



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 5, 2023 – 08:21 PM EDT

PDB ID : 8TGY
Title : Crystal structure of Gdx-Clo from Small Multidrug Resistance family of transporters in complex with guanylyurea
Authors : Lucero, R.M.; Stockbridge, R.B.
Deposited on : 2023-07-13
Resolution : 2.18 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

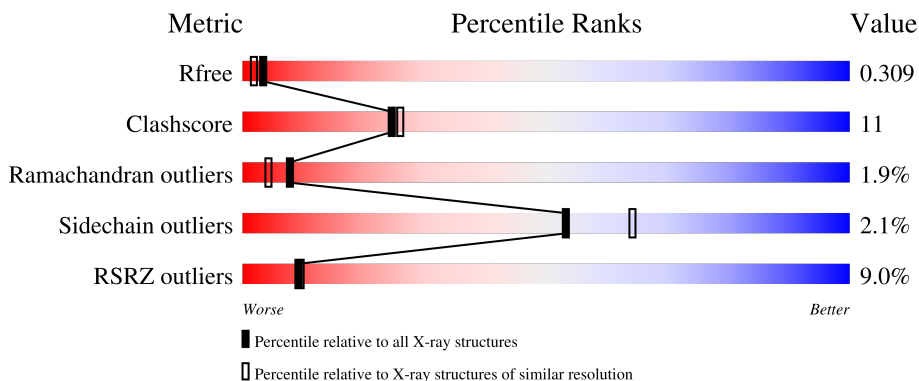
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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.



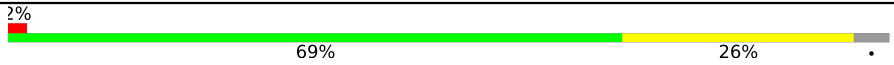


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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	105	 10% 71% 27% ..
1	B	105	 10% 78% 20% ..
1	E	105	 7% 66% 31% ..
1	G	105	 9% 70% 28% ..
2	C	91	 10% 66% 32% .

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Mol	Chain	Length	Quality of chain
2	D	91	 2% 69% 26% .
2	F	91	 13% 67% 26% . .
2	H	91	 11% 70% 20% . . 8%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6058 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.

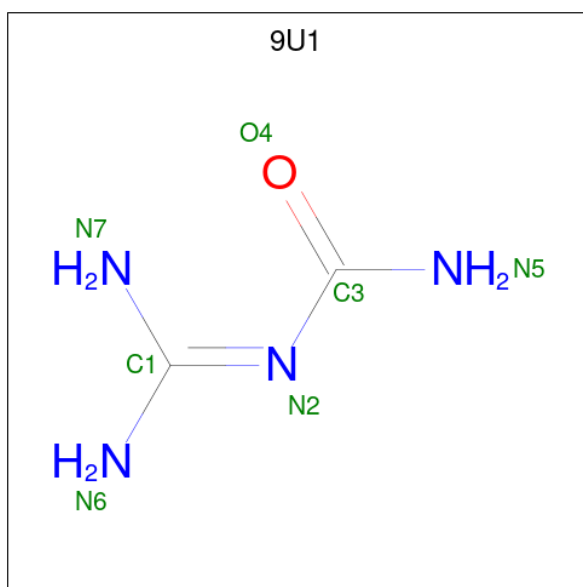
- Molecule 1 is a protein called Multidrug resistance protein, SMR family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	104	792	541	119	127	5	0	0	0
1	E	104	792	541	119	127	5	0	0	0
1	G	104	792	541	119	127	5	0	0	0
1	B	104	792	541	119	127	5	0	0	0

- Molecule 2 is a protein called L10 Monobody.

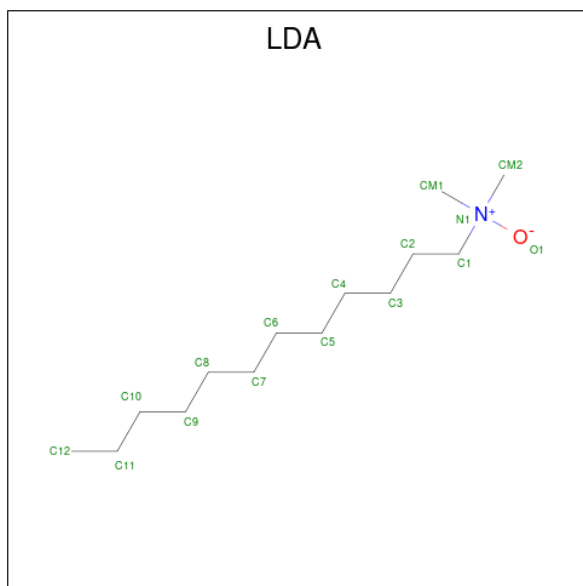
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	89	691	448	106	137	0	0	0
2	D	87	675	437	104	134	0	0	0
2	F	87	677	441	103	133	0	0	0
2	H	84	662	432	100	130	0	0	0

- Molecule 3 is N-(diaminomethylidene)urea (three-letter code: 9U1) (formula: C₂H₆N₄O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	E	1	7	2	4	1	0	0

- Molecule 4 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: $C_{14}H_{31}NO$).



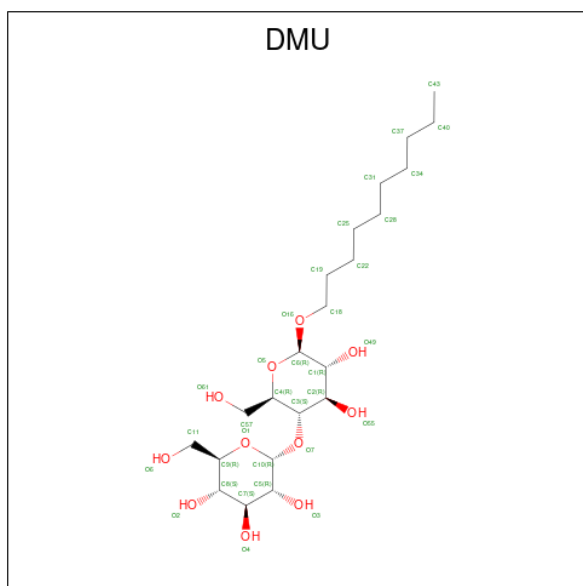
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	E	1	16	14	1	1	0	0
4	B	1	16	14	1	1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			16	14	1	1		
4	C	1	Total	C	N	O	0	0
			13	11	1	1		
4	C	1	Total	C	N	O	0	0
			16	14	1	1		
4	D	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 5 is DECYL-BETA-D-MALTOPYRANOSIDE (three-letter code: DMU) (formula: C₂₂H₄₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			33	22	11		
5	D	1	Total	C	O	0	0
			25	14	11		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	4	Total	O	0	0
			4	4		
6	E	3	Total	O	0	0
			3	3		
6	G	3	Total	O	0	0
			3	3		

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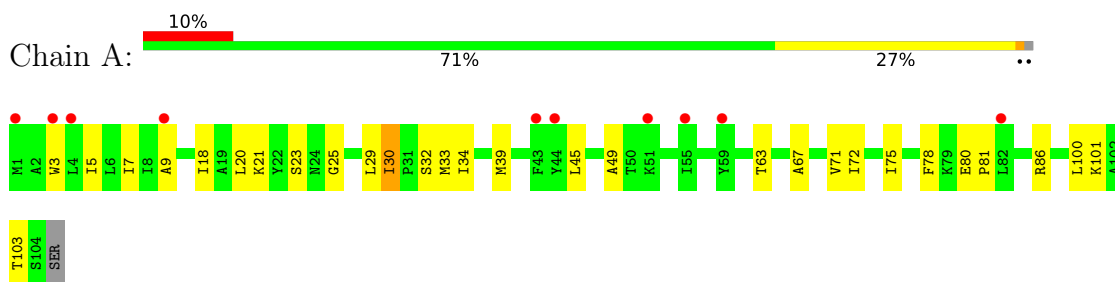
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total O 1 1	0	0
6	C	7	Total O 7 7	0	0
6	D	6	Total O 6 6	0	0
6	F	1	Total O 1 1	0	0
6	H	2	Total O 2 2	0	0

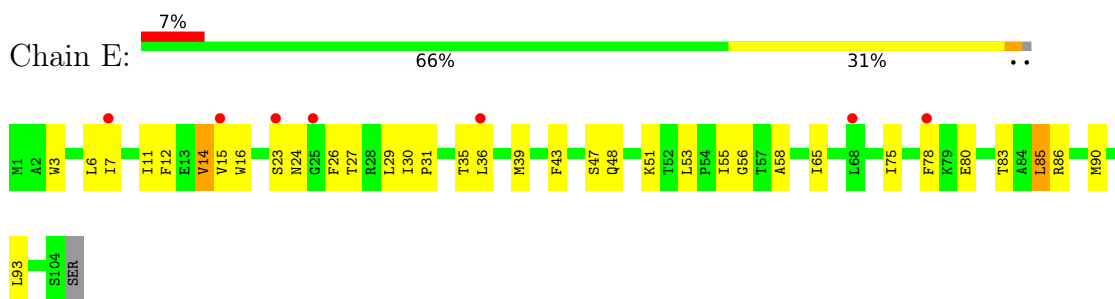
3 Residue-property plots [i](#)

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.

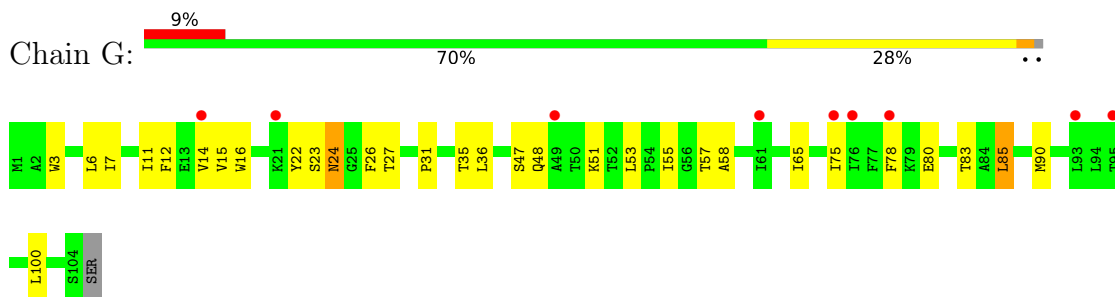
- Molecule 1: Multidrug resistance protein, SMR family



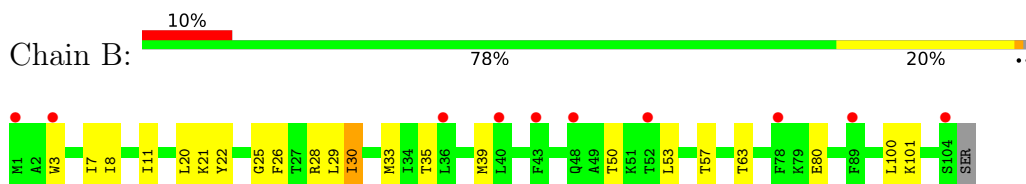
- Molecule 1: Multidrug resistance protein, SMR family



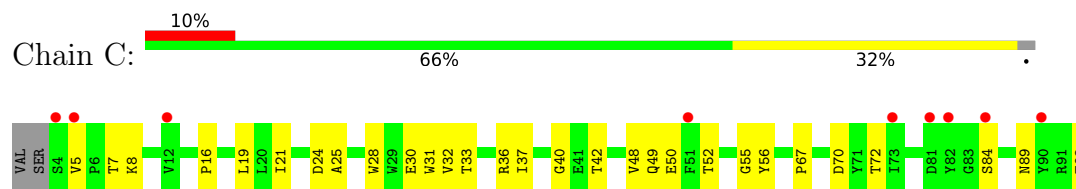
- Molecule 1: Multidrug resistance protein, SMR family



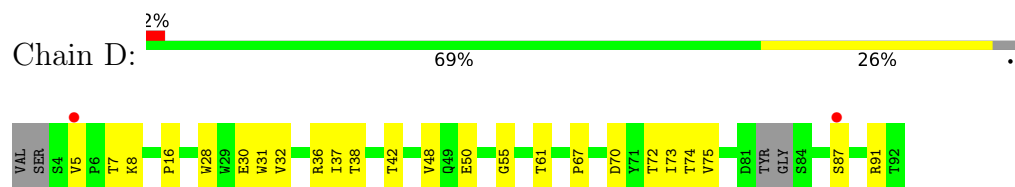
- Molecule 1: Multidrug resistance protein, SMR family



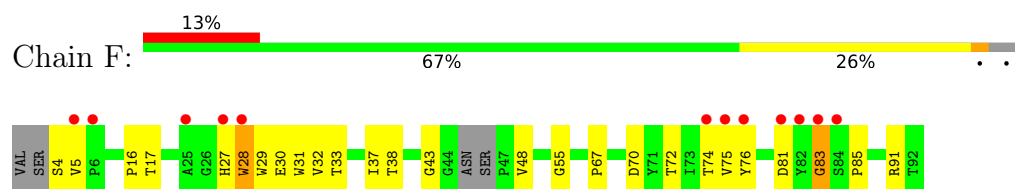
- Molecule 2: L10 Monobody



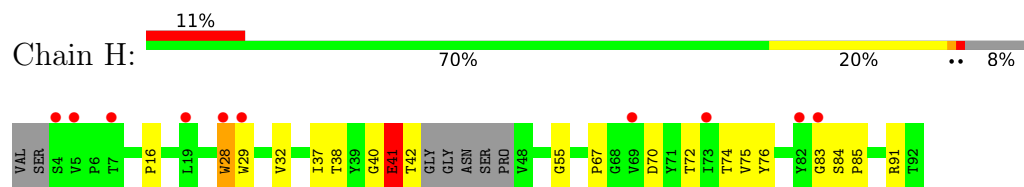
- Molecule 2: L10 Monobody



- Molecule 2: L10 Monobody



- Molecule 2: L10 Monobody



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	49.91Å 74.37Å 107.21Å 86.77° 90.01° 70.26°	Depositor
Resolution (Å)	35.66 – 2.18 35.66 – 2.18	Depositor EDS
% Data completeness (in resolution range)	81.1 (35.66-2.18) 38.8 (35.66-2.18)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.18Å)	Xtrriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.280 , 0.309 0.280 , 0.309	Depositor DCC
R_{free} test set	1360 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtrriage
Anisotropy	0.058	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 25.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtrriage
Estimated twinning fraction	0.327 for h,h-k,-l	Xtrriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	6058	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.06% 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: LDA, 9U1, DMU

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	0.24	0/809	0.39	0/1101
1	B	0.24	0/809	0.40	0/1101
1	E	0.25	0/809	0.39	0/1101
1	G	0.24	0/809	0.39	0/1101
2	C	0.29	0/715	0.49	0/988
2	D	0.25	0/697	0.46	0/962
2	F	0.25	0/700	0.45	0/965
2	H	0.24	0/684	0.47	0/944
All	All	0.25	0/6032	0.43	0/8263

There are no bond length outliers.

There are no bond angle outliers.

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	792	0	873	21	0
1	B	792	0	873	16	0
1	E	792	0	873	26	0
1	G	792	0	873	22	0
2	C	691	0	656	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	675	0	643	16	0
2	F	677	0	645	16	0
2	H	662	0	631	11	0
3	E	7	0	0	0	0
4	B	32	0	62	2	0
4	C	29	0	53	5	0
4	D	16	0	31	2	0
4	E	16	0	31	0	0
5	C	33	0	42	4	0
5	D	25	0	23	1	0
6	A	4	0	0	0	0
6	B	1	0	0	0	0
6	C	7	0	0	3	0
6	D	6	0	0	1	0
6	E	3	0	0	0	0
6	F	1	0	0	0	0
6	G	3	0	0	0	0
6	H	2	0	0	0	0
All	All	6058	0	6309	135	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 11.

All (135) 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)
2:C:52:THR:HG21	5:C:501:DMU:H6	1.61	0.82
2:D:7:THR:HG22	2:D:8:LYS:HG3	1.65	0.79
2:C:7:THR:HG22	2:C:8:LYS:HG3	1.66	0.75
2:C:25:ALA:O	6:C:601:HOH:O	2.04	0.75
1:G:7:ILE:HA	1:G:65:ILE:HD13	1.71	0.72
2:C:84:SER:O	6:C:602:HOH:O	2.08	0.71
2:H:38:THR:OG1	2:H:74:THR:OG1	2.11	0.67
1:B:33:MET:HE1	4:D:102:LDA:H61	1.76	0.65
2:D:5:VAL:HG11	2:D:30:GLU:HB3	1.81	0.61
1:E:23:SER:HB2	1:E:31:PRO:HB2	1.82	0.61
1:G:83:THR:HG22	1:G:85:LEU:H	1.64	0.61
5:C:501:DMU:H21	4:C:502:LDA:H62	1.80	0.61
5:C:501:DMU:O55	6:C:603:HOH:O	2.16	0.61
2:C:7:THR:HB	2:C:24:ASP:HB3	1.82	0.60
1:E:39:MET:HE3	1:E:43:PHE:HE2	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:70:ASP:HA	2:F:91:ARG:HA	1.84	0.60
1:E:83:THR:HB	1:E:86:ARG:HG3	1.83	0.59
1:E:83:THR:HG22	1:E:85:LEU:H	1.67	0.59
1:A:30:ILE:H	1:A:33:MET:HG3	1.67	0.58
2:C:33:THR:HB	5:C:501:DMU:H20	1.84	0.58
2:D:42:THR:HB	2:D:72:THR:HG23	1.86	0.57
1:E:48:GLN:OE1	1:E:51:LYS:NZ	2.38	0.57
1:G:48:GLN:OE1	1:G:51:LYS:NZ	2.38	0.56
2:C:40:GLY:HA3	2:C:48:VAL:HG12	1.87	0.56
2:C:42:THR:HB	2:C:72:THR:HG23	1.87	0.56
2:D:37:ILE:HG12	2:D:75:VAL:HG22	1.88	0.56
2:H:16:PRO:HB3	2:H:67:PRO:HB3	1.85	0.56
1:G:23:SER:HB2	1:G:31:PRO:HB2	1.87	0.56
1:A:80:GLU:OE2	1:E:55:ILE:HG22	2.05	0.55
1:G:55:ILE:HG22	1:B:80:GLU:OE2	2.07	0.55
1:E:27:THR:O	2:F:33:THR:N	2.39	0.55
2:D:16:PRO:HB3	2:D:67:PRO:HB3	1.88	0.55
1:A:20:LEU:HD11	1:E:47:SER:HB3	1.89	0.54
1:A:21:LYS:NZ	1:E:53:LEU:O	2.36	0.54
2:H:37:ILE:HG13	2:H:75:VAL:HG22	1.90	0.54
2:C:16:PRO:HB3	2:C:67:PRO:HB3	1.90	0.53
2:F:17:THR:HG22	2:F:67:PRO:HD3	1.91	0.53
1:A:81:PRO:O	1:A:86:ARG:NH1	2.41	0.53
1:G:6:LEU:HD11	1:G:58:ALA:HA	1.91	0.53
2:C:7:THR:O	2:C:24:ASP:N	2.25	0.52
1:G:3:TRP:O	1:G:7:ILE:HG12	2.10	0.52
1:A:29:LEU:HB2	2:C:32:VAL:O	2.09	0.52
2:F:37:ILE:HG13	2:F:75:VAL:HG22	1.91	0.52
2:H:41:GLU:O	2:H:72:THR:HG22	2.10	0.52
1:E:75:ILE:HA	1:E:80:GLU:HB3	1.92	0.52
2:F:28:TRP:HB3	2:F:29:TRP:CE3	2.45	0.51
2:C:56:TYR:HB2	4:C:503:LDA:H41	1.93	0.51
1:G:11:ILE:O	1:G:14:VAL:HG12	2.11	0.51
2:C:5:VAL:HG11	2:C:30:GLU:HB3	1.91	0.51
1:B:63:THR:HG21	1:B:101:LYS:NZ	2.25	0.51
2:D:70:ASP:HA	2:D:91:ARG:HA	1.93	0.51
2:F:38:THR:OG1	2:F:74:THR:OG1	2.22	0.50
1:B:8:ILE:HA	1:B:11:ILE:HD12	1.93	0.50
2:F:76:TYR:CD1	2:F:85:PRO:HB3	2.46	0.50
1:E:3:TRP:O	1:E:7:ILE:HG12	2.12	0.50
1:G:75:ILE:HA	1:G:80:GLU:HB3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:503:LDA:H81	4:C:503:LDA:H122	1.93	0.49
1:A:9:ALA:HB2	1:A:45:LEU:HB2	1.94	0.49
1:B:35:THR:O	1:B:39:MET:HG2	2.13	0.49
1:E:7:ILE:HA	1:E:65:ILE:HD13	1.95	0.49
2:H:70:ASP:HA	2:H:91:ARG:HA	1.94	0.49
2:D:42:THR:HG21	2:D:70:ASP:OD2	2.13	0.49
2:D:50:GLU:OE2	5:D:101:DMU:O3	2.30	0.48
2:C:42:THR:HG21	2:C:70:ASP:OD2	2.14	0.48
2:H:28:TRP:HE3	2:H:29:TRP:HB3	1.78	0.48
1:A:25:GLY:HA2	1:E:47:SER:HB2	1.96	0.48
2:C:32:VAL:O	2:C:55:GLY:HA3	2.13	0.47
1:G:57:THR:HG23	1:G:100:LEU:HD13	1.95	0.47
2:D:74:THR:HG23	2:D:87:SER:HB3	1.96	0.47
2:F:76:TYR:HD1	2:F:85:PRO:HB3	1.78	0.47
2:C:28:TRP:HA	2:C:31:TRP:CD2	2.49	0.47
1:E:31:PRO:O	1:E:35:THR:HG23	2.14	0.47
1:G:31:PRO:O	1:G:35:THR:HG23	2.15	0.47
1:A:23:SER:HB2	1:A:32:SER:OG	2.15	0.47
1:E:6:LEU:HD11	1:E:58:ALA:HA	1.97	0.47
1:B:3:TRP:CE2	1:B:53:LEU:HD22	2.50	0.47
1:E:30:ILE:HD11	2:F:31:TRP:HE3	1.80	0.46
1:G:16:TRP:HB2	1:G:35:THR:HA	1.96	0.46
1:G:47:SER:HB2	1:B:25:GLY:HA2	1.98	0.46
2:C:21:ILE:HD11	2:C:37:ILE:HD13	1.97	0.46
2:F:5:VAL:HG21	2:F:30:GLU:HB3	1.98	0.45
1:B:26:PHE:HB2	4:B:601:LDA:H51	1.99	0.45
2:H:32:VAL:O	2:H:55:GLY:HA3	2.16	0.45
1:A:30:ILE:O	1:A:34:ILE:HG12	2.16	0.45
2:D:28:TRP:HA	2:D:31:TRP:CD2	2.52	0.45
2:C:19:LEU:HD21	2:C:92:THR:HG22	1.99	0.45
1:G:24:ASN:HB2	1:G:27:THR:OG1	2.17	0.45
1:E:12:PHE:HA	1:E:15:VAL:HG22	1.99	0.45
1:B:29:LEU:O	1:B:30:ILE:HB	2.16	0.44
1:G:11:ILE:O	1:G:15:VAL:HG13	2.17	0.44
1:G:36:LEU:HD12	1:G:36:LEU:HA	1.84	0.44
1:G:12:PHE:HA	1:G:15:VAL:HG22	1.99	0.44
2:D:38:THR:HG23	2:D:48:VAL:HB	1.99	0.44
1:A:3:TRP:O	1:A:7:ILE:HG12	2.17	0.44
1:B:57:THR:OG1	4:B:602:LDA:H121	2.18	0.44
2:H:41:GLU:HB3	2:H:42:THR:H	1.62	0.44
1:A:29:LEU:O	1:A:30:ILE:HB	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:47:SER:HB3	1:B:20:LEU:HD11	2.00	0.44
1:B:22:TYR:O	1:B:28:ARG:NH2	2.51	0.44
2:H:28:TRP:HB3	2:H:29:TRP:HE3	1.83	0.43
1:A:71:VAL:HA	1:E:56:GLY:HA3	2.01	0.43
2:D:42:THR:HG22	2:D:70:ASP:O	2.19	0.43
1:G:53:LEU:O	1:B:21:LYS:NZ	2.51	0.43
1:G:90:MET:HG3	1:B:100:LEU:HD21	2.00	0.43
1:A:67:ALA:O	1:A:71:VAL:HG23	2.18	0.43
1:A:18:ILE:HD12	1:A:78:PHE:HE2	1.84	0.43
1:A:100:LEU:HD21	1:E:90:MET:HG3	2.01	0.43
1:G:22:TYR:HD1	1:G:22:TYR:HA	1.73	0.43
2:H:76:TYR:HD1	2:H:85:PRO:HB3	1.83	0.43
1:E:29:LEU:HB2	2:F:32:VAL:O	2.19	0.43
1:A:5:ILE:HG13	1:A:49:ALA:HB2	2.00	0.42
2:C:36:ARG:HD2	2:C:50:GLU:OE2	2.19	0.42
1:E:30:ILE:HD11	2:F:31:TRP:CE3	2.55	0.42
1:A:72:ILE:O	1:A:75:ILE:HG13	2.19	0.42
1:E:26:PHE:HZ	1:E:36:LEU:HD13	1.84	0.42
2:H:70:ASP:HB2	2:H:91:ARG:NH2	2.34	0.42
1:A:63:THR:HG21	1:A:101:LYS:NZ	2.34	0.42
2:F:27:HIS:HA	2:F:31:TRP:NE1	2.34	0.42
2:F:32:VAL:O	2:F:55:GLY:HA3	2.20	0.42
1:B:7:ILE:O	1:B:11:ILE:HG13	2.20	0.41
2:D:36:ARG:NH1	6:D:205:HOH:O	2.53	0.41
1:E:30:ILE:HB	1:E:31:PRO:HD3	2.02	0.41
1:E:11:ILE:O	1:E:14:VAL:HG13	2.20	0.41
1:E:16:TRP:HB2	1:E:35:THR:HA	2.03	0.41
2:F:16:PRO:HB2	2:F:67:PRO:HB3	2.02	0.41
1:G:24:ASN:O	1:G:26:PHE:N	2.48	0.41
2:D:31:TRP:CE2	4:D:102:LDA:HM21	2.55	0.41
2:C:31:TRP:CZ3	4:C:503:LDA:H21	2.55	0.41
2:D:38:THR:O	2:D:73:ILE:HA	2.20	0.41
1:A:103:THR:OG1	1:E:86:ARG:NH1	2.53	0.41
4:C:503:LDA:H22	4:C:503:LDA:HM11	1.79	0.41
2:F:4:SER:N	2:F:83:GLY:HA3	2.36	0.41
1:B:3:TRP:CE3	1:B:53:LEU:HD13	2.56	0.40
2:D:32:VAL:O	2:D:55:GLY:HA3	2.21	0.40
1:A:29:LEU:HB2	2:C:32:VAL:C	2.42	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	102/105 (97%)	100 (98%)	1 (1%)	1 (1%)	15	12
1	B	102/105 (97%)	99 (97%)	2 (2%)	1 (1%)	15	12
1	E	102/105 (97%)	99 (97%)	1 (1%)	2 (2%)	7	4
1	G	102/105 (97%)	95 (93%)	5 (5%)	2 (2%)	7	4
2	C	87/91 (96%)	83 (95%)	4 (5%)	0	100	100
2	D	83/91 (91%)	80 (96%)	3 (4%)	0	100	100
2	F	83/91 (91%)	76 (92%)	4 (5%)	3 (4%)	3	1
2	H	80/91 (88%)	70 (88%)	5 (6%)	5 (6%)	1	0
All	All	741/784 (94%)	702 (95%)	25 (3%)	14 (2%)	8	4

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	ILE
1	B	30	ILE
1	E	24	ASN
2	F	83	GLY
2	H	28	TRP
2	H	41	GLU
2	H	40	GLY
1	G	24	ASN
2	F	28	TRP
1	E	78	PHE
1	G	78	PHE
2	F	43	GLY
2	H	83	GLY
2	H	84	SER

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	83/84 (99%)	82 (99%)	1 (1%)	71	81
1	B	83/84 (99%)	82 (99%)	1 (1%)	71	81
1	E	83/84 (99%)	80 (96%)	3 (4%)	35	42
1	G	83/84 (99%)	82 (99%)	1 (1%)	71	81
2	C	76/78 (97%)	74 (97%)	2 (3%)	46	55
2	D	75/78 (96%)	74 (99%)	1 (1%)	69	79
2	F	74/78 (95%)	71 (96%)	3 (4%)	30	36
2	H	73/78 (94%)	72 (99%)	1 (1%)	67	78
All	All	630/648 (97%)	617 (98%)	13 (2%)	53	64

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	MET
1	E	14	VAL
1	E	85	LEU
1	E	93	LEU
1	G	85	LEU
1	B	50	THR
2	C	49	GLN
2	C	89	ASN
2	D	61	THR
2	F	48	VAL
2	F	72	THR
2	F	81	ASP
2	H	41	GLU

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

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 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
5	DMU	D	101	-	26,26,34	1.62	8 (30%)	37,37,45	1.42	6 (16%)
4	LDA	C	502	-	9,12,15	0.38	0	11,14,17	0.21	0
4	LDA	B	602	-	12,15,15	0.34	0	14,17,17	0.22	0
5	DMU	C	501	-	34,34,34	1.41	6 (17%)	45,45,45	1.69	8 (17%)
3	9U1	E	201	-	6,6,6	0.22	0	6,7,7	0.25	0
4	LDA	B	601	-	12,15,15	0.34	0	14,17,17	0.19	0
4	LDA	C	503	-	12,15,15	0.35	0	14,17,17	0.23	0
4	LDA	D	102	-	12,15,15	0.38	0	14,17,17	0.69	1 (7%)
4	LDA	E	202	-	12,15,15	0.36	0	14,17,17	0.34	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	DMU	D	101	-	-	7/11/51/59	0/2/2/2
4	LDA	C	502	-	-	2/10/10/13	-
4	LDA	B	602	-	-	6/13/13/13	-
5	DMU	C	501	-	-	10/19/59/59	0/2/2/2
3	9U1	E	201	-	-	1/3/4/4	-
4	LDA	B	601	-	-	7/13/13/13	-
4	LDA	C	503	-	-	8/13/13/13	-
4	LDA	D	102	-	-	3/13/13/13	-
4	LDA	E	202	-	-	3/13/13/13	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	101	DMU	O1-C9	3.50	1.52	1.44
5	C	501	DMU	O1-C9	3.43	1.52	1.44
5	C	501	DMU	C11-C9	-3.05	1.41	1.51
5	D	101	DMU	C11-C9	-3.03	1.41	1.51
5	D	101	DMU	O4-C7	2.82	1.49	1.43
5	C	501	DMU	O4-C7	2.76	1.49	1.43
5	C	501	DMU	C7-C5	-2.34	1.46	1.52
5	D	101	DMU	C7-C5	-2.33	1.46	1.52
5	D	101	DMU	O3-C5	2.28	1.48	1.43
5	C	501	DMU	O3-C5	2.27	1.48	1.43
5	C	501	DMU	O1-C10	2.23	1.47	1.41
5	D	101	DMU	O5-C6	2.20	1.47	1.41
5	D	101	DMU	O1-C10	2.08	1.47	1.41
5	D	101	DMU	O5-C4	2.05	1.49	1.44

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	501	DMU	C10-C5-C7	4.29	118.93	110.00
5	C	501	DMU	C1-C2-C3	4.17	119.19	109.68
5	C	501	DMU	C6-C1-C2	4.11	118.55	110.00
5	D	101	DMU	C8-C7-C5	3.69	117.26	110.82
5	C	501	DMU	O1-C10-C5	3.63	118.02	110.35
5	C	501	DMU	C10-O7-C3	-3.56	109.14	117.96
5	D	101	DMU	C10-C5-C7	3.48	117.25	110.00
5	C	501	DMU	O1-C9-C8	-3.11	104.05	109.69
5	C	501	DMU	C8-C7-C5	2.86	115.81	110.82
5	D	101	DMU	C10-O7-C3	-2.73	111.21	117.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	101	DMU	O1-C10-C5	2.52	115.69	110.35
5	D	101	DMU	O1-C9-C8	-2.32	105.49	109.69
5	D	101	DMU	C7-C8-C9	2.28	114.31	110.24
5	C	501	DMU	O5-C6-C1	2.13	114.85	110.35
4	D	102	LDA	CM1-N1-C1	2.09	114.64	110.23

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	201	9U1	O4-C3-N2-C1
4	E	202	LDA	N1-C1-C2-C3
4	B	601	LDA	N1-C1-C2-C3
4	B	602	LDA	N1-C1-C2-C3
5	D	101	DMU	C1-C6-O16-C18
5	D	101	DMU	O5-C6-O16-C18
5	D	101	DMU	O6-C11-C9-O1
5	C	501	DMU	O6-C11-C9-O1
5	D	101	DMU	O5-C4-C57-O61
5	C	501	DMU	O6-C11-C9-C8
5	D	101	DMU	O6-C11-C9-C8
5	C	501	DMU	O5-C4-C57-O61
5	D	101	DMU	C3-C4-C57-O61
4	B	601	LDA	C5-C6-C7-C8
5	C	501	DMU	C31-C34-C37-C40
5	C	501	DMU	C19-C22-C25-C28
4	C	502	LDA	C5-C6-C7-C8
4	D	102	LDA	C2-C3-C4-C5
4	D	102	LDA	C7-C8-C9-C10
5	C	501	DMU	C18-C19-C22-C25
4	C	503	LDA	C7-C8-C9-C10
4	B	602	LDA	C11-C10-C9-C8
4	D	102	LDA	C1-C2-C3-C4
4	C	503	LDA	N1-C1-C2-C3
4	C	503	LDA	C6-C7-C8-C9
4	C	502	LDA	C6-C7-C8-C9
4	C	503	LDA	C4-C5-C6-C7
4	C	503	LDA	C9-C10-C11-C12
5	C	501	DMU	O5-C6-O16-C18
4	B	601	LDA	C2-C1-N1-CM1
4	B	601	LDA	C2-C1-N1-CM2
4	B	602	LDA	C5-C6-C7-C8

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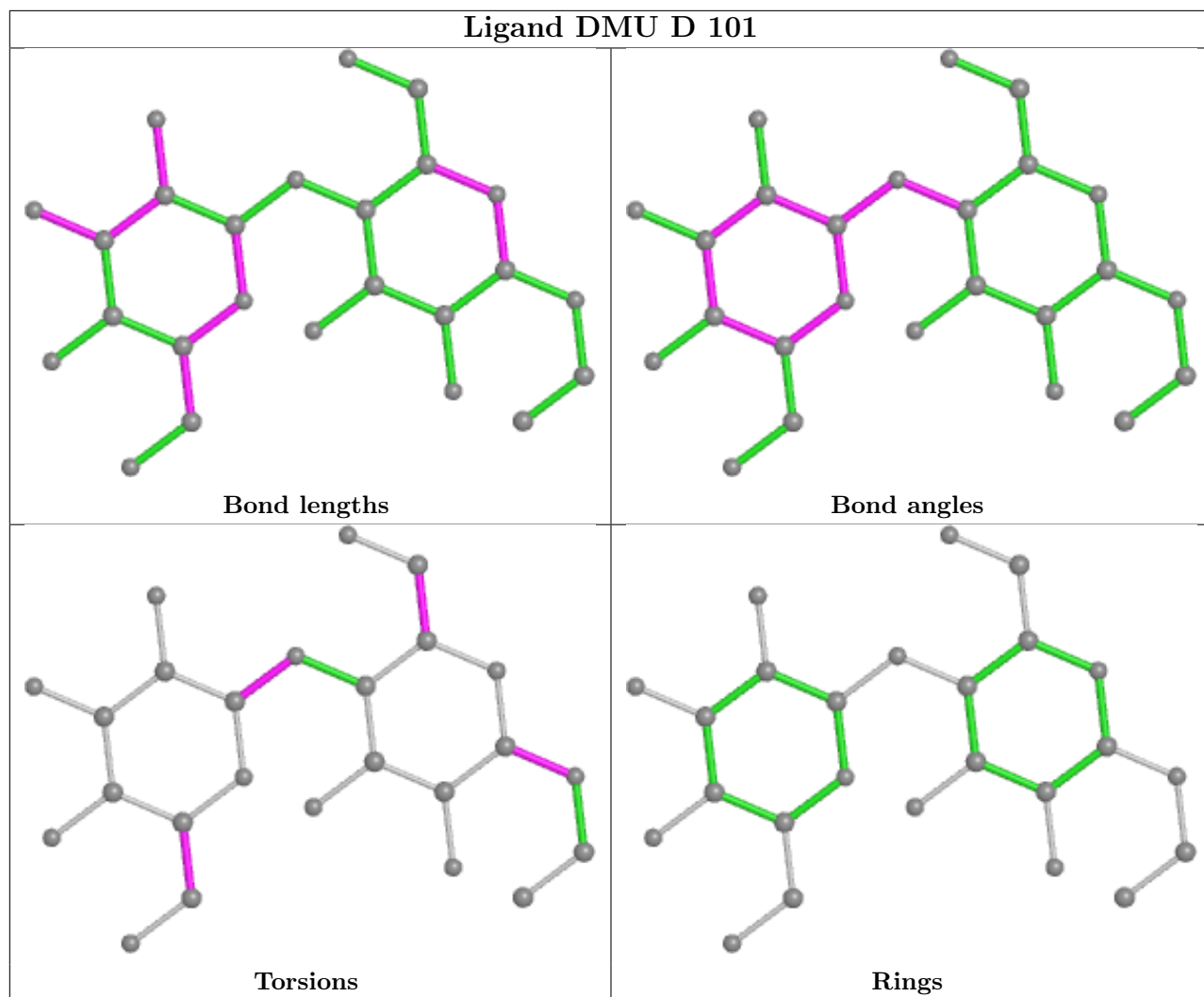
Mol	Chain	Res	Type	Atoms
4	B	601	LDA	C2-C1-N1-O1
5	C	501	DMU	C3-C4-C57-O61
4	B	601	LDA	C4-C5-C6-C7
4	E	202	LDA	C4-C5-C6-C7
4	B	602	LDA	C4-C5-C6-C7
5	C	501	DMU	O16-C18-C19-C22
4	B	601	LDA	C2-C3-C4-C5
4	B	602	LDA	C2-C3-C4-C5
4	C	503	LDA	C3-C4-C5-C6
4	B	602	LDA	C3-C4-C5-C6
5	C	501	DMU	C34-C37-C40-C43
4	C	503	LDA	C2-C3-C4-C5
5	D	101	DMU	C5-C10-O7-C3
4	E	202	LDA	C5-C6-C7-C8
4	C	503	LDA	C5-C6-C7-C8

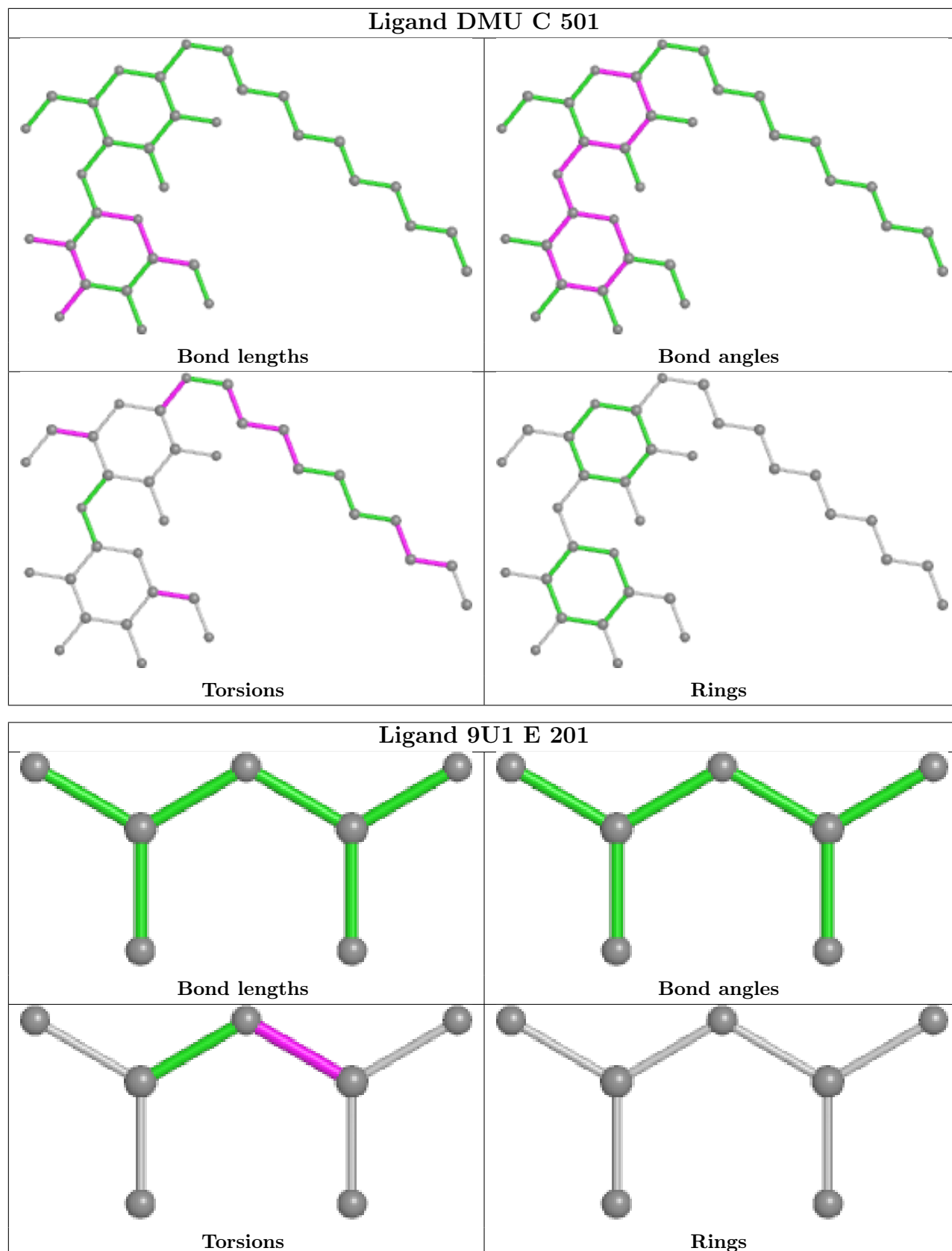
There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	101	DMU	1	0
4	C	502	LDA	1	0
4	B	602	LDA	1	0
5	C	501	DMU	4	0
4	B	601	LDA	1	0
4	C	503	LDA	4	0
4	D	102	LDA	2	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	104/105 (99%)	0.76	10 (9%) 8 8	24, 42, 69, 97	0
1	B	104/105 (99%)	0.83	10 (9%) 8 8	26, 41, 74, 86	0
1	E	104/105 (99%)	0.73	7 (6%) 17 18	24, 38, 70, 88	0
1	G	104/105 (99%)	0.75	9 (8%) 10 10	25, 41, 68, 97	0
2	C	89/91 (97%)	0.67	9 (10%) 7 7	26, 41, 69, 95	0
2	D	87/91 (95%)	0.42	2 (2%) 60 61	26, 41, 63, 71	0
2	F	87/91 (95%)	0.90	12 (13%) 2 3	29, 54, 77, 93	0
2	H	84/91 (92%)	0.88	10 (11%) 4 4	29, 52, 80, 88	0
All	All	763/784 (97%)	0.75	69 (9%) 9 9	24, 42, 72, 97	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	5	VAL	5.6
2	F	28	TRP	5.4
1	A	55	ILE	5.1
1	A	1	MET	4.7
2	H	4	SER	4.7
2	F	76	TYR	4.6
2	H	29	TRP	4.5
1	E	7	ILE	4.3
2	H	83	GLY	4.2
1	A	51	LYS	3.9
1	B	3	TRP	3.8
2	F	74	THR	3.7
2	H	73	ILE	3.7
1	E	78	PHE	3.6
2	F	75	VAL	3.3
1	B	43	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
2	C	81	ASP	3.1
2	F	27	HIS	3.1
2	H	28	TRP	3.1
2	F	82	TYR	3.0
1	B	40	LEU	3.0
2	C	5	VAL	3.0
1	A	44	TYR	2.9
1	B	104	SER	2.9
1	G	21	LYS	2.9
1	G	95	THR	2.8
1	E	36	LEU	2.7
1	B	89	PHE	2.7
2	F	25	ALA	2.7
2	F	83	GLY	2.7
1	G	76	ILE	2.7
1	B	52	THR	2.7
1	A	3	TRP	2.6
2	H	69	VAL	2.6
2	C	4	SER	2.6
1	A	9	ALA	2.6
2	C	82	TYR	2.5
2	D	5	VAL	2.5
1	G	93	LEU	2.5
1	E	15	VAL	2.4
1	G	61	ILE	2.4
1	A	82	LEU	2.4
2	F	81	ASP	2.3
1	G	75	ILE	2.3
2	F	6	PRO	2.3
2	D	87	SER	2.3
1	G	78	PHE	2.3
2	H	7	THR	2.3
2	H	82	TYR	2.3
1	A	4	LEU	2.3
1	B	78	PHE	2.3
1	G	14	VAL	2.3
1	B	36	LEU	2.2
1	B	48	GLN	2.2
1	E	23	SER	2.2
2	C	84	SER	2.2
2	F	5	VAL	2.2
2	C	73	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	1	MET	2.1
2	C	12	VAL	2.1
2	C	90	TYR	2.1
2	C	51	PHE	2.1
1	E	68	LEU	2.1
1	A	43	PHE	2.1
1	A	59	TYR	2.1
1	G	49	ALA	2.0
2	F	84	SER	2.0
1	E	25	GLY	2.0
2	H	19	LEU	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 monosaccharides in this entry.

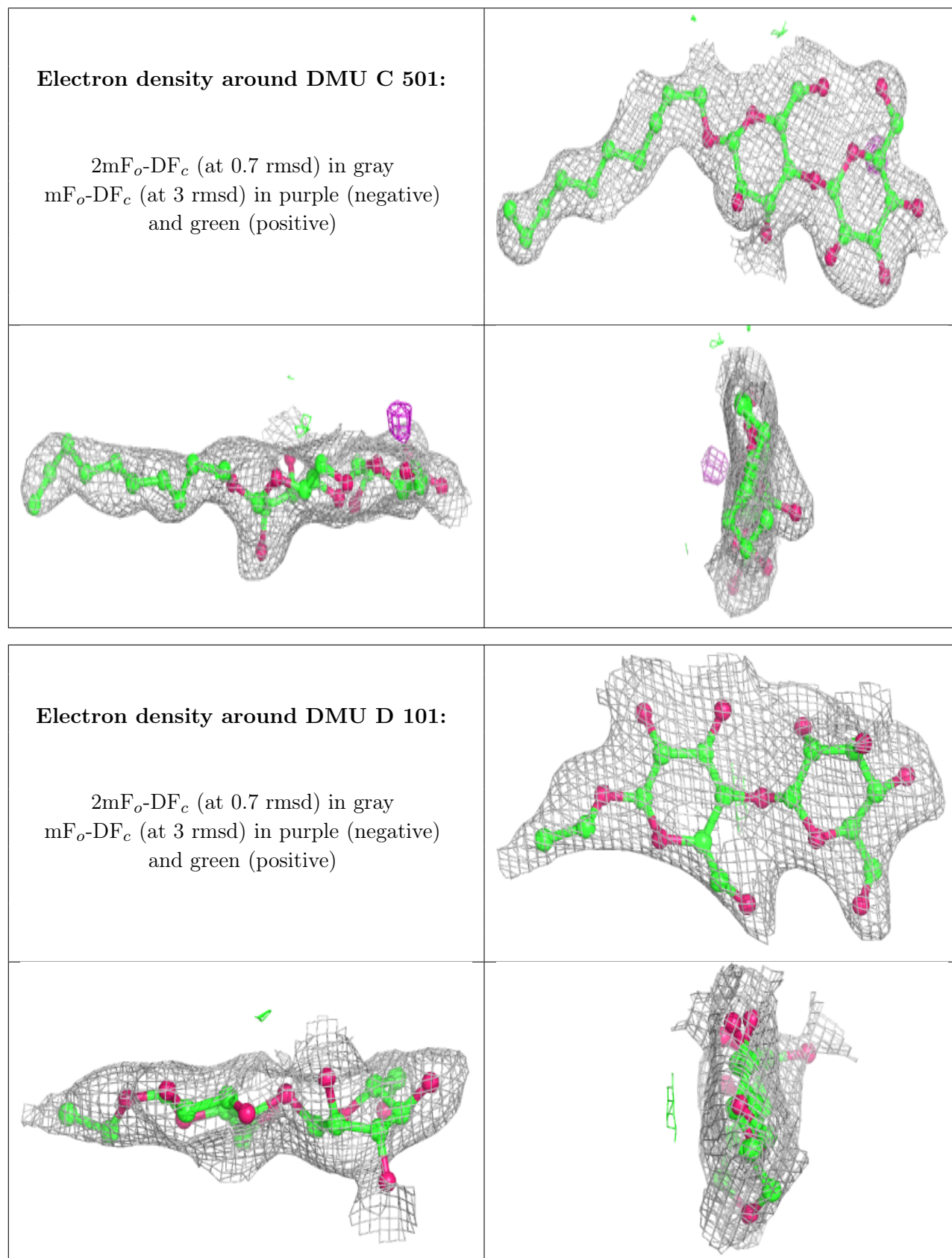
6.4 Ligands [i](#)

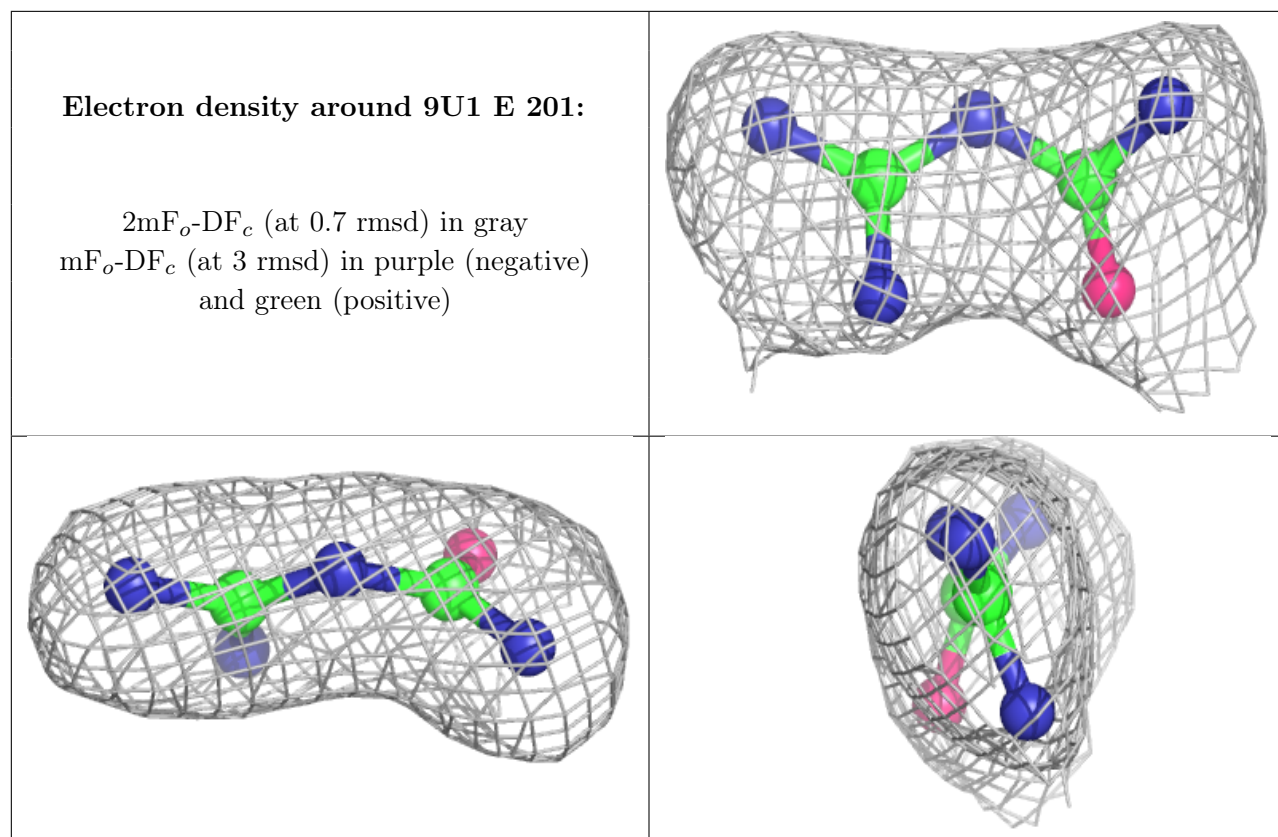
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	LDA	E	202	16/16	0.82	0.27	33,45,62,73	0
5	DMU	C	501	33/33	0.82	0.23	28,70,86,89	0
5	DMU	D	101	25/33	0.85	0.17	54,71,82,88	0
4	LDA	C	503	16/16	0.86	0.24	39,54,76,82	0
4	LDA	C	502	13/16	0.89	0.22	27,33,57,89	0
4	LDA	B	602	16/16	0.91	0.20	41,55,69,74	0
4	LDA	D	102	16/16	0.91	0.10	20,43,54,58	0
4	LDA	B	601	16/16	0.92	0.13	20,42,50,59	0
3	9U1	E	201	7/7	0.93	0.16	30,44,58,59	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.