



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2026 – 04:15 PM EDT

PDB ID : 9ECK / pdb_00009eck
Title : Estrogen receptor alpha
Authors : Lee, C.C.
Deposited on : 2024-11-14
Resolution : 3.20 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

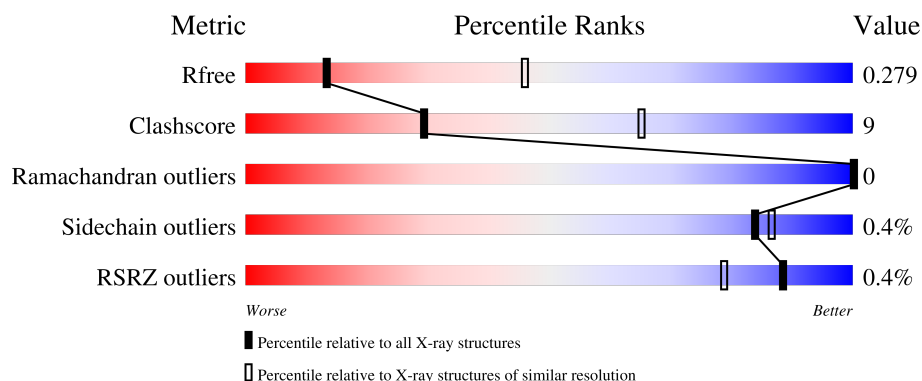
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	259	
1	B	259	
1	C	259	
1	D	259	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13867 atoms, of which 6860 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 Estrogen receptor.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	232	Total	C	H	N	O	S	0	0	0
			3417	1104	1692	296	312	13			
1	B	220	Total	C	H	N	O	S	0	0	0
			3292	1068	1629	279	301	15			
1	C	233	Total	C	H	N	O	S	0	0	0
			3506	1129	1736	306	321	14			
1	D	223	Total	C	H	N	O	S	0	0	0
			3432	1105	1711	286	314	16			

There are 24 discrepancies between the modelled and reference sequences:

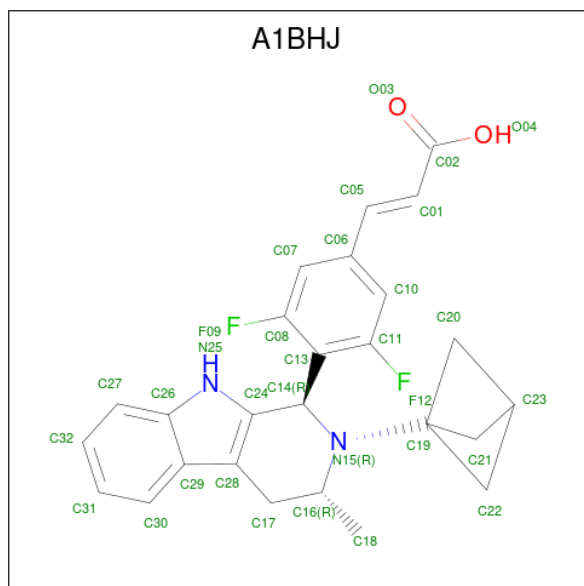
Chain	Residue	Modelled	Actual	Comment	Reference
A	295	SER	-	expression tag	UNP P03372
A	296	ASN	-	expression tag	UNP P03372
A	297	ALA	-	expression tag	UNP P03372
A	381	SER	CYS	engineered mutation	UNP P03372
A	417	SER	CYS	engineered mutation	UNP P03372
A	536	SER	LEU	engineered mutation	UNP P03372
B	295	SER	-	expression tag	UNP P03372
B	296	ASN	-	expression tag	UNP P03372
B	297	ALA	-	expression tag	UNP P03372
B	381	SER	CYS	engineered mutation	UNP P03372
B	417	SER	CYS	engineered mutation	UNP P03372
B	536	SER	LEU	engineered mutation	UNP P03372
C	295	SER	-	expression tag	UNP P03372
C	296	ASN	-	expression tag	UNP P03372
C	297	ALA	-	expression tag	UNP P03372
C	381	SER	CYS	engineered mutation	UNP P03372
C	417	SER	CYS	engineered mutation	UNP P03372
C	536	SER	LEU	engineered mutation	UNP P03372
D	295	SER	-	expression tag	UNP P03372
D	296	ASN	-	expression tag	UNP P03372
D	297	ALA	-	expression tag	UNP P03372

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Chain	Residue	Modelled	Actual	Comment	Reference
D	381	SER	CYS	engineered mutation	UNP P03372
D	417	SER	CYS	engineered mutation	UNP P03372
D	536	SER	LEU	engineered mutation	UNP P03372

- Molecule 2 is (2E)-3-{4-[(1R,3R)-2-(bicyclo[1.1.1]pentan-1-yl)-3-methyl-2,3,4,9-tetrahydro-1H-pyrido[3,4-b]indol-1-yl]-3,5-difluorophenyl}prop-2-enoic acid (CCD ID: A1BHJ) (formula: C₂₆H₂₄F₂N₂O₂) (labeled as "Ligand of Interest" by depositor).

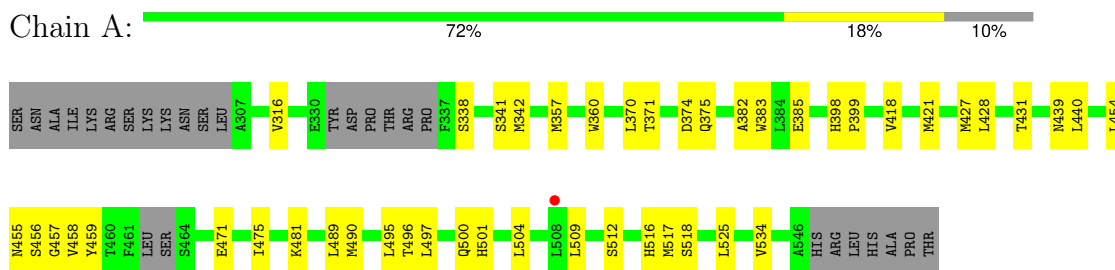


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 55	C 26	F 2	H 23	N 2	O 2	0	0
2	B	1	Total 55	C 26	F 2	H 23	N 2	O 2	0	0
2	C	1	Total 55	C 26	F 2	H 23	N 2	O 2	0	0
2	D	1	Total 55	C 26	F 2	H 23	N 2	O 2	0	0

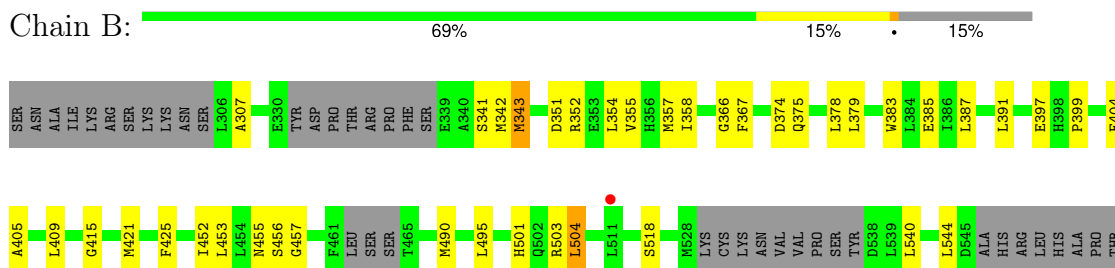
3 Residue-property plots

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.

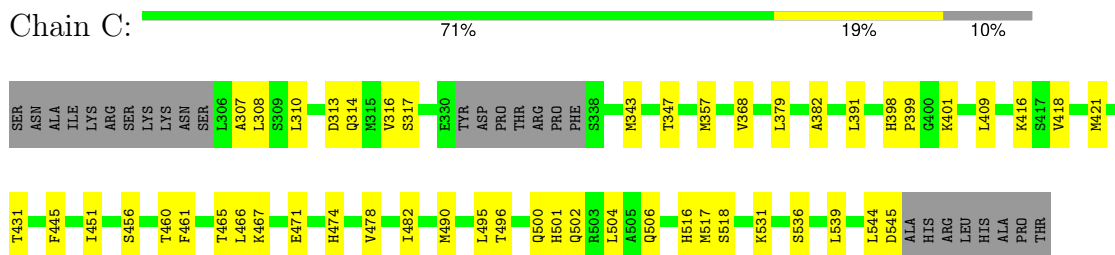
• Molecule 1: Estrogen receptor



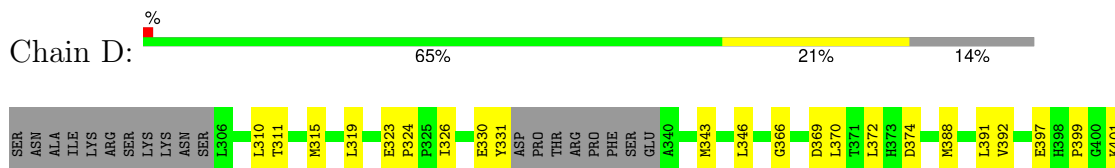
• Molecule 1: Estrogen receptor

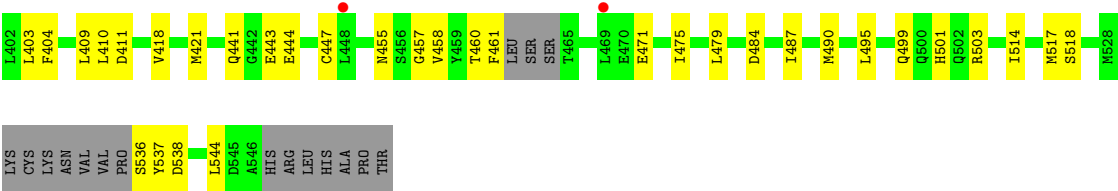


• Molecule 1: Estrogen receptor



• Molecule 1: Estrogen receptor





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.62Å 59.41Å 100.85Å 94.05° 93.83° 115.36°	Depositor
Resolution (Å)	23.77 – 3.20 23.77 – 3.20	Depositor EDS
% Data completeness (in resolution range)	88.8 (23.77-3.20) 88.6 (23.77-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 3.23Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.212 , 0.293 0.227 , 0.279	Depositor DCC
R_{free} test set	955 reflections (4.48%)	wwPDB-VP
Wilson B-factor (Å ²)	98.4	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 90.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13867	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.16% 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: A1BHJ

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.17	4/1751 (0.2%)	1.14	5/2379 (0.2%)
1	B	1.16	2/1690 (0.1%)	1.06	7/2292 (0.3%)
1	C	1.31	7/1800 (0.4%)	1.16	11/2441 (0.5%)
1	D	1.08	1/1749 (0.1%)	1.10	3/2368 (0.1%)
All	All	1.18	14/6990 (0.2%)	1.12	26/9480 (0.3%)

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	478	VAL	C-O	-6.63	1.16	1.24
1	C	516	HIS	CA-C	-5.76	1.45	1.52
1	C	518	SER	CA-C	-5.75	1.45	1.52
1	C	445	PHE	CA-C	-5.62	1.45	1.52
1	C	451	ILE	C-O	-5.50	1.17	1.24
1	A	509	LEU	C-O	-5.49	1.17	1.24
1	C	391	LEU	CA-C	-5.47	1.45	1.52
1	A	512	SER	CA-C	-5.29	1.45	1.52
1	C	313	ASP	C-O	-5.27	1.17	1.24
1	D	518	SER	CA-C	-5.25	1.46	1.52
1	A	518	SER	CA-C	-5.20	1.46	1.52
1	B	457	GLY	C-O	-5.19	1.20	1.24
1	B	518	SER	CA-C	-5.15	1.46	1.52
1	A	481	LYS	C-O	-5.08	1.18	1.24

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	459	TYR	N-CA-C	11.96	124.06	111.14
1	C	307	ALA	N-CA-C	-8.43	102.04	111.82
1	B	504	LEU	N-CA-C	-6.73	103.04	111.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	516	HIS	N-CA-C	-6.47	104.15	111.07
1	D	311	THR	N-CA-C	-6.47	102.17	110.19
1	C	504	LEU	N-CA-C	-6.42	104.19	111.07
1	D	447	CYS	N-CA-C	-6.25	104.39	111.07
1	B	343	MET	N-CA-C	-6.16	104.48	111.07
1	C	506	GLN	N-CA-C	-6.05	104.60	111.07
1	A	504	LEU	N-CA-C	-5.97	104.68	111.07
1	A	385	GLU	N-CA-C	-5.83	104.92	111.28
1	D	444	GLU	N-CA-C	-5.82	105.02	111.36
1	C	516	HIS	N-CA-C	-5.62	105.06	111.07
1	C	451	ILE	N-CA-C	-5.59	104.92	110.62
1	C	382	ALA	N-CA-C	5.36	119.80	112.88
1	C	502	GLN	N-CA-C	-5.36	105.34	111.07
1	C	461	PHE	N-CA-C	5.27	116.89	110.41
1	B	341	SER	N-CA-C	-5.17	105.64	111.28
1	B	501	HIS	CB-CA-C	-5.17	102.06	110.85
1	B	405	ALA	CA-C-N	-5.17	114.29	119.56
1	B	405	ALA	C-N-CA	-5.17	114.29	119.56
1	C	531	LYS	N-CA-C	5.07	121.59	110.80
1	C	456	SER	N-CA-C	5.05	117.17	111.11
1	A	360	TRP	N-CA-C	-5.04	105.24	111.33
1	B	409	LEU	N-CA-C	-5.03	100.98	109.07
1	C	482	ILE	N-CA-C	-5.01	104.82	111.09

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	1725	1692	1691	26	0
1	B	1663	1629	1629	25	0
1	C	1770	1736	1736	31	0
1	D	1721	1711	1709	39	0
2	A	32	23	0	0	0
2	B	32	23	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	32	23	0	0	0
2	D	32	23	0	0	0
All	All	7007	6860	6765	119	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 9.

All (119) 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:497:LEU:HD11	1:A:501:HIS:CE1	2.09	0.86
1:C:418:VAL:HG13	1:C:421:MET:HB2	1.63	0.80
1:D:484:ASP:O	1:D:487:ILE:HG22	1.84	0.77
1:C:501:HIS:NE2	1:D:501:HIS:NE2	2.38	0.72
1:D:495:LEU:HD22	1:D:499:GLN:HB3	1.73	0.71
1:D:326:ILE:O	1:D:326:ILE:HG13	1.91	0.69
1:C:343:MET:O	1:C:347:THR:HG23	1.94	0.67
1:C:544:LEU:HD23	1:C:544:LEU:O	1.96	0.66
1:A:525:LEU:HD21	1:A:534:VAL:HG11	1.77	0.65
1:B:367:PHE:CE2	1:B:453:LEU:HD21	2.32	0.64
1:D:403:LEU:HD13	1:D:409:LEU:HD12	1.79	0.63
1:C:460:THR:O	1:C:460:THR:HG22	1.98	0.63
1:A:496:THR:O	1:A:500:GLN:HG3	2.01	0.61
1:B:351:ASP:HA	1:B:354:LEU:HD12	1.82	0.60
1:D:460:THR:O	1:D:461:PHE:C	2.44	0.60
1:A:382:ALA:HB2	1:A:456:SER:CB	2.33	0.59
1:D:346:LEU:HD22	1:D:404:PHE:CD2	2.37	0.59
1:D:372:LEU:HD11	1:D:544:LEU:HD13	1.85	0.58
1:D:401:LYS:HA	1:D:410:LEU:O	2.04	0.57
1:A:497:LEU:O	1:A:500:GLN:HB2	2.05	0.57
1:D:331:TYR:CD2	1:D:331:TYR:O	2.58	0.57
1:D:404:PHE:CD1	1:D:410:LEU:HD13	2.40	0.57
1:B:540:LEU:H	1:B:540:LEU:HD12	1.68	0.57
1:D:475:ILE:O	1:D:479:LEU:HD13	2.05	0.57
1:D:391:LEU:HD13	1:D:404:PHE:HA	1.86	0.56
1:D:410:LEU:HD12	1:D:410:LEU:N	2.20	0.56
1:B:383:TRP:O	1:B:387:LEU:HD13	2.07	0.55
1:D:310:LEU:HB2	1:D:315:MET:HE2	1.87	0.55
1:A:371:THR:HG23	1:A:471:GLU:OE2	2.07	0.55
1:C:368:VAL:HG12	1:C:368:VAL:O	2.07	0.54
1:C:418:VAL:HG13	1:C:421:MET:CB	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:343:MET:CA	1:C:343:MET:HE2	2.38	0.54
1:C:490:MET:HB3	1:C:495:LEU:HD12	1.90	0.53
1:D:403:LEU:CD1	1:D:409:LEU:HD12	2.38	0.53
1:A:428:LEU:O	1:A:431:THR:OG1	2.27	0.53
1:B:379:LEU:HD22	1:B:453:LEU:HD11	1.90	0.53
1:C:465:THR:HG22	1:C:466:LEU:N	2.24	0.53
1:D:490:MET:SD	1:D:503:ARG:HD3	2.48	0.52
1:B:367:PHE:CE1	1:B:375:GLN:HG2	2.44	0.52
1:C:316:VAL:HG23	1:C:317:SER:N	2.24	0.52
1:A:497:LEU:CD1	1:A:501:HIS:CE1	2.89	0.51
1:B:490:MET:HB3	1:B:495:LEU:HD22	1.92	0.51
1:C:544:LEU:HD23	1:C:544:LEU:C	2.36	0.50
1:A:490:MET:HB3	1:A:495:LEU:HD12	1.94	0.50
1:D:479:LEU:HD12	1:D:479:LEU:N	2.26	0.49
1:C:357:MET:HE3	1:C:379:LEU:HD22	1.94	0.49
1:D:536:SER:O	1:D:537:TYR:C	2.56	0.49
1:B:540:LEU:HD12	1:B:540:LEU:N	2.28	0.49
1:D:475:ILE:O	1:D:479:LEU:CD1	2.61	0.49
1:D:315:MET:O	1:D:319:LEU:HG	2.12	0.49
1:D:397:GLU:C	1:D:399:PRO:HD3	2.38	0.49
1:D:323:GLU:HG3	1:D:324:PRO:HD2	1.94	0.48
1:A:338:SER:H	1:A:341:SER:HB3	1.77	0.48
1:C:471:GLU:O	1:C:474:HIS:HB3	2.13	0.48
1:D:331:TYR:O	1:D:331:TYR:CG	2.67	0.48
1:A:316:VAL:HG21	1:A:489:LEU:HD21	1.95	0.48
1:D:441:GLN:HB3	1:D:443:GLU:OE1	2.13	0.47
1:A:357:MET:HE1	1:A:383:TRP:O	2.15	0.47
1:D:388:MET:O	1:D:392:VAL:HG23	2.13	0.47
1:B:503:ARG:O	1:B:504:LEU:C	2.58	0.47
1:C:501:HIS:CE1	1:D:501:HIS:HE2	2.33	0.47
1:A:370:LEU:O	1:A:375:GLN:NE2	2.38	0.47
1:B:540:LEU:H	1:B:540:LEU:CD1	2.28	0.47
1:D:411:ASP:OD1	1:D:411:ASP:C	2.58	0.47
1:C:401:LYS:HB2	1:C:409:LEU:HD11	1.98	0.46
1:B:374:ASP:O	1:B:378:LEU:HG	2.16	0.46
1:B:540:LEU:O	1:B:544:LEU:HD13	2.16	0.46
1:B:455:ASN:O	1:B:456:SER:C	2.59	0.46
1:A:398:HIS:O	1:A:399:PRO:C	2.59	0.45
1:D:370:LEU:HD23	1:D:374:ASP:HB3	1.96	0.45
1:A:427:MET:O	1:A:431:THR:HG23	2.16	0.45
1:C:431:THR:HG21	1:C:517:MET:CE	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:374:ASP:OD2	1:D:471:GLU:OE1	2.34	0.45
1:B:357:MET:O	1:B:358:ILE:C	2.58	0.45
1:D:404:PHE:HD1	1:D:410:LEU:HD13	1.82	0.45
1:B:375:GLN:O	1:B:379:LEU:HD23	2.16	0.44
1:C:460:THR:O	1:C:460:THR:CG2	2.64	0.44
1:B:383:TRP:O	1:B:387:LEU:CD1	2.65	0.44
1:B:421:MET:CE	1:B:425:PHE:CE1	3.00	0.44
1:C:431:THR:HG21	1:C:517:MET:HE1	2.00	0.44
1:A:439:ASN:O	1:A:440:LEU:C	2.59	0.44
1:C:418:VAL:CG1	1:C:421:MET:HB2	2.42	0.44
1:D:343:MET:HE1	1:D:421:MET:HE1	1.99	0.44
1:B:385:GLU:HB3	1:B:452:ILE:HG21	2.00	0.44
1:A:454:LEU:HB3	1:A:475:ILE:HG23	1.99	0.43
1:A:370:LEU:HB2	1:A:375:GLN:HG3	1.99	0.43
1:B:415:GLY:HA3	1:B:425:PHE:CD2	2.54	0.43
1:C:536:SER:CB	1:C:539:LEU:HB3	2.49	0.43
1:B:352:ARG:O	1:B:355:VAL:HG22	2.19	0.42
1:C:398:HIS:O	1:C:399:PRO:C	2.62	0.42
1:A:455:ASN:O	1:A:456:SER:C	2.61	0.42
1:B:342:MET:O	1:B:343:MET:C	2.58	0.42
1:C:465:THR:CG2	1:C:466:LEU:N	2.81	0.42
1:A:371:THR:O	1:A:375:GLN:HG3	2.20	0.42
1:A:457:GLY:O	1:A:458:VAL:C	2.61	0.42
1:A:525:LEU:HG	1:A:534:VAL:HG21	2.02	0.42
1:D:457:GLY:O	1:D:458:VAL:C	2.59	0.42
1:D:495:LEU:HD22	1:D:499:GLN:CB	2.47	0.42
1:B:495:LEU:HD12	1:B:495:LEU:N	2.35	0.42
1:C:467:LYS:O	1:C:471:GLU:HG3	2.20	0.42
1:C:545:ASP:N	1:C:545:ASP:OD1	2.53	0.42
1:D:536:SER:C	1:D:538:ASP:N	2.72	0.42
1:C:343:MET:HE2	1:C:343:MET:HA	2.02	0.42
1:C:496:THR:O	1:C:500:GLN:HG3	2.20	0.42
1:A:342:MET:HE2	1:A:342:MET:HB3	1.89	0.41
1:A:418:VAL:HB	1:A:421:MET:HG3	2.01	0.41
1:A:431:THR:HG21	1:A:517:MET:CE	2.50	0.41
1:C:316:VAL:CG2	1:C:317:SER:N	2.84	0.41
1:C:465:THR:HG22	1:C:467:LYS:H	1.85	0.41
1:D:330:GLU:O	1:D:331:TYR:HB3	2.20	0.41
1:D:401:LYS:CG	1:D:409:LEU:HD21	2.51	0.41
1:A:374:ASP:CB	1:A:471:GLU:OE1	2.68	0.41
1:B:397:GLU:C	1:B:399:PRO:HD3	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:514:ILE:HA	1:D:517:MET:CE	2.51	0.41
1:C:416:LYS:C	1:C:418:VAL:N	2.79	0.40
1:D:366:GLY:HA2	1:D:369:ASP:OD2	2.21	0.40
1:B:307:ALA:CB	1:B:366:GLY:N	2.84	0.40
1:B:391:LEU:HD13	1:B:404:PHE:HA	2.04	0.40
1:C:310:LEU:HD22	1:C:314:GLN:CB	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	226/259 (87%)	222 (98%)	4 (2%)	0	100	100
1	B	212/259 (82%)	209 (99%)	3 (1%)	0	100	100
1	C	229/259 (88%)	224 (98%)	5 (2%)	0	100	100
1	D	215/259 (83%)	214 (100%)	1 (0%)	0	100	100
All	All	882/1036 (85%)	869 (98%)	13 (2%)	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	177/233 (76%)	177 (100%)	0	100	100
1	B	174/233 (75%)	174 (100%)	0	100	100
1	C	185/233 (79%)	184 (100%)	1 (0%)	81	85
1	D	185/233 (79%)	183 (99%)	2 (1%)	65	79
All	All	721/932 (77%)	718 (100%)	3 (0%)	84	86

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	308	LEU
1	D	418	VAL
1	D	455	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	375	GLN
1	B	439	ASN
1	C	377	HIS
1	C	455	ASN
1	C	519	ASN
1	D	348	ASN
1	D	500	GLN
1	D	516	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	A1BHJ	B	601	-	33,37,37	1.53	7 (21%)	48,58,58	1.65	10 (20%)
2	A1BHJ	D	601	-	33,37,37	1.42	5 (15%)	48,58,58	1.67	9 (18%)
2	A1BHJ	A	601	-	33,37,37	1.58	4 (12%)	48,58,58	1.74	13 (27%)
2	A1BHJ	C	601	-	33,37,37	1.61	6 (18%)	48,58,58	1.68	10 (20%)

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	A1BHJ	B	601	-	-	2/9/46/46	0/7/6/6
2	A1BHJ	D	601	-	-	2/9/46/46	0/7/6/6
2	A1BHJ	A	601	-	-	0/9/46/46	0/7/6/6
2	A1BHJ	C	601	-	-	4/9/46/46	0/7/6/6

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	A1BHJ	C14-C24	4.47	1.56	1.50
2	A	601	A1BHJ	C14-C24	3.76	1.55	1.50
2	B	601	A1BHJ	C14-N15	3.45	1.52	1.47
2	D	601	A1BHJ	C18-C16	3.36	1.63	1.52
2	A	601	A1BHJ	C13-C11	3.10	1.42	1.39
2	C	601	A1BHJ	C18-C16	3.07	1.62	1.52
2	B	601	A1BHJ	C13-C11	2.98	1.42	1.39
2	A	601	A1BHJ	C07-C08	2.85	1.42	1.37
2	C	601	A1BHJ	C07-C08	2.77	1.42	1.37
2	D	601	A1BHJ	C14-C24	2.72	1.54	1.50
2	B	601	A1BHJ	C18-C16	2.71	1.61	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	A1BHJ	C14-C24	2.69	1.54	1.50
2	D	601	A1BHJ	C07-C08	2.64	1.42	1.37
2	D	601	A1BHJ	C13-C11	2.56	1.42	1.39
2	B	601	A1BHJ	C07-C08	2.42	1.41	1.37
2	D	601	A1BHJ	C17-C28	-2.41	1.47	1.50
2	A	601	A1BHJ	C18-C16	2.39	1.60	1.52
2	C	601	A1BHJ	C13-C08	2.24	1.41	1.39
2	B	601	A1BHJ	C24-C28	2.18	1.39	1.36
2	C	601	A1BHJ	C24-C28	2.15	1.39	1.36
2	B	601	A1BHJ	C10-C11	2.08	1.41	1.37
2	C	601	A1BHJ	C27-C26	2.05	1.43	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	A1BHJ	C14-C24-N25	4.92	129.11	123.78
2	C	601	A1BHJ	C14-C24-N25	4.46	128.61	123.78
2	B	601	A1BHJ	C14-C24-N25	4.06	128.17	123.78
2	D	601	A1BHJ	C14-C24-N25	4.02	128.13	123.78
2	D	601	A1BHJ	C16-C17-C28	3.71	115.03	110.17
2	C	601	A1BHJ	C06-C10-C11	3.65	122.11	119.27
2	A	601	A1BHJ	C06-C10-C11	3.62	122.09	119.27
2	D	601	A1BHJ	C06-C10-C11	3.39	121.91	119.27
2	A	601	A1BHJ	C16-C17-C28	3.35	114.56	110.17
2	B	601	A1BHJ	C22-C19-C20	-3.35	85.03	88.28
2	C	601	A1BHJ	C22-C19-C21	-3.32	85.06	88.28
2	B	601	A1BHJ	C16-C17-C28	3.26	114.44	110.17
2	D	601	A1BHJ	C07-C06-C05	3.01	129.60	120.61
2	A	601	A1BHJ	C07-C06-C05	2.99	129.56	120.61
2	C	601	A1BHJ	F09-C08-C13	2.97	121.72	118.14
2	B	601	A1BHJ	C06-C10-C11	2.95	121.56	119.27
2	B	601	A1BHJ	C07-C06-C05	2.91	129.31	120.61
2	B	601	A1BHJ	F09-C08-C13	2.90	121.63	118.14
2	C	601	A1BHJ	C07-C06-C05	2.82	129.03	120.61
2	D	601	A1BHJ	C21-C19-C20	-2.78	85.59	88.28
2	A	601	A1BHJ	F09-C08-C13	2.73	121.43	118.14
2	A	601	A1BHJ	C28-C24-N25	-2.70	107.80	110.22
2	D	601	A1BHJ	C10-C06-C05	-2.70	112.53	120.61
2	D	601	A1BHJ	C28-C24-N25	-2.68	107.82	110.22
2	B	601	A1BHJ	C10-C06-C05	-2.62	112.77	120.61
2	D	601	A1BHJ	F09-C08-C13	2.54	121.19	118.14
2	C	601	A1BHJ	C06-C05-C01	-2.52	121.41	126.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	A1BHJ	C10-C06-C05	-2.48	113.21	120.61
2	C	601	A1BHJ	C10-C06-C05	-2.47	113.23	120.61
2	B	601	A1BHJ	C28-C24-N25	-2.34	108.12	110.22
2	C	601	A1BHJ	C17-C28-C24	-2.34	118.90	121.71
2	A	601	A1BHJ	C14-C24-C28	-2.31	123.07	125.50
2	C	601	A1BHJ	C28-C24-N25	-2.27	108.19	110.22
2	C	601	A1BHJ	C14-C24-C28	-2.26	123.12	125.50
2	D	601	A1BHJ	F12-C11-C13	2.21	120.80	118.14
2	A	601	A1BHJ	C06-C05-C01	-2.19	122.12	126.92
2	A	601	A1BHJ	C21-C19-C20	-2.14	86.20	88.28
2	A	601	A1BHJ	C13-C14-N15	-2.12	106.78	110.70
2	B	601	A1BHJ	F09-C08-C07	-2.10	114.43	118.64
2	A	601	A1BHJ	F09-C08-C07	-2.09	114.46	118.64
2	A	601	A1BHJ	C22-C19-C21	-2.07	86.27	88.28
2	B	601	A1BHJ	C06-C05-C01	-2.02	122.50	126.92

There are no chirality outliers.

All (8) torsion outliers are listed below:

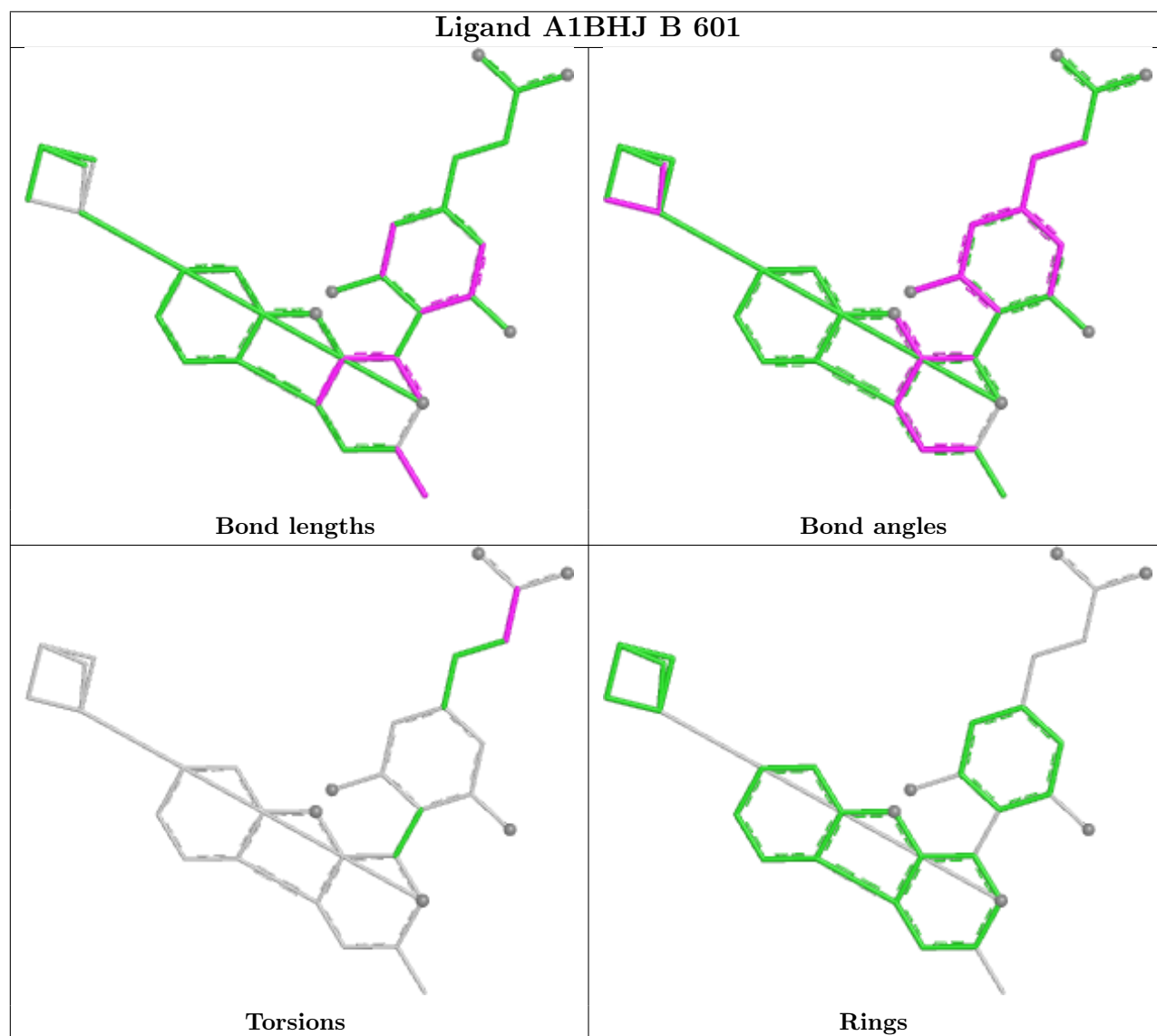
Mol	Chain	Res	Type	Atoms
2	D	601	A1BHJ	C05-C01-C02-O03
2	D	601	A1BHJ	C05-C01-C02-O04
2	B	601	A1BHJ	C05-C01-C02-O03
2	B	601	A1BHJ	C05-C01-C02-O04
2	C	601	A1BHJ	C05-C01-C02-O03
2	C	601	A1BHJ	C01-C05-C06-C10
2	C	601	A1BHJ	C05-C01-C02-O04
2	C	601	A1BHJ	C01-C05-C06-C07

There are no ring outliers.

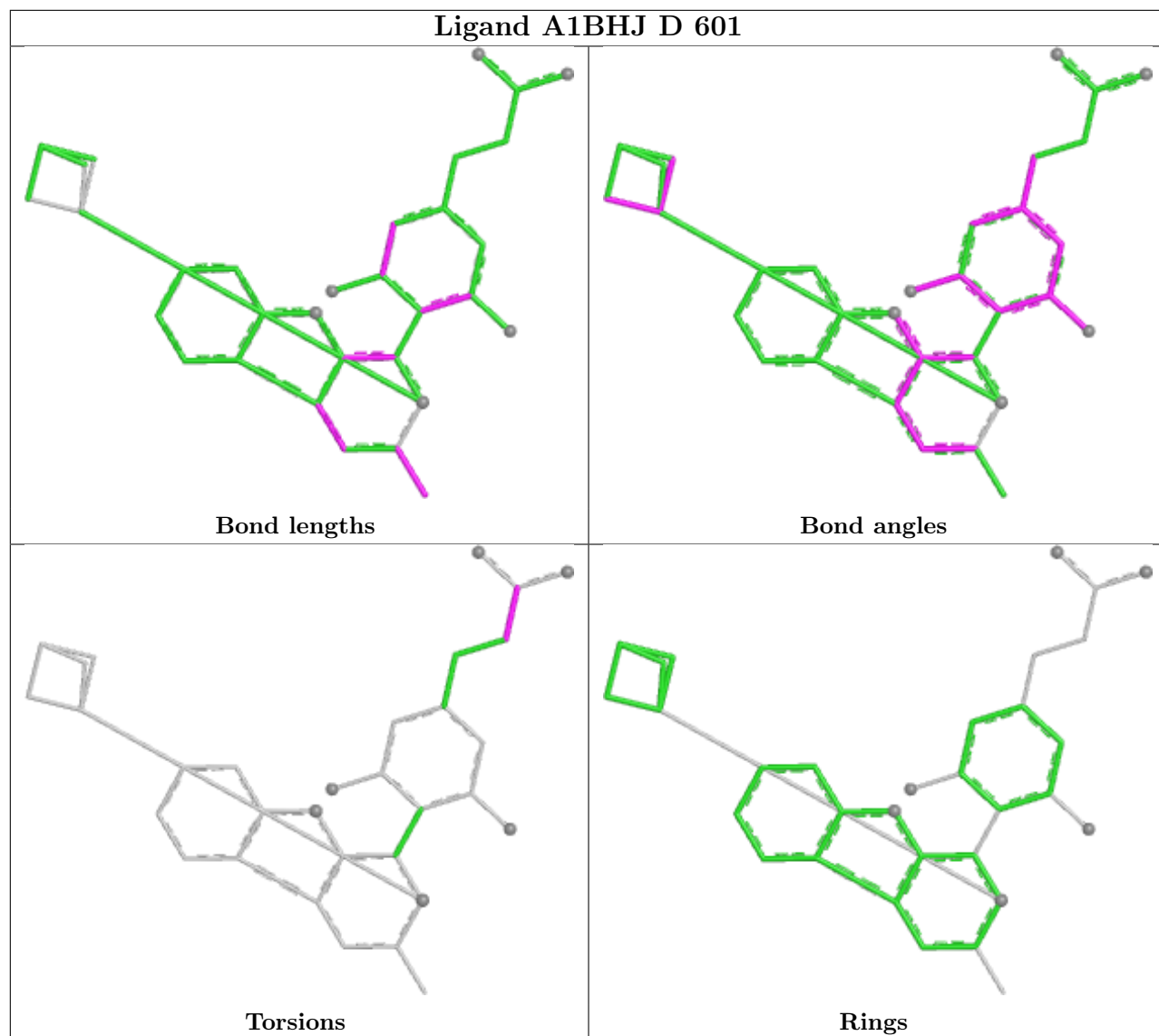
No monomer is involved in short contacts.

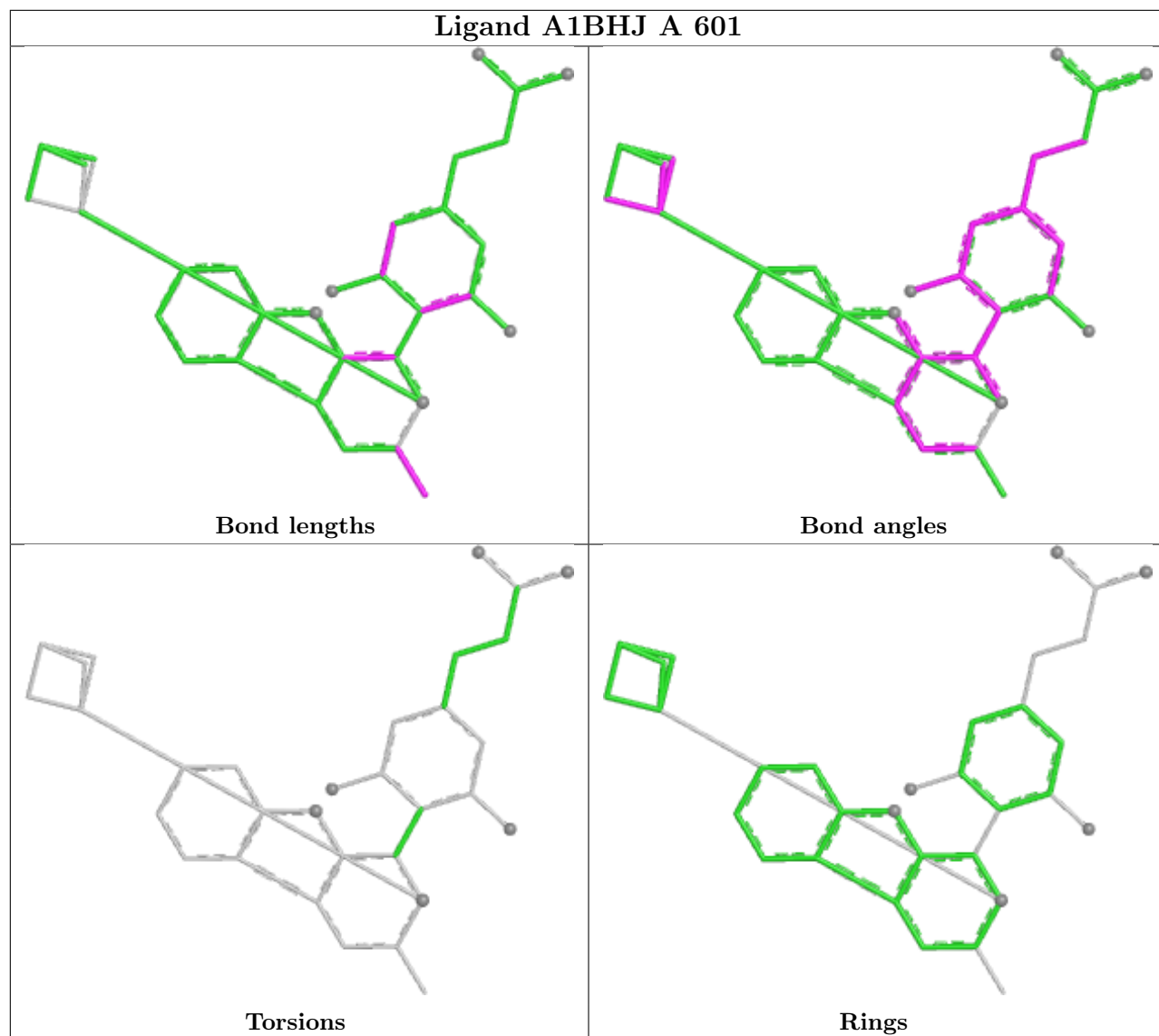
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

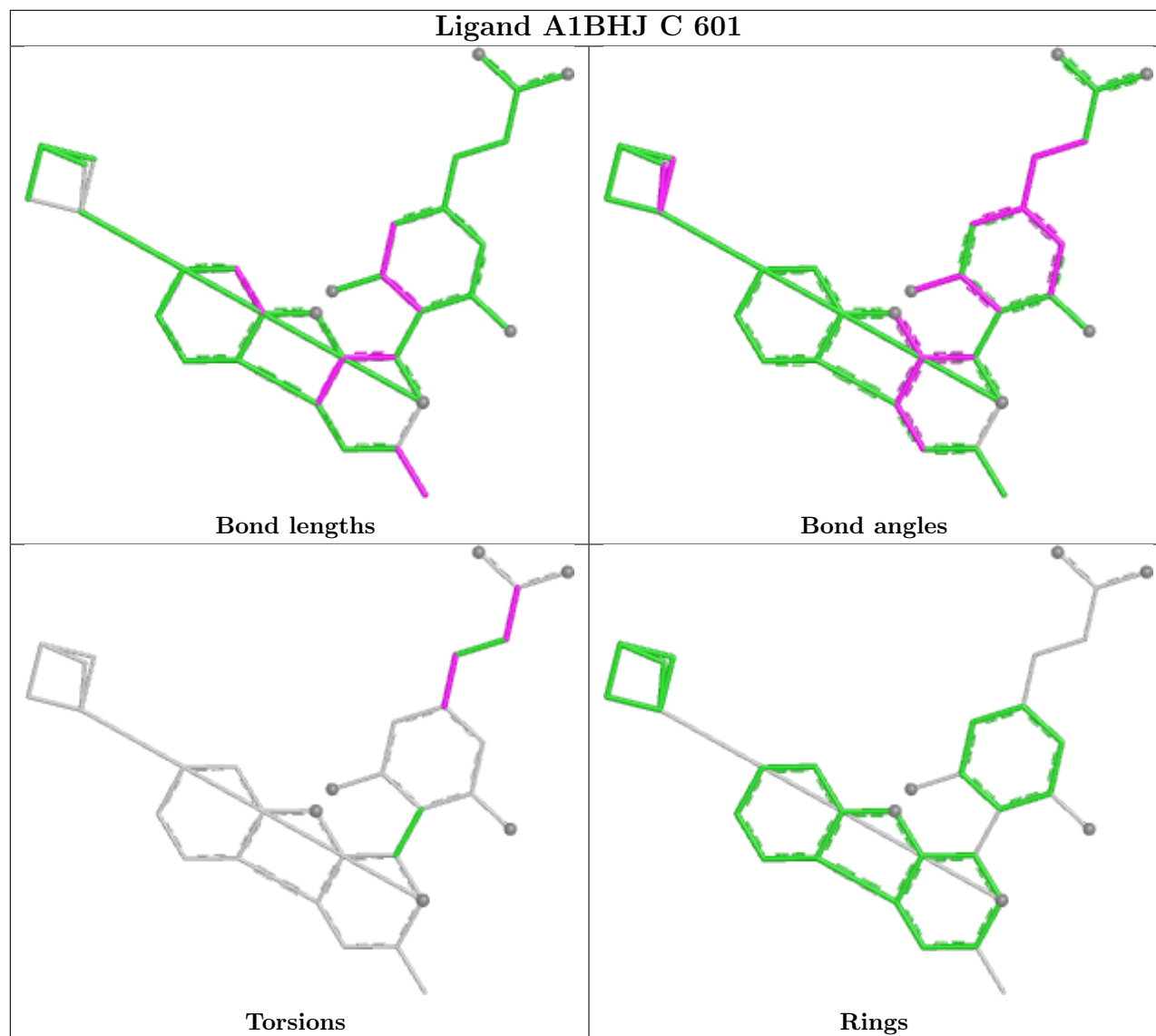
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand A1BHJ D 601







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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	232/259 (89%)	-0.31	1 (0%) 88 79	85, 104, 172, 202	0
1	B	220/259 (84%)	-0.37	1 (0%) 87 76	90, 118, 139, 168	1 (0%)
1	C	233/259 (89%)	-0.21	0 100 100	78, 108, 144, 175	3 (1%)
1	D	223/259 (86%)	-0.12	2 (0%) 81 64	97, 124, 163, 209	5 (2%)
All	All	908/1036 (87%)	-0.25	4 (0%) 88 79	78, 114, 156, 209	9 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	508	LEU	2.6
1	D	469	LEU	2.5
1	B	511	LEU	2.1
1	D	448	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](#)

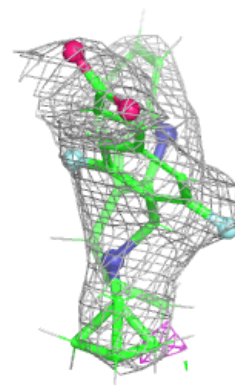
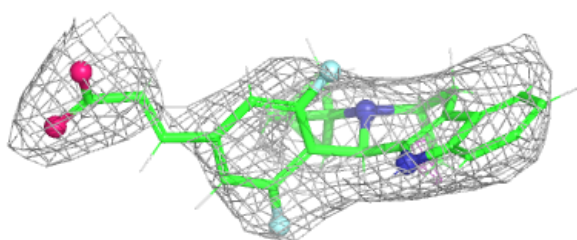
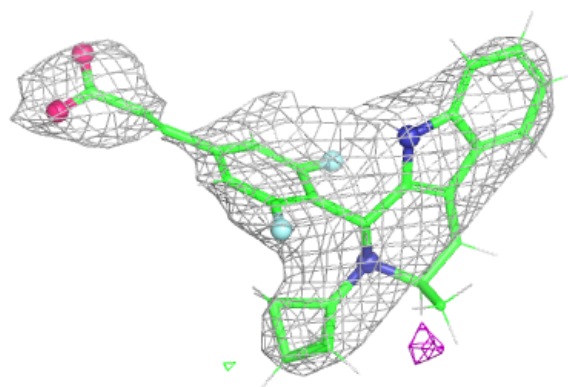
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	A1BHJ	B	601	32/32	0.89	0.11	74,95,118,122	0
2	A1BHJ	A	601	32/32	0.92	0.13	74,89,115,133	0
2	A1BHJ	C	601	32/32	0.92	0.10	74,87,105,118	0
2	A1BHJ	D	601	32/32	0.93	0.12	75,95,115,126	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.

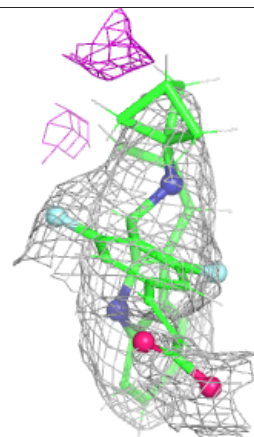
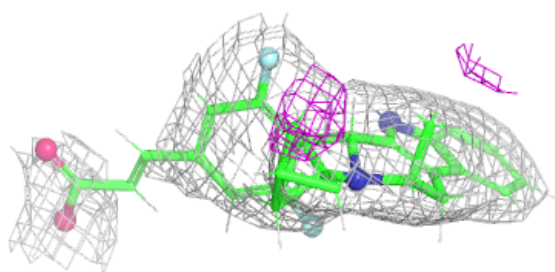
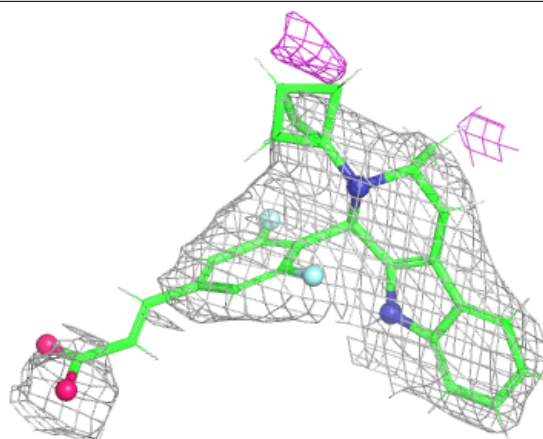
Electron density around A1BHJ B 601:

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



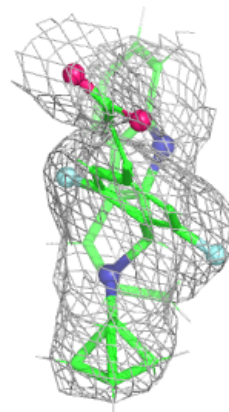
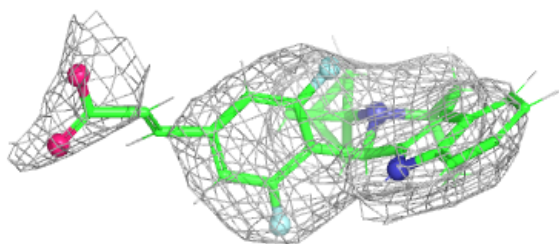
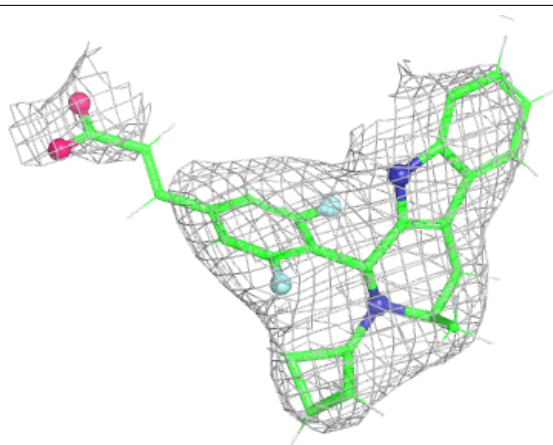
Electron density around A1BHJ A 601:

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



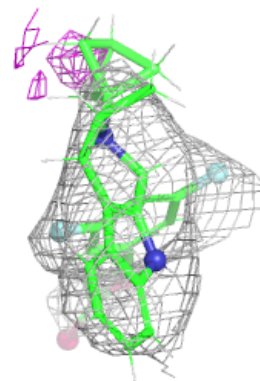
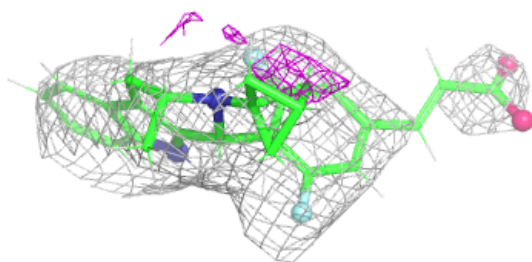
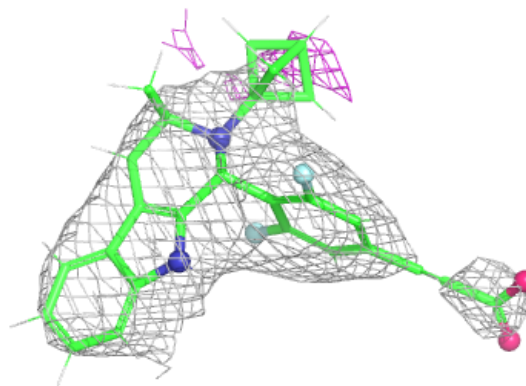
Electron density around A1BHJ C 601:

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



Electron density around A1BHJ D 601:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.