



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 04:37 pm BST

PDB ID : 3MEC
Title : HIV-1 Reverse Transcriptase in Complex with TMC125
Authors : Lansdon, E.B.
Deposited on : 2010-03-31
Resolution : 2.30 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

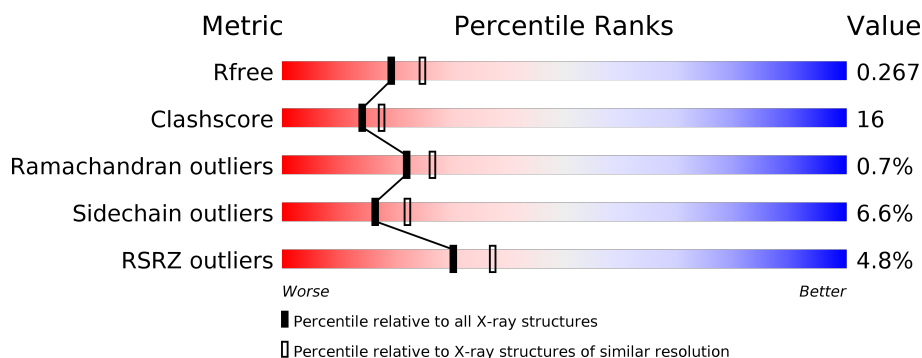
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	560	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>27%</div> <div>5%</div> </div> </div>
2	B	440	<div> <div>6%</div> <div> <div></div> <div>63%</div> <div>26%</div> <div>9%</div> </div> </div>

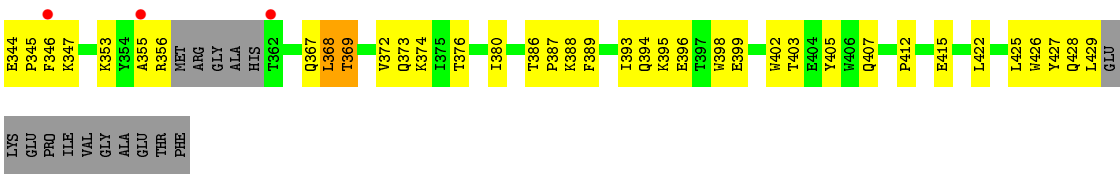
- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	125	Total	O	0	0
			125	125		
5	B	90	Total	O	0	0
			90	90		



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.95Å 154.12Å 154.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.91 – 2.30 47.16 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.6 (29.91-2.30) 90.7 (47.16-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.32Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.226 , 0.270 0.224 , 0.267	Depositor DCC
R_{free} test set	2878 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.0	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	8072	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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: 65B, 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	0.33	0/4598	0.50	0/6246
2	B	0.34	0/3413	0.50	0/4637
All	All	0.33	0/8011	0.50	0/10883

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
2	B	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
2	B	91	GLN	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	4483	0	4537	150	0
2	B	3321	0	3352	103	0
3	A	28	0	15	2	0
4	A	20	0	0	0	0
4	B	5	0	0	1	0
5	A	125	0	0	7	0
5	B	90	0	0	2	0
All	All	8072	0	7904	246	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 16.

All (246) 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:90:VAL:HG23	1:A:91:GLN:H	1.30	0.97
2:B:260:LEU:HD13	2:B:279:LEU:HD21	1.53	0.87
1:A:52:PRO:HA	1:A:53:GLU:C	1.95	0.87
1:A:92:LEU:H	1:A:92:LEU:HD23	1.44	0.82
1:A:451:LYS:HB3	1:A:471:ASP:HA	1.62	0.82
1:A:544:GLY:HA2	2:B:286:THR:HG22	1.61	0.80
1:A:275:LYS:H	1:A:306:ASN:HD21	1.25	0.80
1:A:194:GLU:H	1:A:194:GLU:CD	1.89	0.76
1:A:520:GLN:O	1:A:523:GLU:HG3	1.87	0.75
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.68	0.75
2:B:296:THR:HG22	2:B:298:GLU:H	1.52	0.74
1:A:503:LEU:HD13	1:A:533:LEU:HD12	1.69	0.74
2:B:297:GLU:HG3	2:B:298:GLU:H	1.53	0.74
2:B:259:LYS:HE2	2:B:425:LEU:HD21	1.70	0.73
2:B:337:TRP:HE1	2:B:367:GLN:HE21	1.34	0.72
1:A:199:ARG:HG2	1:A:199:ARG:HH11	1.52	0.72
1:A:60:VAL:HG12	1:A:75:VAL:HG22	1.69	0.72
1:A:298:GLU:H	1:A:298:GLU:CD	1.93	0.72
1:A:64:LYS:NZ	1:A:69:THR:HG22	2.05	0.72
2:B:369:THR:HG21	2:B:405:TYR:HB2	1.73	0.71
1:A:96:HIS:HD1	1:A:98:ALA:H	1.39	0.70
2:B:7:THR:HG22	2:B:119:PRO:HB2	1.73	0.70
2:B:69:THR:O	2:B:70:LYS:HG3	1.92	0.70
2:B:257:ILE:HB	2:B:283:LEU:HD21	1.74	0.70
1:A:450:THR:HG22	1:A:452:LEU:HB2	1.73	0.69
2:B:85:GLN:HA	2:B:88:TRP:NE1	2.08	0.69
1:A:393:ILE:HB	1:A:423:VAL:CG1	2.22	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:81:ASN:ND2	2:B:154:LYS:H	1.91	0.68
2:B:295:LEU:H	2:B:295:LEU:HD22	1.58	0.68
2:B:376:THR:HG23	2:B:386:THR:HG23	1.75	0.68
1:A:503:LEU:HD23	2:B:422:LEU:HD21	1.76	0.68
1:A:64:LYS:HZ3	1:A:69:THR:HA	1.60	0.67
1:A:191:SER:HB2	1:A:193:LEU:HD23	1.78	0.66
1:A:479:LEU:HB3	1:A:517:LEU:HD13	1.79	0.65
2:B:312:GLU:HB3	2:B:313:PRO:HD2	1.79	0.64
2:B:373:GLN:HE22	2:B:407:GLN:H	1.44	0.64
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.80	0.64
1:A:406:TRP:HE3	1:A:407:GLN:HE21	1.46	0.63
1:A:27:THR:HG22	1:A:29:GLU:H	1.62	0.63
1:A:486:LEU:HB3	1:A:524:GLN:HG3	1.81	0.63
2:B:252:TRP:CD1	2:B:295:LEU:HD11	2.33	0.63
1:A:90:VAL:HG23	1:A:91:GLN:N	2.09	0.63
1:A:175:ASN:HB3	1:A:178:ILE:CG2	2.29	0.62
2:B:373:GLN:NE2	2:B:407:GLN:H	1.97	0.62
1:A:169:GLU:HB2	1:A:170:PRO:HD3	1.82	0.62
2:B:297:GLU:HG3	2:B:298:GLU:N	2.15	0.61
1:A:64:LYS:HZ1	1:A:69:THR:HG22	1.63	0.61
1:A:334:GLN:HB3	5:A:611:HOH:O	1.99	0.61
1:A:509:GLN:N	1:A:510:PRO:HD3	2.16	0.61
1:A:175:ASN:HD22	1:A:201:LYS:NZ	1.99	0.60
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.02	0.60
2:B:293:ILE:HG13	2:B:294:PRO:HD2	1.82	0.60
1:A:197:GLN:O	1:A:201:LYS:HG2	2.02	0.60
1:A:175:ASN:ND2	1:A:201:LYS:NZ	2.49	0.60
2:B:13:LYS:HB2	2:B:16:MET:SD	2.41	0.60
2:B:295:LEU:HD22	2:B:295:LEU:N	2.17	0.59
1:A:120:LEU:HD22	1:A:121:ASP:H	1.66	0.59
1:A:242:GLN:O	1:A:244:ILE:HD12	2.02	0.59
2:B:32:LYS:HB2	2:B:32:LYS:NZ	2.17	0.59
1:A:199:ARG:HG2	1:A:199:ARG:NH1	2.12	0.59
1:A:238:LYS:HD2	1:A:315:HIS:CG	2.38	0.59
1:A:287:LYS:HE3	1:A:291:GLU:OE2	2.02	0.59
1:A:54:ASN:ND2	1:A:56:TYR:H	2.01	0.59
2:B:81:ASN:HD21	2:B:154:LYS:H	1.50	0.58
2:B:275:LYS:H	2:B:306:ASN:HD21	1.51	0.58
2:B:393:ILE:HG12	2:B:394:GLN:N	2.19	0.58
1:A:143:ARG:HH11	1:A:143:ARG:HG2	1.68	0.58
2:B:296:THR:HG22	2:B:298:GLU:N	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:441:TYR:CD2	1:A:544:GLY:HA3	2.39	0.57
2:B:80:LEU:HD22	2:B:84:THR:HG23	1.87	0.57
2:B:393:ILE:HG12	2:B:394:GLN:H	1.69	0.57
2:B:345:PRO:O	2:B:346:PHE:HB2	2.04	0.56
1:A:362:THR:HG22	1:A:366:LYS:HD2	1.87	0.56
2:B:81:ASN:HD21	2:B:153:TRP:HA	1.70	0.56
2:B:12:LEU:HD23	2:B:84:THR:HG22	1.87	0.56
1:A:195:ILE:O	1:A:199:ARG:HG3	2.06	0.56
2:B:91:GLN:C	2:B:91:GLN:HE21	2.09	0.56
1:A:328:GLU:HG2	1:A:390:LYS:HB2	1.88	0.56
2:B:184:MET:HE2	5:B:498:HOH:O	2.06	0.55
1:A:19:PRO:HG3	1:A:80:LEU:HA	1.88	0.55
1:A:312:GLU:CD	1:A:312:GLU:H	2.09	0.55
1:A:483:TYR:O	1:A:487:GLN:HG2	2.06	0.55
2:B:305:GLU:O	2:B:309:ILE:HG13	2.06	0.55
2:B:206:ARG:CZ	2:B:206:ARG:HB3	2.37	0.55
1:A:278:GLN:HG3	1:A:298:GLU:HB2	1.89	0.55
1:A:129:ALA:HA	1:A:144:TYR:O	2.06	0.55
1:A:393:ILE:HB	1:A:423:VAL:HG13	1.88	0.55
2:B:259:LYS:NZ	2:B:425:LEU:HD11	2.23	0.54
1:A:486:LEU:HB3	1:A:524:GLN:CG	2.38	0.54
1:A:120:LEU:HD22	1:A:121:ASP:N	2.21	0.54
1:A:368:LEU:HD22	1:A:423:VAL:HG11	1.90	0.54
2:B:282:LEU:HB3	2:B:293:ILE:HD13	1.90	0.54
2:B:344:GLU:HB2	2:B:347:LYS:HD3	1.88	0.54
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.88	0.54
1:A:92:LEU:H	1:A:92:LEU:CD2	2.17	0.54
1:A:350:LYS:HE3	1:A:378:GLU:OE2	2.08	0.53
2:B:183:TYR:CD2	2:B:380:ILE:HD13	2.43	0.53
1:A:503:LEU:CD1	1:A:533:LEU:HD12	2.35	0.53
1:A:536:VAL:HG13	1:A:542:ILE:HG13	1.89	0.53
1:A:24:TRP:CD2	1:A:25:PRO:HD2	2.44	0.53
1:A:255:ASN:HB2	1:A:289:LEU:HG	1.91	0.53
1:A:393:ILE:CB	1:A:423:VAL:HG13	2.40	0.52
1:A:27:THR:CG2	1:A:29:GLU:HG2	2.39	0.52
1:A:27:THR:HG21	5:A:610:HOH:O	2.09	0.52
2:B:253:THR:O	2:B:257:ILE:HG12	2.09	0.52
1:A:205:LEU:HD22	1:A:209:LEU:HD22	1.92	0.52
1:A:107:THR:HG22	1:A:198:HIS:NE2	2.24	0.52
1:A:27:THR:O	1:A:31:ILE:HG13	2.10	0.52
2:B:395:LYS:O	2:B:399:GLU:HG3	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:425:LEU:HD12	2:B:428:GLN:OE1	2.09	0.52
1:A:143:ARG:NH1	1:A:143:ARG:HG2	2.26	0.51
1:A:200:THR:O	1:A:204:GLU:HG3	2.10	0.51
1:A:278:GLN:HB3	1:A:299:ALA:HA	1.92	0.51
1:A:369:THR:O	1:A:373:GLN:HG2	2.09	0.51
2:B:65:LYS:HB2	2:B:68:SER:HB2	1.92	0.51
2:B:259:LYS:HZ3	2:B:425:LEU:HD11	1.75	0.51
1:A:27:THR:HG21	1:A:29:GLU:HG2	1.92	0.51
1:A:107:THR:HG21	1:A:202:ILE:HG13	1.91	0.51
1:A:8:VAL:O	1:A:10:VAL:HG23	2.10	0.51
1:A:517:LEU:HA	1:A:520:GLN:HE21	1.76	0.51
1:A:542:ILE:HD11	2:B:261:VAL:HG11	1.92	0.50
1:A:500:GLN:NE2	2:B:422:LEU:HD23	2.27	0.50
1:A:393:ILE:HG13	1:A:423:VAL:HG13	1.92	0.50
1:A:450:THR:O	1:A:451:LYS:HB2	2.12	0.50
1:A:90:VAL:CG2	1:A:91:GLN:H	2.13	0.50
2:B:306:ASN:O	2:B:310:LEU:HD22	2.11	0.50
2:B:175:ASN:HD21	2:B:201:LYS:NZ	2.10	0.50
1:A:420:PRO:HA	1:A:421:PRO:C	2.32	0.49
1:A:175:ASN:HD22	1:A:201:LYS:HZ1	1.59	0.49
1:A:529:GLU:O	1:A:530:LYS:HG3	2.11	0.49
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.47	0.49
2:B:30:LYS:HE3	2:B:403:THR:HG21	1.93	0.49
1:A:207:GLN:NE2	1:A:207:GLN:HA	2.28	0.49
1:A:547:GLN:O	1:A:551:LEU:HD13	2.13	0.49
2:B:394:GLN:NE2	2:B:396:GLU:HB2	2.27	0.49
1:A:126:LYS:HA	1:A:145:GLN:NE2	2.28	0.49
2:B:214:LEU:HD13	2:B:214:LEU:O	2.12	0.49
1:A:178:ILE:HG12	1:A:179:VAL:N	2.27	0.48
1:A:451:LYS:HD2	1:A:451:LYS:N	2.28	0.48
1:A:52:PRO:HA	1:A:54:ASN:N	2.27	0.48
2:B:165:THR:HG22	5:B:459:HOH:O	2.12	0.48
2:B:167:ILE:HG12	2:B:212:TRP:CD2	2.48	0.48
2:B:296:THR:O	2:B:300:GLU:HG2	2.13	0.48
2:B:24:TRP:CZ3	2:B:399:GLU:HB3	2.48	0.48
2:B:355:ALA:O	2:B:356:ARG:HB3	2.13	0.48
1:A:178:ILE:C	1:A:178:ILE:HD13	2.34	0.48
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.49	0.48
2:B:24:TRP:N	2:B:24:TRP:CD1	2.81	0.48
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.49	0.48
1:A:52:PRO:CA	1:A:53:GLU:C	2.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:SER:HB2	1:A:193:LEU:CD2	2.42	0.48
2:B:376:THR:O	2:B:380:ILE:HG13	2.14	0.47
1:A:450:THR:CG2	1:A:452:LEU:HB2	2.41	0.47
2:B:394:GLN:HE21	2:B:396:GLU:HB2	1.78	0.47
1:A:230:MET:HE3	5:A:620:HOH:O	2.14	0.47
1:A:517:LEU:HA	1:A:520:GLN:NE2	2.29	0.47
2:B:237:ASP:C	2:B:239:TRP:H	2.18	0.47
1:A:126:LYS:HD2	5:A:652:HOH:O	2.15	0.47
1:A:393:ILE:CG1	1:A:423:VAL:HG13	2.45	0.47
2:B:169:GLU:HB3	2:B:170:PRO:HD3	1.97	0.47
2:B:154:LYS:HG2	2:B:184:MET:CE	2.45	0.47
1:A:301:LEU:HD12	1:A:301:LEU:HA	1.74	0.47
2:B:169:GLU:OE1	2:B:173:LYS:HE3	2.15	0.47
2:B:58:THR:HG23	2:B:76:ASP:O	2.15	0.47
1:A:191:SER:CB	1:A:193:LEU:HD23	2.45	0.46
1:A:92:LEU:CD1	2:B:22:LYS:HE2	2.44	0.46
2:B:388:LYS:HE2	2:B:415:GLU:OE1	2.14	0.46
1:A:53:GLU:O	1:A:54:ASN:C	2.54	0.46
2:B:30:LYS:HE3	2:B:403:THR:CG2	2.44	0.46
1:A:457:TYR:HA	1:A:548:VAL:HG21	1.96	0.46
1:A:226:PRO:HB3	1:A:235:HIS:CE1	2.51	0.46
1:A:349:LEU:HD12	1:A:349:LEU:N	2.30	0.46
1:A:130:PHE:CZ	1:A:144:TYR:HB2	2.50	0.46
1:A:392:PRO:O	1:A:423:VAL:HG12	2.15	0.46
1:A:27:THR:HG22	1:A:29:GLU:N	2.29	0.46
2:B:13:LYS:HE3	2:B:16:MET:CE	2.46	0.46
1:A:92:LEU:HD12	2:B:22:LYS:HE2	1.98	0.46
2:B:398:TRP:O	2:B:402:TRP:HD1	1.99	0.46
1:A:220:LYS:HD2	1:A:220:LYS:HA	1.62	0.46
1:A:538:ALA:HA	1:A:545:ASN:HD21	1.81	0.46
2:B:80:LEU:HD13	2:B:153:TRP:CD1	2.51	0.45
2:B:320:ASP:OD2	2:B:323:LYS:HD3	2.16	0.45
2:B:239:TRP:NE1	4:B:441:SO4:O4	2.47	0.45
1:A:298:GLU:CD	1:A:298:GLU:N	2.68	0.45
1:A:406:TRP:CE3	1:A:407:GLN:NE2	2.85	0.45
1:A:345:PRO:HA	1:A:346:PHE:HA	1.53	0.45
1:A:513:SER:H	1:A:519:ASN:HD21	1.64	0.45
2:B:282:LEU:HB3	2:B:293:ILE:CD1	2.45	0.45
2:B:175:ASN:HD21	2:B:201:LYS:HZ2	1.65	0.45
1:A:494:ASN:HB3	2:B:289:LEU:HD12	1.99	0.45
1:A:524:GLN:NE2	5:A:589:HOH:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:ASN:HD22	1:A:54:ASN:C	2.20	0.45
2:B:254:VAL:O	2:B:258:GLN:HG3	2.16	0.45
2:B:422:LEU:HD22	2:B:422:LEU:N	2.32	0.45
1:A:358:ARG:NH1	2:B:394:GLN:HG3	2.33	0.44
2:B:214:LEU:HD13	2:B:214:LEU:C	2.37	0.44
2:B:278:GLN:NE2	2:B:298:GLU:HB3	2.32	0.44
1:A:175:ASN:ND2	1:A:201:LYS:HZ1	2.14	0.44
1:A:169:GLU:CB	1:A:170:PRO:HD3	2.46	0.44
1:A:332:GLN:NE2	1:A:338:THR:HG21	2.33	0.44
2:B:387:PRO:HG2	2:B:389:PHE:CE1	2.52	0.44
1:A:54:ASN:HD22	1:A:56:TYR:H	1.66	0.44
2:B:64:LYS:O	2:B:65:LYS:C	2.56	0.44
1:A:536:VAL:CG1	1:A:542:ILE:HG13	2.46	0.44
1:A:100:LEU:O	1:A:318:TYR:HB3	2.18	0.44
1:A:484:LEU:HD23	1:A:484:LEU:HA	1.81	0.44
1:A:60:VAL:HG11	1:A:130:PHE:CD2	2.53	0.44
1:A:411:ILE:HG13	1:A:414:TRP:CD1	2.53	0.44
3:A:561:65B:N12	3:A:561:65B:H26	2.33	0.43
2:B:368:LEU:O	2:B:372:VAL:HG23	2.19	0.43
2:B:32:LYS:HZ3	2:B:32:LYS:HB2	1.83	0.43
1:A:538:ALA:HA	1:A:545:ASN:ND2	2.33	0.43
1:A:126:LYS:HG2	1:A:127:TYR:N	2.34	0.42
1:A:175:ASN:ND2	1:A:201:LYS:HZ3	2.18	0.42
1:A:178:ILE:HD11	1:A:180:ILE:HG13	2.00	0.42
1:A:279:LEU:O	1:A:282:LEU:HB2	2.18	0.42
2:B:386:THR:CG2	2:B:412:PRO:HB3	2.49	0.42
2:B:303:LEU:HD23	2:B:303:LEU:N	2.35	0.42
1:A:27:THR:CG2	1:A:28:GLU:N	2.82	0.42
1:A:442:VAL:HB	1:A:481:ALA:HB1	2.01	0.42
1:A:211:ARG:NH2	5:A:572:HOH:O	2.52	0.42
2:B:214:LEU:HD13	2:B:215:THR:O	2.20	0.42
1:A:96:HIS:CE1	1:A:350:LYS:HD2	2.55	0.42
1:A:242:GLN:HB2	1:A:243:PRO:HD2	2.02	0.42
1:A:77:PHE:O	1:A:78:ARG:C	2.58	0.42
2:B:405:TYR:O	2:B:407:GLN:HG3	2.20	0.42
1:A:413:GLU:HA	5:A:605:HOH:O	2.20	0.41
1:A:188:TYR:HB3	3:A:561:65B:C3	2.49	0.41
1:A:508:ALA:O	1:A:509:GLN:HB3	2.20	0.41
1:A:58:THR:CG2	1:A:59:PRO:HD2	2.49	0.41
2:B:422:LEU:HB3	2:B:426:TRP:CH2	2.55	0.41
2:B:266:TRP:CH2	2:B:427:TYR:CZ	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:105:SER:O	2:B:190:GLY:HA2	2.20	0.41
2:B:237:ASP:C	2:B:239:TRP:N	2.74	0.41
2:B:30:LYS:HD2	2:B:62:ALA:HB3	2.02	0.41
2:B:373:GLN:HE22	2:B:407:GLN:N	2.15	0.41
1:A:3:SER:OG	1:A:5:ILE:HG22	2.20	0.41
2:B:244:ILE:HG12	2:B:266:TRP:HZ3	1.85	0.41
1:A:548:VAL:O	1:A:552:VAL:HG22	2.21	0.41
2:B:24:TRP:H	2:B:24:TRP:HD1	1.69	0.41
1:A:473:THR:O	1:A:477:THR:HG23	2.21	0.41
1:A:63:ILE:HD11	1:A:74:LEU:HD12	2.03	0.41
2:B:69:THR:C	2:B:70:LYS:HG3	2.40	0.41
1:A:509:GLN:N	1:A:510:PRO:CD	2.83	0.41
1:A:139:THR:HB	1:A:140:PRO:HD2	2.02	0.40
1:A:388:LYS:HE2	1:A:388:LYS:HB3	1.83	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	546/560 (98%)	511 (94%)	29 (5%)	6 (1%)	14	15
2	B	393/440 (89%)	378 (96%)	14 (4%)	1 (0%)	41	50
All	All	939/1000 (94%)	889 (95%)	43 (5%)	7 (1%)	22	26

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	GLU
1	A	90	VAL
1	A	218	ASP
2	B	250	ASP

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Mol	Chain	Res	Type
1	A	543	GLY
1	A	78	ARG
1	A	509	GLN

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	A	492/500 (98%)	457 (93%)	35 (7%)	14	19
2	B	366/400 (92%)	344 (94%)	22 (6%)	19	26
All	All	858/900 (95%)	801 (93%)	57 (7%)	16	22

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LYS
1	A	54	ASN
1	A	67	ASP
1	A	92	LEU
1	A	94	ILE
1	A	107	THR
1	A	120	LEU
1	A	126	LYS
1	A	178	ILE
1	A	194	GLU
1	A	205	LEU
1	A	209	LEU
1	A	218	ASP
1	A	219	LYS
1	A	220	LYS
1	A	264	LEU
1	A	275	LYS
1	A	282	LEU
1	A	283	LEU
1	A	301	LEU

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Mol	Chain	Res	Type
1	A	303	LEU
1	A	310	LEU
1	A	312	GLU
1	A	347	LYS
1	A	350	LYS
1	A	402	TRP
1	A	423	VAL
1	A	448	ARG
1	A	449	GLU
1	A	474	ASN
1	A	479	LEU
1	A	517	LEU
1	A	523	GLU
1	A	529	GLU
1	A	533	LEU
2	B	30	LYS
2	B	48	SER
2	B	80	LEU
2	B	91	GLN
2	B	92	LEU
2	B	113	ASP
2	B	165	THR
2	B	184	MET
2	B	194	GLU
2	B	205	LEU
2	B	209	LEU
2	B	234	LEU
2	B	250	ASP
2	B	260	LEU
2	B	295	LEU
2	B	310	LEU
2	B	323	LYS
2	B	353	LYS
2	B	368	LEU
2	B	369	THR
2	B	374	LYS
2	B	429	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN

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Mol	Chain	Res	Type
1	A	57	ASN
1	A	91	GLN
1	A	145	GLN
1	A	174	GLN
1	A	175	ASN
1	A	207	GLN
1	A	258	GLN
1	A	269	GLN
1	A	306	ASN
1	A	332	GLN
1	A	407	GLN
1	A	500	GLN
1	A	509	GLN
1	A	519	ASN
1	A	520	GLN
1	A	545	ASN
2	B	81	ASN
2	B	85	GLN
2	B	91	GLN
2	B	147	ASN
2	B	151	GLN
2	B	175	ASN
2	B	208	HIS
2	B	255	ASN
2	B	258	GLN
2	B	278	GLN
2	B	306	ASN
2	B	336	GLN
2	B	348	ASN
2	B	367	GLN
2	B	373	GLN
2	B	394	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

6 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
4	SO4	A	564	-	4,4,4	0.23	0	6,6,6	0.09	0
4	SO4	A	563	-	4,4,4	0.25	0	6,6,6	0.14	0
3	65B	A	561	-	29,30,30	2.68	15 (51%)	38,42,42	1.47	4 (10%)
4	SO4	A	565	-	4,4,4	0.27	0	6,6,6	0.09	0
4	SO4	A	562	-	4,4,4	0.25	0	6,6,6	0.10	0
4	SO4	B	441	-	4,4,4	0.49	0	6,6,6	0.41	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	65B	A	561	-	-	0/12/12/12	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	561	65B	C13-N12	6.19	1.41	1.32
3	A	561	65B	C15-N16	5.46	1.42	1.35
3	A	561	65B	C26-C21	5.19	1.48	1.39
3	A	561	65B	C1-C6	3.83	1.46	1.40
3	A	561	65B	C1-C2	3.70	1.46	1.40
3	A	561	65B	C26-C25	3.29	1.44	1.38
3	A	561	65B	C5-C6	2.96	1.43	1.39
3	A	561	65B	C3-C2	2.80	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	561	65B	C22-C21	2.57	1.43	1.39
3	A	561	65B	C11-N16	2.54	1.42	1.34
3	A	561	65B	C25-C24	2.32	1.44	1.39
3	A	561	65B	C23-C24	2.32	1.44	1.39
3	A	561	65B	C23-C22	2.26	1.42	1.38
3	A	561	65B	C5-C4	2.20	1.43	1.39
3	A	561	65B	C6'-C6	2.08	1.55	1.51

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	561	65B	N16-C11-N12	-5.17	118.06	126.23
3	A	561	65B	C11-N12-C13	3.96	122.27	115.18
3	A	561	65B	C14-C13-N12	-2.74	120.53	123.72
3	A	561	65B	C11-N16-C15	2.06	120.74	116.52

There are no chirality outliers.

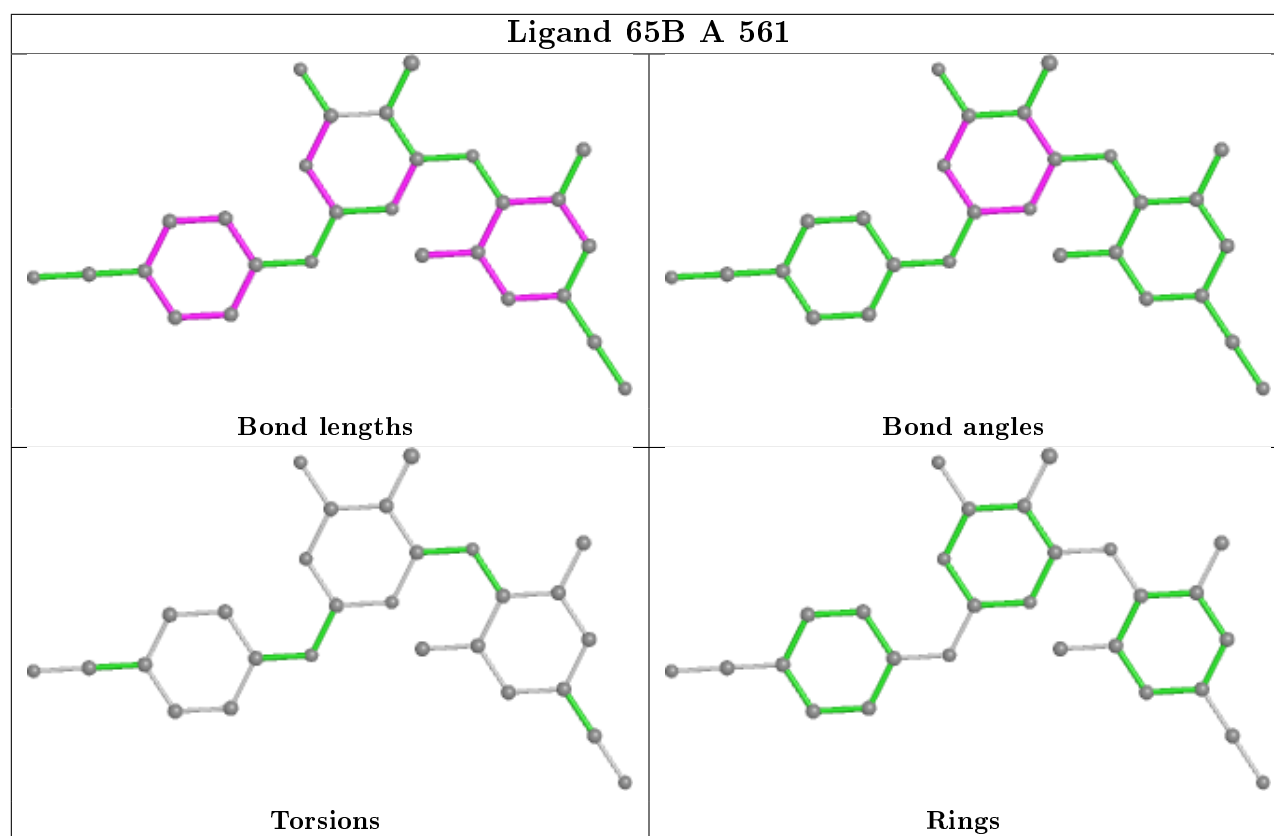
There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	561	65B	2	0
4	B	441	SO4	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	A	550/560 (98%)	0.24	21 (3%)	40	47	30, 50, 90, 122	0
2	B	401/440 (91%)	0.27	25 (6%)	20	26	27, 49, 89, 104	0
All	All	951/1000 (95%)	0.26	46 (4%)	30	37	27, 50, 89, 122	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	ASP	4.8
2	B	87	PHE	3.6
1	A	220	LYS	3.5
1	A	286	THR	3.3
2	B	277	ARG	3.2
2	B	295	LEU	3.2
2	B	14	PRO	2.9
2	B	90	VAL	2.9
1	A	219	LYS	2.9
2	B	92	LEU	2.9
2	B	15	GLY	2.8
2	B	10	VAL	2.8
2	B	276	VAL	2.7
2	B	362	THR	2.7
2	B	283	LEU	2.7
1	A	66	LYS	2.7
2	B	284	ARG	2.6
2	B	298	GLU	2.5
1	A	223	LYS	2.5
1	A	426	TRP	2.5
1	A	67	ASP	2.5
2	B	24	TRP	2.4
2	B	355	ALA	2.4
1	A	552	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	70	LYS	2.4
2	B	281	LYS	2.4
1	A	24	TRP	2.3
1	A	547	GLN	2.3
1	A	406	TRP	2.3
1	A	53	GLU	2.3
2	B	346	PHE	2.3
2	B	238	LYS	2.2
1	A	425	LEU	2.2
2	B	266	TRP	2.2
2	B	12	LEU	2.2
1	A	357	MET	2.2
2	B	70	LYS	2.2
1	A	402	TRP	2.2
2	B	252	TRP	2.1
2	B	279	LEU	2.1
1	A	199	ARG	2.1
1	A	63	ILE	2.1
1	A	194	GLU	2.1
2	B	278	GLN	2.1
2	B	88	TRP	2.1
1	A	22	LYS	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.

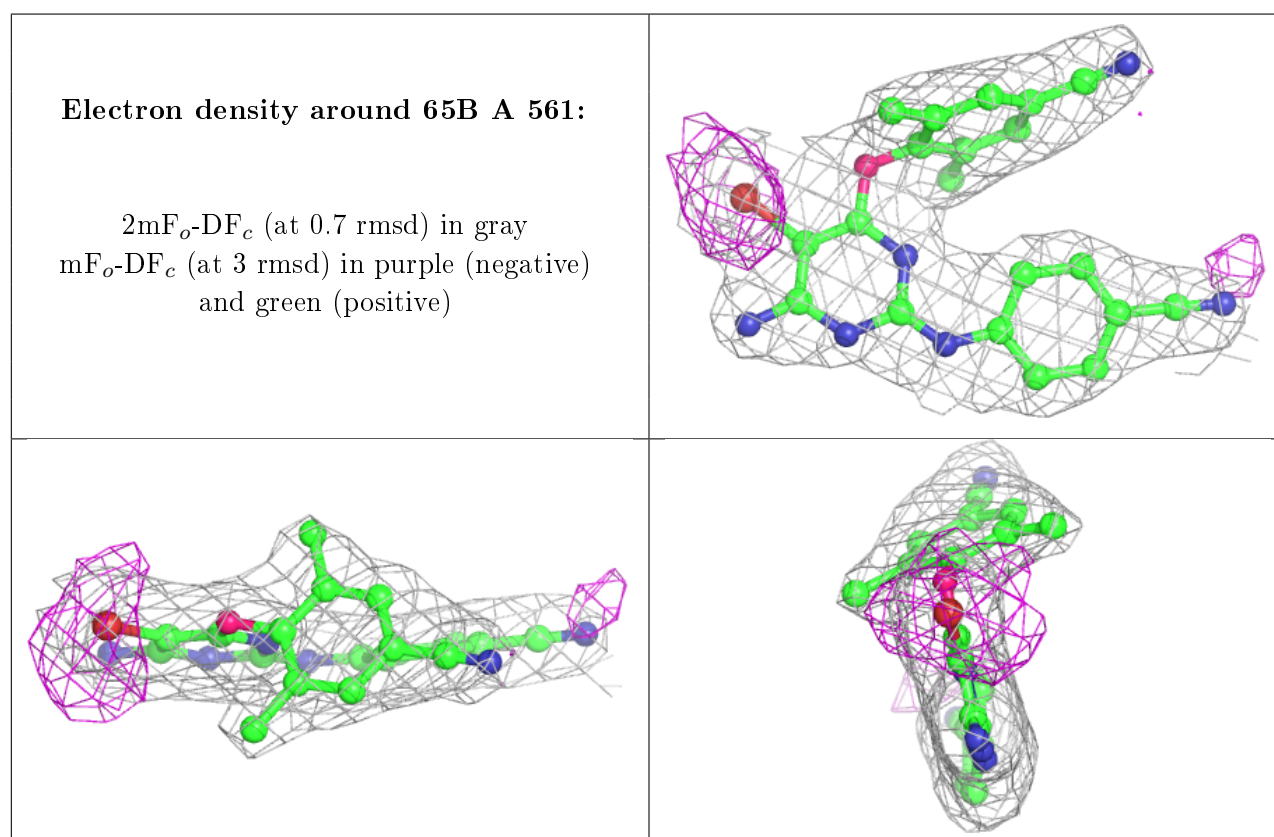
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	A	564	5/5	0.76	0.20	138,138,139,139	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	A	565	5/5	0.91	0.15	105,106,106,106	0
3	65B	A	561	28/28	0.94	0.15	35,38,42,70	0
4	SO4	B	441	5/5	0.94	0.23	96,96,97,97	0
4	SO4	A	563	5/5	0.95	0.14	73,75,76,76	0
4	SO4	A	562	5/5	0.97	0.13	71,72,73,74	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.