



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 02:37 am BST

PDB ID : 3I5V
Title : Crystal structure of beta toxin 275-280 from Staphylococcus aureus
Authors : Huseby, M.; Shi, K.; Kruse, A.C.; Ohlendorf, D.H.
Deposited on : 2009-07-06
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

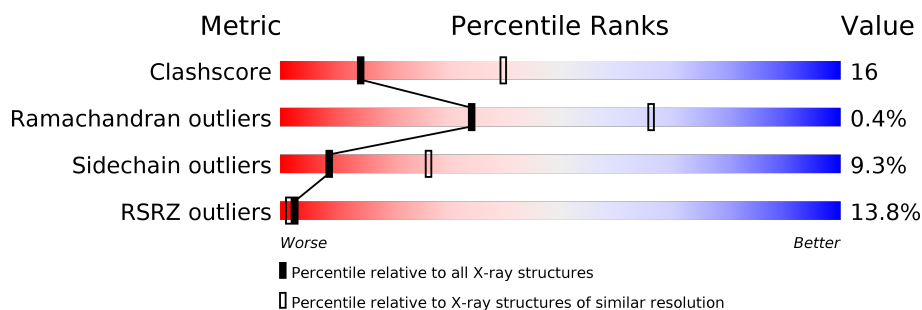
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	313	<div> <div>14%</div> <div> <div></div> <div>62%</div> <div>25%</div> <div>5%</div> <div>9%</div> </div> </div>
1	B	313	<div> <div>13%</div> <div> <div></div> <div>62%</div> <div>24%</div> <div>•</div> <div>9%</div> </div> </div>
1	C	313	<div> <div>14%</div> <div> <div></div> <div>63%</div> <div>23%</div> <div>•</div> <div>9%</div> </div> </div>
1	D	313	<div> <div>10%</div> <div> <div></div> <div>61%</div> <div>25%</div> <div>5%</div> <div>9%</div> </div> </div>

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	DGA	D	401	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9255 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 Beta-hemolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2283	1443	388	447	5			
1	B	285	Total	C	N	O	S	0	0	0
			2283	1443	388	447	5			
1	C	284	Total	C	N	O	S	0	0	0
			2273	1437	386	445	5			
1	D	285	Total	C	N	O	S	0	0	0
			2283	1443	388	447	5			

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP A7LAI8
A	-18	ARG	-	EXPRESSION TAG	UNP A7LAI8
A	-17	SER	-	EXPRESSION TAG	UNP A7LAI8
A	-16	SER	-	EXPRESSION TAG	UNP A7LAI8
A	-15	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	-14	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	-13	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	-12	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	-11	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	-10	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	-9	SER	-	EXPRESSION TAG	UNP A7LAI8
A	-8	SER	-	EXPRESSION TAG	UNP A7LAI8
A	-7	GLY	-	EXPRESSION TAG	UNP A7LAI8
A	-6	LEU	-	EXPRESSION TAG	UNP A7LAI8
A	-5	VAL	-	EXPRESSION TAG	UNP A7LAI8
A	-4	PRO	-	EXPRESSION TAG	UNP A7LAI8
A	-3	ARG	-	EXPRESSION TAG	UNP A7LAI8
A	-2	GLY	-	EXPRESSION TAG	UNP A7LAI8
A	-1	SER	-	EXPRESSION TAG	UNP A7LAI8
A	0	HIS	-	EXPRESSION TAG	UNP A7LAI8
A	1	MET	-	EXPRESSION TAG	UNP A7LAI8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	TYR	DELETION	UNP A7LAI8
A	?	-	ALA	DELETION	UNP A7LAI8
A	?	-	PHE	DELETION	UNP A7LAI8
A	?	-	PRO	DELETION	UNP A7LAI8
A	?	-	TYR	DELETION	UNP A7LAI8
A	?	-	TYR	DELETION	UNP A7LAI8
A	275	ASP	-	SEE REMARK 999	UNP A7LAI8
A	276	GLY	-	SEE REMARK 999	UNP A7LAI8
B	-19	MET	-	EXPRESSION TAG	UNP A7LAI8
B	-18	ARG	-	EXPRESSION TAG	UNP A7LAI8
B	-17	SER	-	EXPRESSION TAG	UNP A7LAI8
B	-16	SER	-	EXPRESSION TAG	UNP A7LAI8
B	-15	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	-14	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	-13	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	-12	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	-11	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	-10	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	-9	SER	-	EXPRESSION TAG	UNP A7LAI8
B	-8	SER	-	EXPRESSION TAG	UNP A7LAI8
B	-7	GLY	-	EXPRESSION TAG	UNP A7LAI8
B	-6	LEU	-	EXPRESSION TAG	UNP A7LAI8
B	-5	VAL	-	EXPRESSION TAG	UNP A7LAI8
B	-4	PRO	-	EXPRESSION TAG	UNP A7LAI8
B	-3	ARG	-	EXPRESSION TAG	UNP A7LAI8
B	-2	GLY	-	EXPRESSION TAG	UNP A7LAI8
B	-1	SER	-	EXPRESSION TAG	UNP A7LAI8
B	0	HIS	-	EXPRESSION TAG	UNP A7LAI8
B	1	MET	-	EXPRESSION TAG	UNP A7LAI8
B	?	-	TYR	DELETION	UNP A7LAI8
B	?	-	ALA	DELETION	UNP A7LAI8
B	?	-	PHE	DELETION	UNP A7LAI8
B	?	-	PRO	DELETION	UNP A7LAI8
B	?	-	TYR	DELETION	UNP A7LAI8
B	?	-	TYR	DELETION	UNP A7LAI8
B	275	ASP	-	SEE REMARK 999	UNP A7LAI8
B	276	GLY	-	SEE REMARK 999	UNP A7LAI8
C	-19	MET	-	EXPRESSION TAG	UNP A7LAI8
C	-18	ARG	-	EXPRESSION TAG	UNP A7LAI8
C	-17	SER	-	EXPRESSION TAG	UNP A7LAI8
C	-16	SER	-	EXPRESSION TAG	UNP A7LAI8
C	-15	HIS	-	EXPRESSION TAG	UNP A7LAI8

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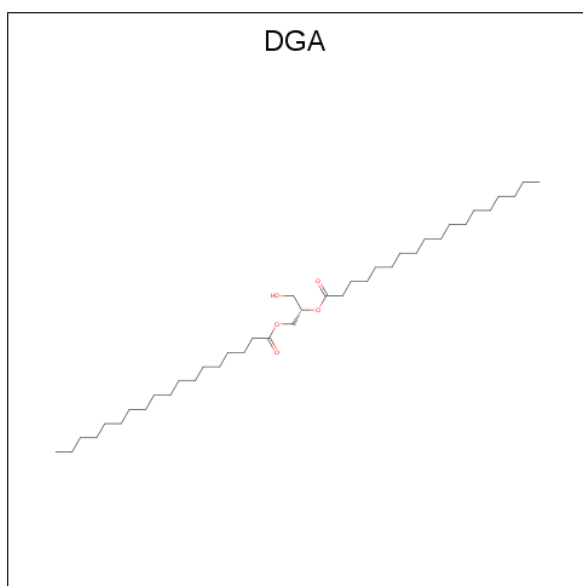
Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	HIS	-	EXPRESSION TAG	UNP A7LAI8
C	-13	HIS	-	EXPRESSION TAG	UNP A7LAI8
C	-12	HIS	-	EXPRESSION TAG	UNP A7LAI8
C	-11	HIS	-	EXPRESSION TAG	UNP A7LAI8
C	-10	HIS	-	EXPRESSION TAG	UNP A7LAI8
C	-9	SER	-	EXPRESSION TAG	UNP A7LAI8
C	-8	SER	-	EXPRESSION TAG	UNP A7LAI8
C	-7	GLY	-	EXPRESSION TAG	UNP A7LAI8
C	-6	LEU	-	EXPRESSION TAG	UNP A7LAI8
C	-5	VAL	-	EXPRESSION TAG	UNP A7LAI8
C	-4	PRO	-	EXPRESSION TAG	UNP A7LAI8
C	-3	ARG	-	EXPRESSION TAG	UNP A7LAI8
C	-2	GLY	-	EXPRESSION TAG	UNP A7LAI8
C	-1	SER	-	EXPRESSION TAG	UNP A7LAI8
C	0	HIS	-	EXPRESSION TAG	UNP A7LAI8
C	1	MET	-	EXPRESSION TAG	UNP A7LAI8
C	?	-	TYR	DELETION	UNP A7LAI8
C	?	-	ALA	DELETION	UNP A7LAI8
C	?	-	PHE	DELETION	UNP A7LAI8
C	?	-	PRO	DELETION	UNP A7LAI8
C	?	-	TYR	DELETION	UNP A7LAI8
C	?	-	TYR	DELETION	UNP A7LAI8
C	275	ASP	-	SEE REMARK 999	UNP A7LAI8
C	276	GLY	-	SEE REMARK 999	UNP A7LAI8
D	-19	MET	-	EXPRESSION TAG	UNP A7LAI8
D	-18	ARG	-	EXPRESSION TAG	UNP A7LAI8
D	-17	SER	-	EXPRESSION TAG	UNP A7LAI8
D	-16	SER	-	EXPRESSION TAG	UNP A7LAI8
D	-15	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	-14	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	-13	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	-12	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	-11	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	-10	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	-9	SER	-	EXPRESSION TAG	UNP A7LAI8
D	-8	SER	-	EXPRESSION TAG	UNP A7LAI8
D	-7	GLY	-	EXPRESSION TAG	UNP A7LAI8
D	-6	LEU	-	EXPRESSION TAG	UNP A7LAI8
D	-5	VAL	-	EXPRESSION TAG	UNP A7LAI8
D	-4	PRO	-	EXPRESSION TAG	UNP A7LAI8
D	-3	ARG	-	EXPRESSION TAG	UNP A7LAI8
D	-2	GLY	-	EXPRESSION TAG	UNP A7LAI8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	SER	-	EXPRESSION TAG	UNP A7LAI8
D	0	HIS	-	EXPRESSION TAG	UNP A7LAI8
D	1	MET	-	EXPRESSION TAG	UNP A7LAI8
D	?	-	TYR	DELETION	UNP A7LAI8
D	?	-	ALA	DELETION	UNP A7LAI8
D	?	-	PHE	DELETION	UNP A7LAI8
D	?	-	PRO	DELETION	UNP A7LAI8
D	?	-	TYR	DELETION	UNP A7LAI8
D	?	-	TYR	DELETION	UNP A7LAI8
D	275	ASP	-	SEE REMARK 999	UNP A7LAI8
D	276	GLY	-	SEE REMARK 999	UNP A7LAI8

- Molecule 2 is DIACYL GLYCEROL (three-letter code: DGA) (formula: $C_{39}H_{76}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			17	13	4		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	33	Total	O	0	0
			33	33		
3	B	26	Total	O	0	0
			26	26		
3	C	31	Total	O	0	0
			31	31		

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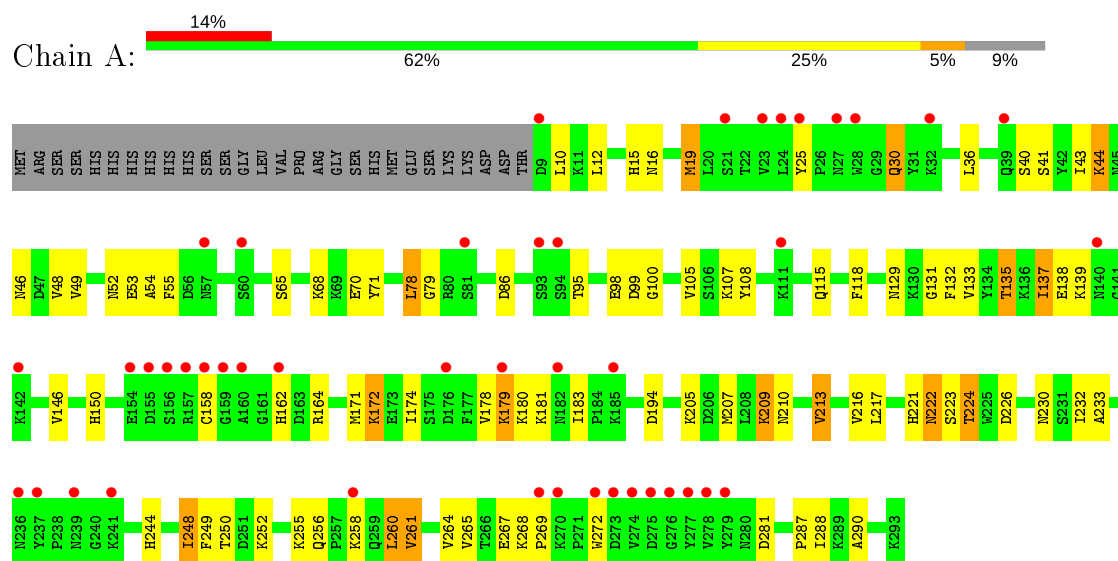
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	26	Total	O	0	0
			26	26		

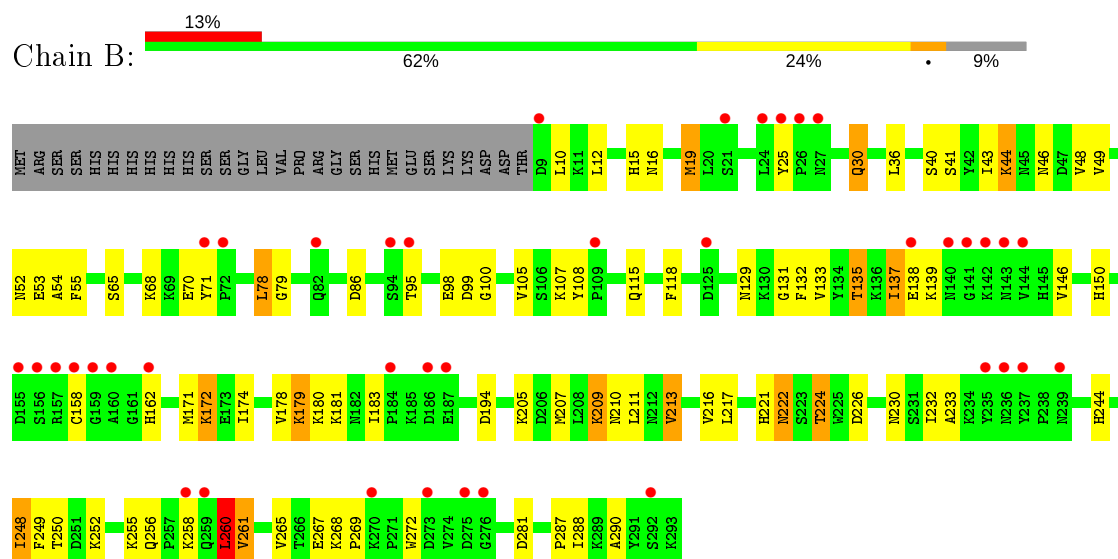
3 Residue-property plots [i](#)

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.

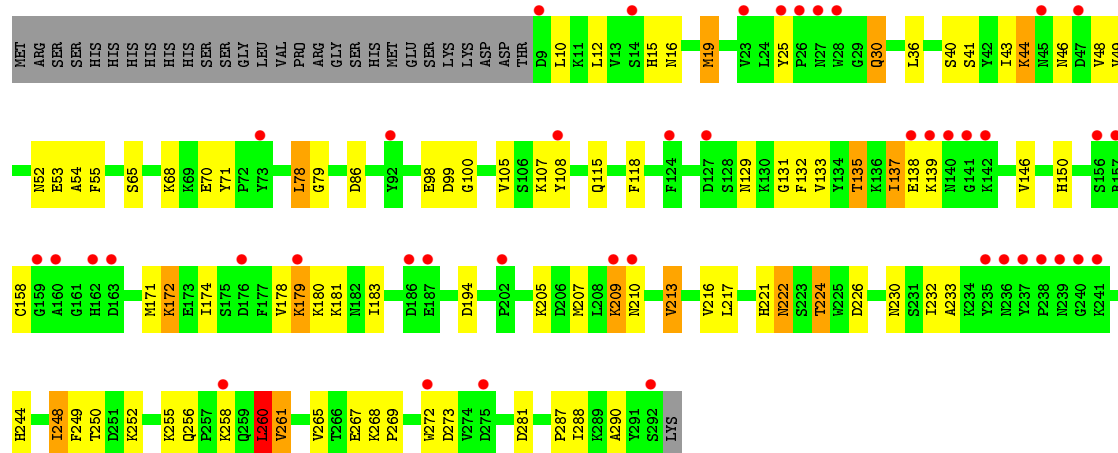
• Molecule 1: Beta-hemolysin



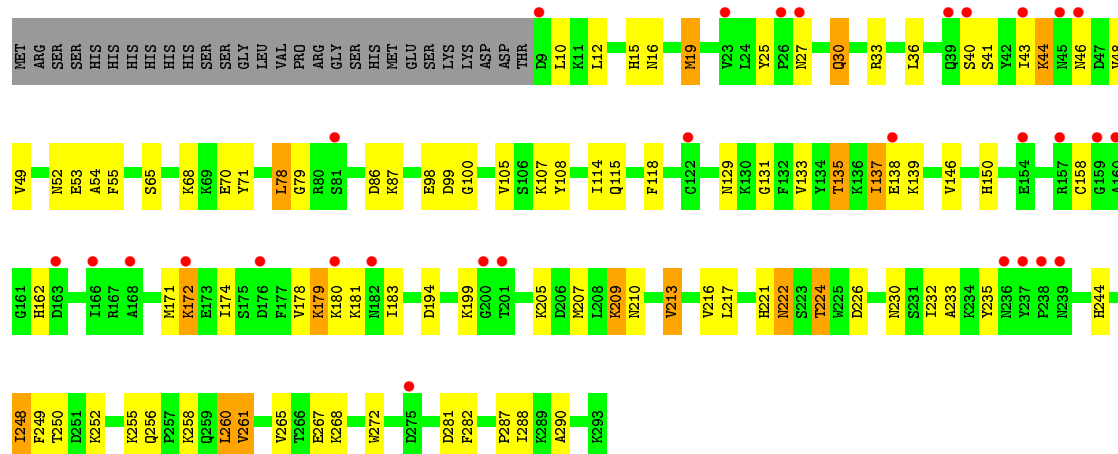
• Molecule 1: Beta-hemolysin



• Molecule 1: Beta-hemolysin



• Molecule 1: Beta-hemolysin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.18Å 69.05Å 75.65Å 93.06° 94.60° 92.13°	Depositor
Resolution (Å)	38.40 – 2.80 38.40 – 2.79	Depositor EDS
% Data completeness (in resolution range)	95.0 (38.40-2.80) 77.7 (38.40-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.48 (at 2.77Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.214 , 0.267 0.232 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	1.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 68.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9255	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.25	0/2339	0.45	2/3162 (0.1%)
1	B	0.25	0/2339	0.45	2/3162 (0.1%)
1	C	0.25	0/2329	0.45	2/3151 (0.1%)
1	D	0.25	0/2339	0.45	2/3162 (0.1%)
All	All	0.25	0/9346	0.45	8/12637 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	260	LEU	CB-CG-CD2	5.96	121.14	111.00
1	B	260	LEU	CB-CG-CD2	5.95	121.12	111.00
1	D	260	LEU	CB-CG-CD2	5.66	120.62	111.00
1	A	260	LEU	CB-CG-CD2	5.64	120.59	111.00
1	D	260	LEU	CB-CG-CD1	5.63	120.57	111.00
1	A	260	LEU	CB-CG-CD1	5.58	120.48	111.00
1	B	260	LEU	CB-CG-CD1	5.55	120.44	111.00
1	C	260	LEU	CB-CG-CD1	5.54	120.41	111.00

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	2283	0	2205	80	0
1	B	2283	0	2205	80	0
1	C	2273	0	2192	63	0
1	D	2283	0	2205	69	0
2	D	17	0	17	7	0
3	A	33	0	0	0	0
3	B	26	0	0	0	0
3	C	31	0	0	3	0
3	D	26	0	0	2	0
All	All	9255	0	8824	279	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 16.

All (279) 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:D:27:ASN:O	2:D:401:DGA:HG12	1.36	1.21
3:C:325:HOH:O	2:D:401:DGA:HB42	1.63	0.98
1:B:30:GLN:HE21	1:B:30:GLN:H	1.13	0.96
1:A:30:GLN:H	1:A:30:GLN:HE21	1.13	0.96
1:D:30:GLN:HE21	1:D:30:GLN:H	1.14	0.95
1:C:30:GLN:HE21	1:C:30:GLN:H	1.14	0.92
1:A:230:ASN:HD22	1:A:233:ALA:H	1.20	0.90
1:D:230:ASN:HD22	1:D:233:ALA:H	1.20	0.90
1:B:230:ASN:HD22	1:B:233:ALA:H	1.20	0.88
3:C:325:HOH:O	2:D:401:DGA:CB4	2.16	0.87
1:C:230:ASN:HD22	1:C:233:ALA:H	1.20	0.86
1:D:108:TYR:HB2	1:D:137:ILE:HD12	1.63	0.81
1:B:108:TYR:HB2	1:B:137:ILE:HD12	1.63	0.81
1:C:78:LEU:HD13	1:C:100:GLY:HA3	1.62	0.81
1:D:27:ASN:O	2:D:401:DGA:CG1	2.25	0.81
1:A:108:TYR:HB2	1:A:137:ILE:HD12	1.63	0.80
1:B:78:LEU:HD13	1:B:100:GLY:HA3	1.63	0.80
1:A:95:THR:HG21	1:B:162:HIS:CD2	2.17	0.80
1:C:108:TYR:HB2	1:C:137:ILE:HD12	1.63	0.79
1:A:78:LEU:HD13	1:A:100:GLY:HA3	1.63	0.79
1:A:30:GLN:HE21	1:A:30:GLN:N	1.82	0.78
1:D:78:LEU:HD13	1:D:100:GLY:HA3	1.62	0.78
1:B:30:GLN:HE21	1:B:30:GLN:N	1.82	0.78
1:A:162:HIS:CD2	1:B:95:THR:HG21	2.18	0.77
1:D:30:GLN:HE21	1:D:30:GLN:N	1.83	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:30:GLN:HE21	1:C:30:GLN:N	1.82	0.76
1:A:162:HIS:CE1	1:B:95:THR:CG2	2.72	0.73
1:A:95:THR:CG2	1:B:162:HIS:CE1	2.73	0.71
1:A:164:ARG:HH12	1:C:273:ASP:HB3	1.55	0.69
1:A:162:HIS:CG	1:B:95:THR:HG21	2.28	0.67
1:A:230:ASN:ND2	1:A:233:ALA:H	1.93	0.66
1:D:230:ASN:ND2	1:D:233:ALA:H	1.94	0.66
1:C:230:ASN:ND2	1:C:233:ALA:H	1.93	0.65
1:C:30:GLN:NE2	1:C:30:GLN:H	1.92	0.65
1:D:235:TYR:OH	2:D:401:DGA:HB22	1.97	0.65
1:B:30:GLN:H	1:B:30:GLN:NE2	1.91	0.63
1:C:230:ASN:HB3	1:C:233:ALA:HB3	1.81	0.63
1:A:95:THR:CG2	1:B:162:HIS:NE2	2.62	0.62
1:D:162:HIS:HE1	3:D:316:HOH:O	1.83	0.62
1:B:213:VAL:HG11	1:B:249:PHE:CG	2.36	0.61
1:B:230:ASN:HB3	1:B:233:ALA:HB3	1.81	0.61
1:B:230:ASN:ND2	1:B:233:ALA:H	1.94	0.61
1:A:230:ASN:HB3	1:A:233:ALA:HB3	1.82	0.60
1:D:213:VAL:HG11	1:D:249:PHE:CG	2.36	0.60
1:C:213:VAL:HG11	1:C:249:PHE:CG	2.36	0.60
1:D:230:ASN:HB3	1:D:233:ALA:HB3	1.82	0.60
1:A:213:VAL:HG11	1:A:249:PHE:CG	2.36	0.60
1:B:209:LYS:N	1:B:209:LYS:HD2	2.17	0.60
1:D:30:GLN:NE2	1:D:30:GLN:H	1.92	0.59
1:D:209:LYS:N	1:D:209:LYS:HD2	2.17	0.59
1:A:95:THR:HG21	1:B:162:HIS:CG	2.36	0.59
1:A:30:GLN:H	1:A:30:GLN:NE2	1.91	0.59
1:C:172:LYS:HE2	1:C:172:LYS:HA	1.85	0.59
1:D:172:LYS:HA	1:D:172:LYS:HE2	1.85	0.59
1:A:209:LYS:HD2	1:A:209:LYS:N	2.17	0.59
1:A:172:LYS:HA	1:A:172:LYS:HE2	1.85	0.59
1:A:12:LEU:HD22	1:A:48:VAL:HB	1.85	0.58
1:D:10:LEU:HB3	1:D:12:LEU:HD21	1.86	0.58
1:A:162:HIS:NE2	1:B:95:THR:CG2	2.66	0.58
1:C:12:LEU:HD22	1:C:48:VAL:HB	1.85	0.58
1:C:209:LYS:HD2	1:C:209:LYS:N	2.17	0.58
1:D:12:LEU:HD22	1:D:48:VAL:HB	1.85	0.58
1:B:172:LYS:HE2	1:B:172:LYS:HA	1.85	0.58
1:A:78:LEU:HD23	1:A:115:GLN:HB2	1.86	0.57
1:B:10:LEU:HB3	1:B:12:LEU:HD21	1.86	0.57
1:A:10:LEU:HB3	1:A:12:LEU:HD21	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:SER:OG	1:A:43:ILE:HG13	2.04	0.57
1:A:171:MET:SD	1:A:207:MET:HG3	2.44	0.57
1:D:78:LEU:HD23	1:D:115:GLN:HB2	1.87	0.57
1:A:78:LEU:CD1	1:A:100:GLY:HA3	2.35	0.57
1:D:40:SER:OG	1:D:43:ILE:HG13	2.05	0.57
1:C:40:SER:OG	1:C:43:ILE:HG13	2.05	0.57
1:B:12:LEU:HD22	1:B:48:VAL:HB	1.86	0.57
1:C:10:LEU:HB3	1:C:12:LEU:HD21	1.86	0.57
1:C:171:MET:SD	1:C:207:MET:HG3	2.45	0.57
1:B:78:LEU:HD23	1:B:115:GLN:HB2	1.86	0.56
1:C:78:LEU:HD23	1:C:115:GLN:HB2	1.87	0.56
1:C:78:LEU:CD1	1:C:100:GLY:HA3	2.35	0.56
1:B:171:MET:SD	1:B:207:MET:HG3	2.45	0.56
1:D:171:MET:SD	1:D:207:MET:HG3	2.46	0.56
1:B:40:SER:OG	1:B:43:ILE:HG13	2.05	0.56
1:B:78:LEU:CD1	1:B:100:GLY:HA3	2.35	0.55
1:B:41:SER:HA	1:B:44:LYS:HE2	1.89	0.55
1:D:41:SER:HA	1:D:44:LYS:HE2	1.89	0.55
1:A:41:SER:HA	1:A:44:LYS:HE2	1.89	0.54
1:D:16:ASN:HA	1:D:52:ASN:HB2	1.90	0.54
1:C:41:SER:HA	1:C:44:LYS:HE2	1.89	0.54
1:A:16:ASN:HA	1:A:52:ASN:HB2	1.90	0.54
1:A:179:LYS:HD2	1:A:180:LYS:N	2.23	0.53
1:D:135:THR:HG22	1:D:146:VAL:HB	1.91	0.53
1:C:179:LYS:HD2	1:C:180:LYS:N	2.24	0.53
1:C:248:ILE:HG23	1:C:288:ILE:HD11	1.90	0.53
1:D:179:LYS:HD2	1:D:180:LYS:N	2.23	0.53
1:C:16:ASN:HA	1:C:52:ASN:HB2	1.90	0.53
1:A:135:THR:HG22	1:A:146:VAL:HB	1.91	0.53
1:A:49:VAL:HB	1:A:105:VAL:HG22	1.91	0.53
1:C:10:LEU:HD13	1:C:12:LEU:HD21	1.90	0.53
1:B:10:LEU:HD13	1:B:12:LEU:HD21	1.91	0.52
1:D:10:LEU:HD13	1:D:12:LEU:HD21	1.91	0.52
1:B:179:LYS:HD2	1:B:180:LYS:N	2.23	0.52
1:C:19:MET:HB2	1:C:54:ALA:HA	1.92	0.52
1:A:10:LEU:HD13	1:A:12:LEU:HD21	1.91	0.52
1:A:248:ILE:HG23	1:A:288:ILE:HD11	1.91	0.52
1:B:16:ASN:HA	1:B:52:ASN:HB2	1.90	0.52
1:B:248:ILE:HG23	1:B:288:ILE:HD11	1.91	0.52
1:D:248:ILE:HG23	1:D:288:ILE:HD11	1.91	0.52
1:B:118:PHE:CE2	1:B:131:GLY:HA2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:THR:HG22	1:C:146:VAL:HB	1.91	0.52
1:A:118:PHE:CE2	1:A:131:GLY:HA2	2.45	0.52
1:D:49:VAL:HB	1:D:105:VAL:HG22	1.92	0.52
1:D:108:TYR:CZ	1:D:139:LYS:HD2	2.45	0.52
1:A:108:TYR:CZ	1:A:139:LYS:HD2	2.45	0.51
1:A:15:HIS:CE1	1:A:287:PRO:HG3	2.45	0.51
1:B:10:LEU:HB3	1:B:12:LEU:CD2	2.40	0.51
1:B:49:VAL:HB	1:B:105:VAL:HG22	1.91	0.51
1:D:15:HIS:CE1	1:D:287:PRO:HG3	2.46	0.51
1:C:118:PHE:CE2	1:C:131:GLY:HA2	2.46	0.51
1:D:19:MET:HB2	1:D:54:ALA:HA	1.92	0.51
1:A:19:MET:HB2	1:A:54:ALA:HA	1.92	0.51
1:C:78:LEU:HD22	1:C:79:GLY:N	2.26	0.51
1:D:10:LEU:HB3	1:D:12:LEU:CD2	2.40	0.51
1:A:78:LEU:HD22	1:A:79:GLY:N	2.25	0.51
1:A:10:LEU:HB3	1:A:12:LEU:CD2	2.40	0.51
1:C:10:LEU:HB3	1:C:12:LEU:CD2	2.40	0.51
1:D:78:LEU:CD1	1:D:100:GLY:HA3	2.34	0.51
1:D:118:PHE:CE2	1:D:131:GLY:HA2	2.46	0.51
1:A:162:HIS:CD2	1:B:95:THR:CG2	2.92	0.51
1:B:135:THR:HG22	1:B:146:VAL:HB	1.92	0.50
1:B:19:MET:HB2	1:B:54:ALA:HA	1.92	0.50
1:B:78:LEU:HD22	1:B:79:GLY:N	2.26	0.50
1:C:49:VAL:HB	1:C:105:VAL:HG22	1.92	0.50
1:C:108:TYR:CZ	1:C:139:LYS:HD2	2.45	0.50
1:D:78:LEU:HD22	1:D:79:GLY:N	2.26	0.50
1:B:15:HIS:CE1	1:B:287:PRO:HG3	2.45	0.50
1:C:15:HIS:CE1	1:C:287:PRO:HG3	2.46	0.50
1:B:108:TYR:CZ	1:B:139:LYS:HD2	2.46	0.50
1:A:95:THR:HG22	1:B:162:HIS:CE1	2.47	0.50
1:D:137:ILE:HG13	1:D:138:GLU:N	2.26	0.50
1:A:137:ILE:HG13	1:A:138:GLU:N	2.26	0.50
1:A:162:HIS:CE1	1:B:95:THR:HG22	2.47	0.49
1:C:150:HIS:CD2	1:C:194:ASP:HB3	2.47	0.49
1:A:150:HIS:CD2	1:A:194:ASP:HB3	2.47	0.49
1:A:95:THR:HG21	1:B:162:HIS:NE2	2.27	0.49
1:B:137:ILE:HG13	1:B:138:GLU:N	2.27	0.49
1:B:118:PHE:CD2	1:B:131:GLY:HA2	2.48	0.49
1:B:150:HIS:CD2	1:B:194:ASP:HB3	2.47	0.49
1:C:118:PHE:CD2	1:C:131:GLY:HA2	2.48	0.49
1:C:232:ILE:HD11	1:C:272:TRP:HH2	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:PHE:CD2	1:A:131:GLY:HA2	2.47	0.48
1:D:118:PHE:CD2	1:D:131:GLY:HA2	2.48	0.48
1:D:150:HIS:CD2	1:D:194:ASP:HB3	2.48	0.48
1:B:268:LYS:HE2	1:B:281:ASP:OD2	2.14	0.48
1:C:43:ILE:HD12	1:C:44:LYS:N	2.28	0.48
1:D:162:HIS:CE1	3:D:316:HOH:O	2.61	0.48
1:D:268:LYS:HE2	1:D:281:ASP:OD2	2.14	0.48
1:A:43:ILE:HD12	1:A:44:LYS:N	2.29	0.48
1:B:43:ILE:HD12	1:B:44:LYS:N	2.29	0.48
1:B:65:SER:HA	1:B:68:LYS:HB2	1.96	0.48
1:C:137:ILE:HG13	1:C:138:GLU:N	2.27	0.48
1:A:268:LYS:HE2	1:A:281:ASP:OD2	2.14	0.47
1:B:49:VAL:HB	1:B:105:VAL:CG2	2.44	0.47
1:C:49:VAL:HB	1:C:105:VAL:CG2	2.44	0.47
1:D:49:VAL:HB	1:D:105:VAL:CG2	2.44	0.47
1:A:78:LEU:HB2	1:A:133:VAL:CG2	2.45	0.47
1:D:222:ASN:HD22	1:D:222:ASN:N	2.12	0.47
1:D:43:ILE:HD12	1:D:44:LYS:N	2.29	0.47
1:C:65:SER:HA	1:C:68:LYS:HB2	1.96	0.47
1:D:78:LEU:HB2	1:D:133:VAL:CG2	2.45	0.47
1:B:232:ILE:HD11	1:B:272:TRP:HH2	1.79	0.47
1:A:49:VAL:HB	1:A:105:VAL:CG2	2.44	0.47
1:B:78:LEU:HB2	1:B:133:VAL:CG2	2.44	0.47
1:C:268:LYS:HE2	1:C:281:ASP:OD2	2.14	0.47
1:D:248:ILE:HG12	1:D:290:ALA:HB2	1.97	0.47
1:A:248:ILE:HG12	1:A:290:ALA:HB2	1.97	0.47
1:B:41:SER:O	1:B:44:LYS:HG2	2.15	0.47
1:C:41:SER:O	1:C:44:LYS:HG2	2.15	0.47
1:B:265:VAL:HG12	1:B:267:GLU:H	1.80	0.47
1:D:41:SER:O	1:D:44:LYS:HG2	2.15	0.47
1:C:248:ILE:HG12	1:C:290:ALA:HB2	1.97	0.46
1:D:232:ILE:HD11	1:D:272:TRP:HH2	1.79	0.46
1:A:232:ILE:HD11	1:A:272:TRP:HH2	1.79	0.46
1:A:265:VAL:HG12	1:A:267:GLU:H	1.80	0.46
1:A:95:THR:HG21	1:B:162:HIS:CE1	2.51	0.46
1:B:221:HIS:HE1	1:B:226:ASP:OD1	1.99	0.46
1:D:65:SER:HA	1:D:68:LYS:HB2	1.96	0.46
1:A:216:VAL:HG12	1:A:217:LEU:O	2.16	0.46
1:B:248:ILE:HG12	1:B:290:ALA:HB2	1.97	0.46
2:D:401:DGA:HG11	2:D:401:DGA:OB1	2.15	0.46
1:A:65:SER:HA	1:A:68:LYS:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:265:VAL:HG12	1:C:267:GLU:H	1.80	0.46
1:B:108:TYR:CE1	1:B:139:LYS:HD2	2.51	0.46
1:C:216:VAL:HG12	1:C:217:LEU:O	2.16	0.46
1:D:216:VAL:HG12	1:D:217:LEU:O	2.15	0.46
1:D:222:ASN:HD22	1:D:222:ASN:H	1.64	0.46
1:A:222:ASN:HD22	1:A:222:ASN:N	2.13	0.46
1:A:41:SER:O	1:A:44:LYS:HG2	2.16	0.46
1:A:108:TYR:CE1	1:A:139:LYS:HD2	2.51	0.45
1:C:108:TYR:CE1	1:C:139:LYS:HD2	2.51	0.45
1:C:261:VAL:O	1:C:290:ALA:HA	2.17	0.45
1:C:78:LEU:HB2	1:C:133:VAL:CG2	2.45	0.45
1:C:222:ASN:HD22	1:C:222:ASN:N	2.13	0.45
1:A:222:ASN:HD22	1:A:222:ASN:H	1.65	0.45
1:B:216:VAL:HG12	1:B:217:LEU:O	2.16	0.45
1:C:221:HIS:HE1	1:C:226:ASP:OD1	1.99	0.45
1:C:53:GLU:HA	1:C:55:PHE:CE1	2.52	0.45
1:D:108:TYR:CE1	1:D:139:LYS:HD2	2.51	0.45
1:D:261:VAL:O	1:D:290:ALA:HA	2.17	0.45
1:D:265:VAL:HG12	1:D:267:GLU:H	1.80	0.45
1:B:53:GLU:HA	1:B:55:PHE:CE1	2.52	0.45
1:B:222:ASN:HD22	1:B:222:ASN:N	2.14	0.45
1:B:261:VAL:O	1:B:290:ALA:HA	2.17	0.44
1:A:222:ASN:HB3	1:D:199:LYS:HD2	1.99	0.44
1:A:53:GLU:HA	1:A:55:PHE:CE1	2.52	0.44
1:B:222:ASN:HD22	1:B:222:ASN:H	1.65	0.44
1:A:261:VAL:O	1:A:290:ALA:HA	2.17	0.44
1:A:46:ASN:HB2	1:A:71:TYR:OH	2.18	0.44
1:D:46:ASN:HB2	1:D:71:TYR:OH	2.18	0.44
1:B:46:ASN:HB2	1:B:71:TYR:OH	2.18	0.44
1:C:46:ASN:HB2	1:C:71:TYR:OH	2.17	0.44
3:C:325:HOH:O	2:D:401:DGA:HB41	2.03	0.44
1:D:53:GLU:HA	1:D:55:PHE:CE1	2.52	0.44
1:A:221:HIS:HE1	1:A:226:ASP:OD1	2.00	0.43
1:C:222:ASN:HD22	1:C:222:ASN:H	1.65	0.43
1:A:162:HIS:CE1	1:B:95:THR:HG21	2.54	0.43
1:D:135:THR:CG2	1:D:146:VAL:HB	2.49	0.43
1:D:221:HIS:HE1	1:D:226:ASP:OD1	2.00	0.43
1:B:70:GLU:O	1:B:107:LYS:HG2	2.19	0.43
1:A:174:ILE:O	1:A:178:VAL:HG23	2.19	0.43
1:A:98:GLU:HG3	1:A:99:ASP:N	2.33	0.43
1:B:172:LYS:HE3	1:B:210:ASN:HA	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:GLU:HG3	1:B:99:ASP:N	2.33	0.43
1:C:224:THR:HB	1:C:244:HIS:HA	2.01	0.43
1:D:216:VAL:HG13	1:D:260:LEU:O	2.19	0.43
1:C:70:GLU:O	1:C:107:LYS:HG2	2.19	0.42
1:D:174:ILE:O	1:D:178:VAL:HG23	2.19	0.42
1:A:216:VAL:HG13	1:A:260:LEU:O	2.19	0.42
1:A:70:GLU:O	1:A:107:LYS:HG2	2.19	0.42
1:B:224:THR:HB	1:B:244:HIS:HA	2.01	0.42
1:C:135:THR:CG2	1:C:146:VAL:HB	2.49	0.42
1:A:135:THR:CG2	1:A:146:VAL:HB	2.49	0.42
1:A:224:THR:HB	1:A:244:HIS:HA	2.00	0.42
1:B:174:ILE:O	1:B:178:VAL:HG23	2.20	0.42
1:B:181:LYS:HB3	1:B:183:ILE:HG13	2.02	0.42
1:B:216:VAL:HG13	1:B:260:LEU:O	2.20	0.42
1:C:174:ILE:O	1:C:178:VAL:HG23	2.19	0.42
1:C:98:GLU:HG3	1:C:99:ASP:N	2.34	0.42
1:D:70:GLU:O	1:D:107:LYS:HG2	2.19	0.42
1:D:172:LYS:HE3	1:D:210:ASN:HA	2.01	0.42
1:C:172:LYS:HE3	1:C:210:ASN:HA	2.01	0.42
1:D:181:LYS:HB3	1:D:183:ILE:HG13	2.02	0.42
1:A:181:LYS:HB3	1:A:183:ILE:HG13	2.02	0.42
1:C:181:LYS:HB3	1:C:183:ILE:HG13	2.02	0.42
1:C:268:LYS:HA	1:C:269:PRO:HD3	1.84	0.42
1:A:162:HIS:NE2	1:B:95:THR:HG23	2.33	0.41
1:B:135:THR:CG2	1:B:146:VAL:HB	2.49	0.41
1:B:213:VAL:HG22	1:B:250:THR:O	2.20	0.41
1:D:98:GLU:HG3	1:D:99:ASP:N	2.34	0.41
1:C:55:PHE:HB3	1:C:98:GLU:CD	2.41	0.41
1:D:213:VAL:HG22	1:D:250:THR:O	2.20	0.41
1:A:172:LYS:HE3	1:A:210:ASN:HA	2.01	0.41
1:B:12:LEU:N	1:B:12:LEU:HD23	2.36	0.41
1:B:268:LYS:HA	1:B:269:PRO:HD3	1.84	0.41
1:C:216:VAL:HG13	1:C:260:LEU:O	2.20	0.41
1:C:213:VAL:HG22	1:C:250:THR:O	2.20	0.41
1:D:224:THR:HB	1:D:244:HIS:HA	2.01	0.41
1:A:268:LYS:HA	1:A:269:PRO:HD3	1.84	0.41
1:A:15:HIS:ND1	1:A:287:PRO:HG3	2.36	0.41
1:B:15:HIS:ND1	1:B:287:PRO:HG3	2.36	0.41
1:D:55:PHE:HB3	1:D:98:GLU:CD	2.41	0.41
1:A:132:PHE:CD1	1:A:174:ILE:HG12	2.57	0.40
1:B:211:LEU:HA	1:B:211:LEU:HD23	1.96	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:87:LYS:HB3	1:D:114:ILE:HG13	2.03	0.40
1:B:132:PHE:CD1	1:B:174:ILE:HG12	2.57	0.40
1:D:33:ARG:HD2	1:D:282:PHE:O	2.22	0.40
1:A:223:SER:O	1:A:264:VAL:HG11	2.22	0.40
1:B:55:PHE:HB3	1:B:98:GLU:CD	2.41	0.40
1:A:213:VAL:HG22	1:A:250:THR:O	2.20	0.40
1:C:132:PHE:CD1	1:C:174:ILE:HG12	2.57	0.40
1:D:12:LEU:N	1:D:12:LEU:HD23	2.36	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	283/313 (90%)	260 (92%)	22 (8%)	1 (0%)	34	66
1	B	283/313 (90%)	260 (92%)	22 (8%)	1 (0%)	34	66
1	C	282/313 (90%)	259 (92%)	22 (8%)	1 (0%)	34	66
1	D	283/313 (90%)	260 (92%)	22 (8%)	1 (0%)	34	66
All	All	1131/1252 (90%)	1039 (92%)	88 (8%)	4 (0%)	34	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	LYS
1	B	258	LYS
1	C	258	LYS
1	D	258	LYS

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	254/280 (91%)	231 (91%)	23 (9%)	9	27
1	B	254/280 (91%)	230 (91%)	24 (9%)	8	26
1	C	253/280 (90%)	229 (90%)	24 (10%)	8	25
1	D	254/280 (91%)	231 (91%)	23 (9%)	9	27
All	All	1015/1120 (91%)	921 (91%)	94 (9%)	9	26

All (94) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	MET
1	A	25	TYR
1	A	30	GLN
1	A	36	LEU
1	A	44	LYS
1	A	78	LEU
1	A	86	ASP
1	A	129	ASN
1	A	135	THR
1	A	137	ILE
1	A	158	CYS
1	A	172	LYS
1	A	179	LYS
1	A	205	LYS
1	A	209	LYS
1	A	213	VAL
1	A	222	ASN
1	A	224	THR
1	A	248	ILE
1	A	252	LYS
1	A	255	LYS
1	A	256	GLN
1	A	261	VAL
1	B	19	MET

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Mol	Chain	Res	Type
1	B	25	TYR
1	B	30	GLN
1	B	36	LEU
1	B	44	LYS
1	B	78	LEU
1	B	86	ASP
1	B	129	ASN
1	B	135	THR
1	B	137	ILE
1	B	158	CYS
1	B	172	LYS
1	B	179	LYS
1	B	205	LYS
1	B	209	LYS
1	B	213	VAL
1	B	222	ASN
1	B	224	THR
1	B	248	ILE
1	B	252	LYS
1	B	255	LYS
1	B	256	GLN
1	B	260	LEU
1	B	261	VAL
1	C	19	MET
1	C	25	TYR
1	C	30	GLN
1	C	36	LEU
1	C	44	LYS
1	C	78	LEU
1	C	86	ASP
1	C	129	ASN
1	C	135	THR
1	C	137	ILE
1	C	158	CYS
1	C	172	LYS
1	C	179	LYS
1	C	205	LYS
1	C	209	LYS
1	C	213	VAL
1	C	222	ASN
1	C	224	THR
1	C	248	ILE

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Mol	Chain	Res	Type
1	C	252	LYS
1	C	255	LYS
1	C	256	GLN
1	C	260	LEU
1	C	261	VAL
1	D	19	MET
1	D	25	TYR
1	D	30	GLN
1	D	36	LEU
1	D	44	LYS
1	D	78	LEU
1	D	86	ASP
1	D	129	ASN
1	D	135	THR
1	D	137	ILE
1	D	158	CYS
1	D	172	LYS
1	D	179	LYS
1	D	205	LYS
1	D	209	LYS
1	D	213	VAL
1	D	222	ASN
1	D	224	THR
1	D	248	ILE
1	D	252	LYS
1	D	255	LYS
1	D	256	GLN
1	D	261	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (54) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	30	GLN
1	A	46	ASN
1	A	52	ASN
1	A	66	ASN
1	A	129	ASN
1	A	140	ASN
1	A	150	HIS
1	A	210	ASN
1	A	212	ASN

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Mol	Chain	Res	Type
1	A	221	HIS
1	A	222	ASN
1	A	230	ASN
1	A	259	GLN
1	B	27	ASN
1	B	30	GLN
1	B	46	ASN
1	B	52	ASN
1	B	66	ASN
1	B	129	ASN
1	B	140	ASN
1	B	150	HIS
1	B	210	ASN
1	B	212	ASN
1	B	221	HIS
1	B	222	ASN
1	B	230	ASN
1	B	259	GLN
1	C	30	GLN
1	C	46	ASN
1	C	52	ASN
1	C	66	ASN
1	C	129	ASN
1	C	140	ASN
1	C	150	HIS
1	C	210	ASN
1	C	212	ASN
1	C	221	HIS
1	C	222	ASN
1	C	230	ASN
1	C	259	GLN
1	D	27	ASN
1	D	30	GLN
1	D	46	ASN
1	D	52	ASN
1	D	66	ASN
1	D	129	ASN
1	D	140	ASN
1	D	150	HIS
1	D	210	ASN
1	D	212	ASN
1	D	221	HIS

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Mol	Chain	Res	Type
1	D	222	ASN
1	D	230	ASN

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	DGA	D	401	-	16,16,43	1.24	1 (6%)	17,17,45	1.92	5 (29%)

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	DGA	D	401	-	1/1/2/3	13/17/17/45	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	DGA	OG2-CB1	4.43	1.46	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	DGA	OG2-CB1-CB2	5.61	123.60	111.50
2	D	401	DGA	OB1-CB1-CB2	-2.43	114.27	123.73
2	D	401	DGA	OXT-CG3-CG2	2.17	117.54	111.78
2	D	401	DGA	CA3-CA2-CA1	-2.11	104.14	113.49
2	D	401	DGA	CA1-OG1-CG1	2.00	122.56	113.61

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	401	DGA	CG2

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	401	DGA	CB2-CB1-OG2-CG2
2	D	401	DGA	CG2-CG1-OG1-CA1
2	D	401	DGA	CG1-CG2-CG3-OXT
2	D	401	DGA	OG2-CG2-CG3-OXT
2	D	401	DGA	OB1-CB1-OG2-CG2
2	D	401	DGA	OG1-CA1-CA2-CA3
2	D	401	DGA	CB2-CB3-CB4-CB5
2	D	401	DGA	OG1-CG1-CG2-CG3
2	D	401	DGA	CA2-CA3-CA4-CA5
2	D	401	DGA	CA2-CA1-OG1-CG1
2	D	401	DGA	OG1-CG1-CG2-OG2
2	D	401	DGA	CG1-CG2-OG2-CB1
2	D	401	DGA	CA1-CA2-CA3-CA4

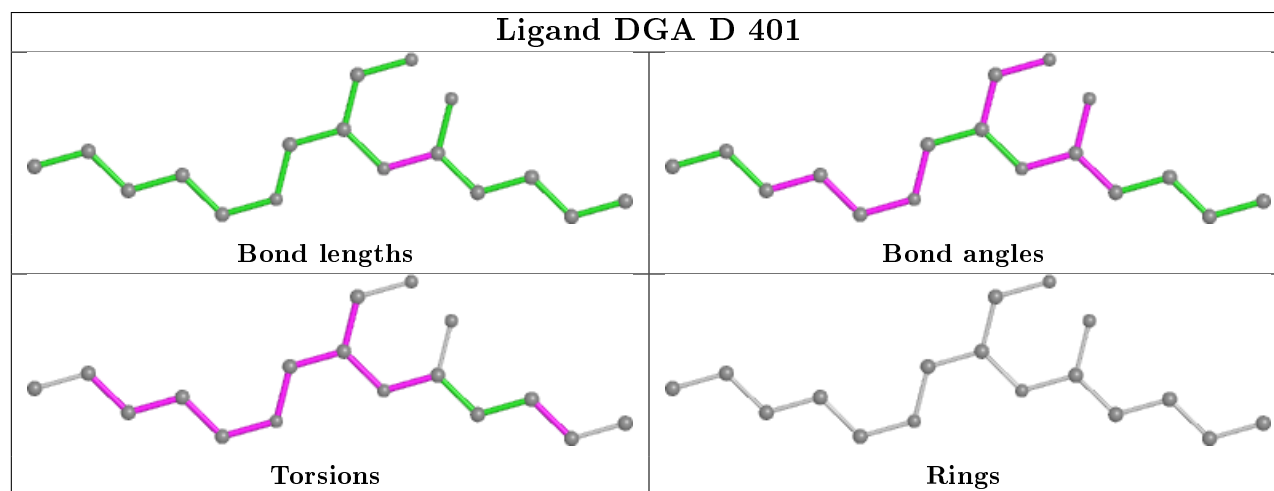
There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	DGA	7	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, 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	285/313 (91%)	1.06	44 (15%)	2 1	51, 82, 141, 251	0
1	B	285/313 (91%)	0.89	40 (14%)	2 1	43, 80, 136, 268	0
1	C	284/313 (90%)	0.86	43 (15%)	2 1	45, 77, 132, 176	0
1	D	285/313 (91%)	0.79	30 (10%)	6 3	44, 78, 129, 186	0
All	All	1139/1252 (90%)	0.90	157 (13%)	2 1	43, 79, 136, 268	0

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	239	ASN	11.4
1	A	278	VAL	11.2
1	C	239	ASN	8.7
1	C	26	PRO	7.9
1	A	155	ASP	7.2
1	B	157	ARG	6.5
1	A	9	ASP	6.5
1	C	156	SER	6.5
1	A	275	ASP	6.4
1	A	94	SER	6.3
1	C	237	TYR	6.1
1	C	157	ARG	6.0
1	D	157	ARG	5.9
1	D	238	PRO	5.6
1	A	274	VAL	5.0
1	A	237	TYR	4.9
1	B	276	GLY	4.9
1	B	160	ALA	4.7
1	A	162	HIS	4.7
1	B	25	TYR	4.7
1	A	273	ASP	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	159	GLY	4.5
1	C	27	ASN	4.5
1	A	27	ASN	4.5
1	A	24	LEU	4.5
1	C	238	PRO	4.4
1	B	239	ASN	4.4
1	B	273	ASP	4.4
1	C	235	TYR	4.4
1	B	184	PRO	4.4
1	B	155	ASP	4.3
1	A	23	VAL	4.2
1	A	279	TYR	4.1
1	B	27	ASN	4.0
1	C	160	ALA	4.0
1	B	140	ASN	4.0
1	B	270	LYS	4.0
1	B	94	SER	4.0
1	A	140	ASN	3.9
1	A	241	LYS	3.8
1	D	200	GLY	3.8
1	A	28	TRP	3.8
1	C	209	LYS	3.7
1	A	156	SER	3.7
1	C	25	TYR	3.7
1	A	276	GLY	3.6
1	A	160	ALA	3.6
1	A	185	LYS	3.6
1	D	160	ALA	3.6
1	D	176	ASP	3.6
1	D	172	LYS	3.6
1	D	23	VAL	3.5
1	A	239	ASN	3.5
1	C	187	GLU	3.5
1	B	237	TYR	3.4
1	B	142	LYS	3.4
1	C	292	SER	3.4
1	A	270	LYS	3.4
1	C	45	ASN	3.3
1	A	179	LYS	3.3
1	B	24	LEU	3.3
1	B	158	CYS	3.3
1	C	240	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	25	TYR	3.2
1	A	157	ARG	3.2
1	B	144	VAL	3.2
1	D	163	ASP	3.2
1	C	124	PHE	3.1
1	A	57	ASN	3.1
1	C	275	ASP	3.1
1	C	236	ASN	3.1
1	C	141	GLY	3.0
1	A	158	CYS	3.0
1	D	154	GLU	2.9
1	D	237	TYR	2.9
1	C	186	ASP	2.9
1	A	277	TYR	2.8
1	C	73	TYR	2.8
1	A	159	GLY	2.8
1	C	47	ASP	2.8
1	D	201	THR	2.8
1	B	235	TYR	2.8
1	C	241	LYS	2.8
1	C	28	TRP	2.8
1	C	9	ASP	2.8
1	C	142	LYS	2.8
1	B	82	GLN	2.7
1	A	182	ASN	2.7
1	B	236	ASN	2.7
1	D	275	ASP	2.7
1	A	258	LYS	2.7
1	B	186	ASP	2.7
1	A	60	SER	2.7
1	B	143	ASN	2.6
1	B	9	ASP	2.6
1	B	275	ASP	2.6
1	C	138	GLU	2.6
1	A	81	SER	2.6
1	B	156	SER	2.6
1	D	81	SER	2.6
1	D	236	ASN	2.6
1	D	40	SER	2.6
1	A	236	ASN	2.6
1	A	93	SER	2.6
1	B	258	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	27	ASN	2.6
1	D	43	ILE	2.5
1	A	39	GLN	2.5
1	D	138	GLU	2.5
1	B	95	THR	2.5
1	B	259	GLN	2.5
1	D	159	GLY	2.5
1	D	45	ASN	2.5
1	C	139	LYS	2.5
1	C	162	HIS	2.5
1	B	138	GLU	2.5
1	A	272	TRP	2.4
1	B	187	GLU	2.4
1	D	168	ALA	2.4
1	A	32	LYS	2.4
1	D	46	ASN	2.4
1	D	26	PRO	2.4
1	B	292	SER	2.4
1	C	176	ASP	2.4
1	A	176	ASP	2.4
1	B	26	PRO	2.3
1	C	140	ASN	2.3
1	D	166	ILE	2.3
1	C	14	SER	2.3
1	C	108	TYR	2.3
1	C	23	VAL	2.3
1	B	72	PRO	2.2
1	D	39	GLN	2.2
1	A	154	GLU	2.2
1	D	182	ASN	2.2
1	C	163	ASP	2.2
1	B	109	PRO	2.2
1	D	122	CYS	2.2
1	D	9	ASP	2.1
1	B	141	GLY	2.1
1	C	210	ASN	2.1
1	A	269	PRO	2.1
1	B	71	TYR	2.1
1	C	92	TYR	2.1
1	A	142	LYS	2.1
1	B	162	HIS	2.1
1	C	127	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	179	LYS	2.1
1	C	202	PRO	2.1
1	B	21	SER	2.1
1	B	125	ASP	2.1
1	A	111	LYS	2.1
1	C	258	LYS	2.1
1	D	180	LYS	2.1
1	A	21	SER	2.0
1	C	159	GLY	2.0
1	C	272	TRP	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](#)

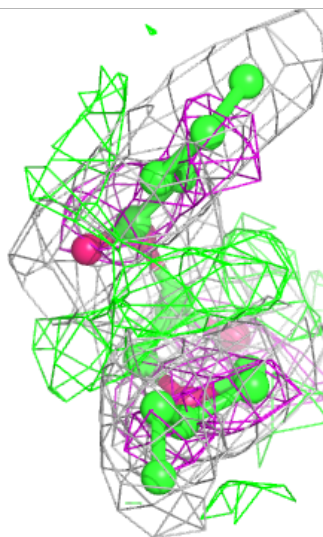
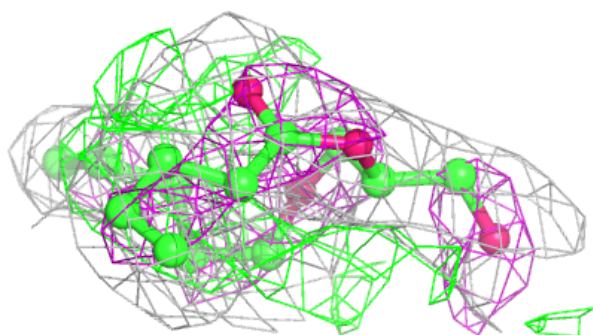
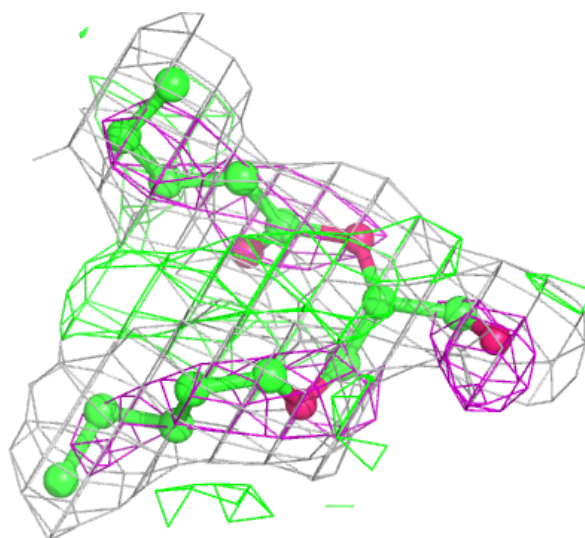
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	DGA	D	401	17/44	0.89	0.28	20,20,20,20	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 DGA D 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.