



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:59 AM UTC

PDB ID : 1VJB / pdb_00001vjb
Title : crystal structure of the ligand-binding domain of the estrogen-related receptor gamma in complex with 4-hydroxytamoxifen
Authors : Greschik, H.; Flaig, R.; Renaud, J.P.; Moras, D.
Deposited on : 2004-02-03
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

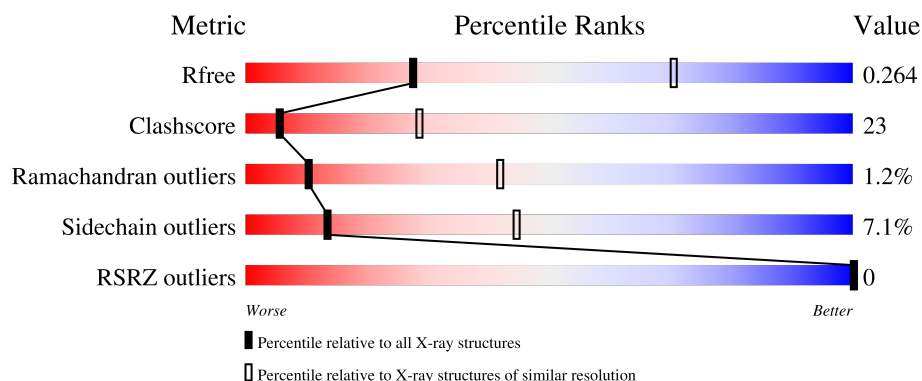
1 Overall quality at a glance

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	251	
1	B	251	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3246 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 Estrogen-related receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1594	1015	260	309	10			
1	B	210	Total	C	N	O	S	0	0	0
			1594	1008	261	315	10			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	208	MET	-	initiating methionine	UNP P62509
A	209	GLY	-	expression tag	UNP P62509
A	210	SER	-	expression tag	UNP P62509
A	211	SER	-	expression tag	UNP P62509
A	212	HIS	-	expression tag	UNP P62509
A	213	HIS	-	expression tag	UNP P62509
A	214	HIS	-	expression tag	UNP P62509
A	215	HIS	-	expression tag	UNP P62509
A	216	HIS	-	expression tag	UNP P62509
A	217	HIS	-	expression tag	UNP P62509
A	218	SER	-	expression tag	UNP P62509
A	219	SER	-	expression tag	UNP P62509
A	220	GLY	-	expression tag	UNP P62509
A	221	LEU	-	expression tag	UNP P62509
A	222	VAL	-	expression tag	UNP P62509
A	223	PRO	-	expression tag	UNP P62509
A	224	ARG	-	expression tag	UNP P62509
A	225	GLY	-	expression tag	UNP P62509
A	226	SER	-	expression tag	UNP P62509
A	227	HIS	-	expression tag	UNP P62509
A	228	MET	-	expression tag	UNP P62509
B	208	MET	-	initiating methionine	UNP P62509
B	209	GLY	-	expression tag	UNP P62509
B	210	SER	-	expression tag	UNP P62509
B	211	SER	-	expression tag	UNP P62509

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Chain	Residue	Modelled	Actual	Comment	Reference
B	212	HIS	-	expression tag	UNP P62509
B	213	HIS	-	expression tag	UNP P62509
B	214	HIS	-	expression tag	UNP P62509
B	215	HIS	-	expression tag	UNP P62509
B	216	HIS	-	expression tag	UNP P62509
B	217	HIS	-	expression tag	UNP P62509
B	218	SER	-	expression tag	UNP P62509
B	219	SER	-	expression tag	UNP P62509
B	220	GLY	-	expression tag	UNP P62509
B	221	LEU	-	expression tag	UNP P62509
B	222	VAL	-	expression tag	UNP P62509
B	223	PRO	-	expression tag	UNP P62509
B	224	ARG	-	expression tag	UNP P62509
B	225	GLY	-	expression tag	UNP P62509
B	226	SER	-	expression tag	UNP P62509
B	227	HIS	-	expression tag	UNP P62509
B	228	MET	-	expression tag	UNP P62509

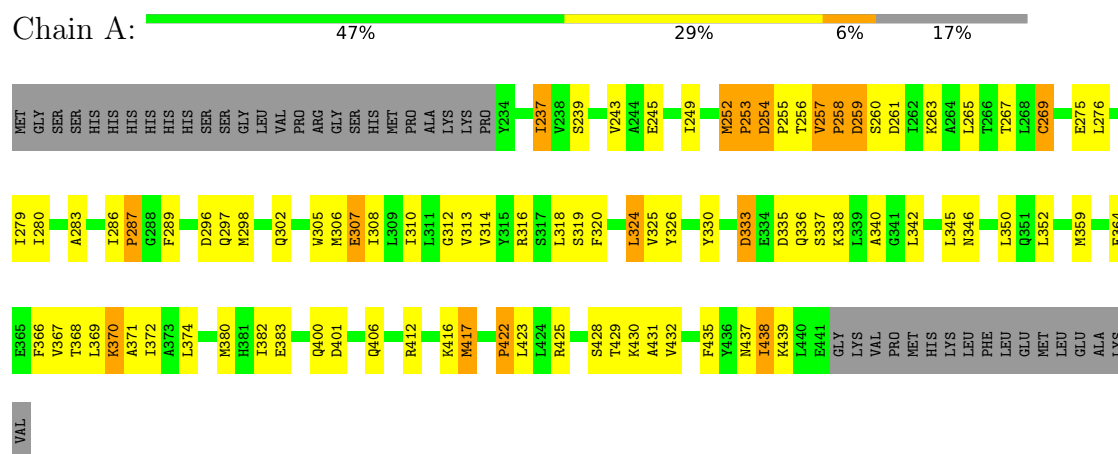
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- The chemical structure of Oxazepam (OHT) is shown with atom labels. The structure consists of a benzodiazepine core. The benzene ring is labeled with C1 through C7. The diazepine ring is labeled with C8 through C11. The nitrogen atom is labeled N24. The carbonyl group is labeled C12 and O13. The methylene group is labeled C14 and C15. The ethyl group is labeled C16 and C17. The methyl group is labeled C18. The hydroxyl group is labeled O19 and H20. The atom labels are: C1, C2, C3, C4, C5, C6, C7, C8, C9, C10, C11, C12, C13, C14, C15, C16, C17, C18, C19, C20, C21, C22, C23, C24, C25, C26, O13, O19, H20, N24.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 29	C 26	N 1	O 2	0	0
2	B	1	Total 29	C 26	N 1	O 2	0	0

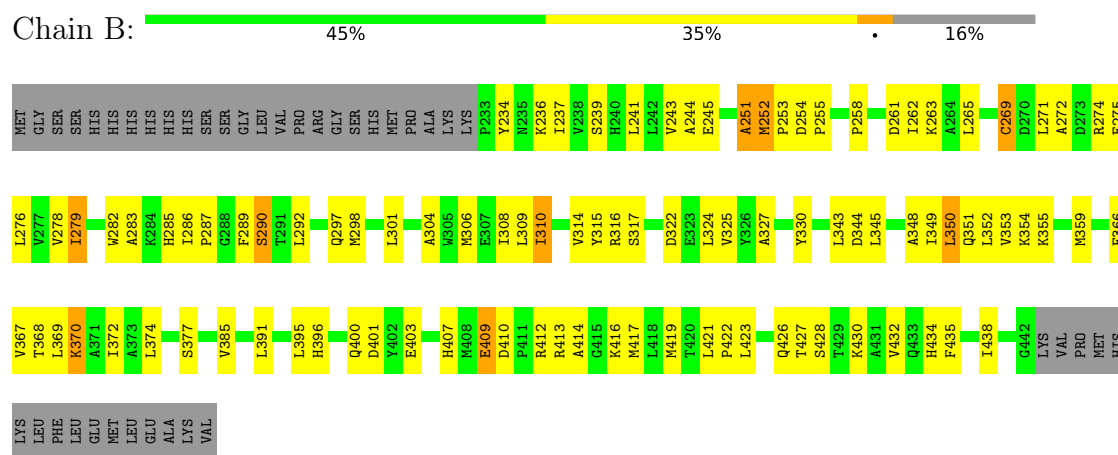
3 Residue-property plots

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-related receptor gamma



- Molecule 1: Estrogen-related receptor gamma



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	74.62Å 74.62Å 144.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.13 – 3.20 24.13 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (24.13-3.20) 99.0 (24.13-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.11 (at 3.23Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.256 , 0.275 0.237 , 0.264	Depositor DCC
R_{free} test set	372 reflections (4.61%)	wwPDB-VP
Wilson B-factor (Å ²)	91.7	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3246	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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: OHT

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	0.60	0/1620	1.08	8/2203 (0.4%)
1	B	0.55	0/1619	1.08	10/2199 (0.5%)
All	All	0.58	0/3239	1.08	18/4402 (0.4%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	257	VAL	N-CA-C	8.86	118.51	108.05
1	A	252	MET	CA-C-N	8.39	130.33	119.84
1	A	252	MET	C-N-CA	8.39	130.33	119.84
1	A	255	PRO	N-CA-CB	6.54	110.11	103.25
1	B	350	LEU	N-CA-C	-6.08	104.73	111.36
1	B	254	ASP	CA-C-N	5.98	125.66	119.56
1	B	254	ASP	C-N-CA	5.98	125.66	119.56
1	B	430	LYS	N-CA-C	-5.80	104.88	111.14
1	B	310	ILE	N-CA-C	5.66	115.86	110.42
1	A	417	MET	N-CA-C	-5.57	105.29	111.36
1	B	401	ASP	N-CA-C	5.53	116.99	111.07
1	B	324	LEU	N-CA-C	5.53	117.53	108.52
1	B	251	ALA	N-CA-C	5.22	117.39	111.02
1	B	385	VAL	N-CA-C	5.12	118.23	111.17
1	B	309	LEU	N-CA-C	-5.08	105.43	110.97
1	A	307	GLU	N-CA-C	-5.08	105.75	111.28
1	A	245	GLU	CA-C-N	5.03	126.13	119.84
1	A	245	GLU	C-N-CA	5.03	126.13	119.84

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	1594	0	1554	72	0
1	B	1594	0	1531	85	0
2	A	29	0	28	3	0
2	B	29	0	28	5	0
All	All	3246	0	3141	150	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 23.

All (150) 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:314:VAL:HG21	1:A:352:LEU:HD23	1.47	0.95
1:A:307:GLU:OE2	1:A:425:ARG:HA	1.78	0.83
1:B:351:GLN:HA	1:B:354:LYS:HE3	1.60	0.82
1:B:304:ALA:HB2	1:B:377:SER:HB3	1.61	0.80
1:B:271:LEU:O	1:B:275:GLU:HG3	1.84	0.77
1:B:352:LEU:HD13	1:B:423:LEU:HG	1.65	0.77
1:A:412:ARG:HG2	1:B:400:GLN:OE1	1.88	0.74
1:A:237:ILE:HG23	1:A:287:PRO:HG2	1.70	0.74
1:B:304:ALA:O	1:B:308:ILE:HG12	1.88	0.72
1:B:269:CYS:HA	2:B:200:OHT:C19	2.21	0.71
1:B:368:THR:O	1:B:372:ILE:HG12	1.92	0.70
1:A:279:ILE:HD12	1:A:308:ILE:HG21	1.72	0.69
1:A:257:VAL:O	1:A:263:LYS:HD2	1.92	0.69
1:A:261:ASP:O	1:A:265:LEU:HG	1.92	0.69
1:B:327:ALA:HB3	1:B:330:TYR:CB	2.24	0.68
1:B:269:CYS:HA	2:B:200:OHT:H19	1.77	0.67
1:A:259:ASP:HA	1:A:263:LYS:HD3	1.75	0.67
1:A:276:LEU:O	1:A:280:ILE:HG12	1.97	0.64
1:A:312:GLY:HA3	1:A:316:ARG:HH21	1.64	0.63
1:B:310:ILE:CD1	1:B:428:SER:HA	2.28	0.63
1:B:325:VAL:HG12	1:B:325:VAL:O	1.98	0.63
1:B:327:ALA:HB3	1:B:330:TYR:HB3	1.81	0.63
1:B:349:ILE:O	1:B:353:VAL:HG23	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:ILE:HD12	1:A:308:ILE:CG2	2.30	0.62
1:A:412:ARG:HG2	1:B:400:GLN:CD	2.25	0.62
1:A:366:PHE:CZ	1:A:370:LYS:HE2	2.35	0.62
1:A:298:MET:O	1:A:302:GLN:HG3	2.01	0.61
1:A:275:GLU:O	1:A:279:ILE:HG12	2.01	0.61
1:A:253:PRO:HG3	1:A:330:TYR:HD1	1.66	0.60
1:B:276:LEU:HA	1:B:279:ILE:HG22	1.83	0.60
1:A:412:ARG:HB3	1:B:396:HIS:CE1	2.37	0.60
1:B:310:ILE:O	1:B:314:VAL:HG23	2.02	0.59
1:B:345:LEU:O	1:B:348:ALA:HB3	2.01	0.59
1:B:372:ILE:HD13	1:B:395:LEU:HB3	1.85	0.59
1:A:324:LEU:HD13	1:A:350:LEU:HG	1.84	0.58
1:A:438:ILE:HG22	1:A:439:LYS:N	2.19	0.58
1:B:274:ARG:O	1:B:278:VAL:HG23	2.03	0.58
1:B:403:GLU:OE1	1:B:413:ARG:HD3	2.04	0.57
1:A:276:LEU:HD21	1:A:305:TRP:HB2	1.86	0.57
1:A:253:PRO:O	1:A:254:ASP:C	2.48	0.57
1:B:276:LEU:HA	1:B:279:ILE:CG2	2.35	0.56
1:B:290:SER:HA	1:B:297:GLN:NE2	2.21	0.56
1:A:412:ARG:HG2	1:B:400:GLN:NE2	2.21	0.56
1:B:251:ALA:HA	1:B:271:LEU:HD13	1.88	0.55
1:A:333:ASP:OD1	1:A:336:GLN:HG3	2.06	0.55
1:A:257:VAL:HG12	1:A:258:PRO:O	2.07	0.55
1:B:310:ILE:HD12	1:B:428:SER:HA	1.87	0.55
1:A:352:LEU:HD13	1:A:423:LEU:HG	1.88	0.54
1:B:435:PHE:HB2	2:B:200:OHT:H15	1.90	0.53
1:A:435:PHE:O	1:A:438:ILE:HG12	2.08	0.53
1:A:269:CYS:HA	2:A:100:OHT:H19	1.90	0.53
1:B:261:ASP:O	1:B:265:LEU:HG	2.07	0.53
1:B:316:ARG:HB3	1:B:325:VAL:CG1	2.39	0.53
1:B:407:HIS:HA	1:B:409:GLU:OE2	2.09	0.52
1:B:269:CYS:SG	2:B:200:OHT:H19	2.50	0.52
1:A:337:SER:O	1:A:340:ALA:N	2.43	0.51
1:B:297:GLN:O	1:B:301:LEU:HG	2.11	0.51
1:A:252:MET:O	1:A:254:ASP:N	2.43	0.51
1:A:286:ILE:HD12	1:A:374:LEU:CD2	2.41	0.51
1:A:425:ARG:HD2	1:B:426:GLN:OE1	2.11	0.51
1:B:253:PRO:O	1:B:255:PRO:HD3	2.10	0.50
1:B:434:HIS:O	1:B:438:ILE:HG13	2.12	0.50
1:B:327:ALA:HB3	1:B:330:TYR:HB2	1.93	0.50
1:B:350:LEU:O	1:B:354:LYS:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:GLU:O	1:B:279:ILE:HG22	2.12	0.50
1:B:241:LEU:CB	1:B:367:VAL:CG1	2.91	0.49
1:B:241:LEU:CB	1:B:367:VAL:HG11	2.42	0.49
1:B:314:VAL:HG21	1:B:352:LEU:HD23	1.93	0.49
1:A:306:MET:HG2	1:A:432:VAL:HG23	1.94	0.49
1:B:352:LEU:HD13	1:B:423:LEU:CG	2.38	0.49
1:A:412:ARG:CG	1:B:400:GLN:HE22	2.25	0.49
1:B:239:SER:O	1:B:243:VAL:HG23	2.13	0.48
1:B:419:MET:O	1:B:422:PRO:HD2	2.13	0.48
1:A:342:LEU:HB3	1:A:346:ASN:HD22	1.78	0.48
1:A:369:LEU:O	1:A:370:LYS:C	2.56	0.48
1:A:422:PRO:O	1:A:425:ARG:HB3	2.12	0.48
1:B:410:ASP:HB3	1:B:413:ARG:HB2	1.94	0.48
1:A:267:THR:HG21	1:A:330:TYR:CD1	2.49	0.48
1:A:324:LEU:HD22	1:A:346:ASN:OD1	2.14	0.47
1:B:306:MET:O	1:B:310:ILE:HG13	2.13	0.47
1:A:249:ILE:HD12	1:A:275:GLU:HG2	1.95	0.47
1:A:324:LEU:CD2	1:A:346:ASN:OD1	2.62	0.47
1:B:252:MET:N	1:B:253:PRO:HD3	2.30	0.47
1:A:306:MET:HG3	1:A:428:SER:HB2	1.97	0.47
1:B:292:LEU:HD11	1:B:391:LEU:HD22	1.96	0.47
1:A:239:SER:O	1:A:243:VAL:HG23	2.14	0.47
1:B:298:MET:N	1:B:298:MET:HE2	2.30	0.47
1:A:412:ARG:O	1:A:416:LYS:HG3	2.15	0.47
1:B:265:LEU:HD22	2:B:200:OHT:H13	1.97	0.47
1:B:413:ARG:O	1:B:416:LYS:HB2	2.14	0.47
1:A:283:ALA:HB1	1:A:289:PHE:CE2	2.50	0.47
1:A:286:ILE:HD12	1:A:374:LEU:HD23	1.97	0.47
1:B:276:LEU:CA	1:B:279:ILE:HG22	2.45	0.46
1:A:260:SER:O	1:A:261:ASP:C	2.58	0.46
1:A:359:MET:SD	1:A:417:MET:HE2	2.55	0.46
1:A:431:ALA:O	2:A:100:OHT:H15	2.16	0.46
1:B:349:ILE:HG12	1:B:427:THR:HG21	1.97	0.46
1:A:254:ASP:O	1:A:256:THR:N	2.44	0.45
1:B:258:PRO:O	1:B:263:LYS:HD2	2.16	0.45
1:A:252:MET:N	1:A:253:PRO:HD3	2.31	0.45
1:B:276:LEU:O	1:B:279:ILE:HG23	2.17	0.45
1:A:269:CYS:HA	2:A:100:OHT:C19	2.47	0.45
1:A:371:ALA:O	1:A:374:LEU:HB3	2.16	0.45
1:B:283:ALA:O	1:B:289:PHE:CD2	2.69	0.45
1:B:297:GLN:C	1:B:298:MET:HE2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:SER:HA	1:B:297:GLN:HE22	1.82	0.45
1:B:410:ASP:C	1:B:412:ARG:H	2.24	0.45
1:B:244:ALA:O	1:B:282:TRP:HD1	2.00	0.45
1:A:296:ASP:O	1:A:297:GLN:C	2.60	0.44
1:B:369:LEU:HD21	1:B:417:MET:HB3	1.99	0.44
1:B:241:LEU:C	1:B:367:VAL:HG11	2.42	0.44
1:B:285:HIS:O	1:B:287:PRO:HD3	2.17	0.44
1:B:286:ILE:O	1:B:287:PRO:C	2.61	0.44
1:B:304:ALA:HB3	1:B:374:LEU:HD12	1.99	0.44
1:A:313:VAL:HG21	1:A:326:TYR:CD2	2.53	0.44
1:A:412:ARG:HG2	1:B:400:GLN:HE22	1.81	0.44
1:A:256:THR:HG22	1:A:257:VAL:HG23	2.00	0.43
1:B:236:LYS:HG3	1:B:237:ILE:H	1.82	0.43
1:B:301:LEU:HD23	1:B:374:LEU:HD11	2.00	0.43
1:A:366:PHE:CE2	1:A:370:LYS:HE2	2.54	0.43
1:A:306:MET:O	1:A:310:ILE:HG12	2.19	0.43
1:B:396:HIS:CD2	1:B:396:HIS:C	2.97	0.43
1:B:343:LEU:O	1:B:344:ASP:C	2.60	0.43
1:B:413:ARG:O	1:B:414:ALA:C	2.62	0.43
1:B:261:ASP:OD1	1:B:261:ASP:N	2.50	0.42
1:A:318:LEU:C	1:A:320:PHE:H	2.27	0.42
1:A:342:LEU:CB	1:A:346:ASN:HD22	2.32	0.42
1:B:372:ILE:CD1	1:B:395:LEU:HB3	2.50	0.42
1:A:337:SER:O	1:A:338:LYS:C	2.63	0.42
1:B:366:PHE:O	1:B:370:LYS:HB2	2.20	0.42
1:A:364:GLU:O	1:A:367:VAL:HG22	2.19	0.42
1:A:382:ILE:HG22	1:A:383:GLU:N	2.34	0.42
1:B:414:ALA:O	1:B:417:MET:HB2	2.19	0.42
1:A:313:VAL:HG13	1:A:325:VAL:O	2.20	0.42
1:B:243:VAL:C	1:B:245:GLU:H	2.28	0.42
1:A:335:ASP:O	1:A:336:GLN:C	2.61	0.41
1:B:316:ARG:HD2	1:B:325:VAL:HG12	2.01	0.41
1:B:421:LEU:N	1:B:422:PRO:CD	2.83	0.41
1:A:342:LEU:HB3	1:A:346:ASN:ND2	2.35	0.41
1:B:315:TYR:CD1	1:B:366:PHE:HB2	2.55	0.41
1:A:286:ILE:CG2	1:A:289:PHE:HB2	2.51	0.41
1:B:310:ILE:HD12	1:B:428:SER:CA	2.51	0.41
1:A:302:GLN:O	1:A:305:TRP:CD1	2.74	0.41
1:B:272:ALA:O	1:B:276:LEU:HG	2.21	0.41
1:A:345:LEU:HD11	1:A:431:ALA:HA	2.03	0.41
1:A:368:THR:O	1:A:372:ILE:HG13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:421:LEU:HD23	1:B:421:LEU:HA	1.91	0.41
1:A:429:THR:O	1:A:430:LYS:C	2.64	0.41
1:B:355:LYS:O	1:B:359:MET:HG3	2.21	0.41
1:A:267:THR:HG21	1:A:330:TYR:CE1	2.56	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	206/251 (82%)	178 (86%)	24 (12%)	4 (2%)	6	33
1	B	208/251 (83%)	194 (93%)	13 (6%)	1 (0%)	24	59
All	All	414/502 (82%)	372 (90%)	37 (9%)	5 (1%)	10	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	438	ILE
1	B	234	TYR
1	A	319	SER
1	A	253	PRO
1	A	254	ASP

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	169/221 (76%)	155 (92%)	14 (8%)	10	38
1	B	167/221 (76%)	157 (94%)	10 (6%)	17	50
All	All	336/442 (76%)	312 (93%)	24 (7%)	13	44

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	237	ILE
1	A	258	PRO
1	A	259	ASP
1	A	269	CYS
1	A	287	PRO
1	A	324	LEU
1	A	333	ASP
1	A	370	LYS
1	A	380	MET
1	A	400	GLN
1	A	401	ASP
1	A	406	GLN
1	A	422	PRO
1	A	437	ASN
1	B	252	MET
1	B	262	ILE
1	B	269	CYS
1	B	279	ILE
1	B	290	SER
1	B	317	SER
1	B	322	ASP
1	B	370	LYS
1	B	409	GLU
1	B	432	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	400	GLN
1	A	437	ASN
1	B	297	GLN
1	B	346	ASN
1	B	396	HIS
1	B	407	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 ⓘ

2 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	OHT	B	200	-	31,31,31	1.94	12 (38%)	41,41,41	1.03	3 (7%)
2	OHT	A	100	-	31,31,31	1.73	10 (32%)	41,41,41	0.93	2 (4%)

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	OHT	B	200	-	-	2/24/24/24	0/3/3/3
2	OHT	A	100	-	-	2/24/24/24	0/3/3/3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	100	OHT	C2-C1	4.50	1.46	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	200	OHT	C7-C8	4.14	1.42	1.35
2	B	200	OHT	C2-C1	4.12	1.45	1.39
2	B	200	OHT	C1-C7	3.47	1.55	1.49
2	A	100	OHT	C16-C11	3.12	1.44	1.39
2	B	200	OHT	C16-C11	3.11	1.44	1.39
2	B	200	OHT	C11-C8	3.09	1.54	1.49
2	A	100	OHT	C7-C8	2.86	1.40	1.35
2	B	200	OHT	C21-C20	2.68	1.43	1.38
2	A	100	OHT	C1-C7	2.67	1.54	1.49
2	B	200	OHT	C22-C17	2.54	1.43	1.39
2	A	100	OHT	C5-C4	2.32	1.43	1.39
2	B	200	OHT	C13-C12	2.26	1.42	1.38
2	A	100	OHT	C6-C1	2.21	1.42	1.39
2	A	100	OHT	C21-C20	2.20	1.42	1.38
2	B	200	OHT	C3-C2	2.17	1.42	1.38
2	B	200	OHT	C15-C16	2.12	1.42	1.38
2	A	100	OHT	C13-C12	2.09	1.42	1.38
2	A	100	OHT	C3-C4	2.09	1.42	1.39
2	A	100	OHT	C22-C17	2.08	1.42	1.39
2	B	200	OHT	C12-C11	2.08	1.42	1.39
2	B	200	OHT	C6-C1	2.05	1.42	1.39

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	200	OHT	C11-C8-C7	2.79	125.10	122.35
2	A	100	OHT	C11-C8-C7	2.69	124.99	122.35
2	B	200	OHT	C1-C7-C8	2.60	125.94	122.83
2	A	100	OHT	C17-C7-C1	-2.10	111.14	115.34
2	B	200	OHT	C17-C7-C1	-2.06	111.22	115.34

There are no chirality outliers.

All (4) torsion outliers are listed below:

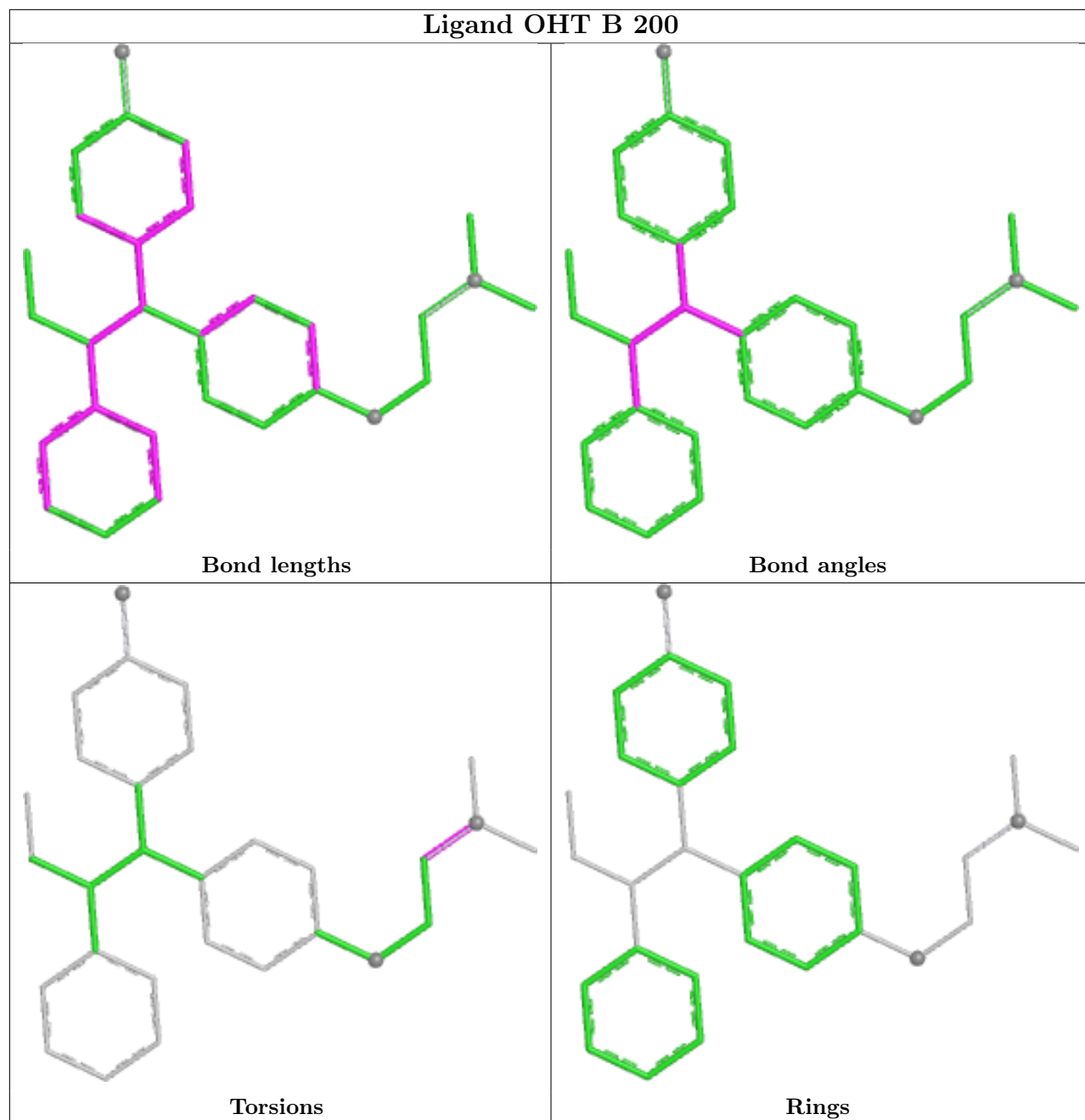
Mol	Chain	Res	Type	Atoms
2	A	100	OHT	C23-C24-N24-C25
2	A	100	OHT	C23-C24-N24-C26
2	B	200	OHT	C23-C24-N24-C25
2	B	200	OHT	C23-C24-N24-C26

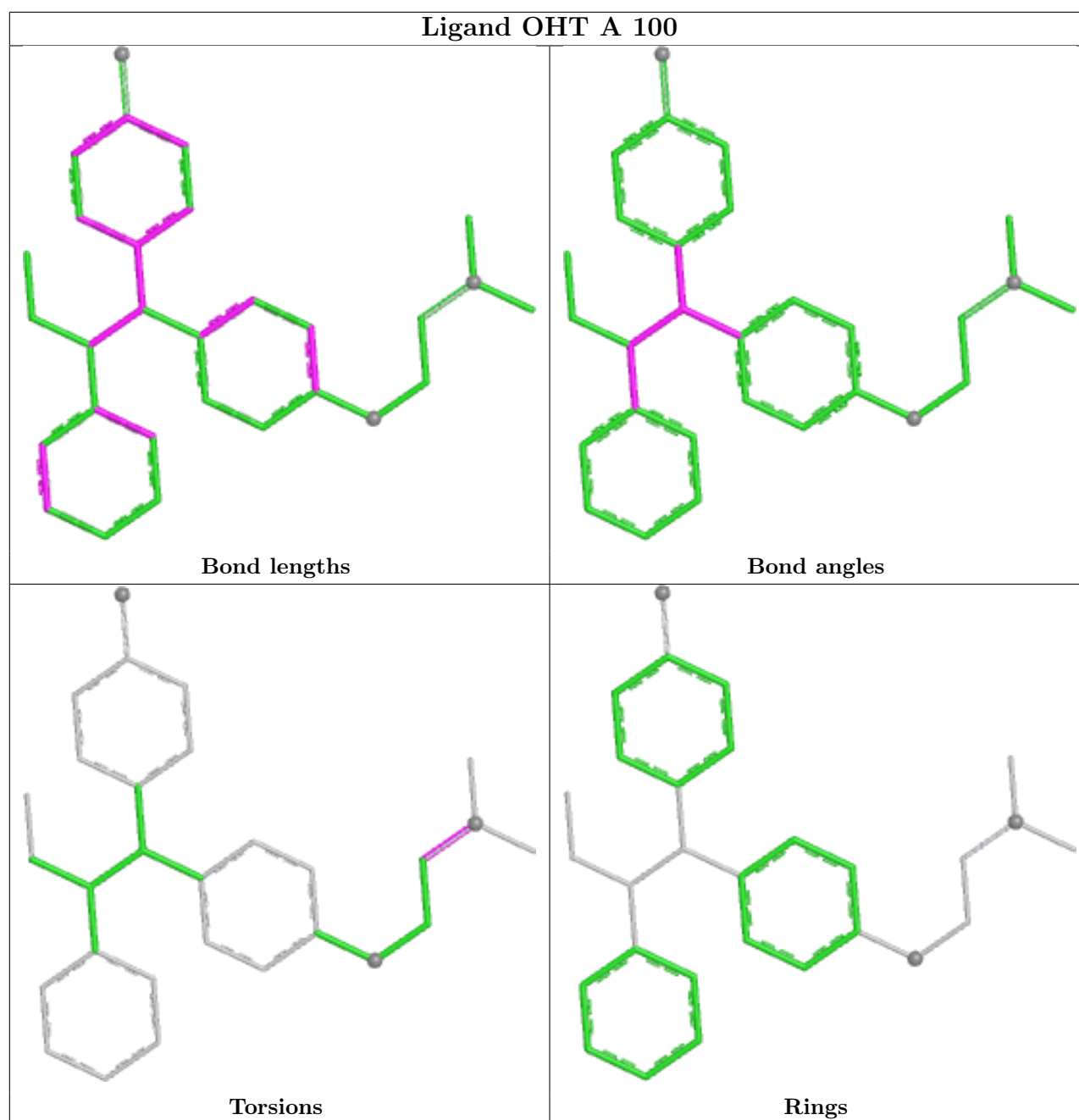
There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	200	OHT	5	0
2	A	100	OHT	3	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 [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	208/251 (82%)	-0.70	0 100 100	40, 64, 96, 100	0
1	B	210/251 (83%)	-0.51	0 100 100	43, 76, 100, 100	0
All	All	418/502 (83%)	-0.60	0 100 100	40, 70, 100, 100	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

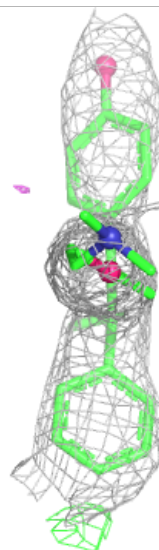
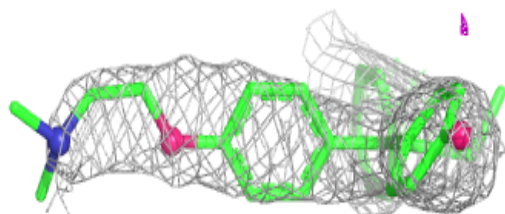
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	OHT	B	200	29/29	0.90	0.11	60,70,94,96	0
2	OHT	A	100	29/29	0.93	0.12	60,64,67,68	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.

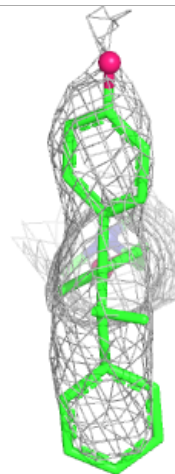
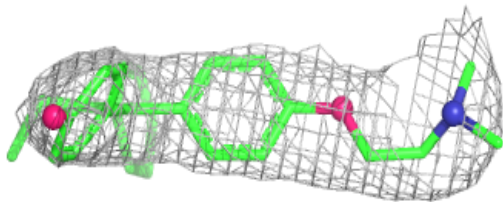
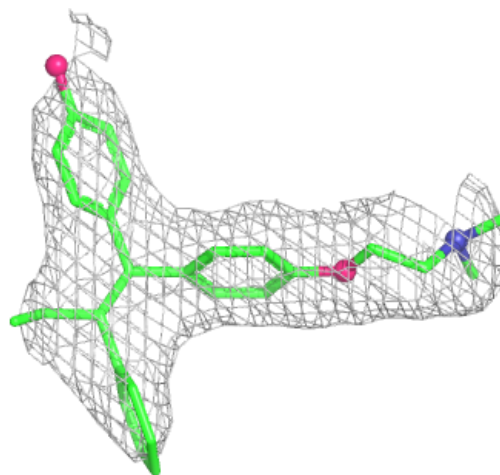
Electron density around OHT B 200:

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



Electron density around OHT A 100:

$2mF_o-DF_c$ (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.