



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 10:17 am BST

PDB ID : 3RA7
Title : Bispecific digoxigenin binding antibodies for targeted payload delivery
Authors : Hopfner, K.P.; Lammens, A.
Deposited on : 2011-03-27
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 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

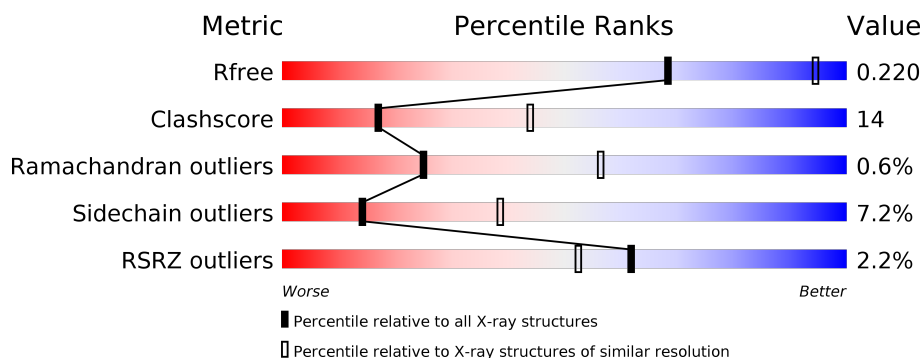
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 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	L	214	<div> <div>2%</div> <div> <div></div> <div>69%</div> <div>27%</div> <div>..</div> </div> </div>
1	M	214	<div> <div>2%</div> <div> <div></div> <div>68%</div> <div>28%</div> <div>..</div> </div> </div>
2	H	228	<div> <div>2%</div> <div> <div></div> <div>67%</div> <div>29%</div> <div>..</div> </div> </div>
2	I	228	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>22%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7061 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.

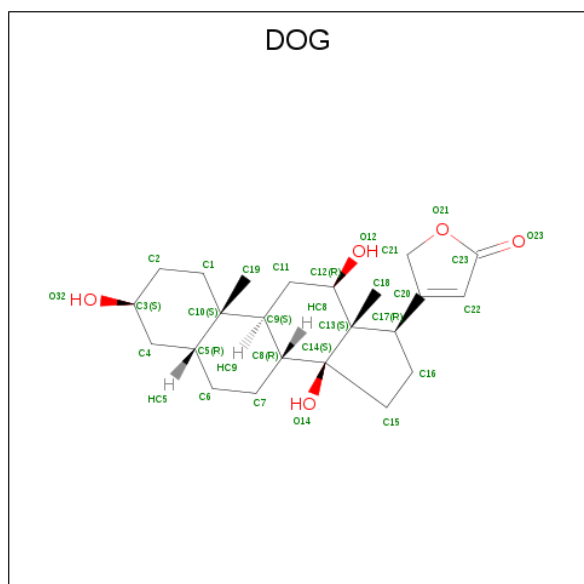
- Molecule 1 is a protein called Fab fragment, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	76	0	0
			1635	1018	277	334	6			
1	M	211	Total	C	N	O	S	47	0	0
			1634	1017	277	334	6			

- Molecule 2 is a protein called Fab fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	225	Total	C	N	O	S	65	0	0
			1694	1075	276	334	9			
2	I	225	Total	C	N	O	S	49	0	0
			1694	1075	276	334	9			

- Molecule 3 is DIGOXIGENIN (three-letter code: DOG) (formula: $C_{23}H_{34}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	L	1	Total	C	O	0	0
			28	23	5		
3	M	1	Total	C	O	0	0
			28	23	5		

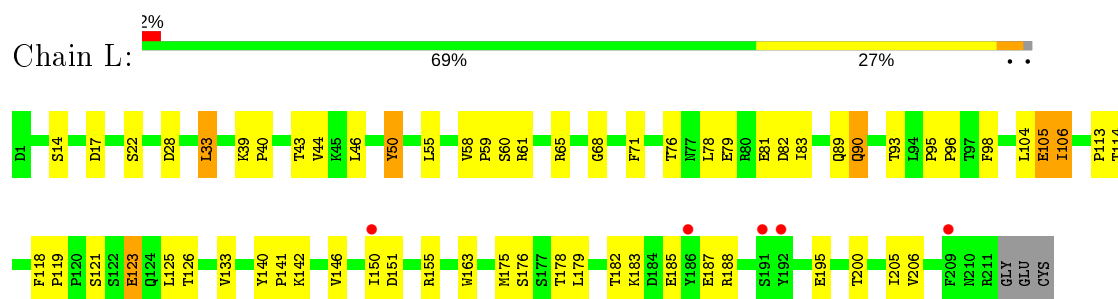
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	66	Total	O	0	0
			66	66		
4	H	97	Total	O	0	0
			97	97		
4	M	79	Total	O	0	0
			79	79		
4	I	106	Total	O	0	0
			106	106		

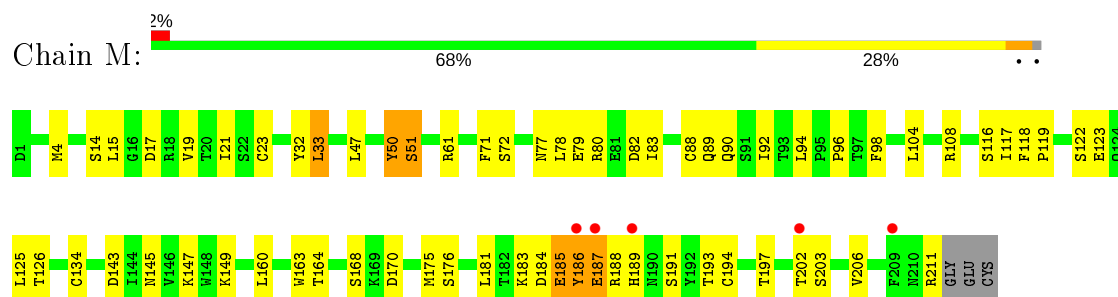
3 Residue-property plots

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.

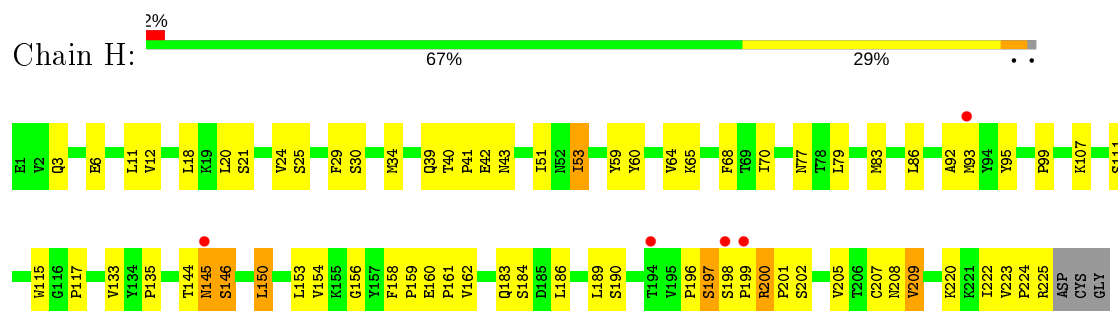
- Molecule 1: Fab fragment, light chain



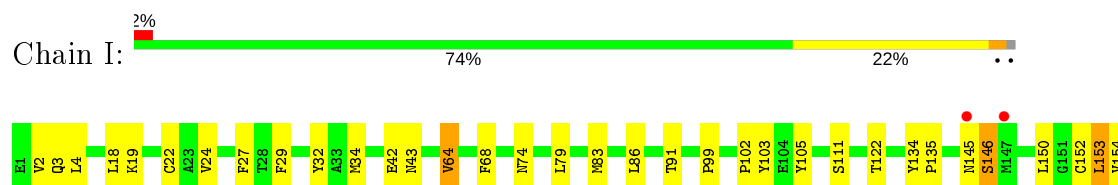
- Molecule 1: Fab fragment, light chain

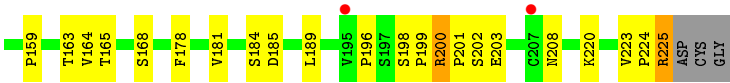


- Molecule 2: Fab fragment, heavy chain



- Molecule 2: Fab fragment, heavy chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	138.01Å 138.01Å 123.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.04 – 2.80 48.79 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.04-2.80) 99.9 (48.79-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.25 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.204 , 0.231 0.200 , 0.220	Depositor DCC
R_{free} test set	1536 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7061	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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: DOG

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	L	0.24	0/1668	0.45	0/2263
1	M	0.46	3/1667 (0.2%)	0.48	0/2261
2	H	0.24	0/1741	0.44	0/2381
2	I	0.24	0/1741	0.46	0/2381
All	All	0.31	3/6817 (0.0%)	0.46	0/9286

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	186	TYR	CD2-CE2	-5.43	1.31	1.39
1	M	186	TYR	CD1-CE1	-5.12	1.31	1.39
1	M	185	GLU	CG-CD	-5.02	1.44	1.51

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	L	1635	0	1590	38	0
1	M	1634	0	1586	39	0
2	H	1694	0	1652	60	0
2	I	1694	0	1652	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	28	0	32	1	0
3	M	28	0	34	1	0
4	H	97	0	0	3	0
4	I	106	0	0	0	0
4	L	66	0	0	1	0
4	M	79	0	0	0	0
All	All	7061	0	6546	178	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 14.

All (178) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:160:GLU:HG3	2:H:161:PRO:HA	1.43	0.98
2:I:200:ARG:HG3	2:I:200:ARG:HH11	1.24	0.97
2:H:200:ARG:HH11	2:H:200:ARG:HG3	1.31	0.93
2:H:40:THR:HG22	2:H:42:GLU:H	1.34	0.93
1:M:186:TYR:O	1:M:187:GLU:HG3	1.70	0.91
2:H:150:LEU:HD23	2:H:205:VAL:HG21	1.55	0.88
1:L:150:ILE:HD11	1:L:179:LEU:HD21	1.54	0.87
2:I:200:ARG:CG	2:I:200:ARG:HH11	1.88	0.85
2:H:200:ARG:HH11	2:H:200:ARG:CG	1.89	0.85
2:I:224:PRO:O	2:I:225:ARG:HB2	1.79	0.82
1:M:78:LEU:HD11	1:M:104:LEU:HD21	1.62	0.81
1:M:186:TYR:O	1:M:187:GLU:CB	2.30	0.79
2:I:200:ARG:HD2	2:I:201:PRO:HA	1.66	0.78
1:M:186:TYR:O	1:M:187:GLU:CG	2.30	0.78
1:M:183:LYS:O	1:M:186:TYR:O	2.05	0.75
2:I:135:PRO:HG3	2:I:220:LYS:HD2	1.72	0.71
1:M:186:TYR:O	1:M:187:GLU:HB2	1.91	0.70
2:H:145:ASN:O	2:H:197:SER:OG	2.11	0.69
2:H:40:THR:HG23	2:H:41:PRO:HD2	1.73	0.69
2:H:196:PRO:O	2:H:199:PRO:HD2	1.93	0.69
1:L:105:GLU:HG3	1:L:106:ILE:N	2.06	0.69
2:H:160:GLU:HG3	2:H:161:PRO:CA	2.20	0.69
2:H:60:TYR:HB2	2:H:65:LYS:HG3	1.74	0.68
1:L:33:LEU:HD22	1:L:71:PHE:CG	2.29	0.67
2:H:200:ARG:NH1	2:H:200:ARG:HG3	2.08	0.67
1:L:195:GLU:HG2	1:L:206:VAL:HG22	1.78	0.66
1:M:33:LEU:HD22	1:M:71:PHE:CG	2.31	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:64:VAL:HG13	2:H:68:PHE:HB2	1.77	0.64
2:H:198:SER:HB2	2:H:199:PRO:HD3	1.80	0.64
2:H:156:GLY:HA2	2:H:186:LEU:HB3	1.79	0.63
1:L:90:GLN:NE2	1:L:93:THR:O	2.32	0.63
2:H:144:THR:HA	2:H:145:ASN:HB2	1.80	0.62
1:M:117:ILE:HD12	1:M:194:CYS:SG	2.40	0.62
1:L:33:LEU:O	1:L:50:TYR:O	2.17	0.61
2:H:200:ARG:HD2	2:H:201:PRO:HA	1.82	0.61
2:I:103:TYR:CE2	2:I:105:TYR:HB2	2.35	0.61
1:M:187:GLU:HA	1:M:211:ARG:CZ	2.31	0.60
2:I:64:VAL:HG13	2:I:68:PHE:HB2	1.83	0.60
2:I:196:PRO:O	2:I:199:PRO:HD2	2.02	0.60
2:H:154:VAL:HB	2:H:189:LEU:HD23	1.83	0.59
1:M:160:LEU:HD21	2:I:181:VAL:HB	1.85	0.58
1:L:182:THR:CG2	1:L:185:GLU:H	2.17	0.58
2:H:18:LEU:HB3	2:H:83:MET:HE2	1.85	0.58
1:M:33:LEU:O	1:M:50:TYR:O	2.22	0.58
2:H:135:PRO:HG3	2:H:220:LYS:HD2	1.86	0.57
2:I:42:GLU:O	2:I:43:ASN:HB2	2.03	0.57
2:I:29:PHE:HE1	2:I:34:MET:HG3	1.70	0.57
2:H:3:GLN:HB2	2:H:25:SER:HB2	1.86	0.57
1:L:90:GLN:HE21	1:L:93:THR:H	1.51	0.56
2:I:165:THR:OG1	2:I:208:ASN:HB2	2.05	0.56
1:M:83:ILE:HD12	1:M:168:SER:HA	1.87	0.56
1:M:189:HIS:O	1:M:211:ARG:HD3	2.06	0.56
2:I:200:ARG:NH1	2:I:200:ARG:CG	2.58	0.56
2:I:200:ARG:NH1	2:I:200:ARG:HG3	2.04	0.55
2:H:196:PRO:HB2	2:H:199:PRO:CD	2.37	0.55
1:M:80:ARG:O	1:M:83:ILE:HG13	2.07	0.54
1:M:122:SER:O	1:M:126:THR:HG23	2.07	0.54
2:H:196:PRO:HB2	2:H:199:PRO:HD3	1.90	0.54
2:H:24:VAL:HG21	2:H:29:PHE:CD1	2.43	0.54
2:H:24:VAL:HG22	2:H:77:ASN:O	2.08	0.54
2:H:93:MET:HG2	2:H:95:TYR:CZ	2.43	0.53
1:L:140:TYR:CG	1:L:141:PRO:HA	2.43	0.53
3:M:3123:DOG:H211	3:M:3123:DOG:H183	1.90	0.53
1:L:150:ILE:O	1:L:151:ASP:HB2	2.08	0.53
1:M:89:GLN:HG3	1:M:98:PHE:CE1	2.43	0.53
2:H:200:ARG:CG	2:H:200:ARG:NH1	2.59	0.53
2:I:64:VAL:HG13	2:I:68:PHE:CG	2.43	0.52
1:L:182:THR:HG23	1:L:185:GLU:H	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:99:PRO:HA	2:H:111:SER:O	2.09	0.52
2:H:144:THR:HA	2:H:145:ASN:CB	2.37	0.52
2:H:64:VAL:CG1	2:H:68:PHE:HB2	2.40	0.52
1:L:14:SER:O	1:L:17:ASP:HB2	2.11	0.51
2:H:160:GLU:CG	2:H:161:PRO:HA	2.30	0.51
2:I:29:PHE:CE1	2:I:34:MET:HG3	2.46	0.51
2:H:183:GLN:O	2:H:184:SER:HB2	2.10	0.51
2:H:223:VAL:HG13	2:H:224:PRO:HD2	1.93	0.51
2:H:29:PHE:CE1	2:H:34:MET:HG3	2.45	0.50
1:M:33:LEU:HD22	1:M:71:PHE:CB	2.41	0.50
2:I:223:VAL:HG13	2:I:224:PRO:HD2	1.93	0.50
2:I:22:CYS:HB3	2:I:79:LEU:HB3	1.94	0.50
1:M:116:SER:O	1:M:134:CYS:HA	2.11	0.50
2:H:93:MET:HG2	2:H:95:TYR:CE1	2.47	0.50
2:I:198:SER:OG	2:I:199:PRO:HD3	2.12	0.50
2:H:53:ILE:HG22	4:H:312:HOH:O	2.12	0.49
2:H:64:VAL:HG13	2:H:68:PHE:CG	2.47	0.49
4:L:266:HOH:O	2:H:117:PRO:HG3	2.12	0.49
2:H:34:MET:HB3	2:H:79:LEU:HD22	1.93	0.49
1:L:133:VAL:HG22	1:L:178:THR:HG23	1.94	0.49
1:L:61:ARG:HH21	1:L:82:ASP:CG	2.16	0.49
1:L:89:GLN:HG3	1:L:98:PHE:CE1	2.47	0.49
2:H:150:LEU:HG	2:H:222:ILE:HG21	1.94	0.48
2:H:145:ASN:O	2:H:146:SER:C	2.52	0.48
2:I:32:TYR:CZ	2:I:102:PRO:HG3	2.47	0.48
1:L:43:THR:HG23	4:H:259:HOH:O	2.12	0.48
1:M:61:ARG:NH2	1:M:82:ASP:OD1	2.47	0.48
1:L:33:LEU:HD22	1:L:71:PHE:CB	2.44	0.48
2:H:144:THR:CA	2:H:145:ASN:HB2	2.43	0.48
2:H:133:VAL:HG21	2:H:209:VAL:HG11	1.96	0.47
1:M:175:MET:HG2	1:M:176:SER:N	2.29	0.47
1:M:94:LEU:HD23	1:M:96:PRO:HD3	1.95	0.47
1:L:28:ASP:OD1	1:L:68:GLY:HA2	2.15	0.47
1:L:39:LYS:HB3	1:L:40:PRO:HD2	1.98	0.46
2:I:196:PRO:HD2	2:I:199:PRO:HG2	1.98	0.46
2:I:164:VAL:HA	2:I:208:ASN:O	2.15	0.46
1:M:145:ASN:HB3	1:M:197:THR:HB	1.96	0.46
1:L:183:LYS:O	1:L:187:GLU:HG3	2.16	0.46
2:H:207:CYS:SG	2:H:220:LYS:HB2	2.56	0.46
2:I:154:VAL:HB	2:I:189:LEU:HB3	1.97	0.45
2:I:34:MET:HB3	2:I:79:LEU:HD22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:39:GLN:O	2:H:92:ALA:HB1	2.16	0.45
2:I:83:MET:HE2	2:I:86:LEU:HD21	1.98	0.45
1:L:163:TRP:CD1	1:L:163:TRP:N	2.85	0.45
2:H:115:TRP:N	2:H:115:TRP:CD1	2.85	0.45
2:I:184:SER:O	2:I:185:ASP:HB2	2.16	0.45
2:I:99:PRO:HA	2:I:111:SER:O	2.17	0.45
1:L:185:GLU:HA	1:L:188:ARG:HG2	1.99	0.45
2:H:11:LEU:HD22	2:H:159:PRO:HG3	1.98	0.45
1:M:32:TYR:CG	1:M:92:ILE:HB	2.52	0.45
1:L:142:LYS:O	1:L:142:LYS:HG2	2.17	0.44
2:H:30:SER:O	2:H:53:ILE:HG21	2.16	0.44
2:I:134:TYR:HA	2:I:135:PRO:HD3	1.78	0.44
1:L:83:ILE:HG23	1:L:104:LEU:O	2.17	0.44
1:M:163:TRP:CD1	1:M:175:MET:HG3	2.52	0.44
2:I:150:LEU:HD11	2:I:200:ARG:HG2	2.00	0.44
1:M:19:VAL:HG21	1:M:78:LEU:HD22	1.99	0.44
1:M:193:THR:CG2	1:M:206:VAL:HG23	2.48	0.44
2:I:168:SER:N	2:I:208:ASN:OD1	2.46	0.43
1:L:118:PHE:HA	1:L:119:PRO:HD3	1.81	0.43
1:M:125:LEU:HD22	1:M:183:LYS:HG3	1.99	0.43
2:I:146:SER:OG	2:I:146:SER:O	2.33	0.43
2:I:64:VAL:HG13	2:I:68:PHE:CB	2.48	0.43
1:M:149:LYS:HB2	1:M:193:THR:HB	2.00	0.43
2:H:40:THR:O	2:H:43:ASN:N	2.49	0.43
1:L:185:GLU:HA	1:L:188:ARG:CG	2.48	0.43
1:L:43:THR:HG22	1:L:44:VAL:N	2.34	0.43
1:L:46:LEU:HB3	1:L:55:LEU:HD22	2.00	0.43
2:H:68:PHE:N	2:H:68:PHE:CD1	2.87	0.43
2:I:29:PHE:HD2	2:I:74:ASN:HA	1.83	0.43
1:L:113:PRO:HG2	1:L:205:ILE:HD12	2.00	0.43
1:M:14:SER:O	1:M:17:ASP:HB2	2.18	0.43
2:H:64:VAL:HG13	2:H:68:PHE:CB	2.46	0.43
1:L:150:ILE:HD12	1:L:155:ARG:HD3	2.01	0.43
1:M:79:GLU:O	1:M:82:ASP:HB2	2.18	0.43
2:H:6:GLU:HA	2:H:21:SER:O	2.18	0.43
2:I:134:TYR:HB2	2:I:153:LEU:HB3	2.01	0.42
1:L:79:GLU:O	1:L:82:ASP:HB2	2.18	0.42
2:I:196:PRO:HB2	2:I:199:PRO:CD	2.49	0.42
2:H:51:ILE:HB	2:H:70:ILE:HD13	2.01	0.42
1:L:58:VAL:HA	1:L:59:PRO:HD3	1.93	0.42
1:M:78:LEU:HD12	1:M:78:LEU:HA	1.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:3123:DOG:H212	3:L:3123:DOG:H161	1.84	0.42
2:H:158:PHE:CG	2:H:159:PRO:HA	2.55	0.42
2:I:83:MET:HB3	2:I:86:LEU:HD21	2.01	0.42
2:H:107:LYS:NZ	4:H:280:HOH:O	2.52	0.42
2:I:4:LEU:CD2	2:I:24:VAL:HG12	2.50	0.42
1:M:118:PHE:HA	1:M:119:PRO:HD3	1.89	0.42
1:M:4:MET:HB3	1:M:88:CYS:SG	2.59	0.42
1:M:108:ARG:HD2	1:M:170:ASP:O	2.19	0.42
1:M:50:TYR:HB3	1:M:51:SER:H	1.73	0.42
1:M:21:ILE:O	1:M:72:SER:HA	2.20	0.42
2:H:83:MET:HE3	2:H:86:LEU:HD11	2.02	0.42
2:I:24:VAL:HG21	2:I:29:PHE:CD1	2.55	0.41
2:I:2:VAL:HG13	2:I:27:PHE:CD1	2.55	0.41
2:H:199:PRO:C	2:H:202:SER:O	2.59	0.41
2:H:68:PHE:N	2:H:68:PHE:HD1	2.18	0.41
2:I:91:THR:HG23	2:I:122:THR:HA	2.01	0.41
1:L:78:LEU:HD11	1:L:104:LEU:HD21	2.02	0.41
2:H:40:THR:HG23	2:H:41:PRO:CD	2.47	0.41
1:M:185:GLU:O	1:M:188:ARG:HB3	2.21	0.41
1:L:121:SER:HB2	1:L:123:GLU:HG2	2.02	0.41
1:L:89:GLN:HG3	1:L:98:PHE:CD1	2.56	0.41
1:L:95:PRO:HA	1:L:96:PRO:HD3	1.96	0.41
2:I:199:PRO:O	2:I:203:GLU:HB2	2.21	0.40
1:M:183:LYS:O	1:M:187:GLU:HG3	2.21	0.40
2:H:30:SER:O	2:H:53:ILE:CG2	2.70	0.40
1:L:95:PRO:HB3	2:H:59:TYR:OH	2.21	0.40
1:L:175:MET:HG2	1:L:176:SER:N	2.37	0.40
1:M:164:THR:HG23	2:I:178:PHE:CD1	2.56	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	L	209/214 (98%)	199 (95%)	9 (4%)	1 (0%)	29	61
1	M	209/214 (98%)	199 (95%)	9 (4%)	1 (0%)	29	61
2	H	223/228 (98%)	210 (94%)	11 (5%)	2 (1%)	17	46
2	I	223/228 (98%)	214 (96%)	8 (4%)	1 (0%)	34	66
All	All	864/884 (98%)	822 (95%)	37 (4%)	5 (1%)	25	56

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	187	GLU
1	L	200	THR
2	H	145	ASN
2	H	146	SER
2	I	159	PRO

5.3.2 Protein sidechains ⓘ

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	L	189/191 (99%)	174 (92%)	15 (8%)	12	34
1	M	189/191 (99%)	173 (92%)	16 (8%)	10	31
2	H	191/193 (99%)	179 (94%)	12 (6%)	18	46
2	I	191/193 (99%)	179 (94%)	12 (6%)	18	46
All	All	760/768 (99%)	705 (93%)	55 (7%)	14	38

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	22	SER
1	L	33	LEU
1	L	50	TYR
1	L	60	SER
1	L	65	ARG
1	L	76	THR

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Mol	Chain	Res	Type
1	L	81	GLU
1	L	90	GLN
1	L	105	GLU
1	L	106	ILE
1	L	114	THR
1	L	123	GLU
1	L	125	LEU
1	L	126	THR
1	L	146	VAL
2	H	12	VAL
2	H	20	LEU
2	H	53	ILE
2	H	150	LEU
2	H	153	LEU
2	H	162	VAL
2	H	190	SER
2	H	197	SER
2	H	200	ARG
2	H	208	ASN
2	H	209	VAL
2	H	225	ARG
1	M	15	LEU
1	M	23	CYS
1	M	33	LEU
1	M	47	LEU
1	M	50	TYR
1	M	51	SER
1	M	77	ASN
1	M	90	GLN
1	M	123	GLU
1	M	143	ASP
1	M	147	LYS
1	M	181	LEU
1	M	184	ASP
1	M	191	SER
1	M	202	THR
1	M	203	SER
2	I	3	GLN
2	I	18	LEU
2	I	19	LYS
2	I	64	VAL
2	I	145	ASN

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Mol	Chain	Res	Type
2	I	146	SER
2	I	152	CYS
2	I	153	LEU
2	I	163	THR
2	I	200	ARG
2	I	202	SER
2	I	225	ARG

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

Mol	Chain	Res	Type
1	L	27	GLN
1	L	77	ASN

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](#)

2 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	DOG	L	3123	-	32,32,32	3.90	20 (62%)	48,53,53	1.75	11 (22%)
3	DOG	M	3123	-	32,32,32	3.40	16 (50%)	48,53,53	1.76	14 (29%)

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	DOG	L	3123	-	-	0/4/81/81	0/5/5/5
3	DOG	M	3123	-	-	0/4/81/81	0/5/5/5

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	M	3123	DOG	C21-C20	-9.35	1.29	1.50
3	L	3123	DOG	C21-C20	-9.17	1.30	1.50
3	L	3123	DOG	O21-C21	-8.11	1.30	1.44
3	M	3123	DOG	O21-C21	-7.94	1.30	1.44
3	L	3123	DOG	C14-C13	-7.43	1.48	1.56
3	M	3123	DOG	C22-C23	-7.37	1.29	1.46
3	L	3123	DOG	C22-C23	-7.19	1.29	1.46
3	M	3123	DOG	C14-C13	-5.44	1.50	1.56
3	L	3123	DOG	C10-C5	-4.84	1.47	1.55
3	L	3123	DOG	O14-C14	-4.50	1.36	1.44
3	L	3123	DOG	C10-C9	-4.44	1.47	1.56
3	M	3123	DOG	C15-C14	-4.35	1.47	1.54
3	L	3123	DOG	O12-C12	-4.31	1.36	1.43
3	L	3123	DOG	C15-C14	-4.03	1.48	1.54
3	L	3123	DOG	C1-C10	-3.78	1.47	1.54
3	M	3123	DOG	O21-C23	-3.74	1.26	1.36
3	M	3123	DOG	O14-C14	-3.57	1.38	1.44
3	L	3123	DOG	C13-C12	-3.55	1.47	1.55
3	L	3123	DOG	C19-C10	-3.51	1.48	1.54
3	L	3123	DOG	C9-C8	-3.48	1.47	1.54
3	L	3123	DOG	O21-C23	-3.27	1.27	1.36
3	M	3123	DOG	C7-C8	-3.23	1.48	1.53
3	L	3123	DOG	C18-C13	-3.10	1.47	1.53
3	M	3123	DOG	O12-C12	-2.98	1.38	1.43
3	L	3123	DOG	C11-C12	-2.92	1.48	1.53
3	M	3123	DOG	C19-C10	-2.91	1.49	1.54
3	L	3123	DOG	C13-C17	-2.73	1.53	1.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	3123	DOG	C11-C9	-2.65	1.49	1.53
3	L	3123	DOG	C7-C8	-2.50	1.49	1.53
3	L	3123	DOG	C1-C2	-2.48	1.48	1.53
3	M	3123	DOG	C11-C9	-2.36	1.49	1.53
3	M	3123	DOG	C13-C17	-2.29	1.54	1.57
3	M	3123	DOG	C18-C13	-2.24	1.49	1.53
3	M	3123	DOG	C11-C12	-2.19	1.49	1.53
3	M	3123	DOG	C9-C8	-2.15	1.50	1.54
3	M	3123	DOG	C17-C20	2.09	1.54	1.50

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	3123	DOG	C11-C9-C10	-5.56	108.00	113.73
3	M	3123	DOG	C1-C2-C3	5.08	116.99	110.47
3	L	3123	DOG	C9-C10-C5	3.95	114.14	108.58
3	M	3123	DOG	C16-C17-C20	-3.56	106.71	113.68
3	M	3123	DOG	C19-C10-C1	-3.37	102.83	108.26
3	M	3123	DOG	C21-O21-C23	-3.25	104.91	108.85
3	L	3123	DOG	C16-C17-C20	-3.19	107.42	113.68
3	L	3123	DOG	C1-C10-C9	-3.01	106.61	111.35
3	L	3123	DOG	C1-C2-C3	3.00	114.31	110.47
3	M	3123	DOG	O21-C21-C20	2.82	110.26	104.80
3	M	3123	DOG	C2-C1-C10	2.57	117.18	112.78
3	L	3123	DOG	O21-C23-C22	2.57	112.02	108.67
3	M	3123	DOG	O21-C23-C22	2.54	111.99	108.67
3	L	3123	DOG	C4-C3-C2	-2.54	107.53	110.55
3	M	3123	DOG	C9-C10-C5	2.52	112.13	108.58
3	M	3123	DOG	C9-C11-C12	-2.51	110.99	114.30
3	L	3123	DOG	O12-C12-C13	-2.47	105.06	110.85
3	M	3123	DOG	C13-C17-C20	2.37	118.99	115.78
3	M	3123	DOG	C19-C10-C5	-2.34	106.39	110.36
3	M	3123	DOG	C1-C10-C5	2.25	111.10	107.77
3	L	3123	DOG	C20-C22-C23	-2.22	104.24	108.84
3	M	3123	DOG	O23-C23-C22	-2.17	126.42	130.81
3	L	3123	DOG	C9-C11-C12	-2.16	111.44	114.30
3	L	3123	DOG	O12-C12-C11	-2.07	104.91	109.12
3	M	3123	DOG	O12-C12-C11	-2.04	104.97	109.12

There are no chirality outliers.

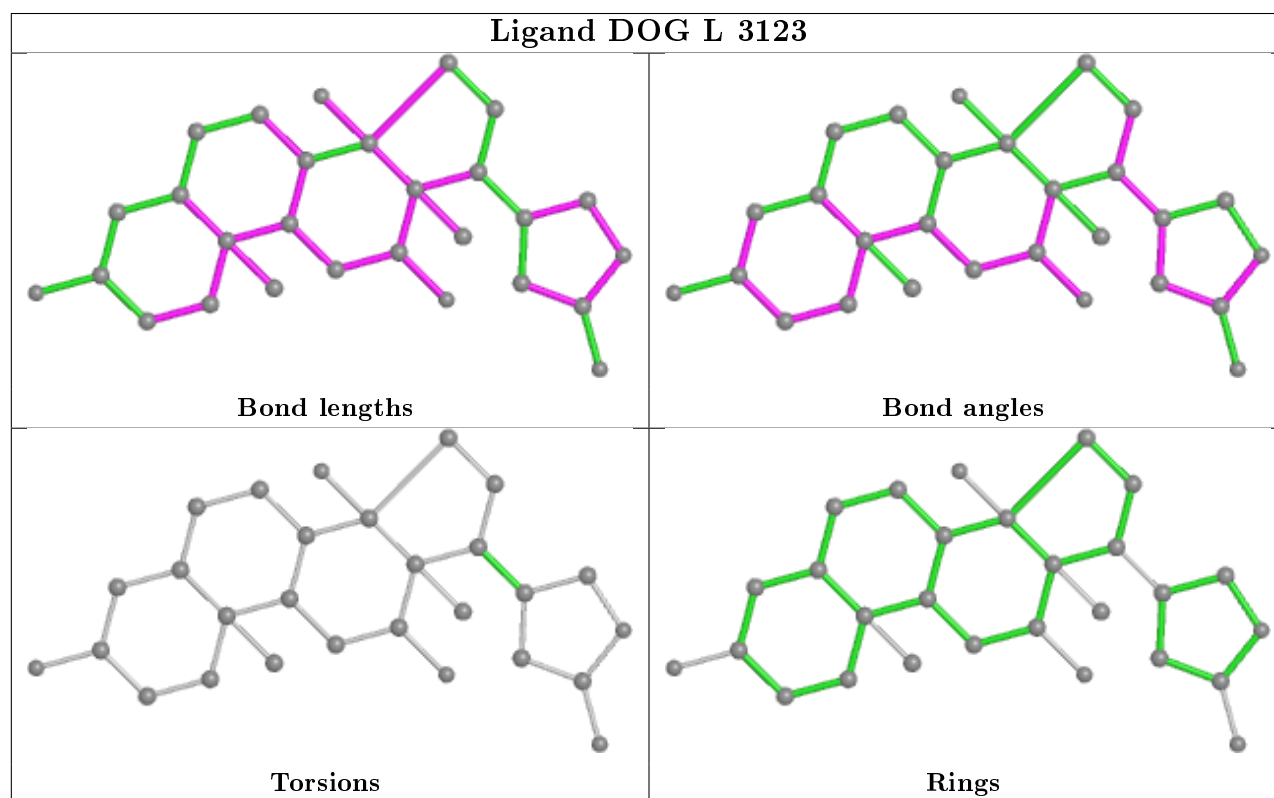
There are no torsion outliers.

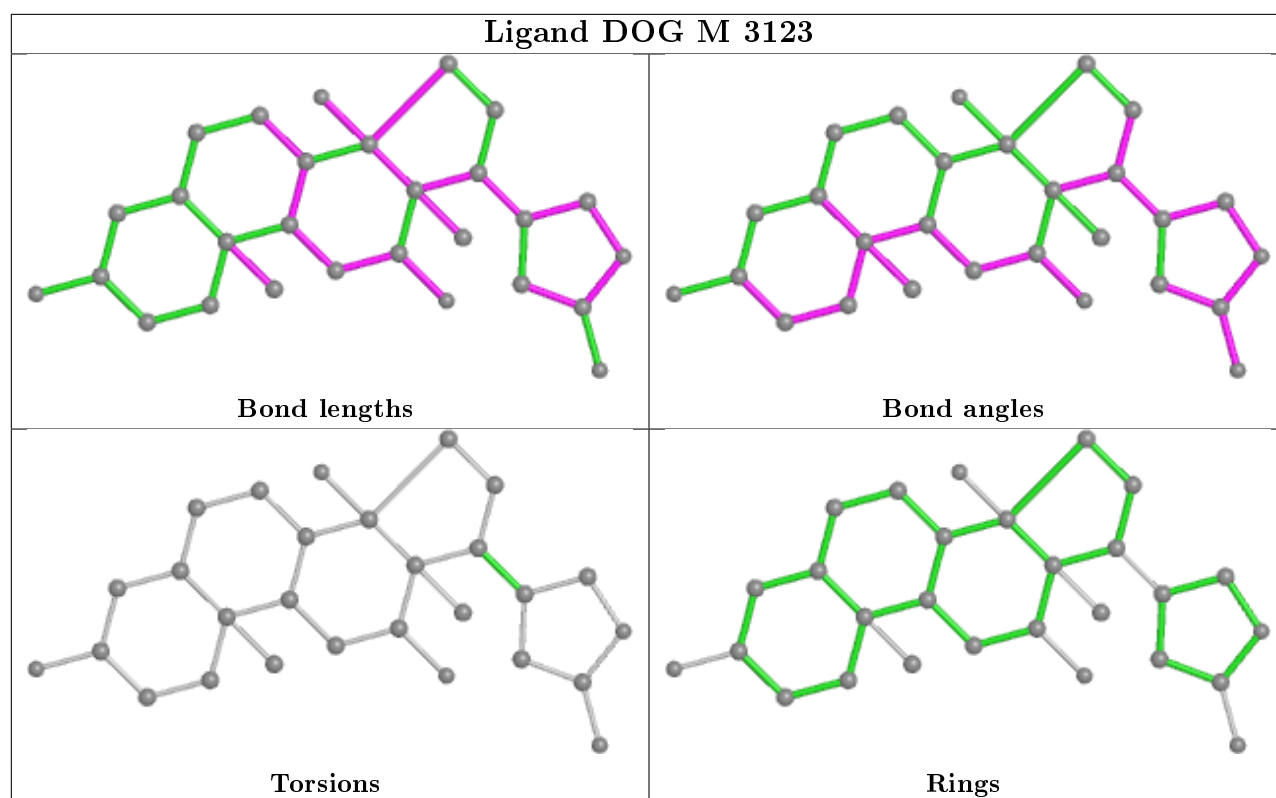
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	3123	DOG	1	0
3	M	3123	DOG	1	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	L	211/214 (98%)	0.23	5 (2%) 59 49	15, 30, 54, 62	35 (16%)
1	M	211/214 (98%)	0.11	5 (2%) 59 49	12, 23, 41, 59	23 (10%)
2	H	221/228 (96%)	0.09	5 (2%) 60 51	13, 23, 42, 62	16 (7%)
2	I	220/228 (96%)	-0.07	4 (1%) 68 61	11, 18, 38, 53	12 (5%)
All	All	863/884 (97%)	0.09	19 (2%) 62 52	11, 23, 46, 62	86 (9%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	209	PHE	4.8
2	H	194	THR	3.4
2	H	199	PRO	3.3
1	L	150	ILE	3.2
1	M	209	PHE	3.1
1	M	187	GLU	3.0
1	M	186	TYR	2.8
1	M	189	HIS	2.6
2	I	145	ASN	2.6
1	M	202	THR	2.5
1	L	186	TYR	2.5
1	L	192	TYR	2.4
2	H	145	ASN	2.4
2	I	207	CYS	2.3
2	I	195	VAL	2.3
1	L	191	SER	2.3
2	H	198	SER	2.2
2	H	93	MET	2.1
2	I	147	MET	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](#)

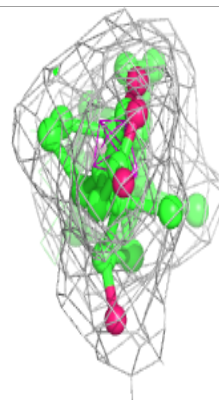
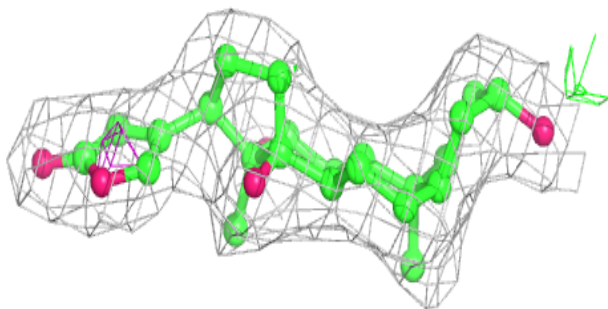
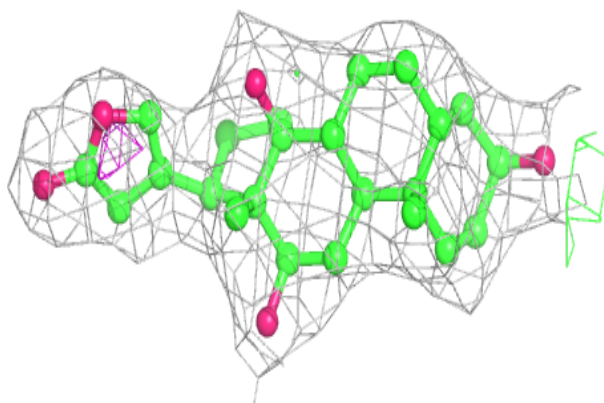
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	DOG	M	3123	28/28	0.96	0.15	12,15,21,25	0
3	DOG	L	3123	28/28	0.98	0.12	14,21,25,31	0

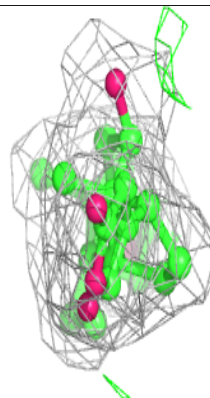
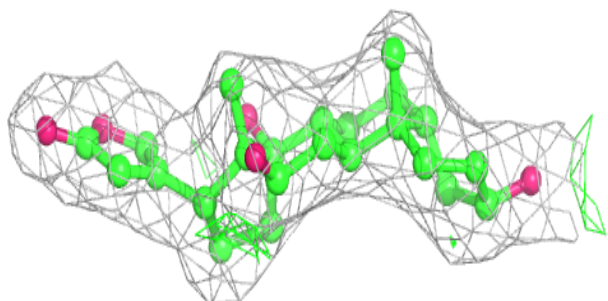
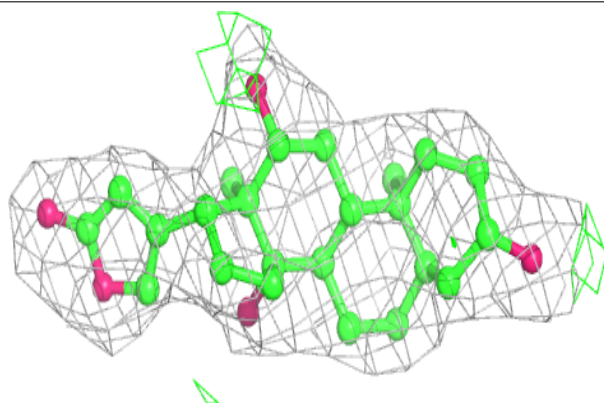
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DOG M 3123:

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

**Electron density around DOG L 3123:**

$2mF_o-DF_c$ (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.