



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 11:03 PM UTC

PDB ID : 5D56 / pdb_00005d56
Title : In meso in situ serial X-ray crystallography structure of diacylglycerol kinase, DgkA, at 100 K
Authors : Huang, C.-Y.; Howe, N.; Olieric, V.; Warshamanage, R.; Diederichs, K.; Wang, M.; Caffrey, M.
Deposited on : 2015-08-10
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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:

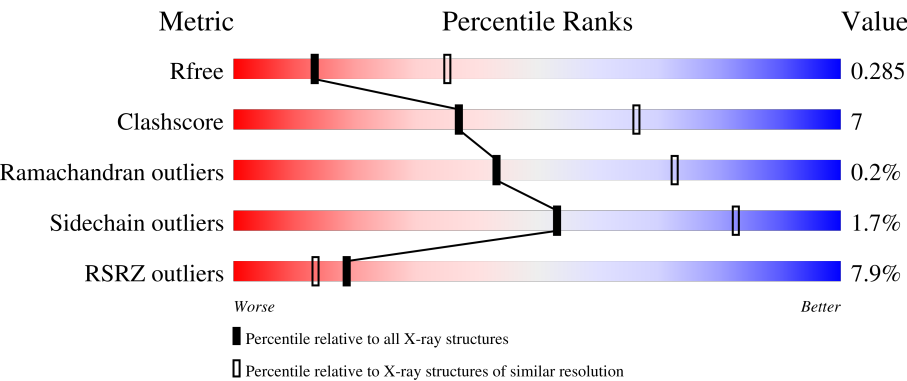
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 i

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

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	130	7% 76% 13% 11%
1	B	130	4% 62% 12% 25%
1	C	130	% 65% 31%
1	D	130	5% 72% 11% 17%

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Mol	Chain	Length	Quality of chain
1	E	130	10% 55% 12% • 32%
1	F	130	8% 54% 11% • 35%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4717 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 Diacylglycerol kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	116	Total	C	N	O	S	0	0	0
			887	580	149	155	3			
1	B	97	Total	C	N	O	S	0	1	0
			747	489	121	134	3			
1	C	90	Total	C	N	O	S	0	0	0
			681	443	112	123	3			
1	D	108	Total	C	N	O	S	0	0	0
			824	538	137	146	3			
1	E	89	Total	C	N	O	S	0	0	0
			670	437	108	122	3			
1	F	85	Total	C	N	O	S	0	0	0
			635	415	101	116	3			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	GLY	-	expression tag	UNP P0ABN1
A	-7	HIS	-	expression tag	UNP P0ABN1
A	-6	HIS	-	expression tag	UNP P0ABN1
A	-5	HIS	-	expression tag	UNP P0ABN1
A	-4	HIS	-	expression tag	UNP P0ABN1
A	-3	HIS	-	expression tag	UNP P0ABN1
A	-2	HIS	-	expression tag	UNP P0ABN1
A	-1	GLU	-	expression tag	UNP P0ABN1
A	0	LEU	-	expression tag	UNP P0ABN1
A	41	CYS	ALA	engineered mutation	UNP P0ABN1
A	46	ALA	CYS	engineered mutation	UNP P0ABN1
A	53	VAL	ILE	engineered mutation	UNP P0ABN1
A	70	LEU	ILE	engineered mutation	UNP P0ABN1
A	96	LEU	MET	engineered mutation	UNP P0ABN1
A	107	ASP	VAL	engineered mutation	UNP P0ABN1
A	113	ALA	CYS	engineered mutation	UNP P0ABN1
B	-8	GLY	-	expression tag	UNP P0ABN1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP P0ABN1
B	-6	HIS	-	expression tag	UNP P0ABN1
B	-5	HIS	-	expression tag	UNP P0ABN1
B	-4	HIS	-	expression tag	UNP P0ABN1
B	-3	HIS	-	expression tag	UNP P0ABN1
B	-2	HIS	-	expression tag	UNP P0ABN1
B	-1	GLU	-	expression tag	UNP P0ABN1
B	0	LEU	-	expression tag	UNP P0ABN1
B	41	CYS	ALA	engineered mutation	UNP P0ABN1
B	46	ALA	CYS	engineered mutation	UNP P0ABN1
B	53	VAL	ILE	engineered mutation	UNP P0ABN1
B	70	LEU	ILE	engineered mutation	UNP P0ABN1
B	96	LEU	MET	engineered mutation	UNP P0ABN1
B	107	ASP	VAL	engineered mutation	UNP P0ABN1
B	113	ALA	CYS	engineered mutation	UNP P0ABN1
C	-8	GLY	-	expression tag	UNP P0ABN1
C	-7	HIS	-	expression tag	UNP P0ABN1
C	-6	HIS	-	expression tag	UNP P0ABN1
C	-5	HIS	-	expression tag	UNP P0ABN1
C	-4	HIS	-	expression tag	UNP P0ABN1
C	-3	HIS	-	expression tag	UNP P0ABN1
C	-2	HIS	-	expression tag	UNP P0ABN1
C	-1	GLU	-	expression tag	UNP P0ABN1
C	0	LEU	-	expression tag	UNP P0ABN1
C	41	CYS	ALA	engineered mutation	UNP P0ABN1
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C	53	VAL	ILE	engineered mutation	UNP P0ABN1
C	70	LEU	ILE	engineered mutation	UNP P0ABN1
C	96	LEU	MET	engineered mutation	UNP P0ABN1
C	107	ASP	VAL	engineered mutation	UNP P0ABN1
C	113	ALA	CYS	engineered mutation	UNP P0ABN1
D	-8	GLY	-	expression tag	UNP P0ABN1
D	-7	HIS	-	expression tag	UNP P0ABN1
D	-6	HIS	-	expression tag	UNP P0ABN1
D	-5	HIS	-	expression tag	UNP P0ABN1
D	-4	HIS	-	expression tag	UNP P0ABN1
D	-3	HIS	-	expression tag	UNP P0ABN1
D	-2	HIS	-	expression tag	UNP P0ABN1
D	-1	GLU	-	expression tag	UNP P0ABN1
D	0	LEU	-	expression tag	UNP P0ABN1
D	41	CYS	ALA	engineered mutation	UNP P0ABN1
D	46	ALA	CYS	engineered mutation	UNP P0ABN1

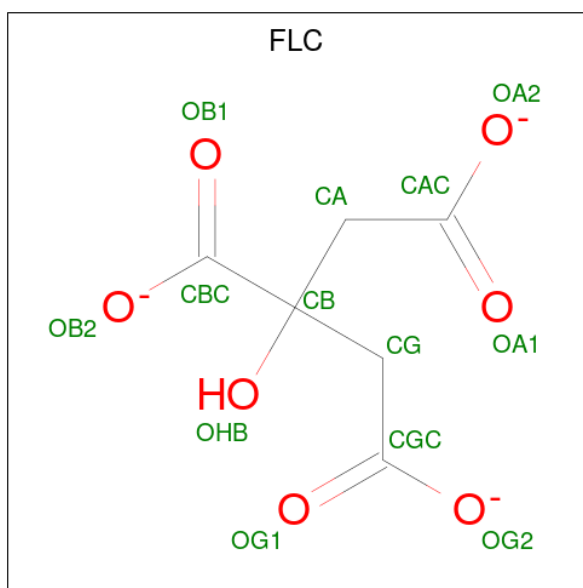
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Chain	Residue	Modelled	Actual	Comment	Reference
D	53	VAL	ILE	engineered mutation	UNP P0ABN1
D	70	LEU	ILE	engineered mutation	UNP P0ABN1
D	96	LEU	MET	engineered mutation	UNP P0ABN1
D	107	ASP	VAL	engineered mutation	UNP P0ABN1
D	113	ALA	CYS	engineered mutation	UNP P0ABN1
E	-8	GLY	-	expression tag	UNP P0ABN1
E	-7	HIS	-	expression tag	UNP P0ABN1
E	-6	HIS	-	expression tag	UNP P0ABN1
E	-5	HIS	-	expression tag	UNP P0ABN1
E	-4	HIS	-	expression tag	UNP P0ABN1
E	-3	HIS	-	expression tag	UNP P0ABN1
E	-2	HIS	-	expression tag	UNP P0ABN1
E	-1	GLU	-	expression tag	UNP P0ABN1
E	0	LEU	-	expression tag	UNP P0ABN1
E	41	CYS	ALA	engineered mutation	UNP P0ABN1
E	46	ALA	CYS	engineered mutation	UNP P0ABN1
E	53	VAL	ILE	engineered mutation	UNP P0ABN1
E	70	LEU	ILE	engineered mutation	UNP P0ABN1
E	96	LEU	MET	engineered mutation	UNP P0ABN1
E	107	ASP	VAL	engineered mutation	UNP P0ABN1
E	113	ALA	CYS	engineered mutation	UNP P0ABN1
F	-8	GLY	-	expression tag	UNP P0ABN1
F	-7	HIS	-	expression tag	UNP P0ABN1
F	-6	HIS	-	expression tag	UNP P0ABN1
F	-5	HIS	-	expression tag	UNP P0ABN1
F	-4	HIS	-	expression tag	UNP P0ABN1
F	-3	HIS	-	expression tag	UNP P0ABN1
F	-2	HIS	-	expression tag	UNP P0ABN1
F	-1	GLU	-	expression tag	UNP P0ABN1
F	0	LEU	-	expression tag	UNP P0ABN1
F	41	CYS	ALA	engineered mutation	UNP P0ABN1
F	46	ALA	CYS	engineered mutation	UNP P0ABN1
F	53	VAL	ILE	engineered mutation	UNP P0ABN1
F	70	LEU	ILE	engineered mutation	UNP P0ABN1
F	96	LEU	MET	engineered mutation	UNP P0ABN1
F	107	ASP	VAL	engineered mutation	UNP P0ABN1
F	113	ALA	CYS	engineered mutation	UNP P0ABN1

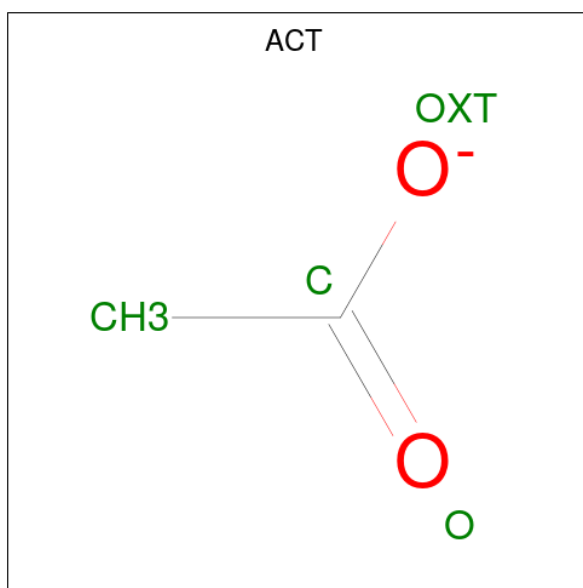
- Molecule 2 is (2S)-2,3-DIHYDROXYPROPYL(7Z)-PENTADEC-7-ENOATE (CCD ID: 78M) (formula: C₁₈H₃₄O₄).

- Molecule 4 is CITRATE ANION (CCD ID: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



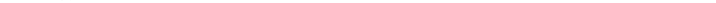
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			4	2	2		

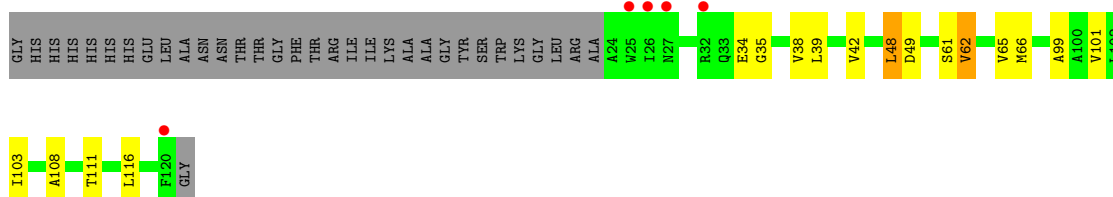
- Molecule 6 is water.

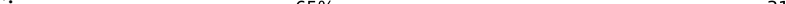
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total 3	O 3	0	0
6	B	1	Total 1	O 1	0	0
6	D	5	Total 5	O 5	0	0
6	E	4	Total 4	O 4	0	0

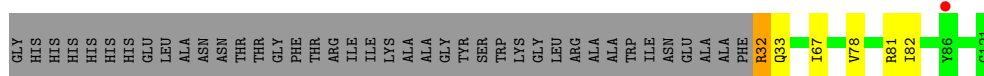
- Molecule 1: Diacylglycerol kinase

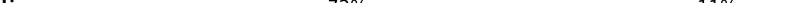


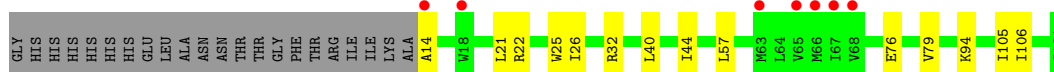
- Chain B:  4% 62% 12% 25%



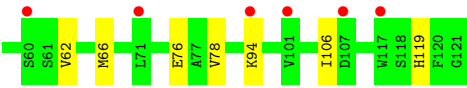
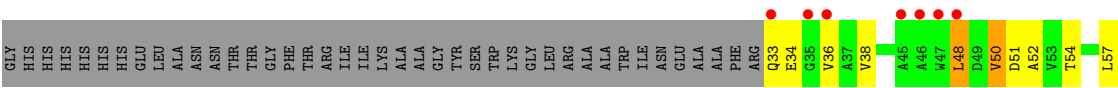
- Chain C:  65% 31% 4%



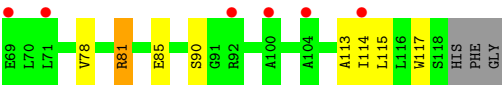
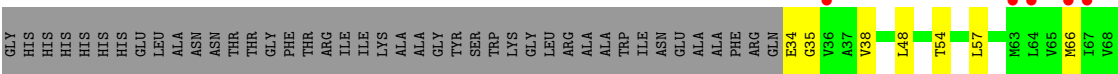
- Chain D:  5% 72% 11% 17%



- Chain E: 10% 55% 12% 32%



● Molecule 1: Diacylglycerol kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.80Å 93.04Å 143.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.52 – 2.80 46.52 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.0 (46.52-2.80) 98.7 (46.52-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.243 , 0.283 0.247 , 0.285	Depositor DCC
R_{free} test set	1288 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	87.5	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4717	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.70% 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: 78M, ZN, ACT, FLC

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	0.32	0/903	0.74	0/1229
1	B	0.31	0/762	0.71	0/1042
1	C	0.30	0/690	0.74	0/941
1	D	0.31	0/839	0.74	0/1144
1	E	0.32	0/679	0.77	0/927
1	F	0.40	0/642	0.78	0/879
All	All	0.33	0/4515	0.74	0/6162

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	887	0	922	18	0
1	B	747	0	776	12	0
1	C	681	0	716	10	0
1	D	824	0	850	11	0
1	E	670	0	703	12	0
1	F	635	0	676	10	0
2	A	88	0	136	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	66	0	102	6	0
2	C	22	0	34	0	0
2	D	66	0	102	3	0
3	D	1	0	0	0	0
4	D	13	0	5	0	0
5	D	4	0	3	0	0
6	A	3	0	0	0	0
6	B	1	0	0	0	0
6	D	5	0	0	1	0
6	E	4	0	0	0	0
All	All	4717	0	5025	69	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 (69) 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:F:54:THR:HA	1:F:57:LEU:HB2	1.71	0.72
1:A:88:GLU:OE2	1:A:88:GLU:N	2.26	0.68
1:A:54:THR:HG23	1:A:115:LEU:HD13	1.79	0.62
1:F:54:THR:HB	1:F:115:LEU:HD13	1.81	0.62
1:F:34:GLU:N	1:F:34:GLU:OE1	2.33	0.62
1:C:32:ARG:HD3	1:C:33:GLN:H	1.64	0.61
1:E:52:ALA:HB1	1:F:114:ILE:HG22	1.84	0.59
1:A:82:ILE:HD12	1:C:81:ARG:HH21	1.68	0.58
1:E:50:VAL:HG22	1:E:51:ASP:H	1.67	0.58
2:A:202:78M:H8	2:A:204:78M:H153	1.85	0.58
1:A:82:ILE:HD12	1:C:81:ARG:NH2	2.20	0.56
2:A:204:78M:H18	1:B:101:VAL:HB	1.88	0.56
1:D:14:ALA:N	6:D:302:HOH:O	2.39	0.55
2:B:201:78M:H51C	1:D:106:ILE:HG12	1.89	0.55
1:A:82:ILE:HD13	1:C:82:ILE:HG12	1.88	0.54
1:E:50:VAL:HG21	1:E:119:HIS:CE1	2.44	0.52
1:A:66:MET:HG2	2:A:201:78M:H101	1.92	0.52
1:B:42:VAL:HG11	2:D:205:78M:H153	1.91	0.51
1:C:32:ARG:HH21	1:C:33:GLN:H	1.58	0.51
1:A:17:SER:OG	2:A:204:78M:O19	2.29	0.50
2:A:202:78M:H102	2:A:204:78M:H131	1.94	0.49
1:A:13:ALA:HB1	2:A:204:78M:H202	1.94	0.49
1:A:25:TRP:CD2	1:A:32:ARG:HG2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:LEU:HD21	1:B:116:LEU:HD11	1.96	0.48
2:B:203:78M:H31C	1:D:40:LEU:HD22	1.94	0.48
1:F:38:VAL:HG21	1:F:66:MET:HE2	1.95	0.48
1:A:90:SER:HA	1:C:81:ARG:NH1	2.29	0.48
1:A:13:ALA:HB1	2:A:204:78M:H171	1.96	0.48
1:E:78:VAL:HG22	1:F:90:SER:HB3	1.96	0.48
1:B:39:LEU:HD11	1:D:44:ILE:HD13	1.95	0.47
1:A:55:ARG:NH2	2:A:202:78M:H201	2.29	0.47
1:A:43:VAL:HG13	1:A:47:TRP:CZ3	2.49	0.47
1:B:49:ASP:HB2	2:B:201:78M:H18	1.96	0.46
1:D:21:LEU:HD22	2:D:206:78M:H7	1.96	0.46
1:C:32:ARG:HD3	1:C:33:GLN:N	2.31	0.46
1:A:88:GLU:H	1:A:88:GLU:CD	2.23	0.46
1:D:76:GLU:HG3	1:D:94:LYS:HE2	1.97	0.46
1:E:34:GLU:OE1	1:E:34:GLU:N	2.50	0.45
1:E:76:GLU:HG2	1:E:94:LYS:HG2	1.98	0.45
1:B:61:SER:HB2	1:B:108:ALA:HA	1.99	0.45
2:A:203:78M:H62C	2:A:203:78M:H91C	1.69	0.44
2:A:203:78M:H22C	1:B:65:VAL:HG11	1.99	0.44
1:D:57:LEU:HD21	1:E:57:LEU:HD21	1.99	0.44
1:E:48:LEU:HD21	1:E:50:VAL:HG12	1.98	0.44
2:B:201:78M:H201	1:D:105:ILE:HD12	1.99	0.44
1:E:38:VAL:HG13	1:E:62:VAL:HG22	1.99	0.44
2:A:202:78M:H18	2:B:202:78M:H18	1.98	0.44
1:B:35:GLY:HA2	1:B:66:MET:HE3	2.00	0.43
1:B:38:VAL:HG22	1:B:62:VAL:HG12	2.00	0.43
2:A:201:78M:H141	2:A:202:78M:H121	2.00	0.43
2:A:204:78M:H92C	2:A:204:78M:H62C	1.87	0.43
1:D:25:TRP:CE2	1:D:32:ARG:HG2	2.52	0.43
1:A:22:ARG:HE	1:A:22:ARG:HB3	1.63	0.43
1:E:50:VAL:HG13	1:E:54:THR:OG1	2.18	0.43
1:A:55:ARG:CZ	2:A:202:78M:H171	2.49	0.42
1:F:81:ARG:NH1	1:F:85:GLU:HB2	2.34	0.42
2:A:203:78M:H171	1:B:34:GLU:HG2	2.01	0.42
1:B:99:ALA:O	1:B:103:ILE:HG12	2.19	0.42
2:D:206:78M:H111	1:E:106:ILE:HG21	2.01	0.42
1:F:34:GLU:HG2	1:F:35:GLY:N	2.35	0.42
1:A:100:ALA:HB1	1:C:67:ILE:HG23	2.02	0.42
1:E:33:GLN:HB3	1:E:36:VAL:HB	2.01	0.42
1:F:113:ALA:O	1:F:117:TRP:HB2	2.20	0.41
1:D:79:VAL:HG13	1:F:78:VAL:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:VAL:O	1:C:82:ILE:HG13	2.20	0.41
2:B:202:78M:H92C	2:B:202:78M:H62C	1.85	0.41
1:A:56:VAL:HG11	1:B:111:THR:HG23	2.03	0.40
1:C:82:ILE:HG13	1:C:82:ILE:H	1.67	0.40
1:D:22:ARG:O	1:D:26:ILE:HG12	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	114/130 (88%)	114 (100%)	0	0	100	100
1	B	96/130 (74%)	96 (100%)	0	0	100	100
1	C	88/130 (68%)	88 (100%)	0	0	100	100
1	D	106/130 (82%)	106 (100%)	0	0	100	100
1	E	87/130 (67%)	85 (98%)	1 (1%)	1 (1%)	11	36
1	F	83/130 (64%)	83 (100%)	0	0	100	100
All	All	574/780 (74%)	572 (100%)	1 (0%)	1 (0%)	43	72

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	50	VAL

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	90/102 (88%)	89 (99%)	1 (1%)	65	87
1	B	79/102 (78%)	77 (98%)	2 (2%)	42	76
1	C	73/102 (72%)	72 (99%)	1 (1%)	59	85
1	D	84/102 (82%)	84 (100%)	0	100	100
1	E	72/102 (71%)	70 (97%)	2 (3%)	38	73
1	F	69/102 (68%)	67 (97%)	2 (3%)	37	73
All	All	467/612 (76%)	459 (98%)	8 (2%)	53	83

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	B	48	LEU
1	B	62	VAL
1	C	32	ARG
1	E	48	LEU
1	E	66	MET
1	F	48	LEU
1	F	81	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	72	ASN
1	D	87	HIS
1	D	119	HIS
1	E	119	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 1 is monoatomic - leaving 13 for Mogul analysis.

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	78M	B	203	-	21,21,21	0.90	1 (4%)	22,22,22	0.78	1 (4%)
5	ACT	D	203	-	3,3,3	0.83	0	3,3,3	1.36	0
2	78M	B	202	-	21,21,21	0.91	1 (4%)	22,22,22	0.83	1 (4%)
2	78M	A	201	-	21,21,21	0.86	1 (4%)	22,22,22	0.81	1 (4%)
2	78M	A	204	-	21,21,21	0.88	1 (4%)	22,22,22	0.83	1 (4%)
4	FLC	D	202	-	12,12,12	1.12	0	17,17,17	1.84	5 (29%)
2	78M	A	203	-	21,21,21	0.86	1 (4%)	22,22,22	0.84	1 (4%)
2	78M	C	201	-	21,21,21	0.87	1 (4%)	22,22,22	0.81	1 (4%)
2	78M	A	202	-	21,21,21	0.89	1 (4%)	22,22,22	0.87	1 (4%)
2	78M	B	201	-	21,21,21	0.90	1 (4%)	22,22,22	0.87	1 (4%)
2	78M	D	205	-	21,21,21	0.87	1 (4%)	22,22,22	0.81	1 (4%)
2	78M	D	204	-	21,21,21	0.87	1 (4%)	22,22,22	0.84	1 (4%)
2	78M	D	206	-	21,21,21	0.89	1 (4%)	22,22,22	0.80	1 (4%)

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	78M	B	203	-	-	10/21/21/21	-
2	78M	B	202	-	-	8/21/21/21	-
2	78M	A	201	-	-	9/21/21/21	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	78M	A	204	-	-	11/21/21/21	-
4	FLC	D	202	-	-	8/16/16/16	-
2	78M	A	203	-	-	6/21/21/21	-
2	78M	C	201	-	-	10/21/21/21	-
2	78M	A	202	-	-	7/21/21/21	-
2	78M	B	201	-	-	10/21/21/21	-
2	78M	D	205	-	-	7/21/21/21	-
2	78M	D	204	-	-	6/21/21/21	-
2	78M	D	206	-	-	12/21/21/21	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	203	78M	O2-C1	2.85	1.41	1.33
2	B	202	78M	O2-C1	2.81	1.41	1.33
2	A	202	78M	O2-C1	2.78	1.41	1.33
2	B	201	78M	O2-C1	2.74	1.41	1.33
2	A	204	78M	O2-C1	2.74	1.41	1.33
2	D	206	78M	O2-C1	2.73	1.41	1.33
2	C	201	78M	O2-C1	2.70	1.41	1.33
2	D	204	78M	O2-C1	2.68	1.41	1.33
2	A	201	78M	O2-C1	2.68	1.41	1.33
2	D	205	78M	O2-C1	2.62	1.41	1.33
2	A	203	78M	O2-C1	2.55	1.40	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	202	FLC	OB2-CBC-CB	4.24	121.28	113.14
2	A	202	78M	O2-C1-C2	3.09	121.25	111.83
2	B	201	78M	O2-C1-C2	2.97	120.90	111.83
2	D	204	78M	O2-C1-C2	2.95	120.83	111.83
2	B	202	78M	O2-C1-C2	2.87	120.59	111.83
2	A	204	78M	O2-C1-C2	2.85	120.52	111.83
2	A	201	78M	O2-C1-C2	2.77	120.28	111.83
2	D	205	78M	O2-C1-C2	2.71	120.09	111.83
2	B	203	78M	O2-C1-C2	2.70	120.06	111.83
4	D	202	FLC	CB-CA-CAC	2.69	121.28	113.92
2	D	206	78M	O2-C1-C2	2.67	119.99	111.83
2	C	201	78M	O2-C1-C2	2.62	119.83	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	203	78M	O2-C1-C2	2.60	119.77	111.83
4	D	202	FLC	OA2-CAC-CA	2.53	122.37	114.35
4	D	202	FLC	OA2-CAC-OA1	-2.35	117.29	123.33
4	D	202	FLC	OG2-CGC-CG	2.01	120.73	114.35

There are no chirality outliers.

All (104) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	78M	O2-C17-C18-C20
2	A	201	78M	O2-C17-C18-O19
2	C	201	78M	O2-C17-C18-C20
2	C	201	78M	O2-C17-C18-O19
2	D	205	78M	C17-C18-C20-O21
2	D	206	78M	C17-C18-C20-O21
2	D	206	78M	O2-C17-C18-C20
2	D	206	78M	O2-C17-C18-O19
4	D	202	FLC	CA-CB-CG-CGC
4	D	202	FLC	OHB-CB-CG-CGC
2	B	203	78M	O1-C1-O2-C17
2	A	203	78M	C2-C1-O2-C17
2	A	202	78M	C2-C1-O2-C17
2	A	204	78M	C2-C1-O2-C17
2	B	203	78M	C2-C1-O2-C17
2	A	203	78M	O1-C1-O2-C17
2	A	203	78M	O2-C17-C18-O19
2	A	204	78M	O1-C1-O2-C17
2	A	202	78M	O1-C1-O2-C17
2	B	201	78M	C2-C1-O2-C17
2	C	201	78M	C2-C1-O2-C17
2	D	204	78M	C2-C1-O2-C17
2	B	201	78M	O2-C17-C18-C20
2	D	205	78M	O2-C17-C18-C20
2	D	205	78M	O2-C17-C18-O19
2	B	201	78M	O1-C1-O2-C17
2	C	201	78M	O1-C1-O2-C17
2	D	204	78M	O1-C1-O2-C17
2	B	201	78M	O2-C17-C18-O19
4	D	202	FLC	CBC-CB-CG-CGC
2	D	205	78M	O19-C18-C20-O21
2	D	206	78M	O19-C18-C20-O21
2	B	202	78M	C9-C10-C11-C12

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Mol	Chain	Res	Type	Atoms
2	D	206	78M	C2-C3-C4-C5
2	B	202	78M	C2-C1-O2-C17
2	D	205	78M	C2-C1-O2-C17
2	A	202	78M	C2-C3-C4-C5
2	D	206	78M	C11-C10-C9-C8
2	B	202	78M	O1-C1-O2-C17
2	A	204	78M	C11-C12-C13-C14
2	A	201	78M	C2-C3-C4-C5
2	B	202	78M	C10-C11-C12-C13
2	A	202	78M	C11-C10-C9-C8
2	A	204	78M	C11-C10-C9-C8
2	A	201	78M	C2-C1-O2-C17
2	B	203	78M	C5-C6-C7-C8
2	D	205	78M	O1-C1-O2-C17
2	B	202	78M	C4-C5-C6-C7
2	C	201	78M	C4-C5-C6-C7
2	D	204	78M	C2-C3-C4-C5
2	B	203	78M	C1-C2-C3-C4
2	B	202	78M	C11-C12-C13-C14
2	A	202	78M	C9-C10-C11-C12
2	A	204	78M	C2-C3-C4-C5
2	A	204	78M	C1-C2-C3-C4
2	A	201	78M	O1-C1-O2-C17
2	B	201	78M	C11-C10-C9-C8
4	D	202	FLC	CAC-CA-CB-OHB
2	A	203	78M	C2-C3-C4-C5
2	A	201	78M	C3-C4-C5-C6
2	D	206	78M	C1-C2-C3-C4
2	A	202	78M	C3-C4-C5-C6
2	B	201	78M	C2-C3-C4-C5
2	B	202	78M	C12-C13-C14-C15
2	B	201	78M	C3-C4-C5-C6
2	A	201	78M	C12-C13-C14-C15
2	D	204	78M	C12-C13-C14-C15
2	C	201	78M	C11-C10-C9-C8
2	B	201	78M	C10-C11-C12-C13
2	B	203	78M	C3-C4-C5-C6
2	B	202	78M	O19-C18-C20-O21
2	B	203	78M	C2-C3-C4-C5
2	C	201	78M	C12-C13-C14-C15
2	D	206	78M	C9-C10-C11-C12
2	A	203	78M	C10-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
2	B	203	78M	O2-C17-C18-O19
2	D	206	78M	C4-C5-C6-C7
4	D	202	FLC	CAC-CA-CB-CBC
2	D	204	78M	C11-C10-C9-C8
2	B	203	78M	C7-C8-C9-C10
2	C	201	78M	C2-C3-C4-C5
2	A	203	78M	O2-C17-C18-C20
2	A	204	78M	C4-C5-C6-C7
2	A	202	78M	C7-C8-C9-C10
2	A	204	78M	C3-C4-C5-C6
2	A	204	78M	O2-C17-C18-O19
2	A	201	78M	C7-C8-C9-C10
2	B	201	78M	C5-C6-C7-C8
2	D	206	78M	C7-C8-C9-C10
2	D	205	78M	C7-C8-C9-C10
2	D	206	78M	C10-C11-C12-C13
2	B	201	78M	C7-C8-C9-C10
2	C	201	78M	C5-C6-C7-C8
2	D	206	78M	C12-C13-C14-C15
2	A	204	78M	C7-C8-C9-C10
2	D	204	78M	C7-C8-C9-C10
2	B	203	78M	C11-C10-C9-C8
2	B	203	78M	C9-C10-C11-C12
4	D	202	FLC	CB-CA-CAC-OA1
4	D	202	FLC	CB-CA-CAC-OA2
2	A	204	78M	C10-C11-C12-C13
4	D	202	FLC	CAC-CA-CB-CG
2	C	201	78M	C7-C8-C9-C10
2	A	201	78M	C11-C10-C9-C8

There are no ring outliers.

9 monomers are involved in 23 short contacts:

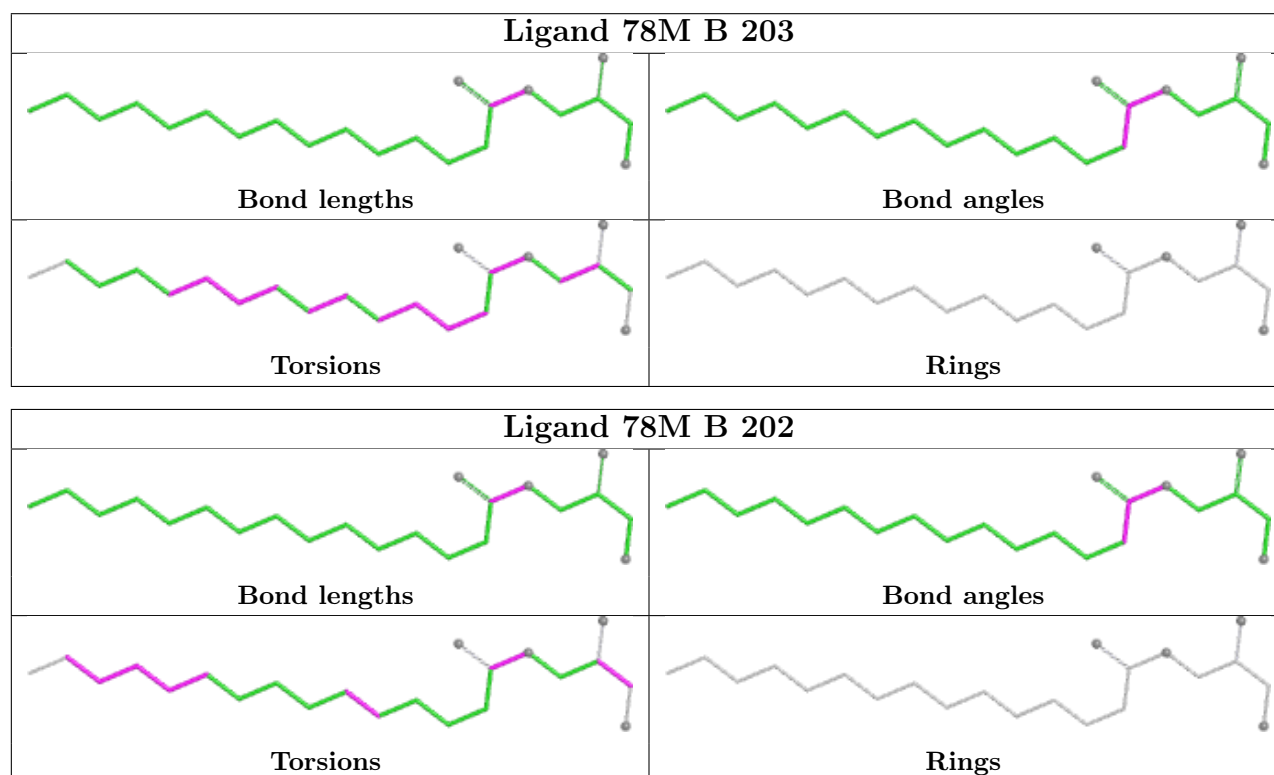
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	203	78M	1	0
2	B	202	78M	2	0
2	A	201	78M	2	0
2	A	204	78M	7	0
2	A	203	78M	3	0
2	A	202	78M	6	0
2	B	201	78M	3	0
2	D	205	78M	1	0

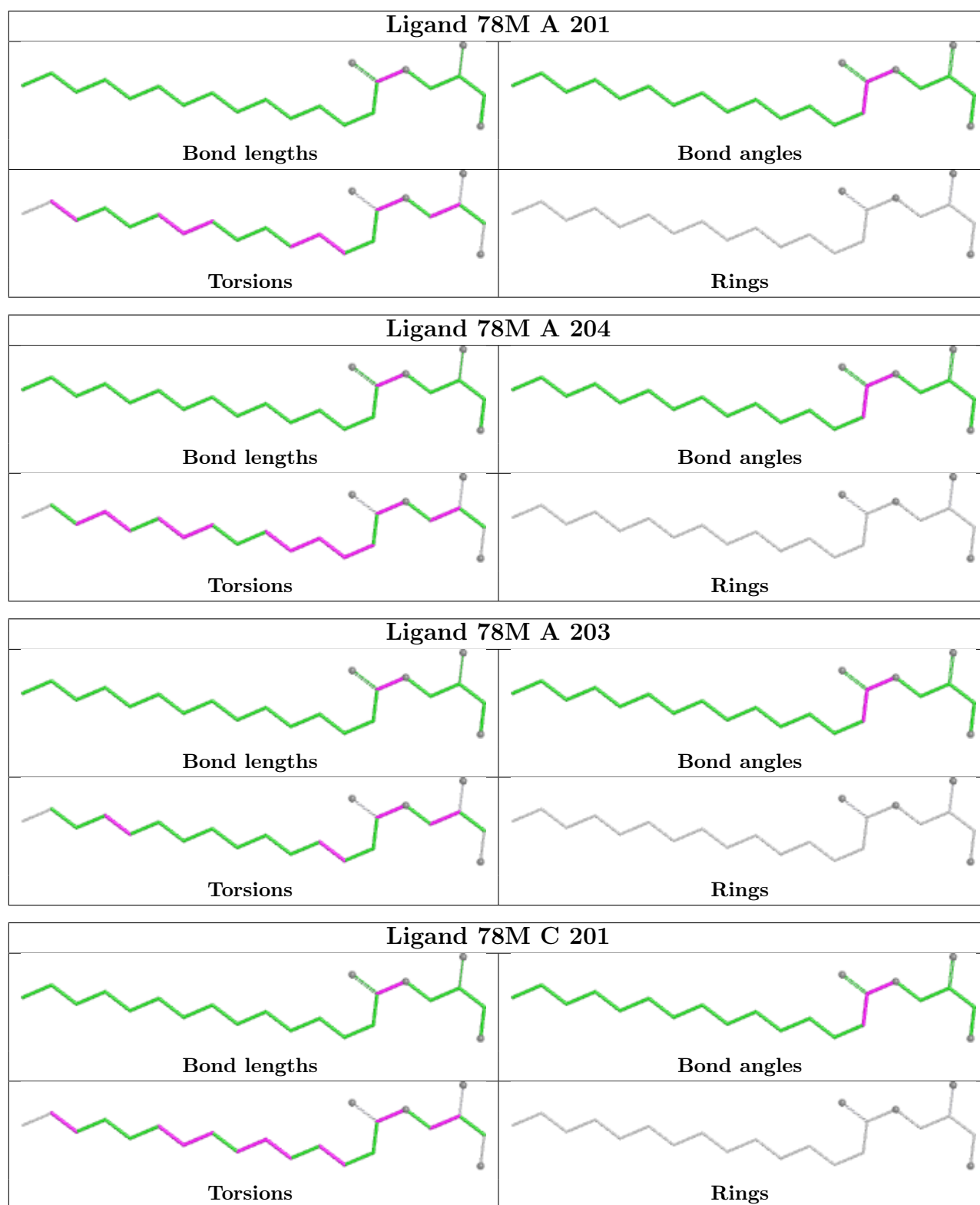
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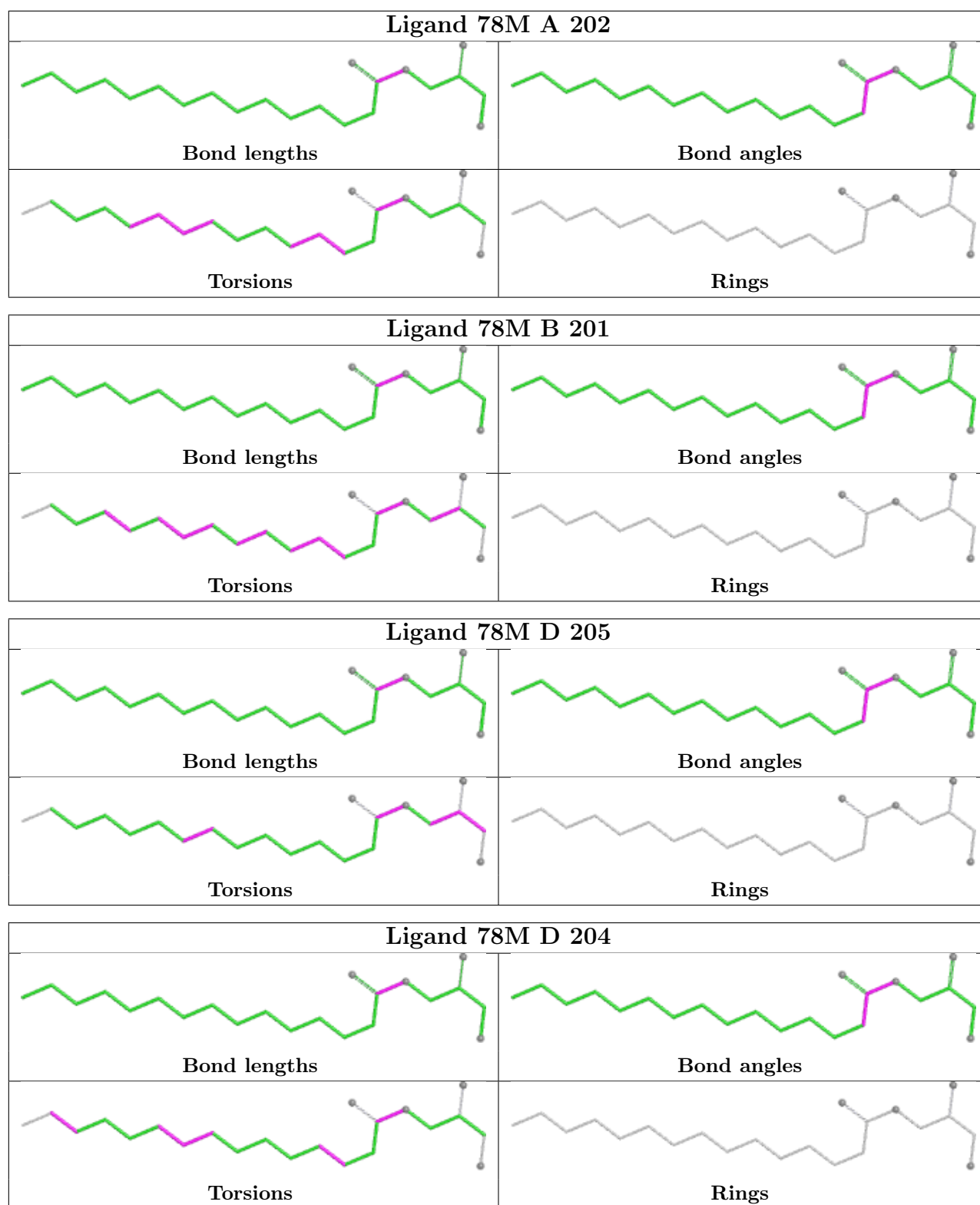
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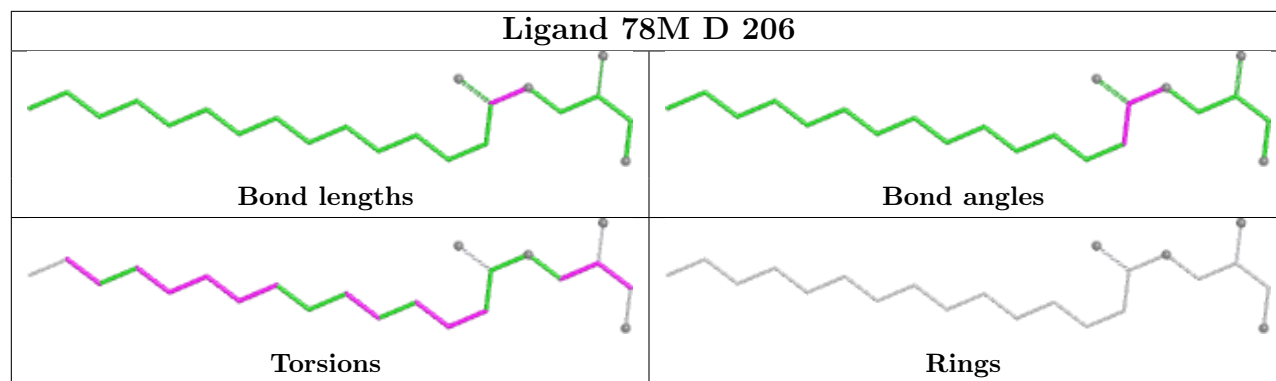
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	206	78M	2	0

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	116/130 (89%)	0.55	9 (7%)	19 14	61, 81, 134, 159	0
1	B	97/130 (74%)	0.31	5 (5%)	33 25	47, 77, 131, 142	1 (1%)
1	C	90/130 (69%)	0.30	1 (1%)	78 70	61, 81, 124, 144	0
1	D	108/130 (83%)	0.55	7 (6%)	25 18	65, 83, 117, 140	0
1	E	89/130 (68%)	0.86	13 (14%)	6 4	66, 98, 144, 175	3 (3%)
1	F	85/130 (65%)	0.87	11 (12%)	7 6	70, 97, 150, 169	0
All	All	585/780 (75%)	0.56	46 (7%)	18 13	47, 85, 138, 175	4 (0%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	100	ALA	3.9
1	D	67	ILE	3.7
1	D	18	TRP	3.5
1	B	27	ASN	3.5
1	E	46	ALA	3.3
1	F	71	LEU	3.3
1	F	67	ILE	3.3
1	E	45	ALA	3.3
1	D	14	ALA	3.2
1	F	104	ALA	3.1
1	E	48	LEU	3.0
1	F	63	MET	3.0
1	F	36	VAL	3.0
1	A	120	PHE	3.0
1	E	35	GLY	3.0
1	E	36	VAL	3.0
1	A	39	LEU	2.9
1	F	66	MET	2.9
1	B	25	TRP	2.8

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Mol	Chain	Res	Type	RSRZ
1	E	71	LEU	2.8
1	B	26	ILE	2.8
1	B	32	ARG	2.8
1	F	69	GLU	2.8
1	D	63	MET	2.8
1	A	47	TRP	2.7
1	E	33	GLN	2.6
1	D	66	MET	2.6
1	B	120	PHE	2.6
1	E	47	TRP	2.6
1	A	117	TRP	2.5
1	D	65	VAL	2.5
1	E	117	TRP	2.5
1	A	71	LEU	2.4
1	A	89	LEU	2.4
1	A	96	LEU	2.4
1	E	101	VAL	2.4
1	A	100	ALA	2.4
1	F	92	ARG	2.3
1	E	60	SER	2.3
1	D	68	VAL	2.3
1	E	94	LYS	2.2
1	F	64	LEU	2.2
1	A	6	GLY	2.2
1	F	114	ILE	2.1
1	C	86	TYR	2.1
1	E	107	ASP	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,

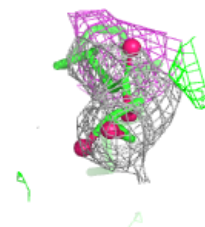
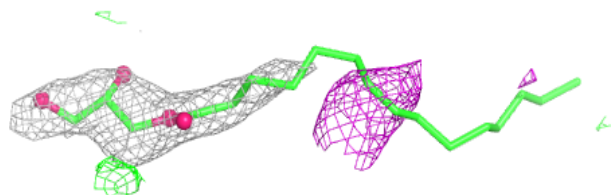
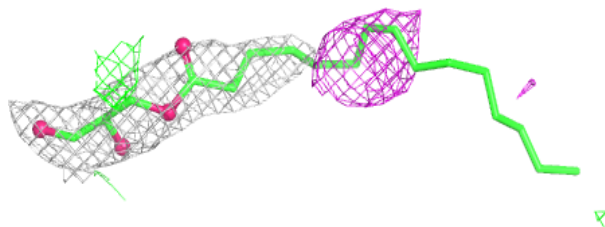
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	FLC	D	202	13/13	0.31	0.20	161,180,191,191	0
2	78M	D	206	22/22	0.80	0.33	74,98,117,122	0
2	78M	A	202	22/22	0.83	0.26	75,110,129,132	0
2	78M	B	201	22/22	0.84	0.28	120,135,145,156	0
2	78M	C	201	22/22	0.84	0.22	64,101,134,144	0
2	78M	B	203	22/22	0.85	0.30	73,97,127,135	22
2	78M	A	203	22/22	0.85	0.20	81,99,129,142	0
2	78M	A	201	22/22	0.85	0.19	66,93,115,119	0
2	78M	B	202	22/22	0.85	0.18	81,113,126,128	0
5	ACT	D	203	4/4	0.87	0.23	113,133,135,138	0
2	78M	D	204	22/22	0.88	0.28	102,121,127,128	0
2	78M	A	204	22/22	0.88	0.20	76,90,107,113	22
2	78M	D	205	22/22	0.92	0.24	107,131,144,155	0
3	ZN	D	201	1/1	0.93	0.07	120,120,120,120	1

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.

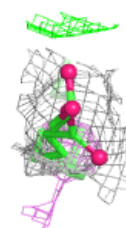
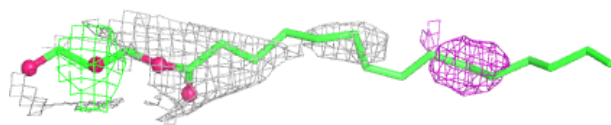
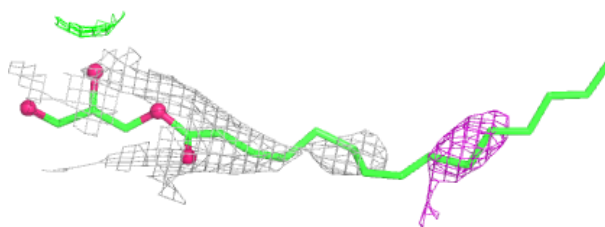
Electron density around 78M D 206:

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

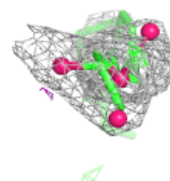
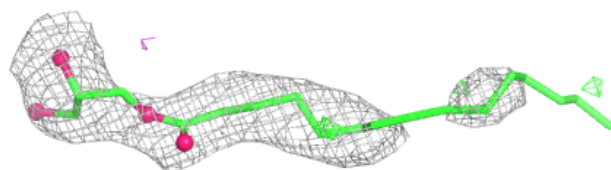
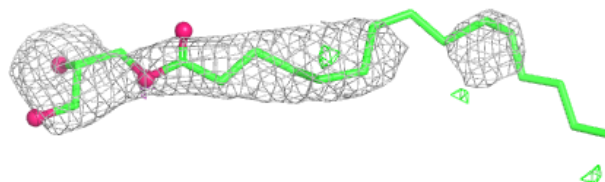


Electron density around 78M A 202:

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

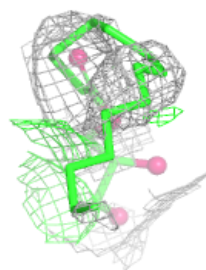
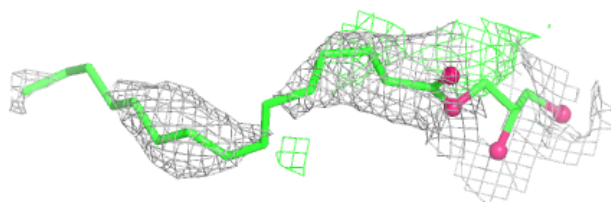
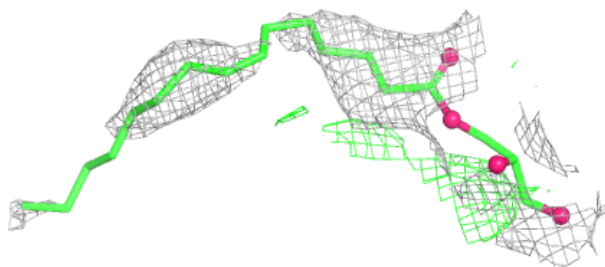
**Electron density around 78M B 201:**

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

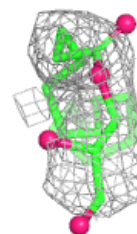
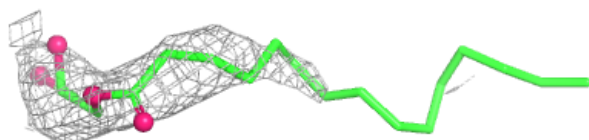
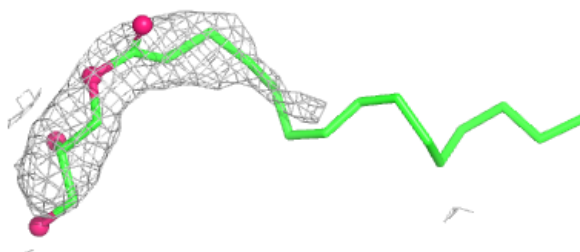


Electron density around 78M C 201:

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

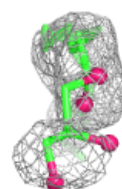
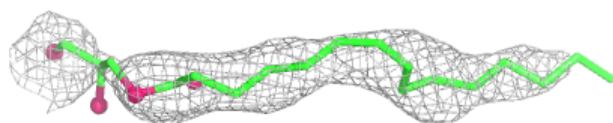
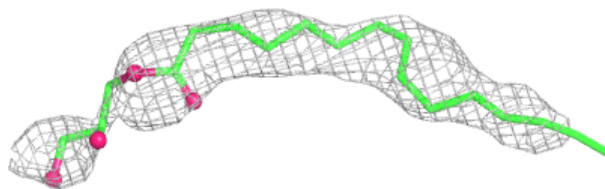
**Electron density around 78M B 203:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

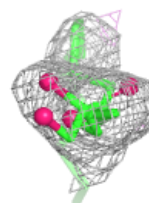
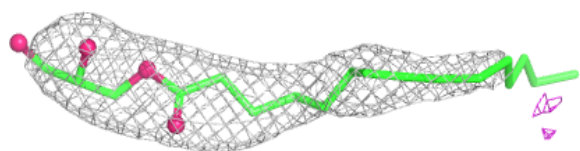
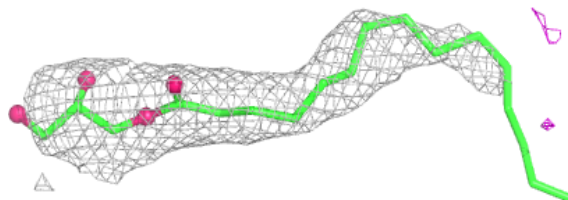


Electron density around 78M A 203:

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

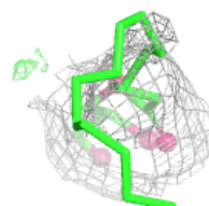
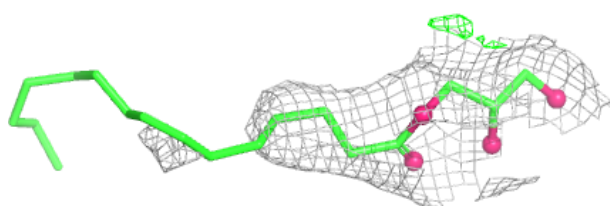
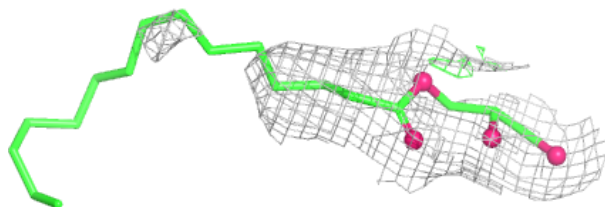
**Electron density around 78M A 201:**

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

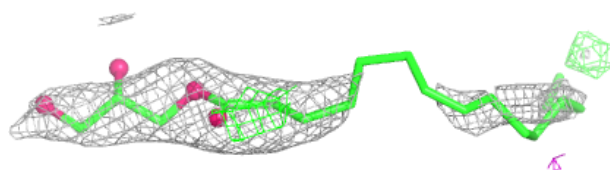
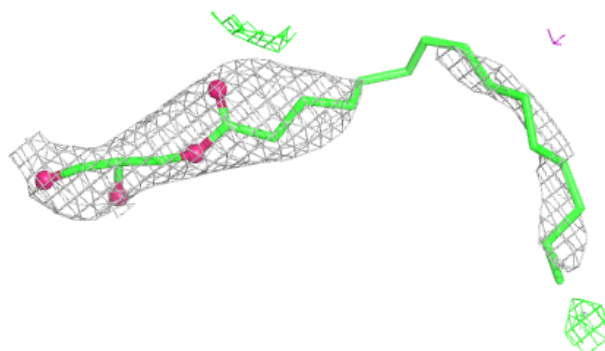


Electron density around 78M B 202:

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

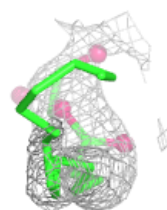
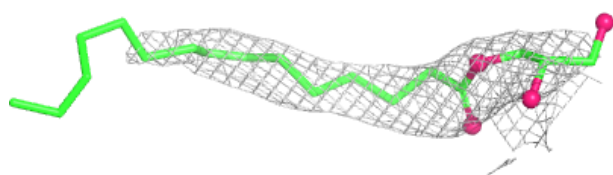
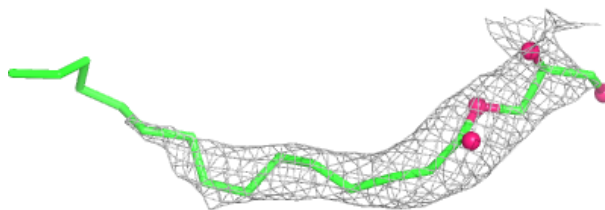
**Electron density around 78M D 204:**

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

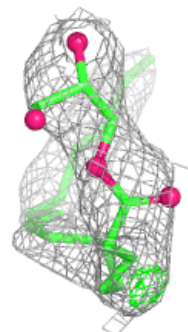
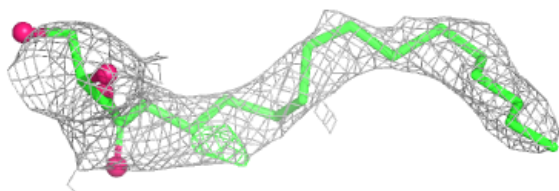
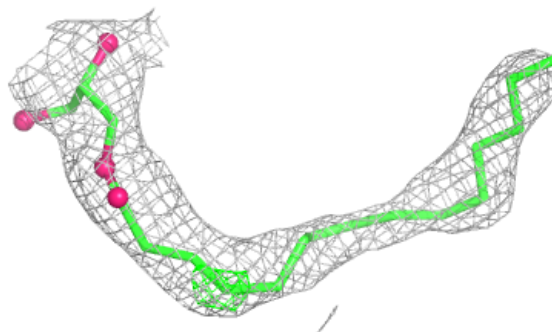


Electron density around 78M A 204:

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

**Electron density around 78M D 205:**

$2mF_o-DF_c$ (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.