



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 11:33 pm BST

PDB ID : 6NJU
Title : Mouse endonuclease G mutant H97A bound to A-DNA
Authors : Vander Zanden, C.M.; Ho, E.N.; Czarny, R.S.; Robertson, A.B.; Ho, P.S.
Deposited on : 2019-01-04
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

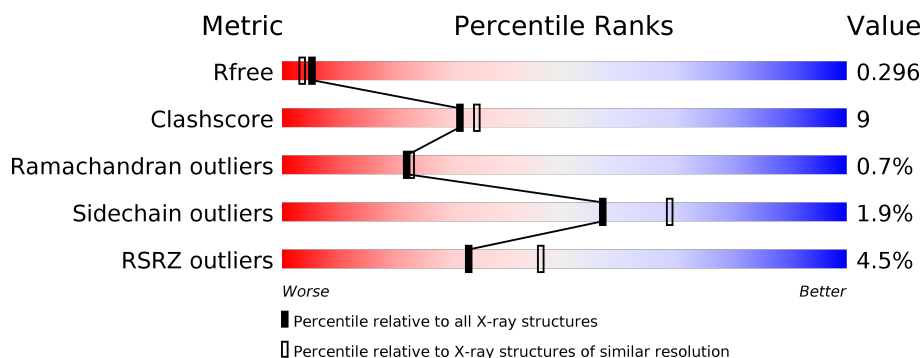
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	253	<div> <div>4%</div> <div>68% 19% • 12%</div> </div>
1	B	253	<div> <div>5%</div> <div>71% 18% • 10%</div> </div>
1	C	253	<div> <div>4%</div> <div>72% 16% • 11%</div> </div>
1	D	253	<div> <div>4%</div> <div>68% 22% • 9%</div> </div>
2	E	10	<div> <div>80% 20%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7736 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 Endonuclease G, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	3	0
			1803	1139	332	328	4			
1	B	228	Total	C	N	O	S	0	2	0
			1825	1150	337	334	4			
1	C	224	Total	C	N	O	S	0	4	0
			1809	1146	329	330	4			
1	D	229	Total	C	N	O	S	0	2	0
			1838	1159	339	336	4			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP O08600
A	97	ALA	HIS	engineered mutation	UNP O08600
B	1	SER	-	expression tag	UNP O08600
B	97	ALA	HIS	engineered mutation	UNP O08600
C	1	SER	-	expression tag	UNP O08600
C	97	ALA	HIS	engineered mutation	UNP O08600
D	1	SER	-	expression tag	UNP O08600
D	97	ALA	HIS	engineered mutation	UNP O08600

- Molecule 2 is a DNA chain called DNA (5'-D(CCGGCGCCGG)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	10	Total	C	N	O	P	0	0	0
			202	95	40	58	9			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		

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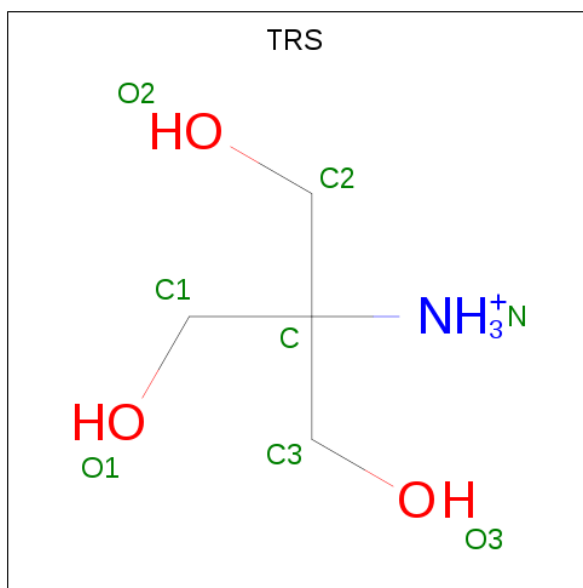
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	C	2	Total	Cl	0	0
			2	2		

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	E	1	Total	C	N	O	0	0
			8	4	1	3		

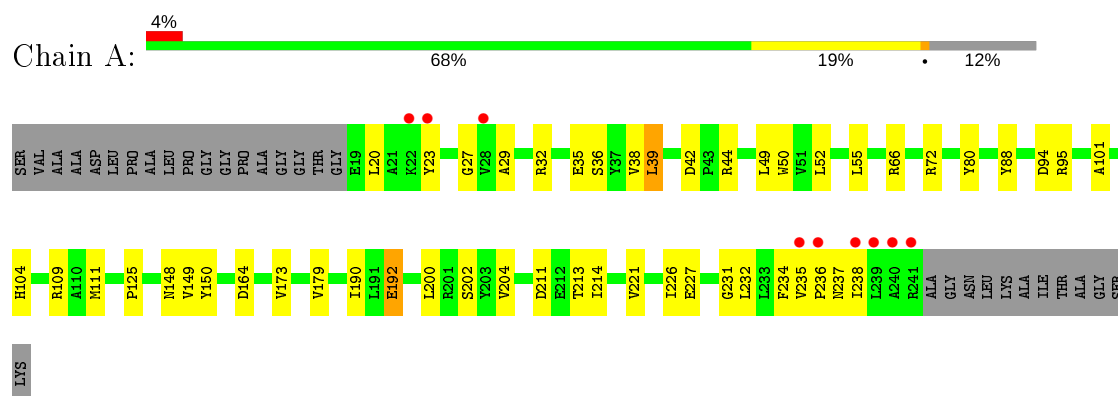
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	48	Total 48	O 48	0	0
6	B	47	Total 47	O 47	0	0
6	C	75	Total 75	O 75	0	0
6	D	62	Total 63	O 63	0	1
6	E	10	Total 10	O 10	0	0

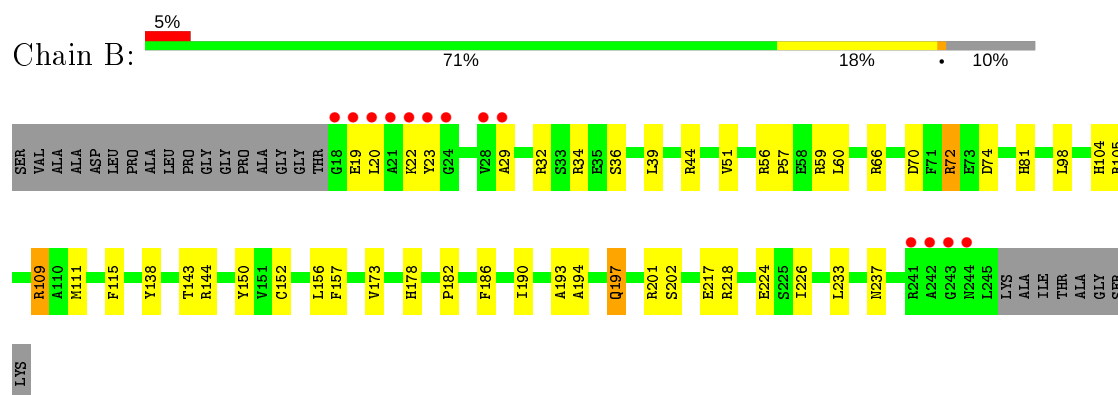
3 Residue-property plots [i](#)

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

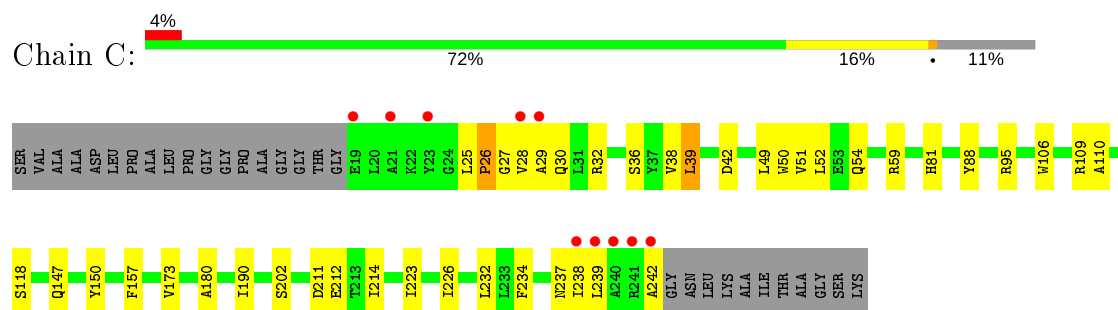
- Molecule 1: Endonuclease G, mitochondrial



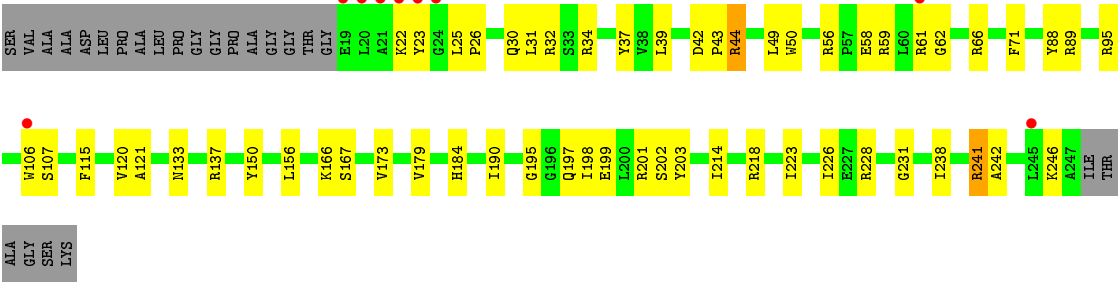
- Molecule 1: Endonuclease G, mitochondrial



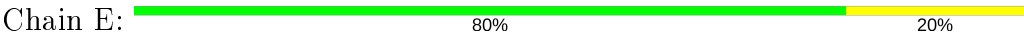
- Molecule 1: Endonuclease G, mitochondrial



- Molecule 1: Endonuclease G, mitochondrial



• Molecule 2: DNA (5'-D(CCGGCGCCGG)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	107.98Å 107.98Å 357.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	53.99 – 2.35 53.99 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (53.99-2.35) 100.0 (53.99-2.35)	Depositor EDS
R_{merge}	0.81	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.34Å)	Xtrriage
Refinement program	PHENIX 1.13-2998	Depositor
R, R_{free}	0.222 , 0.291 0.232 , 0.296	Depositor DCC
R_{free} test set	2540 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	23.2	Xtrriage
Anisotropy	0.279	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7736	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.33 % 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 1.7915e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TRS, MG, CL

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	0.45	0/1856	0.60	1/2522 (0.0%)
1	B	0.44	0/1875	0.58	0/2547
1	C	0.44	0/1865	0.63	1/2536 (0.0%)
1	D	0.49	0/1888	0.64	0/2564
2	E	0.90	0/226	0.94	1/347 (0.3%)
All	All	0.47	0/7710	0.63	3/10516 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	E	9	DG	O4'-C1'-N9	6.04	112.23	108.00
1	A	39	LEU	CA-CB-CG	5.37	127.66	115.30
1	C	39	LEU	CA-CB-CG	5.08	126.99	115.30

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	1803	0	1775	39	0
1	B	1825	0	1791	42	0
1	C	1809	0	1785	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1838	0	1807	41	0
2	E	202	0	112	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	C	2	0	0	0	0
4	D	1	0	0	0	0
5	E	8	0	12	1	0
6	A	48	0	0	0	0
6	B	47	0	0	0	0
6	C	75	0	0	2	0
6	D	63	0	0	3	0
6	E	10	0	0	0	0
All	All	7736	0	7282	133	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 9.

All (133) 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:138:TYR:OH	1:B:201:ARG:NH1	2.17	0.77
1:C:157:PHE:O	1:D:44[A]:ARG:NH2	2.17	0.77
1:A:29:ALA:HB2	1:B:29:ALA:HB2	1.66	0.76
1:B:109:ARG:HE	1:B:109:ARG:H	1.35	0.74
1:C:25:LEU:O	1:C:27:GLY:N	2.20	0.71
1:A:44:ARG:HH12	1:B:44:ARG:HH21	1.39	0.70
1:C:28:VAL:HG12	1:C:29:ALA:H	1.56	0.70
1:D:22:LYS:HD2	1:D:23:TYR:HE2	1.58	0.69
1:D:223:ILE:H	1:D:246:LYS:HE2	1.56	0.69
1:B:20:LEU:HD23	1:B:20:LEU:O	1.94	0.67
1:A:211:ASP:HB3	1:A:214:ILE:HD12	1.77	0.66
1:B:34[B]:ARG:NH1	1:B:74:ASP:OD2	2.29	0.65
1:B:193:ALA:HB3	1:B:197:GLN:HG2	1.80	0.64
1:D:166:LYS:NZ	6:D:401:HOH:O	2.31	0.63
1:D:89:ARG:NH2	6:D:402:HOH:O	2.31	0.63
1:A:148:ASN:HB2	1:A:192:GLU:HG2	1.81	0.62
1:A:80:TYR:OH	1:B:224:GLU:HG3	1.98	0.62
1:C:42:ASP:HB2	1:C:49[B]:LEU:HG	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ARG:HH22	1:B:72:ARG:HH11	1.47	0.61
1:D:223:ILE:HG22	1:D:246:LYS:HD3	1.83	0.60
1:C:42:ASP:OD2	1:D:44[B]:ARG:NH2	2.36	0.59
1:B:202[A]:SER:HB3	1:B:226:ILE:HD12	1.84	0.59
1:B:202[B]:SER:HB3	1:B:226:ILE:HD12	1.84	0.59
1:A:38:VAL:HG12	1:B:20:LEU:CD2	2.33	0.58
1:A:42:ASP:HB2	1:A:49[A]:LEU:HG	1.86	0.58
1:B:66:ARG:HH22	1:B:72:ARG:NH1	2.02	0.57
1:D:202:SER:HB2	1:D:226:ILE:HD12	1.87	0.57
1:D:59:ARG:HH21	1:D:106:TRP:HE1	1.52	0.57
1:B:150:TYR:HB2	1:B:190:ILE:HB	1.87	0.57
1:D:150:TYR:HB2	1:D:190:ILE:HB	1.87	0.56
1:D:22:LYS:HD2	1:D:23:TYR:CE2	2.40	0.56
1:A:148:ASN:ND2	1:A:192:GLU:OE2	2.39	0.55
1:C:212[B]:GLU:H	1:C:212[B]:GLU:CD	2.09	0.55
1:A:49[B]:LEU:HD12	1:A:232:LEU:HD21	1.88	0.54
1:A:235:VAL:HG23	1:A:236:PRO:HD3	1.89	0.54
1:A:42:ASP:HB2	1:A:49[B]:LEU:HD23	1.89	0.53
1:A:237:ASN:HB3	1:B:23:TYR:O	2.08	0.53
1:A:35:GLU:HG3	1:A:109:ARG:HG2	1.91	0.53
1:B:32:ARG:HB2	1:B:39:LEU:HB3	1.91	0.53
1:D:44[B]:ARG:HD2	6:D:439:HOH:O	2.09	0.53
1:D:22:LYS:C	1:D:23:TYR:HD2	2.13	0.53
1:C:39:LEU:HD12	1:C:51[B]:VAL:HG12	1.91	0.53
1:B:156:LEU:HD11	1:B:186:PHE:HB3	1.91	0.53
1:A:202:SER:HB2	1:A:226:ILE:HD12	1.90	0.52
1:B:56:ARG:O	1:B:60:LEU:HD13	2.10	0.52
6:C:410:HOH:O	1:D:44[A]:ARG:HD2	2.10	0.51
1:C:232:LEU:HB3	1:D:26:PRO:HB2	1.93	0.50
1:A:38:VAL:HG12	1:B:20:LEU:HD21	1.92	0.50
1:D:199:GLU:OE2	1:D:201:ARG:NH2	2.31	0.50
1:A:204:VAL:HG23	1:A:221:VAL:HG21	1.94	0.50
1:C:150:TYR:HB2	1:C:190:ILE:HB	1.94	0.50
1:D:133:ASN:O	1:D:137:ARG:HG3	2.11	0.49
1:A:235:VAL:CG2	1:A:236:PRO:HD3	2.42	0.49
1:C:202:SER:HB2	1:C:226:ILE:HD12	1.95	0.49
1:A:44:ARG:HH12	1:B:44:ARG:NH2	2.09	0.49
1:D:34:ARG:HG3	1:D:37:TYR:CE1	2.48	0.49
2:E:3:DG:N7	5:E:101:TRS:H11	2.28	0.49
1:D:34:ARG:HG3	1:D:37:TYR:CZ	2.48	0.48
1:A:211:ASP:OD1	1:A:213:THR:OG1	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:ARG:HA	1:B:111:MET:HE3	1.94	0.48
1:B:109:ARG:H	1:B:109:ARG:NE	2.07	0.47
1:A:36:SER:HB2	1:A:104:HIS:CE1	2.50	0.47
1:B:19:GLU:CG	1:B:22:LYS:HD3	2.45	0.47
1:C:54:GLN:NE2	1:C:150:TYR:OH	2.47	0.47
1:A:150:TYR:HB2	1:A:190:ILE:HB	1.96	0.47
1:C:59:ARG:HA	1:C:106:TRP:HB3	1.97	0.47
1:C:223:ILE:HD11	1:C:234:PHE:CD1	2.50	0.47
1:C:38[A]:VAL:HG12	1:C:52:LEU:HB3	1.96	0.46
1:C:39:LEU:HD12	1:C:51[A]:VAL:HG22	1.98	0.46
1:D:201:ARG:HG2	1:D:203:TYR:CE1	2.51	0.46
1:A:234:PHE:O	1:A:238:ILE:HG12	2.16	0.45
1:A:29:ALA:CB	1:B:29:ALA:HB2	2.40	0.45
1:C:211:ASP:HB3	1:C:214:ILE:HD12	1.98	0.45
1:D:214:ILE:HG23	1:D:218:ARG:HD3	1.98	0.45
1:D:195:GLY:N	1:D:197:GLN:OE1	2.50	0.45
1:C:81:HIS:CD2	1:D:231:GLY:HA2	2.51	0.45
1:B:66:ARG:HD3	1:B:115:PHE:CD1	2.52	0.44
1:B:59:ARG:HH12	1:B:104:HIS:CE1	2.34	0.44
1:D:42:ASP:HB2	1:D:49:LEU:HG	1.99	0.44
1:A:55:LEU:HB2	1:A:149:VAL:HB	1.98	0.44
1:C:173:VAL:HG12	1:C:180:ALA:HA	1.99	0.44
1:A:38:VAL:HG13	1:A:50:TRP:HZ3	1.83	0.44
1:D:37:TYR:HE1	1:D:39:LEU:HD23	1.82	0.44
1:B:36:SER:HA	1:B:59:ARG:HH22	1.84	0.43
1:C:237:ASN:ND2	1:D:25:LEU:O	2.51	0.43
1:C:238:ILE:HD13	1:C:238:ILE:HA	1.74	0.43
1:A:38:VAL:HG12	1:B:20:LEU:HD22	1.99	0.43
1:B:66:ARG:NH2	1:B:70:ASP:OD1	2.50	0.43
1:C:38[B]:VAL:HG22	1:C:52:LEU:HB3	2.01	0.43
1:B:233:LEU:HA	1:B:233:LEU:HD23	1.71	0.43
1:B:51:VAL:O	1:B:152:CYS:HA	2.18	0.43
1:C:38[A]:VAL:HG13	1:C:50:TRP:HZ3	1.82	0.43
1:D:32:ARG:HB2	1:D:39:LEU:HG	2.01	0.43
1:B:157:PHE:CE2	1:B:182:PRO:HA	2.54	0.43
1:C:239:LEU:HA	1:C:242:ALA:HB3	2.01	0.43
1:D:58:GLU:HG2	1:D:58:GLU:H	1.60	0.43
1:C:36:SER:HB3	1:C:110:ALA:HB1	2.00	0.42
1:D:71:PHE:HA	1:D:115:PHE:HB3	2.00	0.42
1:D:173:VAL:HA	1:D:179:VAL:O	2.20	0.42
1:A:101:ALA:O	1:A:111:MET:HE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:LEU:HD23	1:A:20:LEU:O	2.19	0.42
1:B:57:PRO:HD3	1:B:143:THR:HB	2.01	0.42
1:C:147:GLN:NE2	6:C:401:HOH:O	2.24	0.42
1:D:66:ARG:HD3	1:D:115:PHE:CD1	2.55	0.42
1:A:227:GLU:OE2	1:A:235:VAL:HG22	2.20	0.42
1:D:37:TYR:CE1	1:D:39:LEU:HD23	2.55	0.42
1:A:88:TYR:CD2	1:A:95:ARG:HG3	2.55	0.41
1:D:61:ARG:HD2	1:D:62:GLY:N	2.35	0.41
1:A:32:ARG:HB2	1:A:39:LEU:HB3	2.02	0.41
1:A:94:ASP:OD2	1:A:125:PRO:HA	2.21	0.41
1:A:80:TYR:CZ	1:B:224:GLU:HG3	2.55	0.41
1:D:156:LEU:HB2	1:D:184:HIS:HB2	2.02	0.41
1:D:88:TYR:CD2	1:D:95:ARG:HG3	2.55	0.41
1:B:173:VAL:HB	1:B:178:HIS:HB3	2.02	0.41
1:B:104:HIS:O	1:B:111:MET:HB2	2.20	0.41
1:C:88:TYR:CD2	1:C:95:ARG:HG3	2.56	0.41
1:A:192:GLU:OE1	1:B:23:TYR:OH	2.37	0.41
1:A:190:ILE:HG12	1:A:200:LEU:HD11	2.03	0.41
1:B:217:GLU:CD	1:B:217:GLU:H	2.24	0.41
1:A:231:GLY:HA2	1:B:81:HIS:CD2	2.56	0.41
1:A:38:VAL:CG1	1:A:52:LEU:HB3	2.51	0.40
1:C:39:LEU:HD11	1:C:118:SER:O	2.21	0.40
1:C:30:GLN:HB3	1:C:32:ARG:HH12	1.86	0.40
1:D:238:ILE:O	1:D:241:ARG:HB3	2.22	0.40
1:A:173:VAL:HA	1:A:179:VAL:O	2.20	0.40
1:B:98:LEU:HA	1:B:98:LEU:HD23	1.98	0.40
1:C:26:PRO:HG2	1:D:50:TRP:CD2	2.56	0.40
1:C:28:VAL:HG11	1:D:31:LEU:CD2	2.51	0.40
1:D:30:GLN:OE1	1:D:43:PRO:HG3	2.22	0.40
1:B:19:GLU:HG3	1:B:22:LYS:HB2	2.03	0.40
1:D:120:VAL:HG12	1:D:121:ALA:N	2.36	0.40
1:D:198:ILE:HG21	1:D:238:ILE:HG21	2.03	0.40
1:A:23:TYR:O	1:B:237:ASN:HB3	2.21	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	224/253 (88%)	213 (95%)	9 (4%)	2 (1%)	17	17
1	B	228/253 (90%)	217 (95%)	10 (4%)	1 (0%)	34	38
1	C	226/253 (89%)	219 (97%)	6 (3%)	1 (0%)	34	38
1	D	229/253 (90%)	216 (94%)	11 (5%)	2 (1%)	17	17
All	All	907/1012 (90%)	865 (95%)	36 (4%)	6 (1%)	22	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	241	ARG
1	D	242	ALA
1	A	164	ASP
1	C	26	PRO
1	B	194	ALA
1	A	27	GLY

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	190/203 (94%)	187 (98%)	3 (2%)	62	75
1	B	191/203 (94%)	186 (97%)	5 (3%)	46	56
1	C	191/203 (94%)	190 (100%)	1 (0%)	88	94
1	D	192/203 (95%)	186 (97%)	6 (3%)	40	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	764/812 (94%)	749 (98%)	15 (2%)	57	66

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	ARG
1	A	72	ARG
1	A	192	GLU
1	B	72	ARG
1	B	109	ARG
1	B	144	ARG
1	B	197	GLN
1	B	218	ARG
1	C	109	ARG
1	D	44[A]	ARG
1	D	44[B]	ARG
1	D	56	ARG
1	D	107	SER
1	D	167	SER
1	D	228	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

Of 9 ligands modelled in this entry, 8 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	TRS	E	101	-	7,7,7	0.48	0	9,9,9	0.95	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
5	TRS	E	101	-	-	9/9/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	101	TRS	N-C-C1-O1
5	E	101	TRS	C2-C-C1-O1
5	E	101	TRS	C3-C-C1-O1
5	E	101	TRS	C1-C-C2-O2
5	E	101	TRS	C3-C-C2-O2
5	E	101	TRS	N-C-C3-O3
5	E	101	TRS	C1-C-C3-O3
5	E	101	TRS	C2-C-C3-O3
5	E	101	TRS	N-C-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	101	TRS	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	223/253 (88%)	0.03	9 (4%)	38	51	23, 34, 61, 86	0
1	B	228/253 (90%)	0.13	13 (5%)	23	34	23, 35, 71, 91	0
1	C	224/253 (88%)	0.11	10 (4%)	33	46	17, 29, 67, 86	0
1	D	229/253 (90%)	0.12	9 (3%)	39	52	18, 28, 63, 86	0
2	E	10/10 (100%)	-0.49	0	100	100	28, 35, 41, 49	0
All	All	914/1022 (89%)	0.09	41 (4%)	33	46	17, 32, 66, 91	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	240	ALA	7.1
1	B	23	TYR	7.1
1	D	23	TYR	7.0
1	C	240	ALA	6.1
1	C	19	GLU	5.9
1	D	19	GLU	5.5
1	D	21	ALA	5.2
1	B	21	ALA	5.1
1	A	239	LEU	4.9
1	C	29	ALA	4.8
1	C	21	ALA	4.7
1	B	242	ALA	4.4
1	B	20	LEU	4.3
1	C	23	TYR	4.3
1	B	28	VAL	4.0
1	C	242	ALA	4.0
1	B	244	ASN	3.9
1	D	20	LEU	3.9
1	D	22	LYS	3.7
1	C	238	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	241	ARG	3.6
1	C	28	VAL	3.5
1	C	239	LEU	3.4
1	B	243	GLY	3.4
1	D	24	GLY	3.4
1	B	22	LYS	3.3
1	B	19	GLU	3.1
1	A	241	ARG	3.1
1	B	18	GLY	3.0
1	A	235	VAL	2.9
1	A	28	VAL	2.8
1	B	29	ALA	2.7
1	A	23	TYR	2.7
1	B	24	GLY	2.7
1	D	106	TRP	2.6
1	B	241	ARG	2.5
1	A	236	PRO	2.5
1	D	61	ARG	2.3
1	D	245	LEU	2.3
1	A	22	LYS	2.2
1	A	238	ILE	2.1

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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CL	C	302	1/1	0.82	0.15	57,57,57,57	0
4	CL	A	302	1/1	0.87	0.12	54,54,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CL	D	302	1/1	0.92	0.13	51,51,51,51	0
5	TRS	E	101	8/8	0.94	0.12	27,33,41,45	0
4	CL	C	303	1/1	0.95	0.07	53,53,53,53	0
3	MG	C	301	1/1	0.95	0.05	35,35,35,35	0
3	MG	A	301	1/1	0.96	0.06	33,33,33,33	0
3	MG	D	301	1/1	0.97	0.09	34,34,34,34	0
3	MG	B	301	1/1	0.99	0.07	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.