



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 02:30 am BST

PDB ID : 4EBN
Title : BlaC Amoxicillin Acyl-Intermediate Complex
Authors : Mire, J.A.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 2012-03-23
Resolution : 2.85 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

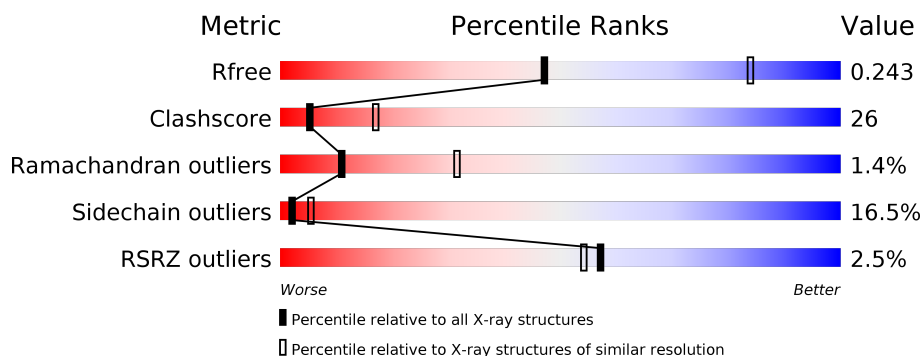
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	265	<div> <div>3%</div> <div> <div></div> <div>60%</div> <div>34%</div> <div>6%</div> </div> </div>
1	B	265	<div> <div>3%</div> <div> <div></div> <div>62%</div> <div>31%</div> <div>8%</div> </div> </div>
1	C	265	<div> <div>2%</div> <div> <div></div> <div>60%</div> <div>31%</div> <div>9%</div> </div> </div>
1	D	265	<div> <div>2%</div> <div> <div></div> <div>55%</div> <div>35%</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PO4	A	301	-	X	-	-
2	PO4	A	303	-	X	-	-
2	PO4	C	302	-	X	-	-
2	PO4	D	301	-	X	-	-
2	PO4	D	302	-	X	-	-
3	AXL	B	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7864 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-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			1974	1236	348	384	6			
1	B	265	Total	C	N	O	S	0	0	0
			1920	1205	338	371	6			
1	C	265	Total	C	N	O	S	0	0	0
			1943	1222	344	371	6			
1	D	265	Total	C	N	O	S	0	0	0
			1947	1221	344	376	6			

There are 4 discrepancies between the modelled and reference sequences:

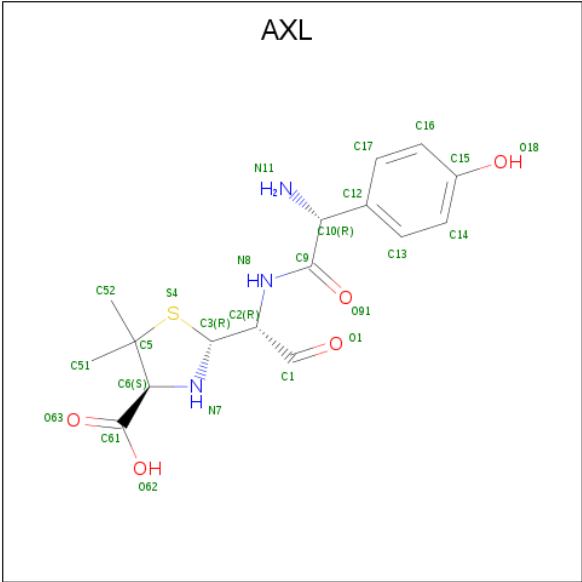
Chain	Residue	Modelled	Actual	Comment	Reference
A	166	ALA	GLU	ENGINEERED MUTATION	UNP P0C5C1
B	166	ALA	GLU	ENGINEERED MUTATION	UNP P0C5C1
C	166	ALA	GLU	ENGINEERED MUTATION	UNP P0C5C1
D	166	ALA	GLU	ENGINEERED MUTATION	UNP P0C5C1

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is 2-{1-[2-AMINO-2-(4-HYDROXY-PHENYL)-ACETYLAMINO]-2-OXO-ETHYL}-5,5-DIMETHYL-THIAZOLIDINE-4-CARBOXYLIC ACID (three-letter code: AXL) (formula: C₁₆H₂₁N₃O₅S).

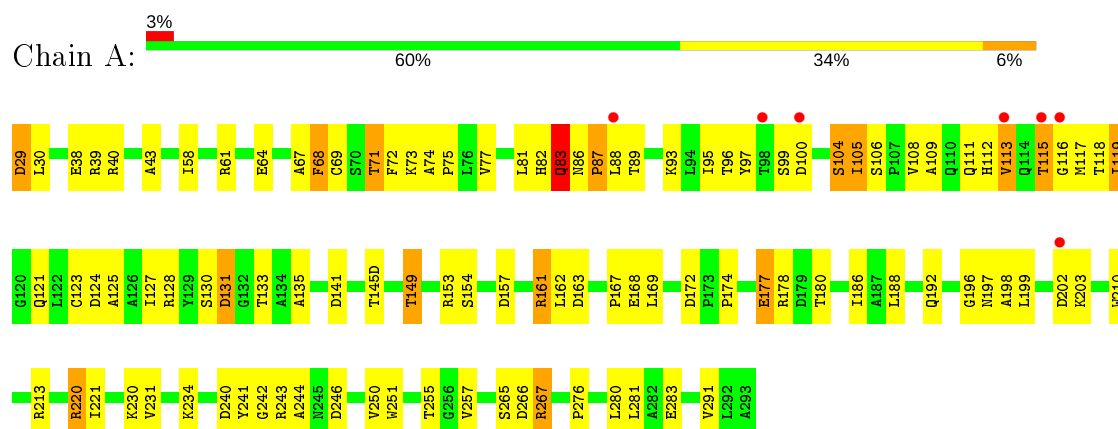


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	S	0	0
			25	16	3	5	1		
3	C	1	Total	C	N	O	S	0	0
			25	16	3	5	1		

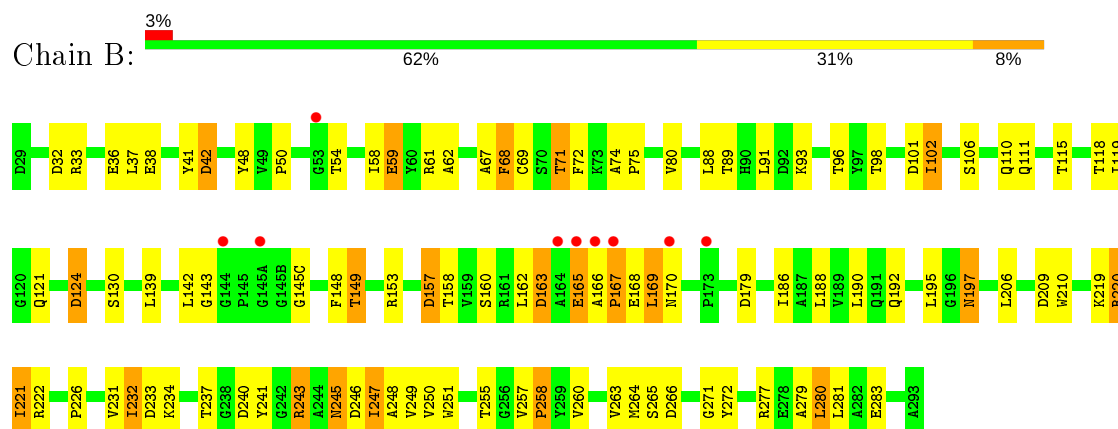
3 Residue-property plots

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

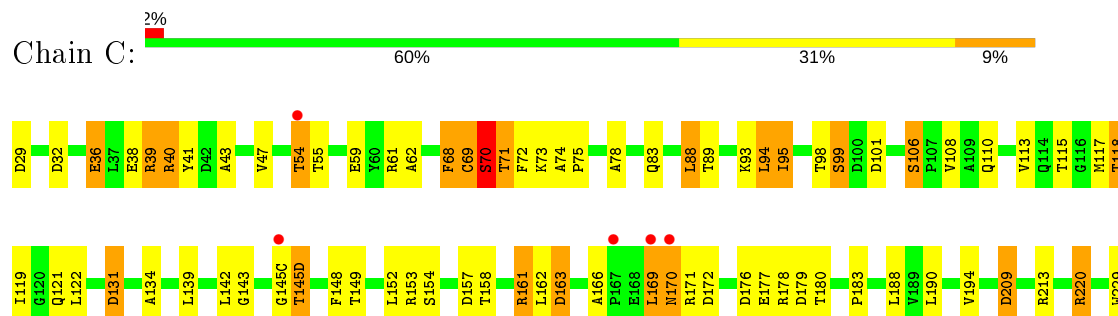
• Molecule 1: Beta-lactamase



• Molecule 1: Beta-lactamase

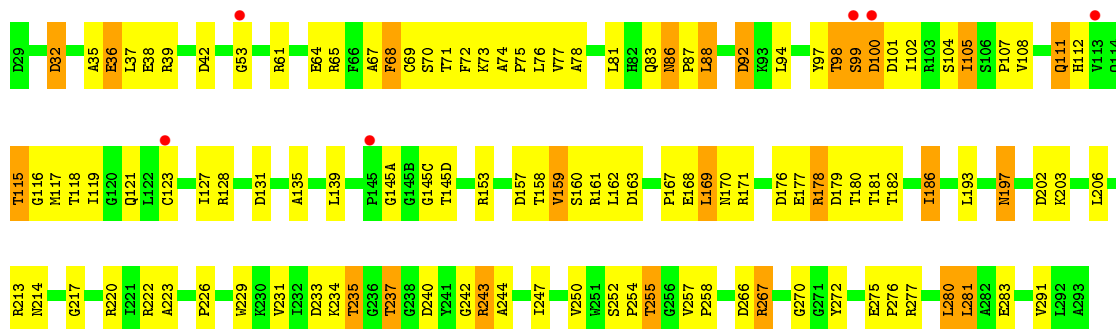


• Molecule 1: Beta-lactamase





• Molecule 1: Beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.57Å 96.67Å 109.81Å 90.00° 107.52° 90.00°	Depositor
Resolution (Å)	47.26 – 2.85 47.26 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.26-2.85) 99.1 (47.26-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.213 , 0.244 0.211 , 0.243	Depositor DCC
R_{free} test set	1842 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	60.5	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.3	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	0.93	EDS
Total number of atoms	7864	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.58% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, AXL

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	0.85	1/2013 (0.0%)	0.88	1/2750 (0.0%)
1	B	0.78	0/1958	0.85	1/2682 (0.0%)
1	C	0.71	0/1982	0.79	2/2711 (0.1%)
1	D	0.82	0/1986	0.87	1/2718 (0.0%)
All	All	0.79	1/7939 (0.0%)	0.85	5/10861 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	29	ASP	CB-CG	5.95	1.64	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	71	THR	N-CA-CB	6.64	122.92	110.30
1	B	69	CYS	C-N-CA	5.80	136.20	121.70
1	C	70	SER	C-N-CA	-5.76	107.31	121.70
1	D	243	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	A	161	ARG	NE-CZ-NH2	-5.29	117.65	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1974	0	1938	89	0
1	B	1920	0	1857	87	0
1	C	1943	0	1902	107	0
1	D	1947	0	1894	118	0
2	A	15	0	0	1	0
2	C	5	0	0	0	0
2	D	10	0	0	1	0
3	B	25	0	18	17	0
3	C	25	0	18	8	0
All	All	7864	0	7627	410	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 26.

All (410) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:THR:HG23	1:C:121:GLN:CD	1.69	1.11
1:C:118:THR:HG23	1:C:121:GLN:NE2	1.67	1.08
1:D:178:ARG:HH11	1:D:178:ARG:CG	1.69	1.06
1:C:163:ASP:OD2	1:C:178:ARG:HG2	1.56	1.05
1:D:112:HIS:HB3	1:D:115:THR:HG22	1.41	1.03
1:D:178:ARG:HH11	1:D:178:ARG:HG3	0.89	1.03
1:D:178:ARG:NH1	1:D:178:ARG:HG3	1.59	1.01
1:C:39:ARG:HG2	1:C:39:ARG:HH21	1.24	1.00
1:C:39:ARG:CG	1:C:39:ARG:HH21	1.73	1.00
1:C:118:THR:CG2	1:C:121:GLN:CD	2.30	0.99
1:C:170:ASN:C	1:C:170:ASN:HD22	1.65	0.99
1:C:237:THR:OG1	3:C:301:AXL:HC6	1.64	0.97
1:C:240:ASP:O	1:C:243:ARG:HG3	1.64	0.95
1:C:118:THR:CG2	1:C:121:GLN:NE2	2.30	0.95
1:D:73:LYS:HE2	1:D:135:ALA:HB2	1.49	0.95
1:B:170:ASN:O	1:B:240:ASP:HB2	1.67	0.94
1:C:235:THR:HG22	1:C:246:ASP:OD1	1.70	0.92
1:A:83:GLN:HG3	1:A:83:GLN:O	1.70	0.90
1:B:166:ALA:N	1:B:167:PRO:HD2	1.86	0.89
1:A:83:GLN:HG3	1:A:86:ASN:ND2	1.86	0.89
1:D:112:HIS:CB	1:D:115:THR:HG22	2.04	0.87
1:A:198:ALA:O	1:A:199:LEU:HD23	1.76	0.85
1:B:221:ILE:HG13	1:B:246:ASP:OD2	1.77	0.85
1:A:82:HIS:O	1:A:82:HIS:CG	2.30	0.84
1:C:74:ALA:HB3	1:C:75:PRO:HD3	1.59	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:170:ASN:C	1:C:170:ASN:ND2	2.30	0.83
1:D:112:HIS:HB3	1:D:115:THR:CG2	2.08	0.82
3:C:301:AXL:HC3	3:C:301:AXL:O63	1.77	0.82
1:A:242:GLY:HA3	1:A:267:ARG:HG3	1.62	0.82
1:C:39:ARG:HG2	1:C:39:ARG:NH2	1.84	0.82
1:A:106:SER:HB3	1:A:109:ALA:HB3	1.61	0.81
1:C:169:LEU:N	1:C:169:LEU:HD23	1.96	0.81
1:A:104:SER:HB3	1:A:167:PRO:HG3	1.62	0.81
3:C:301:AXL:H512	3:C:301:AXL:C14	2.10	0.81
1:A:220:ARG:HD3	1:A:281:LEU:HD12	1.61	0.81
1:D:169:LEU:HD12	1:D:169:LEU:C	2.01	0.80
1:C:170:ASN:ND2	1:C:171:ARG:N	2.30	0.79
1:A:149:THR:HG23	1:A:161:ARG:HA	1.65	0.78
1:A:83:GLN:CG	1:A:86:ASN:ND2	2.47	0.78
1:C:39:ARG:CB	1:C:39:ARG:HH21	1.96	0.78
1:D:74:ALA:HB3	1:D:75:PRO:HD3	1.66	0.77
1:D:169:LEU:HD12	1:D:170:ASN:N	2.00	0.76
1:A:39:ARG:HG3	1:A:39:ARG:HH21	1.50	0.76
1:B:166:ALA:N	1:B:167:PRO:CD	2.49	0.75
1:D:139:LEU:O	1:D:145(C):GLY:HA3	1.86	0.75
1:C:70:SER:O	1:C:73:LYS:HG3	1.86	0.75
1:B:162:LEU:O	1:B:163:ASP:CB	2.35	0.75
1:A:149:THR:CG2	1:A:161:ARG:HA	2.18	0.73
1:A:61:ARG:HG2	1:A:64:GLU:CD	2.08	0.73
1:C:240:ASP:O	1:C:243:ARG:CG	2.35	0.73
1:D:115:THR:HG23	1:D:116:GLY:O	1.89	0.73
1:D:169:LEU:CD1	1:D:169:LEU:C	2.56	0.73
1:B:170:ASN:HD21	3:B:301:AXL:H10	1.52	0.72
1:D:145(A):GLY:O	1:D:145(D):THR:HG23	1.89	0.72
1:A:119:ILE:HD11	1:A:141:ASP:OD1	1.88	0.72
1:B:220:ARG:HD3	1:B:281:LEU:HD12	1.72	0.72
1:D:220:ARG:HG2	1:D:281:LEU:HD12	1.72	0.72
1:B:221:ILE:HG13	1:B:246:ASP:CG	2.09	0.72
1:C:69:CYS:O	1:C:71:THR:N	2.23	0.72
3:C:301:AXL:O63	3:C:301:AXL:C3	2.37	0.71
1:D:104:SER:HB3	1:D:167:PRO:HB3	1.70	0.71
1:A:97:TYR:CE1	1:A:113:VAL:HG22	2.24	0.71
1:C:169:LEU:H	1:C:169:LEU:CD2	2.04	0.70
1:A:112:HIS:HA	1:A:115:THR:HG23	1.73	0.70
1:A:118:THR:H	1:A:121:GLN:NE2	1.91	0.69
1:C:240:ASP:OD2	1:C:272:TYR:CZ	2.45	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLN:HG3	1:A:86:ASN:HD21	1.55	0.69
1:B:221:ILE:CG1	1:B:246:ASP:OD2	2.39	0.69
1:C:131:ASP:C	1:C:131:ASP:OD1	2.30	0.69
1:D:197:ASN:H	1:D:197:ASN:ND2	1.91	0.69
1:A:83:GLN:O	1:A:86:ASN:CG	2.30	0.69
1:C:39:ARG:HB3	1:C:39:ARG:HH21	1.58	0.69
1:B:157:ASP:OD1	1:B:157:ASP:C	2.31	0.68
1:C:72:PHE:O	1:C:75:PRO:HD2	1.93	0.68
1:A:97:TYR:CE1	1:A:113:VAL:CG2	2.77	0.68
1:A:82:HIS:O	1:A:83:GLN:HB3	1.93	0.67
1:D:112:HIS:CA	1:D:115:THR:HG22	2.25	0.67
1:B:170:ASN:OD1	3:B:301:AXL:N11	2.27	0.66
1:C:169:LEU:N	1:C:169:LEU:CD2	2.58	0.66
1:C:75:PRO:O	1:C:78:ALA:HB3	1.96	0.66
3:B:301:AXL:H512	3:B:301:AXL:C14	2.24	0.66
1:C:170:ASN:O	1:C:243:ARG:NH1	2.28	0.66
1:D:112:HIS:O	1:D:116:GLY:N	2.23	0.66
1:C:38:GLU:CD	1:C:61:ARG:HH21	1.98	0.66
1:B:166:ALA:O	1:B:168:GLU:N	2.29	0.65
1:C:237:THR:OG1	3:C:301:AXL:C6	2.41	0.65
1:A:131:ASP:C	1:A:131:ASP:OD1	2.35	0.65
1:B:168:GLU:O	1:B:169:LEU:CB	2.44	0.65
3:B:301:AXL:O63	3:B:301:AXL:H523	1.97	0.65
1:C:194:VAL:HG11	1:C:249:VAL:HG21	1.78	0.65
1:A:82:HIS:O	1:A:82:HIS:ND1	2.30	0.65
1:D:153:ARG:NH1	1:D:157:ASP:O	2.30	0.65
1:D:65:ARG:HA	1:D:182:THR:HG22	1.79	0.65
1:D:36:GLU:O	1:D:38:GLU:N	2.29	0.65
1:C:62:ALA:O	1:C:183:PRO:HD2	1.96	0.65
1:A:112:HIS:O	1:A:116:GLY:N	2.24	0.64
1:D:98:THR:O	1:D:100:ASP:N	2.30	0.64
1:B:148:PHE:CD1	1:B:162:LEU:HD21	2.33	0.64
1:C:163:ASP:OD1	1:C:163:ASP:N	2.30	0.64
1:A:83:GLN:O	1:A:86:ASN:ND2	2.30	0.64
1:B:197:ASN:H	1:B:197:ASN:HD22	1.46	0.64
1:D:70:SER:O	1:D:73:LYS:HG3	1.97	0.63
1:C:231:VAL:HG22	1:C:250:VAL:HG12	1.79	0.63
1:B:149:THR:O	1:B:153:ARG:HG2	1.99	0.63
1:C:170:ASN:HD22	1:C:171:ARG:N	1.91	0.63
1:B:222:ARG:NH2	1:B:226:PRO:O	2.33	0.62
1:D:83:GLN:C	1:D:86:ASN:ND2	2.53	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:72:PHE:HE1	1:D:76:LEU:HD22	1.64	0.62
1:B:166:ALA:C	1:B:168:GLU:H	2.03	0.62
1:C:169:LEU:H	1:C:169:LEU:HD23	1.60	0.62
1:A:39:ARG:HG3	1:A:39:ARG:NH2	2.15	0.61
1:C:38:GLU:HG2	1:C:43:ALA:O	1.99	0.61
1:D:267:ARG:NH2	1:D:275:GLU:O	2.32	0.61
1:A:131:ASP:OD1	1:A:133:THR:N	2.33	0.61
1:C:94:LEU:HG	1:C:95:ILE:N	2.13	0.61
1:D:108:VAL:HB	1:D:131:ASP:OD2	2.00	0.61
1:B:240:ASP:O	1:B:241:TYR:HB2	1.99	0.61
1:D:280:LEU:C	1:D:280:LEU:CD2	2.68	0.61
1:A:97:TYR:HE1	1:A:113:VAL:HG23	1.66	0.61
1:B:170:ASN:CG	3:B:301:AXL:N11	2.54	0.61
3:C:301:AXL:C51	3:C:301:AXL:C14	2.79	0.61
1:D:77:VAL:O	1:D:81:LEU:HD12	2.01	0.60
1:D:162:LEU:HD12	1:D:179:ASP:OD1	2.00	0.60
1:A:97:TYR:CD1	1:A:113:VAL:HG22	2.36	0.60
1:B:197:ASN:H	1:B:197:ASN:ND2	2.00	0.60
1:B:59:GLU:HB3	1:B:62:ALA:HB2	1.83	0.60
1:A:83:GLN:CG	1:A:86:ASN:HD21	2.12	0.60
1:D:220:ARG:HH11	1:D:244:ALA:HB1	1.66	0.60
1:A:61:ARG:HD3	1:A:64:GLU:OE2	2.02	0.59
1:D:76:LEU:O	1:D:76:LEU:HD12	2.02	0.59
1:A:240:ASP:O	1:A:243:ARG:HG3	2.03	0.58
1:A:128:ARG:HD2	1:A:213:ARG:O	2.04	0.58
1:B:170:ASN:ND2	3:B:301:AXL:H10	2.17	0.58
1:B:249:VAL:HG22	1:B:260:VAL:HG22	1.86	0.58
1:C:152:LEU:O	1:C:157:ASP:HB3	2.03	0.58
1:C:118:THR:CG2	1:C:121:GLN:OE1	2.51	0.58
1:B:170:ASN:HD21	3:B:301:AXL:C10	2.17	0.58
1:D:153:ARG:NH1	1:D:158:THR:HA	2.19	0.58
1:D:220:ARG:HD3	1:D:235:THR:HG22	1.86	0.57
1:A:38:GLU:CD	1:A:61:ARG:HH21	2.07	0.57
1:D:105:ILE:HG22	1:D:107:PRO:HD3	1.85	0.57
1:D:214:ASN:HD21	1:D:217:GLY:N	2.03	0.57
1:C:70:SER:HB2	1:C:236:GLY:HA2	1.87	0.57
1:C:118:THR:H	1:C:121:GLN:NE2	2.03	0.57
1:D:157:ASP:OD1	1:D:157:ASP:C	2.42	0.57
1:A:221:ILE:HG13	1:A:246:ASP:OD2	2.05	0.57
1:C:32:ASP:O	1:C:36:GLU:HG2	2.04	0.57
1:D:128:ARG:HD2	1:D:213:ARG:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:ASP:O	1:B:243:ARG:HG3	2.05	0.56
1:C:54:THR:HG23	1:C:55:THR:HG23	1.87	0.56
1:D:61:ARG:HG2	1:D:64:GLU:OE1	2.05	0.56
1:D:280:LEU:C	1:D:280:LEU:HD22	2.25	0.56
1:C:148:PHE:CD1	1:C:162:LEU:HD22	2.40	0.56
1:A:61:ARG:HG2	1:A:64:GLU:OE1	2.06	0.56
1:B:231:VAL:HG22	1:B:250:VAL:HG12	1.87	0.56
1:B:197:ASN:N	1:B:197:ASN:HD22	2.04	0.56
1:B:33:ARG:O	1:B:37:LEU:HD12	2.06	0.56
1:A:240:ASP:O	1:A:241:TYR:HB2	2.05	0.55
1:A:221:ILE:HG13	1:A:246:ASP:HB3	1.88	0.55
1:D:176:ASP:OD2	1:D:178:ARG:HG3	2.06	0.55
1:D:186:ILE:O	1:D:186:ILE:HD12	2.05	0.55
1:A:202:ASP:OD1	1:A:203:LYS:N	2.39	0.55
1:D:240:ASP:O	1:D:243:ARG:HG3	2.07	0.55
1:A:69:CYS:O	1:A:72:PHE:HD2	1.89	0.55
1:C:241:TYR:HA	1:C:271:GLY:O	2.06	0.55
1:A:77:VAL:HG13	1:A:123:CYS:SG	2.46	0.55
1:D:69:CYS:O	1:D:72:PHE:HD2	1.90	0.54
1:B:32:ASP:O	1:B:36:GLU:HG3	2.07	0.54
1:B:41:TYR:O	1:B:42:ASP:HB2	2.06	0.54
1:D:163:ASP:OD1	1:D:178:ARG:HA	2.07	0.54
1:C:68:PHE:C	1:C:69:CYS:O	2.45	0.54
1:A:77:VAL:O	1:A:81:LEU:HD12	2.07	0.54
1:D:112:HIS:HD2	1:D:115:THR:HG21	1.70	0.54
1:A:97:TYR:HE1	1:A:113:VAL:CG2	2.17	0.54
1:B:118:THR:H	1:B:121:GLN:NE2	2.06	0.54
1:B:247:ILE:HG12	1:B:247:ILE:O	2.07	0.54
1:D:112:HIS:CD2	1:D:115:THR:HG21	2.43	0.54
1:A:192:GLN:HA	1:A:196:GLY:HA3	1.90	0.54
1:C:118:THR:HG22	1:C:121:GLN:NE2	2.19	0.54
1:C:169:LEU:HB2	1:C:179:ASP:CG	2.29	0.54
1:B:102:ILE:HD12	1:B:110:GLN:HA	1.90	0.54
1:B:232:ILE:O	1:B:248:ALA:HB1	2.07	0.54
1:C:143:GLY:O	1:C:145(C):GLY:HA2	2.08	0.54
1:C:118:THR:HG23	1:C:121:GLN:CG	2.38	0.53
1:A:117:MET:HE3	1:A:121:GLN:HB3	1.90	0.53
1:C:71:THR:HG23	1:C:236:GLY:HA3	1.89	0.53
1:B:170:ASN:O	1:B:240:ASP:CB	2.49	0.53
1:D:197:ASN:H	1:D:197:ASN:HD22	1.56	0.53
1:D:280:LEU:CD2	1:D:280:LEU:O	2.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:237:THR:OG1	2:D:301:PO4:O4	2.24	0.53
1:B:102:ILE:CD1	1:B:110:GLN:HA	2.38	0.53
1:B:241:TYR:HA	1:B:271:GLY:O	2.09	0.53
1:C:266:ASP:C	1:C:266:ASP:OD2	2.46	0.53
1:A:105:ILE:O	1:A:105:ILE:HG22	2.07	0.53
1:A:86:ASN:O	1:A:87:PRO:C	2.46	0.52
1:D:36:GLU:HA	1:D:39:ARG:CZ	2.39	0.52
1:D:42:ASP:O	1:D:42:ASP:CG	2.48	0.52
1:C:190:LEU:HD22	1:C:247:ILE:HD12	1.90	0.52
1:B:237:THR:HB	3:B:301:AXL:HN7	1.75	0.52
1:B:245:ASN:OD1	1:B:245:ASN:N	2.39	0.52
1:C:106:SER:HB3	1:C:131:ASP:OD2	2.10	0.52
1:A:58:ILE:HD11	1:A:291:VAL:HG11	1.92	0.51
1:A:117:MET:CE	1:A:121:GLN:HB3	2.40	0.51
1:D:77:VAL:HG13	1:D:123:CYS:SG	2.50	0.51
1:B:190:LEU:HD22	1:B:247:ILE:HG13	1.93	0.51
1:A:231:VAL:HG22	1:A:250:VAL:HG12	1.93	0.51
1:A:73:LYS:HE3	1:A:135:ALA:HB2	1.93	0.51
1:A:39:ARG:HH21	1:A:39:ARG:CG	2.22	0.51
1:D:145(A):GLY:O	1:D:145(D):THR:CG2	2.58	0.51
1:D:98:THR:C	1:D:100:ASP:H	2.14	0.51
1:B:188:LEU:O	1:B:192:GLN:HG3	2.11	0.51
1:D:72:PHE:C	1:D:72:PHE:CD1	2.84	0.51
1:D:72:PHE:O	1:D:72:PHE:CD1	2.64	0.50
1:B:165:GLU:C	1:B:167:PRO:HD2	2.31	0.50
1:B:58:ILE:C	1:B:59:GLU:HG3	2.26	0.50
1:B:170:ASN:OD1	3:B:301:AXL:C10	2.59	0.50
1:B:38:GLU:OE2	1:B:61:ARG:NH2	2.44	0.50
1:D:42:ASP:O	1:D:42:ASP:OD2	2.30	0.50
1:A:112:HIS:CA	1:A:115:THR:HG23	2.41	0.50
1:C:68:PHE:O	1:C:69:CYS:O	2.30	0.50
1:D:159:VAL:HG12	1:D:160:SER:N	2.26	0.50
1:C:279:ALA:O	1:C:283:GLU:HG2	2.12	0.50
1:D:35:ALA:O	1:D:36:GLU:O	2.30	0.50
1:A:83:GLN:HG2	1:A:86:ASN:ND2	2.25	0.49
1:B:170:ASN:ND2	3:B:301:AXL:C10	2.75	0.49
1:D:74:ALA:HB3	1:D:75:PRO:CD	2.40	0.49
1:C:149:THR:O	1:C:153:ARG:HG2	2.12	0.49
1:A:68:PHE:O	1:A:71:THR:OG1	2.30	0.49
1:C:32:ASP:O	1:C:36:GLU:CG	2.61	0.49
1:C:229:TRP:CE2	1:C:254:PRO:HD3	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:THR:HB	3:C:301:AXL:HN7	1.78	0.49
1:D:68:PHE:O	1:D:71:THR:OG1	2.30	0.49
1:A:220:ARG:HD3	1:A:281:LEU:CD1	2.39	0.49
1:B:277:ARG:HG3	1:B:277:ARG:HH11	1.77	0.49
1:B:279:ALA:O	1:B:283:GLU:HG2	2.12	0.49
1:D:222:ARG:NH2	1:D:226:PRO:O	2.46	0.49
1:C:249:VAL:HG22	1:C:260:VAL:HG22	1.94	0.48
1:D:242:GLY:HA3	1:D:267:ARG:HG3	1.94	0.48
1:B:59:GLU:HB3	1:B:62:ALA:CB	2.43	0.48
1:C:99:SER:HB2	1:D:105:ILE:HG13	1.95	0.48
1:A:255:THR:HG22	1:A:257:VAL:HG23	1.96	0.48
1:B:280:LEU:HD22	1:B:280:LEU:O	2.13	0.48
3:B:301:AXL:O63	3:B:301:AXL:HC3	2.12	0.48
1:D:65:ARG:HA	1:D:181:THR:O	2.13	0.48
1:A:255:THR:O	1:A:255:THR:CG2	2.61	0.48
1:B:266:ASP:OD2	1:B:266:ASP:C	2.51	0.48
1:D:118:THR:H	1:D:121:GLN:NE2	2.10	0.48
1:C:108:VAL:HG12	1:C:108:VAL:O	2.13	0.48
1:C:72:PHE:HD2	1:C:72:PHE:H	1.59	0.48
1:A:86:ASN:O	1:A:87:PRO:O	2.30	0.48
1:D:220:ARG:HD3	1:D:235:THR:CG2	2.43	0.48
1:D:67:ALA:HB3	1:D:243:ARG:HD3	1.94	0.48
1:B:197:ASN:N	1:B:197:ASN:ND2	2.61	0.48
1:C:176:ASP:C	1:C:176:ASP:OD2	2.52	0.48
3:B:301:AXL:H512	3:B:301:AXL:C13	2.44	0.47
1:D:36:GLU:O	1:D:39:ARG:N	2.37	0.47
1:B:157:ASP:OD1	1:B:157:ASP:O	2.32	0.47
1:C:68:PHE:HB2	1:C:179:ASP:O	2.14	0.47
1:D:88:LEU:HD12	1:D:88:LEU:HA	1.69	0.47
1:D:98:THR:C	1:D:100:ASP:N	2.66	0.47
1:D:247:ILE:O	1:D:247:ILE:CG1	2.61	0.47
1:D:38:GLU:O	1:D:42:ASP:N	2.47	0.47
1:B:226:PRO:HB3	1:D:223:ALA:O	2.14	0.47
1:D:176:ASP:OD2	1:D:178:ARG:CG	2.62	0.47
1:A:43:ALA:HB1	1:A:266:ASP:O	2.14	0.47
1:B:98:THR:O	1:B:101:ASP:HB2	2.15	0.47
1:B:38:GLU:CD	1:B:61:ARG:HH21	2.17	0.47
1:A:188:LEU:O	1:A:192:GLN:HG3	2.15	0.47
1:A:87:PRO:C	1:A:89:THR:H	2.17	0.47
1:B:143:GLY:O	1:B:145(C):GLY:HA2	2.15	0.47
1:A:87:PRO:C	1:A:89:THR:N	2.69	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:74:ALA:CB	1:C:75:PRO:HD3	2.38	0.47
1:D:272:TYR:CD2	1:D:272:TYR:O	2.68	0.47
1:A:30:LEU:HD22	1:A:291:VAL:HG21	1.97	0.47
1:B:263:VAL:CG1	1:B:281:LEU:HD22	2.45	0.47
1:A:83:GLN:C	1:A:86:ASN:ND2	2.69	0.47
1:D:206:LEU:HD23	1:D:206:LEU:HA	1.61	0.47
1:B:241:TYR:CD2	1:B:271:GLY:C	2.89	0.46
1:A:74:ALA:HB3	1:A:75:PRO:HD3	1.97	0.46
1:C:166:ALA:HA	1:C:169:LEU:CD2	2.45	0.46
1:C:69:CYS:C	1:C:71:THR:H	2.16	0.46
1:D:214:ASN:HD21	1:D:217:GLY:H	1.61	0.46
1:A:149:THR:HG21	1:A:161:ARG:HA	1.95	0.46
1:C:88:LEU:HD12	1:C:88:LEU:HA	1.69	0.46
1:B:67:ALA:O	1:B:245:ASN:ND2	2.49	0.46
1:C:220:ARG:HE	1:C:278:GLU:CG	2.29	0.46
1:A:74:ALA:HB3	1:A:75:PRO:CD	2.45	0.46
1:A:124:ASP:OD1	1:A:124:ASP:C	2.54	0.46
1:C:118:THR:HG21	1:C:121:GLN:OE1	2.16	0.46
1:C:145(D):THR:HG22	1:C:162:LEU:HD23	1.98	0.46
1:C:41:TYR:CD2	1:C:277:ARG:HG2	2.51	0.46
3:C:301:AXL:C51	3:C:301:AXL:H14	2.46	0.46
1:D:252:SER:O	1:D:255:THR:N	2.49	0.46
1:D:168:GLU:OE1	1:D:168:GLU:N	2.46	0.46
1:D:280:LEU:O	1:D:280:LEU:HD23	2.16	0.46
1:D:81:LEU:O	1:D:203:LYS:NZ	2.43	0.46
1:A:112:HIS:HA	1:A:115:THR:CG2	2.45	0.45
1:B:80:VAL:HG23	1:B:142:LEU:HD21	1.97	0.45
1:B:237:THR:HB	3:B:301:AXL:N7	2.32	0.45
1:A:141:ASP:OD2	1:A:141:ASP:O	2.35	0.45
1:B:166:ALA:C	1:B:168:GLU:N	2.69	0.45
1:B:72:PHE:O	1:B:75:PRO:HG2	2.16	0.45
1:C:247:ILE:HG13	1:C:247:ILE:O	2.16	0.45
1:D:197:ASN:HD22	1:D:197:ASN:N	2.14	0.45
1:A:178:ARG:HE	1:A:178:ARG:HB2	1.50	0.45
1:A:61:ARG:CD	1:A:64:GLU:OE2	2.64	0.45
1:C:71:THR:HG22	1:C:235:THR:C	2.36	0.45
1:C:71:THR:HG22	1:C:236:GLY:N	2.31	0.45
1:D:220:ARG:HH21	1:D:235:THR:HG21	1.81	0.45
1:D:231:VAL:HG22	1:D:250:VAL:HG12	1.98	0.45
1:D:280:LEU:O	1:D:280:LEU:HD22	2.17	0.45
1:D:65:ARG:HD2	1:D:177:GLU:HG2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:118:THR:O	1:D:119:ILE:C	2.54	0.45
1:A:67:ALA:HB3	1:A:243:ARG:HD3	1.98	0.44
1:B:50:PRO:HD2	1:B:260:VAL:O	2.18	0.44
1:C:139:LEU:O	1:C:142:LEU:HB2	2.16	0.44
1:C:171:ARG:O	1:C:172:ASP:C	2.55	0.44
1:D:255:THR:O	1:D:255:THR:HG23	2.15	0.44
1:C:110:GLN:O	1:C:113:VAL:HG23	2.16	0.44
1:D:86:ASN:HB3	1:D:87:PRO:HD2	1.99	0.44
1:D:92:ASP:N	1:D:92:ASP:OD1	2.49	0.44
1:B:206:LEU:HA	1:B:206:LEU:HD23	1.87	0.44
1:A:97:TYR:N	1:A:97:TYR:CD2	2.85	0.44
1:D:255:THR:HG22	1:D:257:VAL:HG23	1.99	0.44
1:C:292:LEU:O	1:C:293:ALA:C	2.56	0.44
1:B:48:TYR:HA	1:B:58:ILE:O	2.18	0.44
1:C:38:GLU:OE2	1:C:61:ARG:NH2	2.50	0.44
1:D:266:ASP:C	1:D:266:ASP:OD2	2.56	0.44
1:C:98:THR:O	1:C:101:ASP:HB2	2.17	0.44
1:D:111:GLN:H	1:D:111:GLN:HG2	1.36	0.44
1:D:97:TYR:OH	1:D:112:HIS:HB2	2.17	0.43
1:D:242:GLY:HA2	1:D:276:PRO:HG3	2.00	0.43
1:C:75:PRO:O	1:C:78:ALA:N	2.51	0.43
1:D:161:ARG:O	1:D:179:ASP:HA	2.18	0.43
1:D:229:TRP:CE2	1:D:254:PRO:HD3	2.54	0.43
3:B:301:AXL:C14	3:B:301:AXL:C51	2.93	0.43
1:B:68:PHE:O	1:B:71:THR:OG1	2.36	0.43
1:C:176:ASP:OD2	1:C:177:GLU:N	2.51	0.43
1:B:195:LEU:HA	1:B:195:LEU:HD23	1.76	0.43
3:B:301:AXL:O63	3:B:301:AXL:C3	2.65	0.43
1:C:166:ALA:HA	1:C:169:LEU:HD22	2.00	0.43
1:C:153:ARG:HD3	1:C:157:ASP:O	2.19	0.43
1:D:117:MET:HB3	1:D:121:GLN:HE21	1.83	0.43
1:A:266:ASP:C	1:A:267:ARG:HG2	2.39	0.43
1:C:39:ARG:CB	1:C:39:ARG:NH2	2.75	0.43
1:B:220:ARG:HB2	1:B:246:ASP:OD2	2.19	0.42
2:A:301:PO4:P	1:B:111:GLN:HE22	2.42	0.42
1:C:209:ASP:O	1:C:213:ARG:HG3	2.19	0.42
1:C:220:ARG:HD3	1:C:281:LEU:HD12	1.99	0.42
1:A:127:ILE:HG21	1:A:210:TRP:HB3	2.02	0.42
1:A:177:GLU:O	1:A:180:THR:OG1	2.30	0.42
1:C:131:ASP:OD1	1:C:134:ALA:N	2.45	0.42
1:A:168:GLU:CD	1:A:168:GLU:H	2.21	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:ALA:HB2	1:A:276:PRO:HB3	2.02	0.42
1:A:88:LEU:HA	1:A:88:LEU:HD12	1.78	0.42
1:C:108:VAL:CG1	1:C:108:VAL:O	2.67	0.42
1:C:267:ARG:HD3	1:C:275:GLU:O	2.19	0.42
1:A:153:ARG:HD3	1:A:157:ASP:O	2.20	0.42
1:A:230:LYS:HB3	1:A:251:TRP:HB2	2.01	0.42
1:B:245:ASN:ND2	1:B:264:MET:CE	2.83	0.42
1:C:152:LEU:HD23	1:C:152:LEU:HA	1.70	0.42
1:C:161:ARG:HE	1:C:161:ARG:HB3	1.27	0.42
1:D:193:LEU:HD23	1:D:193:LEU:HA	1.81	0.42
1:D:97:TYR:CE2	1:D:116:GLY:HA2	2.55	0.42
1:B:93:LYS:O	1:B:118:THR:HA	2.19	0.42
1:B:170:ASN:ND2	3:B:301:AXL:N11	2.68	0.42
1:B:41:TYR:O	1:B:42:ASP:CB	2.67	0.42
1:D:160:SER:HA	1:D:180:THR:O	2.19	0.42
1:C:39:ARG:CG	1:C:39:ARG:NH2	2.43	0.41
1:D:163:ASP:OD1	1:D:178:ARG:CA	2.68	0.41
1:A:174:PRO:HD3	1:A:241:TYR:CE2	2.55	0.41
1:C:39:ARG:HB3	1:C:39:ARG:NH2	2.30	0.41
1:A:145(D):THR:HB	1:A:162:LEU:O	2.21	0.41
1:B:157:ASP:OD2	1:B:160:SER:OG	2.30	0.41
1:D:127:ILE:HG21	1:D:127:ILE:HD13	1.70	0.41
1:B:74:ALA:HB3	1:B:75:PRO:HD2	2.01	0.41
1:D:76:LEU:HD12	1:D:76:LEU:C	2.41	0.41
1:B:251:TRP:HA	1:B:257:VAL:O	2.20	0.41
1:A:69:CYS:O	1:A:72:PHE:CD2	2.72	0.41
1:B:240:ASP:O	1:B:241:TYR:CB	2.68	0.41
1:C:161:ARG:O	1:C:179:ASP:HA	2.20	0.41
1:D:78:ALA:HA	1:D:81:LEU:HD12	2.02	0.41
1:A:108:VAL:HG21	1:A:125:ALA:HB1	2.03	0.41
1:B:91:LEU:HA	1:B:119:ILE:HB	2.02	0.41
1:B:233:ASP:CB	1:B:248:ALA:HB2	2.51	0.41
1:B:241:TYR:CD2	1:B:272:TYR:HA	2.56	0.41
1:B:170:ASN:OD1	3:B:301:AXL:H10	2.20	0.41
1:C:69:CYS:C	1:C:71:THR:N	2.73	0.41
1:D:72:PHE:HE2	1:D:169:LEU:HD21	1.85	0.41
1:C:47:VAL:O	1:C:59:GLU:HA	2.20	0.41
1:D:32:ASP:N	1:D:32:ASP:OD1	2.54	0.41
1:D:83:GLN:C	1:D:86:ASN:HD22	2.22	0.41
1:D:252:SER:C	1:D:255:THR:N	2.72	0.41
1:D:77:VAL:HG12	1:D:81:LEU:HD11	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ARG:HB2	1:A:40:ARG:HE	1.64	0.41
1:C:122:LEU:HD22	1:C:134:ALA:HA	2.01	0.41
1:B:251:TRP:NE1	1:B:258:PRO:HB3	2.35	0.40
1:C:40:ARG:HG3	1:C:40:ARG:O	2.21	0.40
1:C:72:PHE:N	1:C:72:PHE:CD2	2.89	0.40
1:D:118:THR:H	1:D:121:GLN:HE21	1.69	0.40
1:D:86:ASN:ND2	1:D:86:ASN:N	2.70	0.40
1:C:188:LEU:HD23	1:C:188:LEU:HA	1.88	0.40
1:D:98:THR:N	1:D:101:ASP:OD2	2.30	0.40
1:D:252:SER:O	1:D:254:PRO:C	2.60	0.40
1:A:104:SER:HB2	1:A:105:ILE:H	1.64	0.40
1:B:186:ILE:O	1:B:186:ILE:HG13	2.21	0.40
1:C:162:LEU:C	1:C:163:ASP:OD1	2.60	0.40
1:D:277:ARG:HA	1:D:277:ARG:HD2	1.84	0.40
1:B:124:ASP:HB2	1:B:210:TRP:NE1	2.37	0.40
1:C:163:ASP:OD2	1:C:178:ARG:CG	2.47	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	263/265 (99%)	249 (95%)	12 (5%)	2 (1%)	19	46
1	B	263/265 (99%)	246 (94%)	12 (5%)	5 (2%)	8	24
1	C	263/265 (99%)	252 (96%)	9 (3%)	2 (1%)	19	46
1	D	263/265 (99%)	242 (92%)	15 (6%)	6 (2%)	6	20
All	All	1052/1060 (99%)	989 (94%)	48 (5%)	15 (1%)	11	31

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	83	GLN
1	B	165	GLU
1	C	70	SER
1	D	36	GLU
1	D	37	LEU
1	A	87	PRO
1	B	163	ASP
1	D	99	SER
1	D	105	ILE
1	C	69	CYS
1	D	270	GLY
1	B	167	PRO
1	B	169	LEU
1	B	258	PRO
1	D	53	GLY

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	199/201 (99%)	168 (84%)	31 (16%)	2	6
1	B	187/201 (93%)	156 (83%)	31 (17%)	2	5
1	C	191/201 (95%)	156 (82%)	35 (18%)	1	4
1	D	192/201 (96%)	162 (84%)	30 (16%)	2	6
All	All	769/804 (96%)	642 (84%)	127 (16%)	2	5

All (127) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	68	PHE
1	A	71	THR
1	A	83	GLN
1	A	93	LYS
1	A	95	ILE
1	A	96	THR

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Mol	Chain	Res	Type
1	A	99	SER
1	A	100	ASP
1	A	104	SER
1	A	105	ILE
1	A	111	GLN
1	A	113	VAL
1	A	115	THR
1	A	119	ILE
1	A	130	SER
1	A	131	ASP
1	A	149	THR
1	A	154	SER
1	A	163	ASP
1	A	169	LEU
1	A	172	ASP
1	A	177	GLU
1	A	186	ILE
1	A	197	ASN
1	A	220	ARG
1	A	234	LYS
1	A	265	SER
1	A	267	ARG
1	A	280	LEU
1	A	283	GLU
1	B	42	ASP
1	B	54	THR
1	B	59	GLU
1	B	68	PHE
1	B	71	THR
1	B	88	LEU
1	B	89	THR
1	B	96	THR
1	B	102	ILE
1	B	106	SER
1	B	115	THR
1	B	124	ASP
1	B	130	SER
1	B	139	LEU
1	B	149	THR
1	B	157	ASP
1	B	158	THR
1	B	179	ASP

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Mol	Chain	Res	Type
1	B	197	ASN
1	B	209	ASP
1	B	219	LYS
1	B	220	ARG
1	B	221	ILE
1	B	232	ILE
1	B	234	LYS
1	B	243	ARG
1	B	245	ASN
1	B	247	ILE
1	B	255	THR
1	B	265	SER
1	B	280	LEU
1	C	29	ASP
1	C	36	GLU
1	C	39	ARG
1	C	40	ARG
1	C	54	THR
1	C	68	PHE
1	C	70	SER
1	C	83	GLN
1	C	88	LEU
1	C	89	THR
1	C	93	LYS
1	C	94	LEU
1	C	95	ILE
1	C	99	SER
1	C	106	SER
1	C	115	THR
1	C	117	MET
1	C	118	THR
1	C	119	ILE
1	C	131	ASP
1	C	145(D)	THR
1	C	154	SER
1	C	158	THR
1	C	161	ARG
1	C	163	ASP
1	C	169	LEU
1	C	170	ASN
1	C	180	THR
1	C	209	ASP

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Mol	Chain	Res	Type
1	C	220	ARG
1	C	243	ARG
1	C	255	THR
1	C	273	ASP
1	C	277	ARG
1	C	280	LEU
1	D	32	ASP
1	D	68	PHE
1	D	86	ASN
1	D	88	LEU
1	D	92	ASP
1	D	94	LEU
1	D	98	THR
1	D	99	SER
1	D	100	ASP
1	D	102	ILE
1	D	111	GLN
1	D	115	THR
1	D	159	VAL
1	D	169	LEU
1	D	171	ARG
1	D	178	ARG
1	D	186	ILE
1	D	197	ASN
1	D	202	ASP
1	D	233	ASP
1	D	234	LYS
1	D	235	THR
1	D	237	THR
1	D	255	THR
1	D	258	PRO
1	D	267	ARG
1	D	280	LEU
1	D	281	LEU
1	D	283	GLU
1	D	291	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	A	90	HIS

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Mol	Chain	Res	Type
1	A	111	GLN
1	A	121	GLN
1	A	192	GLN
1	B	110	GLN
1	B	111	GLN
1	B	121	GLN
1	B	192	GLN
1	B	197	ASN
1	C	111	GLN
1	C	121	GLN
1	C	170	ASN
1	C	184	HIS
1	C	192	GLN
1	D	86	ASN
1	D	112	HIS
1	D	121	GLN
1	D	192	GLN
1	D	197	ASN
1	D	214	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 ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	AXL	C	301	1	17,26,26	0.62	0	22,38,38	0.92	1 (4%)
2	PO4	A	302	-	4,4,4	0.92	0	6,6,6	0.43	0
2	PO4	C	302	-	4,4,4	2.94	4 (100%)	6,6,6	0.42	0
2	PO4	D	301	-	4,4,4	2.95	4 (100%)	6,6,6	0.43	0
2	PO4	A	301	-	4,4,4	2.96	4 (100%)	6,6,6	0.43	0
2	PO4	A	303	-	4,4,4	2.97	4 (100%)	6,6,6	0.43	0
2	PO4	D	302	-	4,4,4	2.94	4 (100%)	6,6,6	0.42	0
3	AXL	B	301	1	17,26,26	0.63	0	22,38,38	0.80	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	AXL	C	301	1	-	2/12/37/37	0/2/2/2
3	AXL	B	301	1	-	1/12/37/37	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	303	PO4	P-O1	4.55	1.61	1.50
2	A	301	PO4	P-O1	4.53	1.61	1.50
2	D	301	PO4	P-O1	4.49	1.61	1.50
2	D	302	PO4	P-O1	4.49	1.61	1.50
2	C	302	PO4	P-O1	4.47	1.61	1.50
2	D	302	PO4	P-O3	2.30	1.61	1.54
2	A	301	PO4	P-O3	2.30	1.61	1.54
2	D	301	PO4	P-O2	2.29	1.61	1.54
2	C	302	PO4	P-O3	2.29	1.61	1.54
2	D	301	PO4	P-O3	2.28	1.61	1.54
2	C	302	PO4	P-O2	2.28	1.61	1.54
2	A	303	PO4	P-O2	2.28	1.61	1.54
2	A	303	PO4	P-O3	2.27	1.61	1.54
2	D	302	PO4	P-O2	2.25	1.61	1.54
2	A	301	PO4	P-O2	2.25	1.61	1.54
2	A	301	PO4	P-O4	-2.07	1.48	1.54
2	D	301	PO4	P-O4	-2.05	1.48	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	302	PO4	P-O4	-2.04	1.48	1.54
2	C	302	PO4	P-O4	-2.04	1.48	1.54
2	A	303	PO4	P-O4	-2.03	1.48	1.54

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	301	AXL	C12-C10-C9	-2.44	106.94	112.00

There are no chirality outliers.

All (3) torsion outliers are listed below:

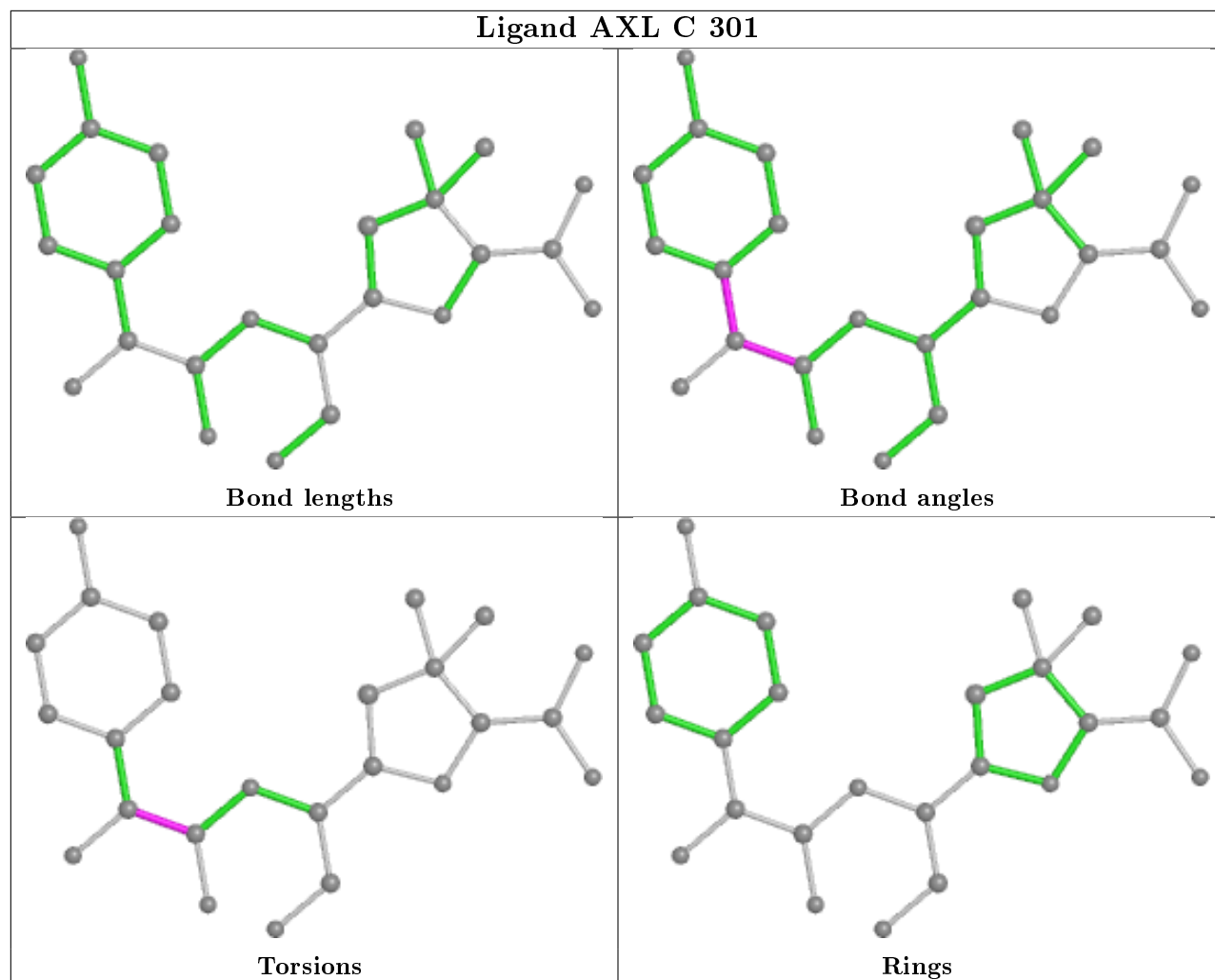
Mol	Chain	Res	Type	Atoms
3	C	301	AXL	N11-C10-C9-O91
3	B	301	AXL	C9-C10-C12-C17
3	C	301	AXL	N11-C10-C9-N8

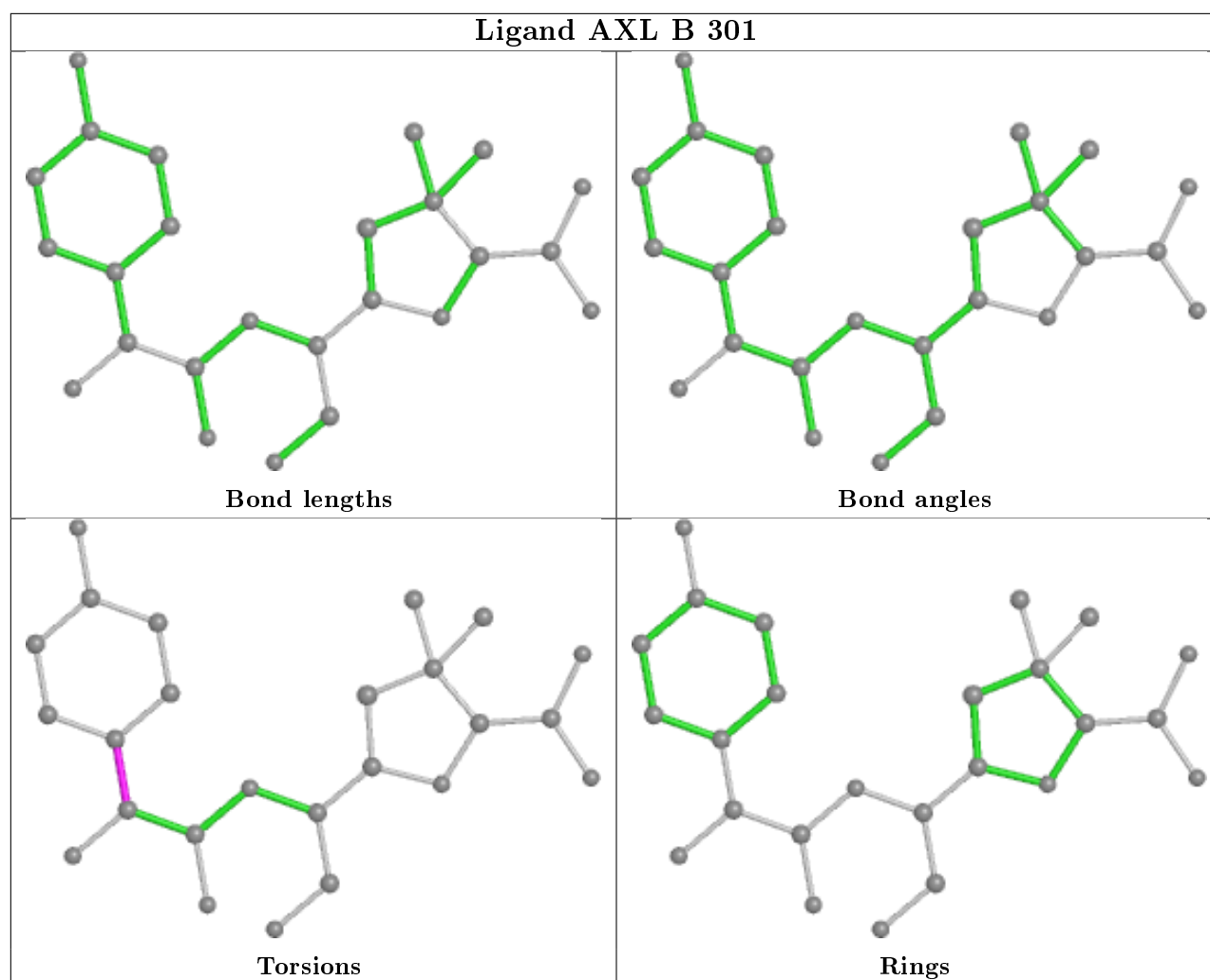
There are no ring outliers.

4 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	301	AXL	8	0
2	D	301	PO4	1	0
2	A	301	PO4	1	0
3	B	301	AXL	17	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	265/265 (100%)	-0.02	7 (2%) 56 52	33, 54, 83, 94	0
1	B	265/265 (100%)	-0.09	9 (3%) 45 39	34, 56, 110, 183	0
1	C	265/265 (100%)	-0.10	5 (1%) 66 64	37, 56, 100, 146	1 (0%)
1	D	265/265 (100%)	-0.04	6 (2%) 60 57	36, 59, 87, 103	0
All	All	1060/1060 (100%)	-0.06	27 (2%) 57 54	33, 56, 93, 183	1 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	164	ALA	5.1
1	B	165	GLU	4.4
1	D	99	SER	3.3
1	C	170	ASN	3.3
1	A	113	VAL	3.2
1	C	167	PRO	3.1
1	B	53	GLY	3.1
1	D	53	GLY	3.0
1	B	166	ALA	3.0
1	C	169	LEU	2.9
1	B	145(A)	GLY	2.8
1	D	145	PRO	2.8
1	B	170	ASN	2.6
1	C	145(C)	GLY	2.6
1	C	54	THR	2.4
1	A	98	THR	2.3
1	D	113	VAL	2.3
1	D	123	CYS	2.2
1	A	202	ASP	2.2
1	A	116	GLY	2.2
1	B	173	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	100	ASP	2.1
1	A	115	THR	2.1
1	A	100	ASP	2.1
1	A	88	LEU	2.1
1	B	144	GLY	2.0
1	B	167	PRO	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](#)

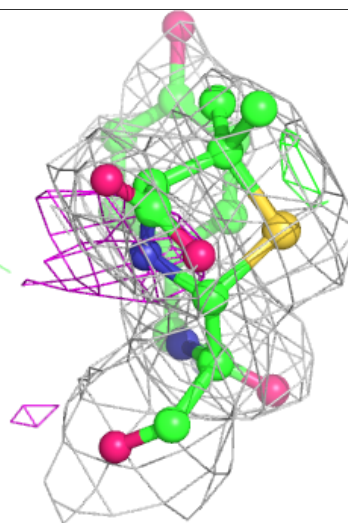
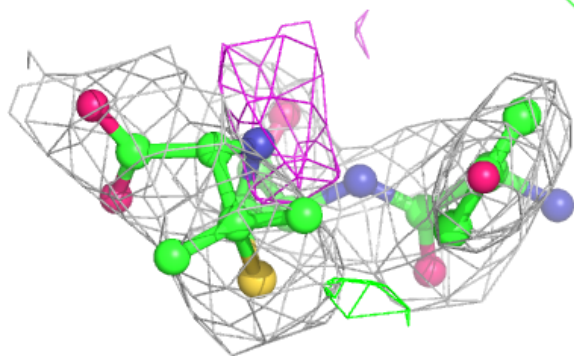
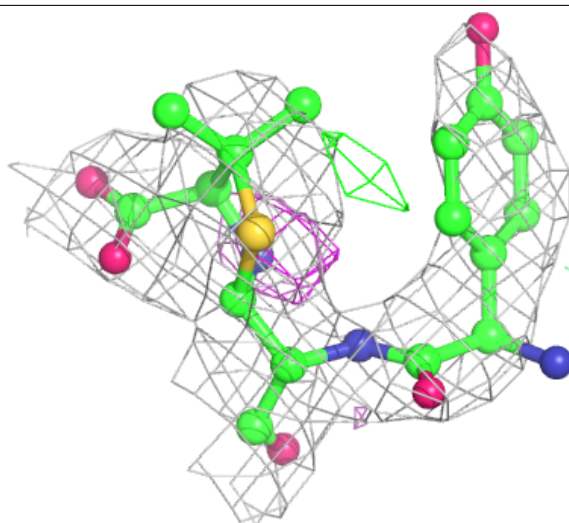
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

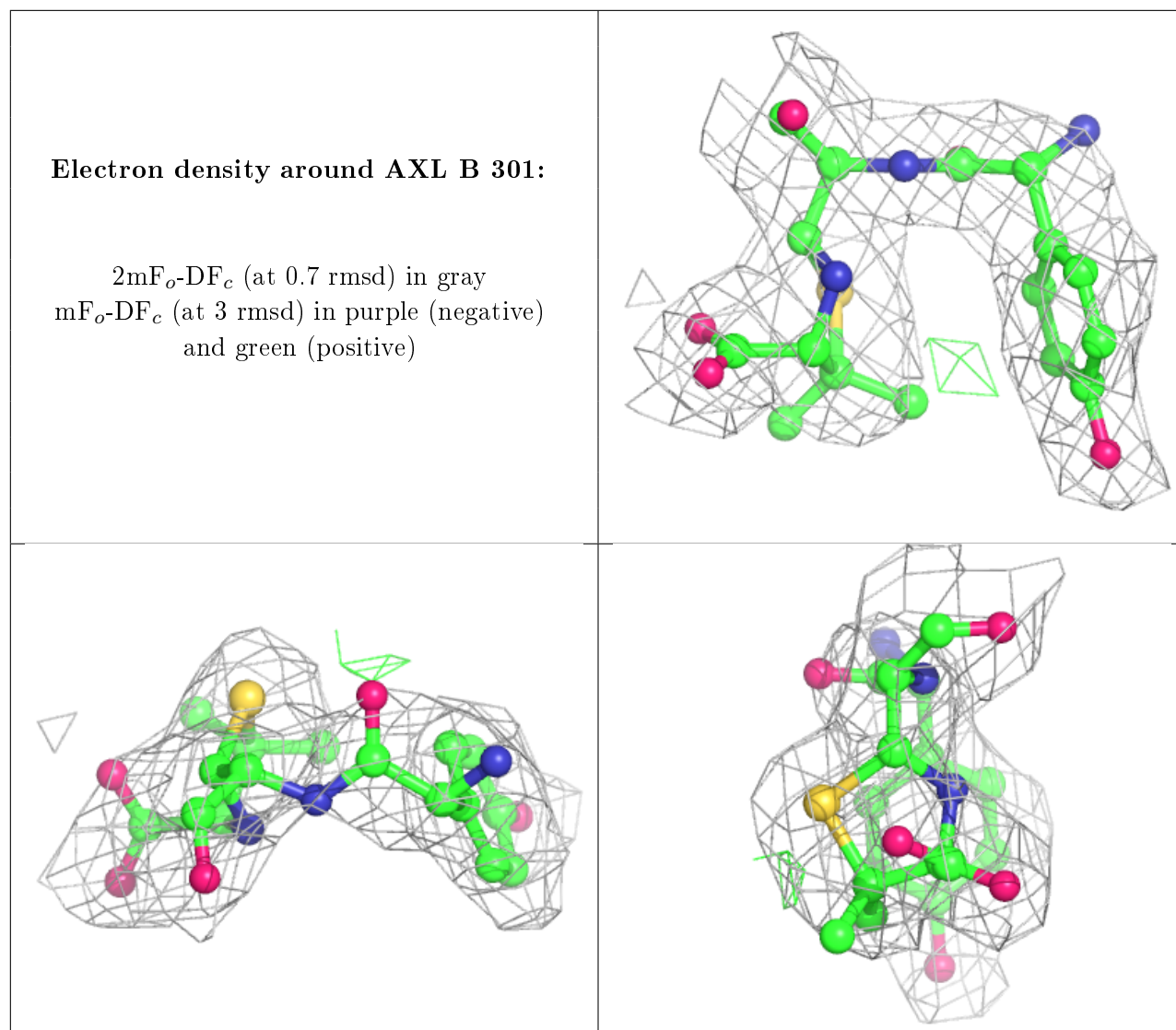
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PO4	A	303	5/5	0.84	0.32	91,107,116,132	0
2	PO4	D	302	5/5	0.84	0.21	110,116,123,133	0
3	AXL	C	301	25/25	0.87	0.23	60,64,66,67	25
2	PO4	C	302	5/5	0.89	0.19	68,80,129,130	0
2	PO4	D	301	5/5	0.89	0.24	95,111,124,125	0
3	AXL	B	301	25/25	0.89	0.24	56,60,62,64	25
2	PO4	A	302	5/5	0.91	0.32	118,118,118,119	0
2	PO4	A	301	5/5	0.96	0.20	66,67,88,135	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 AXL C 301:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.