



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:15 AM UTC

PDB ID : 3K3E / pdb_00003k3e
Title : Crystal structure of the PDE9A catalytic domain in complex with (R)-BAY73-6691
Authors : Wang, H.; Luo, X.; Ye, M.; Hou, J.; Robinson, H.; Ke, H.
Deposited on : 2009-10-02
Resolution : 2.70 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

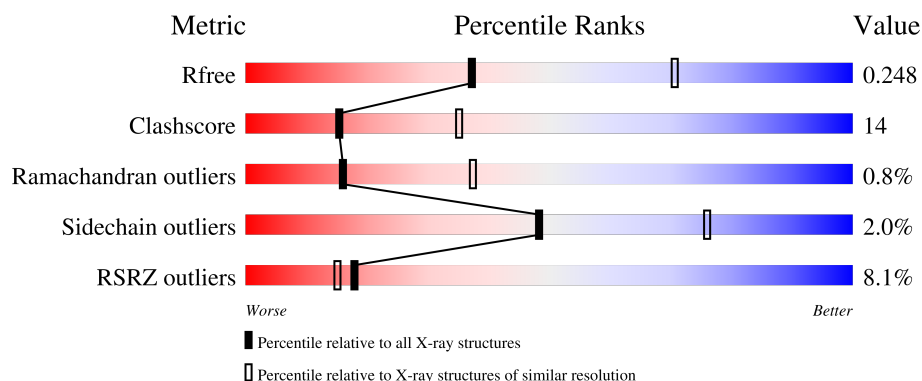
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	326	
1	B	326	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5332 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 High affinity cGMP-specific 3',5'-cyclic phosphodiesterase 9A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	1
			2632	1683	438	481	30			
1	B	320	Total	C	N	O	S	0	0	1
			2632	1683	438	481	30			

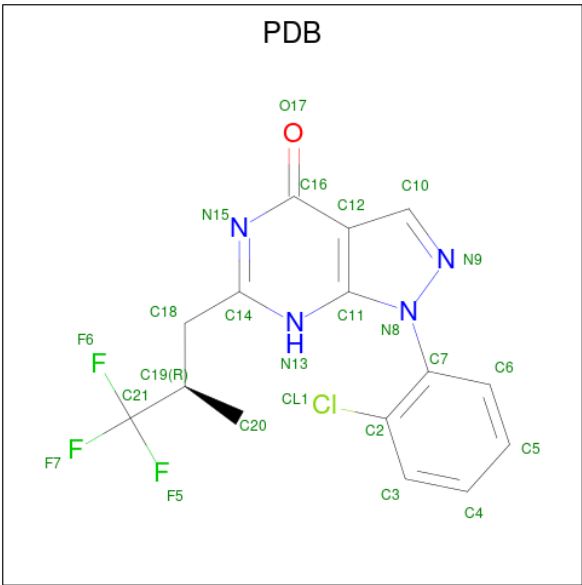
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 1-(2-chlorophenyl)-6-[(2R)-3,3,3-trifluoro-2-methylpropyl]-1,7-dihydro-4H-pyrazolo[3,4-d]pyrimidin-4-one (CCD ID: PDB) (formula: C₁₅H₁₂ClF₃N₄O).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	Cl	F	N	O	0	0
			24	15	1	3	4	1		
4	B	1	Total	C	Cl	F	N	O	0	0
			24	15	1	3	4	1		

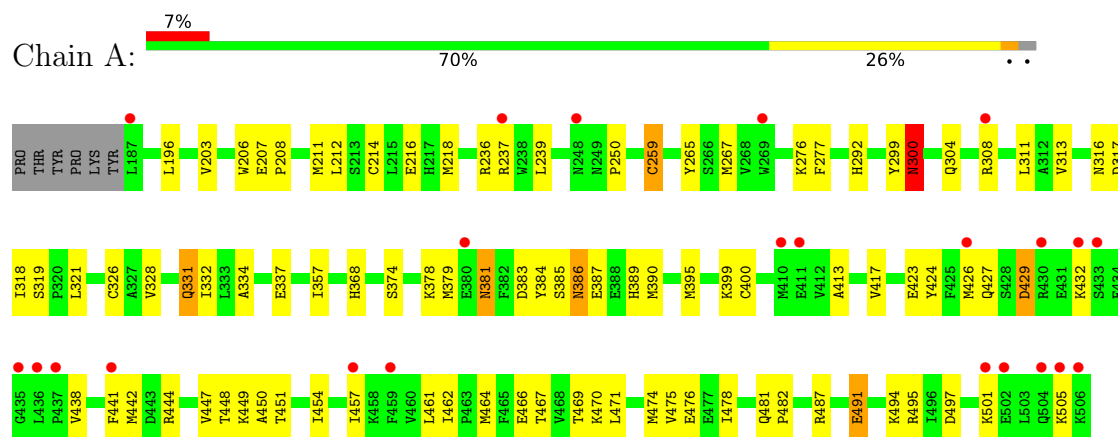
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	10	Total	O	0	0
			10	10		
5	B	6	Total	O	0	0
			6	6		

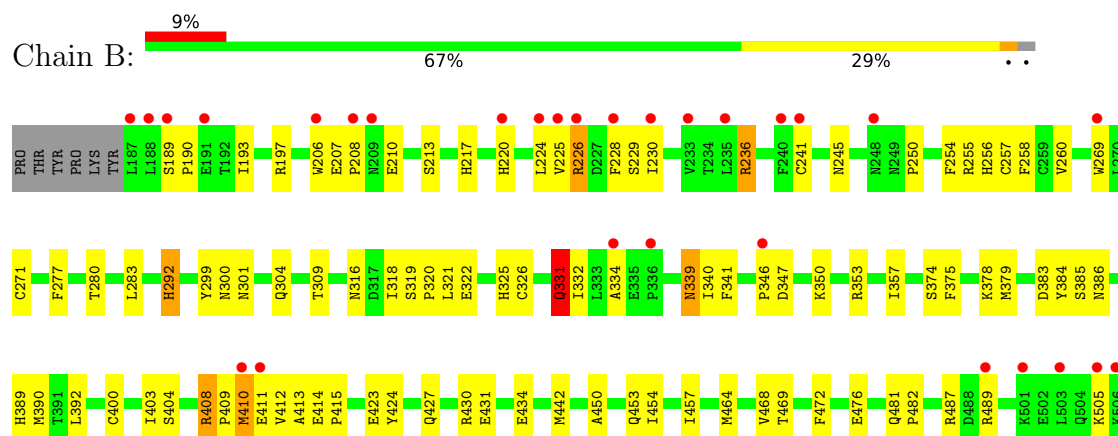
3 Residue-property plots

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: High affinity cGMP-specific 3',5'-cyclic phosphodiesterase 9A



- Molecule 1: High affinity cGMP-specific 3',5'-cyclic phosphodiesterase 9A



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	103.19Å 103.19Å 270.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 30.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.5 (30.00-2.70) 95.4 (30.00-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.37 (at 2.69Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.230 , 0.258 0.222 , 0.248	Depositor DCC
R_{free} test set	4127 reflections (10.04%)	wwPDB-VP
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.542	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5332	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% 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: ZN, PDB, MG

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	0.48	1/2698 (0.0%)	0.95	10/3654 (0.3%)
1	B	0.49	2/2698 (0.1%)	0.96	7/3654 (0.2%)
All	All	0.49	3/5396 (0.1%)	0.95	17/7308 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	410	MET	SD-CE	5.43	1.93	1.79
1	A	505	LYS	C-N	-5.37	1.25	1.33
1	B	505	LYS	C-N	-5.33	1.25	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	412	VAL	N-CA-C	-10.24	100.73	111.58
1	B	408	ARG	CA-C-N	7.81	127.82	119.78
1	B	408	ARG	C-N-CA	7.81	127.82	119.78
1	A	461	LEU	N-CA-C	7.43	119.03	111.07
1	B	241	CYS	N-CA-C	-7.27	103.29	111.07
1	A	206	TRP	N-CA-C	6.78	120.55	109.85
1	B	413	ALA	N-CA-C	6.53	118.19	111.14
1	A	413	ALA	N-CA-C	6.51	118.17	111.14
1	A	478	ILE	N-CA-C	6.20	118.84	111.09
1	A	447	VAL	N-CA-C	6.03	118.31	108.97
1	A	292	HIS	N-CA-C	5.91	120.02	112.34
1	A	331	GLN	N-CA-C	-5.73	105.11	111.36
1	A	429	ASP	N-CA-C	-5.50	104.93	111.03
1	B	331	GLN	N-CA-C	-5.42	105.45	111.36
1	A	300	ASN	N-CA-C	5.26	116.71	110.19
1	A	368	HIS	N-CA-C	5.16	116.59	111.07

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	292	HIS	N-CA-C	5.07	121.59	110.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2632	0	2574	67	0
1	B	2632	0	2572	80	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	24	0	12	0	0
4	B	24	0	12	0	0
5	A	10	0	0	0	0
5	B	6	0	0	0	0
All	All	5332	0	5170	142	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (142) 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:383:ASP:H	1:A:389:HIS:HD2	1.28	0.81
1:A:400:CYS:SG	1:A:464:MET:HE3	2.22	0.79
1:B:207:GLU:HB3	1:B:208:PRO:HD2	1.67	0.77
1:A:454:ILE:HD13	1:A:491:GLU:HG3	1.69	0.74
1:B:206:TRP:H	1:B:206:TRP:CD1	2.04	0.74
1:B:193:ILE:HG23	1:B:217:HIS:NE2	2.02	0.74
1:B:300:ASN:ND2	1:B:442:MET:HE1	2.01	0.74
1:B:411:GLU:CG	1:B:411:GLU:O	2.35	0.73
1:A:386:ASN:HD22	1:A:386:ASN:C	1.96	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:LEU:HD22	1:A:239:LEU:HB3	1.70	0.71
1:A:451:THR:HG22	1:A:494:LYS:HE2	1.73	0.70
1:B:410:MET:SD	1:B:489:ARG:CD	2.81	0.69
1:B:283:LEU:HD21	1:B:341:PHE:HE1	1.59	0.68
1:B:220:HIS:ND1	1:B:225:VAL:HG21	2.09	0.68
1:A:383:ASP:H	1:A:389:HIS:CD2	2.10	0.67
1:B:250:PRO:HG2	1:B:423:GLU:HB2	1.77	0.67
1:A:497:ASP:O	1:A:501:LYS:HG2	1.96	0.66
1:B:300:ASN:HD21	1:B:442:MET:HE1	1.59	0.66
1:A:300:ASN:O	1:A:304:GLN:HG2	1.96	0.66
1:A:457:ILE:HG21	1:A:487:ARG:HB2	1.79	0.65
1:B:374:SER:O	1:B:378:LYS:HG2	1.97	0.65
1:B:410:MET:SD	1:B:489:ARG:HD2	2.38	0.64
1:B:481:GLN:HB3	1:B:482:PRO:CD	2.28	0.64
1:B:378:LYS:HG3	1:B:392:LEU:HD12	1.80	0.63
1:B:245:ASN:ND2	1:B:332:ILE:HD11	2.14	0.62
1:B:220:HIS:CE1	1:B:225:VAL:HG21	2.34	0.61
1:A:196:LEU:HD21	1:A:214:CYS:HB3	1.80	0.61
1:B:277:PHE:CE1	1:B:390:MET:HE3	2.35	0.60
1:B:414:GLU:CD	1:B:489:ARG:HH22	2.10	0.60
1:A:374:SER:O	1:A:378:LYS:HG2	2.03	0.59
1:B:197:ARG:HD3	1:B:269:TRP:CD1	2.37	0.58
1:B:283:LEU:HD21	1:B:341:PHE:CE1	2.39	0.58
1:B:430:ARG:O	1:B:434:GLU:HG3	2.03	0.58
1:A:317:ASP:OD2	1:B:353:ARG:NE	2.36	0.57
1:A:417:VAL:HG11	1:A:449:LYS:HG3	1.86	0.57
1:A:386:ASN:HB3	1:A:389:HIS:HB2	1.86	0.56
1:A:469:THR:HG21	1:A:476:GLU:HB2	1.88	0.56
1:B:226:ARG:HH11	1:B:226:ARG:HB3	1.71	0.56
1:A:432:LYS:HG3	1:A:438:VAL:HG21	1.89	0.55
1:A:308:ARG:HH22	1:B:350:LYS:HG2	1.72	0.55
1:A:462:ILE:O	1:A:466:GLU:HB2	2.07	0.55
1:A:216:GLU:OE2	1:A:236:ARG:NH2	2.40	0.55
1:B:411:GLU:O	1:B:411:GLU:HG2	2.06	0.55
1:B:236:ARG:HH11	1:B:236:ARG:HB2	1.72	0.54
1:B:411:GLU:O	1:B:411:GLU:HG3	2.06	0.54
1:A:386:ASN:HD22	1:A:387:GLU:N	2.04	0.54
1:B:450:ALA:O	1:B:454:ILE:HG13	2.08	0.54
1:B:410:MET:SD	1:B:489:ARG:NE	2.81	0.54
1:A:379:MET:C	1:A:381:ASN:H	2.16	0.54
1:B:271:CYS:HB3	1:B:472:PHE:CD2	2.43	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:TYR:HD2	1:B:427:GLN:HG3	1.73	0.53
1:B:346:PRO:HG2	1:B:347:ASP:H	1.72	0.53
1:A:379:MET:SD	1:A:471:LEU:HD13	2.48	0.53
1:A:300:ASN:C	1:A:300:ASN:HD22	2.15	0.53
1:B:383:ASP:C	1:B:385:SER:H	2.17	0.52
1:B:326:CYS:SG	1:B:357:ILE:HG12	2.49	0.52
1:A:495:ARG:HH11	1:A:495:ARG:HG3	1.74	0.52
1:A:379:MET:HE2	1:A:379:MET:CA	2.40	0.51
1:B:189:SER:HB2	1:B:190:PRO:HD2	1.93	0.51
1:B:469:THR:HG21	1:B:476:GLU:HB2	1.92	0.51
1:A:454:ILE:HD11	1:A:491:GLU:HA	1.93	0.51
1:B:210:GLU:O	1:B:213:SER:HB2	2.10	0.51
1:B:481:GLN:HB3	1:B:482:PRO:HD3	1.93	0.51
1:B:457:ILE:HG21	1:B:487:ARG:HB2	1.92	0.50
1:B:260:VAL:HG21	1:B:292:HIS:CE1	2.47	0.50
1:A:203:VAL:HG11	1:A:259:CYS:SG	2.52	0.50
1:A:328:VAL:HA	1:A:331:GLN:OE1	2.12	0.50
1:A:395:MET:O	1:A:399:LYS:HG3	2.11	0.50
1:B:228:PHE:HZ	1:B:280:THR:HA	1.77	0.50
1:B:331:GLN:O	1:B:334:ALA:HB3	2.12	0.49
1:B:236:ARG:HH11	1:B:236:ARG:CB	2.25	0.49
1:A:454:ILE:CD1	1:A:491:GLU:HG3	2.41	0.49
1:B:228:PHE:O	1:B:229:SER:HB2	2.12	0.49
1:B:375:PHE:CZ	1:B:379:MET:HE3	2.47	0.49
1:B:254:PHE:O	1:B:257:CYS:HB3	2.12	0.49
1:A:383:ASP:N	1:A:389:HIS:HD2	2.05	0.49
1:A:457:ILE:CG2	1:A:487:ARG:HB2	2.43	0.49
1:A:328:VAL:O	1:A:332:ILE:HG13	2.12	0.49
1:A:308:ARG:NH2	1:B:350:LYS:HG2	2.28	0.48
1:A:316:ASN:HB3	1:B:357:ILE:HD13	1.94	0.48
1:B:245:ASN:HD22	1:B:332:ILE:HD11	1.75	0.48
1:A:207:GLU:HB3	1:A:208:PRO:HD2	1.95	0.48
1:A:441:PHE:CE1	1:A:442:MET:HG3	2.49	0.47
1:B:400:CYS:O	1:B:404:SER:HB3	2.13	0.47
1:A:379:MET:C	1:A:381:ASN:N	2.72	0.47
1:B:226:ARG:CB	1:B:226:ARG:NH1	2.77	0.47
1:B:379:MET:HA	1:B:379:MET:HE2	1.97	0.47
1:B:304:GLN:NE2	1:B:309:THR:OG1	2.45	0.47
1:A:481:GLN:HB3	1:A:482:PRO:CD	2.45	0.46
1:B:245:ASN:ND2	1:B:332:ILE:CD1	2.78	0.46
1:A:457:ILE:O	1:A:462:ILE:HG13	2.14	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:PHE:CE1	1:A:390:MET:HE3	2.51	0.46
1:A:400:CYS:HA	1:A:464:MET:CE	2.46	0.46
1:A:318:ILE:O	1:A:319:SER:C	2.58	0.46
1:B:339:ASN:C	1:B:341:PHE:H	2.24	0.46
1:A:426:MET:HE3	1:A:426:MET:HB2	1.73	0.45
1:A:331:GLN:O	1:A:334:ALA:HB3	2.16	0.45
1:B:256:HIS:O	1:B:260:VAL:HG23	2.16	0.45
1:A:300:ASN:C	1:A:300:ASN:ND2	2.74	0.45
1:A:299:TYR:HD2	1:A:427:GLN:HG3	1.81	0.45
1:B:224:LEU:O	1:B:228:PHE:HB2	2.16	0.45
1:B:300:ASN:HA	1:B:321:LEU:HD11	1.99	0.45
1:B:414:GLU:N	1:B:415:PRO:HD2	2.31	0.45
1:A:276:LYS:HB3	1:A:384:TYR:CD2	2.52	0.45
1:A:424:TYR:CG	1:A:442:MET:HG2	2.51	0.45
1:A:429:ASP:OD1	1:A:444:ARG:NH1	2.49	0.45
1:B:245:ASN:HD22	1:B:332:ILE:CD1	2.29	0.45
1:B:383:ASP:O	1:B:385:SER:N	2.49	0.45
1:B:225:VAL:HA	1:B:230:ILE:HB	1.99	0.44
1:B:408:ARG:HB3	1:B:409:PRO:HD2	1.97	0.44
1:A:379:MET:HE2	1:A:379:MET:HA	1.99	0.44
1:A:386:ASN:C	1:A:386:ASN:ND2	2.64	0.44
1:B:226:ARG:HH11	1:B:226:ARG:CB	2.31	0.44
1:A:450:ALA:O	1:A:454:ILE:HG13	2.17	0.44
1:B:403:ILE:HD11	1:B:464:MET:HE1	2.00	0.44
1:A:218:MET:HG2	1:A:265:TYR:CD1	2.52	0.44
1:B:318:ILE:O	1:B:319:SER:C	2.61	0.44
1:B:386:ASN:HB3	1:B:389:HIS:HB2	1.99	0.44
1:B:408:ARG:HB3	1:B:409:PRO:CD	2.48	0.44
1:A:208:PRO:HA	1:A:211:MET:HE3	1.99	0.43
1:A:237:ARG:HD3	1:A:337:GLU:HB2	2.01	0.43
1:A:250:PRO:HG2	1:A:423:GLU:HB2	2.00	0.43
1:B:193:ILE:O	1:B:193:ILE:HG22	2.19	0.43
1:B:255:ARG:O	1:B:258:PHE:HB3	2.19	0.43
1:A:383:ASP:N	1:A:389:HIS:CD2	2.84	0.42
1:B:383:ASP:C	1:B:385:SER:N	2.77	0.42
1:B:453:GLN:O	1:B:457:ILE:HG13	2.20	0.42
1:A:467:THR:HA	1:A:470:LYS:HD2	2.01	0.41
1:B:322:GLU:H	1:B:322:GLU:CD	2.28	0.41
1:A:326:CYS:SG	1:A:357:ILE:HG12	2.60	0.41
1:B:269:TRP:CE3	1:B:269:TRP:HA	2.55	0.41
1:A:383:ASP:OD2	1:A:385:SER:OG	2.37	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:GLU:HA	1:B:325:HIS:CD2	2.56	0.41
1:B:424:TYR:HB3	1:B:442:MET:HG2	2.02	0.41
1:B:301:ASN:OD1	1:B:320:PRO:HG2	2.21	0.41
1:B:427:GLN:O	1:B:431:GLU:HG3	2.19	0.41
1:A:300:ASN:HA	1:A:321:LEU:HD11	2.02	0.41
1:B:383:ASP:N	1:B:389:HIS:HD2	2.19	0.41
1:A:357:ILE:HD13	1:B:316:ASN:HB3	2.02	0.40
1:A:448:THR:O	1:A:449:LYS:C	2.65	0.40
1:A:236:ARG:HG3	1:A:236:ARG:NH1	2.37	0.40
1:A:417:VAL:CG1	1:A:449:LYS:HG3	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	318/326 (98%)	295 (93%)	22 (7%)	1 (0%)	36	60
1	B	318/326 (98%)	288 (91%)	26 (8%)	4 (1%)	9	25
All	All	636/652 (98%)	583 (92%)	48 (8%)	5 (1%)	16	37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	381	ASN
1	B	339	ASN
1	B	384	TYR
1	B	226	ARG
1	B	340	ILE

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	296/303 (98%)	287 (97%)	9 (3%)	36	66
1	B	296/303 (98%)	293 (99%)	3 (1%)	68	86
All	All	592/606 (98%)	580 (98%)	12 (2%)	48	76

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	259	CYS
1	A	267	MET
1	A	300	ASN
1	A	311	LEU
1	A	313	VAL
1	A	386	ASN
1	A	474	MET
1	A	475	VAL
1	A	491	GLU
1	B	236	ARG
1	B	331	GLN
1	B	468	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	248	ASN
1	A	274	GLN
1	A	300	ASN
1	A	304	GLN
1	A	351	GLN
1	A	354	GLN
1	A	381	ASN
1	A	386	ASN
1	A	389	HIS
1	B	245	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	304	GLN
1	B	351	GLN
1	B	389	HIS
1	B	453	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PDB	B	600	-	23,26,26	2.42	2 (8%)	23,39,39	2.65	5 (21%)
4	PDB	A	600	-	23,26,26	2.41	2 (8%)	23,39,39	2.65	5 (21%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PDB	B	600	-	-	2/12/14/14	0/3/3/3
4	PDB	A	600	-	-	5/12/14/14	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	600	PDB	N8-N9	-9.56	1.25	1.38
4	A	600	PDB	N8-N9	-9.48	1.25	1.38
4	B	600	PDB	C7-N8	-5.47	1.35	1.43
4	A	600	PDB	C7-N8	-5.44	1.35	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	600	PDB	C10-N9-N8	8.53	111.40	104.59
4	A	600	PDB	C10-N9-N8	8.50	111.37	104.59
4	B	600	PDB	C7-N8-N9	5.69	124.69	119.12
4	A	600	PDB	C7-N8-N9	5.60	124.61	119.12
4	A	600	PDB	C12-C10-N9	-4.95	106.30	112.45
4	B	600	PDB	C12-C10-N9	-4.89	106.38	112.45
4	A	600	PDB	N13-C11-N8	4.39	133.45	126.84
4	B	600	PDB	N13-C11-N8	4.30	133.31	126.84
4	A	600	PDB	C10-C12-C11	2.18	105.26	103.74
4	B	600	PDB	C10-C12-C11	2.09	105.19	103.74

There are no chirality outliers.

All (7) torsion outliers are listed below:

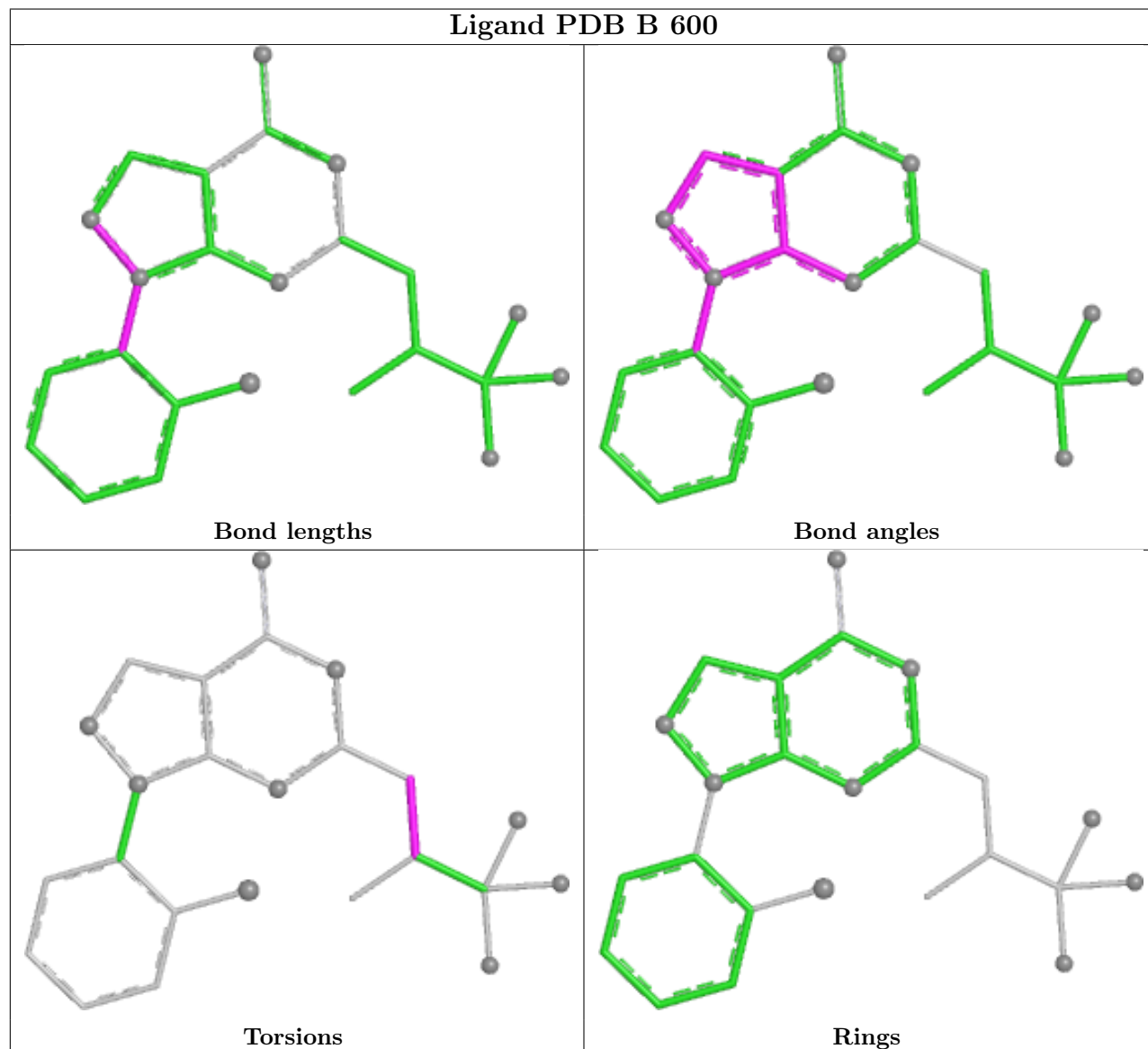
Mol	Chain	Res	Type	Atoms
4	A	600	PDB	C18-C19-C21-F5
4	A	600	PDB	C14-C18-C19-C21
4	B	600	PDB	C14-C18-C19-C21
4	A	600	PDB	C18-C19-C21-F6
4	A	600	PDB	C18-C19-C21-F7
4	A	600	PDB	C14-C18-C19-C20
4	B	600	PDB	C14-C18-C19-C20

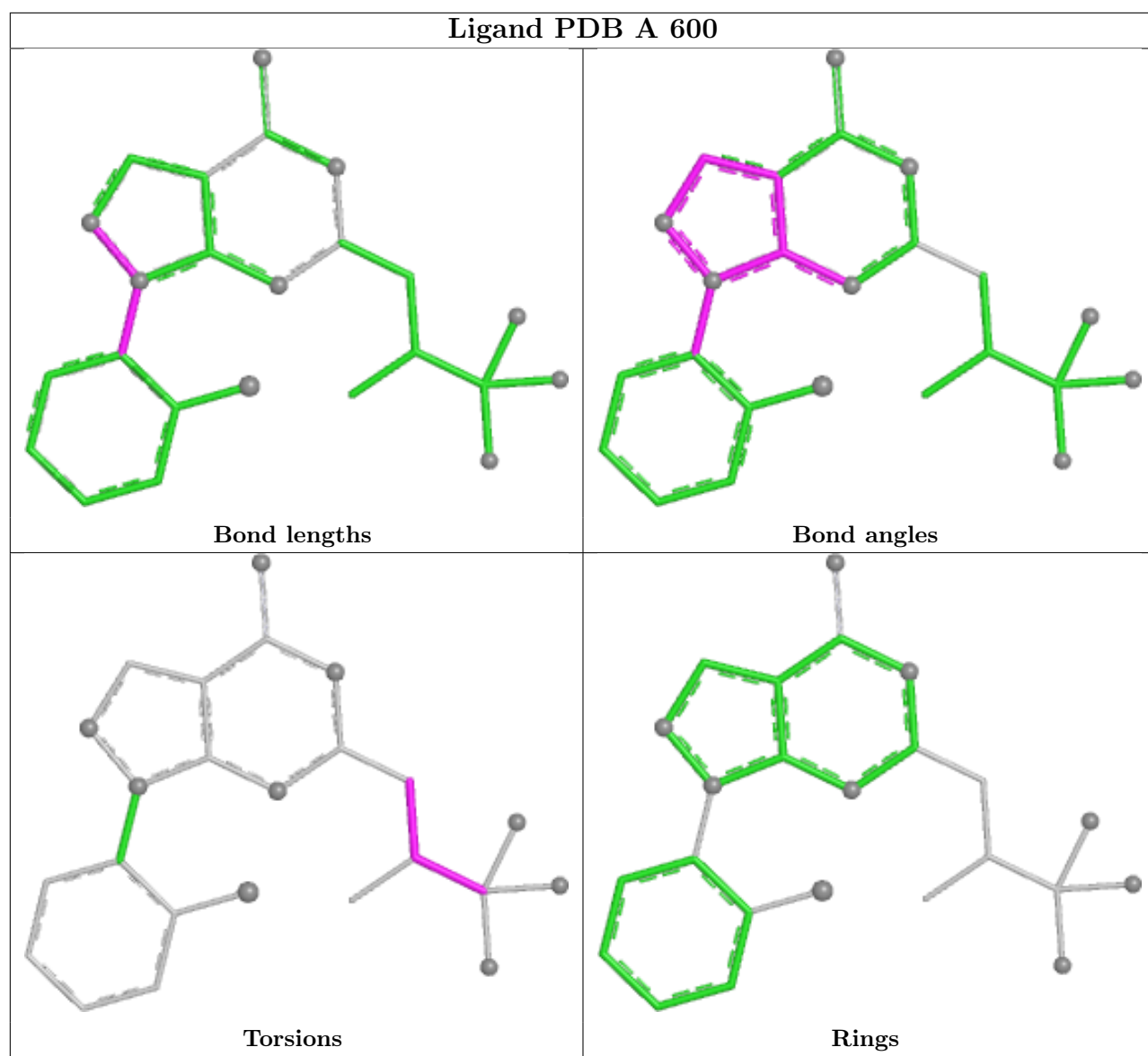
There are no ring outliers.

No monomer is involved in short contacts.

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	320/326 (98%)	0.42	23 (7%) 21 18	33, 54, 77, 92	0
1	B	320/326 (98%)	0.43	29 (9%) 15 12	32, 57, 82, 105	0
All	All	640/652 (98%)	0.43	52 (8%) 18 15	32, 56, 79, 105	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	410	MET	6.6
1	A	505	LYS	6.4
1	A	410	MET	4.9
1	B	209	ASN	4.7
1	B	187	LEU	4.4
1	A	506	LYS	4.1
1	B	411	GLU	3.9
1	B	505	LYS	3.9
1	B	191	GLU	3.6
1	B	269	TRP	3.6
1	B	241	CYS	3.3
1	A	187	LEU	3.2
1	A	308	ARG	3.0
1	A	432	LYS	3.0
1	B	501	LYS	3.0
1	B	336	PRO	2.9
1	B	228	PHE	2.9
1	B	233	VAL	2.8
1	A	441	PHE	2.8
1	A	269	TRP	2.8
1	A	380	GLU	2.7
1	A	459	PHE	2.7
1	B	226	ARG	2.6
1	B	506	LYS	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	435	GLY	2.6
1	B	206	TRP	2.5
1	B	230	ILE	2.5
1	A	237	ARG	2.5
1	B	248	ASN	2.5
1	A	411	GLU	2.5
1	B	188	LEU	2.5
1	A	248	ASN	2.5
1	B	224	LEU	2.5
1	B	235	LEU	2.4
1	A	502	GLU	2.4
1	B	240	PHE	2.4
1	A	457	ILE	2.4
1	B	208	PRO	2.4
1	B	225	VAL	2.4
1	B	503	LEU	2.3
1	B	334	ALA	2.3
1	A	501	LYS	2.3
1	A	504	GLN	2.2
1	B	220	HIS	2.2
1	B	346	PRO	2.2
1	A	426	MET	2.2
1	A	430	ARG	2.1
1	A	437	PRO	2.1
1	A	433	SER	2.1
1	B	189	SER	2.1
1	B	489	ARG	2.0
1	A	436	LEU	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.

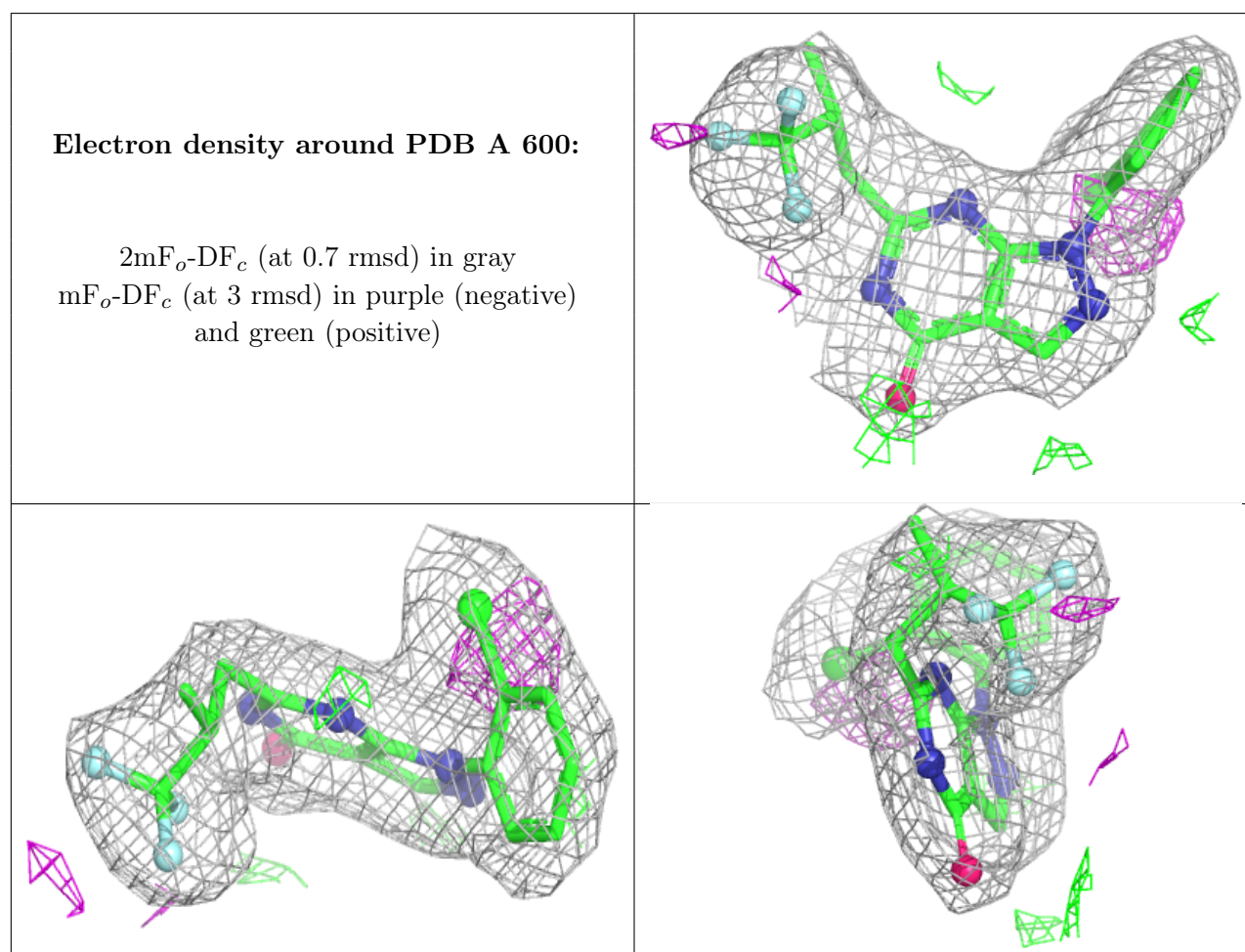
6.4 Ligands [i](#)

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.

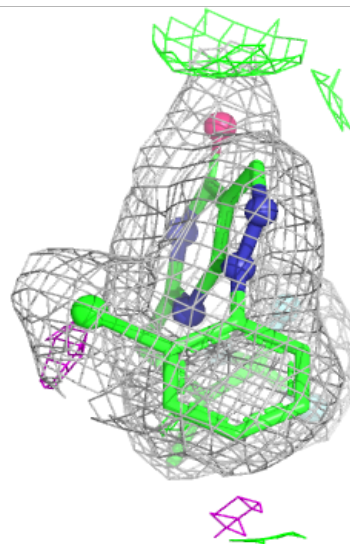
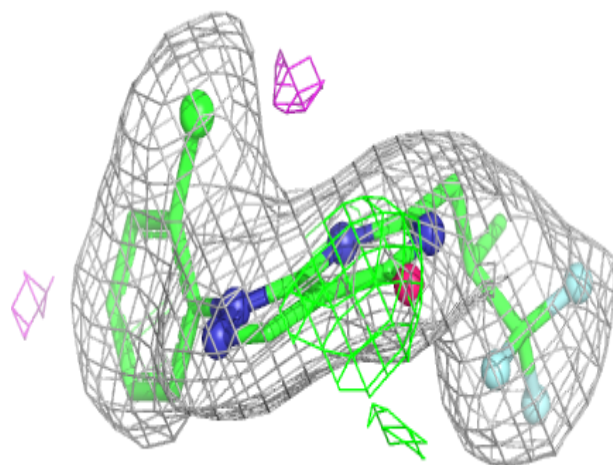
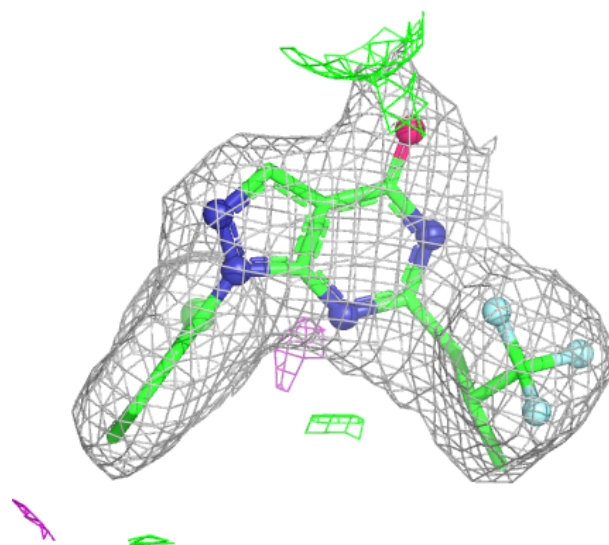
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	B	602	1/1	0.90	0.09	50,50,50,50	0
4	PDB	A	600	24/24	0.95	0.09	41,47,51,53	0
4	PDB	B	600	24/24	0.97	0.07	40,44,45,47	0
3	MG	A	602	1/1	0.98	0.04	37,37,37,37	0
2	ZN	A	601	1/1	0.99	0.02	44,44,44,44	0
2	ZN	B	601	1/1	1.00	0.04	47,47,47,47	0

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 PDB B 600:

$2mF_o-DF_c$ (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.