



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:53 PM UTC

PDB ID : 5C1M / pdb_00005c1m
Title : Crystal structure of active mu-opioid receptor bound to the agonist BU72
Authors : Huang, W.J.; Manglik, A.; Venkatakrisnan, A.J.; Laeremans, T.; Feinberg, E.N.; Sanborn, A.L.; Kato, H.E.; Livingston, K.E.; Thorsen, T.S.; Kling, R.; Granier, S.; Gmeiner, P.; Husbands, S.M.; Traynor, J.R.; Weis, W.I.; Steyaert, J.; Dror, R.O.; Kobilka, B.K.
Deposited on : 2015-06-15
Resolution : 2.07 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](#) ①) were used in the production of this report:

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

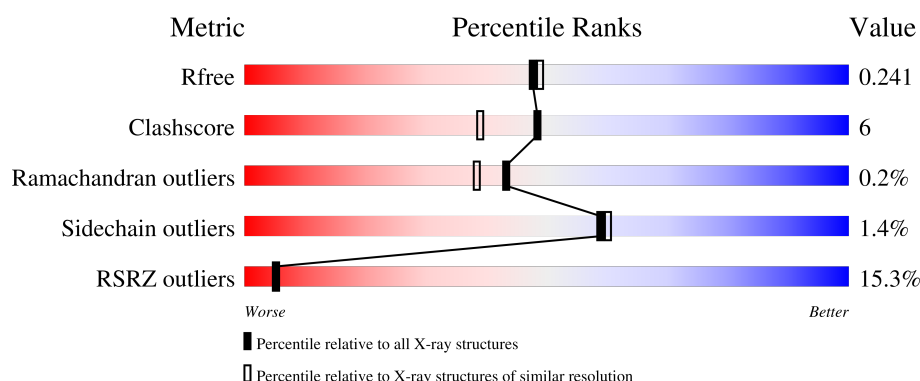
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	296	14% 89% 10% .
2	B	125	18% 81% 13% 6%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 3518 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.

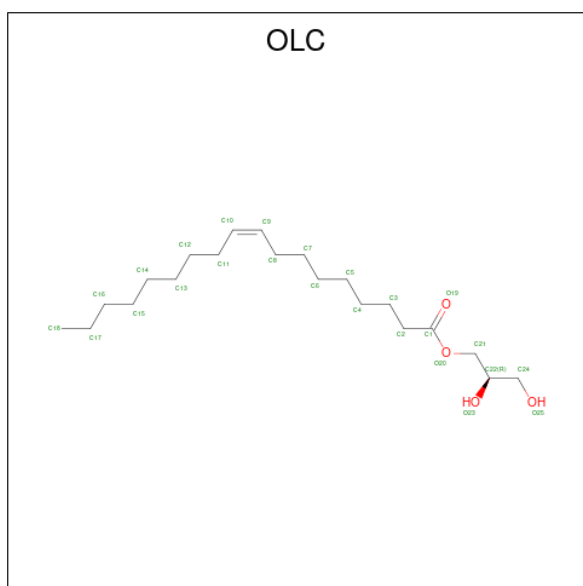
- Molecule 1 is a protein called Mu-type opioid receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	3	0
			2374	1569	381	398	26			

- Molecule 2 is a protein called Nanobody 39.

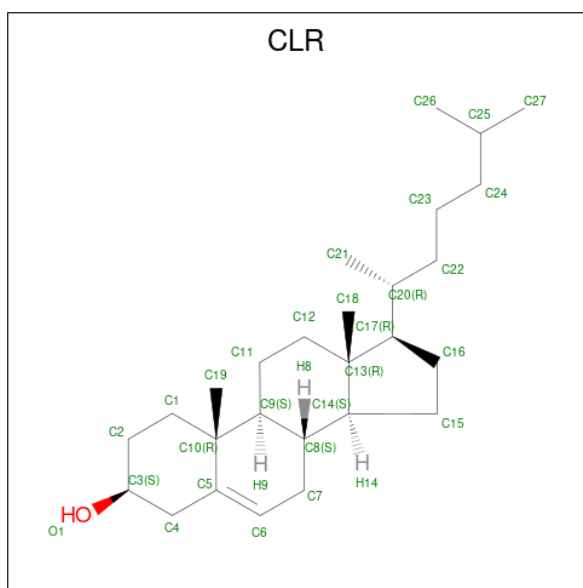
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	118	Total	C	N	O	S	0	1	0
			919	571	162	182	4			

- Molecule 3 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLC) (formula: $C_{21}H_{40}O_4$).



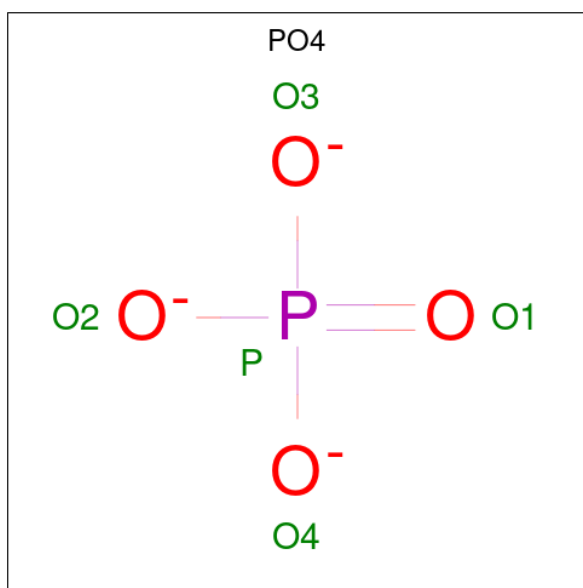
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			16	12	4		
3	A	1	Total	C	O	0	0
			18	14	4		

- Molecule 4 is CHOLESTEROL (CCD ID: CLR) (formula: $C_{27}H_{46}O$).



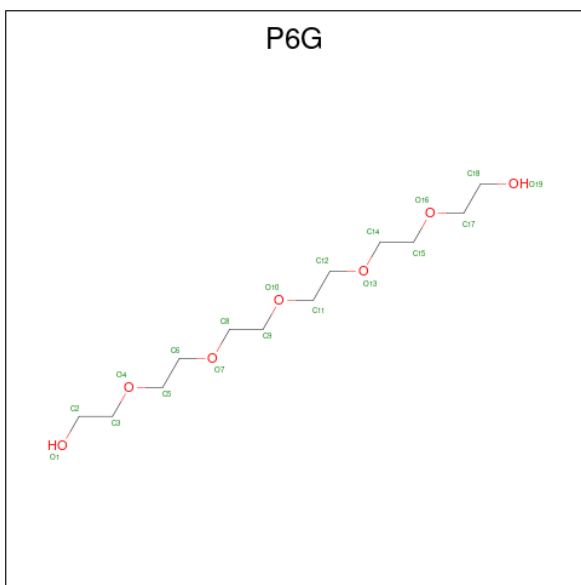
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			28	27	1		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



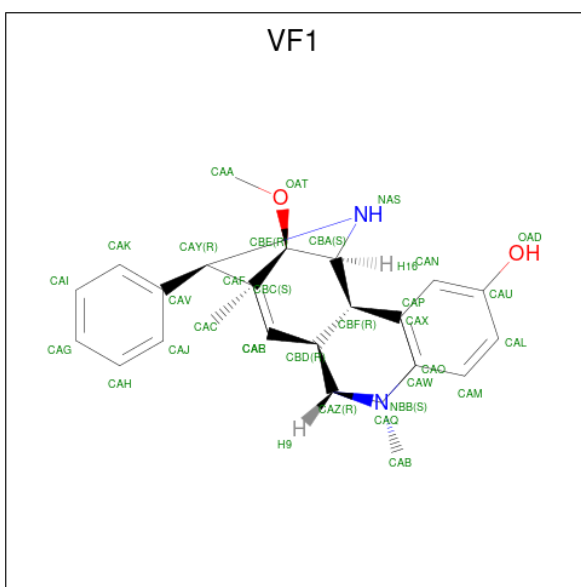
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is HEXAETHYLENE GLYCOL (CCD ID: P6G) (formula: $C_{12}H_{26}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			19	12	7		
6	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 7 is (2R,3S,3aR,5aR,6R,11bR,11cS)-3a-methoxy-3,14-dimethyl-2-phenyl-2,3,3a,6,7,11c-hexahydro-1H-6,11b-(epiminoethano)-3,5a-methanonaphtho[2,1-g]indol-10-ol (CCD ID: VF1) (formula: C₂₈H₃₂N₂O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			32	28	2	2		

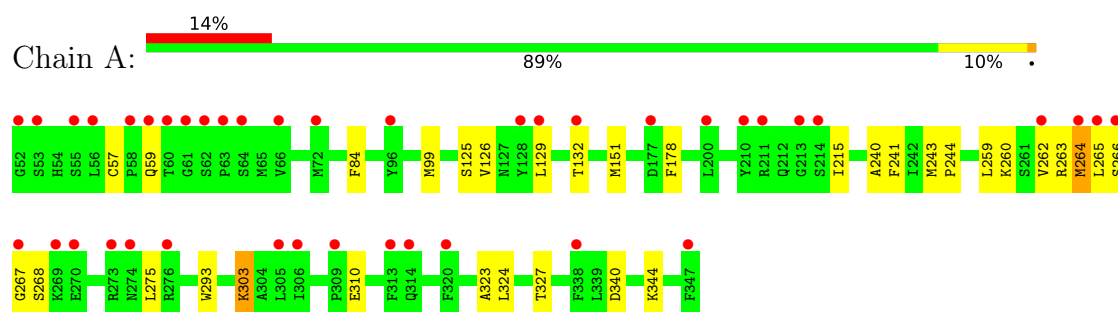
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	67	Total 67	O 67	0	0
8	B	27	Total 27	O 27	0	0

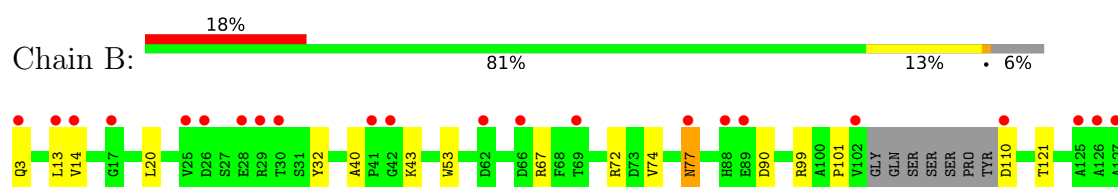
3 Residue-property plots

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: Mu-type opioid receptor



- Molecule 2: Nanobody 39



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.43Å 144.00Å 209.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.47 – 2.07 43.47 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.47-2.07) 99.8 (43.47-2.07)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.94 (at 1.79Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660, PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.199 , 0.229 (Not available) , 0.241	Depositor DCC
R_{free} test set	3144 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3518	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: VF1, PO4, YCM, OLC, P6G, CLR

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	0.22	1/2424 (0.0%)	0.34	0/3304
2	B	0.12	0/939	0.32	0/1277
All	All	0.20	1/3363 (0.0%)	0.33	0/4581

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	263	ARG	CA-C	-5.27	1.46	1.52

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2374	0	2447	28	0
2	B	919	0	865	10	0
3	A	34	0	44	0	0
4	A	28	0	46	1	0
5	A	5	0	0	0	0
6	A	32	0	43	3	0
7	A	32	0	0	0	0
8	A	67	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	27	0	0	1	0
All	All	3518	0	3445	38	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 6.

All (38) 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:A:264:MET:SD	1:A:264:MET:N	2.38	0.93
2:B:32:TYR:HD1	2:B:77[A]:ASN:HD21	1.39	0.70
1:A:132:THR:HG23	1:A:215:ILE:H	1.58	0.68
2:B:99:ARG:NH1	2:B:110:ASP:O	2.28	0.66
2:B:13:LEU:HD23	2:B:121:THR:HB	1.79	0.64
1:A:264:MET:HE1	1:A:268:SER:HA	1.82	0.61
1:A:264:MET:HE1	1:A:268:SER:CA	2.31	0.60
1:A:57:YCM:HB2	1:A:59:GLN:HE21	1.66	0.60
1:A:264:MET:CE	1:A:268:SER:HA	2.32	0.59
1:A:266:SER:HB3	2:B:101:PRO:HG2	1.85	0.59
2:B:3:GLN:N	8:B:201:HOH:O	2.35	0.58
1:A:264:MET:HE3	1:A:267:GLY:C	2.30	0.57
2:B:40:ALA:HB3	2:B:43:LYS:HB2	1.88	0.56
1:A:264:MET:O	1:A:264:MET:HE2	2.08	0.54
1:A:178:PHE:HB2	6:A:405:P6G:H31	1.90	0.53
1:A:151:MET:HE2	1:A:240:ALA:HB1	1.91	0.52
1:A:151:MET:HE3	1:A:293:TRP:CZ3	2.45	0.52
2:B:53:TRP:O	2:B:72:ARG:NH1	2.43	0.52
1:A:84:PHE:HD1	6:A:406:P6G:H121	1.74	0.50
2:B:72:ARG:NH1	2:B:74:VAL:HG12	2.28	0.49
1:A:264:MET:CE	1:A:267:GLY:C	2.87	0.47
1:A:260:LYS:HA	1:A:275:LEU:HD11	1.96	0.47
1:A:126:VAL:HA	1:A:129:LEU:HD13	1.96	0.47
1:A:125:SER:O	1:A:129:LEU:HD12	2.17	0.45
1:A:241:PHE:C	1:A:244:PRO:HD2	2.42	0.44
2:B:67:ARG:NH1	2:B:90:ASP:OD2	2.44	0.43
1:A:340:ASP:O	1:A:344:LYS:HB2	2.18	0.43
1:A:84:PHE:CD1	6:A:406:P6G:H121	2.52	0.43
2:B:14:VAL:HG11	2:B:20:LEU:HG	2.01	0.43
1:A:264:MET:CE	1:A:268:SER:N	2.81	0.43
1:A:324:LEU:O	1:A:327[B]:THR:OG1	2.33	0.43
1:A:57:YCM:HD3	1:A:57:YCM:HA	1.79	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:MET:HE3	1:A:267:GLY:O	2.18	0.42
1:A:57:YCM:HZ22	1:A:303:LYS:HE2	1.84	0.42
1:A:243:MET:HB3	1:A:244:PRO:HD3	2.02	0.41
1:A:323:ALA:O	1:A:327[B]:THR:HG23	2.20	0.41
1:A:259:LEU:O	1:A:262:VAL:HG12	2.20	0.41
4:A:403:CLR:H22	4:A:403:CLR:H192	1.75	0.41

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	296/296 (100%)	282 (95%)	13 (4%)	1 (0%)	36	30
2	B	115/125 (92%)	113 (98%)	2 (2%)	0	100	100
All	All	411/421 (98%)	395 (96%)	15 (4%)	1 (0%)	43	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	265	LEU

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	267/267 (100%)	263 (98%)	4 (2%)	57	58
2	B	97/102 (95%)	95 (98%)	2 (2%)	47	45
All	All	364/369 (99%)	358 (98%)	6 (2%)	59	55

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

Mol	Chain	Res	Type
1	A	99	MET
1	A	264	MET
1	A	303	LYS
1	A	310	GLU
2	B	77[A]	ASN
2	B	77[B]	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	127	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	YCM	A	57	1	7,9,10	1.19	0	5,10,12	0.71	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
1	YCM	A	57	1	-	5/6/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	57	YCM	N-CA-CB-SG
1	A	57	YCM	C-CA-CB-SG
1	A	57	YCM	CA-CB-SG-CD
1	A	57	YCM	CE-CD-SG-CB
1	A	57	YCM	SG-CD-CE-OZ1

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	57	YCM	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	OLC	A	402	-	17,17,24	0.97	1 (5%)	18,18,25	1.20	3 (16%)
3	OLC	A	401	-	15,15,24	0.53	0	16,16,25	0.86	0
6	P6G	A	405	-	18,18,18	0.57	0	17,17,17	0.43	0
7	VF1	A	407	-	34,38,38	2.84	8 (23%)	46,64,64	2.24	11 (23%)
4	CLR	A	403	-	31,31,31	2.79	7 (22%)	48,48,48	2.21	14 (29%)
6	P6G	A	406	-	12,12,18	0.52	0	11,11,17	0.54	0
5	PO4	A	404	-	4,4,4	0.97	0	6,6,6	0.47	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	OLC	A	402	-	-	7/17/17/24	-
3	OLC	A	401	-	-	2/15/15/24	-
6	P6G	A	405	-	-	7/16/16/16	-
7	VF1	A	407	-	-	0/7/86/86	0/1/7/7
4	CLR	A	403	-	-	1/10/68/68	0/4/4/4
6	P6G	A	406	-	-	5/10/10/16	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	407	VF1	CAE-CAF	12.86	1.51	1.32
4	A	403	CLR	C7-C8	-8.36	1.39	1.53
4	A	403	CLR	C4-C3	-6.68	1.40	1.52
4	A	403	CLR	C10-C5	-6.40	1.40	1.52
4	A	403	CLR	C4-C5	-5.77	1.40	1.51
7	A	407	VF1	CBF-CAX	-5.37	1.42	1.51
4	A	403	CLR	C7-C6	-5.17	1.39	1.50
7	A	407	VF1	CAQ-CAW	-4.94	1.42	1.51
7	A	407	VF1	CBD-CBF	-3.83	1.51	1.55
4	A	403	CLR	C6-C5	3.39	1.40	1.33
3	A	402	OLC	C9-C10	3.29	1.53	1.29
7	A	407	VF1	CBE-CBC	-2.95	1.52	1.57
7	A	407	VF1	CBD-CAE	2.82	1.55	1.50
7	A	407	VF1	CBD-CAZ	-2.82	1.52	1.54
7	A	407	VF1	CAV-CAY	2.38	1.55	1.52
4	A	403	CLR	C10-C9	-2.25	1.52	1.56

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	407	VF1	CBF-CBD-CAE	7.60	111.47	107.52
4	A	403	CLR	C3-C4-C5	7.52	124.04	112.05
7	A	407	VF1	CBD-CAZ-NBB	6.05	111.63	108.52
7	A	407	VF1	CBE-CBC-CAY	-5.70	93.73	100.24
7	A	407	VF1	CAZ-CBD-CAE	-5.36	109.16	114.02
4	A	403	CLR	C8-C7-C6	5.33	120.15	112.76
4	A	403	CLR	C4-C5-C10	4.14	121.72	116.42
4	A	403	CLR	C15-C14-C8	-3.98	112.75	119.10
4	A	403	CLR	C17-C13-C14	-3.57	96.00	100.10
4	A	403	CLR	C2-C1-C10	-3.15	106.08	112.78
4	A	403	CLR	C10-C9-C8	-2.89	108.50	112.71
7	A	407	VF1	CAP-CBF-CBA	-2.83	107.45	112.47
7	A	407	VF1	CAQ-CAZ-CBD	-2.83	108.73	111.54
4	A	403	CLR	C16-C17-C13	-2.82	100.52	103.84
3	A	402	OLC	C8-C9-C10	-2.76	109.58	130.48
4	A	403	CLR	C19-C10-C9	-2.70	108.63	111.66
7	A	407	VF1	CBF-CBD-CAZ	2.70	106.96	105.72
7	A	407	VF1	CAR-CBD-CAE	2.63	107.17	105.02
4	A	403	CLR	C13-C14-C8	-2.58	110.75	114.41
4	A	403	CLR	C16-C17-C20	-2.54	108.33	112.18
7	A	407	VF1	CAB-NBB-CAZ	-2.45	110.12	112.70
4	A	403	CLR	C1-C10-C9	2.37	111.87	108.74
3	A	402	OLC	C21-O20-C1	-2.17	109.20	117.12
3	A	402	OLC	C11-C10-C9	-2.15	109.65	126.43
7	A	407	VF1	CAV-CAY-NAS	2.12	116.69	113.48
4	A	403	CLR	C7-C6-C5	-2.11	121.45	125.02
7	A	407	VF1	CBC-CAY-CAV	-2.10	113.03	116.17
4	A	403	CLR	C10-C5-C6	-2.01	119.99	122.93

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	405	P6G	O4-C5-C6-O7
3	A	402	OLC	C2-C1-O20-C21
3	A	402	OLC	O19-C1-O20-C21
6	A	406	P6G	O7-C8-C9-O10
3	A	402	OLC	C2-C3-C4-C5
3	A	402	OLC	C3-C4-C5-C6
3	A	402	OLC	C6-C7-C8-C9
3	A	401	OLC	C2-C1-O20-C21

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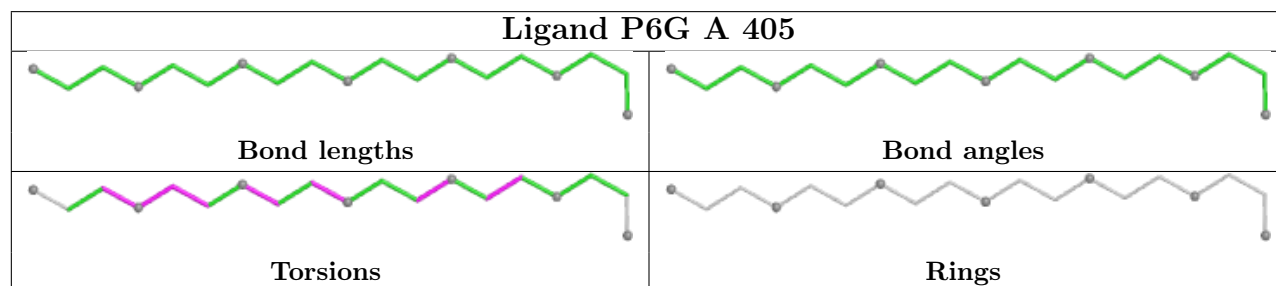
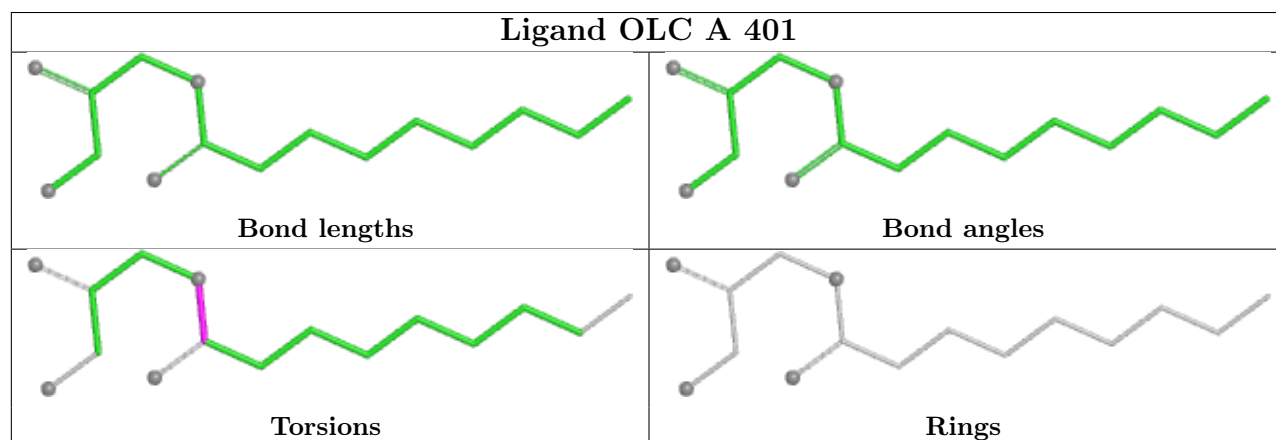
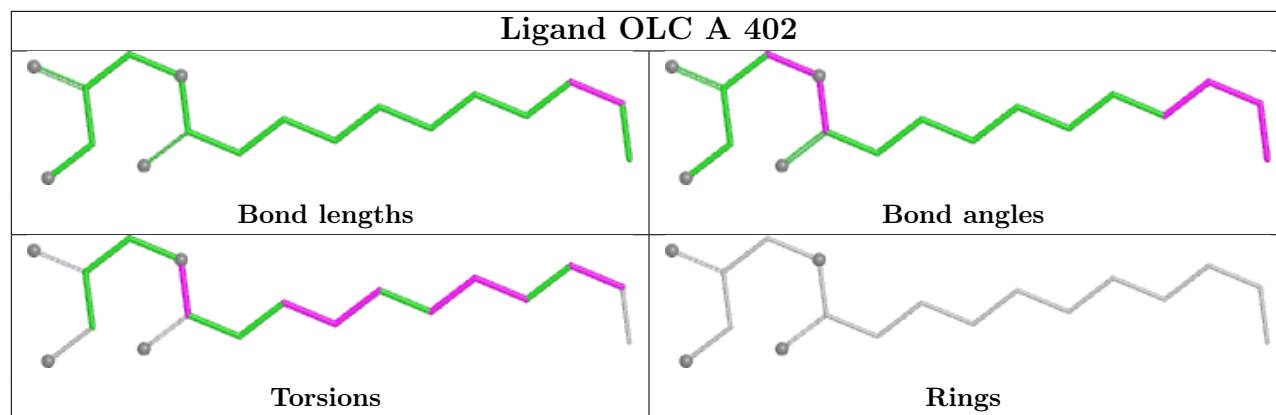
Mol	Chain	Res	Type	Atoms
6	A	405	P6G	O13-C14-C15-O16
3	A	401	OLC	O19-C1-O20-C21
3	A	402	OLC	C11-C10-C9-C8
6	A	405	P6G	C11-C12-O13-C14
6	A	405	P6G	C9-C8-O7-C6
6	A	405	P6G	C18-C17-O16-C15
6	A	405	P6G	C12-C11-O10-C9
3	A	402	OLC	C5-C6-C7-C8
4	A	403	CLR	C22-C23-C24-C25
6	A	406	P6G	O10-C11-C12-O13
6	A	406	P6G	O13-C14-C15-O16
6	A	406	P6G	C11-C12-O13-C14
6	A	406	P6G	C12-C11-O10-C9
6	A	405	P6G	C14-C15-O16-C17

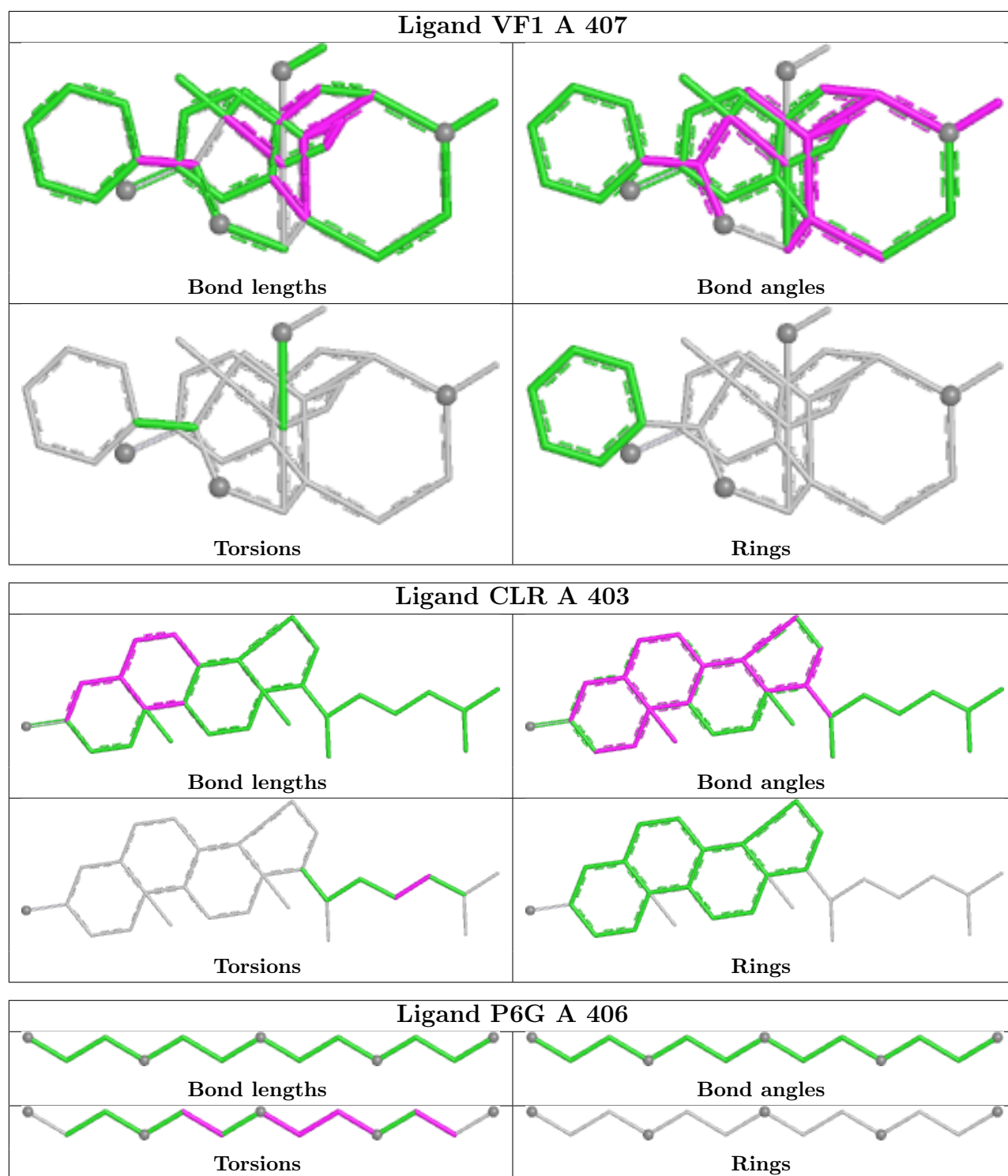
There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	405	P6G	1	0
4	A	403	CLR	1	0
6	A	406	P6G	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

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	295/296 (99%)	0.73	41 (13%) 6 6	19, 46, 110, 159	3 (1%)
2	B	118/125 (94%)	1.11	22 (18%) 3 3	27, 57, 97, 121	1 (0%)
All	All	413/421 (98%)	0.84	63 (15%) 5 5	19, 50, 109, 159	4 (0%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	102	VAL	7.0
1	A	265	LEU	5.8
1	A	347	PHE	5.3
2	B	66	ASP	5.1
1	A	270	GLU	4.8
1	A	129	LEU	4.6
2	B	3	GLN	4.5
1	A	177	ASP	4.2
2	B	13	LEU	4.2
1	A	213	GLY	4.0
2	B	77[A]	ASN	3.9
1	A	264	MET	3.9
1	A	210	TYR	3.8
1	A	58	PRO	3.5
2	B	14	VAL	3.3
2	B	29	ARG	3.3
1	A	63	PRO	3.3
1	A	211	ARG	3.2
1	A	267	GLY	3.2
1	A	313	PHE	3.1
1	A	266	SER	3.1
1	A	269	LYS	3.0
1	A	262	VAL	3.0
2	B	126	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	60	THR	2.9
2	B	28	GLU	2.8
2	B	41	PRO	2.8
2	B	88	HIS	2.7
1	A	52	GLY	2.7
2	B	89	GLU	2.6
2	B	62	ASP	2.6
2	B	127	ALA	2.5
2	B	25	VAL	2.5
1	A	61	GLY	2.5
1	A	273	ARG	2.5
1	A	66	VAL	2.4
1	A	306	ILE	2.4
1	A	132	THR	2.4
1	A	56	LEU	2.4
1	A	72	MET	2.4
1	A	276	ARG	2.4
1	A	53	SER	2.3
1	A	214	SER	2.3
2	B	26	ASP	2.3
1	A	55	SER	2.2
1	A	62	SER	2.2
2	B	42	GLY	2.2
1	A	320	PHE	2.2
1	A	309	PRO	2.2
2	B	125	ALA	2.2
1	A	338	PHE	2.2
2	B	30	THR	2.2
1	A	59	GLN	2.1
1	A	314	GLN	2.1
1	A	305	LEU	2.1
1	A	274	ASN	2.1
1	A	128	TYR	2.1
2	B	17	GLY	2.1
1	A	64	SER	2.1
1	A	96	TYR	2.1
2	B	110	ASP	2.0
1	A	200	LEU	2.0
2	B	69	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	YCM	A	57	10/11	0.73	0.16	90,100,111,113	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

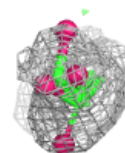
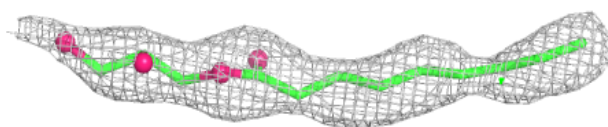
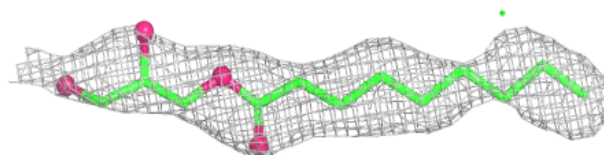
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	PO4	A	404	5/5	0.75	0.12	110,118,120,121	0
3	OLC	A	401	16/25	0.80	0.20	70,88,92,95	0
6	P6G	A	406	13/19	0.85	0.18	83,89,92,92	0
4	CLR	A	403	28/28	0.86	0.19	59,85,96,105	0
6	P6G	A	405	19/19	0.87	0.21	71,77,89,90	0
3	OLC	A	402	18/25	0.88	0.19	73,80,96,97	0
7	VF1	A	407	32/32	0.95	0.07	30,39,45,48	0

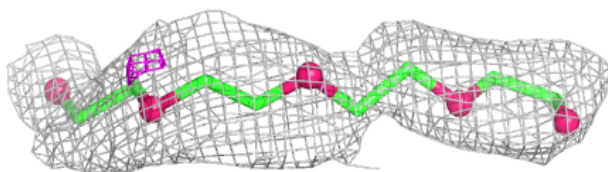
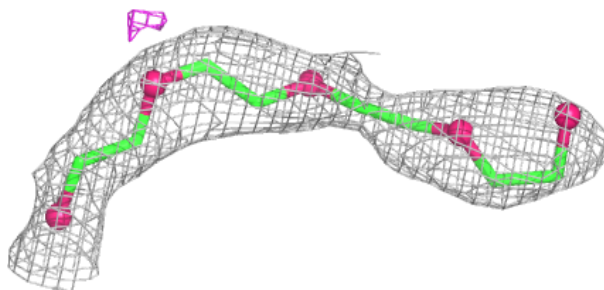
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around OLC A 401:

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

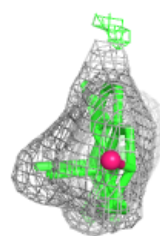
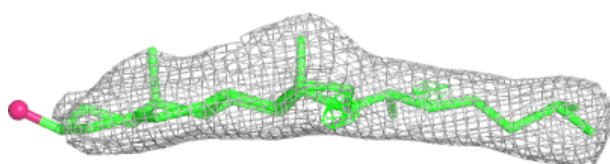
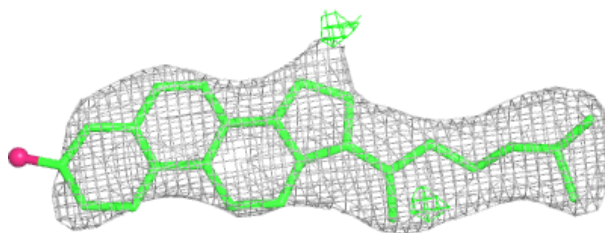
**Electron density around P6G A 406:**

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

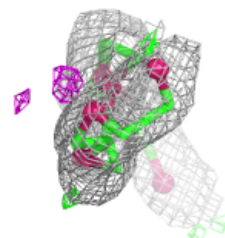
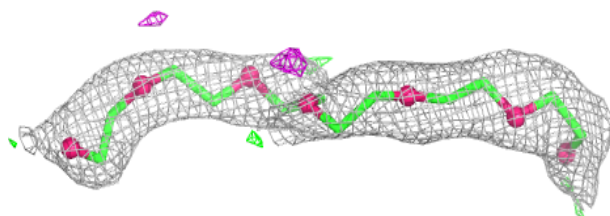
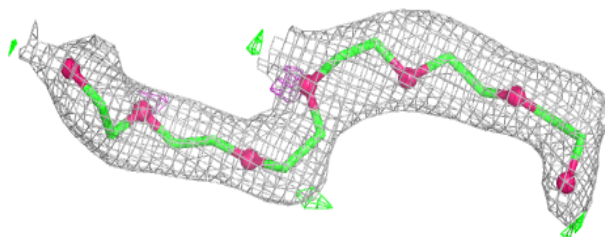


Electron density around CLR A 403:

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

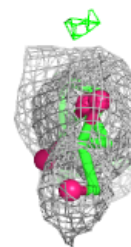
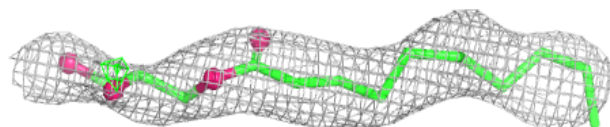
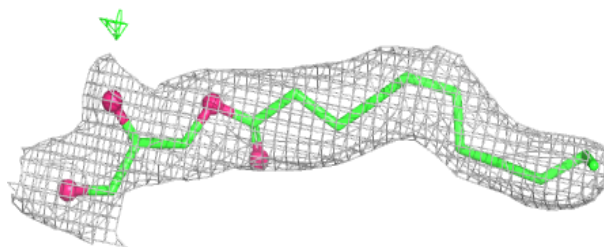
**Electron density around P6G A 405:**

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



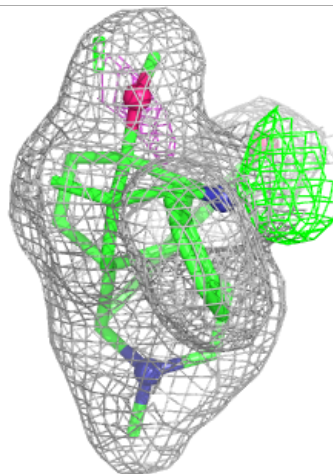
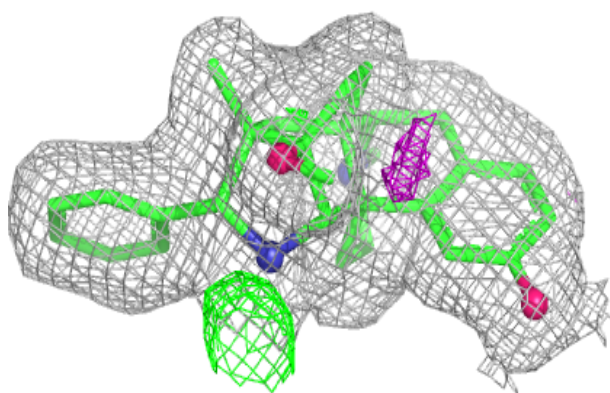
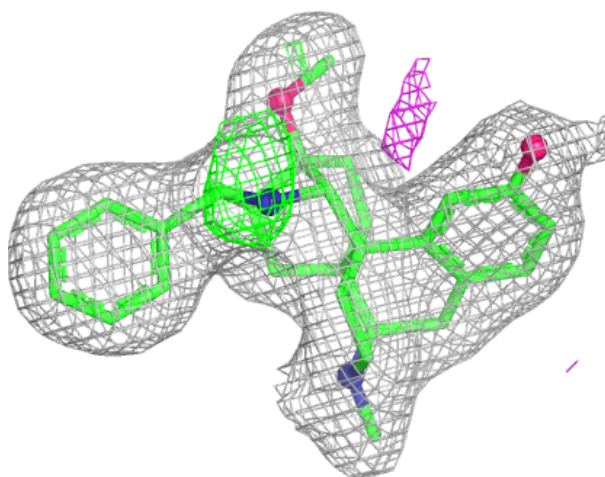
Electron density around OLC A 402:

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



Electron density around VF1 A 407:

$2mF_o-DF_c$ (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.