



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 07:31 am BST

PDB ID : 4RDY
Title : Crystal structure of VmoLac bound to 3-oxo-C10 AHL
Authors : Hiblot, J.; Bzdrenga, J.; Champion, C.; Gotthard, G.; Gonzalez, D.; Chabriere, E.; Elias, M.
Deposited on : 2014-09-20
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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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

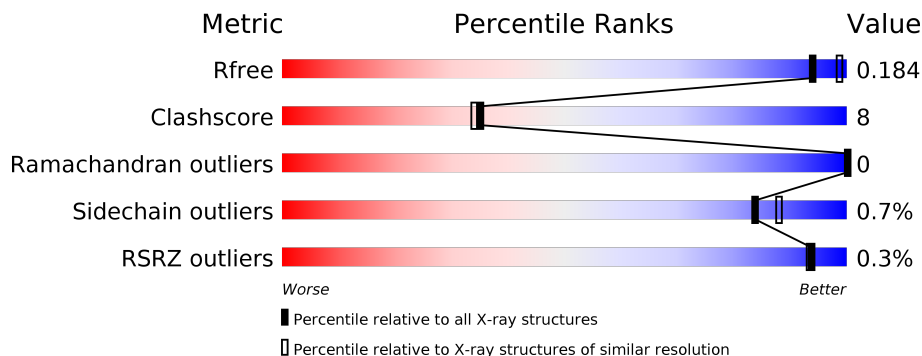
1 Overall quality at a glance

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 $\leq 5\%$. 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	316	 89% 9%
1	B	316	 87% 12%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	3M5	B	403	-	-	X	-
4	GOL	A	404	-	-	X	-
4	GOL	A	406	-	X	-	-
4	GOL	A	408	-	-	X	-
4	GOL	A	411	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5831 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 Parathion hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	313	2503	1604	432	460	7	0	3	0
1	B	314	2539	1632	431	468	8	0	9	0

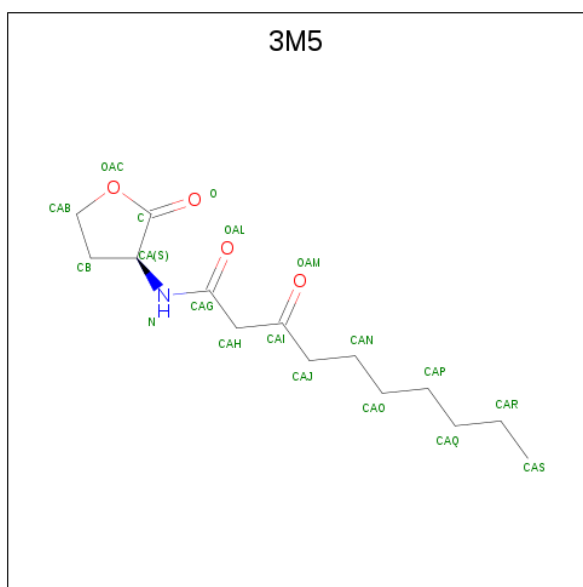
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP F0QXN6
B	0	GLY	-	EXPRESSION TAG	UNP F0QXN6

- Molecule 2 is COBALT (II) ION (three-letter code: CO) (formula: Co).

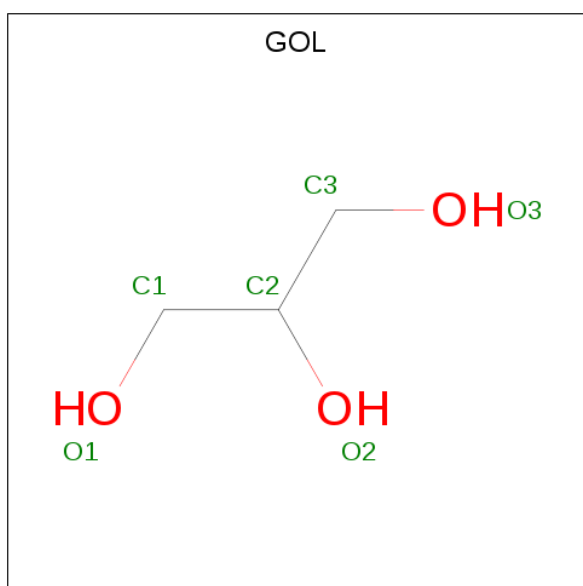
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Co	0	0
			2	2		
2	A	2	Total	Co	0	0
			2	2		

- Molecule 3 is 3-oxo-N-[(3S)-2-oxotetrahydrofuran-3-yl]decanamide (three-letter code: 3M5) (formula: C₁₄H₂₃NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			19	14	1	4		
3	B	1	Total	C	N	O	0	0
			19	14	1	4		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



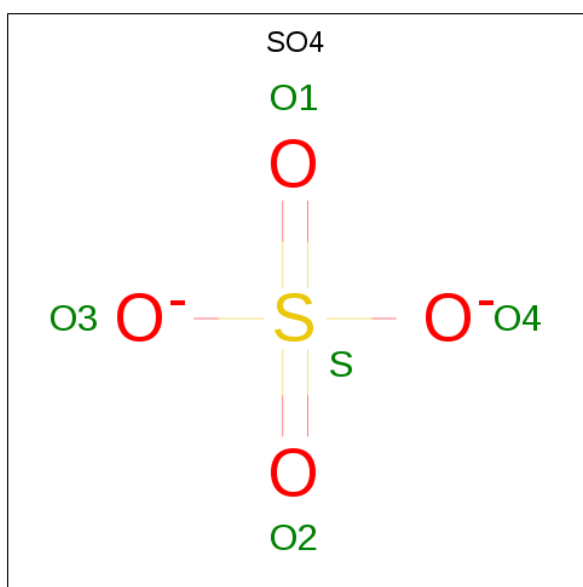
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

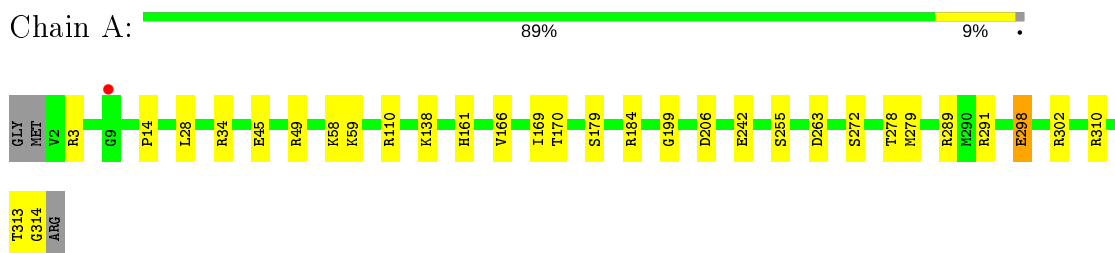
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	330	Total	O	0	0
			330	330		
6	B	329	Total	O	0	0
			329	329		

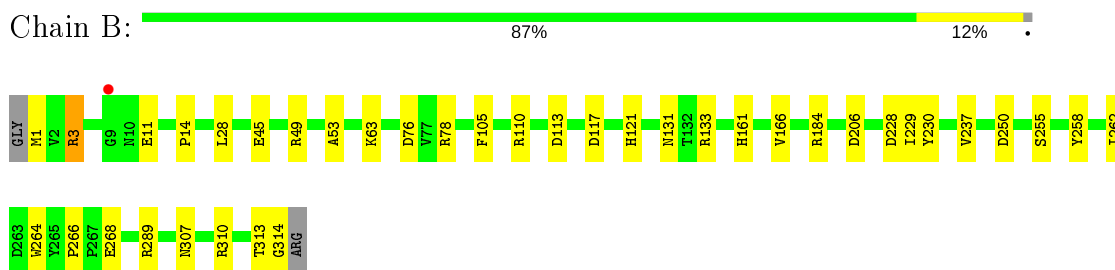
3 Residue-property plots [i](#)

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: Parathion hydrolase



- Molecule 1: Parathion hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	174.74Å 174.74Å 61.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.75 – 2.00 47.75 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.75-2.00) 99.8 (47.75-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 2.00Å)	Xtrriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.139 , 0.174 0.151 , 0.184	Depositor DCC
R_{free} test set	3630 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtrriage
Anisotropy	0.037	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5831	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% 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: GOL, 3M5, CO, SO4, KCX

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	1.09	3/2549 (0.1%)	1.02	5/3455 (0.1%)
1	B	1.09	1/2603 (0.0%)	1.04	7/3527 (0.2%)
All	All	1.09	4/5152 (0.1%)	1.03	12/6982 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	242	GLU	CD-OE1	5.99	1.32	1.25
1	B	258	TYR	CG-CD2	-5.94	1.31	1.39
1	A	272	SER	CB-OG	-5.16	1.35	1.42
1	A	298	GLU	CD-OE2	-5.04	1.20	1.25

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	291	ARG	NE-CZ-NH1	6.95	123.78	120.30
1	B	110	ARG	NE-CZ-NH2	-6.71	116.95	120.30
1	A	263	ASP	CB-CG-OD1	6.24	123.92	118.30
1	B	206	ASP	CB-CG-OD1	6.17	123.85	118.30
1	A	206	ASP	CB-CG-OD1	5.86	123.57	118.30
1	B	78	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	B	228	ASP	CB-CG-OD1	5.45	123.20	118.30
1	A	272	SER	N-CA-CB	-5.41	102.38	110.50
1	A	263	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	B	76	ASP	CB-CG-OD1	5.38	123.14	118.30
1	B	250	ASP	CB-CG-OD2	-5.37	113.46	118.30
1	B	117	ASP	CB-CG-OD1	5.20	122.98	118.30

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	2503	0	2541	34	0
1	B	2539	0	2596	35	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	19	0	23	8	0
3	B	19	0	23	9	0
4	A	48	0	64	20	0
4	B	30	0	39	9	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
6	A	330	0	0	15	4
6	B	329	0	0	19	4
All	All	5831	0	5286	86	4

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 8.

All (86) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:ASN:HB3	6:B:815:HOH:O	1.58	1.03
1:B:28:LEU:HD11	3:B:403:3M5:H3	1.53	0.90
1:A:59:LYS:HD2	6:A:613:HOH:O	1.74	0.85
1:B:28:LEU:CD1	3:B:403:3M5:H3	2.06	0.85
1:B:268[A]:GLU:HG2	6:B:820:HOH:O	1.79	0.82
1:A:161:HIS:HD2	1:A:166:VAL:O	1.65	0.79
1:B:161:HIS:HD2	1:B:166:VAL:O	1.64	0.79
1:A:28:LEU:HD11	3:A:403:3M5:H3	1.64	0.78
1:B:45[A]:GLU:HG3	1:B:49:ARG:NH1	1.98	0.77
3:A:403:3M5:H2	3:A:403:3M5:OAL	1.87	0.74
1:A:298:GLU:OE1	6:A:787:HOH:O	2.06	0.73
3:A:403:3M5:OAL	3:A:403:3M5:CB	2.35	0.73
1:A:45:GLU:HG3	1:A:49:ARG:NH1	2.07	0.70
1:B:63:LYS:HE2	6:B:823:HOH:O	1.90	0.70
1:A:3:ARG:HG3	6:A:823:HOH:O	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:403:3M5:OAC	6:A:830:HOH:O	2.11	0.68
3:B:403:3M5:OAL	3:B:403:3M5:H2	1.92	0.68
3:B:403:3M5:OAL	3:B:403:3M5:CB	2.41	0.67
1:A:310[A]:ARG:NH2	4:A:409:GOL:O2	2.29	0.65
1:B:1:MET:N	6:B:811:HOH:O	2.25	0.63
1:A:310[A]:ARG:HD2	6:A:781:HOH:O	1.99	0.62
1:A:28:LEU:CD1	3:A:403:3M5:H3	2.30	0.61
1:B:14:PRO:HG2	1:B:313:THR:HG22	1.82	0.61
1:B:310[A]:ARG:NH2	4:B:404:GOL:O2	2.34	0.60
1:A:110:ARG:HE	4:A:408:GOL:H32	1.66	0.60
1:A:184:ARG:HD3	4:A:405:GOL:O3	2.01	0.60
1:A:314:GLY:C	6:A:790:HOH:O	2.40	0.60
4:B:407:GOL:H2	6:B:709:HOH:O	2.01	0.60
1:A:45:GLU:OE1	6:A:803:HOH:O	2.17	0.58
1:B:310[A]:ARG:HH22	4:B:404:GOL:C2	2.16	0.58
4:B:408:GOL:O2	6:B:576:HOH:O	1.98	0.57
3:A:403:3M5:CAB	3:A:403:3M5:OAL	2.53	0.57
4:A:406:GOL:H11	1:B:105:PHE:CE2	2.39	0.57
1:B:28:LEU:CD1	3:B:403:3M5:CB	2.82	0.56
1:B:310[B]:ARG:HD3	6:B:824:HOH:O	2.05	0.56
1:A:161:HIS:HE1	6:A:536:HOH:O	1.89	0.55
1:A:34:ARG:HE	4:A:411:GOL:C2	2.20	0.55
3:A:403:3M5:H4	3:A:403:3M5:OAL	2.08	0.54
1:A:58:LYS:NZ	6:A:775:HOH:O	2.41	0.54
1:B:229:ILE:HG23	6:B:789:HOH:O	2.07	0.54
3:B:403:3M5:OAL	3:B:403:3M5:H4	2.08	0.54
1:A:34:ARG:HG3	4:A:411:GOL:H32	1.90	0.54
1:B:45[A]:GLU:CG	1:B:49:ARG:NH1	2.69	0.54
1:B:11:GLU:OE2	1:B:133:ARG:NH2	2.41	0.53
1:A:110:ARG:HH21	4:A:408:GOL:H12	1.72	0.53
1:B:161:HIS:HE1	6:B:538:HOH:O	1.91	0.53
4:A:410:GOL:H11	6:A:698:HOH:O	2.09	0.52
1:B:121:HIS:HE1	6:B:636:HOH:O	1.93	0.52
3:B:403:3M5:OAL	3:B:403:3M5:CAB	2.58	0.52
4:A:411:GOL:C1	6:A:685:HOH:O	2.58	0.52
1:B:237:VAL:HG11	1:B:289:ARG:HD2	1.92	0.51
1:B:314:GLY:C	6:B:794:HOH:O	2.49	0.51
1:B:53:ALA:HA	4:B:405:GOL:H12	1.94	0.50
3:B:403:3M5:C	6:B:829:HOH:O	2.61	0.49
1:A:28:LEU:CD1	3:A:403:3M5:CB	2.91	0.48
4:B:407:GOL:H32	6:B:709:HOH:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:408:GOL:C1	6:A:795:HOH:O	2.61	0.48
4:A:408:GOL:H11	1:B:266:PRO:HB3	1.95	0.48
1:A:302:ARG:HD2	6:A:617:HOH:O	2.14	0.48
1:A:184:ARG:CD	4:A:405:GOL:O3	2.62	0.48
1:B:113:ASP:HB2	6:B:825:HOH:O	2.14	0.47
1:A:289[B]:ARG:NH2	6:A:705:HOH:O	2.15	0.47
1:A:110:ARG:NE	4:A:408:GOL:H32	2.30	0.47
1:B:184:ARG:HD2	6:B:800:HOH:O	2.14	0.46
6:A:561:HOH:O	4:B:406:GOL:H12	2.14	0.46
1:A:138:KCX:HA	1:A:169:ILE:O	2.15	0.46
1:B:229:ILE:HD12	1:B:230:TYR:CD2	2.50	0.45
1:B:289:ARG:HG2	6:B:819:HOH:O	2.15	0.45
1:B:3:ARG:HA	1:B:3:ARG:HD2	1.79	0.45
1:A:14:PRO:HG2	1:A:313:THR:HG22	1.99	0.45
1:A:161:HIS:CD2	1:A:166:VAL:O	2.57	0.44
1:B:184:ARG:CD	6:B:800:HOH:O	2.65	0.44
1:B:310[A]:ARG:NH2	4:B:404:GOL:C2	2.80	0.44
1:B:3:ARG:HG3	6:B:787:HOH:O	2.17	0.44
1:A:170:THR:O	1:A:199:GLY:HA3	2.18	0.43
1:A:278:THR:HB	4:A:404:GOL:C3	2.48	0.42
1:A:110:ARG:HE	4:A:408:GOL:H12	1.83	0.42
1:B:264:TRP:CH2	3:B:403:3M5:H1	2.55	0.42
1:B:262:ILE:HG22	4:B:407:GOL:H11	2.01	0.42
1:A:34:ARG:HE	4:A:411:GOL:C1	2.33	0.41
1:A:34:ARG:HE	4:A:411:GOL:H2	1.84	0.41
1:A:279:MET:H	4:A:404:GOL:H31	1.86	0.41
1:A:279:MET:H	4:A:404:GOL:C3	2.33	0.41
1:B:3:ARG:CG	6:B:787:HOH:O	2.69	0.41
1:A:278:THR:HB	4:A:404:GOL:H32	2.03	0.40
1:B:307:ASN:OD1	1:B:310[B]:ARG:NH2	2.48	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:812:HOH:O	6:B:812:HOH:O[3_455]	1.82	0.38
6:A:815:HOH:O	6:B:699:HOH:O[6_665]	2.05	0.15
6:A:791:HOH:O	6:B:574:HOH:O[6_665]	2.09	0.11
6:A:636:HOH:O	6:B:806:HOH:O[6_665]	2.14	0.06

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	313/316 (99%)	308 (98%)	5 (2%)	0	100	100
1	B	320/316 (101%)	315 (98%)	5 (2%)	0	100	100
All	All	633/632 (100%)	623 (98%)	10 (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	271/270 (100%)	269 (99%)	2 (1%)	84	88
1	B	278/270 (103%)	276 (99%)	2 (1%)	84	88
All	All	549/540 (102%)	545 (99%)	4 (1%)	84	88

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

Mol	Chain	Res	Type
1	A	179	SER
1	A	255	SER
1	B	3	ARG
1	B	255	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	90	ASN
1	A	161	HIS
1	A	251	GLN
1	B	36	ASN
1	B	90	ASN
1	B	161	HIS
1	B	251	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	KCX	B	138	1,2	7,11,12	1.34	1 (14%)	4,12,14	1.00	0
1	KCX	A	138	1,2	7,11,12	0.74	0	4,12,14	1.31	1 (25%)

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	KCX	B	138	1,2	-	1/7/10/12	-
1	KCX	A	138	1,2	-	0/7/10/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	138	KCX	CB-CA	-3.13	1.49	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	KCX	CE-NZ-CX	2.56	127.28	122.95

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	138	KCX	O-C-CA-CB

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	138	KCX	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 4 are monoatomic - leaving 17 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
4	GOL	B	405	-	5,5,5	1.35	1 (20%)	5,5,5	0.69	0
4	GOL	B	406	-	5,5,5	0.49	0	5,5,5	1.92	2 (40%)
4	GOL	A	404	-	5,5,5	0.94	0	5,5,5	1.57	1 (20%)
4	GOL	A	409	-	5,5,5	0.76	0	5,5,5	0.84	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	A	412	-	4,4,4	0.64	0	6,6,6	0.78	0
4	GOL	A	407	-	5,5,5	0.28	0	5,5,5	0.58	0
4	GOL	B	407	-	5,5,5	0.60	0	5,5,5	0.50	0
3	3M5	B	403	2	19,19,19	1.99	3 (15%)	18,23,23	2.66	6 (33%)
3	3M5	A	403	2	19,19,19	1.97	5 (26%)	18,23,23	2.73	6 (33%)
4	GOL	A	410	-	5,5,5	0.66	0	5,5,5	0.87	0
4	GOL	A	408	-	5,5,5	0.80	0	5,5,5	2.17	1 (20%)
5	SO4	B	409	-	4,4,4	0.57	0	6,6,6	0.40	0
4	GOL	B	404	-	5,5,5	0.68	0	5,5,5	1.38	1 (20%)
4	GOL	B	408	-	5,5,5	0.39	0	5,5,5	0.88	0
4	GOL	A	406	-	5,5,5	1.08	0	5,5,5	1.45	2 (40%)
4	GOL	A	405	-	5,5,5	0.36	0	5,5,5	1.05	0
4	GOL	A	411	-	5,5,5	0.93	0	5,5,5	1.90	1 (20%)

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
4	GOL	B	405	-	-	2/4/4/4	-
4	GOL	B	406	-	-	0/4/4/4	-
4	GOL	A	404	-	-	0/4/4/4	-
4	GOL	A	409	-	-	0/4/4/4	-
4	GOL	A	407	-	-	4/4/4/4	-
4	GOL	B	407	-	-	2/4/4/4	-
3	3M5	B	403	2	-	4/15/25/25	0/1/1/1
3	3M5	A	403	2	-	3/15/25/25	0/1/1/1
4	GOL	A	410	-	-	2/4/4/4	-
4	GOL	A	408	-	-	4/4/4/4	-
4	GOL	B	404	-	-	4/4/4/4	-
4	GOL	B	408	-	-	0/4/4/4	-
4	GOL	A	406	-	-	4/4/4/4	-
4	GOL	A	405	-	-	2/4/4/4	-
4	GOL	A	411	-	-	2/4/4/4	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	3M5	CA-C	-6.99	1.37	1.52
3	A	403	3M5	CA-C	-6.05	1.39	1.52
3	B	403	3M5	OAC-CAB	-3.27	1.38	1.46
3	A	403	3M5	CA-N	2.92	1.52	1.45
3	A	403	3M5	OAL-CAG	-2.40	1.18	1.23
3	A	403	3M5	OAC-CAB	-2.36	1.40	1.46
4	B	405	GOL	C1-C2	-2.31	1.42	1.51
3	B	403	3M5	OAL-CAG	-2.12	1.19	1.23
3	A	403	3M5	CAH-CAG	-2.07	1.49	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	403	3M5	OAC-CAB-CB	-8.61	91.63	105.73
3	A	403	3M5	OAC-CAB-CB	-8.46	91.87	105.73
4	A	408	GOL	O2-C2-C1	-4.20	90.64	109.12
4	A	411	GOL	O3-C3-C2	3.49	126.94	110.20
3	A	403	3M5	CAB-OAC-C	-3.47	107.18	110.39
3	A	403	3M5	CAH-CAG-N	3.47	122.23	116.31
3	A	403	3M5	CAO-CAN-CAJ	-3.40	100.98	113.19
3	B	403	3M5	CB-CA-N	-3.10	108.00	114.96
3	A	403	3M5	OAL-CAG-CAH	-2.98	116.47	121.33
3	B	403	3M5	CAB-OAC-C	-2.97	107.65	110.39
4	A	404	GOL	C3-C2-C1	-2.85	100.63	111.70
3	B	403	3M5	OAC-C-O	2.83	124.36	121.42
4	B	406	GOL	O2-C2-C1	-2.47	98.25	109.12
3	B	403	3M5	OAL-CAG-CAH	-2.42	117.38	121.33
4	B	406	GOL	C3-C2-C1	-2.26	102.91	111.70
3	A	403	3M5	CAG-CAH-CAI	-2.24	104.34	113.71
4	A	406	GOL	O3-C3-C2	2.18	120.65	110.20
4	B	404	GOL	O2-C2-C3	-2.08	99.97	109.12
3	B	403	3M5	CAO-CAN-CAJ	-2.08	105.72	113.19
4	A	406	GOL	O1-C1-C2	2.02	119.89	110.20

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	405	GOL	O1-C1-C2-C3
3	B	403	3M5	CB-CA-N-CAG
3	A	403	3M5	CB-CA-N-CAG
4	A	410	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
4	A	408	GOL	O1-C1-C2-C3
4	A	408	GOL	C1-C2-C3-O3
4	B	404	GOL	O1-C1-C2-C3
4	B	404	GOL	C1-C2-C3-O3
4	A	406	GOL	C1-C2-C3-O3
4	A	411	GOL	C1-C2-C3-O3
4	B	405	GOL	O1-C1-C2-O2
4	A	408	GOL	O1-C1-C2-O2
4	B	404	GOL	O1-C1-C2-O2
3	B	403	3M5	CAJ-CAN-CAO-CAP
4	A	407	GOL	O1-C1-C2-C3
4	A	407	GOL	C1-C2-C3-O3
4	B	407	GOL	O1-C1-C2-C3
4	A	405	GOL	O1-C1-C2-C3
4	B	404	GOL	O2-C2-C3-O3
4	A	405	GOL	O1-C1-C2-O2
4	A	411	GOL	O2-C2-C3-O3
3	B	403	3M5	CAO-CAP-CAQ-CAR
4	A	407	GOL	O1-C1-C2-O2
4	A	410	GOL	O2-C2-C3-O3
4	A	408	GOL	O2-C2-C3-O3
4	A	407	GOL	O2-C2-C3-O3
4	A	406	GOL	O2-C2-C3-O3
4	A	406	GOL	O1-C1-C2-O2
3	A	403	3M5	CAN-CAO-CAP-CAQ
4	A	406	GOL	O1-C1-C2-C3
4	B	407	GOL	O1-C1-C2-O2
3	B	403	3M5	CAP-CAQ-CAR-CAS
3	A	403	3M5	CAG-CAH-CAI-CAJ

There are no ring outliers.

14 monomers are involved in 46 short contacts:

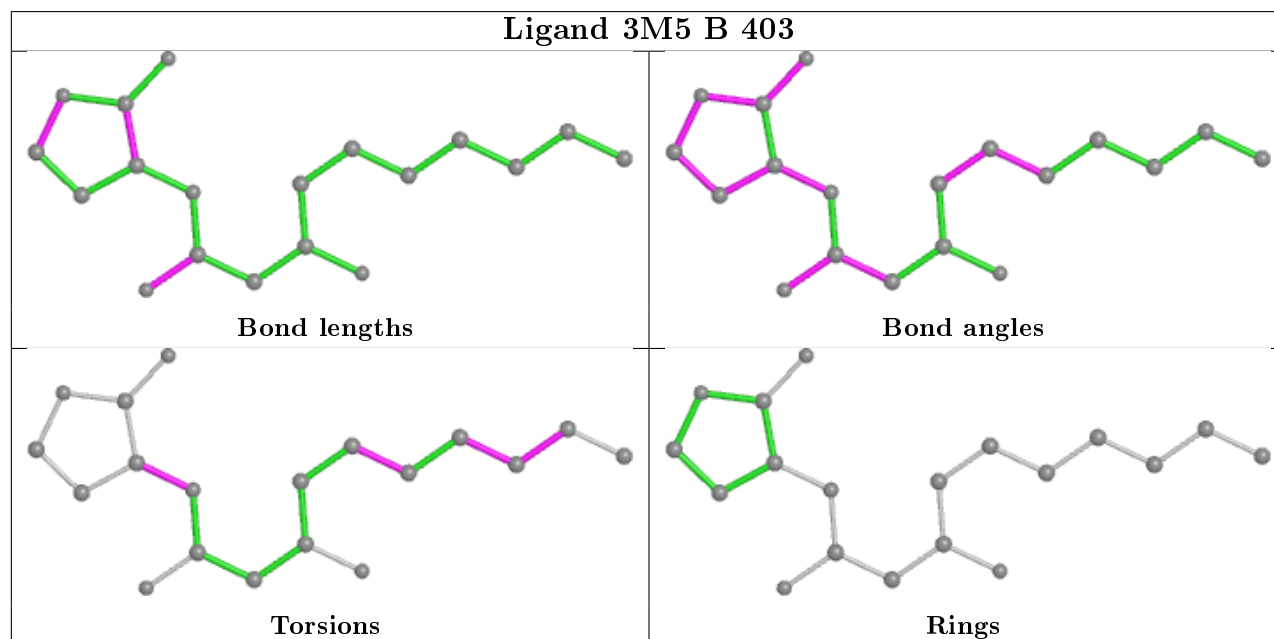
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	405	GOL	1	0
4	B	406	GOL	1	0
4	A	404	GOL	4	0
4	A	409	GOL	1	0
4	B	407	GOL	3	0
3	B	403	3M5	9	0
3	A	403	3M5	8	0
4	A	410	GOL	1	0

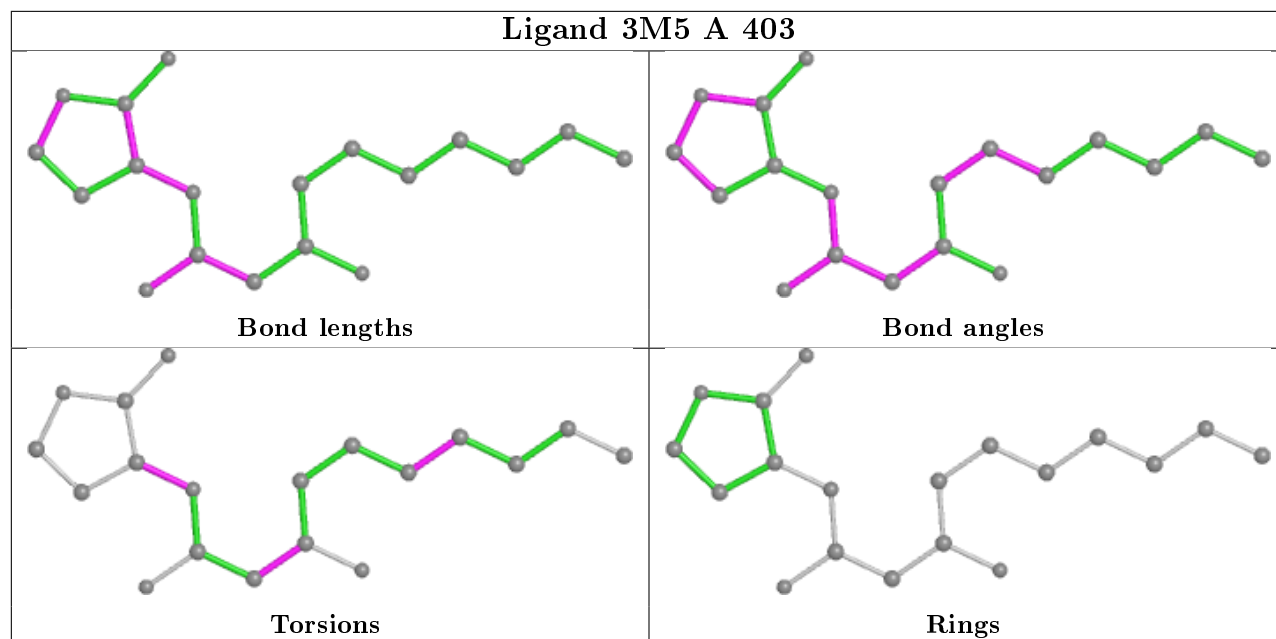
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	408	GOL	6	0
4	B	404	GOL	3	0
4	B	408	GOL	1	0
4	A	406	GOL	1	0
4	A	405	GOL	2	0
4	A	411	GOL	5	0

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 [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	312/316 (98%)	-0.67	1 (0%) 94 93	13, 20, 36, 51	0
1	B	313/316 (99%)	-0.70	1 (0%) 94 93	13, 20, 38, 67	0
All	All	625/632 (98%)	-0.69	2 (0%) 94 93	13, 20, 37, 67	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	9	GLY	2.8
1	B	9	GLY	2.3

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	KCX	B	138	12/13	0.98	0.09	12,14,17,18	0
1	KCX	A	138	12/13	0.99	0.09	13,14,15,15	0

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,

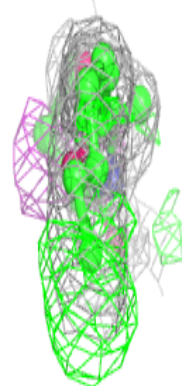
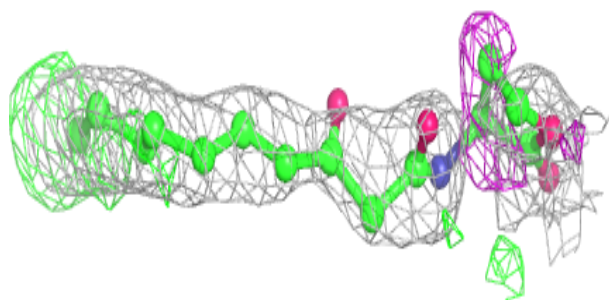
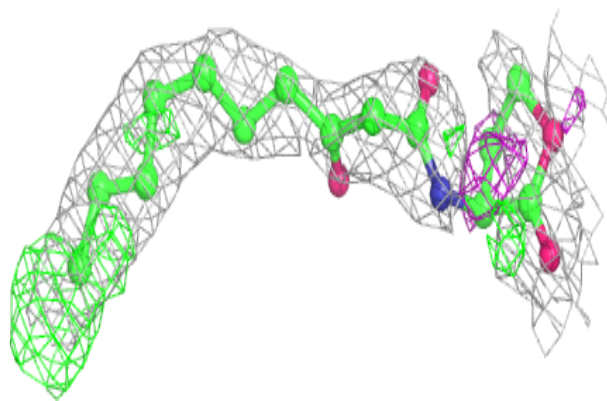
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	408	6/6	0.76	0.20	52,59,62,62	0
3	3M5	A	403	19/19	0.85	0.26	25,33,49,61	19
3	3M5	B	403	19/19	0.85	0.29	29,42,55,69	19
4	GOL	A	411	6/6	0.86	0.15	38,49,56,60	0
4	GOL	A	408	6/6	0.87	0.23	40,51,59,59	0
4	GOL	A	410	6/6	0.88	0.20	38,49,53,55	0
4	GOL	A	405	6/6	0.89	0.28	55,59,63,67	0
4	GOL	A	407	6/6	0.92	0.28	60,66,70,75	0
4	GOL	A	409	6/6	0.92	0.19	33,43,48,48	0
5	SO4	B	409	5/5	0.93	0.24	64,64,74,82	0
4	GOL	B	407	6/6	0.94	0.15	38,41,59,66	0
4	GOL	B	406	6/6	0.94	0.12	37,46,49,50	0
4	GOL	B	404	6/6	0.94	0.16	37,40,43,52	0
4	GOL	A	404	6/6	0.95	0.19	26,34,43,47	0
4	GOL	A	406	6/6	0.96	0.13	27,40,41,51	0
5	SO4	A	412	5/5	0.96	0.16	41,47,58,61	0
4	GOL	B	405	6/6	0.96	0.13	27,32,41,47	0
2	CO	A	402	1/1	1.00	0.07	18,18,18,18	0
2	CO	A	401	1/1	1.00	0.07	17,17,17,17	0
2	CO	B	401	1/1	1.00	0.09	16,16,16,16	0
2	CO	B	402	1/1	1.00	0.07	17,17,17,17	0

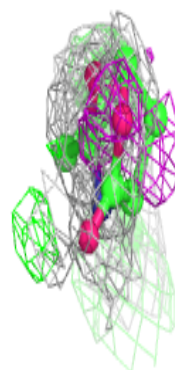
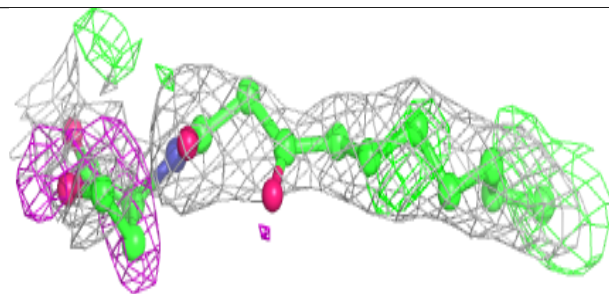
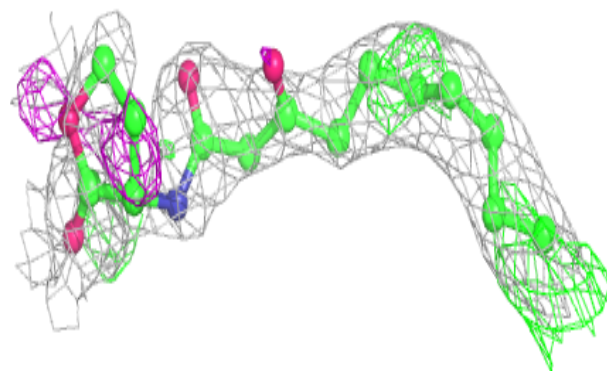
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 3M5 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 3M5 B 403:**

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



6.5 Other polymers

There are no such residues in this entry.