78 (88%)	11 (12%)	4	1
1	H	93/103 (90%)	87 (94%)	6 (6%)	15	8
All	All	733/824 (89%)	657 (90%)	76 (10%)	7	2

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	45[A]	ARG
1	A	45[B]	ARG
1	A	46	MSE
1	A	73	GLU
1	A	81	GLU
1	A	95	ARG
1	A	111	ARG
1	A	117	LEU
1	A	124	LEU
1	A	142	LEU
1	A	152	LEU
1	B	56	SER
1	B	61	GLN
1	B	74	THR
1	B	92[A]	GLU
1	B	92[B]	GLU
1	B	103	GLU
1	B	118	GLN
1	B	133	ASP
1	B	162	THR
1	C	42	GLN
1	C	43	GLU

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Mol	Chain	Res	Type
1	C	67	LEU
1	C	93	GLN
1	C	116	GLN
1	C	117	LEU
1	C	131[A]	ASP
1	C	131[B]	ASP
1	C	144	LEU
1	C	149	LEU
1	C	152	LEU
1	C	162	THR
1	D	55	VAL
1	D	67	LEU
1	D	78	GLN
1	D	80	LEU
1	D	115	LEU
1	D	132	ASN
1	D	133	ASP
1	D	151	GLN
1	D	161	GLU
1	E	53	ILE
1	E	62	LEU
1	E	70	LEU
1	E	78	GLN
1	E	86	ASP
1	E	104	GLN
1	E	133	ASP
1	E	137	GLU
1	E	145	ARG
1	F	43	GLU
1	F	67	LEU
1	F	78	GLN
1	F	115	LEU
1	F	133	ASP
1	F	137	GLU
1	F	144	LEU
1	F	162	THR
1	G	46	MSE
1	G	48	HIS
1	G	53	ILE
1	G	86	ASP
1	G	89	ARG
1	G	97	ASN

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Mol	Chain	Res	Type
1	G	99	VAL
1	G	121	THR
1	G	137	GLU
1	G	152	LEU
1	G	154	LEU
1	H	72	ARG
1	H	137	GLU
1	H	142	LEU
1	H	144	LEU
1	H	154	LEU
1	H	159	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	GLN
1	A	66	GLN
1	A	132	ASN
1	B	61	GLN
1	B	120	HIS
1	B	147	GLN
1	C	97	ASN
1	D	57	GLN
1	D	58	GLN
1	D	66	GLN
1	D	120	HIS
1	D	147	GLN
1	E	44	GLN
1	E	97	ASN
1	F	78	GLN
1	F	120	HIS
1	F	147	GLN
1	G	42	GLN
1	G	84	GLN
1	G	104	GLN
1	H	85	ASN
1	H	97	ASN
1	H	150	GLN

5.3.3 RNA [i](#)

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	108/130 (83%)	0.22	9 (8%) 17 17	16, 33, 70, 100	4 (3%)
1	B	102/130 (78%)	0.55	13 (12%) 8 7	20, 45, 91, 127	1 (0%)
1	C	121/130 (93%)	0.06	4 (3%) 49 49	12, 31, 68, 101	7 (5%)
1	D	105/130 (80%)	0.41	9 (8%) 16 17	17, 37, 83, 101	3 (2%)
1	E	101/130 (77%)	0.88	14 (13%) 6 6	22, 50, 87, 104	0
1	F	113/130 (86%)	0.51	11 (9%) 13 13	18, 39, 87, 102	3 (2%)
1	G	109/130 (83%)	0.50	13 (11%) 9 9	20, 37, 78, 108	0
1	H	110/130 (84%)	0.20	3 (2%) 56 58	17, 34, 79, 90	3 (2%)
All	All	869/1040 (83%)	0.41	76 (8%) 16 16	12, 38, 85, 127	21 (2%)

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	41	ALA	5.1
1	G	132	ASN	5.0
1	D	102	ALA	4.3
1	D	122	PRO	4.1
1	C	162	THR	4.0
1	E	80	LEU	3.8
1	B	96	ALA	3.8
1	H	75	PRO	3.7
1	G	99	VAL	3.7
1	C	41	ALA	3.5
1	A	124	LEU	3.5
1	D	97	ASN	3.3
1	B	122	PRO	3.2
1	E	132	ASN	3.2
1	E	117	LEU	3.2
1	G	133	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	F	136	SER	3.1
1	F	45	ARG	3.1
1	E	158	SER	3.0
1	D	131	ASP	3.0
1	D	162	THR	3.0
1	A	152	LEU	2.9
1	E	154	LEU	2.9
1	B	115	LEU	2.9
1	G	124	LEU	2.8
1	H	45	ARG	2.8
1	G	75	PRO	2.8
1	H	72	ARG	2.8
1	B	164	ALA	2.8
1	E	68	VAL	2.8
1	A	50	TYR	2.8
1	E	115	LEU	2.8
1	F	42	GLN	2.8
1	A	123	ALA	2.8
1	B	162	THR	2.8
1	G	122	PRO	2.8
1	F	85	ASN	2.7
1	B	160	ALA	2.7
1	F	162	THR	2.6
1	F	75	PRO	2.6
1	F	122	PRO	2.6
1	A	43	GLU	2.6
1	G	42	GLN	2.5
1	F	161	GLU	2.5
1	A	48	HIS	2.5
1	B	49	HIS	2.5
1	B	51	ALA	2.4
1	E	62	LEU	2.4
1	G	157	ILE	2.4
1	E	157	ILE	2.4
1	B	54	GLU	2.4
1	G	74	THR	2.4
1	D	115	LEU	2.3
1	C	163	SER	2.3
1	D	98	THR	2.3
1	B	94	GLY	2.3
1	F	50	TYR	2.3
1	B	72	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	F	82[A]	ARG	2.2
1	B	50	TYR	2.2
1	E	81	GLU	2.2
1	F	72	ARG	2.2
1	A	74	THR	2.2
1	G	86	ASP	2.2
1	E	120	HIS	2.2
1	E	67	LEU	2.2
1	A	75	PRO	2.1
1	E	98	THR	2.2
1	B	158	SER	2.1
1	D	132	ASN	2.1
1	D	50	TYR	2.1
1	G	48	HIS	2.1
1	C	161	GLU	2.1
1	G	156	GLY	2.1
1	A	132	ASN	2.0
1	E	77	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.