



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 19, 2019 – 09:42 PM EDT

PDB ID : 6T6D
Title : Crystal structure of the ACVR1 (ALK2) kinase in complex with the compound M4K2149
Authors : Adamson, R.J.; Williams, E.P.; Smil, D.; Burgess-Brown, N.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Bullock, A.N.
Deposited on : 2019-10-18
Resolution : 2.56 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 : 1.8.0 (224370), CSD as540be (2019)
Xtriage (Phenix) : 1.13
EDS : 2.4
buster-report : 1.1.7 (2018)
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.4

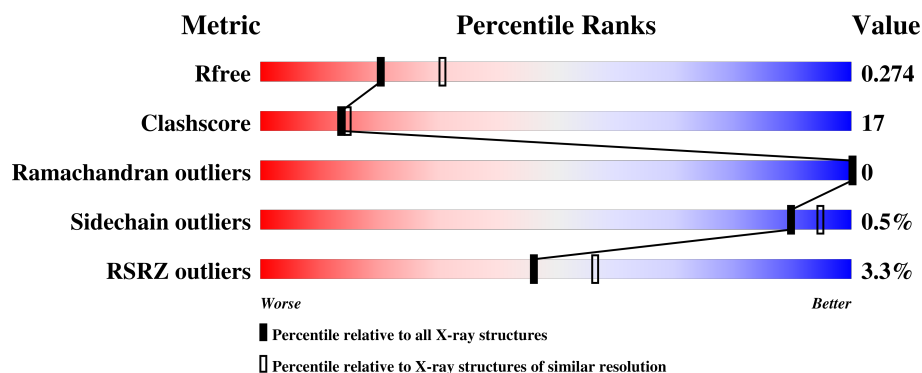
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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}	111664	1056 (2.58-2.54)
Clashscore	122126	1102 (2.58-2.54)
Ramachandran outliers	120053	1092 (2.58-2.54)
Sidechain outliers	120020	1092 (2.58-2.54)
RSRZ outliers	108989	1048 (2.58-2.54)

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. 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	301	<div> <div>%</div> <div> <div></div> <div>68%</div> <div>29%</div> <div>.</div> </div> </div>
1	B	301	<div> <div>3%</div> <div> <div></div> <div>70%</div> <div>29%</div> <div>.</div> </div> </div>
1	C	301	<div> <div>5%</div> <div> <div></div> <div>66%</div> <div>26%</div> <div>8%</div> </div> </div>
1	D	301	<div> <div>4%</div> <div> <div></div> <div>61%</div> <div>33%</div> <div>5%</div> </div> </div>

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	SO4	A	504	-	-	X	-
3	SO4	B	503	-	-	X	-
3	SO4	D	503	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9250 atoms, of which 104 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.

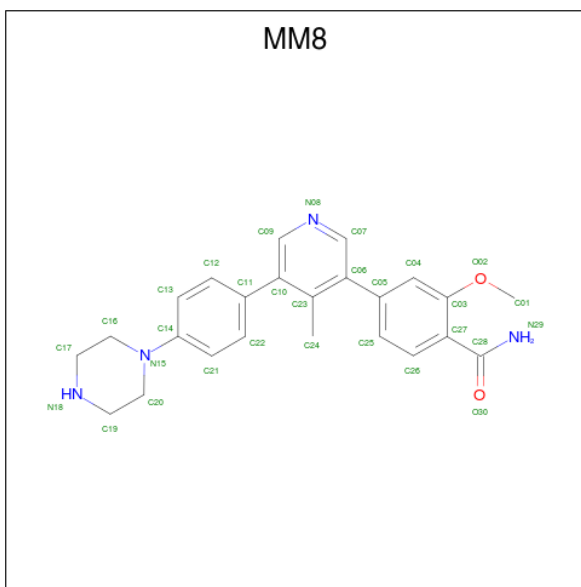
- Molecule 1 is a protein called Activin receptor type I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	1	0
			2282	1460	386	421	15			
1	B	297	Total	C	N	O	S	0	1	0
			2288	1464	393	417	14			
1	C	278	Total	C	N	O	S	0	0	0
			2136	1371	365	386	14			
1	D	285	Total	C	N	O	S	0	2	0
			2235	1430	378	412	15			

There are 12 discrepancies between the modelled and reference sequences:

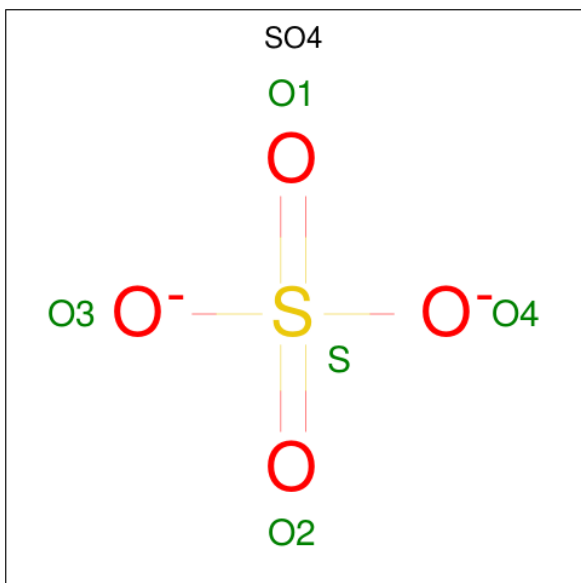
Chain	Residue	Modelled	Actual	Comment	Reference
A	199	SER	-	expression tag	UNP Q04771
A	200	MET	-	expression tag	UNP Q04771
A	207	ASP	GLN	engineered mutation	UNP Q04771
B	199	SER	-	expression tag	UNP Q04771
B	200	MET	-	expression tag	UNP Q04771
B	207	ASP	GLN	engineered mutation	UNP Q04771
C	199	SER	-	expression tag	UNP Q04771
C	200	MET	-	expression tag	UNP Q04771
C	207	ASP	GLN	engineered mutation	UNP Q04771
D	199	SER	-	expression tag	UNP Q04771
D	200	MET	-	expression tag	UNP Q04771
D	207	ASP	GLN	engineered mutation	UNP Q04771

- Molecule 2 is 2-methoxy-4-[4-methyl-5-(4-piperazin-1-ylphenyl)pyridin-3-yl]benzamide (three-letter code: MM8) (formula: C₂₄H₂₆N₄O₂) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			56	24	26	4	2		
2	B	1	Total	C	H	N	O	0	0
			56	24	26	4	2		
2	C	1	Total	C	H	N	O	0	0
			56	24	26	4	2		
2	D	1	Total	C	H	N	O	0	0
			56	24	26	4	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

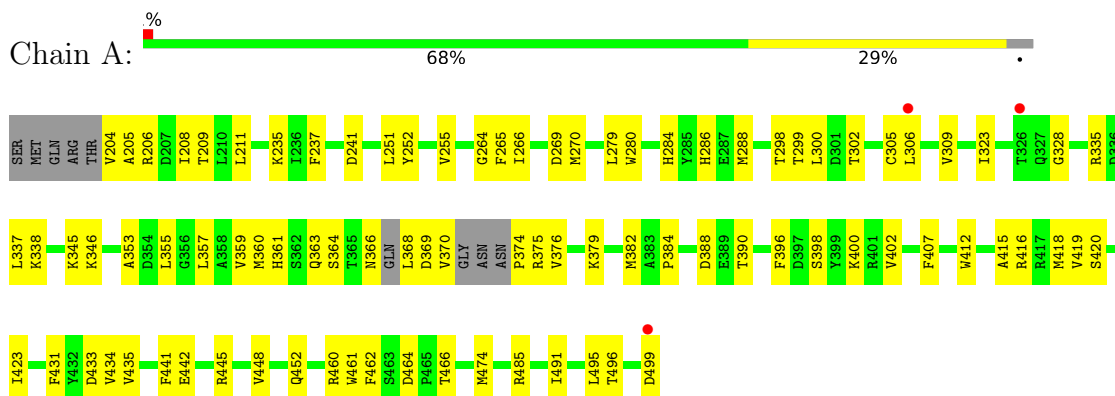
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	16	Total O 16 16	0	0
4	B	4	Total O 4 4	0	0
4	C	4	Total O 4 4	0	0
4	D	6	Total O 6 6	0	0

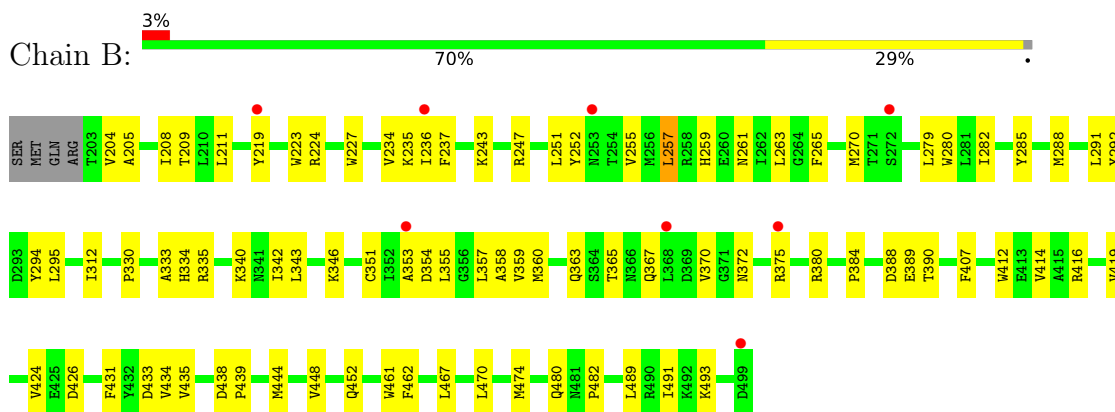
3 Residue-property plots [i](#)

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.

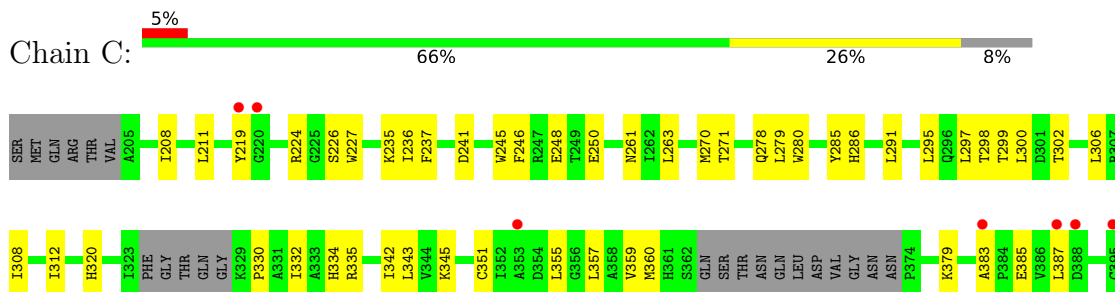
• Molecule 1: Activin receptor type I

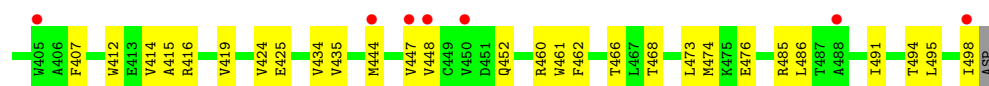


• Molecule 1: Activin receptor type I

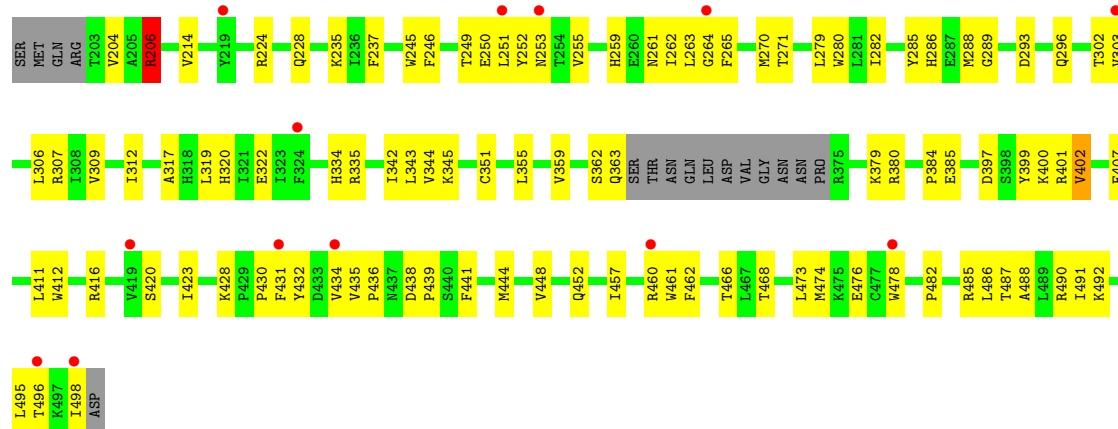


• Molecule 1: Activin receptor type I





● Molecule 1: Activin receptor type I



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	198.09Å 101.88Å 84.27Å 90.00° 111.43° 90.00°	Depositor
Resolution (Å)	78.45 – 2.56 89.18 – 2.56	Depositor EDS
% Data completeness (in resolution range)	99.3 (78.45-2.56) 99.7 (89.18-2.56)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.55Å)	Xtriage
Refinement program	PHENIX 1.16_3549, PHENIX 1.16_3549	Depositor
R, R_{free}	0.222 , 0.272 0.221 , 0.274	Depositor DCC
R_{free} test set	2493 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	56.0	Xtriage
Anisotropy	0.941	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 57.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9250	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: MM8, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.24	0/2332	0.42	0/3171
1	B	0.23	0/2340	0.42	0/3191
1	C	0.23	0/2185	0.41	0/2977
1	D	0.24	0/2288	0.41	0/3113
All	All	0.23	0/9145	0.41	0/12452

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	D	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	206	ARG	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	2282	0	2222	79	0
1	B	2288	0	2196	77	0
1	C	2136	0	2053	64	0
1	D	2235	0	2167	91	0
2	A	30	26	0	0	0
2	B	30	26	0	0	0
2	C	30	26	0	0	0
2	D	30	26	0	4	0
3	A	15	0	0	3	0
3	B	15	0	0	2	0
3	C	15	0	0	0	0
3	D	10	0	0	3	0
4	A	16	0	0	0	0
4	B	4	0	0	0	0
4	C	4	0	0	0	0
4	D	6	0	0	0	0
All	All	9146	104	8638	304	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 17.

All (304) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:VAL:HG23	1:A:206:ARG:H	1.26	0.99
1:A:302:THR:HA	1:A:418:MET:HE1	1.45	0.98
1:D:270:MET:HG3	1:D:279:LEU:HD23	1.48	0.96
1:B:270:MET:HG3	1:B:279:LEU:HD23	1.51	0.91
1:D:286:HIS:NE2	3:D:503:SO4:O1	2.06	0.88
1:A:360:MET:H	1:A:368:LEU:HD23	1.40	0.87
1:A:286:HIS:NE2	3:A:504:SO4:O3	2.08	0.86
1:C:419:VAL:HA	1:C:424:VAL:HG12	1.54	0.86
1:A:264:GLY:N	3:A:504:SO4:O1	2.08	0.85
1:A:270:MET:HG3	1:A:279:LEU:HD23	1.57	0.83
1:D:270:MET:HG3	1:D:279:LEU:CD2	2.07	0.83
1:B:234:VAL:HG12	1:B:282:ILE:HD12	1.61	0.81
1:D:264:GLY:N	3:D:503:SO4:O4	2.13	0.81
1:A:407:PHE:HE2	1:A:491:ILE:HG21	1.46	0.81
1:B:435:VAL:HG21	1:B:444:MET:HE1	1.61	0.80
1:D:462:PHE:HA	1:D:468:THR:CG2	2.11	0.80
1:D:312:ILE:HD13	1:D:342:ILE:HD13	1.64	0.79
1:B:489:LEU:HG	1:B:493:LYS:HE3	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:259:HIS:HD2	1:D:261:ASN:H	1.30	0.78
1:A:270:MET:HG3	1:A:279:LEU:CD2	2.13	0.78
1:B:270:MET:HG3	1:B:279:LEU:CD2	2.15	0.77
1:D:204:VAL:HG23	1:D:206:ARG:H	1.50	0.76
1:D:462:PHE:HA	1:D:468:THR:HG23	1.69	0.74
1:D:487:THR:HG23	1:D:490:ARG:H	1.53	0.74
1:B:416:ARG:NH1	1:B:426:ASP:O	2.20	0.73
1:A:208:ILE:HG21	1:A:280:TRP:HZ3	1.52	0.73
1:A:286:HIS:CE1	1:A:345:LYS:HG2	2.22	0.73
1:D:206:ARG:HH12	1:D:271:THR:HG22	1.54	0.72
1:D:462:PHE:HD1	1:D:468:THR:HG22	1.54	0.71
1:C:444:MET:O	1:C:448:VAL:HG12	1.91	0.71
1:D:420:SER:O	1:D:423:ILE:HG22	1.90	0.71
1:C:476:GLU:HB3	1:C:486:LEU:CD1	2.21	0.71
1:C:211:LEU:HD11	1:C:226:SER:OG	1.91	0.70
1:A:302:THR:O	1:A:306:LEU:HD13	1.90	0.70
1:A:375:ARG:HD2	1:A:441:PHE:CE1	2.26	0.70
1:D:476:GLU:HB3	1:D:486:LEU:CD1	2.22	0.70
1:A:496:THR:O	1:D:307:ARG:NH1	2.26	0.69
1:C:412:TRP:CZ2	1:C:416:ARG:HD2	2.28	0.68
1:D:259:HIS:CD2	1:D:261:ASN:H	2.11	0.68
1:B:412:TRP:O	1:B:416:ARG:HG3	1.94	0.68
1:C:219:TYR:HA	1:C:241:ASP:OD2	1.94	0.67
1:B:431:PHE:O	1:B:435:VAL:HG22	1.95	0.67
1:A:412:TRP:HD1	1:A:474:MET:HE1	1.59	0.67
1:C:419:VAL:CA	1:C:424:VAL:HG12	2.26	0.66
1:D:309:VAL:HG13	1:D:411:LEU:HD11	1.78	0.66
1:C:224:ARG:HD3	1:C:285:TYR:CZ	2.31	0.66
1:A:300:LEU:O	1:A:419:VAL:HG12	1.96	0.65
1:C:270:MET:HB2	1:C:279:LEU:HD23	1.78	0.65
1:D:206:ARG:HH12	1:D:271:THR:CG2	2.10	0.65
1:A:252:TYR:CG	1:A:265:PHE:HB2	2.32	0.64
1:B:288:MET:CE	1:B:346:LYS:HA	2.27	0.64
1:D:492:LYS:O	1:D:496:THR:HG23	1.97	0.64
1:D:462:PHE:CD1	1:D:468:THR:HG22	2.33	0.63
1:D:431:PHE:O	1:D:435:VAL:HG12	1.98	0.63
1:A:306:LEU:HD11	1:A:466:THR:CG2	2.29	0.63
1:C:476:GLU:HB3	1:C:486:LEU:HD13	1.78	0.63
1:A:390:THR:HG21	1:B:363:GLN:NE2	2.15	0.62
1:B:334:HIS:NE2	1:B:354:ASP:O	2.32	0.62
1:A:363:GLN:OE1	1:B:390:THR:HG21	2.00	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:435:VAL:HG22	1:D:436:PRO:HD2	1.82	0.61
1:A:418:MET:O	1:A:460:ARG:NH2	2.29	0.61
1:B:419:VAL:HG22	1:B:424:VAL:HB	1.81	0.61
1:A:357:LEU:HD13	1:A:376:VAL:HG13	1.83	0.61
1:B:439:PRO:HG3	1:B:444:MET:HE3	1.81	0.61
1:D:385:GLU:OE2	1:D:485:ARG:NH2	2.34	0.61
1:A:448:VAL:O	1:A:452:GLN:HA	2.01	0.60
1:D:476:GLU:HB3	1:D:486:LEU:HD11	1.83	0.60
1:C:237:PHE:HB2	1:C:279:LEU:HB2	1.83	0.60
1:A:209:THR:O	1:A:211:LEU:HD12	2.02	0.60
1:D:420:SER:HB3	1:D:460:ARG:HH11	1.67	0.60
1:D:252:TYR:CD1	1:D:265:PHE:HB2	2.38	0.59
1:D:466:THR:HG22	1:D:498:ILE:HG22	1.84	0.59
1:B:353:ALA:O	1:B:355:LEU:HD12	2.02	0.59
1:D:235:LYS:HD3	1:D:237:PHE:CZ	2.37	0.59
1:D:473:LEU:HD21	1:D:491:ILE:HG23	1.83	0.59
1:A:370:VAL:HG11	1:A:374:PRO:HD3	1.85	0.59
1:C:300:LEU:O	1:C:419:VAL:HG12	2.03	0.58
1:C:434:VAL:HG23	1:C:435:VAL:HG13	1.84	0.58
1:B:243:LYS:HD2	1:B:372:ASN:O	2.04	0.58
1:A:286:HIS:ND1	3:A:502:SO4:O2	2.36	0.58
1:B:236:ILE:HD12	1:B:280:TRP:NE1	2.19	0.58
1:B:461:TRP:CD1	1:B:467:LEU:HD13	2.39	0.58
1:D:435:VAL:CG2	1:D:439:PRO:HB3	2.33	0.58
1:A:305:CYS:HB3	1:A:418:MET:HE2	1.85	0.58
1:D:397:ASP:O	1:D:401:ARG:HG3	2.04	0.57
1:D:288:MET:HB2	1:D:344:VAL:O	2.04	0.57
1:D:206:ARG:HD3	1:D:280:TRP:CE3	2.39	0.57
1:D:286:HIS:ND1	3:D:502:SO4:O4	2.38	0.57
1:D:309:VAL:HG12	1:D:495:LEU:HD13	1.87	0.57
1:A:364:SER:O	1:A:366:ASN:ND2	2.38	0.56
1:D:214:VAL:HG11	2:D:501:MM8:C22	2.35	0.56
1:D:302:THR:O	1:D:306:LEU:HD23	2.05	0.56
1:B:435:VAL:HG21	1:B:444:MET:CE	2.32	0.56
1:D:363:GLN:OE1	1:D:363:GLN:HA	2.06	0.56
1:C:308:ILE:O	1:C:312:ILE:HG13	2.05	0.56
1:D:309:VAL:CG1	1:D:411:LEU:HD11	2.35	0.56
1:B:434:VAL:HG23	1:B:435:VAL:HG13	1.88	0.56
1:C:379:LYS:O	1:C:444:MET:HG3	2.05	0.56
1:C:476:GLU:HB3	1:C:486:LEU:HD11	1.88	0.56
1:D:435:VAL:HG22	1:D:439:PRO:HB3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:448:VAL:O	1:D:452:GLN:HA	2.06	0.56
1:D:288:MET:HA	1:D:288:MET:CE	2.36	0.56
1:B:412:TRP:CD1	1:B:474:MET:HE1	2.41	0.56
1:D:286:HIS:CE1	1:D:345:LYS:HG2	2.41	0.55
1:A:309:VAL:HG12	1:A:495:LEU:HD13	1.87	0.55
1:D:312:ILE:CD1	1:D:342:ILE:HD13	2.35	0.55
1:A:298:THR:HG22	1:A:299:THR:O	2.06	0.55
1:B:312:ILE:HD13	1:B:342:ILE:HD13	1.88	0.55
1:C:302:THR:O	1:C:306:LEU:HD23	2.06	0.55
1:C:235:LYS:HD3	1:C:237:PHE:CZ	2.41	0.55
1:A:376:VAL:HG22	1:A:376:VAL:O	2.07	0.55
1:C:412:TRP:HB2	1:C:474:MET:HE3	1.89	0.55
1:B:295:LEU:HD21	1:B:414:VAL:HG22	1.89	0.55
1:C:473:LEU:HD21	1:C:494:THR:HG21	1.89	0.55
1:A:208:ILE:CG2	1:A:280:TRP:HZ3	2.18	0.54
1:A:412:TRP:CZ2	1:A:416:ARG:HD2	2.43	0.54
1:B:412:TRP:HD1	1:B:474:MET:HE1	1.72	0.54
1:B:335:ARG:HD3	1:B:357:LEU:O	2.07	0.54
1:A:442:GLU:HG3	1:A:445:ARG:NH2	2.23	0.54
1:C:208:ILE:CG2	1:C:227:TRP:HB2	2.38	0.54
1:A:412:TRP:HB2	1:A:474:MET:HE3	1.89	0.54
1:C:335:ARG:HD3	1:C:357:LEU:O	2.08	0.54
1:D:379:LYS:HE2	1:D:441:PHE:CZ	2.42	0.54
1:A:204:VAL:HG23	1:A:206:ARG:N	2.10	0.54
1:D:263:LEU:HD13	1:D:343:LEU:HD12	1.90	0.54
1:D:309:VAL:HG13	1:D:411:LEU:CD1	2.38	0.54
1:A:305:CYS:HB3	1:A:418:MET:CE	2.37	0.54
1:A:407:PHE:CE2	1:A:491:ILE:HG21	2.34	0.54
1:B:251:LEU:HB3	1:B:257:LEU:HD12	1.88	0.54
1:A:412:TRP:CD1	1:A:474:MET:HE1	2.42	0.53
1:C:278:GLN:O	1:C:278:GLN:HG3	2.08	0.53
1:A:431:PHE:O	1:A:435:VAL:HG22	2.09	0.53
1:B:223:TRP:HB2	1:B:234:VAL:HG22	1.91	0.53
1:A:251:LEU:HD13	1:A:355:LEU:HD23	1.91	0.53
1:B:247:ARG:NH2	1:B:360:MET:SD	2.81	0.53
1:C:412:TRP:HD1	1:C:474:MET:HE1	1.73	0.53
1:D:322:GLU:OE2	1:D:362:SER:HB2	2.08	0.53
1:D:476:GLU:HB3	1:D:486:LEU:HD13	1.91	0.53
1:A:306:LEU:HD11	1:A:466:THR:HG21	1.89	0.53
1:B:439:PRO:CG	1:B:444:MET:HE3	2.39	0.53
1:A:370:VAL:CG1	1:A:374:PRO:HD3	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:306:LEU:HD21	1:C:466:THR:HG23	1.91	0.52
1:D:249:THR:O	1:D:253:ASN:N	2.41	0.52
1:B:208:ILE:CD1	1:B:227:TRP:HB2	2.39	0.52
1:C:298:THR:HG22	1:C:299:THR:O	2.09	0.52
1:D:251:LEU:HD23	1:D:355:LEU:HD13	1.91	0.52
1:D:251:LEU:O	1:D:255:VAL:HB	2.10	0.52
1:C:495:LEU:HD23	1:C:498:ILE:HD12	1.92	0.52
1:B:259:HIS:HE1	1:B:261:ASN:HD22	1.58	0.52
1:C:245:TRP:CZ2	1:C:270:MET:HB3	2.45	0.52
1:D:412:TRP:CZ2	1:D:416:ARG:HD2	2.45	0.52
1:B:235:LYS:HD3	1:B:237:PHE:CZ	2.45	0.51
1:A:306:LEU:HD11	1:A:466:THR:HG23	1.91	0.51
1:C:271:THR:OG1	1:C:278:GLN:HG2	2.10	0.51
1:A:442:GLU:HG3	1:A:445:ARG:HH21	1.76	0.51
1:D:399:TYR:O	1:D:402:VAL:HG13	2.10	0.51
1:B:252:TYR:CD1	1:B:265:PHE:HB2	2.46	0.51
1:B:480:GLN:O	1:B:482:PRO:HD3	2.11	0.51
1:C:462:PHE:HA	1:C:468:THR:OG1	2.10	0.51
1:D:401:ARG:HD3	1:D:482:PRO:O	2.10	0.51
1:B:263:LEU:HD13	1:B:343:LEU:HD12	1.93	0.51
1:A:323:ILE:HG22	1:A:328:GLY:HA2	1.92	0.51
1:C:383:ALA:O	1:C:387:LEU:HD22	2.10	0.51
1:B:407:PHE:CE2	1:B:491:ILE:HG21	2.46	0.50
1:C:237:PHE:CG	1:C:245:TRP:HB2	2.46	0.50
1:D:206:ARG:HD3	1:D:280:TRP:CZ3	2.46	0.50
1:C:208:ILE:HB	1:C:226:SER:O	2.11	0.50
1:C:385:GLU:OE2	1:C:485:ARG:NH1	2.37	0.50
1:D:457:ILE:HG23	1:D:461:TRP:CE3	2.47	0.49
1:B:251:LEU:HD11	1:B:358:ALA:HB3	1.93	0.49
1:D:359:VAL:HG21	1:D:399:TYR:CD2	2.48	0.49
1:B:470:LEU:O	1:B:474:MET:HG3	2.13	0.49
1:A:335:ARG:HD3	1:A:357:LEU:O	2.12	0.49
1:A:379:LYS:HA	1:A:382:MET:HG3	1.94	0.49
1:C:330:PRO:HG3	1:C:360:MET:HE3	1.94	0.49
1:A:288:MET:CE	1:A:346:LYS:HA	2.42	0.49
1:B:448:VAL:O	1:B:452:GLN:HA	2.13	0.49
1:C:407:PHE:HE2	1:C:491:ILE:HG21	1.78	0.49
1:B:261:ASN:O	1:B:351:CYS:HA	2.12	0.48
1:B:255:VAL:HG13	1:B:330:PRO:HD2	1.95	0.48
1:C:286:HIS:CE1	1:C:345:LYS:HG2	2.48	0.48
1:B:223:TRP:HB2	1:B:234:VAL:CG2	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:214:VAL:HG11	2:D:501:MM8:C11	2.44	0.48
1:A:396:PHE:CE2	1:A:400:LYS:HE3	2.49	0.48
1:B:224:ARG:HD3	1:B:285:TYR:CZ	2.49	0.48
1:C:263:LEU:HD13	1:C:343:LEU:HD12	1.95	0.48
1:A:398:SER:O	1:A:402:VAL:HG23	2.12	0.48
1:B:251:LEU:HD13	1:B:355:LEU:HD23	1.94	0.48
1:A:266:ILE:HD11	1:A:284:HIS:CD2	2.48	0.48
1:B:292:TYR:CE2	1:B:340:LYS:HE2	2.49	0.48
1:C:448:VAL:O	1:C:452:GLN:HA	2.13	0.48
1:C:425:GLU:OE1	1:C:460:ARG:NH2	2.40	0.48
1:C:407:PHE:CE2	1:C:491:ILE:HG21	2.49	0.47
1:D:428:LYS:HB3	1:D:432:TYR:CD1	2.49	0.47
1:B:489:LEU:O	1:B:493:LYS:HG3	2.14	0.47
1:D:317:ALA:HA	1:D:488:ALA:HB1	1.96	0.47
1:D:412:TRP:O	1:D:416:ARG:HG3	2.14	0.47
1:D:407:PHE:CE2	1:D:411:LEU:HD13	2.49	0.47
1:B:390:THR:O	1:B:390:THR:HG22	2.15	0.47
1:B:412:TRP:HB2	1:B:474:MET:HE3	1.97	0.47
1:A:375:ARG:HD2	1:A:441:PHE:CZ	2.49	0.47
1:A:462:PHE:CE2	1:C:297:LEU:HD21	2.49	0.47
1:B:209:THR:O	1:B:211:LEU:HD12	2.15	0.47
1:B:365:THR:O	1:B:367:GLN:N	2.48	0.47
1:A:337:LEU:HD23	1:A:338:LYS:N	2.30	0.47
1:C:412:TRP:CD1	1:C:474:MET:HE1	2.49	0.47
1:B:294:TYR:HD2	1:B:295:LEU:HD12	1.79	0.47
1:A:353:ALA:O	1:A:355:LEU:HD12	2.15	0.47
1:A:420:SER:O	1:A:423:ILE:HG22	2.16	0.46
1:A:384:PRO:HD2	1:A:485:ARG:NH2	2.31	0.46
1:C:236:ILE:HG12	1:C:280:TRP:NE1	2.29	0.46
1:D:435:VAL:CG2	1:D:436:PRO:HD2	2.42	0.46
1:A:433:ASP:OD1	1:A:434:VAL:HG13	2.16	0.46
1:B:288:MET:HE3	1:B:346:LYS:HA	1.98	0.46
1:D:289:GLY:HA2	2:D:501:MM8:C13	2.45	0.46
1:A:335:ARG:NH2	1:A:369:ASP:OD2	2.49	0.46
1:B:291:LEU:HD12	1:B:295:LEU:HD13	1.98	0.46
1:A:384:PRO:HG3	1:A:448:VAL:CG1	2.45	0.46
1:B:204:VAL:HG12	1:B:205:ALA:N	2.31	0.46
1:B:388:ASP:O	1:B:389:GLU:HB2	2.16	0.46
1:A:374:PRO:HG2	1:A:375:ARG:H	1.80	0.46
1:A:361:HIS:HB2	1:A:396:PHE:CD1	2.51	0.46
1:D:435:VAL:HG21	1:D:444:MET:SD	2.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:VAL:O	1:B:370:VAL:HG13	2.16	0.45
1:B:435:VAL:HG11	1:B:444:MET:CE	2.47	0.45
1:A:396:PHE:CD2	1:A:400:LYS:HE3	2.51	0.45
1:A:499:ASP:CB	1:D:303:VAL:HG21	2.47	0.45
1:D:224:ARG:HB2	1:D:285:TYR:CE2	2.52	0.45
1:B:204:VAL:HG12	1:B:205:ALA:H	1.81	0.45
1:D:245:TRP:CZ2	1:D:270:MET:HB2	2.52	0.45
1:B:288:MET:HE1	1:B:346:LYS:HA	1.99	0.45
1:C:208:ILE:HG22	1:C:227:TRP:CD1	2.52	0.45
1:D:407:PHE:CZ	1:D:411:LEU:HD13	2.51	0.45
1:A:252:TYR:CD1	1:A:265:PHE:HB2	2.51	0.45
1:B:334:HIS:O	1:B:335:ARG:HB2	2.17	0.44
1:D:401:ARG:HD3	1:D:482:PRO:C	2.38	0.44
1:A:205:ALA:HB3	1:A:269:ASP:OD2	2.18	0.44
1:D:282:ILE:N	1:D:282:ILE:HD12	2.32	0.44
1:D:380:ARG:HD3	1:D:430:PRO:O	2.17	0.44
1:A:335:ARG:HD2	1:A:359:VAL:HG13	1.99	0.44
1:C:291:LEU:O	1:C:295:LEU:HB2	2.18	0.44
1:D:259:HIS:HB3	1:D:262[A]:ILE:HG12	1.99	0.44
1:C:320:HIS:HD2	1:C:332:ILE:O	2.02	0.43
1:C:208:ILE:HG22	1:C:227:TRP:HB2	2.00	0.43
1:B:252:TYR:CG	1:B:265:PHE:HB2	2.53	0.43
1:A:235:LYS:HD3	1:A:237:PHE:CZ	2.53	0.43
1:B:462:PHE:CE2	1:D:228:GLN:HA	2.54	0.43
1:C:246:PHE:O	1:C:250:GLU:HB2	2.18	0.43
1:A:370:VAL:HG22	1:A:374:PRO:N	2.34	0.43
1:A:415:ALA:HB3	1:A:461:TRP:CH2	2.53	0.43
1:B:219:TYR:HD1	1:B:219:TYR:H	1.67	0.43
1:C:248:GLU:HG3	1:C:355:LEU:HB2	1.99	0.43
1:C:224:ARG:HB2	1:C:285:TYR:CE1	2.53	0.43
1:A:300:LEU:HB2	1:A:418:MET:HA	2.00	0.43
1:A:255:VAL:HG12	1:A:255:VAL:O	2.19	0.43
1:B:234:VAL:CG1	1:B:282:ILE:HD12	2.40	0.43
2:D:501:MM8:C24	2:D:501:MM8:C12	2.97	0.43
1:D:309:VAL:HG12	1:D:495:LEU:CD1	2.48	0.42
1:D:334:HIS:O	1:D:335:ARG:HB2	2.19	0.42
1:C:312:ILE:HD13	1:C:342:ILE:HD13	2.01	0.42
1:D:438:ASP:N	1:D:439:PRO:HD3	2.34	0.42
1:C:335:ARG:HD2	1:C:359:VAL:HG22	2.01	0.42
1:A:464:ASP:OD1	1:A:466:THR:N	2.52	0.42
1:B:292:TYR:CD2	1:B:340:LYS:HE2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:LEU:HD21	1:C:414:VAL:HA	2.02	0.42
1:C:419:VAL:HB	1:C:424:VAL:CG1	2.50	0.42
1:A:364:SER:CA	1:B:390:THR:HG23	2.50	0.42
1:B:438:ASP:N	3:B:503:SO4:O3	2.48	0.42
1:C:270:MET:HA	1:C:278:GLN:O	2.20	0.42
1:C:236:ILE:HG12	1:C:280:TRP:CD1	2.55	0.41
1:D:261:ASN:O	1:D:351:CYS:HA	2.20	0.41
1:C:415:ALA:HB3	1:C:461:TRP:CH2	2.55	0.41
1:D:246:PHE:O	1:D:250:GLU:HB2	2.20	0.41
1:A:270:MET:HG3	1:A:279:LEU:HD21	1.98	0.41
1:A:364:SER:N	1:B:390:THR:HG23	2.35	0.41
1:D:335:ARG:HH11	1:D:359:VAL:HG13	1.85	0.41
1:D:474:MET:HE3	1:D:478:TRP:HH2	1.85	0.41
1:B:333:ALA:HB3	1:B:359:VAL:HG22	2.02	0.41
1:D:420:SER:HB3	1:D:460:ARG:NH1	2.34	0.41
1:C:219:TYR:CD2	1:C:235:LYS:HE3	2.56	0.41
1:B:330:PRO:HG3	1:B:360:MET:CE	2.50	0.41
1:D:380:ARG:HA	1:D:444:MET:CE	2.50	0.41
1:C:261:ASN:O	1:C:351:CYS:HA	2.20	0.41
1:D:319:LEU:HD12	1:D:319:LEU:HA	1.90	0.41
1:D:474:MET:HE3	1:D:478:TRP:CH2	2.55	0.41
1:A:388:ASP:HB3	1:A:390:THR:OG1	2.20	0.41
1:B:360:MET:HE2	1:B:360:MET:HB3	1.94	0.41
1:B:384:PRO:HG3	1:B:448:VAL:HG12	2.01	0.41
1:C:270:MET:HG3	1:C:270:MET:O	2.21	0.41
1:D:434:VAL:O	1:D:434:VAL:HG12	2.21	0.41
1:B:461:TRP:HB3	1:B:467:LEU:HB3	2.02	0.41
1:C:330:PRO:HG3	1:C:360:MET:CE	2.51	0.41
1:D:293:ASP:O	1:D:296:GLN:HB2	2.21	0.41
1:D:320:HIS:HB3	1:D:400:LYS:HE3	2.02	0.41
1:C:334:HIS:O	1:C:335:ARG:HB2	2.21	0.41
1:D:436:PRO:O	1:D:439:PRO:HG3	2.21	0.40
1:C:434:VAL:HG21	1:C:447:VAL:HG11	2.03	0.40
1:D:384:PRO:HG3	1:D:448:VAL:CG1	2.51	0.40
1:B:251:LEU:HD21	1:B:360:MET:SD	2.61	0.40
1:B:489:LEU:CG	1:B:493:LYS:HE3	2.45	0.40
1:A:252:TYR:HB3	1:A:265:PHE:CG	2.56	0.40
1:B:357:LEU:HD11	1:B:375:ARG:HD3	2.03	0.40
1:B:380:ARG:NE	3:B:503:SO4:O4	2.48	0.40
1:A:205:ALA:O	1:A:208:ILE:HG22	2.22	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	287/301 (95%)	279 (97%)	8 (3%)	0	100	100
1	B	296/301 (98%)	282 (95%)	14 (5%)	0	100	100
1	C	272/301 (90%)	261 (96%)	11 (4%)	0	100	100
1	D	283/301 (94%)	273 (96%)	10 (4%)	0	100	100
All	All	1138/1204 (94%)	1095 (96%)	43 (4%)	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	245/270 (91%)	244 (100%)	1 (0%)	92	96
1	B	237/270 (88%)	235 (99%)	2 (1%)	83	90
1	C	221/270 (82%)	221 (100%)	0	100	100
1	D	238/270 (88%)	236 (99%)	2 (1%)	83	90
All	All	941/1080 (87%)	936 (100%)	5 (0%)	90	95

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	241	ASP
1	B	257	LEU

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Mol	Chain	Res	Type
1	B	433	ASP
1	D	206	ARG
1	D	402	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

15 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	MM8	A	501	-	33,33,33	2.16	9 (27%)	43,46,46	1.53	3 (6%)
3	SO4	A	502	-	4,4,4	0.15	0	6,6,6	0.11	0
3	SO4	A	503	-	4,4,4	0.15	0	6,6,6	0.07	0
3	SO4	A	504	-	4,4,4	0.17	0	6,6,6	0.15	0
2	MM8	B	501	-	33,33,33	2.16	9 (27%)	43,46,46	1.45	7 (16%)
3	SO4	B	502	-	4,4,4	0.15	0	6,6,6	0.09	0
3	SO4	B	503	-	4,4,4	0.16	0	6,6,6	0.05	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	B	504	-	4,4,4	0.16	0	6,6,6	0.17	0
2	MM8	C	501	-	33,33,33	2.15	8 (24%)	43,46,46	1.36	7 (16%)
3	SO4	C	502	-	4,4,4	0.14	0	6,6,6	0.08	0
3	SO4	C	503	-	4,4,4	0.15	0	6,6,6	0.06	0
3	SO4	C	504	-	4,4,4	0.16	0	6,6,6	0.06	0
2	MM8	D	501	-	33,33,33	2.16	8 (24%)	43,46,46	1.45	2 (4%)
3	SO4	D	502	-	4,4,4	0.16	0	6,6,6	0.06	0
3	SO4	D	503	-	4,4,4	0.16	0	6,6,6	0.13	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MM8	A	501	-	-	3/18/26/26	0/4/4/4
2	MM8	B	501	-	-	5/18/26/26	0/4/4/4
2	MM8	C	501	-	-	0/18/26/26	0/4/4/4
2	MM8	D	501	-	-	2/18/26/26	0/4/4/4

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	MM8	C28-N29	6.36	1.45	1.33
2	B	501	MM8	C28-N29	6.33	1.44	1.33
2	A	501	MM8	C28-N29	6.32	1.44	1.33
2	C	501	MM8	C28-N29	6.27	1.44	1.33
2	D	501	MM8	O02-C03	5.09	1.45	1.37
2	A	501	MM8	O02-C03	5.09	1.45	1.37
2	C	501	MM8	O02-C03	5.05	1.45	1.37
2	B	501	MM8	O02-C03	5.01	1.45	1.37
2	D	501	MM8	C14-N15	3.97	1.49	1.38
2	C	501	MM8	C14-N15	3.92	1.49	1.38
2	B	501	MM8	C14-N15	3.91	1.49	1.38
2	C	501	MM8	C06-C05	-3.83	1.42	1.49
2	A	501	MM8	C14-N15	3.79	1.49	1.38
2	B	501	MM8	C06-C05	-3.64	1.43	1.49
2	D	501	MM8	C06-C05	-3.60	1.43	1.49
2	A	501	MM8	C06-C05	-3.57	1.43	1.49
2	B	501	MM8	C10-C11	-3.46	1.43	1.49
2	A	501	MM8	C10-C11	-3.40	1.43	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	MM8	C10-C11	-3.37	1.43	1.49
2	D	501	MM8	C10-C11	-3.15	1.43	1.49
2	C	501	MM8	O30-C28	-3.14	1.18	1.24
2	A	501	MM8	O30-C28	-3.13	1.18	1.24
2	D	501	MM8	O30-C28	-3.12	1.18	1.24
2	B	501	MM8	O30-C28	-3.11	1.18	1.24
2	B	501	MM8	C27-C28	2.76	1.53	1.50
2	C	501	MM8	C27-C28	2.69	1.53	1.50
2	A	501	MM8	C27-C28	2.61	1.53	1.50
2	D	501	MM8	C27-C28	2.45	1.53	1.50
2	D	501	MM8	C16-N15	2.44	1.50	1.46
2	B	501	MM8	C16-N15	2.39	1.50	1.46
2	A	501	MM8	C16-N15	2.39	1.50	1.46
2	C	501	MM8	C16-N15	2.26	1.50	1.46
2	A	501	MM8	C10-C23	2.06	1.43	1.40
2	B	501	MM8	C20-N15	2.01	1.49	1.46

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	MM8	C03-C27-C28	-7.29	119.97	125.11
2	D	501	MM8	C03-C27-C28	-6.70	120.38	125.11
2	B	501	MM8	C03-C27-C28	-5.61	121.15	125.11
2	C	501	MM8	C03-C27-C28	-4.46	121.96	125.11
2	B	501	MM8	C01-O02-C03	-2.55	113.77	117.53
2	B	501	MM8	O02-C03-C04	-2.52	119.83	124.14
2	C	501	MM8	O02-C03-C04	-2.52	119.84	124.14
2	A	501	MM8	C09-N08-C07	2.41	120.81	117.48
2	C	501	MM8	O02-C03-C27	2.37	120.05	116.53
2	C	501	MM8	C09-N08-C07	2.36	120.74	117.48
2	D	501	MM8	C09-N08-C07	2.29	120.65	117.48
2	B	501	MM8	C09-N08-C07	2.27	120.61	117.48
2	B	501	MM8	O02-C03-C27	2.27	119.90	116.53
2	C	501	MM8	C27-C28-N29	2.23	121.68	118.25
2	C	501	MM8	O30-C28-N29	-2.22	119.44	122.60
2	B	501	MM8	O30-C28-N29	-2.18	119.50	122.60
2	A	501	MM8	C10-C09-N08	-2.17	121.26	124.51
2	B	501	MM8	C10-C09-N08	-2.11	121.35	124.51
2	C	501	MM8	C20-N15-C16	-2.09	107.05	111.53

There are no chirality outliers.

All (10) torsion outliers are listed below:

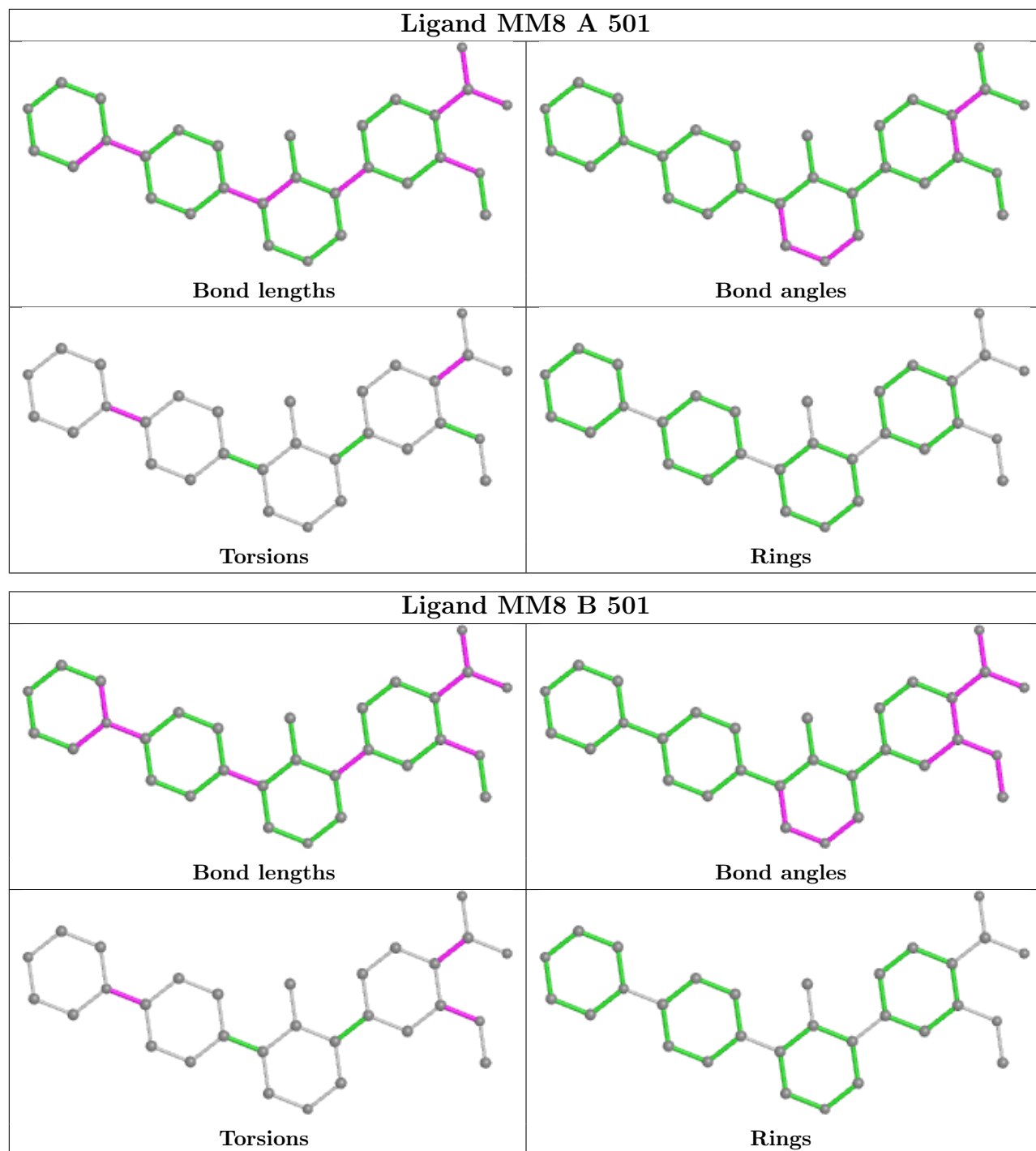
Mol	Chain	Res	Type	Atoms
2	B	501	MM8	C21-C14-N15-C16
2	B	501	MM8	C13-C14-N15-C16
2	B	501	MM8	C04-C03-O02-C01
2	B	501	MM8	C27-C03-O02-C01
2	D	501	MM8	C03-C27-C28-N29
2	D	501	MM8	C03-C27-C28-O30
2	A	501	MM8	C21-C14-N15-C16
2	A	501	MM8	C13-C14-N15-C16
2	A	501	MM8	C26-C27-C28-N29
2	B	501	MM8	C03-C27-C28-N29

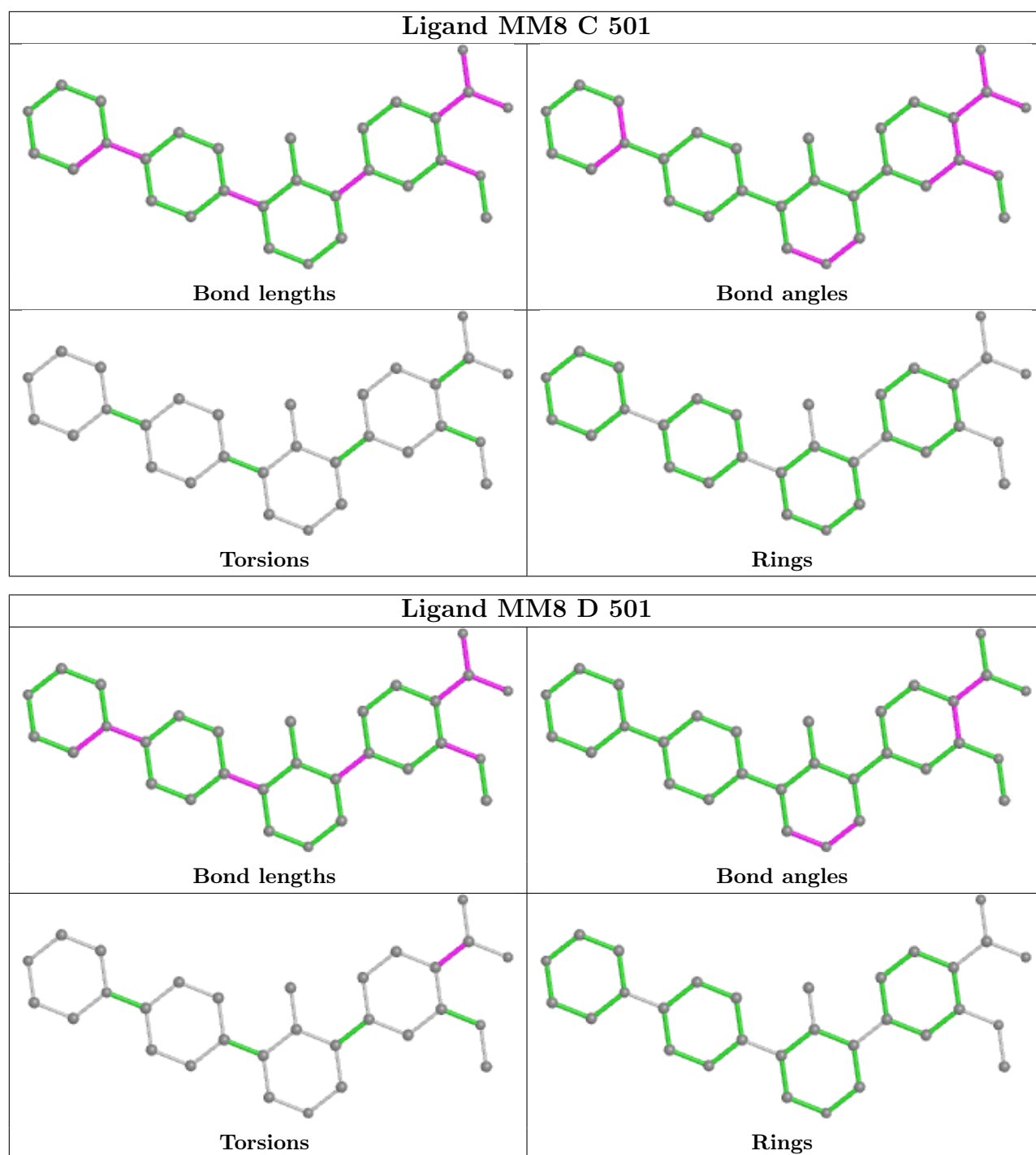
There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	MM8	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	292/301 (97%)	0.32	3 (1%) 82 87	49, 66, 96, 111	0
1	B	297/301 (98%)	0.42	8 (2%) 54 64	53, 69, 98, 121	0
1	C	278/301 (92%)	0.49	14 (5%) 29 36	56, 75, 99, 112	0
1	D	285/301 (94%)	0.62	13 (4%) 32 42	57, 76, 99, 116	0
All	All	1152/1204 (95%)	0.46	38 (3%) 46 56	49, 72, 99, 121	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	499	ASP	4.8
1	A	499	ASP	3.6
1	D	498	ILE	3.6
1	C	219	TYR	3.5
1	C	488	ALA	3.4
1	D	478	TRP	3.3
1	C	383	ALA	3.3
1	C	450	VAL	3.1
1	C	447	VAL	3.1
1	D	496	THR	3.0
1	B	353	ALA	2.9
1	D	419	VAL	2.8
1	C	388	ASP	2.7
1	D	434	VAL	2.7
1	C	448	VAL	2.7
1	D	264	GLY	2.6
1	C	498	ILE	2.6
1	B	368	LEU	2.6
1	D	219	TYR	2.6
1	C	387	LEU	2.5
1	D	324	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	236	ILE	2.5
1	C	405	TRP	2.4
1	D	251	LEU	2.4
1	D	460	ARG	2.4
1	B	253	ASN	2.3
1	C	353	ALA	2.3
1	A	326	THR	2.3
1	D	431	PHE	2.3
1	B	375	ARG	2.3
1	C	395	CYS	2.3
1	C	444	MET	2.2
1	A	306	LEU	2.2
1	C	220	GLY	2.2
1	B	219	TYR	2.1
1	B	272	SER	2.1
1	D	253	ASN	2.1
1	D	303	VAL	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.

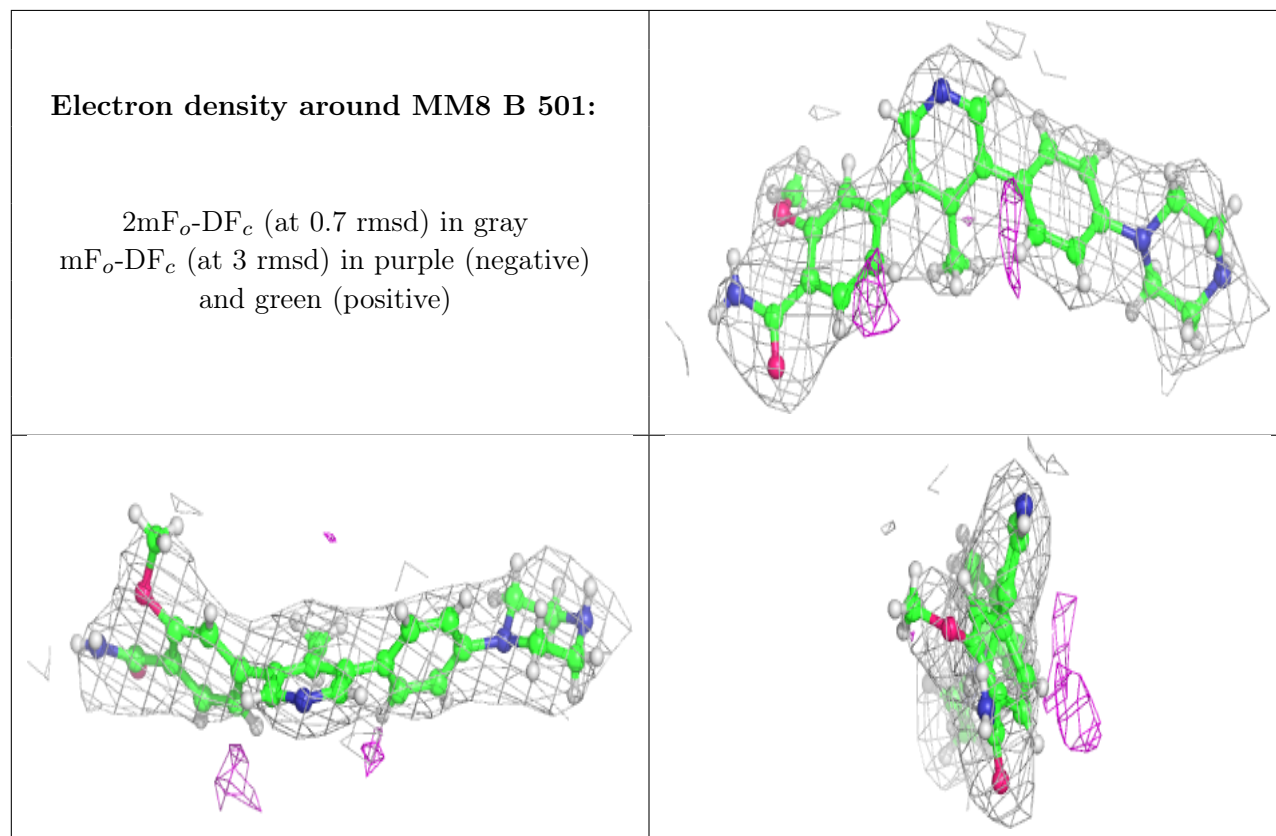
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	D	503	5/5	0.87	0.42	78,87,94,109	0
3	SO4	C	504	5/5	0.88	0.28	78,85,105,107	0
3	SO4	A	504	5/5	0.89	0.47	81,87,99,115	0
2	MM8	B	501	30/30	0.90	0.21	64,80,119,124	0
3	SO4	C	503	5/5	0.92	0.26	103,113,122,126	0
3	SO4	B	503	5/5	0.92	0.12	105,106,113,120	0

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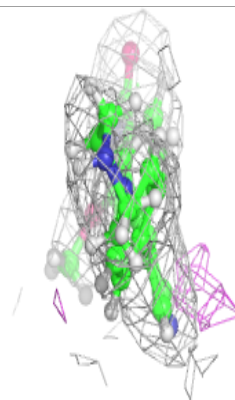
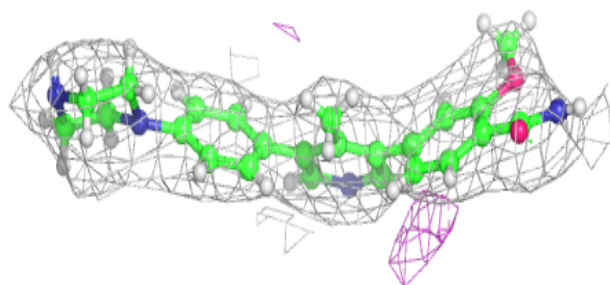
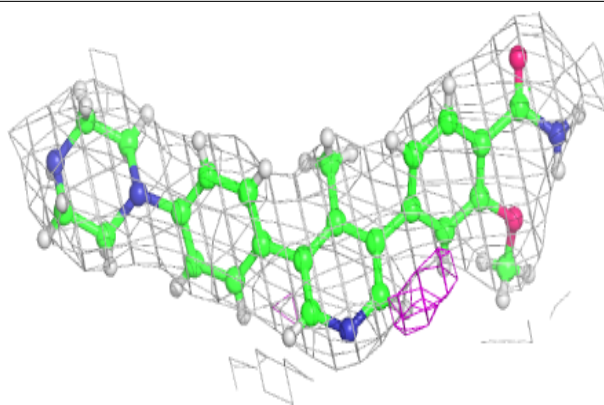
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	B	502	5/5	0.93	0.24	73,78,93,98	0
3	SO4	C	502	5/5	0.93	0.13	71,73,82,87	0
2	MM8	A	501	30/30	0.94	0.20	48,68,103,105	0
3	SO4	B	504	5/5	0.94	0.13	70,73,82,85	0
2	MM8	D	501	30/30	0.94	0.21	55,73,98,103	0
2	MM8	C	501	30/30	0.95	0.25	49,68,91,93	0
3	SO4	D	502	5/5	0.95	0.25	67,78,84,97	0
3	SO4	A	503	5/5	0.95	0.11	75,82,94,99	0
3	SO4	A	502	5/5	0.97	0.13	73,73,80,87	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.

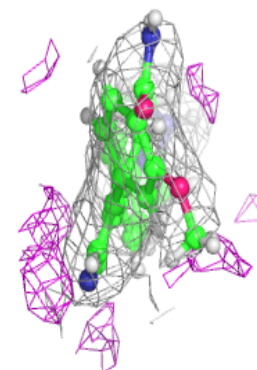
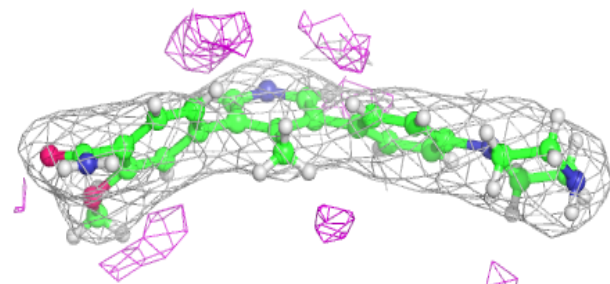
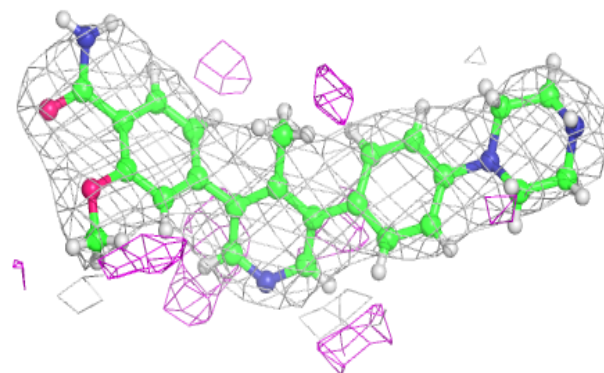


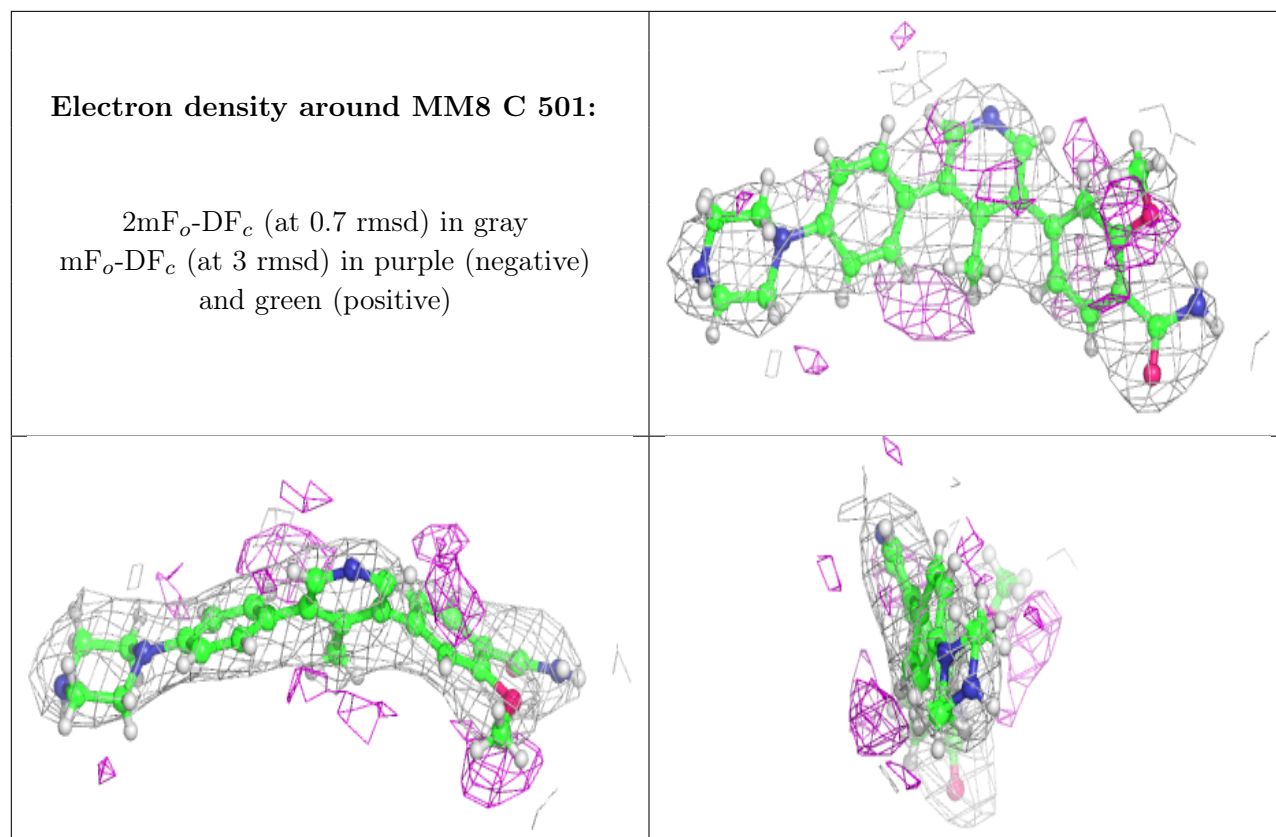
Electron density around MM8 A 501:

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

**Electron density around MM8 D 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 [i](#)

There are no such residues in this entry.