Full wwPDB X-ray Structure Validation Report (i)

May 27, 2020 – 12:21 am BST

PDB ID	:	6OPB
Title	:	Structure of Aedes aegypti OBP22 in the complex with arachidic acid
Authors	:	Jones, D.N.; Wang, J.
Deposited on	:	2019-04-24
Resolution	:	2.00 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 (i) symbol.

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

|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance ⓘ

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

The reported resolution of this entry is 2.00 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.
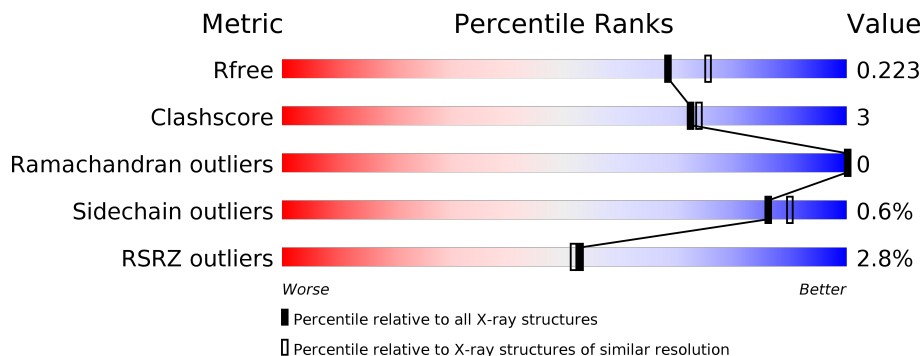
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	123	2% / 93% / 6% ·
1	B	123	2% / 86% / 13% ·
1	C	123	88% / 11% ·
1	D	123	2% / 91% / 8% ·
1	E	123	8% / 88% / 9% ··
1	F	123	8% / 87% / 11% ·

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Mol	Chain	Length	Quality of chain
1	G	123	90% 8% ·
1	H	123	% 94% 5% ·
1	I	123	2% 88% 10% ·

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 17442 atoms, of which 8431 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 AAEL005772-PA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	E	120	Total 1881	C 609	H 916	N 164	O 184	S 8	0	0	0
1	A	121	Total 1911	C 618	H 929	N 169	O 187	S 8	0	2	0
1	B	122	Total 1916	C 619	H 933	N 167	O 189	S 8	0	0	0
1	F	120	Total 1829	C 600	H 881	N 156	O 184	S 8	0	0	0
1	C	121	Total 1914	C 618	H 931	N 170	O 187	S 8	0	2	0
1	D	122	Total 1916	C 619	H 933	N 167	O 189	S 8	0	0	0
1	G	121	Total 1902	C 615	H 927	N 166	O 186	S 8	0	0	0
1	H	122	Total 1915	C 619	H 932	N 167	O 189	S 8	0	0	0
1	I	120	Total 1900	C 614	H 926	N 165	O 187	S 8	0	1	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	1	MET	-	initiating methionine	UNP Q1HRL7
A	1	MET	-	initiating methionine	UNP Q1HRL7
B	1	MET	-	initiating methionine	UNP Q1HRL7
F	1	MET	-	initiating methionine	UNP Q1HRL7
C	1	MET	-	initiating methionine	UNP Q1HRL7
D	1	MET	-	initiating methionine	UNP Q1HRL7
G	1	MET	-	initiating methionine	UNP Q1HRL7
H	1	MET	-	initiating methionine	UNP Q1HRL7
I	1	MET	-	initiating methionine	UNP Q1HRL7

- Molecule 2 is icosanoic acid (three-letter code: DCR) (formula: $C_{20}H_{40}O_2$) (labeled as "Lig-

and of Interest" by author).
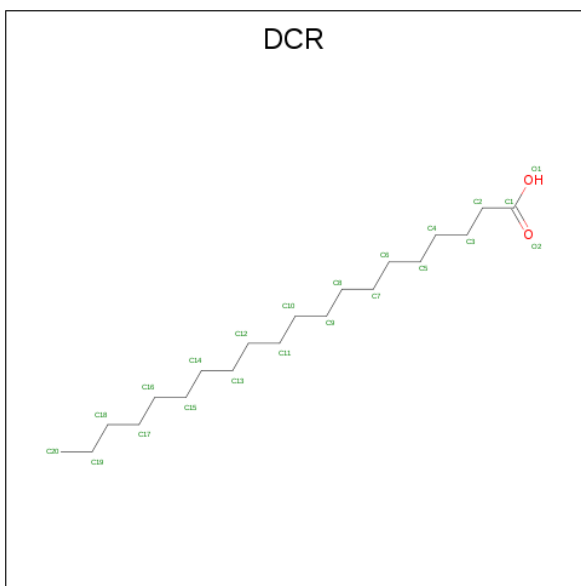
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	E	1	Total	C	H	O	0	0
			61	20	39	2		
2	F	1	Total	C	H	O	0	0
			61	20	39	2		
2	I	1	Total	C	H	O	0	0
			61	20	39	2		

- Molecule 3 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).
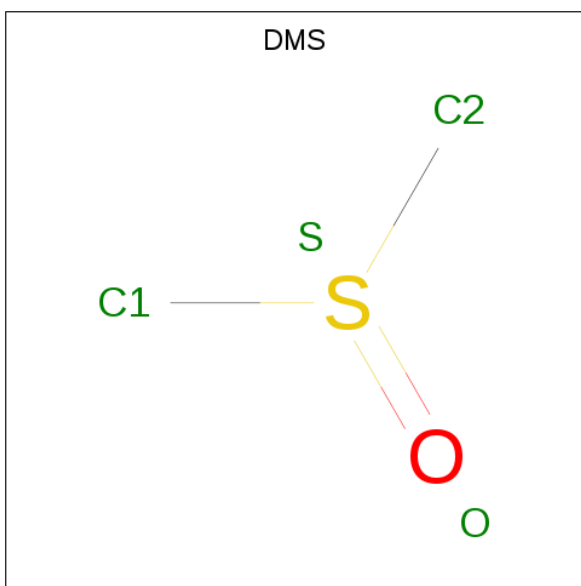
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total 10	C 2	H 6	O 1	S 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	7	Total 7	O 7	0	0
4	A	15	Total 15	O 15	0	0
4	B	22	Total 22	O 22	0	0
4	F	8	Total 8	O 8	0	0
4	C	14	Total 14	O 14	0	0
4	D	23	Total 23	O 23	0	0
4	G	14	Total 14	O 14	0	0
4	H	26	Total 26	O 26	0	0
4	I	36	Total 36	O 36	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: AAEL005772-PA
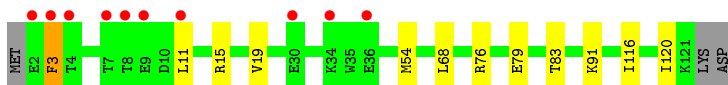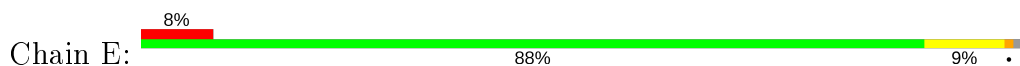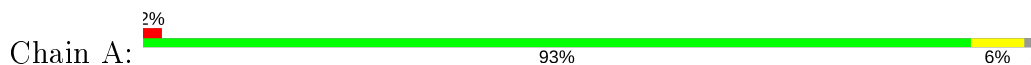
Chain E:

- Molecule 1: AAEL005772-PA

Chain A:

- Molecule 1: AAEL005772-PA
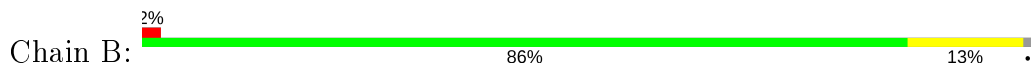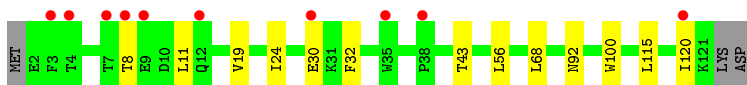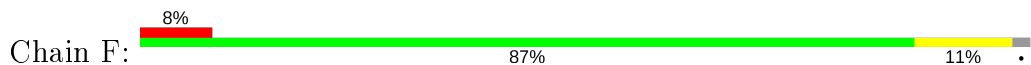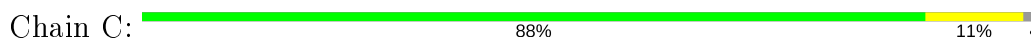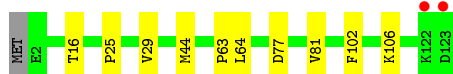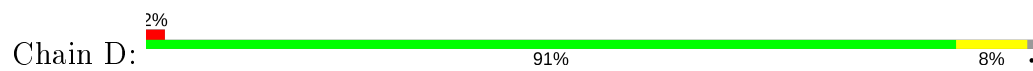
Chain B:

- Molecule 1: AAEL005772-PA

Chain F:

- Molecule 1: AAEL005772-PA

Chain C:

- Molecule 1: AAEL005772-PA

Chain D:

Chain D residue bar: MET E2 T16 P25 V29 M44 P63 L64 D77 V81 F102 K106 K122 D123

- Molecule 1: AAEL005772-PA

Chain G:

Chain G residue bar: MET E2 M44 G62 V65 L72 E79 T83 W100 A101 F102 N111 L115 K122 ASP

- Molecule 1: AAEL005772-PA

Chain H:

Chain H residue bar: MET E2 M44 L72 E80 E84 F102 L115 K122 D123

- Molecule 1: AAEL005772-PA

Chain I:

Chain I residue bar: MET E2 F3 T8 V29 K33 P38 E39 D40 M44 L68 A73 H74 G75 V81 V89 S119 I120 K121 LYS ASP
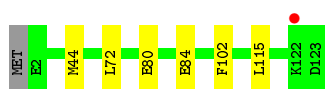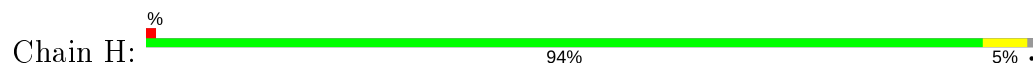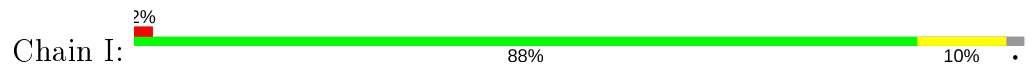
4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	154.26Å 154.26Å 58.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.56 – 2.00 38.56 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.1 (38.56-2.00) 99.9 (38.56-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	2.94 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.176 , 0.218 0.187 , 0.223	Depositor DCC
R_{free} test set	5097 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å^2)	23.2	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.46 , 42.7	EDS
L-test for twinning[2]	$<|L|> = 0.50, <L^2> = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -h,-k,l 0.023 for h,-h-k,-l 0.140 for -k,-h,-l	Xtriage
Reported twinning fraction	0.505 for H, K, L 0.495 for -K, -H, -L	Depositor
Outliers	0 of 104289 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	17442	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.10 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.1428e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $<|L|>$, $<L^2>$ 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: DMS, DCR

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	0.26	0/1009	0.51	0/1365
1	B	0.28	0/1003	0.57	0/1356
1	C	0.25	0/1010	0.55	0/1365
1	D	0.28	0/1003	0.56	0/1356
1	E	0.26	0/985	0.54	0/1332
1	F	0.27	0/968	0.52	0/1312
1	G	0.27	0/995	0.55	0/1345
1	H	0.29	0/1003	0.61	0/1356
1	I	0.28	0/997	0.56	0/1348
All	All	0.27	0/8973	0.55	0/12135

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	982	929	921	6	0
1	B	983	933	930	11	0
1	C	983	931	923	8	0
1	D	983	933	930	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	965	916	916	9	0
1	F	948	881	882	7	0
1	G	975	927	926	5	0
1	H	983	932	930	3	0
1	I	974	926	926	10	0
2	E	22	39	39	4	0
2	F	22	39	39	1	0
2	I	22	39	39	4	0
3	B	4	6	6	1	0
4	A	15	0	0	0	0
4	B	22	0	0	0	0
4	C	14	0	0	0	0
4	D	23	0	0	1	0
4	E	7	0	0	0	0
4	F	8	0	0	0	0
4	G	14	0	0	1	0
4	H	26	0	0	0	0
4	I	36	0	0	1	0
All	All	9011	8431	8407	58	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 3.

All (58) 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:I:68:LEU:HD13	2:I:201:DCR:H20B	1.69	0.75
1:C:65:VAL:O	1:C:69:VAL:HG23	1.92	0.69
1:A:118:ALA:HB3	1:B:109:GLN:OE1	1.96	0.66
1:F:19:VAL:HA	1:F:24:ILE:HD12	1.79	0.62
1:I:120:ILE:HD11	2:I:201:DCR:C8	2.30	0.61
1:I:29:VAL:HG12	1:I:33:LYS:HE2	1.82	0.60
1:E:79:GLU:O	1:E:83:THR:HG23	2.03	0.59
1:E:68:LEU:HD22	2:E:201:DCR:H20B	1.87	0.56
1:G:44:MET:HG2	1:G:102:PHE:CG	2.41	0.56
1:D:25:PRO:O	1:D:29:VAL:HG23	2.06	0.55
1:C:40:ASP:O	1:C:44:MET:HG3	2.07	0.55
1:E:15:ARG:O	1:E:19:VAL:HG23	2.06	0.54
1:I:73:ALA:HB2	1:I:81:VAL:HG21	1.91	0.53
1:B:11:LEU:HD21	3:B:201:DMS:O	2.09	0.52
1:A:82:ARG:HG2	1:A:86:LEU:HD12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:115:LEU:HD21	1:D:16:THR:HG21	1.92	0.51
1:D:77:ASP:O	1:D:81:VAL:HG23	2.11	0.50
1:B:76:ARG:NH2	1:B:84:GLU:OE1	2.45	0.49
1:A:65:VAL:O	1:A:69:VAL:HG23	2.12	0.49
1:F:56:LEU:HD22	1:F:68:LEU:HD23	1.95	0.49
1:C:15[B]:ARG:O	1:C:19:VAL:HG23	2.12	0.49
1:F:92:ASN:HB2	1:F:100:TRP:CZ3	2.48	0.48
1:E:116:ILE:HG23	2:E:201:DCR:C14	2.44	0.47
1:A:34:LYS:HD3	1:B:7:THR:HG21	1.97	0.47
1:C:79:GLU:O	1:C:83:THR:HG23	2.15	0.47
1:E:116:ILE:HG23	2:E:201:DCR:H14	1.96	0.47
1:D:63:PRO:HG3	4:D:205:HOH:O	2.14	0.46
1:B:44:MET:HG2	1:B:102:PHE:CG	2.51	0.46
1:C:120:ILE:HD13	1:D:106:LYS:HG3	1.97	0.46
1:F:11:LEU:HD11	2:F:201:DCR:H3	1.97	0.46
1:F:32:PHE:CE1	1:F:43:THR:HG23	2.50	0.46
1:G:65:VAL:HG23	4:G:203:HOH:O	2.14	0.46
1:C:15[A]:ARG:O	1:C:19:VAL:HG23	2.16	0.45
1:B:8:THR:O	1:B:12:GLN:HG3	2.16	0.45
1:I:75:GLY:HA2	4:I:332:HOH:O	2.16	0.45
1:H:80:GLU:O	1:H:84:GLU:HG3	2.17	0.45
1:I:40:ASP:O	1:I:44:MET:HG3	2.18	0.44
1:B:77:ASP:O	1:B:81:VAL:HG23	2.19	0.43
1:H:72:LEU:HD21	1:H:115:LEU:HD23	2.00	0.43
1:I:119:SER:OG	1:I:120:ILE:HD12	2.18	0.43
1:F:8:THR:OG1	1:F:120:ILE:HG22	2.18	0.43
1:E:120:ILE:HD11	2:E:201:DCR:H10A	2.01	0.43
1:G:62:GLY:HA3	1:G:100:TRP:CH2	2.55	0.42
1:D:44:MET:HG2	1:D:102:PHE:CG	2.54	0.42
1:I:120:ILE:HD11	2:I:201:DCR:H8A	1.99	0.42
1:E:3:PHE:HE2	1:E:54:MET:O	2.03	0.41
1:E:11:LEU:HD22	1:E:120:ILE:HD13	2.02	0.41
1:I:120:ILE:HD12	1:I:120:ILE:N	2.36	0.41
1:I:89:VAL:HG22	2:I:201:DCR:H20	2.03	0.41
1:E:76:ARG:HB3	1:B:19:VAL:HG12	2.03	0.41
1:G:72:LEU:HD21	1:G:115:LEU:HD23	2.03	0.41
1:C:68:LEU:O	1:C:72:LEU:HG	2.20	0.41
1:A:8:THR:OG1	1:B:34:LYS:HA	2.21	0.40
1:C:92:ASN:HB2	1:C:100:TRP:CE3	2.56	0.40
1:B:95:ASN:OD1	1:D:64:LEU:HB3	2.20	0.40
1:G:79:GLU:O	1:G:83:THR:HG23	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:ALA:HA	1:B:35:TRP:HB3	2.03	0.40
1:H:44:MET:HG2	1:H:102:PHE:CG	2.57	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	121/123 (98%)	119 (98%)	2 (2%)	0	100	100
1	B	120/123 (98%)	116 (97%)	4 (3%)	0	100	100
1	C	121/123 (98%)	121 (100%)	0	0	100	100
1	D	120/123 (98%)	118 (98%)	2 (2%)	0	100	100
1	E	118/123 (96%)	112 (95%)	6 (5%)	0	100	100
1	F	118/123 (96%)	110 (93%)	8 (7%)	0	100	100
1	G	119/123 (97%)	119 (100%)	0	0	100	100
1	H	120/123 (98%)	118 (98%)	2 (2%)	0	100	100
1	I	119/123 (97%)	119 (100%)	0	0	100	100
All	All	1076/1107 (97%)	1052 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	109/114 (96%)	109 (100%)	0	100	100
1	B	109/114 (96%)	108 (99%)	1 (1%)	78	83
1	C	109/114 (96%)	109 (100%)	0	100	100
1	D	109/114 (96%)	109 (100%)	0	100	100
1	E	107/114 (94%)	105 (98%)	2 (2%)	57	61
1	F	104/114 (91%)	103 (99%)	1 (1%)	76	81
1	G	108/114 (95%)	107 (99%)	1 (1%)	78	83
1	H	109/114 (96%)	109 (100%)	0	100	100
1	I	109/114 (96%)	108 (99%)	1 (1%)	78	83
All	All	973/1026 (95%)	967 (99%)	6 (1%)	86	90

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	3	PHE
1	E	91	LYS
1	B	41	ASP
1	F	30	GLU
1	G	111	ASN
1	I	3	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	112	ASN
1	A	111	ASN
1	C	55	GLN
1	C	109	GLN
1	H	74	HIS

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 carbohydrates 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	DCR	I	201	-	18,21,21	0.16	0	17,21,21	0.80	0
3	DMS	B	201	-	3,3,3	0.28	0	3,3,3	0.17	0
2	DCR	F	201	-	18,21,21	0.28	0	17,21,21	0.48	0
2	DCR	E	201	-	18,21,21	0.24	0	17,21,21	0.57	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
2	DCR	I	201	-	-	11/17/19/19	-
2	DCR	F	201	-	-	10/17/19/19	-
2	DCR	E	201	-	-	13/17/19/19	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	201	DCR	C16-C17-C18-C19
2	E	201	DCR	C11-C10-C9-C8
2	I	201	DCR	C13-C14-C15-C16

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Mol	Chain	Res	Type	Atoms
2	F	201	DCR	C9-C10-C11-C12
2	F	201	DCR	C15-C16-C17-C18
2	I	201	DCR	C7-C8-C9-C10
2	I	201	DCR	C11-C12-C13-C14
2	F	201	DCR	C7-C8-C9-C10
2	I	201	DCR	C10-C11-C12-C13
2	F	201	DCR	C14-C15-C16-C17
2	E	201	DCR	C3-C4-C5-C6
2	E	201	DCR	C10-C11-C12-C13
2	F	201	DCR	C4-C5-C6-C7
2	F	201	DCR	C13-C14-C15-C16
2	E	201	DCR	C12-C13-C14-C15
2	I	201	DCR	C12-C13-C14-C15
2	E	201	DCR	C4-C5-C6-C7
2	E	201	DCR	C11-C12-C13-C14
2	F	201	DCR	C3-C4-C5-C6
2	F	201	DCR	C11-C10-C9-C8
2	E	201	DCR	C5-C6-C7-C8
2	E	201	DCR	C9-C10-C11-C12
2	I	201	DCR	C17-C18-C19-C20
2	I	201	DCR	C15-C16-C17-C18
2	E	201	DCR	C13-C14-C15-C16
2	I	201	DCR	C11-C10-C9-C8
2	I	201	DCR	C6-C7-C8-C9
2	F	201	DCR	C11-C12-C13-C14
2	I	201	DCR	C3-C4-C5-C6
2	E	201	DCR	C6-C7-C8-C9
2	E	201	DCR	C7-C8-C9-C10
2	F	201	DCR	C6-C7-C8-C9
2	E	201	DCR	C15-C16-C17-C18
2	E	201	DCR	C17-C18-C19-C20

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	201	DCR	4	0
3	B	201	DMS	1	0
2	F	201	DCR	1	0
2	E	201	DCR	4	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 then 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.
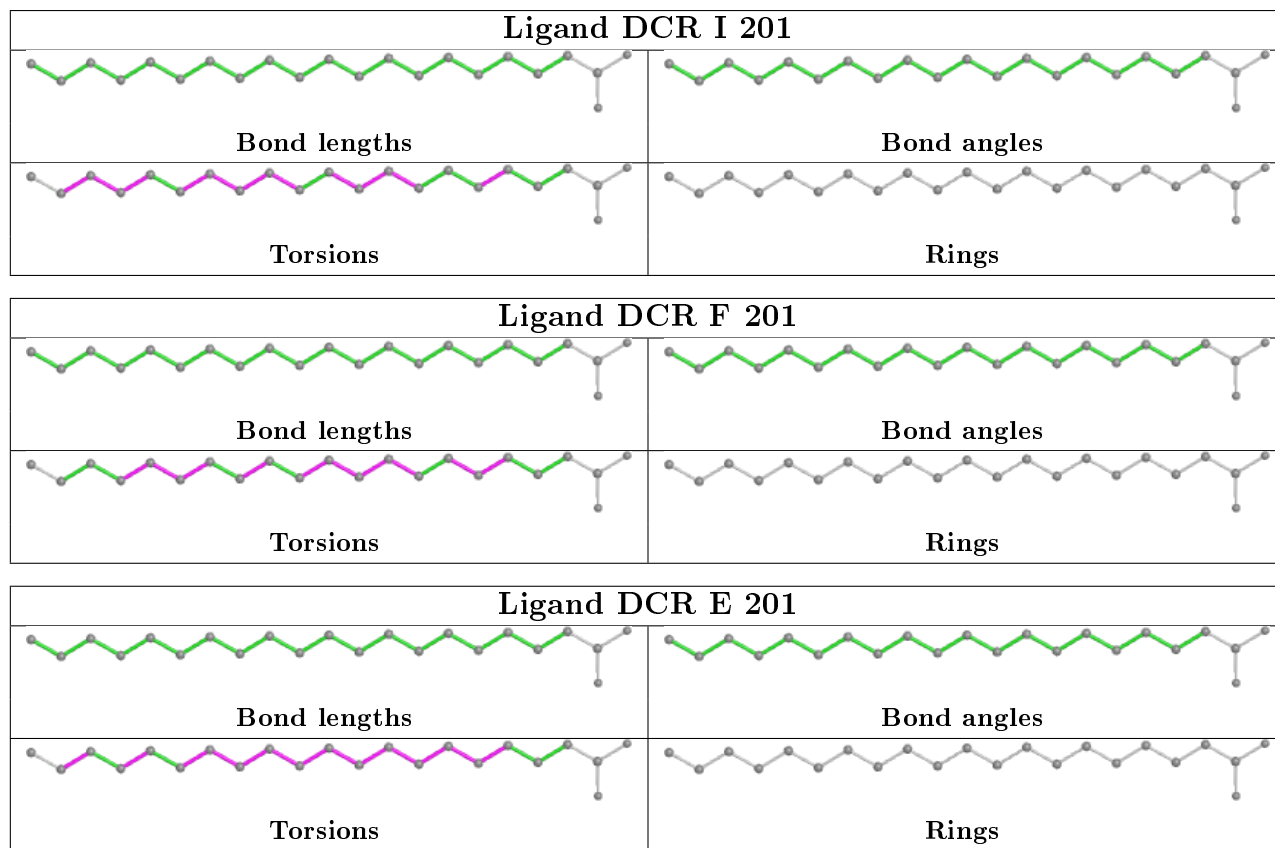
5.7 Other polymers (i)

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	121/123 (98%)	0.02	2 (1%)	70	68	21, 32, 50, 61	0
1	B	122/123 (99%)	0.17	3 (2%)	57	56	15, 25, 38, 52	0
1	C	121/123 (98%)	0.02	0	100	100	20, 30, 44, 50	0
1	D	122/123 (99%)	0.00	2 (1%)	72	70	15, 27, 47, 58	0
1	E	120/123 (97%)	0.11	10 (8%)	11	10	19, 30, 49, 55	0
1	F	120/123 (97%)	0.23	10 (8%)	11	10	21, 32, 48, 54	0
1	G	121/123 (98%)	-0.06	0	100	100	17, 25, 36, 45	0
1	H	122/123 (99%)	-0.02	1 (0%)	86	85	13, 23, 43, 51	0
1	I	120/123 (97%)	0.01	2 (1%)	70	68	13, 22, 37, 45	0
All	All	1089/1107 (98%)	0.06	30 (2%)	53	51	13, 28, 46, 61	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	4	THR	4.5
1	B	122	LYS	3.3
1	E	34	LYS	3.2
1	E	7	THR	3.1
1	D	123	ASP	3.1
1	H	122	LYS	2.9
1	A	122	LYS	2.8
1	F	4	THR	2.7
1	F	35	TRP	2.7
1	I	8	THR	2.7
1	E	11	LEU	2.7
1	E	3	PHE	2.5
1	B	25	PRO	2.5
1	F	30	GLU	2.5
1	B	123	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	122	LYS	2.4
1	F	38	PRO	2.3
1	E	30	GLU	2.3
1	F	3	PHE	2.2
1	I	38	PRO	2.2
1	E	8	THR	2.2
1	E	36	GLU	2.2
1	F	12	GLN	2.2
1	F	7	THR	2.1
1	F	8	THR	2.1
1	F	9	GLU	2.1
1	F	120	ILE	2.0
1	E	2	GLU	2.0
1	A	27	ASP	2.0
1	E	9	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

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

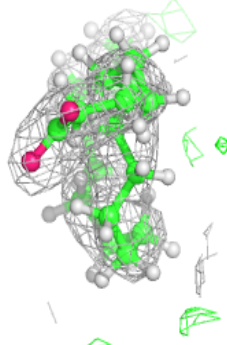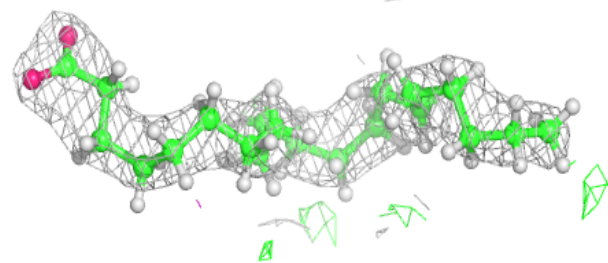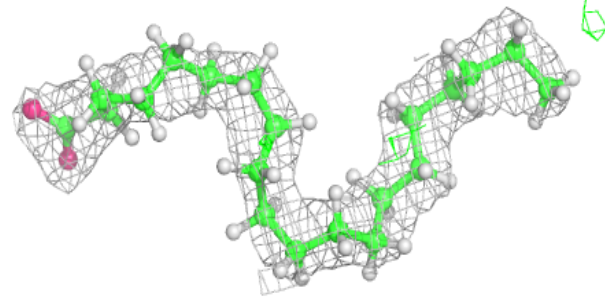| Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	DCR	F	201	22/22	0.82	0.19	30,41,57,61	0
2	DCR	E	201	22/22	0.84	0.18	39,43,49,51	0
2	DCR	I	201	22/22	0.86	0.14	27,29,30,35	0
3	DMS	B	201	4/4	0.91	0.18	35,37,39,39	0

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.
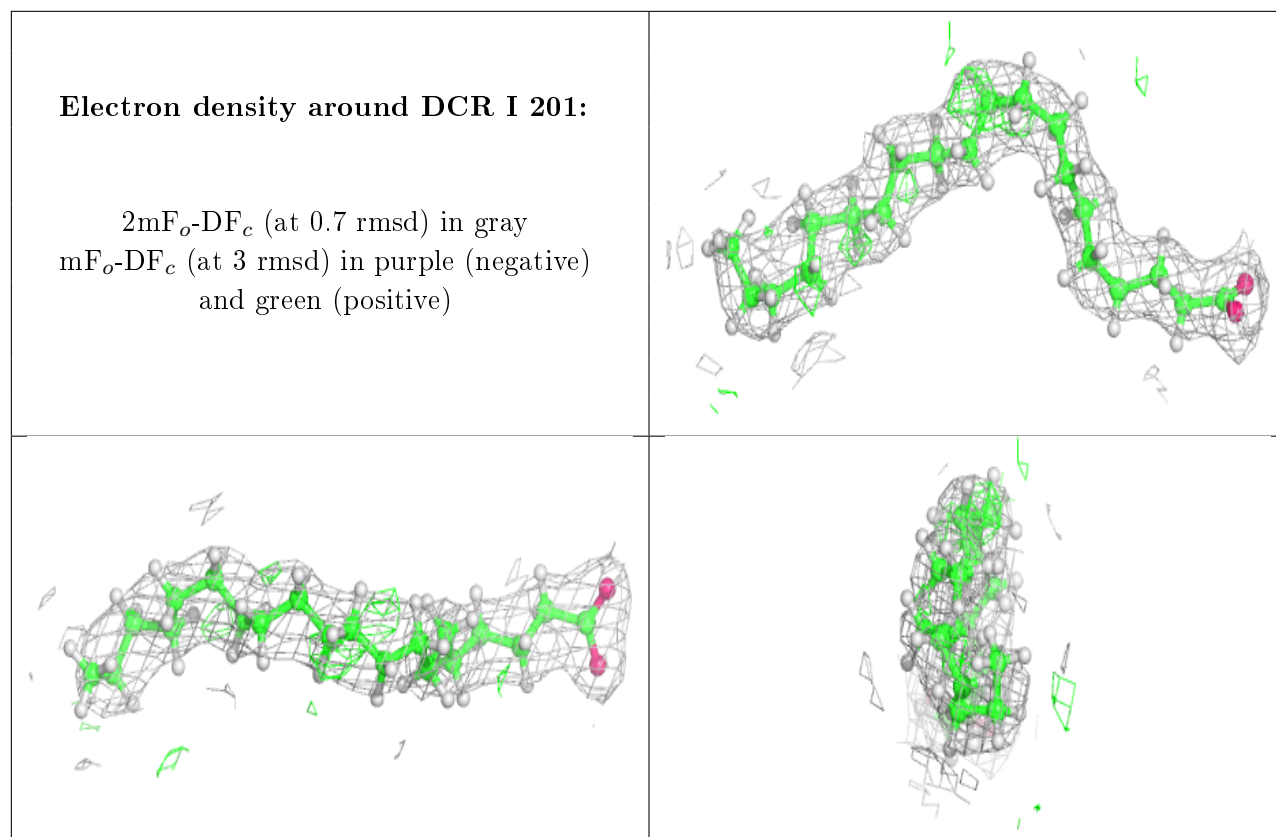
Electron density around DCR F 201:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
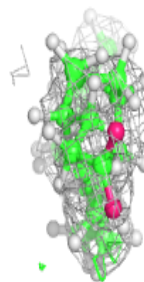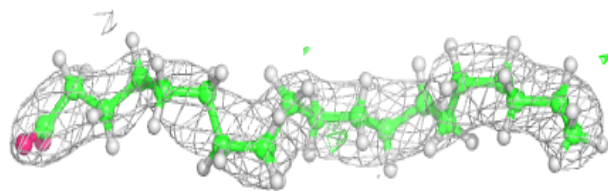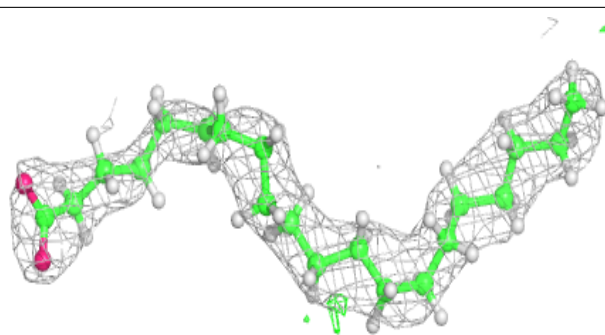
Electron density around DCR E 201:

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

Electron density around DCR I 201:

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

6.5 Other polymers (i)

There are no such residues in this entry.