



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:09 AM GMT

PDB ID : 2PKH
Title : Structural Genomics, the crystal structure of the C-terminal domain of histidine utilization repressor from *Pseudomonas syringae* pv. tomato str. DC3000
Authors : Tan, K.; Zhou, M.; Clancy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2007-04-17
Resolution : 1.95 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

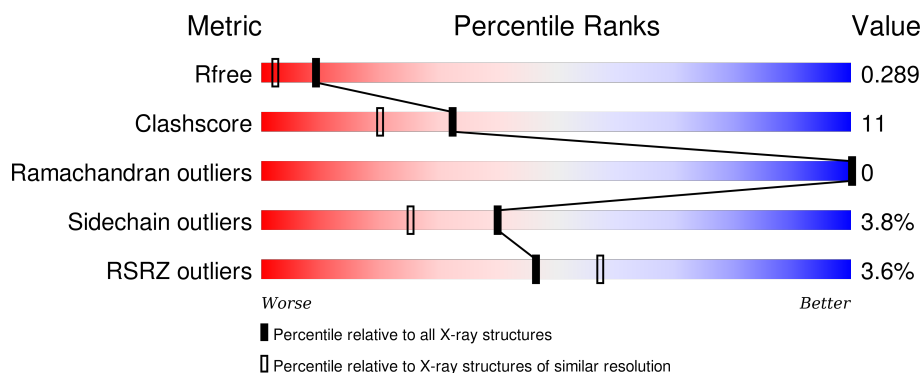
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	91344	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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. 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	148	 4% 76% 16% •• 5%
1	B	148	 2% 76% 16% • •
1	C	148	 6% 77% 18% •• •
1	D	148	 3% 77% 16% • 5%
1	E	148	 2% 75% 19% •• •

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Mol	Chain	Length	Quality of chain
1	F	148	
1	G	148	
1	H	148	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	EDO	B	1010	-	-	-	X
2	EDO	B	1026	-	-	X	-
2	EDO	B	1029	-	-	-	X
2	EDO	C	1017	-	-	X	X
2	EDO	D	1003	-	-	-	X
2	EDO	E	1005	-	-	-	X
2	EDO	F	1028	-	-	X	X
2	EDO	H	1025	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10115 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 Histidine utilization repressor.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	Se	0	0	0
			1127	708	208	206	3	2			
1	B	142	Total	C	N	O	S	Se	0	0	0
			1137	714	211	207	3	2			
1	C	142	Total	C	N	O	S	Se	0	0	0
			1137	714	211	207	3	2			
1	D	141	Total	C	N	O	S	Se	0	0	0
			1127	708	208	206	3	2			
1	E	142	Total	C	N	O	S	Se	0	0	0
			1137	714	211	207	3	2			
1	F	141	Total	C	N	O	S	Se	0	0	0
			1127	708	208	206	3	2			
1	G	141	Total	C	N	O	S	Se	0	0	0
			1127	708	208	206	3	2			
1	H	142	Total	C	N	O	S	Se	0	0	0
			1137	714	211	207	3	2			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
A	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
A	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
A	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
A	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
B	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
B	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
B	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
B	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
B	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
C	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
C	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
C	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
C	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
D	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
D	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
D	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
D	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
D	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
E	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
E	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
E	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
E	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
E	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
F	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
F	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
F	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
F	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
F	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
G	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
G	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
G	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
G	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
G	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
H	102	SER	-	CLONING ARTIFACT	UNP Q87UX0
H	103	ASN	-	CLONING ARTIFACT	UNP Q87UX0
H	104	ALA	-	CLONING ARTIFACT	UNP Q87UX0
H	115	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0
H	132	MSE	MET	MODIFIED RESIDUE	UNP Q87UX0

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	F	1	Total	C	O	0	0
			4	2	2		
2	H	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	E	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	F	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	123	Total O 123 123	0	0

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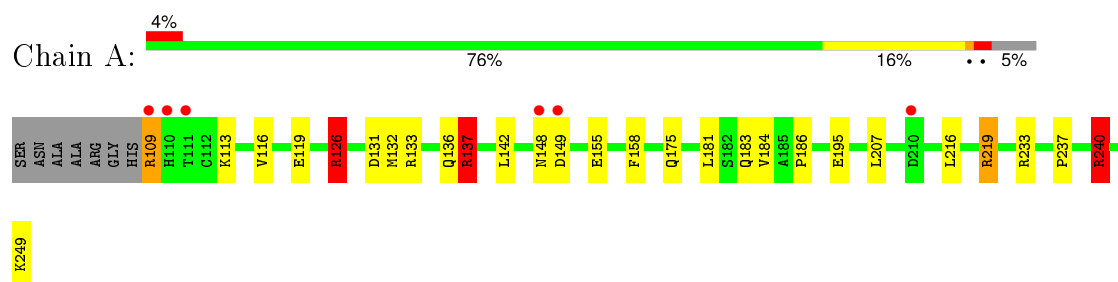
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	116	Total 116	O 116	0	0
3	C	106	Total 106	O 106	0	0
3	D	101	Total 101	O 101	0	0
3	E	121	Total 121	O 121	0	0
3	F	123	Total 123	O 123	0	0
3	G	133	Total 133	O 133	0	0
3	H	108	Total 108	O 108	0	0

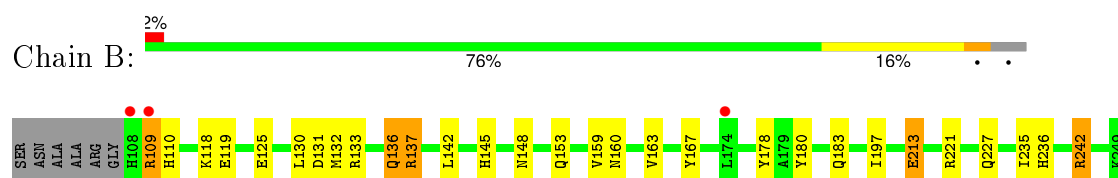
3 Residue-property plots

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

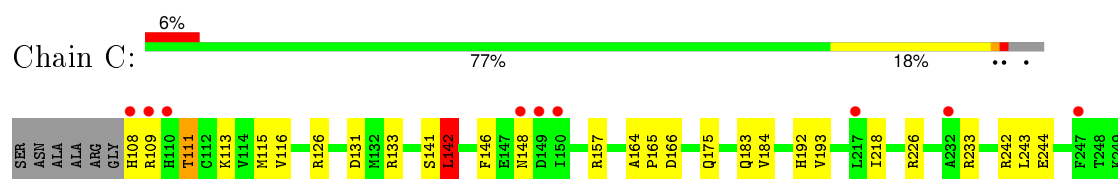
- Molecule 1: Histidine utilization repressor



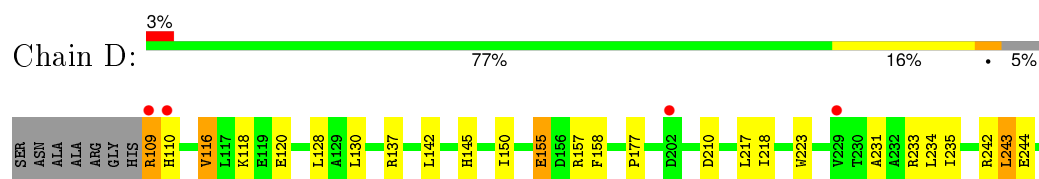
- Molecule 1: Histidine utilization repressor



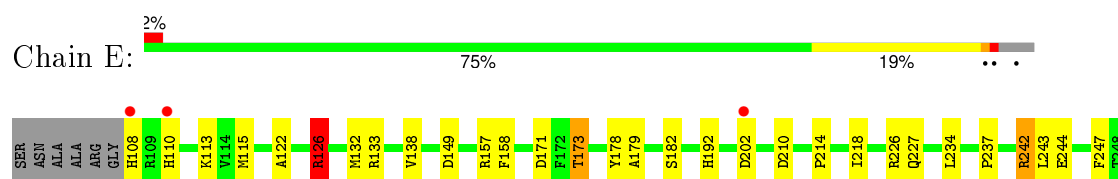
- Molecule 1: Histidine utilization repressor



- Molecule 1: Histidine utilization repressor



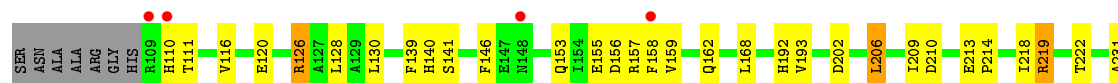
- Molecule 1: Histidine utilization repressor




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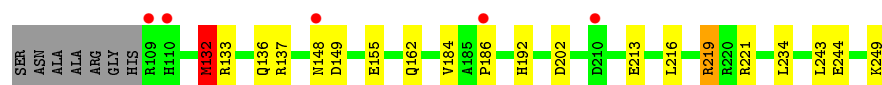
- Molecule 1: Histidine utilization repressor

Chain F:  3% 69% 24% 5%




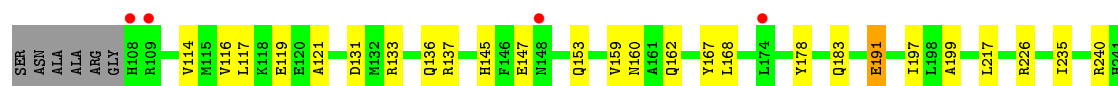
- Molecule 1: Histidine utilization repressor

Chain G:  3% 82% 12% 5%



- Molecule 1: Histidine utilization repressor

Chain H:  3% 77% 18% 2%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	100.84Å 100.95Å 130.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.87 – 1.95 39.87 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.87-1.95) 99.1 (39.87-1.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.219 , 0.277 0.232 , 0.289	Depositor DCC
R_{free} test set	4836 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.1	EDS
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	8 of 96704 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10115	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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 62.56 % 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 1.1050e-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, 0.375 respectively for untwinned datasets, and 0.333, 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: EDO

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	2/1147 (0.2%)	1.36	11/1549 (0.7%)
1	B	1.10	1/1158 (0.1%)	0.99	1/1564 (0.1%)
1	C	1.06	0/1158	1.03	5/1564 (0.3%)
1	D	1.08	1/1147 (0.1%)	1.01	2/1549 (0.1%)
1	E	1.06	0/1158	1.06	5/1564 (0.3%)
1	F	1.09	1/1147 (0.1%)	1.06	6/1549 (0.4%)
1	G	1.17	3/1147 (0.3%)	1.07	4/1549 (0.3%)
1	H	1.18	4/1158 (0.3%)	0.99	2/1564 (0.1%)
All	All	1.10	12/9220 (0.1%)	1.08	36/12452 (0.3%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	155	GLU	CD-OE1	-7.66	1.17	1.25
1	F	155	GLU	CB-CG	7.39	1.66	1.52
1	H	191	GLU	CB-CG	-6.46	1.39	1.52
1	G	184	VAL	CB-CG2	-6.26	1.39	1.52
1	H	121	ALA	CA-CB	6.25	1.65	1.52
1	G	155	GLU	CG-CD	5.92	1.60	1.51
1	H	119	GLU	CB-CG	-5.76	1.41	1.52
1	B	213	GLU	CB-CG	5.47	1.62	1.52
1	A	126	ARG	CD-NE	-5.09	1.37	1.46
1	D	155	GLU	CB-CG	5.03	1.61	1.52
1	A	155	GLU	CD-OE2	-5.03	1.20	1.25
1	H	199	ALA	CA-CB	-5.03	1.41	1.52

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	240	ARG	NE-CZ-NH1	19.22	129.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ARG	NE-CZ-NH2	-17.87	111.36	120.30
1	A	240	ARG	NE-CZ-NH2	-15.44	112.58	120.30
1	A	126	ARG	NE-CZ-NH1	13.99	127.30	120.30
1	G	219	ARG	NE-CZ-NH2	-9.81	115.40	120.30
1	F	206	LEU	CA-CB-CG	8.20	134.17	115.30
1	A	137	ARG	NE-CZ-NH2	-8.02	116.29	120.30
1	A	219	ARG	NE-CZ-NH1	7.54	124.07	120.30
1	G	219	ARG	NE-CZ-NH1	7.41	124.01	120.30
1	E	157	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	D	157	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	E	210	ASP	CB-CG-OD1	6.58	124.22	118.30
1	F	202	ASP	CB-CG-OD1	6.54	124.19	118.30
1	H	242	ARG	NE-CZ-NH2	-6.53	117.03	120.30
1	F	157	ARG	NE-CZ-NH1	6.43	123.52	120.30
1	A	240	ARG	CD-NE-CZ	6.23	132.32	123.60
1	F	202	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	G	132	MSE	CA-CB-CG	6.21	123.85	113.30
1	E	126	ARG	NE-CZ-NH1	-6.12	117.24	120.30
1	C	157	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	G	137	ARG	NE-CZ-NH1	-6.00	117.30	120.30
1	C	226	ARG	NE-CZ-NH1	-5.86	117.37	120.30
1	C	233	ARG	NE-CZ-NH1	-5.83	117.38	120.30
1	C	142	LEU	CB-CG-CD1	5.83	120.91	111.00
1	B	242	ARG	NE-CZ-NH1	5.71	123.16	120.30
1	A	216	LEU	CB-CG-CD2	-5.47	101.70	111.00
1	E	242	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	A	137	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	H	240	ARG	NE-CZ-NH2	-5.33	117.63	120.30
1	E	157	ARG	NE-CZ-NH2	-5.31	117.64	120.30
1	A	219	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	F	126	ARG	NE-CZ-NH2	5.20	122.90	120.30
1	C	142	LEU	CA-CB-CG	5.13	127.11	115.30
1	F	168	LEU	CB-CG-CD2	5.11	119.69	111.00
1	D	210	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	A	233	ARG	NE-CZ-NH1	5.06	122.83	120.30

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	1127	0	1127	20	0
1	B	1137	0	1134	32	0
1	C	1137	0	1134	32	0
1	D	1127	0	1127	21	0
1	E	1137	0	1134	35	0
1	F	1127	0	1127	38	0
1	G	1127	0	1127	14	0
1	H	1137	0	1134	19	0
2	A	20	0	30	0	0
2	B	24	0	36	7	0
2	C	12	0	18	9	0
2	D	12	0	18	2	0
2	E	20	0	30	0	0
2	F	12	0	18	5	0
2	G	8	0	12	2	0
2	H	20	0	30	1	0
3	A	123	0	0	3	0
3	B	116	0	0	3	0
3	C	106	0	0	3	0
3	D	101	0	0	3	0
3	E	121	0	0	1	0
3	F	123	0	0	8	0
3	G	133	0	0	4	0
3	H	108	0	0	6	0
All	All	10115	0	9236	194	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 11.

All (194) 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:C:113:LYS:HG2	1:C:115:MSE:CE	1.62	1.29
1:C:113:LYS:HG2	1:C:115:MSE:HE1	1.22	1.13
1:E:113:LYS:HG2	1:E:115:MSE:HE1	1.29	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:202:ASP:HB2	3:E:1108:HOH:O	1.55	1.07
1:C:113:LYS:HG2	1:C:115:MSE:HE3	1.44	1.00
1:H:131:ASP:HB3	3:H:1129:HOH:O	1.63	0.99
1:E:113:LYS:HG2	1:E:115:MSE:CE	1.99	0.92
1:E:113:LYS:CG	1:E:115:MSE:HE1	2.00	0.91
1:C:111:THR:HG22	1:C:146:PHE:HB2	1.54	0.88
1:F:111:THR:HG22	3:F:1130:HOH:O	1.79	0.82
2:B:1026:EDO:H21	3:B:1144:HOH:O	1.79	0.82
1:A:148:ASN:O	1:A:149:ASP:HB2	1.80	0.82
2:G:1023:EDO:H11	3:G:1150:HOH:O	1.79	0.80
1:A:119:GLU:OE2	1:A:137:ARG:CD	2.29	0.80
1:E:122:ALA:HB2	1:E:132:MSE:CE	2.11	0.79
1:E:113:LYS:HB3	1:E:115:MSE:HE3	1.65	0.78
1:C:115:MSE:HE2	1:C:115:MSE:HA	1.66	0.77
1:B:197:ILE:HD11	1:B:235:ILE:HD12	1.66	0.77
1:F:213:GLU:HB2	2:F:1028:EDO:O1	1.83	0.77
1:A:119:GLU:OE2	1:A:137:ARG:HD3	1.85	0.76
1:H:114:VAL:HG12	3:H:1073:HOH:O	1.85	0.75
1:B:110:HIS:NE2	1:B:145:HIS:HD2	1.85	0.74
1:C:133:ARG:HG2	1:H:242:ARG:HD3	1.70	0.74
1:E:122:ALA:CB	1:E:132:MSE:CE	2.66	0.73
1:B:130:LEU:HB3	1:B:132:MSE:HE3	1.69	0.73
1:E:122:ALA:HB2	1:E:132:MSE:HE1	1.71	0.71
1:C:175:GLN:HE22	1:C:183:GLN:NE2	1.88	0.71
1:C:113:LYS:CG	1:C:115:MSE:CE	2.58	0.71
1:C:113:LYS:CG	1:C:115:MSE:HE1	2.12	0.71
1:G:221:ARG:HD2	3:G:1090:HOH:O	1.91	0.70
1:A:237:PRO:HB2	1:A:240:ARG:HG3	1.73	0.70
1:B:242:ARG:HD3	1:E:133:ARG:HG2	1.72	0.70
1:C:113:LYS:CG	1:C:115:MSE:HE3	2.21	0.69
1:E:244:GLU:HG2	1:F:193:VAL:HG22	1.75	0.69
1:H:133:ARG:NH2	1:H:136:GLN:NE2	2.42	0.68
1:G:148:ASN:O	1:G:149:ASP:HB2	1.93	0.68
1:B:133:ARG:NH2	1:B:136:GLN:OE1	2.28	0.67
1:B:109:ARG:CG	1:B:109:ARG:HH11	2.08	0.65
1:E:126:ARG:HD3	1:E:158:PHE:CD1	2.31	0.65
1:C:184:VAL:HG11	2:C:1017:EDO:H11	1.79	0.65
1:B:132:MSE:HE1	1:B:160:ASN:HD22	1.61	0.65
1:F:153:GLN:HE21	1:F:236:HIS:CE1	2.15	0.65
1:F:111:THR:CG2	1:F:146:PHE:HB2	2.27	0.64
1:A:119:GLU:OE2	1:A:137:ARG:HD2	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:219:ARG:HD3	3:F:1091:HOH:O	1.97	0.64
1:C:116:VAL:HG13	1:C:142:LEU:HB3	1.80	0.64
1:D:116:VAL:HG13	1:D:142:LEU:HB3	1.79	0.64
1:G:132:MSE:HE2	1:G:136:GLN:HB3	1.80	0.63
1:F:158:PHE:CZ	1:F:231:ALA:HB3	2.33	0.63
1:A:186:PRO:HG3	1:D:110:HIS:HB2	1.80	0.62
1:E:113:LYS:CG	1:E:115:MSE:CE	2.69	0.62
1:B:137:ARG:HG3	1:B:137:ARG:HH21	1.64	0.62
1:B:153:GLN:HE21	1:B:236:HIS:CE1	2.17	0.61
1:D:242:ARG:NE	1:D:244:GLU:OE2	2.34	0.61
1:A:113:LYS:HD3	3:A:1147:HOH:O	2.00	0.61
1:D:109:ARG:HH11	1:D:109:ARG:HB3	1.65	0.61
1:H:116:VAL:C	3:H:1073:HOH:O	2.39	0.60
1:B:160:ASN:OD1	1:B:163:VAL:HG23	2.02	0.60
1:D:158:PHE:CZ	1:D:231:ALA:HB3	2.36	0.60
1:A:249:LYS:HE2	1:B:178:TYR:OH	2.02	0.60
1:F:111:THR:HG23	1:F:146:PHE:HB2	1.83	0.59
1:C:184:VAL:HG22	2:C:1017:EDO:H22	1.85	0.59
1:A:219:ARG:HH11	1:A:219:ARG:HG2	1.68	0.59
1:D:150:ILE:CD1	1:G:162:GLN:NE2	2.65	0.59
1:F:110:HIS:HB2	1:G:186:PRO:HG3	1.84	0.58
1:C:242:ARG:HD2	1:C:244:GLU:OE2	2.03	0.58
1:D:116:VAL:HG13	1:D:142:LEU:CB	2.33	0.58
1:E:122:ALA:HB2	1:E:132:MSE:HE3	1.86	0.58
1:B:119:GLU:OE2	1:B:137:ARG:NH2	2.35	0.58
1:F:120:GLU:OE2	1:F:126:ARG:NH2	2.33	0.58
1:A:116:VAL:HG13	1:A:142:LEU:HB3	1.86	0.57
1:A:132:MSE:HE2	1:A:136:GLN:HB3	1.87	0.57
2:D:1013:EDO:H22	3:D:1084:HOH:O	2.05	0.57
1:A:240:ARG:HD3	3:A:1136:HOH:O	2.06	0.56
1:B:180:TYR:O	1:B:183:GLN:HB2	2.04	0.56
1:G:216:LEU:HD21	1:G:243:LEU:HG	1.88	0.56
1:H:197:ILE:HD11	1:H:235:ILE:HD12	1.88	0.55
1:F:233:ARG:HG2	1:F:235:ILE:HD11	1.88	0.55
1:B:110:HIS:NE2	1:B:145:HIS:CD2	2.71	0.55
1:E:122:ALA:CB	1:E:132:MSE:HE1	2.35	0.54
1:C:165:PRO:HD2	2:C:1017:EDO:H12	1.90	0.54
1:B:132:MSE:HB2	1:B:136:GLN:HE21	1.72	0.54
1:A:175:GLN:OE1	1:A:183:GLN:NE2	2.41	0.53
1:A:126:ARG:HD3	1:A:158:PHE:CD1	2.43	0.53
1:E:179:ALA:O	1:E:182:SER:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:HG3	1:B:109:ARG:NH1	2.24	0.53
1:C:164:ALA:HA	2:C:1017:EDO:H12	1.91	0.53
1:H:197:ILE:HD13	1:H:217:LEU:HD22	1.91	0.53
1:C:108:HIS:N	3:C:1094:HOH:O	2.42	0.53
1:D:118:LYS:HE3	1:D:120:GLU:OE1	2.09	0.52
1:C:218:ILE:HD11	1:C:243:LEU:HD21	1.91	0.52
1:C:192:HIS:CD2	1:D:247:PHE:HE1	2.28	0.52
1:B:109:ARG:HD3	1:B:148:ASN:OD1	2.08	0.52
1:E:178:TYR:OH	1:F:249:LYS:HE2	2.10	0.52
1:E:113:LYS:CB	1:E:115:MSE:HE3	2.36	0.52
1:H:117:LEU:HG	1:H:168:LEU:HD22	1.92	0.52
1:C:111:THR:HB	3:C:1057:HOH:O	2.09	0.51
1:E:132:MSE:HE1	1:E:138:VAL:CG1	2.41	0.51
1:E:247:PHE:CE2	1:F:192:HIS:HD2	2.28	0.51
1:F:219:ARG:HG2	3:F:1091:HOH:O	2.10	0.51
1:B:153:GLN:HE21	1:B:236:HIS:HE1	1.57	0.51
1:F:162:GLN:HG2	3:F:1030:HOH:O	2.11	0.51
1:B:109:ARG:CG	1:B:109:ARG:NH1	2.71	0.51
1:E:122:ALA:HB1	1:E:132:MSE:CE	2.41	0.51
1:C:115:MSE:HE2	1:C:115:MSE:CA	2.39	0.50
1:B:213:GLU:OE1	2:B:1026:EDO:H11	2.11	0.50
1:F:219:ARG:CD	3:F:1091:HOH:O	2.57	0.50
1:D:130:LEU:HD11	1:D:158:PHE:CD1	2.47	0.50
1:H:145:HIS:CD2	1:H:153:GLN:HE21	2.29	0.50
1:E:108:HIS:HB2	1:E:110:HIS:NE2	2.27	0.50
1:F:210:ASP:H	2:F:1028:EDO:C1	2.25	0.50
1:F:153:GLN:HE21	1:F:236:HIS:HE1	1.58	0.50
1:E:108:HIS:HB2	1:E:110:HIS:CD2	2.47	0.49
1:C:175:GLN:HE22	1:C:183:GLN:HE22	1.58	0.49
1:H:133:ARG:NH2	1:H:136:GLN:HE21	2.11	0.49
1:D:110:HIS:CE1	1:D:145:HIS:ND1	2.81	0.49
1:E:242:ARG:HD2	1:E:244:GLU:OE2	2.13	0.49
1:B:109:ARG:HG3	1:B:109:ARG:HH11	1.74	0.49
1:C:109:ARG:NH1	1:C:148:ASN:HD22	2.11	0.48
1:B:213:GLU:OE1	2:B:1026:EDO:C1	2.62	0.48
1:F:139:PHE:HB2	1:F:159:VAL:HG12	1.96	0.48
1:H:160:ASN:OD1	1:H:162:GLN:N	2.47	0.48
1:B:142:LEU:C	1:B:142:LEU:HD23	2.33	0.48
1:E:113:LYS:CD	1:E:115:MSE:HE1	2.43	0.47
1:A:131:ASP:OD1	1:F:242:ARG:NH2	2.46	0.47
1:F:110:HIS:CB	1:G:186:PRO:HG3	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:116:VAL:O	1:C:141:SER:HA	2.13	0.47
1:G:192:HIS:CG	3:G:1082:HOH:O	2.67	0.47
1:H:162:GLN:NE2	3:H:1089:HOH:O	2.42	0.47
1:B:131:ASP:O	2:B:1029:EDO:H21	2.15	0.46
1:H:137:ARG:HD3	3:H:1085:HOH:O	2.15	0.46
2:C:1002:EDO:H21	1:D:223:TRP:CE2	2.51	0.46
1:E:132:MSE:HE1	1:E:138:VAL:HG13	1.97	0.46
2:D:1003:EDO:H12	3:D:1053:HOH:O	2.14	0.46
1:E:247:PHE:HE2	1:F:192:HIS:HD2	1.63	0.46
1:E:249:LYS:HE3	1:F:222:THR:HG23	1.98	0.46
1:A:207:LEU:O	1:A:240:ARG:NH2	2.49	0.45
1:B:227:GLN:HG3	2:B:1029:EDO:H12	1.98	0.45
1:F:111:THR:CG2	3:F:1130:HOH:O	2.51	0.45
1:B:137:ARG:HH21	1:B:137:ARG:CG	2.28	0.45
1:A:181:LEU:HA	1:A:184:VAL:HG22	1.98	0.45
1:E:115:MSE:HA	1:E:115:MSE:HE2	1.99	0.45
1:E:113:LYS:HB3	1:E:115:MSE:CE	2.43	0.45
1:F:111:THR:HG22	1:F:146:PHE:HB2	1.99	0.45
1:H:117:LEU:HB2	3:H:1073:HOH:O	2.16	0.45
1:C:164:ALA:HA	2:C:1017:EDO:C1	2.47	0.45
1:C:193:VAL:HG22	1:D:244:GLU:HG2	1.98	0.45
3:D:1062:HOH:O	1:G:136:GLN:HG2	2.16	0.45
1:C:184:VAL:CG2	2:C:1017:EDO:H22	2.46	0.45
1:F:120:GLU:OE2	1:F:126:ARG:NH1	2.49	0.44
1:C:192:HIS:HD2	1:D:247:PHE:HE1	1.62	0.44
1:E:113:LYS:CB	1:E:115:MSE:CE	2.96	0.44
1:A:195:GLU:OE2	1:B:242:ARG:NH1	2.39	0.44
1:D:217:LEU:HD11	1:D:233:ARG:HG3	1.98	0.44
1:G:132:MSE:CE	1:G:136:GLN:HB3	2.45	0.44
1:F:116:VAL:O	1:F:141:SER:HA	2.18	0.44
1:G:244:GLU:OE1	2:G:1018:EDO:H12	2.18	0.43
1:C:113:LYS:CB	1:C:115:MSE:HE3	2.49	0.43
1:D:218:ILE:HD11	1:D:243:LEU:HD11	2.00	0.43
1:E:171:ASP:OD1	1:E:173:THR:HB	2.18	0.43
2:B:1026:EDO:C2	3:B:1144:HOH:O	2.53	0.43
1:F:219:ARG:HG2	1:F:219:ARG:H	1.75	0.43
2:F:1028:EDO:H21	3:F:1149:HOH:O	2.19	0.42
1:C:131:ASP:OD2	1:E:226:ARG:NH1	2.45	0.42
1:G:249:LYS:HE3	1:H:178:TYR:OH	2.20	0.42
3:C:1085:HOH:O	1:D:242:ARG:NH1	2.43	0.42
1:D:150:ILE:CD1	1:G:162:GLN:HE21	2.31	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:160:ASN:OD1	1:H:160:ASN:C	2.57	0.42
1:H:191:GLU:OE2	2:H:1019:EDO:C2	2.67	0.42
1:F:244:GLU:HB2	1:H:248:THR:HG23	2.02	0.42
1:E:218:ILE:HD11	1:E:243:LEU:HD21	2.02	0.42
1:B:130:LEU:CB	1:B:132:MSE:HE3	2.45	0.42
1:F:130:LEU:HG	1:F:158:PHE:HE2	1.85	0.42
1:A:109:ARG:O	1:A:109:ARG:HG2	2.19	0.42
1:F:219:ARG:CG	3:F:1091:HOH:O	2.67	0.42
1:D:155:GLU:HG3	1:D:234:LEU:HD21	2.01	0.41
1:G:219:ARG:HD2	3:G:1074:HOH:O	2.20	0.41
1:B:118:LYS:NZ	3:B:1143:HOH:O	2.53	0.41
1:A:113:LYS:CD	3:A:1147:HOH:O	2.65	0.41
1:B:118:LYS:NZ	2:B:1016:EDO:H22	2.35	0.41
1:D:110:HIS:HE1	1:D:145:HIS:ND1	2.18	0.41
1:E:192:HIS:CD2	1:F:247:PHE:HE2	2.38	0.41
1:C:184:VAL:HG11	2:C:1017:EDO:C1	2.49	0.41
1:B:183:GLN:HE21	1:B:183:GLN:HB3	1.67	0.41
1:H:159:VAL:HG21	1:H:167:TYR:CE1	2.55	0.41
1:F:140:HIS:NE2	1:F:156:ASP:OD1	2.54	0.41
1:F:193:VAL:HB	1:F:219:ARG:HG3	2.01	0.40
1:F:218:ILE:HD11	1:F:243:LEU:HD21	2.03	0.40
1:E:214:PRO:HB3	1:F:214:PRO:HB3	2.03	0.40
1:F:209:ILE:HA	2:F:1028:EDO:H11	2.03	0.40
1:F:210:ASP:H	2:F:1028:EDO:H12	1.85	0.40
1:D:235:ILE:HD12	1:D:235:ILE:N	2.37	0.40
1:B:159:VAL:HG21	1:B:167:TYR:CE1	2.56	0.40
1:C:166:ASP:N	2:C:1017:EDO:O1	2.46	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	139/148 (94%)	137 (99%)	2 (1%)	0	100	100
1	B	140/148 (95%)	139 (99%)	1 (1%)	0	100	100
1	C	140/148 (95%)	139 (99%)	1 (1%)	0	100	100
1	D	139/148 (94%)	137 (99%)	2 (1%)	0	100	100
1	E	140/148 (95%)	138 (99%)	2 (1%)	0	100	100
1	F	139/148 (94%)	137 (99%)	2 (1%)	0	100	100
1	G	139/148 (94%)	138 (99%)	1 (1%)	0	100	100
1	H	140/148 (95%)	139 (99%)	1 (1%)	0	100	100
All	All	1116/1184 (94%)	1104 (99%)	12 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	120/122 (98%)	115 (96%)	5 (4%)	36	21
1	B	121/122 (99%)	116 (96%)	5 (4%)	37	22
1	C	121/122 (99%)	118 (98%)	3 (2%)	55	45
1	D	120/122 (98%)	114 (95%)	6 (5%)	30	14
1	E	121/122 (99%)	115 (95%)	6 (5%)	30	14
1	F	120/122 (98%)	116 (97%)	4 (3%)	45	32
1	G	120/122 (98%)	115 (96%)	5 (4%)	36	21
1	H	121/122 (99%)	118 (98%)	3 (2%)	55	45
All	All	964/976 (99%)	927 (96%)	37 (4%)	40	25

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	ARG
1	A	126	ARG

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Mol	Chain	Res	Type
1	A	133	ARG
1	A	137	ARG
1	A	240	ARG
1	B	109	ARG
1	B	125	GLU
1	B	136	GLN
1	B	137	ARG
1	B	221	ARG
1	C	111	THR
1	C	126	ARG
1	C	142	LEU
1	D	109	ARG
1	D	116	VAL
1	D	128	LEU
1	D	137	ARG
1	D	177	PRO
1	D	243	LEU
1	E	126	ARG
1	E	149	ASP
1	E	173	THR
1	E	227	GLN
1	E	234	LEU
1	E	237	PRO
1	F	128	LEU
1	F	206	LEU
1	F	219	ARG
1	F	249	LYS
1	G	132	MSE
1	G	133	ARG
1	G	202	ASP
1	G	213	GLU
1	G	234	LEU
1	H	147	GLU
1	H	183	GLN
1	H	226	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	175	GLN
1	A	183	GLN

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Mol	Chain	Res	Type
1	A	227	GLN
1	A	241	HIS
1	B	108	HIS
1	B	145	HIS
1	B	236	HIS
1	C	136	GLN
1	C	148	ASN
1	C	153	GLN
1	C	183	GLN
1	C	192	HIS
1	C	241	HIS
1	D	110	HIS
1	D	148	ASN
1	D	153	GLN
1	D	183	GLN
1	D	192	HIS
1	D	241	HIS
1	E	153	GLN
1	E	192	HIS
1	E	227	GLN
1	E	241	HIS
1	F	110	HIS
1	F	148	ASN
1	F	162	GLN
1	F	175	GLN
1	F	192	HIS
1	F	236	HIS
1	F	241	HIS
1	G	153	GLN
1	G	162	GLN
1	G	175	GLN
1	G	183	GLN
1	G	236	HIS
1	G	241	HIS
1	H	153	GLN
1	H	170	GLN
1	H	227	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

32 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	EDO	A	1001	-	3,3,3	0.55	0	2,2,2	0.33	0
2	EDO	A	1004	-	3,3,3	0.47	0	2,2,2	0.75	0
2	EDO	A	1009	-	3,3,3	0.64	0	2,2,2	0.30	0
2	EDO	A	1014	-	3,3,3	0.56	0	2,2,2	0.44	0
2	EDO	A	1024	-	3,3,3	0.52	0	2,2,2	0.18	0
2	EDO	B	1010	-	3,3,3	0.58	0	2,2,2	0.66	0
2	EDO	B	1016	-	3,3,3	0.78	0	2,2,2	0.08	0
2	EDO	B	1021	-	3,3,3	0.71	0	2,2,2	0.09	0
2	EDO	B	1026	-	3,3,3	0.83	0	2,2,2	0.25	0
2	EDO	B	1029	-	3,3,3	1.15	0	2,2,2	1.16	0
2	EDO	B	1032	-	3,3,3	0.63	0	2,2,2	0.23	0
2	EDO	C	1002	-	3,3,3	0.41	0	2,2,2	0.41	0
2	EDO	C	1012	-	3,3,3	0.44	0	2,2,2	0.42	0
2	EDO	C	1017	-	3,3,3	0.49	0	2,2,2	0.49	0
2	EDO	D	1003	-	3,3,3	0.64	0	2,2,2	0.21	0
2	EDO	D	1006	-	3,3,3	0.56	0	2,2,2	0.42	0
2	EDO	D	1013	-	3,3,3	0.67	0	2,2,2	0.64	0
2	EDO	E	1005	-	3,3,3	0.73	0	2,2,2	0.51	0
2	EDO	E	1011	-	3,3,3	0.51	0	2,2,2	0.37	0
2	EDO	E	1020	-	3,3,3	0.46	0	2,2,2	0.55	0
2	EDO	E	1027	-	3,3,3	0.59	0	2,2,2	0.91	0
2	EDO	E	1030	-	3,3,3	0.42	0	2,2,2	0.65	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	F	1007	-	3,3,3	0.42	0	2,2,2	0.70	0
2	EDO	F	1015	-	3,3,3	0.52	0	2,2,2	0.31	0
2	EDO	F	1028	-	3,3,3	0.52	0	2,2,2	0.88	0
2	EDO	G	1018	-	3,3,3	0.54	0	2,2,2	0.74	0
2	EDO	G	1023	-	3,3,3	0.43	0	2,2,2	0.58	0
2	EDO	H	1008	-	3,3,3	0.44	0	2,2,2	0.48	0
2	EDO	H	1019	-	3,3,3	0.78	0	2,2,2	0.46	0
2	EDO	H	1022	-	3,3,3	0.52	0	2,2,2	0.29	0
2	EDO	H	1025	-	3,3,3	0.59	0	2,2,2	0.07	0
2	EDO	H	1031	-	3,3,3	0.44	0	2,2,2	0.54	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	A	1001	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1004	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1009	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1014	-	-	0/1/1/1	0/0/0/0
2	EDO	A	1024	-	-	0/1/1/1	0/0/0/0
2	EDO	B	1010	-	-	0/1/1/1	0/0/0/0
2	EDO	B	1016	-	-	0/1/1/1	0/0/0/0
2	EDO	B	1021	-	-	0/1/1/1	0/0/0/0
2	EDO	B	1026	-	-	0/1/1/1	0/0/0/0
2	EDO	B	1029	-	-	0/1/1/1	0/0/0/0
2	EDO	B	1032	-	-	0/1/1/1	0/0/0/0
2	EDO	C	1002	-	-	0/1/1/1	0/0/0/0
2	EDO	C	1012	-	-	0/1/1/1	0/0/0/0
2	EDO	C	1017	-	-	0/1/1/1	0/0/0/0
2	EDO	D	1003	-	-	0/1/1/1	0/0/0/0
2	EDO	D	1006	-	-	0/1/1/1	0/0/0/0
2	EDO	D	1013	-	-	0/1/1/1	0/0/0/0
2	EDO	E	1005	-	-	0/1/1/1	0/0/0/0
2	EDO	E	1011	-	-	0/1/1/1	0/0/0/0
2	EDO	E	1020	-	-	0/1/1/1	0/0/0/0
2	EDO	E	1027	-	-	0/1/1/1	0/0/0/0
2	EDO	E	1030	-	-	0/1/1/1	0/0/0/0
2	EDO	F	1007	-	-	0/1/1/1	0/0/0/0
2	EDO	F	1015	-	-	0/1/1/1	0/0/0/0
2	EDO	F	1028	-	-	0/1/1/1	0/0/0/0
2	EDO	G	1018	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	G	1023	-	-	0/1/1/1	0/0/0/0
2	EDO	H	1008	-	-	0/1/1/1	0/0/0/0
2	EDO	H	1019	-	-	0/1/1/1	0/0/0/0
2	EDO	H	1022	-	-	0/1/1/1	0/0/0/0
2	EDO	H	1025	-	-	0/1/1/1	0/0/0/0
2	EDO	H	1031	-	-	0/1/1/1	0/0/0/0

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.

11 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1016	EDO	1	0
2	B	1026	EDO	4	0
2	B	1029	EDO	2	0
2	C	1002	EDO	1	0
2	C	1017	EDO	8	0
2	D	1003	EDO	1	0
2	D	1013	EDO	1	0
2	F	1028	EDO	5	0
2	G	1018	EDO	1	0
2	G	1023	EDO	1	0
2	H	1019	EDO	1	0

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	139/148 (93%)	0.26	6 (4%) 39 50	19, 26, 35, 47	0
1	B	140/148 (94%)	0.24	3 (2%) 67 75	17, 25, 35, 48	0
1	C	140/148 (94%)	0.54	9 (6%) 23 33	20, 27, 38, 60	0
1	D	139/148 (93%)	0.36	5 (3%) 46 57	21, 28, 39, 54	0
1	E	140/148 (94%)	0.52	3 (2%) 67 75	21, 27, 37, 64	0
1	F	139/148 (93%)	0.37	5 (3%) 46 57	22, 28, 38, 55	0
1	G	139/148 (93%)	0.26	5 (3%) 46 57	20, 25, 35, 53	0
1	H	140/148 (94%)	0.37	4 (2%) 55 65	19, 26, 36, 49	0
All	All	1116/1184 (94%)	0.37	40 (3%) 46 57	17, 26, 37, 64	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	110	HIS	5.4
1	G	109	ARG	5.0
1	A	109	ARG	4.9
1	E	108	HIS	4.5
1	E	110	HIS	4.3
1	H	108	HIS	4.1
1	C	109	ARG	4.0
1	A	110	HIS	3.5
1	F	247	PHE	3.4
1	F	109	ARG	3.4
1	C	108	HIS	3.4
1	D	109	ARG	3.3
1	D	110	HIS	3.2
1	G	110	HIS	3.0
1	B	174	LEU	3.0
1	C	148	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	148	ASN	2.8
1	C	110	HIS	2.8
1	H	109	ARG	2.8
1	C	247	PHE	2.8
1	F	158	PHE	2.8
1	A	149	ASP	2.7
1	A	148	ASN	2.6
1	A	210	ASP	2.5
1	C	150	ILE	2.5
1	G	186	PRO	2.5
1	E	202	ASP	2.4
1	C	149	ASP	2.4
1	G	210	ASP	2.3
1	B	109	ARG	2.2
1	H	148	ASN	2.2
1	B	108	HIS	2.2
1	A	111	THR	2.2
1	D	202	ASP	2.2
1	C	232	ALA	2.2
1	D	229	VAL	2.2
1	D	247	PHE	2.2
1	H	174	LEU	2.1
1	C	217	LEU	2.1
1	G	148	ASN	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	EDO	E	1005	4/4	0.76	0.30	9.80	43,43,46,48	0
2	EDO	C	1017	4/4	0.92	0.26	7.12	30,31,36,37	0
2	EDO	H	1025	4/4	0.81	0.21	5.72	36,37,39,42	0
2	EDO	D	1003	4/4	0.63	0.17	4.34	57,57,58,59	0
2	EDO	B	1029	4/4	0.64	0.22	3.86	30,33,34,44	0
2	EDO	F	1028	4/4	0.81	0.19	3.51	25,28,34,42	0
2	EDO	B	1010	4/4	0.82	0.16	2.46	44,45,46,46	0
2	EDO	B	1026	4/4	0.89	0.13	1.19	37,38,41,45	0
2	EDO	D	1013	4/4	0.82	0.15	0.81	29,30,30,34	0
2	EDO	A	1004	4/4	0.88	0.18	0.64	43,44,44,45	0
2	EDO	A	1001	4/4	0.86	0.16	0.12	47,49,51,51	0
2	EDO	E	1020	4/4	0.90	0.15	0.11	48,48,49,49	0
2	EDO	D	1006	4/4	0.85	0.12	-0.26	49,51,51,52	0
2	EDO	G	1018	4/4	0.93	0.11	-0.60	35,39,39,40	0
2	EDO	B	1032	4/4	0.83	0.12	-0.80	49,53,54,55	0
2	EDO	E	1011	4/4	0.92	0.10	-0.98	55,57,57,59	0
2	EDO	F	1015	4/4	0.90	0.12	-	50,51,52,53	0
2	EDO	A	1014	4/4	0.87	0.14	-	43,44,45,46	0
2	EDO	H	1008	4/4	0.90	0.17	-	46,46,47,47	0
2	EDO	F	1007	4/4	0.86	0.29	-	49,50,50,51	0
2	EDO	H	1022	4/4	0.90	0.18	-	54,55,55,56	0
2	EDO	C	1002	4/4	0.93	0.17	-	44,46,47,48	0
2	EDO	C	1012	4/4	0.80	0.15	-	44,45,45,46	0
2	EDO	G	1023	4/4	0.90	0.21	-	54,56,57,58	0
2	EDO	E	1027	4/4	0.70	0.22	-	36,37,38,41	0
2	EDO	B	1016	4/4	0.89	0.15	-	26,36,37,43	0
2	EDO	H	1031	4/4	0.84	0.18	-	45,49,49,52	0
2	EDO	E	1030	4/4	0.81	0.14	-	56,57,57,58	0
2	EDO	A	1009	4/4	0.84	0.14	-	38,38,42,43	0
2	EDO	H	1019	4/4	0.82	0.18	-	37,40,42,45	0
2	EDO	B	1021	4/4	0.72	0.22	-	43,45,47,48	0
2	EDO	A	1024	4/4	0.88	0.14	-	45,46,48,49	0

6.5 Other polymers ⓘ

There are no such residues in this entry.