



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:34 PM UTC

PDB ID : 9NST / pdb_00009nst
Title : Bacterial Pictet-Spenglerase KslB in complex with product of L-Trp and α -ketoglutaric acid
Authors : Kim, K.; Kim, W.
Deposited on : 2025-03-17
Resolution : 3.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

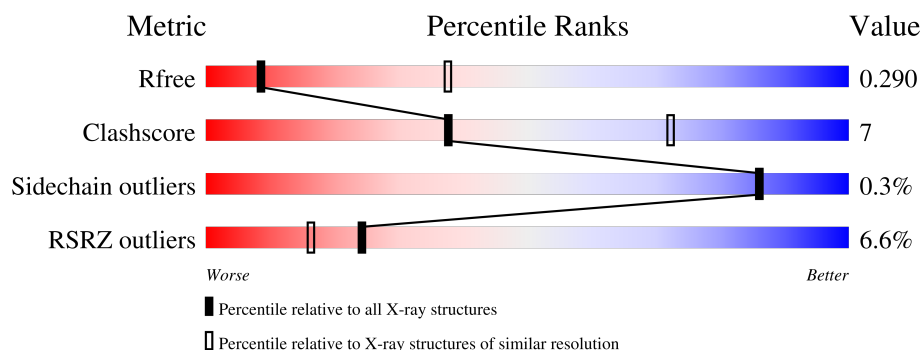
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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 of 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	317	
1	B	317	
1	C	317	
1	D	317	
1	E	317	
1	F	317	

2 Entry composition [i](#)

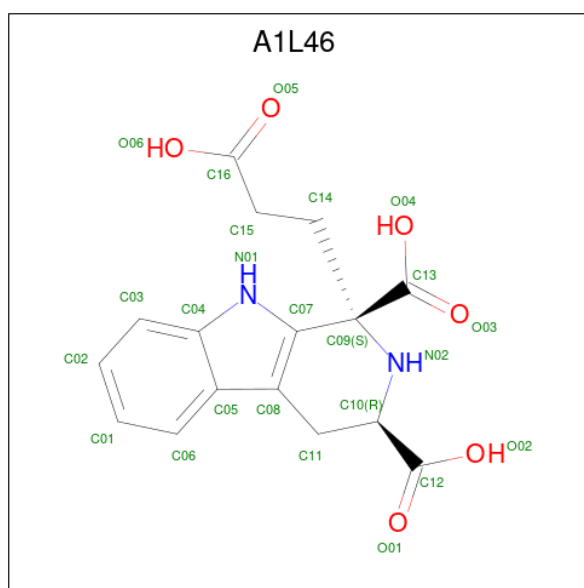
There are 2 unique types of molecules in this entry. The entry contains 14219 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pictet-Spenglerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2394	1531	404	455	4			
1	B	308	Total	C	N	O	S	0	0	0
			2423	1547	407	466	3			
1	C	288	Total	C	N	O	S	0	0	0
			2280	1461	383	433	3			
1	D	303	Total	C	N	O	S	0	0	0
			2393	1530	402	458	3			
1	E	292	Total	C	N	O	S	0	0	0
			2320	1488	390	439	3			
1	F	291	Total	C	N	O	S	0	0	0
			2313	1482	389	438	4			

- Molecule 2 is (1 {S},3 {S})-1-(3-hydroxy-3-oxopropyl)-2,3,4,9-tetrahydropyrido[3,4-b]indole-1,3-dicarboxylic acid (CCD ID: A1L46) (formula: C₁₆H₁₆N₂O₆) (labeled as "Ligand of Interest" by depositor).

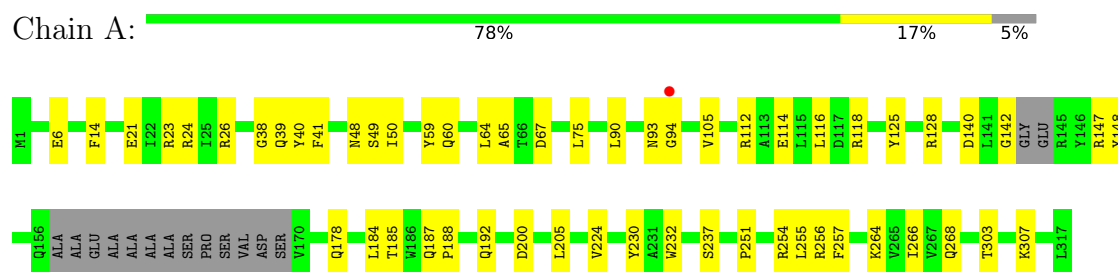


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 24	C 16	N 2	O 6	0	0
2	B	1	Total 24	C 16	N 2	O 6	0	0
2	C	1	Total 24	C 16	N 2	O 6	0	0
2	D	1	Total 24	C 16	N 2	O 6	0	0

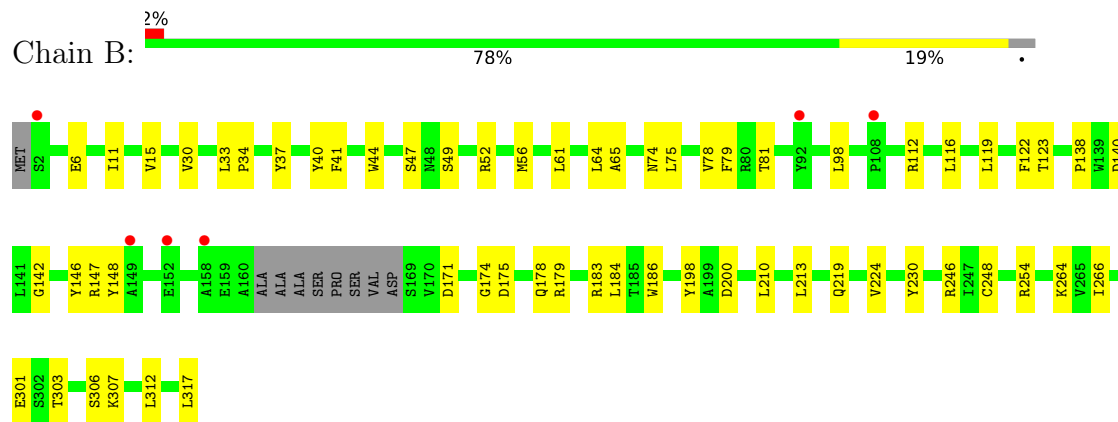
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

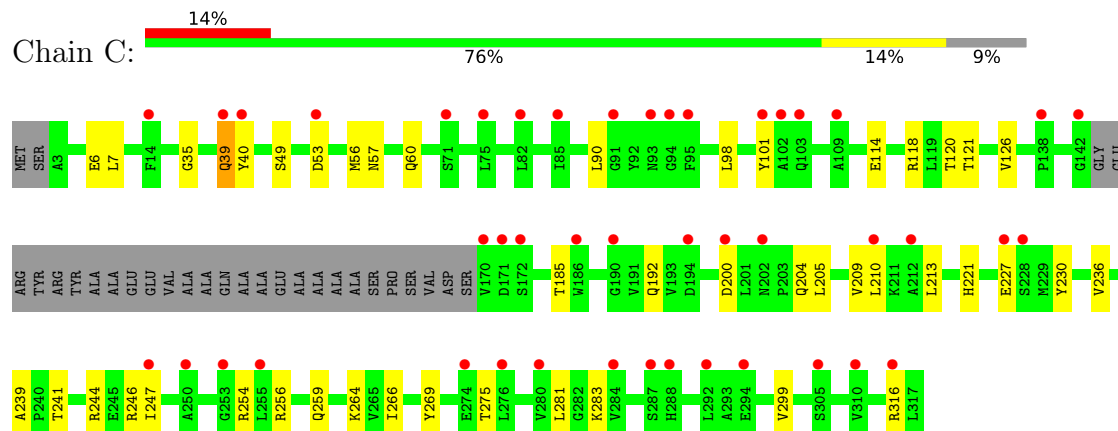
• Molecule 1: Pictet-Spenglerase



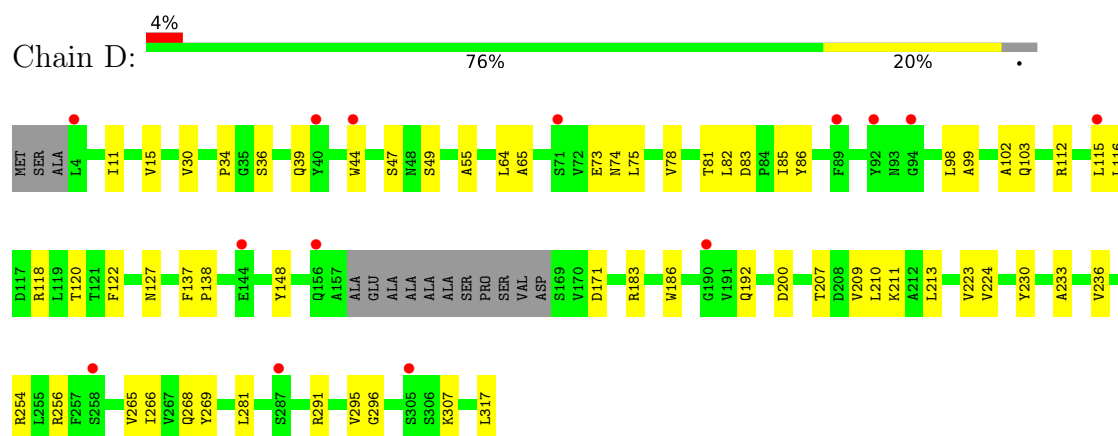
• Molecule 1: Pictet-Spenglerase



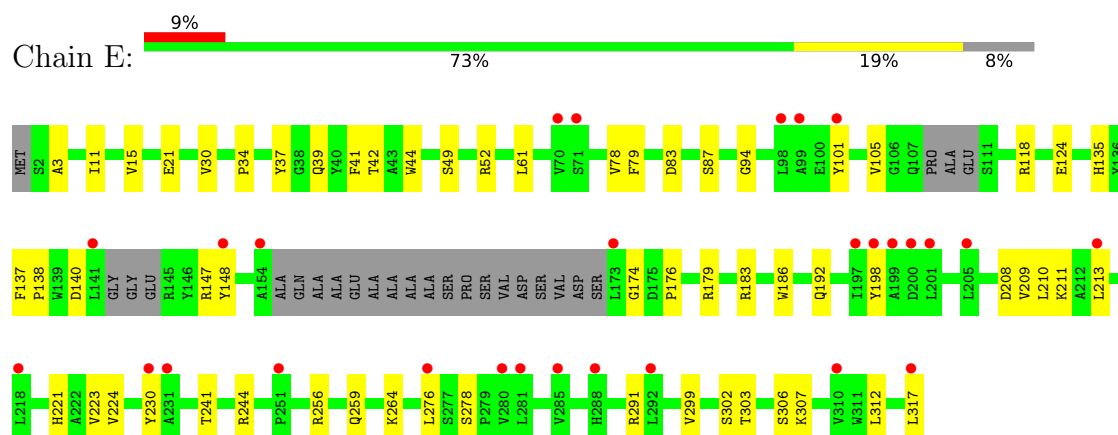
• Molecule 1: Pictet-Spenglerase



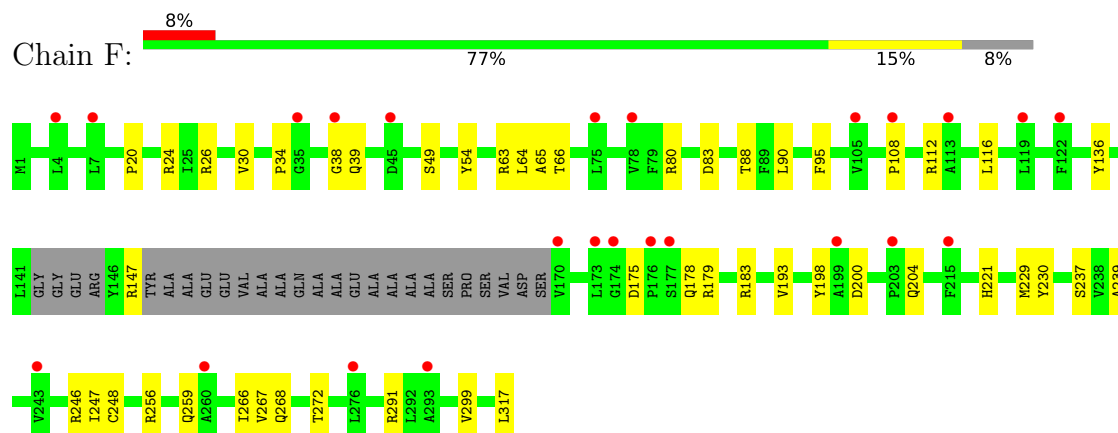
• Molecule 1: Pictet-Spenglerase



• Molecule 1: Pictet-Spenglerase



• Molecule 1: Pictet-Spenglerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	95.57Å 95.57Å 193.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.78 – 3.30 47.78 – 3.30	Depositor EDS
% Data completeness (in resolution range)	97.3 (47.78-3.30) 97.3 (47.78-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.27 (at 3.33Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.194 , 0.213 (Not available) , 0.290	Depositor DCC
R_{free} test set	2020 reflections (6.85%)	wwPDB-VP
Wilson B-factor (Å ²)	62.7	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.075 for -h,-k,l 0.388 for h,-h-k,-l 0.077 for -k,-h,-l	Xtriage
Reported twinning fraction	0.480 for h,-h-k,-l	Depositor
Outliers	0 of 28837 reflections	Xtriage
F_o, F_c correlation	0.80	EDS
Total number of atoms	14219	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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: A1L46

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	0.11	0/2455	0.32	0/3357
1	B	0.15	0/2485	0.33	0/3401
1	C	0.20	0/2340	0.34	0/3204
1	D	0.12	0/2455	0.30	0/3360
1	E	0.11	0/2379	0.30	0/3253
1	F	0.10	0/2373	0.30	0/3246
All	All	0.14	0/14487	0.31	0/19821

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	C	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	316	ARG	Sidechain

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	2394	0	2347	42	0
1	B	2423	0	2360	40	0
1	C	2280	0	2238	30	0
1	D	2393	0	2334	39	0
1	E	2320	0	2271	40	0
1	F	2313	0	2273	31	0
2	A	24	0	0	3	0
2	B	24	0	0	1	0
2	C	24	0	0	1	0
2	D	24	0	0	2	0
All	All	14219	0	13823	195	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 7.

All (195) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:ARG:NH1	1:B:248:CYS:SG	2.57	0.77
1:B:64:LEU:HB3	1:B:75:LEU:HD21	1.70	0.72
1:F:246:ARG:NH1	1:F:248:CYS:SG	2.65	0.70
1:E:176:PRO:HA	1:E:179:ARG:HG3	1.73	0.69
1:E:3:ALA:HB1	1:E:118:ARG:HG2	1.76	0.67
1:F:90:LEU:HB3	1:F:95:PHE:HB2	1.78	0.66
1:A:256:ARG:NH1	1:A:268:GLN:OE1	2.29	0.66
1:F:20:PRO:HB2	1:F:24:ARG:HH21	1.62	0.65
1:F:175:ASP:HB3	1:F:178:GLN:HG3	1.79	0.65
1:C:57:ASN:HA	1:D:82:LEU:HD21	1.77	0.65
1:E:210:LEU:HD23	1:E:213:LEU:HD12	1.80	0.64
1:A:185:THR:HG23	1:A:192:GLN:HG2	1.79	0.63
1:E:208:ASP:HA	1:E:211:LYS:HD2	1.80	0.63
1:B:65:ALA:HB3	1:B:116:LEU:HD13	1.80	0.62
1:D:192:GLN:O	1:D:291:ARG:NH1	2.32	0.62
1:E:105:VAL:HA	1:E:118:ARG:HD2	1.82	0.61
1:E:192:GLN:O	1:E:291:ARG:NH2	2.34	0.61
1:E:78:VAL:HA	1:F:64:LEU:HD21	1.84	0.60
1:D:207:THR:HG22	1:D:211:LYS:HD2	1.84	0.59
1:D:64:LEU:HB3	1:D:75:LEU:HD21	1.83	0.59
1:A:50:ILE:HB	1:A:90:LEU:HD21	1.85	0.59
1:F:256:ARG:NH2	1:F:268:GLN:OE1	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:THR:O	1:B:307:LYS:NZ	2.35	0.58
1:A:41:PHE:HB3	1:B:142:GLY:HA3	1.86	0.58
1:A:230:TYR:HB3	1:A:266:ILE:HG12	1.86	0.57
1:D:256:ARG:NH1	2:D:401:A1L46:O05	2.37	0.57
1:B:184:LEU:HB3	1:B:312:LEU:HD11	1.86	0.57
1:D:115:LEU:HA	1:D:118:ARG:HD2	1.86	0.57
1:C:230:TYR:HB3	1:C:266:ILE:HG12	1.87	0.56
1:E:221:HIS:HB2	1:E:299:VAL:HG12	1.87	0.56
1:A:178:GLN:HG3	1:A:200:ASP:HB3	1.87	0.56
1:F:229:MET:HE3	1:F:267:VAL:HB	1.88	0.56
1:E:37:TYR:O	1:F:147:ARG:NH2	2.39	0.55
1:A:264:LYS:HZ1	2:A:401:A1L46:C13	2.19	0.55
1:D:30:VAL:HG23	1:D:34:PRO:HD2	1.88	0.55
1:A:40:TYR:CZ	1:B:146:TYR:HB3	2.42	0.54
1:F:230:TYR:HB3	1:F:266:ILE:HG12	1.90	0.54
1:A:187:GLN:HE21	1:A:188:PRO:HA	1.73	0.54
1:A:256:ARG:NH2	2:A:401:A1L46:O05	2.40	0.53
1:E:147:ARG:HA	1:F:39:GLN:HA	1.91	0.53
1:A:49:SER:HB3	1:B:49:SER:HB2	1.90	0.53
1:C:120:THR:HG23	1:C:236:VAL:HB	1.90	0.53
1:B:210:LEU:HD23	1:B:213:LEU:HD12	1.91	0.53
1:E:39:GLN:HE22	1:E:42:THR:HG23	1.74	0.53
1:E:174:GLY:HA3	1:E:179:ARG:HH22	1.72	0.53
1:C:221:HIS:HB2	1:C:299:VAL:HG12	1.90	0.53
1:C:200:ASP:OD1	1:C:200:ASP:N	2.36	0.53
1:D:210:LEU:HD23	1:D:213:LEU:HD12	1.90	0.53
1:E:303:THR:O	1:E:307:LYS:NZ	2.38	0.52
1:F:247:ILE:O	1:F:272:THR:OG1	2.27	0.52
1:E:148:TYR:N	1:F:38:GLY:O	2.40	0.51
1:D:233:ALA:HB2	1:D:265:VAL:HG23	1.92	0.51
1:A:21:GLU:HA	1:A:24:ARG:HG2	1.92	0.51
1:E:256:ARG:HD2	1:E:278:SER:HB3	1.93	0.51
1:E:230:TYR:HB2	1:E:264:LYS:HD2	1.91	0.51
1:C:264:LYS:HZ1	2:C:401:A1L46:C13	2.23	0.51
1:D:137:PHE:O	1:D:307:LYS:NZ	2.39	0.51
1:D:209:VAL:HG11	1:D:281:LEU:HD21	1.93	0.51
1:A:232:TRP:CZ2	1:A:264:LYS:HD3	2.46	0.50
1:B:230:TYR:HB3	1:B:266:ILE:HG12	1.93	0.50
1:A:251:PRO:HD2	1:A:254:ARG:HE	1.75	0.50
1:C:241:THR:OG1	1:C:244:ARG:NH1	2.44	0.50
1:F:54:TYR:OH	1:F:83:ASP:OD1	2.24	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:LEU:HD13	1:B:79:PHE:HD1	1.77	0.50
1:F:26:ARG:HD3	1:F:136:TYR:HB3	1.93	0.50
1:A:93:ASN:HD21	1:B:224:VAL:HG11	1.75	0.50
1:A:105:VAL:HA	1:A:118:ARG:HD2	1.94	0.50
1:D:224:VAL:HB	2:D:401:A1L46:O02	2.12	0.49
1:E:11:ILE:HD13	1:E:124:GLU:HB3	1.94	0.49
1:E:61:LEU:HD13	1:E:79:PHE:HD1	1.78	0.49
1:B:264:LYS:NZ	2:B:401:A1L46:O04	2.29	0.48
1:A:147:ARG:NH2	1:B:37:TYR:O	2.46	0.48
1:A:14:PHE:HE2	1:A:125:TYR:HE1	1.61	0.48
1:A:128:ARG:NH2	1:A:232:TRP:O	2.47	0.48
1:E:183:ARG:HB2	1:E:317:LEU:HD11	1.96	0.48
1:A:60:GLN:HB3	1:B:81:THR:HG22	1.95	0.48
1:B:183:ARG:HB2	1:B:317:LEU:HD11	1.95	0.48
1:B:301:GLU:O	1:B:306:SER:N	2.31	0.48
1:C:35:GLY:HA2	1:D:224:VAL:HA	1.94	0.48
1:D:65:ALA:HB3	1:D:116:LEU:HD13	1.95	0.48
1:E:223:VAL:HG23	1:E:224:VAL:HG23	1.96	0.48
1:D:44:TRP:CZ2	1:D:138:PRO:HD3	2.49	0.48
1:A:184:LEU:HD21	1:A:255:LEU:HD21	1.95	0.48
1:F:65:ALA:O	1:F:112:ARG:NH1	2.46	0.47
1:C:254:ARG:NH1	1:C:283:LYS:HB2	2.29	0.47
1:D:230:TYR:HB3	1:D:266:ILE:HG12	1.95	0.47
1:B:44:TRP:CH2	1:B:138:PRO:HD3	2.50	0.47
1:A:67:ASP:O	1:A:112:ARG:NH1	2.48	0.47
1:B:186:TRP:HE1	1:B:219:GLN:HE21	1.63	0.47
1:E:276:LEU:HD22	1:F:88:THR:HB	1.95	0.47
1:B:174:GLY:HA3	1:B:200:ASP:OD2	2.15	0.46
1:C:114:GLU:OE2	1:C:118:ARG:NH2	2.48	0.46
1:E:148:TYR:H	1:F:38:GLY:C	2.23	0.46
1:B:179:ARG:HG3	1:B:198:TYR:HA	1.96	0.46
1:C:6:GLU:H	1:C:6:GLU:CD	2.23	0.46
1:C:6:GLU:HB2	1:C:101:TYR:OH	2.16	0.46
1:E:44:TRP:CH2	1:E:138:PRO:HD3	2.51	0.46
1:E:49:SER:HB3	1:F:49:SER:HB2	1.95	0.46
1:C:204:GLN:HB2	1:C:239:ALA:HB1	1.98	0.46
1:D:11:ILE:O	1:D:15:VAL:HG23	2.16	0.46
1:D:65:ALA:O	1:D:112:ARG:HD3	2.15	0.46
1:D:120:THR:HG23	1:D:236:VAL:HB	1.98	0.46
1:F:237:SER:HB3	1:F:259:GLN:HG2	1.98	0.46
1:D:269:TYR:CZ	1:D:296:GLY:HA3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:52:ARG:NH2	1:F:49:SER:OG	2.49	0.45
1:E:137:PHE:O	1:E:307:LYS:NZ	2.37	0.45
1:E:140:ASP:OD1	1:E:140:ASP:N	2.48	0.45
1:D:200:ASP:OD1	1:D:200:ASP:N	2.37	0.45
1:E:39:GLN:OE1	1:E:41:PHE:N	2.50	0.45
1:B:175:ASP:HB3	1:B:178:GLN:HG3	1.98	0.45
1:D:183:ARG:HB2	1:D:317:LEU:HD11	1.98	0.45
1:C:39:GLN:HE21	1:C:39:GLN:HB2	1.65	0.45
1:D:186:TRP:CD2	1:D:295:VAL:HG11	2.51	0.45
1:F:221:HIS:HB2	1:F:299:VAL:HG12	1.98	0.45
1:A:59:TYR:OH	1:A:237:SER:O	2.23	0.44
1:B:30:VAL:HG23	1:B:34:PRO:HD2	1.99	0.44
1:E:259:GLN:H	1:E:259:GLN:HG2	1.62	0.44
1:A:38:GLY:HA2	1:B:148:TYR:HB2	1.99	0.44
1:A:64:LEU:HB2	1:A:75:LEU:HD21	2.00	0.44
1:B:98:LEU:HG	1:B:122:PHE:HZ	1.82	0.44
1:D:73:GLU:H	1:D:73:GLU:CD	2.25	0.44
1:E:209:VAL:O	1:E:213:LEU:HG	2.17	0.44
1:F:30:VAL:HB	1:F:34:PRO:HD2	1.99	0.44
1:B:52:ARG:O	1:B:56:MET:HG3	2.17	0.44
1:B:171:ASP:CG	1:B:254:ARG:HH22	2.24	0.44
1:C:185:THR:HG23	1:C:192:GLN:HG2	2.00	0.44
1:A:65:ALA:HB3	1:A:116:LEU:HD13	1.99	0.44
1:C:209:VAL:HG11	1:C:281:LEU:HD21	1.98	0.44
1:A:6:GLU:H	1:A:6:GLU:CD	2.26	0.44
1:D:74:ASN:O	1:D:78:VAL:HG13	2.18	0.44
1:D:223:VAL:HG23	1:D:224:VAL:HG23	2.00	0.44
1:E:135:HIS:O	1:E:307:LYS:HD2	2.18	0.43
1:A:205:LEU:HD21	1:A:257:PHE:HB3	2.00	0.43
1:C:227:GLU:HB3	1:C:269:TYR:CZ	2.53	0.43
1:A:224:VAL:HB	2:A:401:A1L46:O02	2.18	0.43
1:C:60:GLN:HB3	1:D:81:THR:HG22	1.99	0.43
1:F:179:ARG:NH1	1:F:200:ASP:OD1	2.50	0.43
1:C:49:SER:HB3	1:D:49:SER:HB2	1.99	0.43
1:C:53:ASP:HB3	1:D:86:TYR:OH	2.19	0.43
1:E:11:ILE:O	1:E:15:VAL:HG13	2.18	0.43
1:A:21:GLU:HB3	1:A:94:GLY:O	2.18	0.43
1:D:171:ASP:CG	1:D:254:ARG:HH22	2.26	0.43
1:B:65:ALA:O	1:B:112:ARG:HD3	2.18	0.43
1:C:7:LEU:HD23	1:C:121:THR:HG22	2.01	0.43
1:C:40:TYR:HB3	1:D:148:TYR:CZ	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:ASP:OD2	1:D:102:ALA:HB1	2.19	0.42
1:A:140:ASP:OD1	1:A:140:ASP:N	2.47	0.42
1:F:183:ARG:HB2	1:F:317:LEU:HD11	2.00	0.42
1:A:41:PHE:CB	1:B:142:GLY:HA3	2.47	0.42
1:A:232:TRP:CE2	1:A:264:LYS:HD3	2.55	0.42
1:B:119:LEU:O	1:B:123:THR:OG1	2.36	0.42
1:C:56:MET:HE3	1:D:85:ILE:HG13	2.01	0.42
1:C:246:ARG:NH1	1:C:275:THR:O	2.52	0.42
1:D:256:ARG:NH2	1:D:268:GLN:OE1	2.48	0.42
1:A:39:GLN:HA	1:B:147:ARG:HA	2.00	0.42
1:F:193:VAL:HG12	1:F:291:ARG:HD2	2.02	0.42
1:A:303:THR:O	1:A:307:LYS:NZ	2.35	0.42
1:D:98:LEU:HG	1:D:122:PHE:HZ	1.84	0.42
1:F:63:ARG:O	1:F:66:THR:OG1	2.31	0.42
1:A:14:PHE:CE2	1:A:125:TYR:HE1	2.38	0.42
1:C:98:LEU:HD12	1:C:126:VAL:HG22	2.02	0.42
1:E:186:TRP:CE2	1:E:312:LEU:HD13	2.55	0.42
1:E:302:SER:HA	1:E:306:SER:HB2	2.02	0.42
1:F:20:PRO:O	1:F:24:ARG:HG3	2.19	0.42
1:B:6:GLU:CD	1:B:6:GLU:H	2.27	0.41
1:C:90:LEU:HD23	1:C:90:LEU:HA	1.86	0.41
1:A:142:GLY:HA3	1:B:41:PHE:HB3	2.02	0.41
1:C:205:LEU:HD13	1:C:259:GLN:HE21	1.85	0.41
1:A:148:TYR:CE1	1:B:33:LEU:HD22	2.55	0.41
1:B:74:ASN:O	1:B:78:VAL:HG13	2.20	0.41
1:B:140:ASP:OD1	1:B:140:ASP:N	2.49	0.41
1:E:21:GLU:HB3	1:E:94:GLY:O	2.21	0.41
1:E:30:VAL:HB	1:E:34:PRO:HD2	2.03	0.41
1:A:114:GLU:OE2	1:A:118:ARG:NH2	2.49	0.41
1:E:241:THR:OG1	1:E:244:ARG:NH1	2.53	0.41
1:C:247:ILE:HG23	1:C:256:ARG:NH1	2.36	0.41
1:B:11:ILE:O	1:B:15:VAL:HG23	2.19	0.41
1:E:83:ASP:OD1	1:E:87:SER:OG	2.38	0.41
1:E:179:ARG:HD3	1:E:198:TYR:CD1	2.56	0.41
1:F:80:ARG:HH22	1:F:108:PRO:HG3	1.85	0.41
1:B:98:LEU:HG	1:B:122:PHE:CZ	2.56	0.41
1:D:98:LEU:HG	1:D:122:PHE:CZ	2.56	0.41
1:D:99:ALA:O	1:D:103:GLN:HG3	2.21	0.41
1:A:65:ALA:O	1:A:112:ARG:HD3	2.21	0.40
1:E:101:TYR:O	1:E:105:VAL:HG23	2.21	0.40
1:A:148:TYR:CZ	1:B:40:TYR:HB3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:TRP:NE1	1:B:312:LEU:HD13	2.36	0.40
1:D:36:SER:O	1:D:39:GLN:HG3	2.21	0.40
1:F:179:ARG:HD2	1:F:198:TYR:CD1	2.57	0.40
1:C:56:MET:O	1:C:60:GLN:HG2	2.21	0.40
1:F:116:LEU:HD12	1:F:116:LEU:HA	1.92	0.40
1:A:23:ARG:HA	1:A:26:ARG:HG2	2.03	0.40
1:D:55:ALA:HB1	1:D:127:ASN:OD1	2.21	0.40
1:F:204:GLN:HB2	1:F:239:ALA:HB1	2.02	0.40
1:C:210:LEU:HA	1:C:213:LEU:HD12	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	260/268 (97%)	259 (100%)	1 (0%)	84	84
1	B	261/268 (97%)	260 (100%)	1 (0%)	84	84
1	C	250/268 (93%)	249 (100%)	1 (0%)	84	84
1	D	259/268 (97%)	258 (100%)	1 (0%)	84	84
1	E	252/268 (94%)	252 (100%)	0	100	100
1	F	254/268 (95%)	254 (100%)	0	100	100
All	All	1536/1608 (96%)	1532 (100%)	4 (0%)	86	86

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	ASN
1	B	47	SER
1	C	39	GLN
1	D	47	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	ASN
1	A	187	GLN
1	A	192	GLN
1	B	74	ASN
1	B	93	ASN
1	B	103	GLN
1	B	219	GLN
1	C	39	GLN
1	C	93	ASN
1	C	192	GLN
1	C	219	GLN
1	E	103	GLN
1	E	219	GLN
1	F	103	GLN
1	F	219	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	A1L46	C	401	-	24,26,26	1.77	4 (16%)	29,39,39	2.31	4 (13%)
2	A1L46	D	401	-	24,26,26	1.72	4 (16%)	29,39,39	2.67	4 (13%)
2	A1L46	A	401	-	24,26,26	1.69	4 (16%)	29,39,39	2.48	3 (10%)
2	A1L46	B	401	-	24,26,26	1.82	5 (20%)	29,39,39	2.50	4 (13%)

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	A1L46	C	401	-	-	7/16/31/31	0/3/3/3
2	A1L46	D	401	-	-	12/16/31/31	0/3/3/3
2	A1L46	A	401	-	-	4/16/31/31	0/3/3/3
2	A1L46	B	401	-	-	9/16/31/31	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	A1L46	C09-C13	-5.97	1.48	1.54
2	C	401	A1L46	C09-C13	-5.80	1.49	1.54
2	D	401	A1L46	C09-C13	-5.46	1.49	1.54
2	A	401	A1L46	C09-C13	-5.33	1.49	1.54
2	A	401	A1L46	C05-C04	-3.17	1.37	1.41
2	B	401	A1L46	C05-C04	-3.13	1.37	1.41
2	D	401	A1L46	C05-C04	-3.10	1.37	1.41
2	C	401	A1L46	C05-C04	-3.07	1.37	1.41
2	A	401	A1L46	C04-N01	-2.68	1.33	1.38
2	C	401	A1L46	C04-N01	-2.67	1.33	1.38
2	B	401	A1L46	C04-N01	-2.65	1.34	1.38
2	D	401	A1L46	C10-N02	2.45	1.49	1.47
2	D	401	A1L46	C04-N01	-2.44	1.34	1.38
2	B	401	A1L46	C07-N01	-2.41	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	A1L46	C07-N01	-2.20	1.34	1.38
2	B	401	A1L46	C10-N02	2.18	1.49	1.47
2	C	401	A1L46	C07-N01	-2.06	1.34	1.38

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	A1L46	C09-C07-N01	9.10	133.09	123.84
2	A	401	A1L46	C10-C11-C08	8.58	117.51	108.24
2	D	401	A1L46	C10-C11-C08	8.07	116.97	108.24
2	B	401	A1L46	C10-C11-C08	8.02	116.91	108.24
2	B	401	A1L46	C09-C07-N01	7.97	131.94	123.84
2	C	401	A1L46	C09-C07-N01	7.58	131.55	123.84
2	A	401	A1L46	C09-C07-N01	7.56	131.52	123.84
2	C	401	A1L46	C10-C11-C08	7.08	115.89	108.24
2	D	401	A1L46	C09-C07-C08	-5.06	118.53	125.65
2	B	401	A1L46	C09-C07-C08	-4.20	119.74	125.65
2	A	401	A1L46	C09-C07-C08	-4.15	119.81	125.65
2	C	401	A1L46	C09-C07-C08	-3.89	120.18	125.65
2	D	401	A1L46	C03-C04-C05	-2.54	119.73	122.19
2	B	401	A1L46	C03-C04-C05	-2.48	119.79	122.19
2	C	401	A1L46	C03-C04-C05	-2.26	120.00	122.19

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	A1L46	N02-C10-C12-O01
2	B	401	A1L46	C07-C09-C13-O04
2	B	401	A1L46	C14-C09-C13-O04
2	B	401	A1L46	N02-C10-C12-O02
2	B	401	A1L46	C09-C14-C15-C16
2	C	401	A1L46	C07-C09-C13-O04
2	C	401	A1L46	C09-C14-C15-C16
2	D	401	A1L46	C07-C09-C13-O04
2	D	401	A1L46	N02-C10-C12-O01
2	C	401	A1L46	N02-C10-C12-O01
2	D	401	A1L46	C09-C14-C15-C16
2	B	401	A1L46	C14-C09-C13-O03
2	D	401	A1L46	C14-C09-C13-O03
2	D	401	A1L46	C14-C09-C13-O04
2	B	401	A1L46	N02-C10-C12-O01

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Mol	Chain	Res	Type	Atoms
2	B	401	A1L46	C13-C09-C14-C15
2	A	401	A1L46	N02-C10-C12-O02
2	C	401	A1L46	N02-C10-C12-O02
2	D	401	A1L46	N02-C10-C12-O02
2	B	401	A1L46	N02-C09-C14-C15
2	D	401	A1L46	C07-C09-C13-O03
2	D	401	A1L46	C11-C10-C12-O01
2	D	401	A1L46	C11-C10-C12-O02
2	C	401	A1L46	C14-C15-C16-O05
2	C	401	A1L46	C14-C15-C16-O06
2	A	401	A1L46	C14-C15-C16-O05
2	A	401	A1L46	C14-C15-C16-O06
2	D	401	A1L46	C14-C15-C16-O06
2	C	401	A1L46	C14-C09-C13-O04
2	D	401	A1L46	C14-C15-C16-O05
2	B	401	A1L46	C07-C09-C14-C15
2	D	401	A1L46	C07-C09-C14-C15

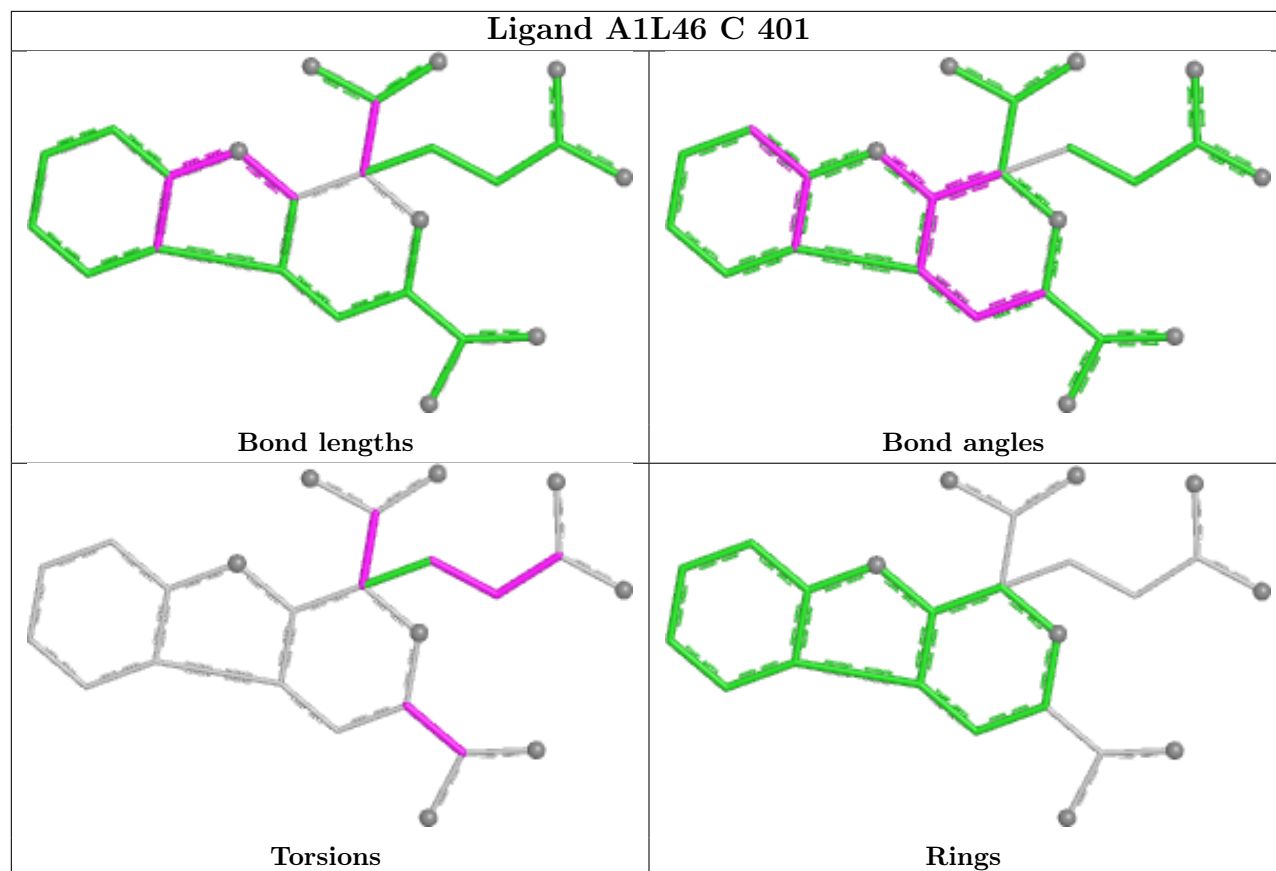
There are no ring outliers.

4 monomers are involved in 7 short contacts:

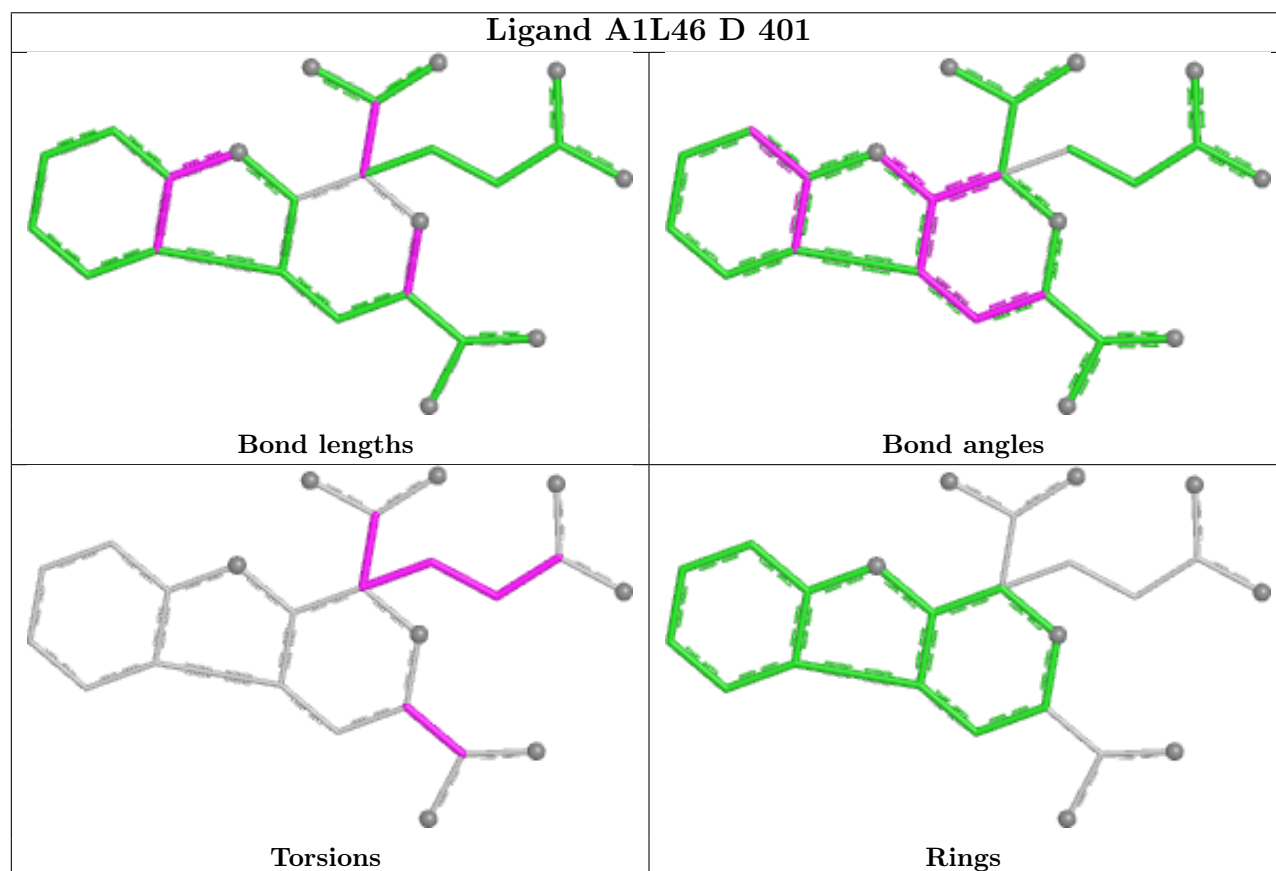
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	A1L46	1	0
2	D	401	A1L46	2	0
2	A	401	A1L46	3	0
2	B	401	A1L46	1	0

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.

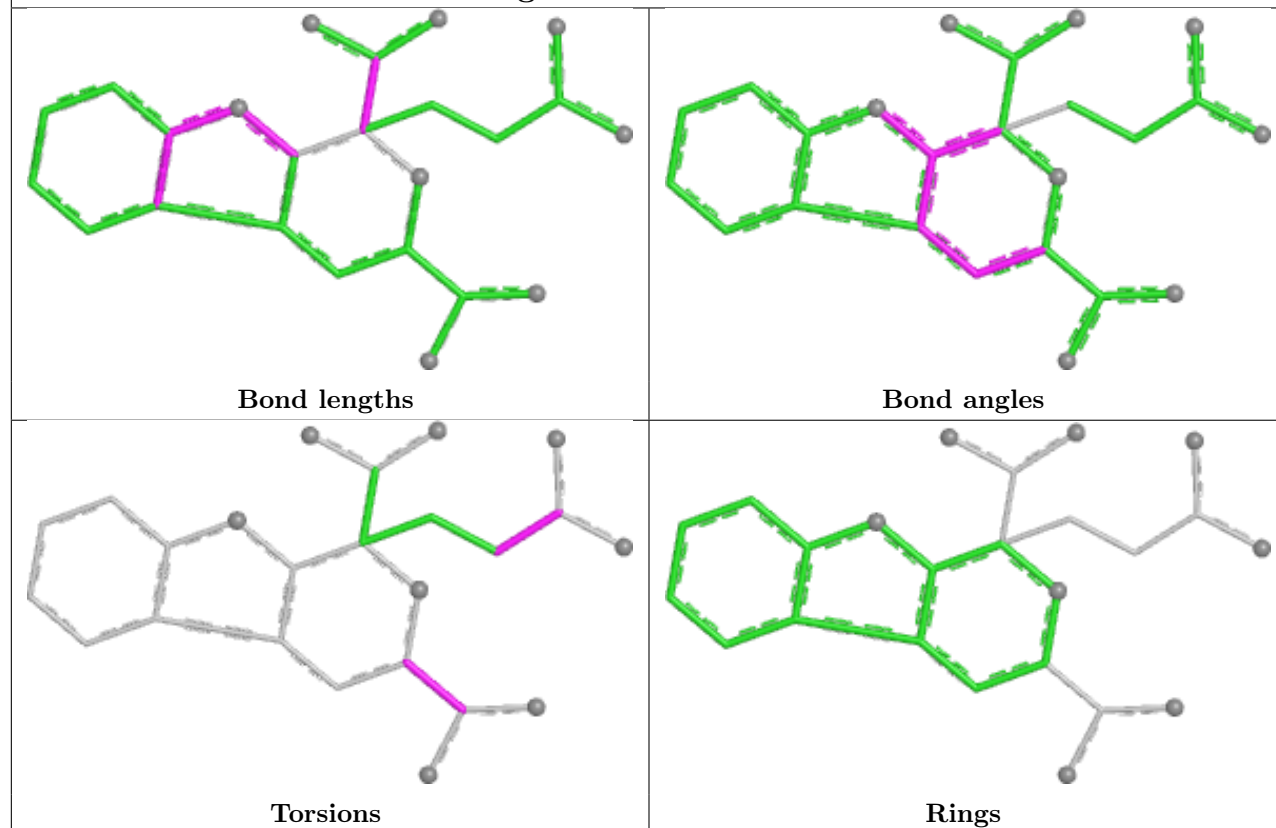
Ligand A1L46 C 401



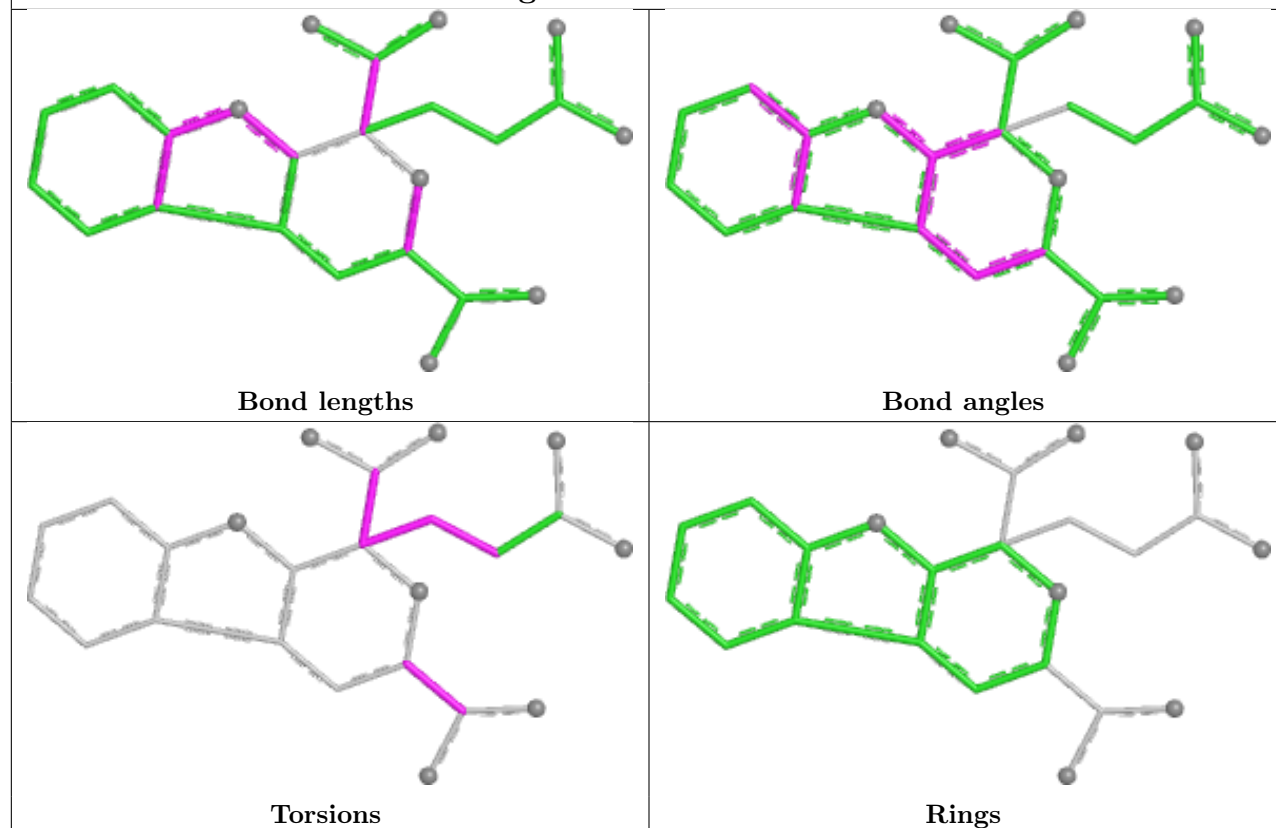
Ligand A1L46 D 401



Ligand A1L46 A 401



Ligand A1L46 B 401



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	302/317 (95%)	0.33	1 (0%) 90 82	22, 34, 53, 65	0
1	B	308/317 (97%)	0.37	6 (1%) 66 48	25, 37, 61, 81	0
1	C	288/317 (90%)	1.34	45 (15%) 5 4	47, 64, 88, 104	0
1	D	303/317 (95%)	0.97	14 (4%) 37 25	41, 60, 80, 94	0
1	E	292/317 (92%)	1.05	28 (9%) 13 11	45, 67, 89, 118	0
1	F	291/317 (91%)	1.00	24 (8%) 17 13	48, 65, 89, 104	0
All	All	1784/1902 (93%)	0.83	118 (6%) 24 16	22, 58, 83, 118	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	94	GLY	4.9
1	C	109	ALA	4.1
1	F	174	GLY	4.0
1	C	250	ALA	3.8
1	C	247	ILE	3.7
1	D	94	GLY	3.7
1	E	288	HIS	3.6
1	C	142	GLY	3.5
1	C	93	ASN	3.5
1	F	7	LEU	3.5
1	E	99	ALA	3.4
1	C	103	GLN	3.4
1	C	255	LEU	3.3
1	C	280	VAL	3.3
1	F	199	ALA	3.2
1	D	40	TYR	3.1
1	E	280	VAL	3.1
1	F	75	LEU	3.1
1	C	316	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	39	GLN	3.0
1	C	274	GLU	3.0
1	B	108	PRO	2.9
1	C	284	VAL	2.9
1	C	253	GLY	2.9
1	C	276	LEU	2.9
1	E	317	LEU	2.9
1	D	92	TYR	2.9
1	F	215	PHE	2.9
1	E	201	LEU	2.9
1	E	230	TYR	2.9
1	D	115	LEU	2.9
1	E	310	VAL	2.8
1	E	205	LEU	2.8
1	E	148	TYR	2.8
1	D	156	GLN	2.7
1	E	292	LEU	2.7
1	D	305	SER	2.7
1	F	108	PRO	2.7
1	C	200	ASP	2.7
1	C	82	LEU	2.7
1	E	98	LEU	2.6
1	C	101	TYR	2.6
1	C	170	VAL	2.6
1	C	210	LEU	2.6
1	F	243	VAL	2.5
1	C	212	ALA	2.5
1	E	101	TYR	2.5
1	C	75	LEU	2.5
1	F	4	LEU	2.5
1	C	228	SER	2.5
1	C	294	GLU	2.5
1	F	119	LEU	2.5
1	C	190	GLY	2.5
1	E	199	ALA	2.4
1	D	287	SER	2.4
1	B	152	GLU	2.4
1	C	202	ASN	2.4
1	E	200	ASP	2.4
1	C	53	ASP	2.4
1	F	260	ALA	2.4
1	C	172	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	213	LEU	2.4
1	F	78	VAL	2.4
1	F	35	GLY	2.3
1	E	281	LEU	2.3
1	F	203	PRO	2.3
1	A	94	GLY	2.3
1	E	197	ILE	2.3
1	F	176	PRO	2.3
1	E	154	ALA	2.3
1	C	305	SER	2.3
1	D	190	GLY	2.3
1	E	218	LEU	2.3
1	F	45	ASP	2.3
1	C	227	GLU	2.2
1	E	173	LEU	2.2
1	C	40	TYR	2.2
1	F	113	ALA	2.2
1	D	44	TRP	2.2
1	F	276	LEU	2.2
1	F	105	VAL	2.2
1	F	170	VAL	2.2
1	F	177	SER	2.2
1	C	288	HIS	2.2
1	C	287	SER	2.2
1	E	231	ALA	2.2
1	C	91	GLY	2.2
1	E	285	VAL	2.1
1	D	89	PHE	2.1
1	C	194	ASP	2.1
1	C	186	TRP	2.1
1	E	70	VAL	2.1
1	B	92	TYR	2.1
1	E	198	TYR	2.1
1	F	173	LEU	2.1
1	C	71	SER	2.1
1	F	38	GLY	2.1
1	C	14	PHE	2.1
1	C	95	PHE	2.1
1	C	171	ASP	2.1
1	B	158	ALA	2.1
1	D	4	LEU	2.1
1	E	141	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	102	ALA	2.1
1	D	144	GLU	2.1
1	E	251	PRO	2.1
1	B	2	SER	2.1
1	D	71	SER	2.1
1	C	310	VAL	2.1
1	C	292	LEU	2.1
1	B	149	ALA	2.1
1	F	293	ALA	2.1
1	C	138	PRO	2.1
1	F	122	PHE	2.1
1	D	258	SER	2.0
1	E	71	SER	2.0
1	E	276	LEU	2.0
1	C	85	ILE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

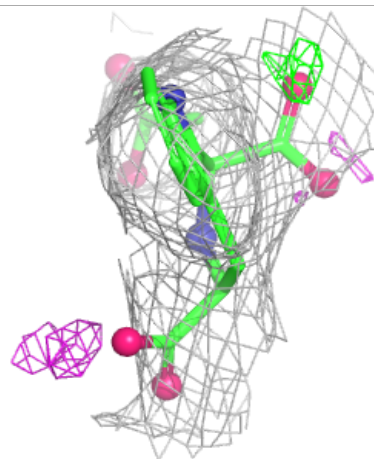
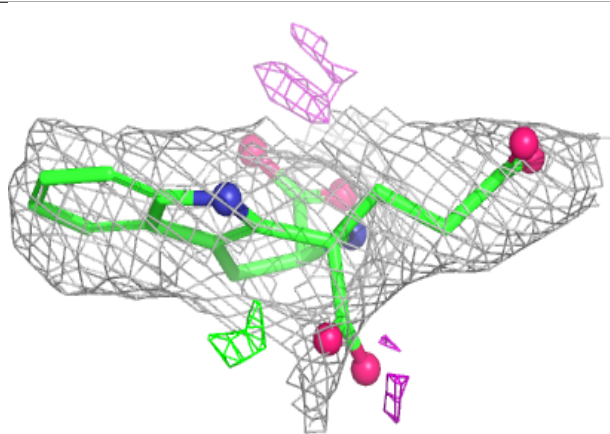
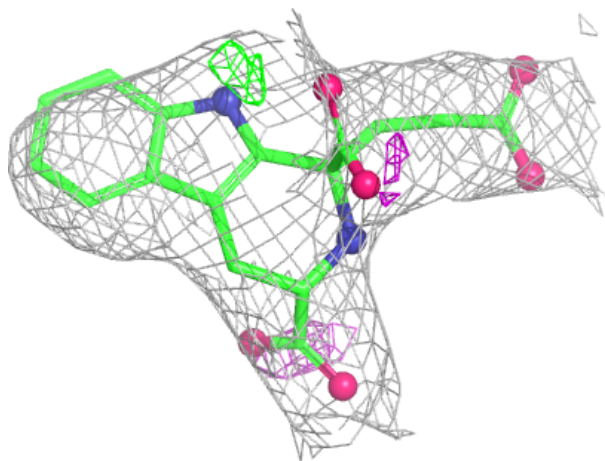
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	A1L46	D	401	24/24	0.75	0.20	55,67,72,73	0
2	A1L46	C	401	24/24	0.83	0.18	68,77,80,83	0
2	A1L46	A	401	24/24	0.83	0.14	33,37,47,50	0
2	A1L46	B	401	24/24	0.89	0.12	35,42,49,51	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.

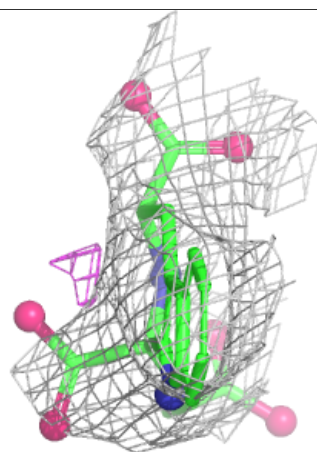
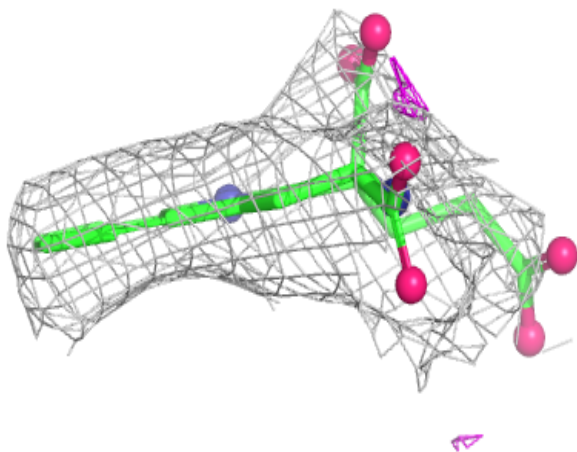
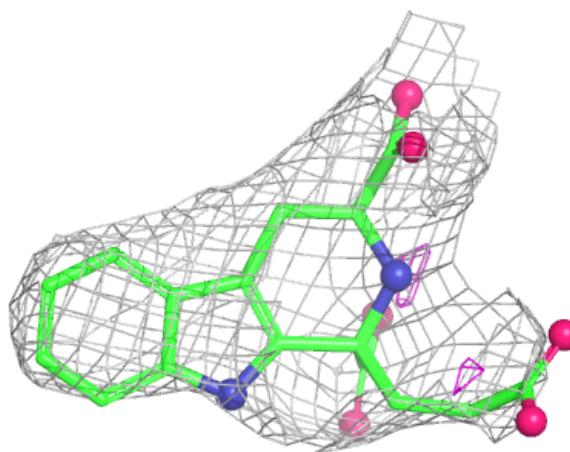
Electron density around A1L46 D 401:

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



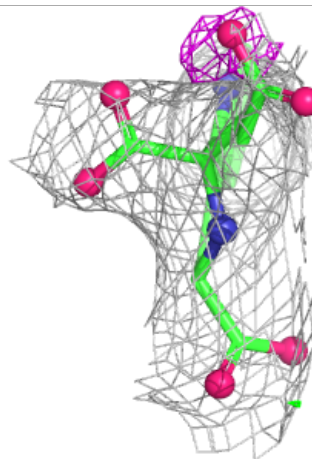
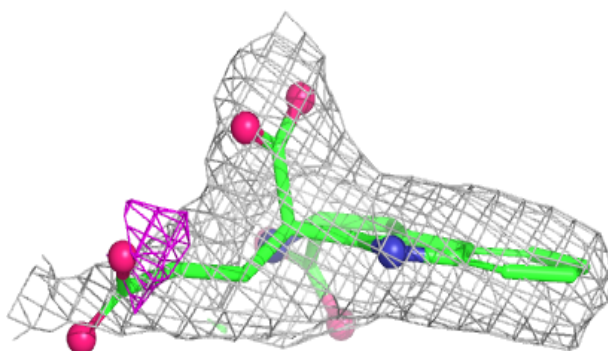
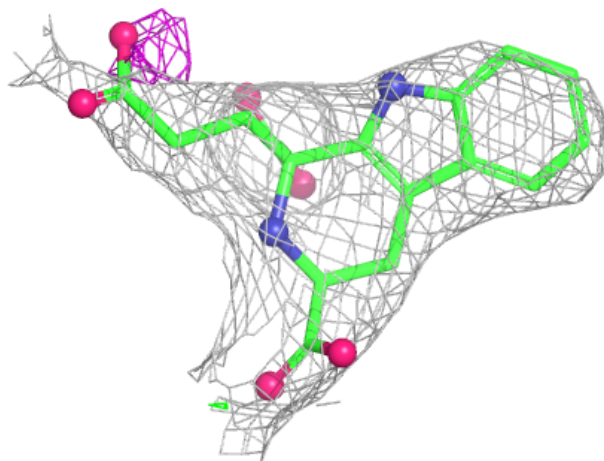
Electron density around A1L46 C 401:

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



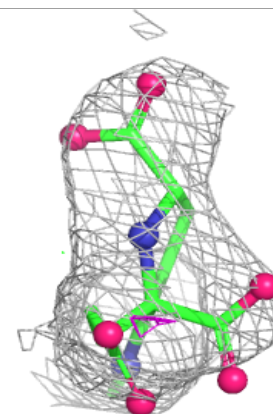
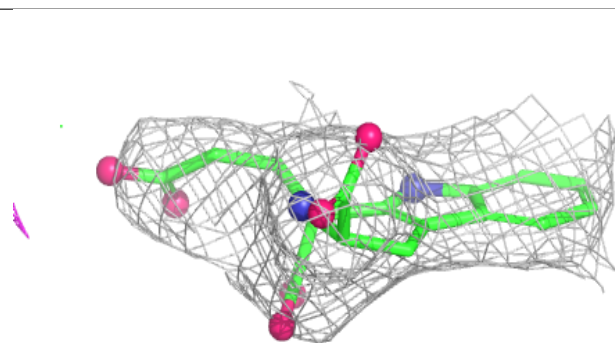
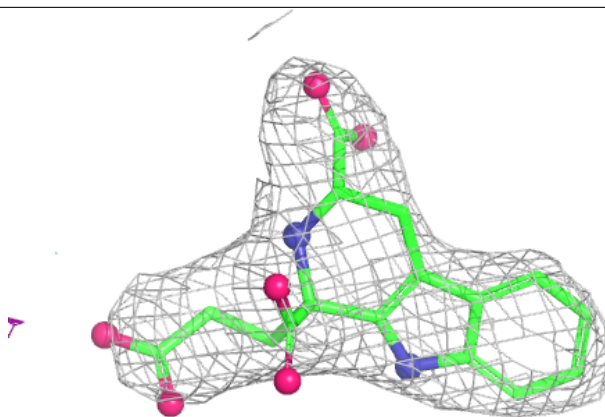
Electron density around A1L46 A 401:

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



Electron density around A1L46 B 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.