



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:34 PM UTC

PDB ID : 7DBL / pdb_00007dbl
Title : Acyl-CoA hydrolase MpaH' mutant S139A in complex with MPA
Authors : Li, S.Y.; You, C.
Deposited on : 2020-10-20
Resolution : 1.84 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

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

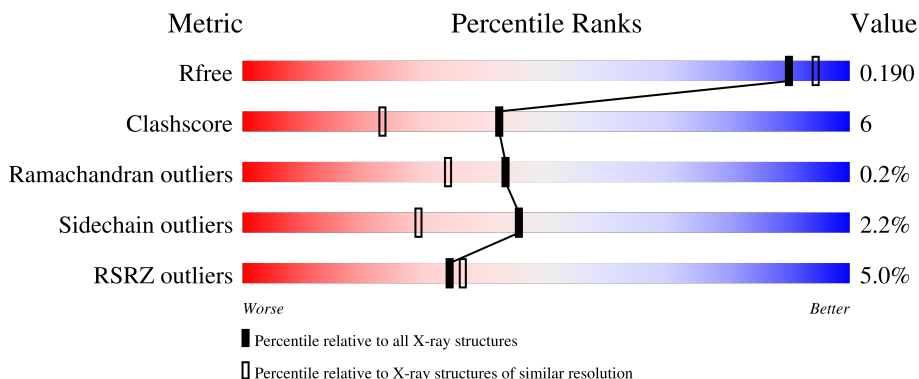
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.84 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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	433	
1	B	433	
1	C	433	
1	D	433	

2 Entry composition i

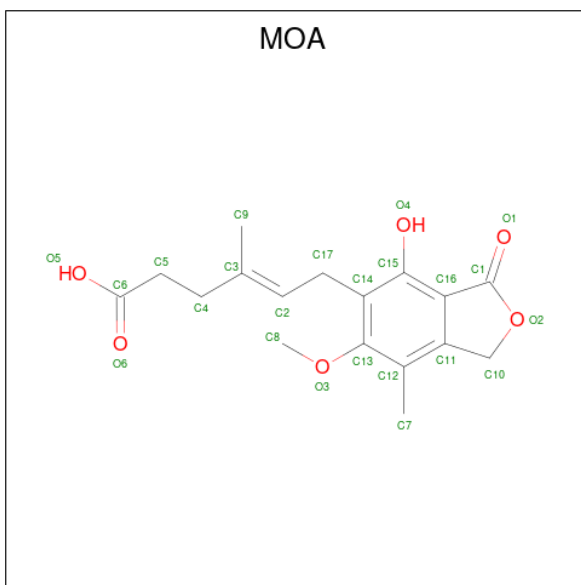
There are 3 unique types of molecules in this entry. The entry contains 14365 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 acyl-CoA hydrolase MpaH'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	422	Total 3369	C 2143	N 593	O 615	S 18	0	0	0
1	B	418	Total 3342	C 2127	N 589	O 609	S 17	0	0	0
1	C	418	Total 3342	C 2127	N 589	O 609	S 17	0	0	0
1	D	420	Total 3351	C 2131	N 591	O 611	S 18	0	0	0

- Molecule 2 is MYCOPHENOLIC ACID (CCD ID: MOA) (formula: C₁₇H₂₀O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 23	C 17	O 6	0	0
2	B	1	Total 23	C 17	O 6	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			23	17	6		
2	D	1	Total	C	O	0	0
			23	17	6		

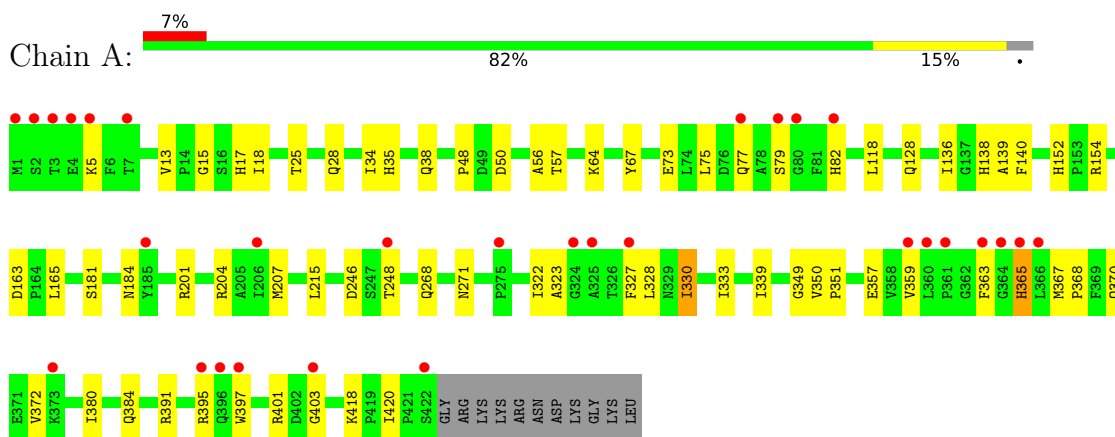
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	214	Total	O	0	0
			214	214		
3	B	224	Total	O	0	0
			224	224		
3	C	210	Total	O	0	0
			210	210		
3	D	221	Total	O	0	0
			221	221		

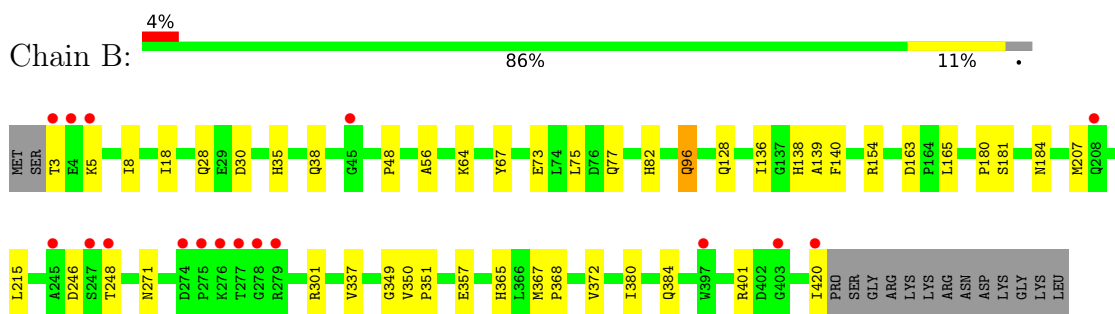
3 Residue-property plots [i](#)

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.

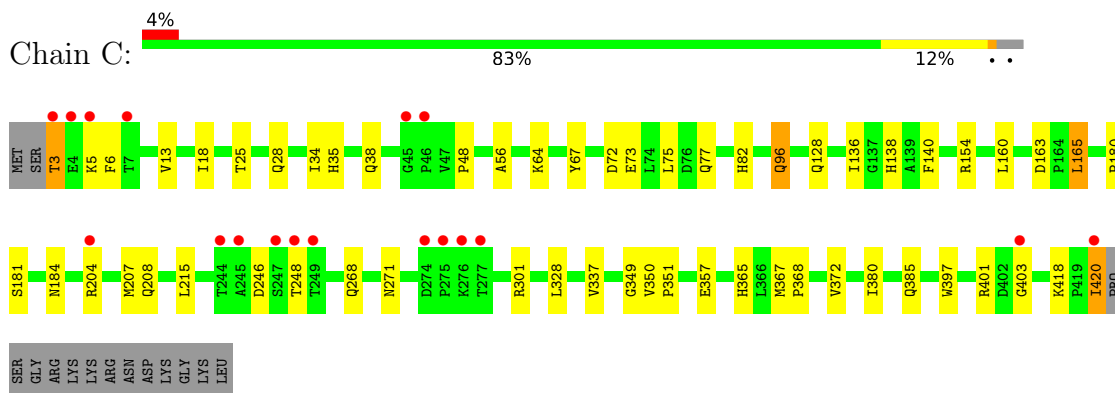
- Molecule 1: acyl-CoA hydrolase MpaH'



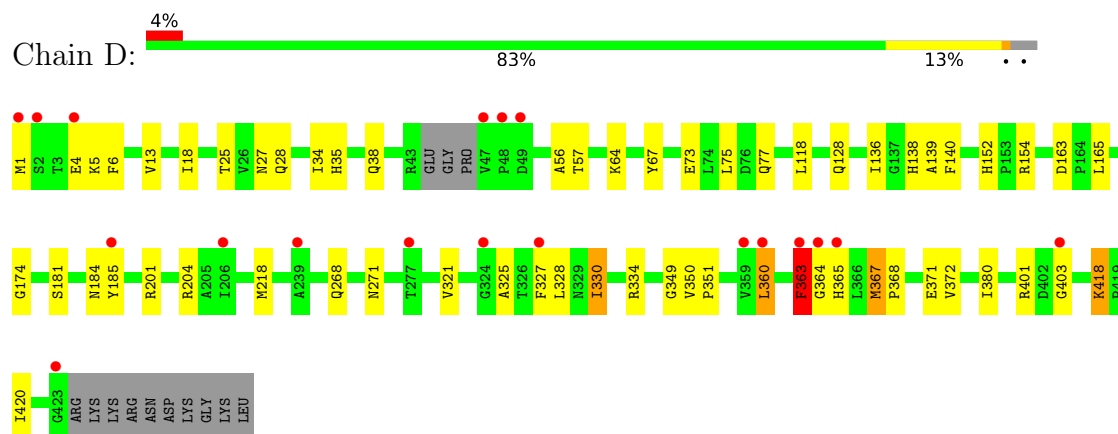
- Molecule 1: acyl-CoA hydrolase MpaH'



- Molecule 1: acyl-CoA hydrolase MpaH'



- Molecule 1: acyl-CoA hydrolase MpaH'



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.82Å 92.78Å 160.76Å 90.00° 95.39° 90.00°	Depositor
Resolution (Å)	48.08 – 1.84 48.08 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.08-1.84) 99.3 (48.08-1.84)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 1.84Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.179 , 0.205 0.188 , 0.190	Depositor DCC
R_{free} test set	8113 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtrriage
Anisotropy	0.453	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14365	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% 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: MOA

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	1.05	2/3460 (0.1%)	1.23	5/4710 (0.1%)
1	B	1.03	1/3432 (0.0%)	1.21	4/4672 (0.1%)
1	C	1.04	0/3432	1.21	4/4672 (0.1%)
1	D	1.06	1/3439 (0.0%)	1.23	9/4678 (0.2%)
All	All	1.05	4/13763 (0.0%)	1.22	22/18732 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	1
1	D	0	3
All	All	0	6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	364	GLY	C-O	13.28	1.41	1.23
1	A	365	HIS	C-O	6.50	1.31	1.24
1	A	17	HIS	CE1-NE2	5.80	1.38	1.32
1	B	30	ASP	C-O	5.23	1.29	1.23

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	363	PHE	CB-CA-C	8.01	123.63	111.85
1	C	96	GLN	CB-CG-CD	-6.24	102.00	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	271	ASN	CA-CB-CG	-6.21	106.39	112.60
1	A	363	PHE	N-CA-C	5.92	120.68	112.45
1	D	349	GLY	CA-C-N	5.80	123.90	120.24
1	D	349	GLY	C-N-CA	5.80	123.90	120.24
1	D	363	PHE	CA-CB-CG	5.75	119.55	113.80
1	C	349	GLY	CA-C-N	5.64	123.79	120.24
1	C	349	GLY	C-N-CA	5.64	123.79	120.24
1	B	349	GLY	CA-C-N	5.62	123.78	120.24
1	B	349	GLY	C-N-CA	5.62	123.78	120.24
1	B	96	GLN	CB-CG-CD	-5.60	103.08	112.60
1	D	163	ASP	CB-CA-C	-5.46	107.32	111.20
1	A	349	GLY	CA-C-N	5.17	123.50	120.24
1	A	349	GLY	C-N-CA	5.17	123.50	120.24
1	A	15	GLY	CA-C-N	5.14	128.07	120.82
1	A	15	GLY	C-N-CA	5.14	128.07	120.82
1	D	6	PHE	CA-C-N	5.05	128.04	120.87
1	D	6	PHE	C-N-CA	5.05	128.04	120.87
1	D	363	PHE	CA-C-N	5.02	131.25	121.41
1	D	363	PHE	C-N-CA	5.02	131.25	121.41
1	B	384	GLN	CB-CA-C	5.00	119.35	110.85

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	323	ALA	Peptide
1	A	403	GLY	Peptide
1	C	403	GLY	Peptide
1	D	363	PHE	Mainchain,Peptide
1	D	403	GLY	Peptide

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	3369	0	3337	53	0
1	B	3342	0	3308	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3342	0	3308	49	0
1	D	3351	0	3317	50	0
2	A	23	0	19	5	0
2	B	23	0	19	2	0
2	C	23	0	18	1	0
2	D	23	0	19	2	0
3	A	214	0	0	2	0
3	B	224	0	0	0	0
3	C	210	0	0	1	0
3	D	221	0	0	3	0
All	All	14365	0	13345	168	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 (168) 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:C:420:ILE:HD12	1:D:185:TYR:CZ	2.17	0.80
1:C:3:THR:HG23	1:C:72:ASP:OD1	1.82	0.80
1:D:330:ILE:CD1	1:D:334:ARG:HD2	2.16	0.74
1:B:246:ASP:OD1	1:B:248:THR:HG22	1.88	0.74
1:A:246:ASP:OD1	1:A:248:THR:HG22	1.88	0.73
1:D:27:ASN:HB3	3:D:738:HOH:O	1.86	0.73
1:C:246:ASP:OD1	1:C:248:THR:HG22	1.89	0.73
1:D:5:LYS:HB3	1:D:75:LEU:HD23	1.71	0.71
1:C:28:GLN:HE22	1:D:154:ARG:HH12	1.39	0.71
1:A:5:LYS:HB3	1:A:75:LEU:HD23	1.73	0.70
1:C:154:ARG:HH12	1:D:28:GLN:HE22	1.41	0.69
1:D:5:LYS:CB	1:D:75:LEU:HD23	2.24	0.68
1:C:420:ILE:HD12	1:D:185:TYR:CE1	2.28	0.68
1:A:327:PHE:HE1	2:A:501:MOA:C13	2.08	0.67
1:A:327:PHE:CE1	2:A:501:MOA:O3	2.48	0.67
1:A:5:LYS:CB	1:A:75:LEU:HD23	2.25	0.66
1:A:328:LEU:HD13	1:A:333:ILE:CD1	2.25	0.66
1:D:330:ILE:HD13	1:D:334:ARG:HD2	1.77	0.66
1:C:5:LYS:HB3	1:C:75:LEU:HD23	1.78	0.66
1:C:207:MET:HE3	1:C:215:LEU:HD22	1.76	0.66
1:B:38:GLN:HE22	1:B:128:GLN:HE22	1.44	0.65
1:C:5:LYS:CB	1:C:75:LEU:HD23	2.26	0.65
1:A:154:ARG:HH12	1:B:28:GLN:HE22	1.43	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:PHE:HE2	1:A:365:HIS:CE1	2.14	0.65
1:A:370:GLN:C	3:A:601:HOH:O	2.38	0.64
1:A:38:GLN:HE22	1:A:128:GLN:HE22	1.45	0.64
1:A:420:ILE:HD13	1:B:180:PRO:HB2	1.79	0.64
1:B:5:LYS:CB	1:B:75:LEU:HD23	2.28	0.64
1:C:67:TYR:OH	1:C:138:HIS:HD2	1.81	0.63
1:A:67:TYR:OH	1:A:138:HIS:HD2	1.82	0.63
1:D:165:LEU:HD11	1:D:327:PHE:CE1	2.34	0.63
1:B:67:TYR:OH	1:B:138:HIS:HD2	1.81	0.63
1:A:28:GLN:HE22	1:B:154:ARG:HH12	1.46	0.63
1:D:35:HIS:HD2	1:D:64:LYS:NZ	1.97	0.62
1:D:67:TYR:OH	1:D:138:HIS:HD2	1.83	0.62
1:B:35:HIS:HD2	1:B:64:LYS:NZ	1.97	0.62
1:A:165:LEU:HD12	1:A:328:LEU:HD21	1.81	0.62
1:C:207:MET:HE3	1:C:215:LEU:CD2	2.30	0.61
1:A:35:HIS:HD2	1:A:64:LYS:NZ	1.99	0.61
1:C:35:HIS:HD2	1:C:64:LYS:NZ	1.98	0.60
1:A:322:ILE:HG21	1:A:330:ILE:HD13	1.83	0.60
1:A:328:LEU:HD13	1:A:333:ILE:HD11	1.84	0.60
1:B:77:GLN:HB2	1:B:380:ILE:CD1	2.32	0.59
1:C:401:ARG:NH1	3:C:601:HOH:O	2.31	0.59
1:D:38:GLN:HE22	1:D:128:GLN:HE22	1.49	0.59
1:A:327:PHE:CE1	2:A:501:MOA:C13	2.85	0.59
1:B:5:LYS:HB3	1:B:75:LEU:HD23	1.84	0.59
1:A:77:GLN:HB2	1:A:380:ILE:CD1	2.34	0.58
1:C:38:GLN:HE22	1:C:128:GLN:HE22	1.52	0.58
1:C:77:GLN:HB2	1:C:380:ILE:CD1	2.34	0.58
1:C:181:SER:H	1:C:184:ASN:ND2	2.03	0.57
1:A:163:ASP:OD2	1:A:367:MET:HE2	2.05	0.57
1:D:165:LEU:HD12	1:D:328:LEU:HD21	1.87	0.57
1:C:3:THR:CG2	1:C:72:ASP:OD1	2.53	0.56
1:B:35:HIS:CD2	1:B:96:GLN:HG2	2.40	0.56
1:D:77:GLN:HB2	1:D:380:ILE:CD1	2.36	0.56
1:D:27:ASN:OD1	3:D:601:HOH:O	2.18	0.55
1:D:181:SER:H	1:D:184:ASN:ND2	2.03	0.55
1:B:181:SER:H	1:B:184:ASN:ND2	2.03	0.55
1:C:3:THR:HG21	1:C:6:PHE:HD1	1.72	0.55
1:A:181:SER:H	1:A:184:ASN:ND2	2.03	0.54
1:C:73:GLU:HG3	1:C:372:VAL:HG12	1.88	0.54
1:B:181:SER:H	1:B:184:ASN:HD22	1.56	0.54
1:A:330:ILE:HD12	1:A:330:ILE:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:GLN:CB	1:B:380:ILE:HD13	2.37	0.54
1:B:77:GLN:HB2	1:B:380:ILE:HD13	1.89	0.54
1:B:73:GLU:HG3	1:B:372:VAL:HG12	1.89	0.54
1:A:207:MET:HE3	1:A:215:LEU:CD2	2.38	0.53
1:D:321:VAL:HG13	1:D:360:LEU:HD22	1.89	0.53
1:D:330:ILE:HD11	1:D:334:ARG:HD2	1.90	0.53
1:A:139:ALA:HB1	2:A:501:MOA:C6	2.39	0.53
1:C:181:SER:H	1:C:184:ASN:HD22	1.55	0.53
1:D:365:HIS:CD2	2:D:501:MOA:O5	2.62	0.53
1:A:181:SER:H	1:A:184:ASN:HD22	1.57	0.53
1:A:201:ARG:NH1	1:A:204:ARG:HG3	2.24	0.52
1:D:165:LEU:HD11	1:D:327:PHE:HE1	1.76	0.51
1:C:154:ARG:HH22	1:D:28:GLN:HE21	1.58	0.51
1:D:181:SER:H	1:D:184:ASN:HD22	1.57	0.51
1:C:418:LYS:HD3	1:D:174:GLY:O	2.10	0.51
1:A:35:HIS:HD2	1:A:64:LYS:HZ2	1.59	0.51
1:A:207:MET:HE3	1:A:215:LEU:HD22	1.92	0.51
1:C:28:GLN:NE2	1:D:154:ARG:HH12	2.06	0.51
1:D:321:VAL:CG1	1:D:360:LEU:HD22	2.41	0.51
1:B:138:HIS:HE1	1:B:365:HIS:O	1.94	0.50
1:C:77:GLN:HB2	1:C:380:ILE:HD13	1.94	0.50
1:A:138:HIS:HE1	1:A:365:HIS:O	1.94	0.50
1:D:139:ALA:HB1	2:D:501:MOA:C6	2.42	0.50
1:C:420:ILE:HD12	1:D:185:TYR:CE2	2.47	0.50
1:C:138:HIS:HE1	1:C:365:HIS:O	1.94	0.50
1:A:154:ARG:HH12	1:B:28:GLN:NE2	2.10	0.49
1:C:180:PRO:HB2	1:D:420:ILE:HD13	1.94	0.49
1:D:13:VAL:HG22	1:D:34:ILE:HG13	1.93	0.49
1:D:165:LEU:HD11	1:D:327:PHE:CZ	2.48	0.49
1:B:5:LYS:HB2	1:B:75:LEU:HD23	1.95	0.48
1:C:77:GLN:CB	1:C:380:ILE:HD13	2.43	0.48
1:C:28:GLN:HE21	1:D:154:ARG:HH22	1.61	0.48
1:C:401:ARG:NH2	1:D:25:THR:O	2.47	0.48
1:C:3:THR:CG2	1:C:6:PHE:HD1	2.27	0.48
1:C:165:LEU:HD13	1:C:328:LEU:HD22	1.96	0.48
1:D:152:HIS:HD2	3:D:730:HOH:O	1.97	0.48
1:A:154:ARG:HH22	1:B:28:GLN:HE21	1.61	0.47
1:C:13:VAL:HG22	1:C:34:ILE:HG13	1.96	0.47
1:C:25:THR:O	1:D:401:ARG:NH2	2.47	0.47
1:C:5:LYS:HB2	1:C:75:LEU:HD23	1.96	0.47
1:A:13:VAL:HG22	1:A:34:ILE:HG13	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:THR:HG21	1:C:6:PHE:CD1	2.49	0.47
1:D:367:MET:N	1:D:368:PRO:CD	2.76	0.47
1:D:35:HIS:HD2	1:D:64:LYS:HZ1	1.61	0.47
1:A:367:MET:N	1:A:368:PRO:CD	2.78	0.46
1:A:73:GLU:HG3	1:A:372:VAL:HG12	1.96	0.46
1:A:327:PHE:HE1	2:A:501:MOA:C12	2.29	0.46
1:B:35:HIS:HD2	1:B:64:LYS:HZ2	1.62	0.46
1:A:77:GLN:HB2	1:A:380:ILE:HD13	1.96	0.46
1:D:77:GLN:HB2	1:D:380:ILE:HD13	1.98	0.46
1:B:207:MET:HE3	1:B:215:LEU:HD22	1.98	0.46
1:B:207:MET:HE3	1:B:215:LEU:CD2	2.46	0.46
1:A:25:THR:O	1:B:401:ARG:NH2	2.49	0.46
1:C:35:HIS:HD2	1:C:64:LYS:HZ2	1.63	0.46
1:D:73:GLU:HG3	1:D:372:VAL:HG12	1.98	0.45
1:C:56:ALA:HA	1:C:136:ILE:O	2.16	0.45
1:A:56:ALA:HA	1:A:136:ILE:O	2.17	0.45
1:B:56:ALA:HA	1:B:136:ILE:O	2.17	0.45
1:C:301:ARG:HG2	2:C:501:MOA:O1	2.17	0.45
1:D:56:ALA:HA	1:D:136:ILE:O	2.17	0.45
1:B:38:GLN:HE22	1:B:128:GLN:NE2	2.12	0.45
1:A:79:SER:OG	1:A:384:GLN:OE1	2.34	0.44
1:A:391:ARG:O	1:A:395:ARG:HG3	2.17	0.44
1:D:201:ARG:NH2	1:D:204:ARG:HG3	2.32	0.44
1:D:77:GLN:CB	1:D:380:ILE:HD13	2.47	0.44
1:A:50:ASP:OD1	1:A:50:ASP:N	2.50	0.43
1:A:152:HIS:HD2	3:A:694:HOH:O	1.99	0.43
1:D:5:LYS:HB2	1:D:75:LEU:HD23	1.99	0.43
1:B:337:VAL:HG11	1:B:357:GLU:HB3	2.01	0.43
1:D:271:ASN:O	1:D:271:ASN:CG	2.61	0.43
1:C:48:PRO:HG2	1:C:82:HIS:CD2	2.53	0.43
1:B:301:ARG:HG2	2:B:501:MOA:O1	2.18	0.43
1:C:154:ARG:HH22	1:D:28:GLN:NE2	2.16	0.43
1:B:48:PRO:HG2	1:B:82:HIS:CG	2.54	0.43
1:C:207:MET:CE	1:C:215:LEU:HD22	2.44	0.43
1:C:154:ARG:HH12	1:D:28:GLN:NE2	2.10	0.43
1:D:325:ALA:O	1:D:363:PHE:HA	2.18	0.43
1:A:28:GLN:HE21	1:B:154:ARG:HH22	1.65	0.43
1:A:28:GLN:NE2	1:B:154:ARG:HH12	2.14	0.42
1:A:77:GLN:CB	1:A:380:ILE:HD13	2.49	0.42
1:A:350:VAL:N	1:A:351:PRO:CD	2.83	0.42
1:C:337:VAL:HG11	1:C:357:GLU:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:THR:HB	1:B:8:ILE:HD11	2.02	0.42
1:B:139:ALA:HB2	1:B:365:HIS:CE1	2.55	0.42
1:A:48:PRO:HG2	1:A:82:HIS:CG	2.54	0.42
1:C:367:MET:N	1:C:368:PRO:CD	2.83	0.42
1:B:350:VAL:N	1:B:351:PRO:CD	2.83	0.42
1:C:38:GLN:HE22	1:C:128:GLN:NE2	2.15	0.42
1:A:271:ASN:CG	1:A:271:ASN:O	2.62	0.41
1:B:207:MET:CE	1:B:215:LEU:HD22	2.51	0.41
1:A:397:TRP:CZ2	1:A:401:ARG:HD2	2.55	0.41
1:B:139:ALA:HB1	2:B:501:MOA:C6	2.50	0.41
1:C:350:VAL:N	1:C:351:PRO:CD	2.83	0.41
1:A:327:PHE:HE2	1:A:365:HIS:ND1	2.19	0.41
1:B:271:ASN:CG	1:B:271:ASN:O	2.64	0.41
1:D:57:THR:CG2	1:D:118:LEU:HD11	2.51	0.41
1:D:350:VAL:N	1:D:351:PRO:CD	2.84	0.41
1:D:418:LYS:HB3	1:D:418:LYS:HE3	1.98	0.41
1:A:5:LYS:HB2	1:A:75:LEU:HD23	2.00	0.41
1:B:367:MET:N	1:B:368:PRO:CD	2.84	0.41
1:C:397:TRP:CZ2	1:C:401:ARG:HD2	2.56	0.41
1:A:327:PHE:CD1	1:A:327:PHE:C	2.99	0.41
1:C:48:PRO:HG2	1:C:82:HIS:CG	2.56	0.40
1:A:57:THR:CG2	1:A:118:LEU:HD11	2.52	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	420/433 (97%)	406 (97%)	13 (3%)	1 (0%)	43 34
1	B	416/433 (96%)	406 (98%)	9 (2%)	1 (0%)	43 34
1	C	416/433 (96%)	407 (98%)	8 (2%)	1 (0%)	43 34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	D	416/433 (96%)	401 (96%)	14 (3%)	1 (0%)	43 34
All	All	1668/1732 (96%)	1620 (97%)	44 (3%)	4 (0%)	43 34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ILE
1	B	18	ILE
1	C	18	ILE
1	D	18	ILE

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	369/378 (98%)	362 (98%)	7 (2%)	50 34
1	B	365/378 (97%)	361 (99%)	4 (1%)	65 54
1	C	365/378 (97%)	354 (97%)	11 (3%)	36 19
1	D	366/378 (97%)	356 (97%)	10 (3%)	39 22
All	All	1465/1512 (97%)	1433 (98%)	32 (2%)	45 29

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	140	PHE
1	A	268	GLN
1	A	330	ILE
1	A	339	ILE
1	A	357	GLU
1	A	359	VAL
1	A	418	LYS
1	B	140	PHE
1	B	163	ASP
1	B	165	LEU

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Mol	Chain	Res	Type
1	B	420	ILE
1	C	3	THR
1	C	96	GLN
1	C	140	PHE
1	C	160	LEU
1	C	163	ASP
1	C	165	LEU
1	C	204	ARG
1	C	208	GLN
1	C	268	GLN
1	C	385	GLN
1	C	420	ILE
1	D	1	MET
1	D	4	GLU
1	D	140	PHE
1	D	218	MET
1	D	268	GLN
1	D	330	ILE
1	D	360	LEU
1	D	367	MET
1	D	371	GLU
1	D	418	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	GLN
1	A	35	HIS
1	A	77	GLN
1	A	128	GLN
1	A	138	HIS
1	A	147	ASN
1	A	152	HIS
1	A	184	ASN
1	A	268	GLN
1	B	28	GLN
1	B	35	HIS
1	B	77	GLN
1	B	96	GLN
1	B	128	GLN
1	B	138	HIS
1	B	147	ASN

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Mol	Chain	Res	Type
1	B	152	HIS
1	B	184	ASN
1	B	208	GLN
1	B	221	HIS
1	C	28	GLN
1	C	35	HIS
1	C	77	GLN
1	C	82	HIS
1	C	95	ASN
1	C	96	GLN
1	C	128	GLN
1	C	138	HIS
1	C	147	ASN
1	C	152	HIS
1	C	184	ASN
1	C	208	GLN
1	C	221	HIS
1	C	268	GLN
1	D	27	ASN
1	D	28	GLN
1	D	35	HIS
1	D	77	GLN
1	D	100	HIS
1	D	128	GLN
1	D	138	HIS
1	D	147	ASN
1	D	152	HIS
1	D	184	ASN
1	D	268	GLN
1	D	365	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 oligosaccharides 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	MOA	D	501	-	24,24,24	3.50	10 (41%)	34,34,34	2.34	15 (44%)
2	MOA	A	501	-	24,24,24	3.18	8 (33%)	34,34,34	2.71	16 (47%)
2	MOA	C	501	-	24,24,24	2.85	9 (37%)	34,34,34	2.59	11 (32%)
2	MOA	B	501	-	24,24,24	2.72	9 (37%)	34,34,34	2.70	11 (32%)

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	MOA	D	501	-	-	5/12/21/21	0/2/2/2
2	MOA	A	501	-	-	2/12/21/21	0/2/2/2
2	MOA	C	501	-	-	0/12/21/21	0/2/2/2
2	MOA	B	501	-	-	0/12/21/21	0/2/2/2

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	MOA	O2-C1	8.17	1.46	1.36
2	A	501	MOA	O2-C1	7.97	1.46	1.36
2	D	501	MOA	C13-C14	7.22	1.51	1.39
2	A	501	MOA	C13-C12	7.11	1.52	1.39
2	D	501	MOA	C13-C12	6.95	1.51	1.39
2	D	501	MOA	C15-C14	6.75	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	MOA	C15-C14	6.04	1.48	1.40
2	C	501	MOA	C13-C12	6.01	1.50	1.39
2	B	501	MOA	C13-C12	5.69	1.49	1.39
2	C	501	MOA	O2-C1	5.50	1.43	1.36
2	B	501	MOA	C13-C14	5.43	1.48	1.39
2	C	501	MOA	C13-C14	5.31	1.48	1.39
2	A	501	MOA	C13-C14	5.25	1.48	1.39
2	B	501	MOA	C15-C14	5.20	1.47	1.40
2	C	501	MOA	C15-C14	4.81	1.47	1.40
2	B	501	MOA	O2-C1	4.80	1.42	1.36
2	B	501	MOA	C11-C12	4.68	1.48	1.40
2	D	501	MOA	C16-C15	4.57	1.50	1.41
2	A	501	MOA	C16-C15	4.41	1.50	1.41
2	C	501	MOA	C11-C12	4.40	1.47	1.40
2	D	501	MOA	C11-C12	4.30	1.47	1.40
2	A	501	MOA	C11-C12	4.13	1.47	1.40
2	B	501	MOA	C16-C11	3.94	1.46	1.39
2	C	501	MOA	C16-C11	3.86	1.46	1.39
2	D	501	MOA	C16-C11	3.71	1.45	1.39
2	C	501	MOA	C16-C15	3.39	1.48	1.41
2	A	501	MOA	C16-C11	3.33	1.45	1.39
2	A	501	MOA	O6-C6	3.27	1.32	1.22
2	B	501	MOA	C10-C11	2.51	1.53	1.50
2	C	501	MOA	C10-C11	2.47	1.53	1.50
2	B	501	MOA	C16-C15	2.47	1.46	1.41
2	D	501	MOA	O6-C6	2.41	1.30	1.22
2	D	501	MOA	C2-C3	2.15	1.38	1.33
2	D	501	MOA	C4-C3	2.13	1.55	1.51
2	B	501	MOA	C16-C1	-2.04	1.44	1.47
2	C	501	MOA	C17-C14	2.03	1.53	1.51

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	MOA	C15-C16-C11	6.81	127.45	121.87
2	C	501	MOA	O2-C1-O1	6.63	128.08	121.09
2	B	501	MOA	O2-C1-O1	6.10	127.53	121.09
2	C	501	MOA	C15-C16-C11	5.86	126.67	121.87
2	C	501	MOA	C16-C11-C12	-5.84	117.60	122.68
2	C	501	MOA	O1-C1-C16	-5.76	120.73	130.99
2	B	501	MOA	O1-C1-C16	-5.62	120.97	130.99
2	A	501	MOA	O2-C1-O1	5.00	126.37	121.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	MOA	C16-C11-C12	-4.97	118.36	122.68
2	A	501	MOA	C14-C17-C2	-4.89	103.72	112.06
2	A	501	MOA	C15-C16-C11	4.56	125.60	121.87
2	D	501	MOA	C14-C17-C2	4.55	119.83	112.06
2	D	501	MOA	C11-C16-C1	-4.30	103.59	108.39
2	A	501	MOA	C9-C3-C4	4.13	122.39	115.23
2	A	501	MOA	C16-C15-C14	-4.10	114.05	121.63
2	A	501	MOA	O1-C1-C16	-3.99	123.89	130.99
2	A	501	MOA	C11-C16-C1	-3.92	104.01	108.39
2	B	501	MOA	O2-C1-C16	3.90	111.33	108.28
2	D	501	MOA	O1-C1-C16	-3.89	124.06	130.99
2	D	501	MOA	O2-C1-O1	3.86	125.16	121.09
2	B	501	MOA	C16-C15-C14	-3.84	114.54	121.63
2	D	501	MOA	C17-C14-C15	-3.66	115.07	120.85
2	C	501	MOA	O2-C1-C16	3.52	111.03	108.28
2	A	501	MOA	O3-C13-C12	3.42	123.75	118.86
2	A	501	MOA	C10-C11-C16	3.29	112.13	107.90
2	A	501	MOA	C4-C3-C2	-3.25	113.87	121.17
2	C	501	MOA	C17-C2-C3	-3.23	122.78	127.42
2	D	501	MOA	O2-C1-C16	3.21	110.79	108.28
2	D	501	MOA	C10-C11-C16	3.11	111.90	107.90
2	A	501	MOA	C7-C12-C13	3.05	126.01	121.24
2	D	501	MOA	C16-C15-C14	-3.00	116.08	121.63
2	B	501	MOA	C11-C16-C1	-2.99	105.05	108.39
2	C	501	MOA	C16-C15-C14	-2.95	116.18	121.63
2	A	501	MOA	C15-C14-C13	2.93	121.59	117.65
2	B	501	MOA	C14-C17-C2	-2.91	107.09	112.06
2	D	501	MOA	C15-C16-C11	2.85	124.20	121.87
2	A	501	MOA	O5-C6-O6	2.83	130.62	123.33
2	D	501	MOA	C9-C3-C2	-2.83	116.36	123.63
2	A	501	MOA	O6-C6-C5	-2.79	114.24	123.09
2	B	501	MOA	O6-C6-C5	-2.77	114.30	123.09
2	C	501	MOA	O3-C13-C14	-2.76	113.57	118.82
2	B	501	MOA	C17-C2-C3	-2.69	123.57	127.42
2	A	501	MOA	O3-C13-C14	-2.61	113.85	118.82
2	D	501	MOA	C7-C12-C13	2.61	125.33	121.24
2	C	501	MOA	C14-C17-C2	-2.57	107.66	112.06
2	C	501	MOA	C11-C16-C1	-2.48	105.62	108.39
2	A	501	MOA	C16-C11-C12	-2.46	120.55	122.68
2	D	501	MOA	O6-C6-C5	-2.43	115.38	123.09
2	D	501	MOA	C16-C11-C12	-2.30	120.68	122.68
2	D	501	MOA	C7-C12-C11	-2.18	117.29	120.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	MOA	C15-C16-C1	2.12	132.21	129.35
2	B	501	MOA	O5-C6-C5	2.12	120.68	114.00
2	C	501	MOA	O3-C13-C12	2.04	121.77	118.86

There are no chirality outliers.

All (7) torsion outliers are listed below:

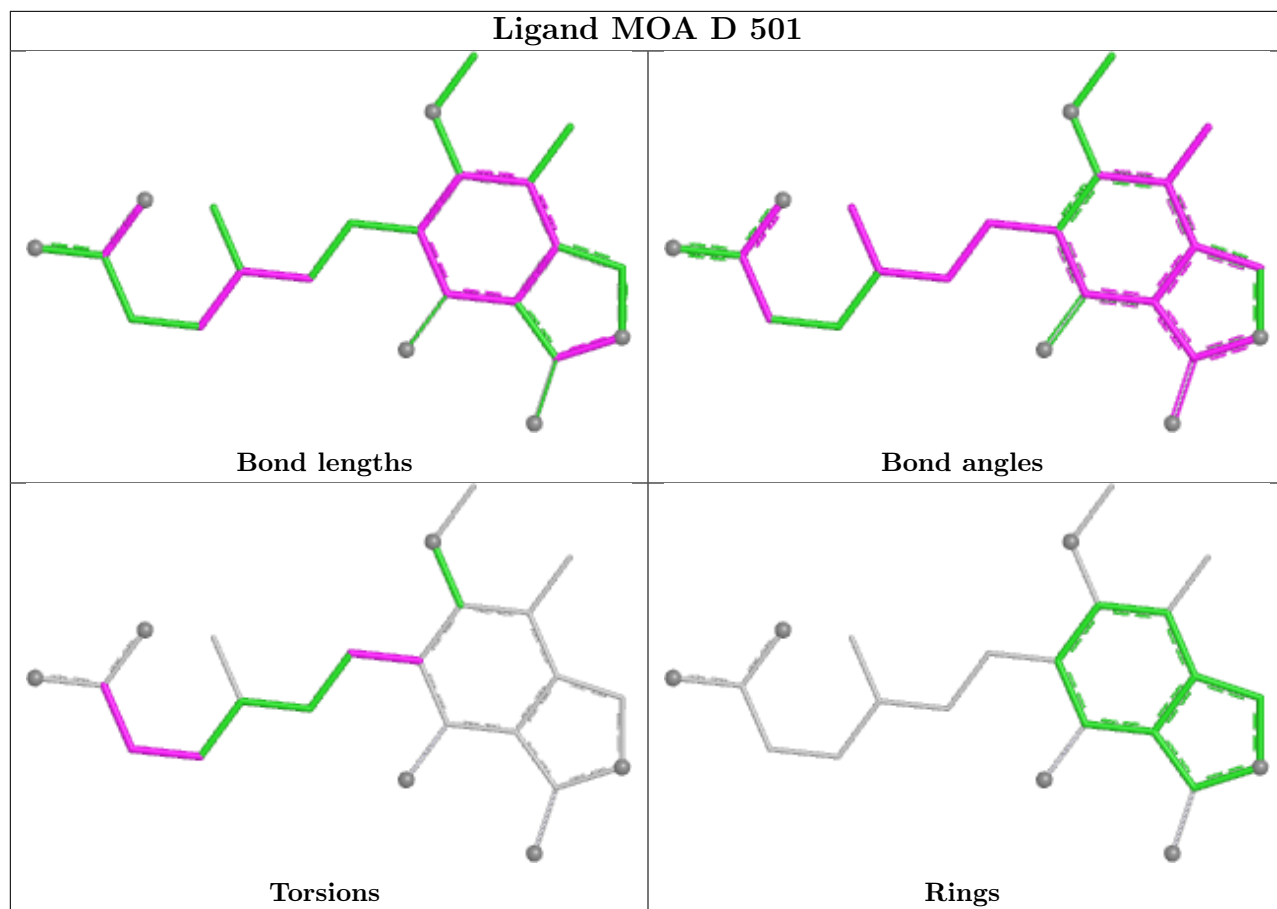
Mol	Chain	Res	Type	Atoms
2	D	501	MOA	C13-C14-C17-C2
2	D	501	MOA	C15-C14-C17-C2
2	D	501	MOA	C3-C4-C5-C6
2	D	501	MOA	C4-C5-C6-O5
2	D	501	MOA	C4-C5-C6-O6
2	A	501	MOA	C4-C5-C6-O6
2	A	501	MOA	C4-C5-C6-O5

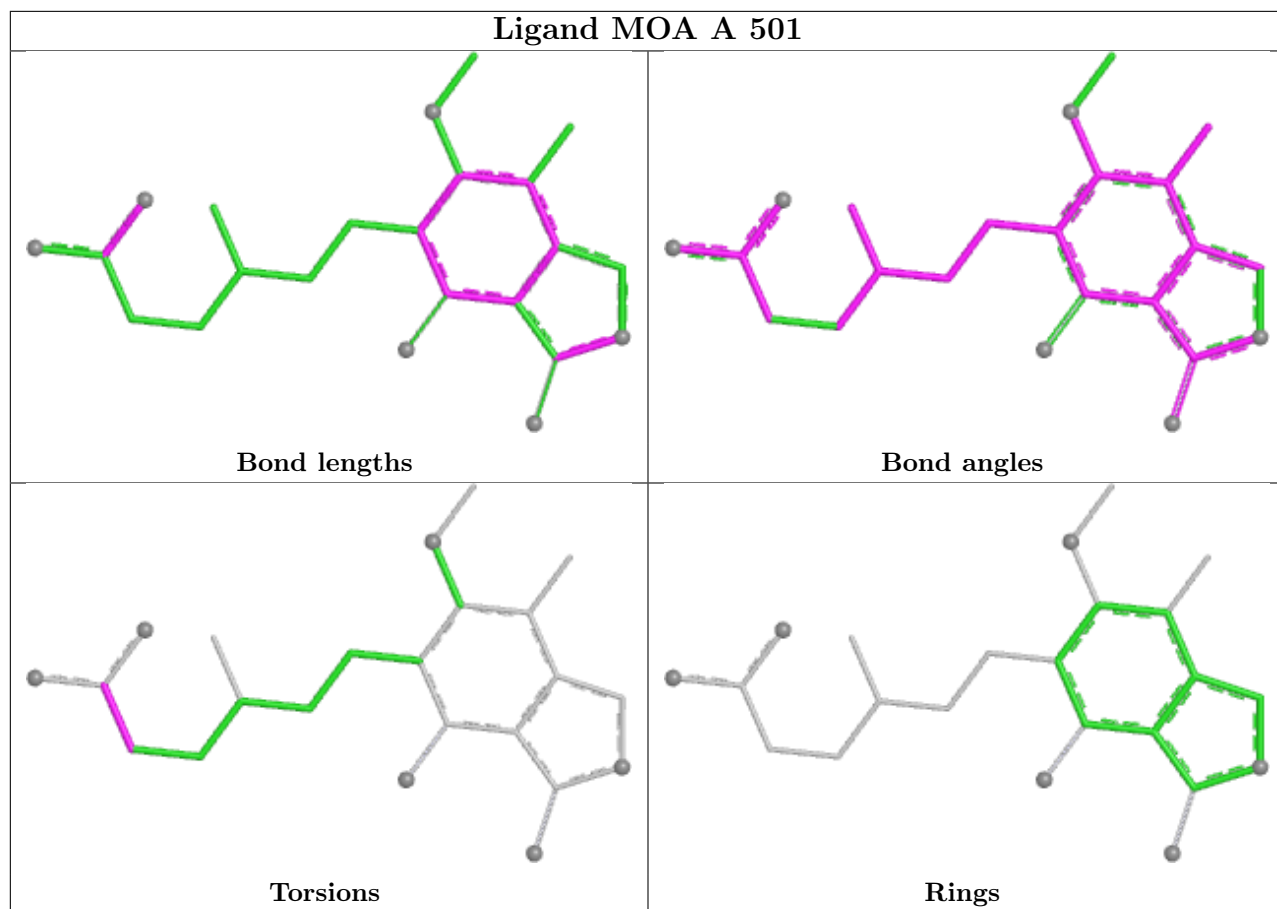
There are no ring outliers.

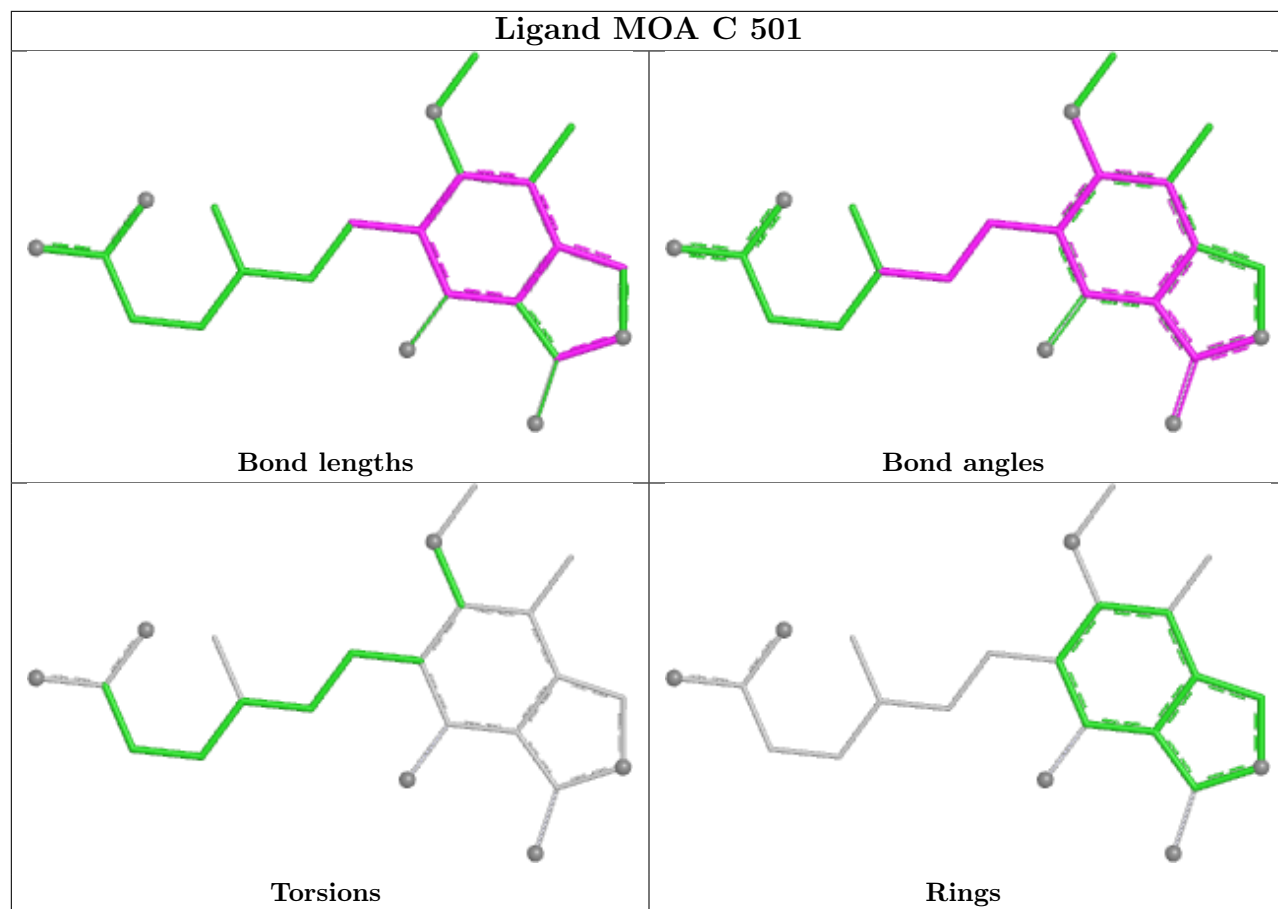
4 monomers are involved in 10 short contacts:

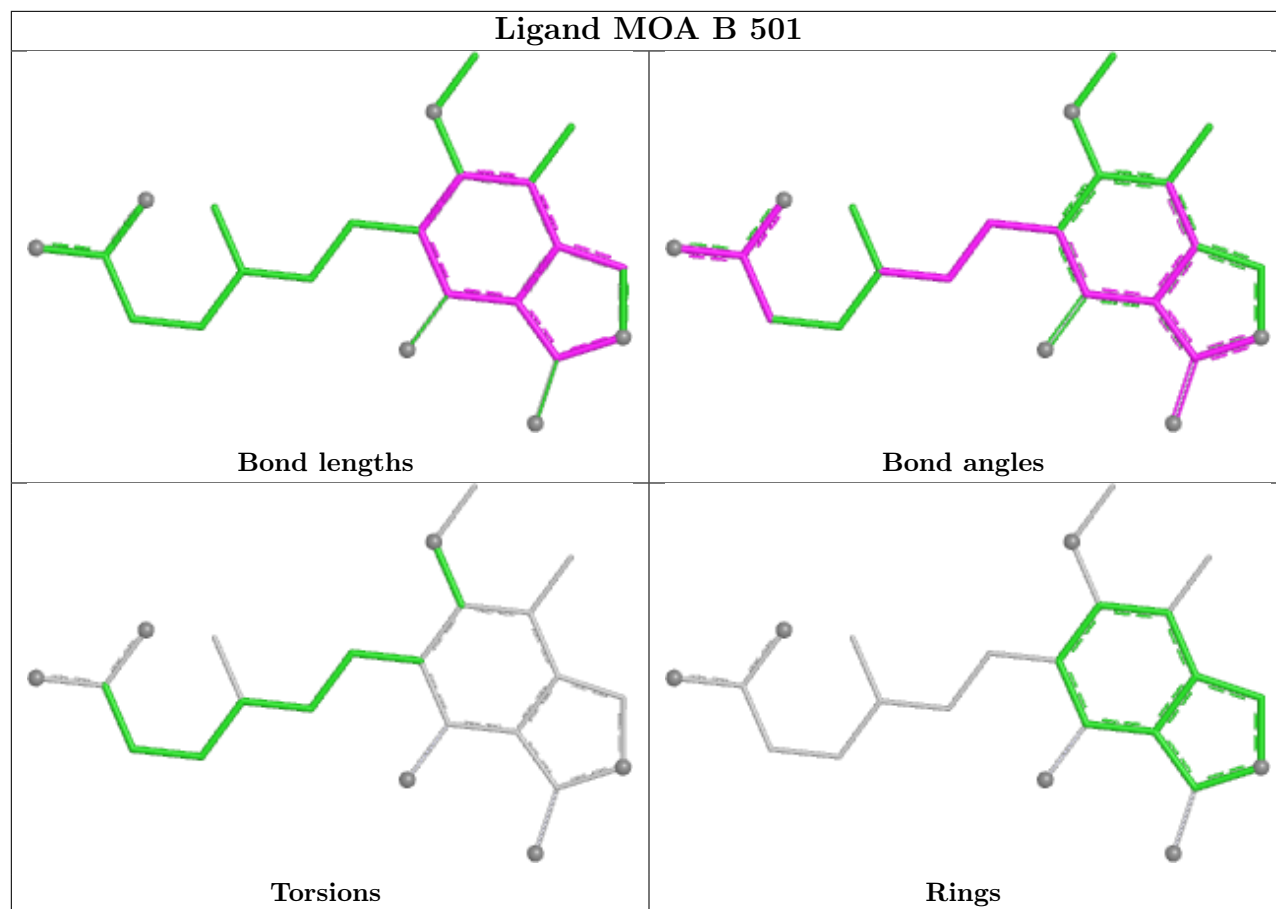
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	MOA	2	0
2	A	501	MOA	5	0
2	C	501	MOA	1	0
2	B	501	MOA	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	422/433 (97%)	0.24	30 (7%) 22 23	17, 24, 48, 76	0
1	B	418/433 (96%)	0.05	17 (4%) 41 44	16, 23, 48, 94	0
1	C	418/433 (96%)	0.02	18 (4%) 40 42	15, 23, 45, 86	0
1	D	420/433 (96%)	0.09	19 (4%) 38 40	16, 23, 43, 67	0
All	All	1678/1732 (96%)	0.10	84 (5%) 34 36	15, 24, 47, 94	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	360	LEU	6.6
1	D	365	HIS	6.0
1	A	359	VAL	6.0
1	A	364	GLY	5.4
1	B	3	THR	5.0
1	A	363	PHE	4.9
1	D	423	GLY	4.5
1	C	420	ILE	4.3
1	D	327	PHE	4.2
1	B	277	THR	4.1
1	A	206	ILE	4.1
1	B	420	ILE	4.0
1	D	364	GLY	4.0
1	B	4	GLU	4.0
1	C	4	GLU	4.0
1	A	365	HIS	4.0
1	B	248	THR	3.8
1	B	275	PRO	3.8
1	D	403	GLY	3.7
1	C	403	GLY	3.7
1	A	361	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	277	THR	3.6
1	A	327	PHE	3.6
1	A	4	GLU	3.5
1	D	360	LEU	3.5
1	D	359	VAL	3.3
1	C	275	PRO	3.3
1	A	403	GLY	3.2
1	A	422	SER	3.2
1	C	276	LYS	3.2
1	D	1	MET	3.1
1	C	248	THR	3.1
1	D	324	GLY	3.0
1	D	206	ILE	3.0
1	D	4	GLU	3.0
1	C	5	LYS	2.9
1	A	397	TRP	2.9
1	D	2	SER	2.9
1	A	275	PRO	2.8
1	D	363	PHE	2.8
1	A	324	GLY	2.8
1	B	5	LYS	2.8
1	A	248	THR	2.7
1	A	79	SER	2.6
1	B	276	LYS	2.6
1	C	46	PRO	2.6
1	C	204	ARG	2.5
1	D	185	TYR	2.5
1	A	396	GLN	2.5
1	C	244	THR	2.5
1	C	274	ASP	2.5
1	B	45	GLY	2.5
1	C	45	GLY	2.5
1	B	247	SER	2.5
1	B	274	ASP	2.5
1	A	80	GLY	2.4
1	A	395	ARG	2.4
1	A	1	MET	2.4
1	B	397	TRP	2.4
1	A	7	THR	2.4
1	A	5	LYS	2.4
1	A	185	TYR	2.3
1	A	2	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	325	ALA	2.3
1	B	278	GLY	2.3
1	C	3	THR	2.3
1	C	249	THR	2.3
1	D	277	THR	2.3
1	B	279	ARG	2.2
1	D	239	ALA	2.2
1	D	47	VAL	2.2
1	A	82	HIS	2.2
1	C	245	ALA	2.2
1	A	373	LYS	2.2
1	B	208	GLN	2.2
1	A	366	LEU	2.1
1	D	48	PRO	2.1
1	B	245	ALA	2.1
1	B	403	GLY	2.1
1	A	77	GLN	2.1
1	A	3	THR	2.0
1	D	49	ASP	2.0
1	C	247	SER	2.0
1	C	7	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

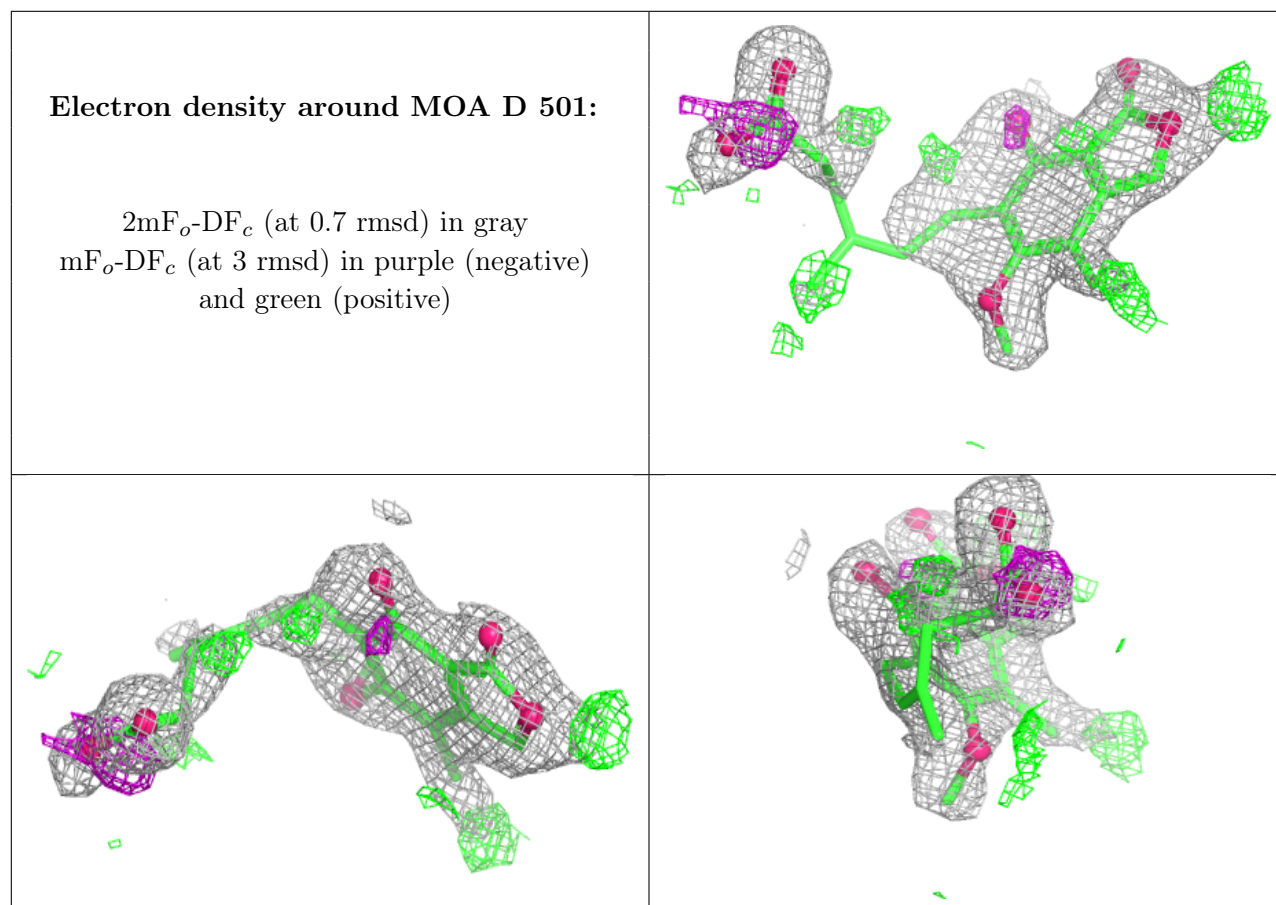
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MOA	D	501	23/23	0.82	0.20	20,57,63,70	0
2	MOA	A	501	23/23	0.83	0.18	20,52,60,68	0

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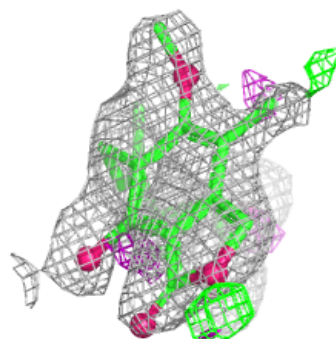
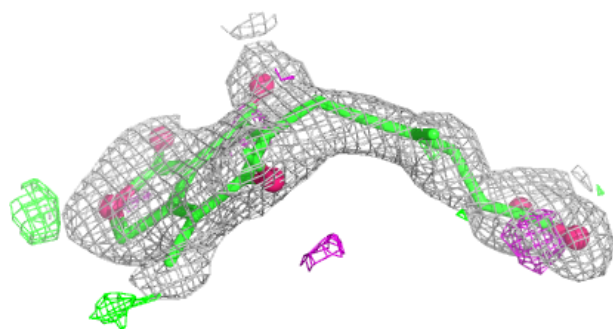
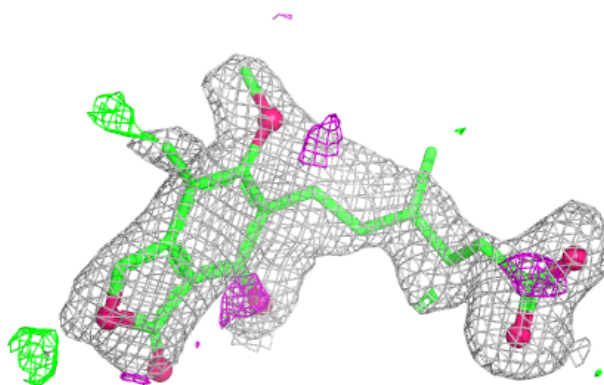
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MOA	C	501	23/23	0.93	0.08	21,28,34,38	0
2	MOA	B	501	23/23	0.93	0.09	21,30,35,38	0

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.

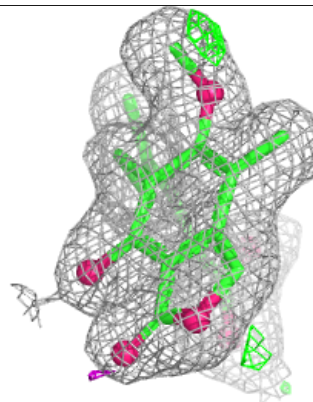
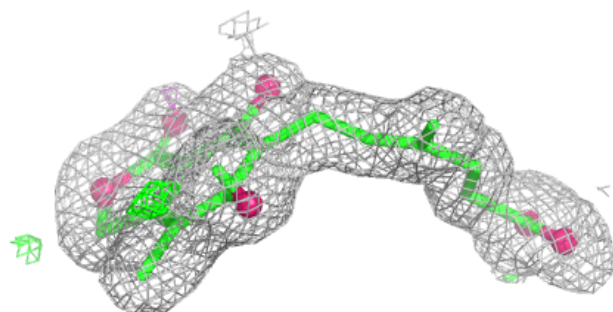
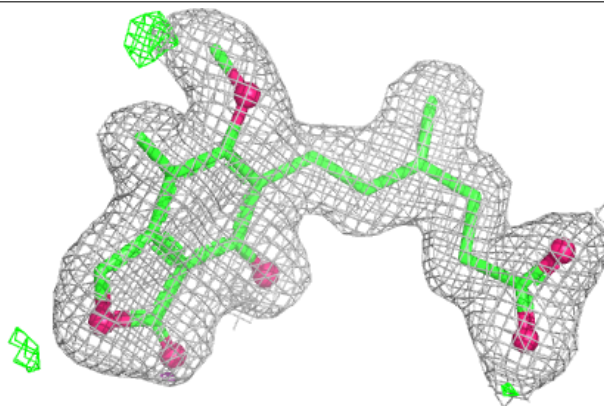


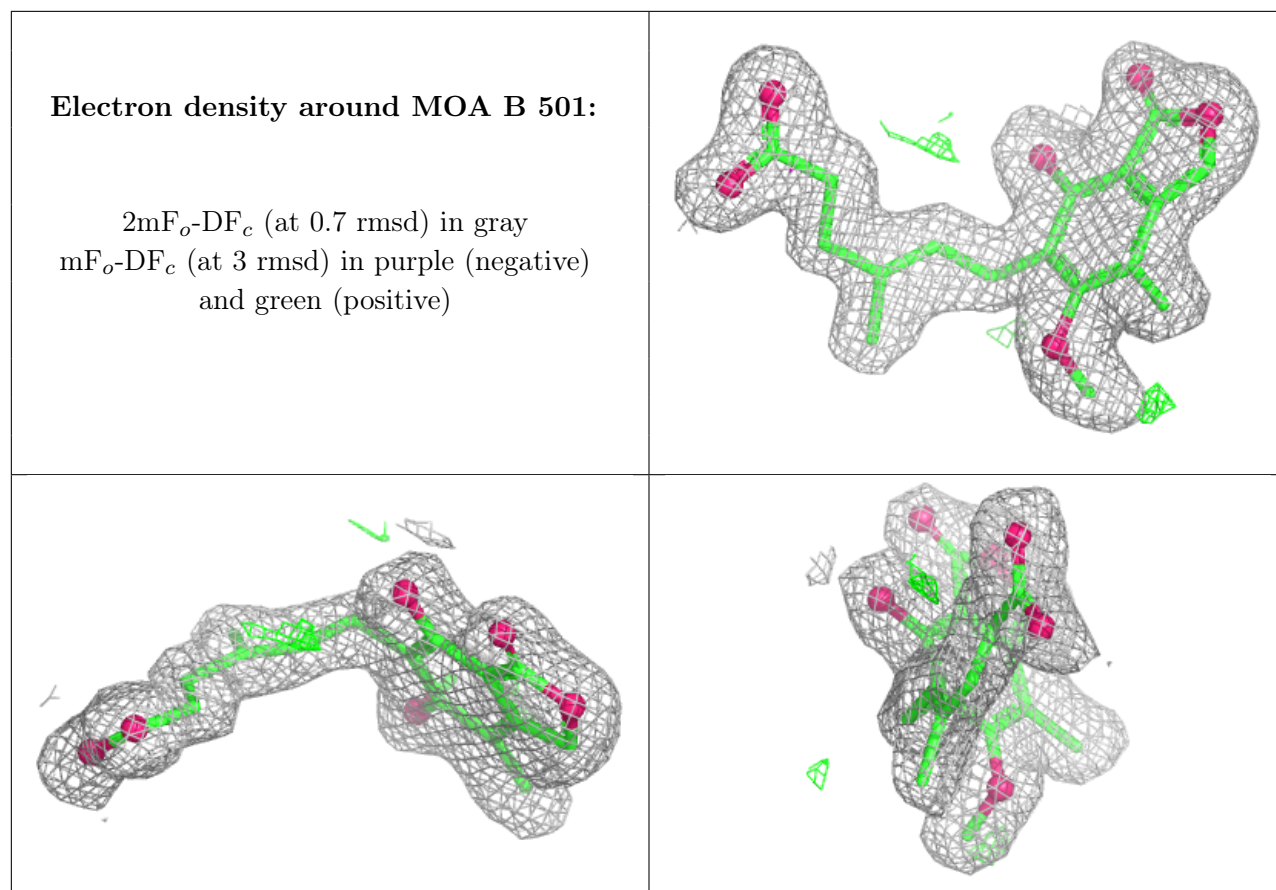
Electron density around MOA A 501:

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

**Electron density around MOA C 501:**

$2mF_o-DF_c$ (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.