



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 02:40 pm BST

PDB ID : 5ZJF  
Title : LDHA-MA  
Authors : Han, C.W.; Jang, S.B.  
Deposited on : 2018-03-20  
Resolution : 2.60 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](mailto: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.

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

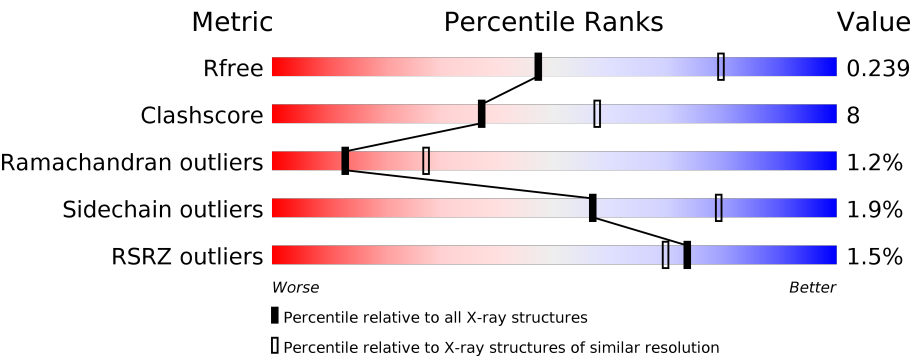
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:  
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	337	<div><div>%</div><div><div></div><div></div><div></div><div></div></div><div>81%16%..</div></div>
1	B	337	<div><div>%</div><div><div></div><div></div><div></div><div></div></div><div>80%18%..</div></div>
1	C	337	<div><div>%</div><div><div></div><div></div><div></div><div></div></div><div>82%15%..</div></div>
1	D	337	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>81%17%..</div></div>
1	E	337	<div><div>3%</div><div><div></div><div></div><div></div><div></div></div><div>80%17%..</div></div>
1	F	337	<div><div>%</div><div><div></div><div></div><div></div><div></div></div><div>83%14%..</div></div>

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Mol	Chain	Length	Quality of chain
1	G	337	
1	H	337	
1	I	337	
1	J	337	
1	K	337	
1	L	337	

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	9G9	C	401	-	-	X	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 31211 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 L-lactate dehydrogenase A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	B	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	C	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	D	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	E	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	F	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	G	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	H	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	I	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	J	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	K	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			
1	L	331	Total	C	N	O	S	0	0	0
			2568	1639	439	477	13			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP P00338
A	-4	HIS	-	expression tag	UNP P00338
A	-3	HIS	-	expression tag	UNP P00338
A	-2	HIS	-	expression tag	UNP P00338
A	-1	HIS	-	expression tag	UNP P00338

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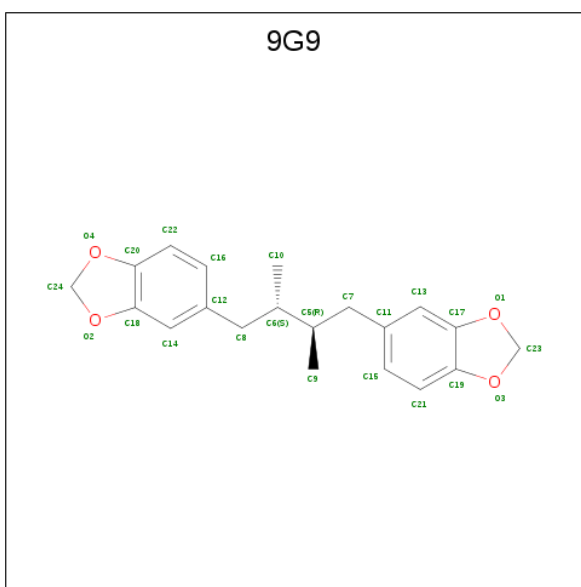
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	expression tag	UNP P00338
B	-5	HIS	-	expression tag	UNP P00338
B	-4	HIS	-	expression tag	UNP P00338
B	-3	HIS	-	expression tag	UNP P00338
B	-2	HIS	-	expression tag	UNP P00338
B	-1	HIS	-	expression tag	UNP P00338
B	0	HIS	-	expression tag	UNP P00338
C	-5	HIS	-	expression tag	UNP P00338
C	-4	HIS	-	expression tag	UNP P00338
C	-3	HIS	-	expression tag	UNP P00338
C	-2	HIS	-	expression tag	UNP P00338
C	-1	HIS	-	expression tag	UNP P00338
C	0	HIS	-	expression tag	UNP P00338
D	-5	HIS	-	expression tag	UNP P00338
D	-4	HIS	-	expression tag	UNP P00338
D	-3	HIS	-	expression tag	UNP P00338
D	-2	HIS	-	expression tag	UNP P00338
D	-1	HIS	-	expression tag	UNP P00338
D	0	HIS	-	expression tag	UNP P00338
E	-5	HIS	-	expression tag	UNP P00338
E	-4	HIS	-	expression tag	UNP P00338
E	-3	HIS	-	expression tag	UNP P00338
E	-2	HIS	-	expression tag	UNP P00338
E	-1	HIS	-	expression tag	UNP P00338
E	0	HIS	-	expression tag	UNP P00338
F	-5	HIS	-	expression tag	UNP P00338
F	-4	HIS	-	expression tag	UNP P00338
F	-3	HIS	-	expression tag	UNP P00338
F	-2	HIS	-	expression tag	UNP P00338
F	-1	HIS	-	expression tag	UNP P00338
F	0	HIS	-	expression tag	UNP P00338
G	-5	HIS	-	expression tag	UNP P00338
G	-4	HIS	-	expression tag	UNP P00338
G	-3	HIS	-	expression tag	UNP P00338
G	-2	HIS	-	expression tag	UNP P00338
G	-1	HIS	-	expression tag	UNP P00338
G	0	HIS	-	expression tag	UNP P00338
H	-5	HIS	-	expression tag	UNP P00338
H	-4	HIS	-	expression tag	UNP P00338
H	-3	HIS	-	expression tag	UNP P00338
H	-2	HIS	-	expression tag	UNP P00338
H	-1	HIS	-	expression tag	UNP P00338

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Chain	Residue	Modelled	Actual	Comment	Reference
H	0	HIS	-	expression tag	UNP P00338
I	-5	HIS	-	expression tag	UNP P00338
I	-4	HIS	-	expression tag	UNP P00338
I	-3	HIS	-	expression tag	UNP P00338
I	-2	HIS	-	expression tag	UNP P00338
I	-1	HIS	-	expression tag	UNP P00338
I	0	HIS	-	expression tag	UNP P00338
J	-5	HIS	-	expression tag	UNP P00338
J	-4	HIS	-	expression tag	UNP P00338
J	-3	HIS	-	expression tag	UNP P00338
J	-2	HIS	-	expression tag	UNP P00338
J	-1	HIS	-	expression tag	UNP P00338
J	0	HIS	-	expression tag	UNP P00338
K	-5	HIS	-	expression tag	UNP P00338
K	-4	HIS	-	expression tag	UNP P00338
K	-3	HIS	-	expression tag	UNP P00338
K	-2	HIS	-	expression tag	UNP P00338
K	-1	HIS	-	expression tag	UNP P00338
K	0	HIS	-	expression tag	UNP P00338
L	-5	HIS	-	expression tag	UNP P00338
L	-4	HIS	-	expression tag	UNP P00338
L	-3	HIS	-	expression tag	UNP P00338
L	-2	HIS	-	expression tag	UNP P00338
L	-1	HIS	-	expression tag	UNP P00338
L	0	HIS	-	expression tag	UNP P00338

- Molecule 2 is 5,5'-[(2R,3S)-2,3-dimethylbutane-1,4-diyl]bis(2H-1,3-benzodioxole) (three-letter code: 9G9) (formula: C<sub>20</sub>H<sub>22</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			24	20	4		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	24	Total O 24 24	0	0
3	B	17	Total O 17 17	0	0
3	C	37	Total O 37 37	0	0
3	D	18	Total O 18 18	0	0
3	E	26	Total O 26 26	0	0
3	F	36	Total O 36 36	0	0
3	G	22	Total O 22 22	0	0
3	H	17	Total O 17 17	0	0
3	I	61	Total O 61 61	0	0
3	J	53	Total O 53 53	0	0
3	K	41	Total O 41 41	0	0

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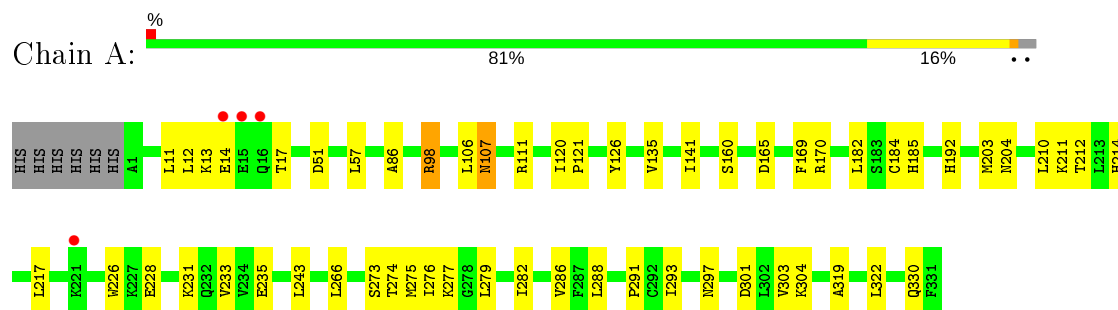
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	19	Total	O	0	0
			19	19		



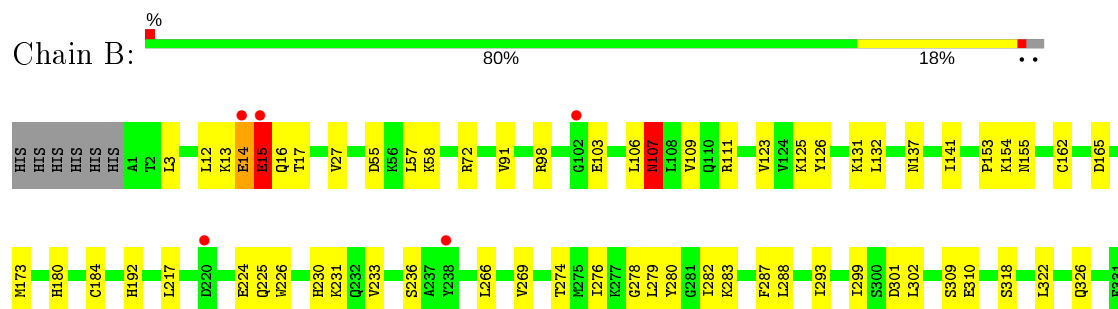
### 3 Residue-property plots [i](#)

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.

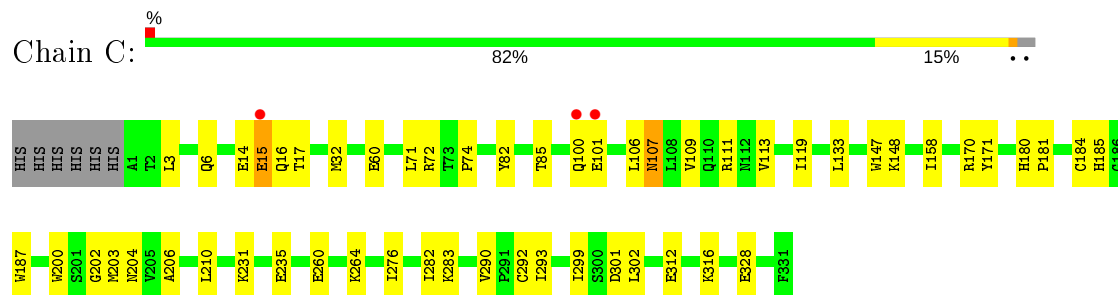
- Molecule 1: L-lactate dehydrogenase A chain



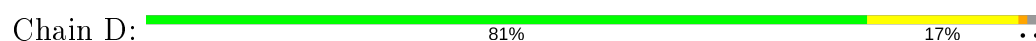
- Molecule 1: L-lactate dehydrogenase A chain

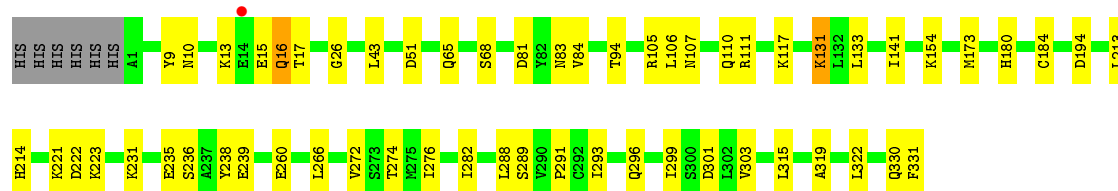


- Molecule 1: L-lactate dehydrogenase A chain

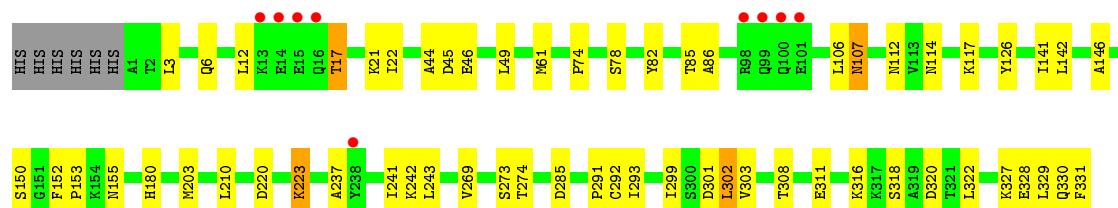
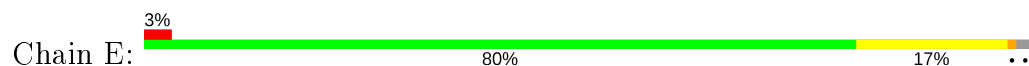


- Molecule 1: L-lactate dehydrogenase A chain

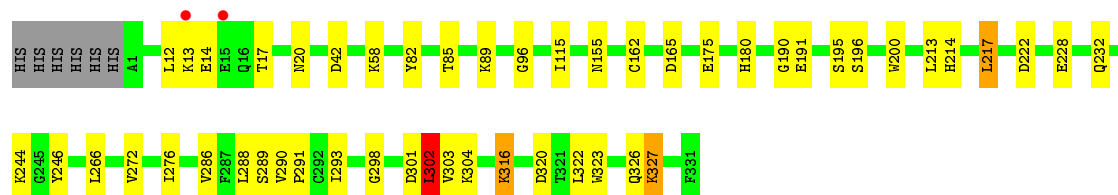
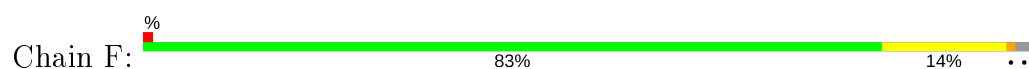




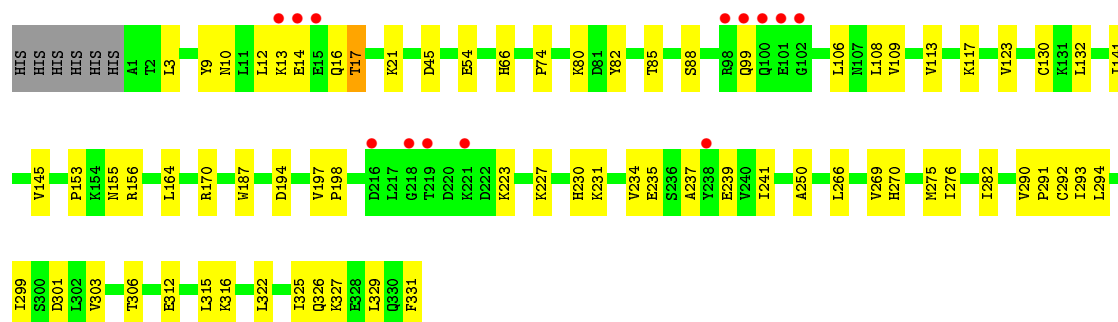
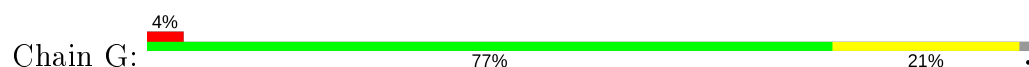
• Molecule 1: L-lactate dehydrogenase A chain



• Molecule 1: L-lactate dehydrogenase A chain

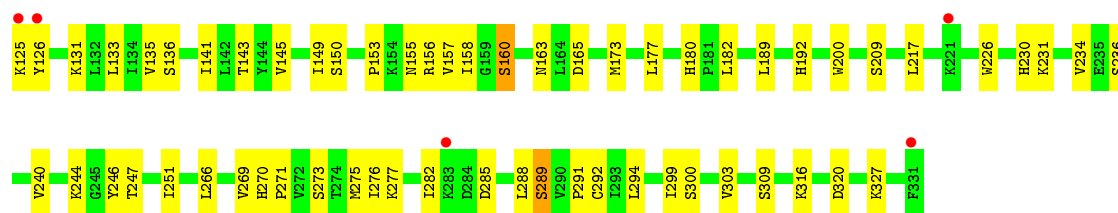


• Molecule 1: L-lactate dehydrogenase A chain



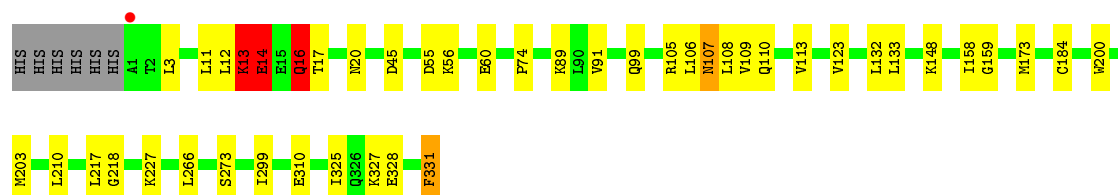
• Molecule 1: L-lactate dehydrogenase A chain





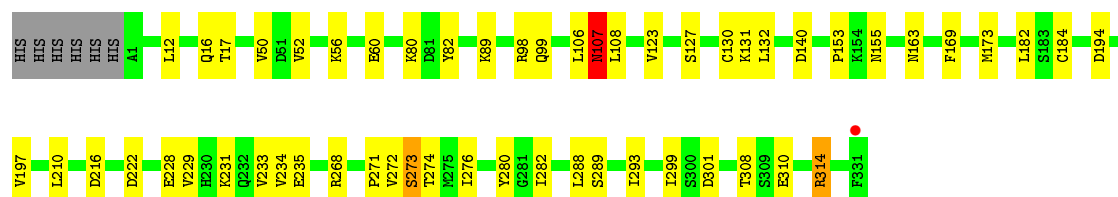
• Molecule 1: L-lactate dehydrogenase A chain

Chain I: 85% 12% ..



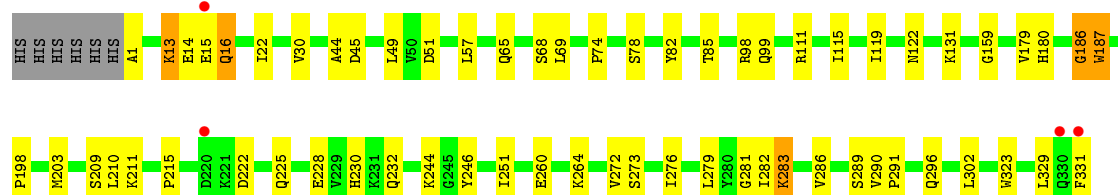
• Molecule 1: L-lactate dehydrogenase A chain

Chain J: 82% 15% ..



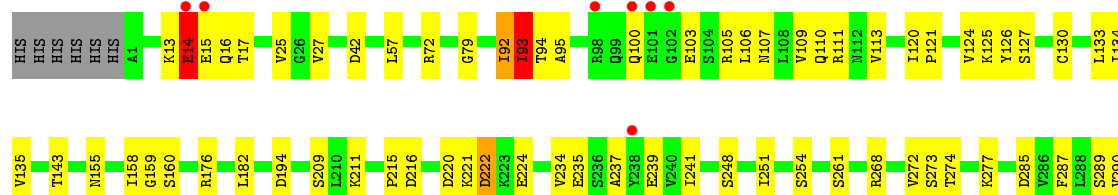
• Molecule 1: L-lactate dehydrogenase A chain

Chain K: 80% 17% ..



• Molecule 1: L-lactate dehydrogenase A chain

Chain L: 75% 22% ..





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	134.23Å 141.50Å 144.08Å 90.00° 110.18° 90.00°	Depositor
Resolution (Å)	37.10 – 2.60 37.10 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.4 (37.10-2.60) 99.4 (37.10-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.46 (at 2.61Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, $R_{free}$	0.181 , 0.239 0.181 , 0.239	Depositor DCC
$R_{free}$ test set	7532 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	50.8	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 45.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	31211	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 9G9

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	0.44	0/2612	0.60	0/3532
1	B	0.42	0/2612	0.59	1/3532 (0.0%)
1	C	0.44	0/2612	0.62	0/3532
1	D	0.42	0/2612	0.60	0/3532
1	E	0.39	0/2612	0.56	0/3532
1	F	0.43	0/2612	0.59	1/3532 (0.0%)
1	G	0.40	0/2612	0.58	0/3532
1	H	0.45	0/2612	0.62	1/3532 (0.0%)
1	I	0.45	0/2612	0.64	0/3532
1	J	0.45	0/2612	0.59	0/3532
1	K	0.47	0/2612	0.63	0/3532
1	L	0.41	0/2612	0.59	1/3532 (0.0%)
All	All	0.43	0/31344	0.60	4/42384 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	2
1	H	0	4
1	I	0	2
1	L	0	2
All	All	0	11

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	302	LEU	CA-CB-CG	6.62	130.53	115.30
1	B	302	LEU	CA-CB-CG	5.90	128.87	115.30
1	H	93	ILE	N-CA-C	5.40	125.58	111.00
1	L	93	ILE	CG1-CB-CG2	-5.12	100.13	111.40

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	14	GLU	Peptide
1	C	14	GLU	Peptide
1	C	15	GLU	Peptide
1	H	13	LYS	Peptide
1	H	15	GLU	Peptide
1	H	92	ILE	Peptide
1	H	93	ILE	Peptide
1	I	13	LYS	Peptide
1	I	16	GLN	Peptide
1	L	13	LYS	Peptide
1	L	92	ILE	Peptide

## 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	2568	0	2656	48	0
1	B	2568	0	2656	36	0
1	C	2568	0	2656	39	0
1	D	2568	0	2656	33	0
1	E	2568	0	2656	45	0
1	F	2568	0	2656	33	0
1	G	2568	0	2656	47	0
1	H	2568	0	2656	86	0
1	I	2568	0	2656	26	0
1	J	2568	0	2656	33	0
1	K	2568	0	2656	35	0
1	L	2568	0	2656	65	0
2	C	24	0	0	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	24	0	0	4	0
3	B	17	0	0	1	0
3	C	37	0	0	0	0
3	D	18	0	0	0	0
3	E	26	0	0	1	0
3	F	36	0	0	1	0
3	G	22	0	0	5	0
3	H	17	0	0	0	0
3	I	61	0	0	0	0
3	J	53	0	0	0	0
3	K	41	0	0	1	0
3	L	19	0	0	0	0
All	All	31211	0	31872	490	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 8.

All (490) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:93:ILE:HD12	1:H:135:VAL:N	1.14	1.43
1:H:93:ILE:HD11	1:H:136:SER:N	1.41	1.32
1:H:93:ILE:CD1	1:H:135:VAL:H	1.44	1.30
1:H:93:ILE:CD1	1:H:135:VAL:N	2.09	1.10
1:H:93:ILE:CD1	1:H:136:SER:N	2.15	1.08
1:F:316:LYS:NZ	1:F:320:ASP:OD1	1.92	1.02
1:H:93:ILE:HD12	1:H:135:VAL:CA	1.92	0.99
1:H:93:ILE:CD1	1:H:136:SER:H	1.72	0.96
1:A:98:ARG:NH1	1:A:111:ARG:HH12	1.66	0.91
1:A:98:ARG:HD3	1:A:98:ARG:H	1.32	0.90
1:A:98:ARG:HH12	1:A:111:ARG:HH12	0.94	0.89
1:L:93:ILE:HD12	1:L:135:VAL:N	1.87	0.88
1:L:93:ILE:CD1	1:L:135:VAL:HB	2.05	0.87
1:A:98:ARG:HH12	1:A:111:ARG:NH1	1.76	0.83
1:C:148:LYS:HE2	1:C:148:LYS:HA	1.57	0.83
1:H:93:ILE:CD1	1:H:135:VAL:CA	2.52	0.82
1:L:93:ILE:HD12	1:L:135:VAL:HB	1.60	0.82
1:H:89:LYS:HE2	1:H:131:LYS:NZ	1.93	0.82
1:H:92:ILE:HG22	1:H:93:ILE:HG22	1.61	0.82
1:H:93:ILE:HD11	1:H:136:SER:H	1.03	0.82
1:L:92:ILE:HA	1:L:93:ILE:HG23	1.60	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:99:GLN:HG2	1:G:108:LEU:HD22	1.61	0.81
1:D:276:ILE:HD11	1:D:288:LEU:HD12	1.62	0.80
1:C:203:MET:HG2	1:C:210:LEU:HD22	1.62	0.79
1:H:93:ILE:HD13	1:H:135:VAL:HB	1.64	0.79
1:C:185:HIS:HD1	2:C:401:9G9:C16	1.96	0.79
1:J:106:LEU:O	1:J:107:ASN:HB2	1.81	0.78
1:E:293:ILE:HD12	1:E:301:ASP:HB2	1.67	0.77
1:F:316:LYS:NZ	1:F:320:ASP:CG	2.38	0.76
1:A:204:ASN:HD21	2:C:401:9G9:C9	1.98	0.76
1:C:185:HIS:ND1	2:C:401:9G9:C16	2.50	0.75
1:H:276:ILE:HD11	1:H:288:LEU:HD12	1.68	0.74
1:G:227:LYS:HD3	1:G:227:LYS:O	1.85	0.74
1:L:293:ILE:HD12	1:L:301:ASP:HB2	1.69	0.74
1:F:190:GLY:HA2	1:F:288:LEU:HD23	1.68	0.73
1:C:204:ASN:HD21	2:C:401:9G9:C10	2.02	0.72
1:F:82:TYR:O	1:F:85:THR:HB	1.88	0.72
1:A:275:MET:HE3	1:A:277:LYS:H	1.54	0.71
1:A:297:ASN:ND2	3:A:401:HOH:O	2.18	0.71
1:H:89:LYS:HE2	1:H:131:LYS:HZ3	1.54	0.70
1:E:316:LYS:NZ	1:E:320:ASP:OD1	2.24	0.70
1:G:14:GLU:OE1	3:G:401:HOH:O	2.09	0.70
1:C:204:ASN:HA	1:C:210:LEU:HD13	1.74	0.69
1:F:316:LYS:HZ2	1:F:320:ASP:CG	1.95	0.69
1:K:260:GLU:HG3	1:K:264:LYS:HD2	1.74	0.69
1:D:131:LYS:NZ	1:D:296:GLN:O	2.23	0.69
1:L:312:GLU:HG2	1:L:316:LYS:HE3	1.76	0.68
1:L:194:ASP:HA	1:L:234:VAL:HG11	1.75	0.68
1:L:93:ILE:HG13	1:L:135:VAL:HG12	1.74	0.68
1:H:93:ILE:CD1	1:H:135:VAL:C	2.61	0.68
1:B:103:GLU:OE2	1:B:111:ARG:NH1	2.26	0.68
1:E:223:LYS:HD3	1:E:223:LYS:H	1.59	0.68
1:G:14:GLU:HB2	1:G:16:GLN:HE22	1.59	0.67
1:H:93:ILE:HG23	1:H:94:THR:N	2.10	0.67
1:H:12:LEU:HG	1:H:14:GLU:HB2	1.76	0.67
1:J:140:ASP:OD1	1:J:273:SER:OG	2.13	0.66
1:F:290:VAL:HG21	1:F:302:LEU:HD12	1.78	0.66
1:J:276:ILE:HD11	1:J:288:LEU:HD12	1.76	0.65
1:H:12:LEU:HA	1:H:14:GLU:OE2	1.95	0.65
1:K:69:LEU:HD12	1:L:182:LEU:HD12	1.78	0.65
1:G:231:LYS:NZ	3:G:403:HOH:O	2.30	0.64
1:C:185:HIS:NE2	2:C:401:9G9:C11	2.60	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:ARG:HD2	1:A:184:CYS:O	1.98	0.64
1:F:42:ASP:OD2	3:F:401:HOH:O	2.14	0.63
1:K:159:GLY:HA3	1:K:273:SER:HB2	1.81	0.63
1:A:98:ARG:N	1:A:98:ARG:HD3	2.09	0.62
1:B:276:ILE:HD11	1:B:288:LEU:HG	1.80	0.62
1:I:106:LEU:O	1:I:107:ASN:HB2	1.99	0.62
1:H:135:VAL:HG11	1:H:251:ILE:HD11	1.80	0.62
1:J:127:SER:HB3	1:J:130:CYS:HB3	1.81	0.62
1:J:293:ILE:HD12	1:J:301:ASP:HB2	1.80	0.62
1:K:82:TYR:O	1:K:85:THR:HB	1.99	0.62
1:L:93:ILE:HD12	1:L:135:VAL:CB	2.30	0.62
1:H:93:ILE:HD11	1:H:136:SER:CA	2.26	0.62
1:B:280:TYR:HB2	1:B:282:ILE:HD12	1.83	0.61
1:L:92:ILE:HG23	1:L:93:ILE:HG12	1.82	0.60
1:C:82:TYR:O	1:C:85:THR:HB	2.01	0.60
1:D:117:LYS:NZ	1:D:331:PHE:OXT	2.32	0.60
1:F:293:ILE:HD12	1:F:301:ASP:HB2	1.85	0.59
1:F:96:GLY:HA2	1:F:115:ILE:HD13	1.84	0.59
1:H:51:ASP:OD1	1:H:52:VAL:N	2.33	0.59
1:L:158:ILE:HG23	1:L:299:ILE:HD11	1.85	0.59
1:G:109:VAL:O	1:G:113:VAL:HG23	2.03	0.58
1:H:93:ILE:HG23	1:H:94:THR:HG23	1.83	0.58
1:L:92:ILE:CA	1:L:93:ILE:HG23	2.30	0.58
1:A:185:HIS:O	1:A:203:MET:HA	2.02	0.58
1:F:180:HIS:HB2	1:H:266:LEU:O	2.02	0.58
1:L:292:CYS:HB3	1:L:299:ILE:HG23	1.85	0.58
1:G:293:ILE:HD12	1:G:301:ASP:HB2	1.84	0.58
1:G:82:TYR:O	1:G:85:THR:HB	2.04	0.58
1:H:135:VAL:HA	1:H:160:SER:HB3	1.85	0.58
1:I:203:MET:HE1	1:I:210:LEU:HD22	1.85	0.58
1:A:276:ILE:HD11	1:A:288:LEU:HD12	1.85	0.58
1:C:231:LYS:O	1:C:235:GLU:HG2	2.04	0.58
1:A:12:LEU:O	1:A:13:LYS:HD2	2.04	0.57
1:E:22:ILE:HD12	1:E:44:ALA:HB2	1.86	0.57
1:K:198:PRO:HD3	1:K:230:HIS:CE1	2.39	0.57
1:L:120:ILE:O	1:L:124:VAL:HG23	2.04	0.57
1:L:14:GLU:OE2	1:L:16:GLN:N	2.37	0.57
1:L:93:ILE:HG21	1:L:133:LEU:C	2.24	0.57
1:G:14:GLU:HA	1:G:16:GLN:OE1	2.05	0.57
1:I:3:LEU:HD21	1:J:210:LEU:HG	1.85	0.57
1:L:93:ILE:HD11	1:L:135:VAL:HB	1.83	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:145:VAL:HG13	1:G:331:PHE:HE1	1.69	0.57
1:I:12:LEU:HD22	1:I:13:LYS:HG2	1.84	0.56
1:C:147:TRP:CZ3	1:C:148:LYS:HE3	2.39	0.56
1:J:99:GLN:HA	1:J:108:LEU:HD13	1.86	0.56
1:C:200:TRP:CE3	1:C:203:MET:HE1	2.41	0.56
1:G:326:GLN:HA	1:G:329:LEU:HD12	1.88	0.56
1:H:96:GLY:HA2	1:H:115:ILE:HD13	1.88	0.56
1:H:93:ILE:HD13	1:H:135:VAL:CB	2.35	0.56
1:H:93:ILE:HG12	1:H:94:THR:HA	1.88	0.56
1:D:293:ILE:HD12	1:D:301:ASP:HB2	1.88	0.56
1:D:231:LYS:O	1:D:235:GLU:HG2	2.06	0.56
1:J:194:ASP:HA	1:J:234:VAL:HG11	1.88	0.56
1:G:291:PRO:HB2	1:G:303:VAL:HB	1.86	0.55
1:H:269:VAL:HA	1:H:292:CYS:O	2.06	0.55
1:L:92:ILE:CG2	1:L:93:ILE:HG12	2.36	0.55
1:E:45:ASP:O	1:E:74:PRO:HD2	2.05	0.55
1:F:316:LYS:HZ1	1:F:320:ASP:CG	2.07	0.55
1:H:292:CYS:HB3	1:H:299:ILE:HG23	1.88	0.55
1:L:107:ASN:HA	1:L:110:GLN:HB2	1.88	0.55
1:B:106:LEU:O	1:B:107:ASN:HB2	2.05	0.55
1:D:141:ILE:HG13	1:D:322:LEU:HD22	1.88	0.55
1:J:276:ILE:HD12	1:J:282:ILE:HD13	1.88	0.55
1:E:21:LYS:HG3	1:E:46:GLU:HG2	1.89	0.55
1:B:14:GLU:HG3	1:B:15:GLU:HG3	1.88	0.55
1:C:276:ILE:HD13	1:C:282:ILE:HD12	1.87	0.55
1:G:327:LYS:NZ	3:G:405:HOH:O	2.39	0.55
1:H:200:TRP:HB3	1:H:217:LEU:HD23	1.89	0.55
1:L:321:THR:O	1:L:325:ILE:HG13	2.07	0.55
1:K:82:TYR:OH	1:K:119:ILE:HG12	2.05	0.54
1:L:93:ILE:HG13	1:L:94:THR:N	2.22	0.54
1:I:107:ASN:HA	1:I:110:GLN:HB3	1.89	0.54
1:A:13:LYS:HA	1:A:13:LYS:HE3	1.89	0.54
1:F:316:LYS:NZ	1:F:320:ASP:OD2	2.41	0.54
1:B:287:PHE:O	1:B:288:LEU:HD23	2.08	0.54
1:G:231:LYS:O	1:G:235:GLU:HG2	2.08	0.54
1:I:327:LYS:HE2	1:I:328:GLU:OE2	2.08	0.54
1:I:45:ASP:O	1:I:74:PRO:HD2	2.06	0.54
1:E:17:THR:N	3:E:401:HOH:O	2.32	0.54
1:H:291:PRO:HG2	1:H:303:VAL:HB	1.90	0.54
1:B:310:GLU:CD	1:B:310:GLU:H	2.11	0.54
1:E:243:LEU:HD13	1:F:58:LYS:HG2	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:302:LEU:HD23	1:G:9:TYR:HB3	1.90	0.53
1:E:327:LYS:HG3	1:E:328:GLU:HG3	1.90	0.53
1:E:146:ALA:O	1:E:150:SER:HB3	2.08	0.53
1:K:290:VAL:HG22	1:K:291:PRO:HD2	1.89	0.53
1:H:89:LYS:CE	1:H:131:LYS:HZ3	2.21	0.53
1:F:14:GLU:N	1:F:14:GLU:OE2	2.42	0.53
1:L:25:VAL:O	1:L:94:THR:OG1	2.25	0.52
1:A:231:LYS:HE2	1:A:235:GLU:OE1	2.10	0.52
1:E:21:LYS:HE3	1:E:86:ALA:O	2.10	0.52
1:G:290:VAL:HG22	1:G:291:PRO:HD2	1.89	0.52
1:J:274:THR:HB	1:J:299:ILE:HD13	1.91	0.52
1:D:107:ASN:HA	1:D:110:GLN:HB3	1.90	0.52
1:F:228:GLU:O	1:F:232:GLN:HG3	2.09	0.52
1:L:109:VAL:O	1:L:113:VAL:HG23	2.10	0.52
1:K:276:ILE:HD13	1:K:282:ILE:HD13	1.91	0.52
1:C:185:HIS:NE2	2:C:401:9G9:C15	2.73	0.52
1:G:276:ILE:HD13	1:G:282:ILE:HD13	1.91	0.52
1:H:93:ILE:CD1	1:H:135:VAL:CB	2.88	0.52
1:G:106:LEU:HD12	1:G:325:ILE:HG23	1.92	0.52
1:K:22:ILE:HG13	1:K:44:ALA:HB2	1.91	0.52
1:D:330:GLN:HG2	1:D:331:PHE:H	1.74	0.51
1:F:20:ASN:HA	1:F:89:LYS:HD2	1.92	0.51
1:E:106:LEU:O	1:E:107:ASN:HB2	2.09	0.51
1:B:274:THR:HB	1:B:299:ILE:HD13	1.92	0.51
1:G:117:LYS:O	3:G:402:HOH:O	2.19	0.51
1:G:237:ALA:O	1:G:241:ILE:HG13	2.11	0.51
1:E:302:LEU:CD2	1:H:11:LEU:HD11	2.40	0.51
1:H:143:THR:HG23	1:H:157:VAL:HG12	1.93	0.51
1:E:12:LEU:HD12	1:H:300:SER:HA	1.92	0.51
1:L:312:GLU:O	1:L:316:LYS:HG3	2.10	0.51
1:H:15:GLU:OE2	1:H:16:GLN:HA	2.11	0.51
1:H:273:SER:HA	1:H:289:SER:HA	1.93	0.51
1:A:304:LYS:HD3	1:D:9:TYR:HB2	1.92	0.51
1:D:173:MET:SD	1:D:184:CYS:HB3	2.50	0.51
1:E:112:ASN:HB3	1:E:142:LEU:HD21	1.92	0.51
1:E:114:ASN:HA	1:E:117:LYS:CD	2.40	0.51
1:E:6:GLN:HG2	1:F:213:LEU:HD22	1.92	0.51
1:I:218:GLY:O	1:I:227:LYS:HD3	2.11	0.51
1:H:110:GLN:O	1:H:114:ASN:ND2	2.44	0.50
1:J:308:THR:OG1	1:J:310:GLU:HG3	2.12	0.50
1:F:286:VAL:HG23	1:F:326:GLN:NE2	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:VAL:HG11	1:B:141:ILE:HG21	1.92	0.50
1:B:224:GLU:O	1:B:225:GLN:HB2	2.10	0.50
1:H:135:VAL:CG1	1:H:251:ILE:HD11	2.40	0.50
1:G:270:HIS:CD2	1:G:294:LEU:HD12	2.47	0.50
1:B:173:MET:SD	1:B:184:CYS:HB3	2.52	0.50
1:G:45:ASP:O	1:G:74:PRO:HD2	2.12	0.50
1:I:123:VAL:HG11	1:I:132:LEU:HD21	1.94	0.49
1:F:272:VAL:O	1:F:289:SER:HA	2.12	0.49
1:I:310:GLU:OE1	1:I:310:GLU:N	2.43	0.49
1:C:202:GLY:O	1:C:204:ASN:ND2	2.42	0.49
1:H:230:HIS:O	1:H:234:VAL:HG23	2.13	0.49
1:E:292:CYS:HB3	1:E:299:ILE:HG23	1.95	0.49
1:F:244:LYS:HE3	1:F:246:TYR:O	2.12	0.49
1:K:1:ALA:N	3:K:401:HOH:O	2.37	0.49
1:K:222:ASP:O	1:K:225:GLN:NE2	2.37	0.49
1:A:165:ASP:OD2	1:A:192:HIS:ND1	2.41	0.49
1:J:56:LYS:NZ	1:J:60:GLU:OE2	2.45	0.49
1:J:123:VAL:HG11	1:J:132:LEU:HD21	1.95	0.49
1:D:223:LYS:HD2	1:D:223:LYS:H	1.77	0.49
1:D:272:VAL:O	1:D:289:SER:HA	2.13	0.49
1:C:6:GLN:HB3	1:D:213:LEU:HD23	1.94	0.49
1:L:105:ARG:HD3	1:L:325:ILE:HD13	1.94	0.48
1:A:214:HIS:HB2	1:B:3:LEU:HD13	1.94	0.48
1:G:170:ARG:NH2	1:H:69:LEU:HD21	2.28	0.48
1:L:272:VAL:O	1:L:289:SER:HA	2.13	0.48
1:F:323:TRP:O	1:F:327:LYS:HG2	2.13	0.48
1:K:279:LEU:HD11	1:K:302:LEU:HD11	1.96	0.48
1:L:120:ILE:HB	1:L:121:PRO:HD3	1.94	0.48
1:G:312:GLU:O	1:G:316:LYS:HG3	2.14	0.48
1:H:141:ILE:O	1:H:145:VAL:HG23	2.14	0.48
1:C:204:ASN:ND2	2:C:401:9G9:C10	2.75	0.48
1:D:291:PRO:HG2	1:D:303:VAL:HB	1.96	0.48
1:G:66:HIS:CE1	1:H:236:SER:HB2	2.49	0.48
1:C:106:LEU:HD11	1:C:328:GLU:HB2	1.96	0.48
1:H:282:ILE:HD11	1:H:316:LYS:HA	1.95	0.48
1:D:15:GLU:O	1:D:16:GLN:HB2	2.14	0.47
1:J:268:ARG:HD3	1:L:182:LEU:HD23	1.96	0.47
1:L:277:LYS:HD3	1:L:285:ASP:OD1	2.13	0.47
1:K:45:ASP:O	1:K:74:PRO:HD2	2.14	0.47
1:A:293:ILE:HD13	1:A:301:ASP:HB2	1.96	0.47
1:B:266:LEU:HD23	1:C:74:PRO:HG3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:292:CYS:HB3	1:C:299:ILE:HG23	1.96	0.47
1:E:114:ASN:HA	1:E:117:LYS:HD2	1.96	0.47
1:E:308:THR:CG2	1:E:311:GLU:H	2.28	0.47
1:L:93:ILE:HB	1:L:135:VAL:H	1.79	0.47
1:A:106:LEU:O	1:A:107:ASN:CB	2.62	0.47
1:B:322:LEU:O	1:B:326:GLN:HG3	2.15	0.47
1:K:329:LEU:C	1:K:331:PHE:H	2.18	0.47
1:B:153:PRO:HB2	1:B:155:ASN:OD1	2.14	0.47
1:G:153:PRO:HB2	1:G:155:ASN:OD1	2.15	0.47
1:I:173:MET:SD	1:I:184:CYS:HB3	2.55	0.47
1:B:125:LYS:HD2	1:B:126:TYR:CE2	2.49	0.47
1:A:243:LEU:HD13	1:B:58:LYS:HG2	1.97	0.47
1:D:330:GLN:HG2	1:D:331:PHE:N	2.30	0.47
1:E:180:HIS:HB2	1:G:266:LEU:O	2.15	0.47
1:G:306:THR:HB	3:G:415:HOH:O	2.15	0.47
1:A:275:MET:CE	1:A:277:LYS:HB3	2.44	0.47
1:C:82:TYR:OH	1:C:119:ILE:HG12	2.15	0.47
1:C:107:ASN:O	1:C:111:ARG:HG3	2.15	0.47
1:F:276:ILE:HG21	1:F:288:LEU:HD13	1.97	0.47
1:F:304:LYS:HD3	1:G:9:TYR:HB2	1.97	0.46
1:A:286:VAL:HG11	1:A:319:ALA:HB1	1.97	0.46
1:G:197:VAL:HG21	1:G:315:LEU:CD1	2.45	0.46
1:I:14:GLU:OE1	1:I:16:GLN:HG3	2.15	0.46
1:K:244:LYS:HE2	1:K:246:TYR:O	2.14	0.46
1:L:100:GLN:HG3	1:L:103:GLU:HB2	1.97	0.46
1:J:52:VAL:O	1:J:80:LYS:HE3	2.16	0.46
1:E:141:ILE:HG13	1:E:322:LEU:HD22	1.96	0.46
1:H:97:ALA:N	1:H:112:ASN:OD1	2.47	0.46
1:E:3:LEU:HD13	1:F:214:HIS:HB2	1.98	0.46
1:G:106:LEU:O	1:G:108:LEU:N	2.46	0.46
1:H:93:ILE:HB	1:H:133:LEU:O	2.16	0.46
1:H:94:THR:O	1:H:95:ALA:HB3	2.15	0.46
1:I:99:GLN:HG3	1:I:108:LEU:HG	1.97	0.46
1:K:30:VAL:HG22	1:K:251:ILE:HG21	1.98	0.46
1:C:312:GLU:HG2	1:C:316:LYS:HE3	1.97	0.46
1:A:228:GLU:OE2	1:A:231:LYS:HD2	2.15	0.46
1:B:162:CYS:HB3	1:B:165:ASP:HB2	1.97	0.46
1:C:185:HIS:O	1:C:203:MET:HA	2.16	0.46
1:E:153:PRO:HB2	1:E:155:ASN:OD1	2.16	0.46
1:E:308:THR:HG23	1:E:311:GLU:H	1.81	0.46
1:A:212:THR:HB	3:A:414:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:50:VAL:HG11	1:J:82:TYR:CZ	2.51	0.46
1:K:14:GLU:O	1:K:16:GLN:N	2.45	0.46
1:L:125:LYS:HE2	1:L:126:TYR:CZ	2.51	0.46
1:C:206:ALA:H	2:C:401:9G9:C15	2.28	0.46
1:J:216:ASP:O	1:J:222:ASP:HB2	2.15	0.46
1:J:197:VAL:HG22	1:J:314:ARG:HD2	1.98	0.46
1:A:274:THR:O	1:A:276:ILE:HG23	2.16	0.46
1:E:330:GLN:OE1	1:E:330:GLN:HA	2.16	0.46
1:L:127:SER:HB3	1:L:130:CYS:HB3	1.98	0.46
1:A:51:ASP:H	1:A:57:LEU:HD22	1.81	0.45
1:B:293:ILE:HD12	1:B:301:ASP:HB2	1.98	0.45
1:C:283:LYS:HA	1:C:283:LYS:HD3	1.87	0.45
1:C:3:LEU:HD13	1:D:214:HIS:HB2	1.97	0.45
1:E:203:MET:HE2	1:E:210:LEU:HD12	1.98	0.45
1:E:237:ALA:O	1:E:241:ILE:HG13	2.16	0.45
1:H:13:LYS:HB2	1:H:14:GLU:OE1	2.17	0.45
1:J:228:GLU:OE2	1:J:231:LYS:HD3	2.16	0.45
1:L:93:ILE:CG2	1:L:134:ILE:HA	2.46	0.45
1:L:235:GLU:O	1:L:239:GLU:HG2	2.16	0.45
1:B:266:LEU:O	1:D:180:HIS:HB2	2.16	0.45
1:B:123:VAL:HG11	1:B:132:LEU:HD21	1.98	0.45
1:I:12:LEU:HD11	1:L:155:ASN:ND2	2.31	0.45
1:A:135:VAL:HA	1:A:160:SER:OG	2.17	0.45
1:A:273:SER:HA	1:A:288:LEU:O	2.17	0.45
1:B:106:LEU:HG	1:B:107:ASN:N	2.31	0.45
1:F:13:LYS:N	1:F:14:GLU:OE2	2.42	0.45
1:E:12:LEU:HD11	1:H:155:ASN:ND2	2.31	0.45
1:A:11:LEU:HD21	1:D:154:LYS:HE2	1.99	0.45
1:L:25:VAL:O	1:L:93:ILE:O	2.34	0.45
1:A:141:ILE:HG13	1:A:322:LEU:HD22	1.98	0.45
1:D:274:THR:HB	1:D:299:ILE:HD13	1.98	0.45
1:G:12:LEU:HD22	1:G:14:GLU:CD	2.37	0.45
1:H:125:LYS:HE3	1:H:126:TYR:CZ	2.51	0.45
1:H:189:LEU:HB3	1:H:289:SER:O	2.17	0.45
1:H:93:ILE:HG23	1:H:94:THR:H	1.80	0.45
1:K:276:ILE:HD11	1:K:286:VAL:HB	1.99	0.45
1:A:111:ARG:HD2	3:A:411:HOH:O	2.17	0.45
1:E:269:VAL:HA	1:E:292:CYS:O	2.16	0.45
1:E:49:LEU:O	1:E:78:SER:HA	2.17	0.45
1:I:91:VAL:HB	1:I:132:LEU:HD23	1.98	0.45
1:B:217:LEU:HD12	1:B:226:TRP:CG	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:LEU:O	1:D:107:ASN:HB3	2.17	0.45
1:G:16:GLN:O	1:G:17:THR:OG1	2.34	0.45
1:H:153:PRO:HB2	1:H:155:ASN:OD1	2.17	0.45
1:H:158:ILE:HG23	1:H:299:ILE:HD11	1.98	0.45
1:J:272:VAL:O	1:J:289:SER:HA	2.17	0.45
1:E:291:PRO:HG2	1:E:303:VAL:HB	1.98	0.44
1:L:94:THR:O	1:L:95:ALA:HB3	2.17	0.44
1:A:13:LYS:CE	1:A:13:LYS:HA	2.46	0.44
1:A:169:PHE:HD2	1:A:233:VAL:HG21	1.83	0.44
1:A:182:LEU:HD11	1:B:72:ARG:NH1	2.32	0.44
1:A:330:GLN:NE2	3:A:406:HOH:O	2.50	0.44
1:B:180:HIS:HB2	1:D:266:LEU:O	2.16	0.44
1:K:13:LYS:HB3	1:K:14:GLU:H	1.47	0.44
1:L:93:ILE:HG13	1:L:94:THR:H	1.81	0.44
1:C:15:GLU:HB2	1:C:16:GLN:NE2	2.31	0.44
1:D:288:LEU:HD13	1:D:315:LEU:HD22	1.99	0.44
1:J:182:LEU:HD12	1:L:268:ARG:HD3	2.00	0.44
1:A:185:HIS:CG	2:C:401:9G9:C5	3.01	0.44
1:E:285:ASP:N	1:E:285:ASP:OD1	2.50	0.44
1:E:61:MET:SD	1:E:78:SER:OG	2.62	0.44
1:H:145:VAL:O	1:H:149:ILE:HG13	2.18	0.44
1:L:134:ILE:HG13	1:L:143:THR:HA	2.00	0.44
1:C:260:GLU:HG3	1:C:264:LYS:HD2	2.00	0.44
1:C:293:ILE:HD13	1:C:301:ASP:HB2	1.99	0.44
1:F:291:PRO:HB2	1:F:303:VAL:HB	2.00	0.44
1:H:165:ASP:OD2	1:H:192:HIS:ND1	2.51	0.44
1:H:294:LEU:HD23	1:H:294:LEU:HA	1.75	0.44
1:J:16:GLN:HB3	1:K:296:GLN:NE2	2.32	0.44
1:J:280:TYR:HB2	1:J:282:ILE:HD12	2.00	0.44
1:K:211:LYS:HE3	1:K:215:PRO:O	2.18	0.44
1:D:106:LEU:HD22	1:D:107:ASN:H	1.83	0.44
1:D:282:ILE:HD13	1:D:319:ALA:HB2	1.99	0.44
1:E:114:ASN:O	1:E:117:LYS:HG2	2.17	0.44
1:G:198:PRO:HD3	1:G:230:HIS:CE1	2.53	0.44
1:G:3:LEU:HD23	1:H:226:TRP:CZ2	2.52	0.44
1:K:65:GLN:O	1:K:68:SER:OG	2.31	0.44
1:L:27:VAL:HG11	1:L:57:LEU:HD12	2.00	0.44
1:A:185:HIS:HD1	2:C:401:9G9:C13	2.30	0.43
1:A:204:ASN:ND2	2:C:401:9G9:C9	2.76	0.43
1:G:10:ASN:ND2	1:G:13:LYS:HE2	2.33	0.43
1:B:15:GLU:N	3:B:403:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:194:ASP:HA	1:L:234:VAL:CG1	2.47	0.43
1:L:94:THR:HG22	1:L:135:VAL:HG11	2.00	0.43
1:B:27:VAL:HG11	1:B:57:LEU:HD12	2.00	0.43
1:J:229:VAL:O	1:J:233:VAL:HG23	2.18	0.43
1:L:93:ILE:HD12	1:L:135:VAL:CA	2.46	0.43
1:H:93:ILE:CG2	1:H:94:THR:N	2.79	0.43
1:L:159:GLY:HA3	1:L:273:SER:HB3	1.99	0.43
1:L:57:LEU:HD21	1:L:79:GLY:H	1.83	0.43
1:H:57:LEU:HD21	1:H:79:GLY:H	1.84	0.43
1:L:93:ILE:HG22	1:L:133:LEU:O	2.18	0.43
1:H:25:VAL:O	1:H:94:THR:N	2.52	0.43
1:A:291:PRO:HG2	1:A:303:VAL:HB	1.99	0.43
1:C:109:VAL:O	1:C:113:VAL:HG23	2.19	0.43
1:K:51:ASP:H	1:K:57:LEU:HD22	1.84	0.43
1:G:292:CYS:SG	1:G:299:ILE:HD13	2.59	0.43
1:H:143:THR:CG2	1:H:157:VAL:HG12	2.49	0.43
1:H:163:ASN:HA	1:H:271:PRO:HG2	2.01	0.43
1:J:153:PRO:HB2	1:J:155:ASN:OD1	2.19	0.43
1:K:286:VAL:HG22	1:K:323:TRP:HB2	2.00	0.43
1:C:100:GLN:HG3	1:C:101:GLU:OE2	2.19	0.43
1:C:290:VAL:HG11	1:C:302:LEU:HD13	2.01	0.43
1:E:21:LYS:HG3	1:E:46:GLU:HB3	2.01	0.43
1:G:164:LEU:HD11	1:G:250:ALA:HB1	2.01	0.43
1:K:99:GLN:HG2	1:K:99:GLN:H	1.55	0.43
1:D:26:GLY:HA2	1:D:51:ASP:OD2	2.18	0.43
1:L:274:THR:O	1:L:287:PHE:HA	2.19	0.43
1:A:120:ILE:HB	1:A:121:PRO:HD3	2.00	0.42
1:D:10:ASN:OD1	1:D:13:LYS:HG3	2.18	0.42
1:E:86:ALA:HA	1:E:126:TYR:O	2.19	0.42
1:G:54:GLU:CD	1:G:80:LYS:HE2	2.39	0.42
1:K:272:VAL:O	1:K:289:SER:HA	2.18	0.42
1:G:194:ASP:HA	1:G:234:VAL:CG1	2.49	0.42
1:H:23:THR:HB	1:H:91:VAL:HG22	2.01	0.42
1:H:316:LYS:NZ	1:H:320:ASP:OD2	2.50	0.42
1:I:11:LEU:HD11	1:L:302:LEU:HD13	2.01	0.42
1:I:266:LEU:O	1:K:180:HIS:HB2	2.19	0.42
1:K:281:GLY:O	1:K:283:LYS:HD3	2.20	0.42
1:A:279:LEU:O	1:A:282:ILE:HG12	2.19	0.42
1:H:270:HIS:CD2	1:H:294:LEU:HD12	2.54	0.42
1:J:194:ASP:HA	1:J:234:VAL:CG1	2.50	0.42
1:B:91:VAL:HG21	1:B:123:VAL:HG13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:LEU:O	1:C:180:HIS:HB2	2.19	0.42
1:A:185:HIS:ND1	2:C:401:9G9:C13	2.82	0.42
1:E:329:LEU:HD23	1:E:329:LEU:HA	1.88	0.42
1:H:277:LYS:HD3	1:H:285:ASP:OD1	2.19	0.42
1:E:21:LYS:HG3	1:E:46:GLU:CG	2.49	0.42
1:G:130:CYS:O	1:G:156:ARG:HD2	2.19	0.42
1:H:107:ASN:O	1:H:111:ARG:N	2.34	0.42
1:L:211:LYS:O	1:L:215:PRO:HA	2.20	0.42
1:D:43:LEU:HD21	1:D:260:GLU:HG3	2.01	0.42
1:A:86:ALA:HA	1:A:126:TYR:HB3	2.01	0.42
1:B:14:GLU:HG3	1:B:15:GLU:N	2.34	0.42
1:A:226:TRP:CZ2	1:B:3:LEU:HD23	2.55	0.42
1:D:141:ILE:CG1	1:D:322:LEU:HD22	2.49	0.42
1:D:65:GLN:O	1:D:68:SER:OG	2.31	0.42
1:H:131:LYS:HG3	1:H:156:ARG:HG2	2.02	0.42
1:H:124:VAL:HG21	1:H:150:SER:HB2	2.02	0.42
1:H:180:HIS:ND1	1:H:182:LEU:HB2	2.35	0.42
1:H:240:VAL:CG1	1:H:247:THR:HG22	2.49	0.42
1:I:105:ARG:O	1:I:109:VAL:HG23	2.20	0.42
1:I:159:GLY:HA3	1:I:273:SER:HB3	2.02	0.42
1:K:203:MET:HE1	1:K:210:LEU:HD12	2.01	0.42
1:B:276:ILE:HD12	1:B:282:ILE:HD13	2.01	0.42
1:F:191:GLU:HG3	1:F:322:LEU:HD21	2.01	0.42
1:E:302:LEU:HD23	1:H:11:LEU:HD11	2.01	0.42
1:L:290:VAL:HG22	1:L:291:PRO:HD2	2.01	0.42
1:H:251:ILE:HD13	1:H:251:ILE:HG21	1.77	0.42
1:I:113:VAL:HG11	1:I:331:PHE:HB3	2.00	0.42
1:L:105:ARG:O	1:L:109:VAL:HG23	2.20	0.42
1:C:106:LEU:CD1	1:C:328:GLU:HB2	2.49	0.41
1:E:274:THR:HB	1:E:299:ILE:HD13	2.02	0.41
1:L:176:ARG:NH1	1:L:224:GLU:O	2.53	0.41
1:J:231:LYS:O	1:J:235:GLU:HG3	2.19	0.41
1:L:106:LEU:HG	1:L:107:ASN:H	1.84	0.41
1:L:133:LEU:HA	1:L:133:LEU:HD12	1.92	0.41
1:F:155:ASN:O	1:F:298:GLY:HA3	2.19	0.41
1:F:266:LEU:O	1:H:180:HIS:HB2	2.20	0.41
1:I:200:TRP:HZ2	1:I:227:LYS:HD2	1.85	0.41
1:K:179:VAL:HG12	1:K:180:HIS:O	2.20	0.41
1:L:14:GLU:OE2	1:L:15:GLU:N	2.53	0.41
1:A:211:LYS:HD2	1:A:217:LEU:HB3	2.02	0.41
1:G:269:VAL:HA	1:G:292:CYS:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:100:GLN:H	1:L:100:GLN:HG2	1.50	0.41
1:B:230:HIS:O	1:B:233:VAL:HB	2.19	0.41
1:B:278:GLY:O	1:B:279:LEU:HD23	2.20	0.41
1:D:81:ASP:O	1:D:84:VAL:HG22	2.20	0.41
1:G:227:LYS:O	1:G:227:LYS:CD	2.62	0.41
1:H:93:ILE:HG21	1:H:135:VAL:CG2	2.51	0.41
1:K:82:TYR:CG	1:K:122:ASN:HB3	2.55	0.41
1:I:105:ARG:HD3	1:I:325:ILE:HD11	2.01	0.41
1:J:163:ASN:HA	1:J:271:PRO:HG2	2.03	0.41
1:K:228:GLU:O	1:K:232:GLN:HG3	2.21	0.41
1:B:269:VAL:HG22	1:B:293:ILE:HG12	2.01	0.41
1:G:21:LYS:HB3	1:G:88:SER:HA	2.03	0.41
1:H:105:ARG:O	1:H:109:VAL:HG23	2.21	0.41
1:H:21:LYS:HB3	1:H:88:SER:HA	2.03	0.41
1:I:148:LYS:HB3	1:I:331:PHE:CE2	2.56	0.41
1:J:106:LEU:O	1:J:107:ASN:CB	2.61	0.41
1:J:89:LYS:O	1:J:131:LYS:HE3	2.21	0.41
1:L:107:ASN:O	1:L:111:ARG:HG3	2.21	0.41
1:D:239:GLU:OE2	1:D:239:GLU:HA	2.20	0.41
1:F:162:CYS:HB3	1:F:165:ASP:HB2	2.03	0.41
1:H:327:LYS:HE3	1:H:327:LYS:HB2	1.90	0.41
1:H:282:ILE:N	1:H:282:ILE:HD13	2.36	0.41
1:C:71:LEU:O	1:C:72:ARG:HD3	2.21	0.41
1:H:93:ILE:HG21	1:H:135:VAL:HG23	2.03	0.41
1:J:169:PHE:HD1	1:J:233:VAL:HG21	1.86	0.41
1:L:93:ILE:CG2	1:L:133:LEU:C	2.88	0.41
1:L:216:ASP:O	1:L:222:ASP:HB2	2.21	0.41
1:L:248:SER:O	1:L:251:ILE:HG22	2.21	0.41
1:F:213:LEU:HD23	1:F:213:LEU:HA	1.89	0.40
1:G:123:VAL:HG11	1:G:132:LEU:HD21	2.04	0.40
1:H:177:LEU:HA	1:H:177:LEU:HD23	1.86	0.40
1:J:173:MET:SD	1:J:184:CYS:HB3	2.61	0.40
1:K:111:ARG:O	1:K:115:ILE:HG13	2.21	0.40
1:C:170:ARG:NE	1:C:184:CYS:O	2.51	0.40
1:A:185:HIS:NE2	2:C:401:9G9:C14	2.84	0.40
1:J:99:GLN:HB2	1:J:108:LEU:HD22	2.03	0.40
1:A:203:MET:HE2	1:A:210:LEU:HD12	2.03	0.40
1:C:133:LEU:HA	1:C:158:ILE:O	2.21	0.40
1:D:111:ARG:HD2	1:D:111:ARG:HH11	1.75	0.40
1:E:150:SER:OG	1:E:152:PHE:CG	2.74	0.40
1:E:302:LEU:HD23	1:H:11:LEU:CD1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:200:TRP:HB3	1:F:217:LEU:HD23	2.03	0.40
1:G:16:GLN:HB2	1:G:16:GLN:HE21	1.75	0.40
1:H:244:LYS:HE2	1:H:246:TYR:O	2.22	0.40
1:I:158:ILE:HG12	1:I:299:ILE:HG13	2.03	0.40
1:K:49:LEU:O	1:K:78:SER:HA	2.20	0.40
1:B:165:ASP:OD2	1:B:192:HIS:ND1	2.51	0.40
1:B:98:ARG:HH12	1:B:137:ASN:HB2	1.86	0.40
1:C:32:MET:CE	1:C:60:GLU:HB3	2.52	0.40
1:E:223:LYS:CD	1:E:223:LYS:H	2.29	0.40
1:C:171:TYR:HA	1:C:181:PRO:HG3	2.04	0.40
1:E:82:TYR:O	1:E:85:THR:HB	2.20	0.40
1:G:141:ILE:HG13	1:G:322:LEU:HD22	2.03	0.40
1:H:93:ILE:HD11	1:H:136:SER:CB	2.51	0.40
1:I:56:LYS:NZ	1:I:60:GLU:OE2	2.54	0.40
1:I:20:ASN:HA	1:I:89:LYS:HD2	2.03	0.40
1:K:186:GLY:O	1:K:187:TRP:HB2	2.21	0.40
1:L:237:ALA:O	1:L:241:ILE:HG13	2.22	0.40
1:L:42:ASP:OD1	1:L:72:ARG:HB2	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	329/337 (98%)	312 (95%)	14 (4%)	3 (1%)	17	35
1	B	329/337 (98%)	310 (94%)	14 (4%)	5 (2%)	10	21
1	C	329/337 (98%)	312 (95%)	14 (4%)	3 (1%)	17	35
1	D	329/337 (98%)	310 (94%)	15 (5%)	4 (1%)	13	27
1	E	329/337 (98%)	313 (95%)	14 (4%)	2 (1%)	25	47
1	F	329/337 (98%)	317 (96%)	9 (3%)	3 (1%)	17	35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	329/337 (98%)	314 (95%)	13 (4%)	2 (1%)	25	47
1	H	329/337 (98%)	306 (93%)	16 (5%)	7 (2%)	7	13
1	I	329/337 (98%)	314 (95%)	9 (3%)	6 (2%)	8	16
1	J	329/337 (98%)	317 (96%)	10 (3%)	2 (1%)	25	47
1	K	329/337 (98%)	315 (96%)	9 (3%)	5 (2%)	10	21
1	L	329/337 (98%)	308 (94%)	17 (5%)	4 (1%)	13	27
All	All	3948/4044 (98%)	3748 (95%)	154 (4%)	46 (1%)	13	27

All (46) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	107	ASN
1	D	16	GLN
1	D	94	THR
1	E	107	ASN
1	H	14	GLU
1	H	15	GLU
1	I	13	LYS
1	I	16	GLN
1	J	107	ASN
1	K	15	GLU
1	L	14	GLU
1	L	222	ASP
1	A	107	ASN
1	B	15	GLU
1	B	55	ASP
1	B	107	ASN
1	D	222	ASP
1	H	16	GLN
1	H	93	ILE
1	I	14	GLU
1	I	55	ASP
1	K	13	LYS
1	K	16	GLN
1	L	93	ILE
1	A	14	GLU
1	C	17	THR
1	E	17	THR
1	F	222	ASP
1	H	13	LYS

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Mol	Chain	Res	Type
1	A	17	THR
1	B	17	THR
1	F	17	THR
1	F	196	SER
1	G	187	TRP
1	H	17	THR
1	I	107	ASN
1	J	17	THR
1	L	17	THR
1	B	16	GLN
1	C	187	TRP
1	D	17	THR
1	G	17	THR
1	K	187	TRP
1	K	186	GLY
1	I	17	THR
1	H	53	ILE

### 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	287/293 (98%)	286 (100%)	1 (0%)	92	98
1	B	287/293 (98%)	276 (96%)	11 (4%)	33	59
1	C	287/293 (98%)	287 (100%)	0	100	100
1	D	287/293 (98%)	279 (97%)	8 (3%)	43	69
1	E	287/293 (98%)	280 (98%)	7 (2%)	49	74
1	F	287/293 (98%)	280 (98%)	7 (2%)	49	74
1	G	287/293 (98%)	284 (99%)	3 (1%)	76	90
1	H	287/293 (98%)	277 (96%)	10 (4%)	36	62
1	I	287/293 (98%)	283 (99%)	4 (1%)	67	85
1	J	287/293 (98%)	282 (98%)	5 (2%)	60	81
1	K	287/293 (98%)	283 (99%)	4 (1%)	67	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	L	287/293 (98%)	280 (98%)	7 (2%)	49 74
All	All	3444/3516 (98%)	3377 (98%)	67 (2%)	57 79

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ARG
1	B	12	LEU
1	B	13	LYS
1	B	15	GLU
1	B	107	ASN
1	B	131	LYS
1	B	154	LYS
1	B	231	LYS
1	B	236	SER
1	B	283	LYS
1	B	309	SER
1	B	318	SER
1	D	83	ASN
1	D	105	ARG
1	D	131	LYS
1	D	133	LEU
1	D	194	ASP
1	D	221	LYS
1	D	236	SER
1	D	238	TYR
1	E	220	ASP
1	E	223	LYS
1	E	242	LYS
1	E	273	SER
1	E	302	LEU
1	E	318	SER
1	E	331	PHE
1	F	12	LEU
1	F	175	GLU
1	F	195	SER
1	F	217	LEU
1	F	302	LEU
1	F	316	LYS
1	F	327	LYS
1	G	223	LYS
1	G	239	GLU

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Mol	Chain	Res	Type
1	G	275	MET
1	H	54	GLU
1	H	56	LYS
1	H	89	LYS
1	H	160	SER
1	H	173	MET
1	H	209	SER
1	H	231	LYS
1	H	275	MET
1	H	289	SER
1	H	309	SER
1	I	14	GLU
1	I	133	LEU
1	I	217	LEU
1	I	331	PHE
1	J	12	LEU
1	J	98	ARG
1	J	107	ASN
1	J	273	SER
1	J	314	ARG
1	K	98	ARG
1	K	131	LYS
1	K	209	SER
1	K	283	LYS
1	L	14	GLU
1	L	160	SER
1	L	209	SER
1	L	220	ASP
1	L	221	LYS
1	L	254	SER
1	L	261	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	297	ASN
1	I	330	GLN
1	J	185	HIS
1	K	110	GLN
1	K	296	GLN
1	L	20	ASN
1	L	185	HIS



### 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 ⓘ

1 ligand is 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	9G9	C	401	-	27,27,27	2.47	9 (33%)	38,38,38	7.83	20 (52%)

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	9G9	C	401	-	-	5/12/24/24	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	9G9	C8-C6	8.08	1.62	1.54
2	C	401	9G9	C7-C5	4.47	1.58	1.54
2	C	401	9G9	C6-C5	4.33	1.66	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	9G9	C20-C18	3.41	1.47	1.39
2	C	401	9G9	C14-C18	2.94	1.44	1.38
2	C	401	9G9	C19-C17	2.83	1.46	1.39
2	C	401	9G9	C15-C11	2.21	1.43	1.38
2	C	401	9G9	C8-C12	2.13	1.56	1.51
2	C	401	9G9	C14-C12	2.11	1.42	1.39

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	9G9	C9-C5-C7	-26.92	81.89	110.91
2	C	401	9G9	C7-C5-C6	23.56	146.43	110.07
2	C	401	9G9	C8-C6-C5	20.44	141.61	110.07
2	C	401	9G9	C10-C6-C8	-12.53	97.40	110.91
2	C	401	9G9	C12-C8-C6	11.96	131.08	113.69
2	C	401	9G9	C22-C16-C12	8.99	133.39	121.03
2	C	401	9G9	C11-C7-C5	6.85	123.65	113.69
2	C	401	9G9	C8-C12-C14	6.49	131.57	120.44
2	C	401	9G9	O2-C18-C20	-5.33	103.71	109.78
2	C	401	9G9	C24-O2-C18	4.22	110.83	105.34
2	C	401	9G9	C16-C22-C20	-3.96	112.52	120.06
2	C	401	9G9	O2-C18-C14	3.58	132.64	127.85
2	C	401	9G9	C8-C12-C16	-3.28	114.38	120.91
2	C	401	9G9	C16-C12-C14	-3.22	114.03	118.54
2	C	401	9G9	C23-O3-C19	2.71	108.87	105.34
2	C	401	9G9	O1-C17-C13	2.47	131.15	127.85
2	C	401	9G9	C18-C14-C12	-2.39	116.47	120.05
2	C	401	9G9	C23-O1-C17	2.37	108.42	105.34
2	C	401	9G9	C21-C15-C11	2.35	124.26	121.03
2	C	401	9G9	C7-C11-C13	2.25	124.30	120.44

There are no chirality outliers.

All (5) torsion outliers are listed below:

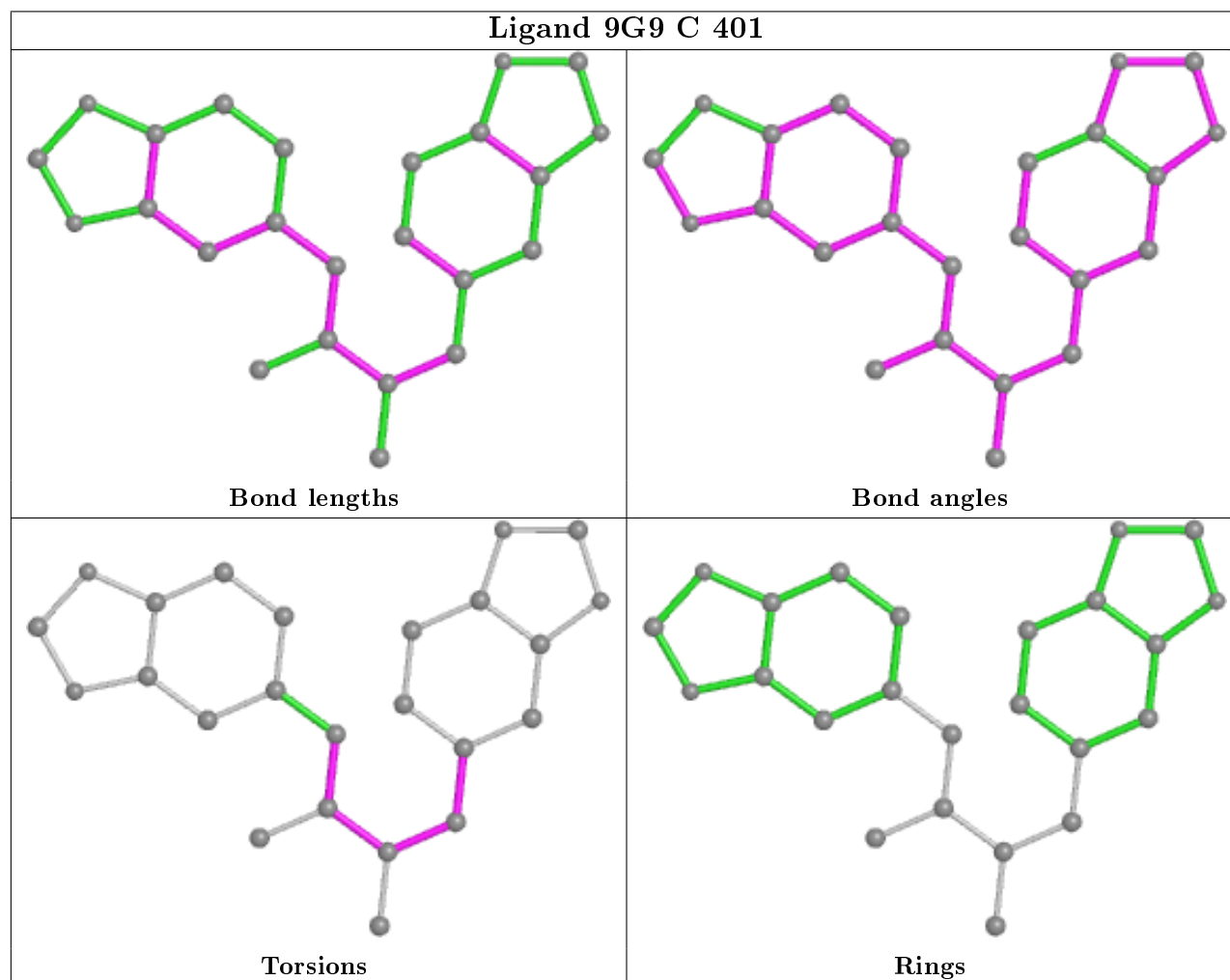
Mol	Chain	Res	Type	Atoms
2	C	401	9G9	C13-C11-C7-C5
2	C	401	9G9	C15-C11-C7-C5
2	C	401	9G9	C9-C5-C6-C10
2	C	401	9G9	C6-C5-C7-C11
2	C	401	9G9	C5-C6-C8-C12

There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	9G9	13	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	331/337 (98%)	-0.34	4 (1%) 79 76	36, 48, 67, 110	0
1	B	331/337 (98%)	-0.15	5 (1%) 73 70	37, 51, 78, 97	0
1	C	331/337 (98%)	-0.25	3 (0%) 84 82	34, 45, 65, 99	0
1	D	331/337 (98%)	-0.32	1 (0%) 94 93	36, 51, 72, 96	0
1	E	331/337 (98%)	-0.15	9 (2%) 54 48	36, 56, 83, 108	0
1	F	331/337 (98%)	-0.34	2 (0%) 89 88	35, 46, 68, 98	0
1	G	331/337 (98%)	-0.13	13 (3%) 39 32	37, 51, 77, 103	0
1	H	331/337 (98%)	-0.02	8 (2%) 59 53	39, 58, 75, 99	0
1	I	331/337 (98%)	-0.37	1 (0%) 94 93	32, 42, 64, 101	0
1	J	331/337 (98%)	-0.47	1 (0%) 94 93	32, 44, 64, 91	0
1	K	331/337 (98%)	-0.40	4 (1%) 79 76	30, 43, 58, 87	0
1	L	331/337 (98%)	-0.16	8 (2%) 59 53	34, 53, 80, 108	0
All	All	3972/4044 (98%)	-0.26	59 (1%) 73 70	30, 49, 74, 110	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	14	GLU	5.8
1	B	15	GLU	4.7
1	G	15	GLU	4.6
1	E	14	GLU	4.2
1	G	102	GLY	3.9
1	A	15	GLU	3.8
1	K	330	GLN	3.7
1	E	101	GLU	3.7
1	L	15	GLU	3.7
1	L	101	GLU	3.6
1	B	14	GLU	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	15	GLU	3.3
1	D	14	GLU	3.2
1	G	221	LYS	3.2
1	C	15	GLU	3.1
1	G	101	GLU	3.1
1	G	14	GLU	3.1
1	C	101	GLU	3.1
1	E	98	ARG	3.1
1	F	13	LYS	3.0
1	H	221	LYS	2.9
1	H	331	PHE	2.9
1	B	102	GLY	2.9
1	J	331	PHE	2.9
1	G	99	GLN	2.9
1	H	15	GLU	2.8
1	H	1	ALA	2.8
1	L	102	GLY	2.8
1	F	15	GLU	2.8
1	K	15	GLU	2.7
1	L	14	GLU	2.7
1	G	100	GLN	2.5
1	E	16	GLN	2.5
1	I	1	ALA	2.4
1	C	100	GLN	2.4
1	A	16	GLN	2.4
1	L	331	PHE	2.4
1	K	331	PHE	2.3
1	G	219	THR	2.3
1	B	220	ASP	2.3
1	A	221	LYS	2.2
1	G	216	ASP	2.2
1	E	100	GLN	2.2
1	L	238	TYR	2.2
1	G	218	GLY	2.2
1	H	125	LYS	2.2
1	H	6	GLN	2.2
1	G	98	ARG	2.2
1	E	238	TYR	2.1
1	G	13	LYS	2.1
1	L	100	GLN	2.1
1	B	238	TYR	2.1
1	E	13	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	283	LYS	2.1
1	L	98	ARG	2.1
1	E	99	GLN	2.1
1	K	220	ASP	2.1
1	G	238	TYR	2.0
1	H	126	TYR	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](#)

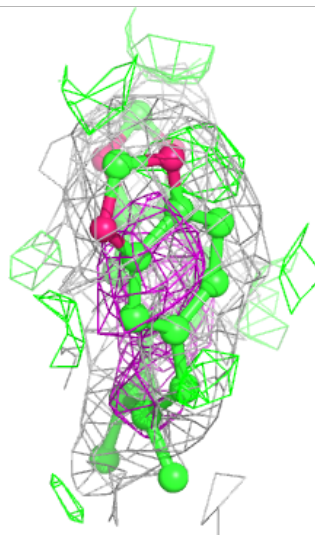
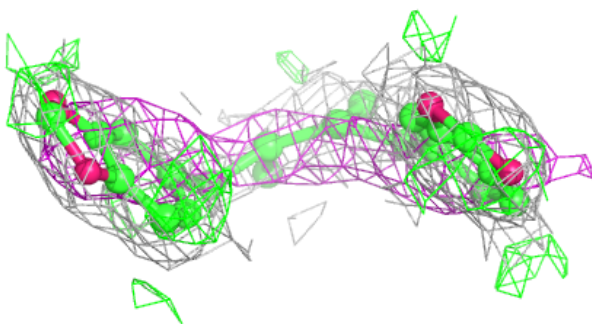
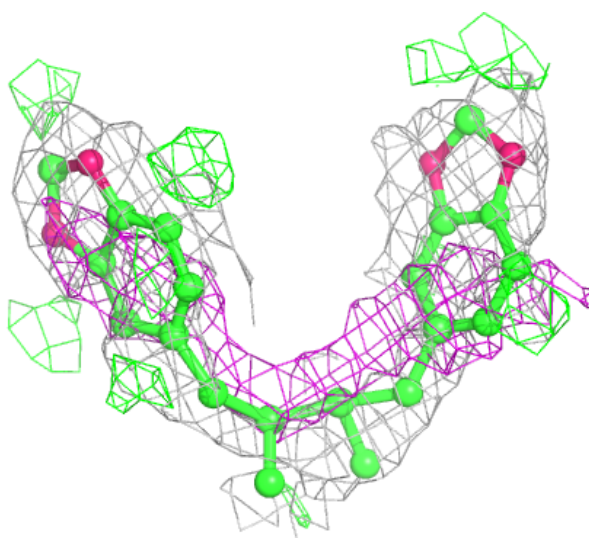
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	9G9	C	401	24/24	0.75	0.33	49,58,62,67	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around 9G9 C 401:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.