



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 04:53 AM EDT

PDB ID : 2B6A / pdb_00002b6a
Title : Crystal structure of HIV-1 reverse transcriptase (RT) in complex with THR-50
Authors : Morningstar, M.L.; Roth, T.; Smith, M.K.; Zajac, M.; Watson, K.; Buckheit, R.W.; Das, K.; Zhang, W.; Arnold, E.; Michejda, C.J.
Deposited on : 2005-09-30
Resolution : 2.65 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

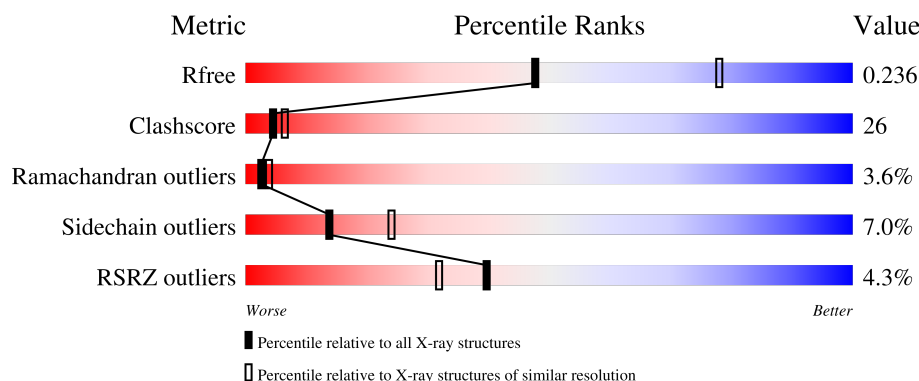
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	560	3% 51% 41% 7% .
2	B	430	6% 52% 37% 9% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7921 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 Reverse transcriptase p66 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	558	Total	C	N	O	S	0	0	0
			4451	2882	740	822	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	280	SER	CYS	engineered mutation	UNP P03366

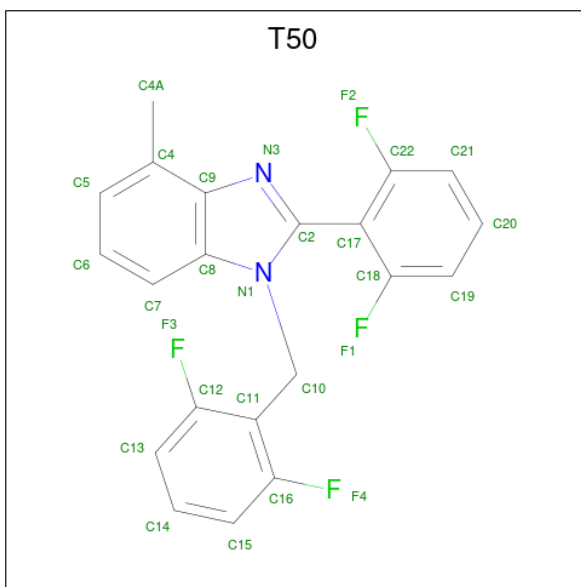
- Molecule 2 is a protein called Reverse transcriptase p51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	427	Total	C	N	O	S	0	0	0
			3443	2241	567	630	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	engineered mutation	UNP P03366

- Molecule 3 is 1-(2,6-DIFLUOROBENZYL)-2-(2,6-DIFLUOROPHENYL)-4-METHYL-1H-BENZIMIDAZOLE (CCD ID: T50) (formula: C₂₁H₁₄F₄N₂).

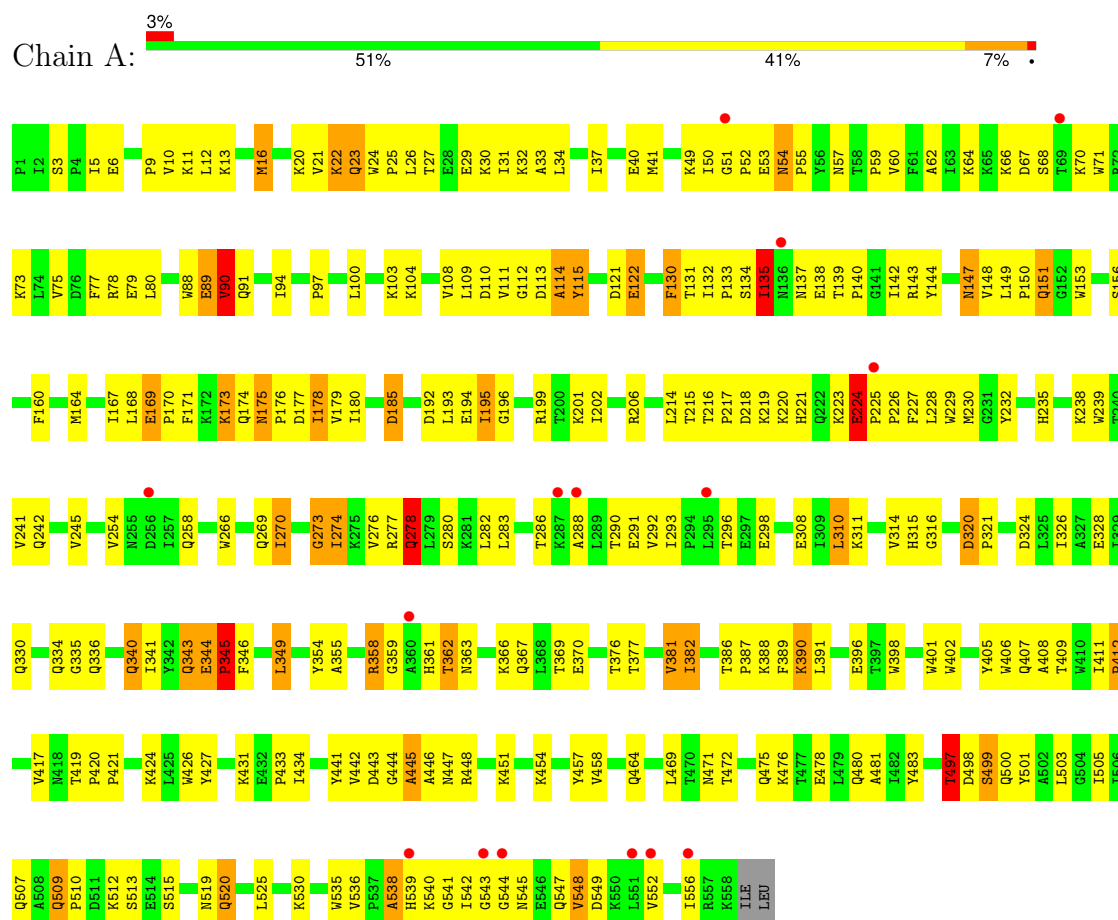


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	F	N	0	0
			27	21	4	2		

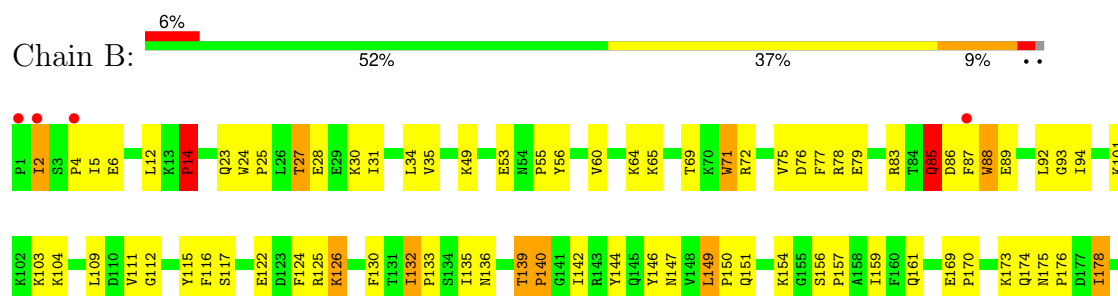
3 Residue-property plots

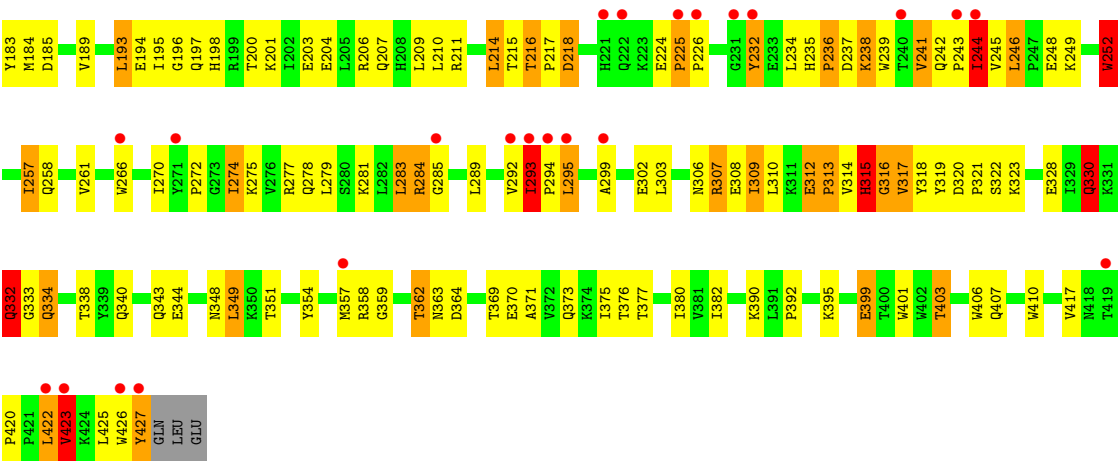
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Reverse transcriptase p66 subunit



• Molecule 2: Reverse transcriptase p51 subunit





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	227.71Å 70.17Å 106.63Å 90.00° 105.82° 90.00°	Depositor
Resolution (Å)	20.00 – 2.65 20.00 – 2.65	Depositor EDS
% Data completeness (in resolution range)	94.6 (20.00-2.65) 94.4 (20.00-2.65)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.226 , 0.277 0.227 , 0.236	Depositor DCC
R_{free} test set	2520 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	66.2	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 75.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7921	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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: T50

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	0.68	1/4566 (0.0%)	1.17	42/6209 (0.7%)
2	B	0.78	0/3541	1.31	45/4822 (0.9%)
All	All	0.73	1/8107 (0.0%)	1.23	87/11031 (0.8%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	361	HIS	N-CA	5.95	1.54	1.46

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	89	GLU	N-CA-C	-11.70	99.22	114.31
2	B	226	PRO	N-CA-CB	11.38	110.86	103.23
1	A	344	GLU	CA-C-N	11.09	133.70	119.84
1	A	344	GLU	C-N-CA	11.09	133.70	119.84
2	B	334	GLN	OE1-CD-NE2	-10.03	112.57	122.60
2	B	332	GLN	OE1-CD-NE2	-9.90	112.70	122.60
2	B	88	TRP	N-CA-C	-9.71	99.31	112.03
2	B	204	GLU	N-CA-C	-9.56	99.73	111.40
2	B	330	GLN	OE1-CD-NE2	-9.47	113.13	122.60
2	B	85	GLN	OE1-CD-NE2	-9.27	113.33	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	GLU	N-CA-C	9.23	125.30	108.69
2	B	151	GLN	N-CA-C	9.16	122.96	111.69
1	A	382	ILE	N-CA-C	8.78	118.78	110.53
1	A	320	ASP	CA-C-N	8.60	128.33	119.56
1	A	320	ASP	C-N-CA	8.60	128.33	119.56
2	B	422	LEU	N-CA-C	-8.57	102.13	112.59
1	A	499	SER	N-CA-C	8.26	121.48	108.67
2	B	244	ILE	N-CA-C	8.05	114.83	106.21
1	A	513	SER	N-CA-C	7.88	120.50	110.33
2	B	423	VAL	N-CA-C	-7.48	105.06	112.17
1	A	273	GLY	N-CA-C	7.08	124.66	114.67
1	A	349	LEU	N-CA-C	-7.08	102.99	111.69
2	B	225	PRO	N-CA-CB	7.06	109.92	103.08
1	A	169	GLU	CA-C-N	-6.99	112.97	119.82
1	A	169	GLU	C-N-CA	-6.99	112.97	119.82
1	A	115	TYR	N-CA-C	6.92	119.71	111.33
1	A	225	PRO	N-CA-C	-6.92	102.26	110.70
2	B	139	THR	CA-C-N	6.88	127.17	119.32
2	B	139	THR	C-N-CA	6.88	127.17	119.32
1	A	509	GLN	CA-C-N	6.87	127.81	120.11
1	A	509	GLN	C-N-CA	6.87	127.81	120.11
2	B	382	ILE	N-CA-C	6.87	116.99	110.53
2	B	332	GLN	CG-CD-NE2	6.85	126.67	116.40
2	B	122	GLU	N-CA-C	6.72	119.44	111.71
2	B	71	TRP	N-CA-C	6.69	119.45	110.35
2	B	315	HIS	N-CA-C	-6.68	103.61	111.03
1	A	361	HIS	N-CA-C	6.63	119.67	107.99
1	A	151	GLN	N-CA-C	6.60	120.79	112.87
2	B	334	GLN	CG-CD-NE2	6.59	126.28	116.40
2	B	85	GLN	CG-CD-NE2	6.55	126.23	116.40
2	B	241	VAL	N-CA-C	6.46	115.64	106.53
2	B	330	GLN	CG-CD-NE2	6.41	126.01	116.40
1	A	539	HIS	N-CA-C	6.39	120.08	112.93
2	B	149	LEU	CA-C-N	6.32	126.50	120.31
2	B	149	LEU	C-N-CA	6.32	126.50	120.31
1	A	70	LYS	N-CA-C	6.31	116.61	108.34
1	A	149	LEU	CA-C-N	6.16	126.11	119.76
1	A	149	LEU	C-N-CA	6.16	126.11	119.76
2	B	246	LEU	N-CA-C	-6.02	101.06	110.14
2	B	401	TRP	N-CA-C	5.96	121.28	113.30
1	A	388	LYS	N-CA-C	-5.95	100.00	109.76
1	A	173	LYS	N-CA-C	-5.93	105.12	112.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	381	VAL	CB-CA-C	-5.93	104.25	112.02
1	A	390	LYS	N-CA-C	-5.88	98.83	108.41
1	A	175	ASN	CA-C-N	5.87	126.28	119.47
1	A	175	ASN	C-N-CA	5.87	126.28	119.47
2	B	349	LEU	N-CA-C	-5.76	106.22	113.18
2	B	216	THR	CA-C-N	5.68	125.61	119.76
2	B	216	THR	C-N-CA	5.68	125.61	119.76
1	A	278	GLN	N-CA-C	5.65	119.35	112.23
1	A	536	VAL	CA-C-N	5.58	126.81	119.84
1	A	536	VAL	C-N-CA	5.58	126.81	119.84
1	A	54	ASN	CA-C-N	5.58	126.81	119.84
1	A	54	ASN	C-N-CA	5.58	126.81	119.84
2	B	112	GLY	N-CA-C	5.51	119.34	112.73
2	B	189	VAL	N-CA-C	5.49	115.69	107.51
1	A	180	ILE	N-CA-C	5.41	115.92	108.12
2	B	293	ILE	CA-C-N	5.38	126.56	119.84
2	B	293	ILE	C-N-CA	5.38	126.56	119.84
2	B	348	ASN	N-CA-C	5.35	117.46	109.59
2	B	218	ASP	N-CA-C	5.27	118.32	110.52
1	A	497	THR	CB-CA-C	-5.23	99.07	109.79
1	A	409	THR	N-CA-C	5.20	116.19	108.86
1	A	169	GLU	N-CA-C	5.19	119.81	112.75
2	B	232	TYR	N-CA-C	5.18	121.84	110.80
1	A	343	GLN	N-CA-C	-5.14	107.39	113.97
1	A	32	LYS	N-CA-C	-5.12	106.19	112.90
2	B	140	PRO	N-CA-C	5.12	121.37	113.75
2	B	178	ILE	N-CA-C	5.11	116.11	108.54
1	A	401	TRP	N-CA-C	5.08	116.63	111.14
1	A	147	ASN	N-CA-C	-5.06	105.80	112.94
2	B	156	SER	CA-C-N	-5.05	113.44	119.05
2	B	156	SER	C-N-CA	-5.05	113.44	119.05
2	B	312	GLU	CA-C-N	5.01	126.10	119.84
2	B	312	GLU	C-N-CA	5.01	126.10	119.84
1	A	168	LEU	N-CA-C	5.00	117.62	111.82
2	B	252	TRP	N-CA-C	5.00	116.43	108.67

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	146	TYR	Sidechain

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	4451	0	4430	244	0
2	B	3443	0	3415	179	0
3	A	27	0	14	2	0
All	All	7921	0	7859	413	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 26.

All (413) 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)
2:B:284:ARG:HH11	2:B:284:ARG:HB2	1.08	1.18
2:B:31:ILE:HG23	2:B:132:ILE:HD11	1.40	1.03
1:A:454:LYS:HD2	1:A:552:VAL:HA	1.41	1.02
2:B:85:GLN:HA	2:B:88:TRP:HB2	1.43	1.00
1:A:442:VAL:HG13	1:A:481:ALA:HB1	1.41	0.99
2:B:373:GLN:HE22	2:B:407:GLN:H	1.10	0.96
2:B:85:GLN:HG3	2:B:154:LYS:CB	1.95	0.96
1:A:417:VAL:HG12	1:A:419:THR:HG23	1.47	0.94
2:B:85:GLN:HG3	2:B:154:LYS:HB3	1.47	0.94
1:A:447:ASN:HD22	1:A:556:ILE:HG12	1.30	0.93
1:A:542:ILE:HG22	1:A:543:GLY:H	1.32	0.93
1:A:132:ILE:HB	1:A:142:ILE:HG23	1.48	0.92
1:A:109:LEU:HD13	1:A:216:THR:HG21	1.52	0.91
2:B:2:ILE:O	2:B:2:ILE:HG13	1.69	0.91
1:A:3:SER:OG	1:A:5:ILE:HG13	1.71	0.90
1:A:51:GLY:H	1:A:52:PRO:HD2	1.38	0.87
2:B:292:VAL:HG12	2:B:294:PRO:HD3	1.59	0.83
2:B:284:ARG:HB2	2:B:284:ARG:NH1	1.92	0.82
1:A:254:VAL:HG13	1:A:283:LEU:HD22	1.60	0.82
2:B:278:GLN:O	2:B:281:LYS:HG2	1.81	0.81
1:A:51:GLY:H	1:A:52:PRO:CD	1.93	0.81
1:A:542:ILE:HG22	1:A:543:GLY:N	1.94	0.81
1:A:64:LYS:HD3	1:A:71:TRP:CE2	2.16	0.80
1:A:545:ASN:O	1:A:549:ASP:HB2	1.83	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:TRP:NE1	2:B:420:PRO:HG3	1.98	0.78
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.67	0.77
1:A:542:ILE:H	1:A:545:ASN:HB2	1.49	0.77
1:A:164:MET:HE1	1:A:214:LEU:HD13	1.67	0.76
2:B:24:TRP:CE3	2:B:25:PRO:HD2	2.20	0.76
2:B:363:ASN:HD22	2:B:364:ASP:H	1.34	0.76
2:B:242:GLN:N	2:B:243:PRO:HD3	2.01	0.76
2:B:249:LYS:HB2	2:B:252:TRP:CZ3	2.21	0.75
1:A:199:ARG:NH1	1:A:220:LYS:HE2	2.02	0.75
1:A:50:ILE:HG13	1:A:50:ILE:O	1.85	0.74
1:A:148:VAL:O	1:A:150:PRO:HD3	1.87	0.74
1:A:60:VAL:HG21	1:A:130:PHE:HD2	1.53	0.74
2:B:363:ASN:HD22	2:B:364:ASP:N	1.85	0.74
1:A:51:GLY:N	1:A:52:PRO:HD2	2.03	0.73
1:A:53:GLU:O	1:A:55:PRO:HD3	1.89	0.73
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.70	0.73
2:B:284:ARG:HH11	2:B:284:ARG:CB	1.96	0.73
1:A:288:ALA:HB1	1:A:291:GLU:HB3	1.72	0.72
2:B:241:VAL:HG12	2:B:351:THR:O	1.89	0.72
1:A:447:ASN:ND2	1:A:556:ILE:HG12	2.05	0.72
2:B:215:THR:HG22	2:B:216:THR:H	1.53	0.71
1:A:57:ASN:OD1	1:A:143:ARG:NH1	2.24	0.71
2:B:303:LEU:HD21	2:B:307:ARG:HH21	1.56	0.71
2:B:206:ARG:HH12	2:B:218:ASP:HA	1.55	0.71
1:A:542:ILE:CG2	1:A:543:GLY:H	2.02	0.71
2:B:252:TRP:CD1	2:B:295:LEU:HG	2.26	0.70
1:A:469:LEU:HD11	1:A:480:GLN:HG2	1.71	0.70
2:B:94:ILE:HG12	2:B:161:GLN:NE2	2.07	0.69
1:A:23:GLN:NE2	1:A:131:THR:O	2.25	0.69
1:A:501:TYR:CZ	1:A:505:ILE:HD11	2.28	0.69
1:A:454:LYS:CD	1:A:552:VAL:HA	2.20	0.69
1:A:34:LEU:HD21	1:A:62:ALA:HB2	1.74	0.69
2:B:85:GLN:HG3	2:B:154:LYS:HB2	1.75	0.68
2:B:206:ARG:NH1	2:B:218:ASP:HA	2.07	0.68
2:B:244:ILE:HD13	2:B:244:ILE:H	1.58	0.68
1:A:111:VAL:HG12	1:A:185:ASP:O	1.94	0.68
1:A:424:LYS:HE3	1:A:426:TRP:CH2	2.29	0.68
2:B:87:PHE:O	2:B:88:TRP:CD1	2.47	0.68
2:B:373:GLN:NE2	2:B:407:GLN:H	1.88	0.68
1:A:389:PHE:HB3	1:A:391:LEU:HD21	1.76	0.68
1:A:130:PHE:CE1	1:A:144:TYR:HB2	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:VAL:HG21	1:A:130:PHE:CD2	2.28	0.67
2:B:275:LYS:HE3	2:B:277:ARG:HD3	1.77	0.67
1:A:52:PRO:HG2	1:A:53:GLU:H	1.58	0.67
2:B:323:LYS:HB2	2:B:343:GLN:OE1	1.94	0.67
1:A:454:LYS:HD2	1:A:552:VAL:CA	2.20	0.67
2:B:257:ILE:O	2:B:261:VAL:HG23	1.95	0.66
1:A:135:ILE:HG13	1:A:138:GLU:CD	2.21	0.66
1:A:66:LYS:C	1:A:68:SER:H	2.03	0.66
1:A:417:VAL:HG12	1:A:419:THR:CG2	2.26	0.66
1:A:90:VAL:O	1:A:91:GLN:HG2	1.96	0.65
2:B:2:ILE:O	2:B:2:ILE:CG1	2.42	0.65
2:B:132:ILE:HD12	2:B:133:PRO:HD2	1.77	0.65
1:A:362:THR:HG21	1:A:367:GLN:HG3	1.79	0.65
1:A:97:PRO:HA	1:A:100:LEU:HD12	1.78	0.65
1:A:442:VAL:HG13	1:A:481:ALA:CB	2.23	0.65
2:B:24:TRP:HE3	2:B:25:PRO:HD2	1.61	0.64
1:A:406:TRP:HE1	2:B:420:PRO:HG3	1.62	0.64
1:A:328:GLU:HG2	1:A:390:LYS:HB2	1.79	0.64
1:A:26:LEU:CB	1:A:30:LYS:HE2	2.27	0.64
2:B:343:GLN:HE21	2:B:349:LEU:HD11	1.63	0.64
2:B:315:HIS:O	2:B:317:VAL:N	2.31	0.63
2:B:274:ILE:HB	2:B:306:ASN:OD1	1.98	0.63
2:B:215:THR:HG22	2:B:216:THR:N	2.13	0.63
2:B:425:LEU:HG	2:B:425:LEU:O	1.99	0.63
1:A:202:ILE:O	1:A:206:ARG:HG3	1.98	0.63
1:A:132:ILE:HB	1:A:142:ILE:CG2	2.27	0.62
1:A:427:TYR:CE2	1:A:525:LEU:HD13	2.34	0.62
2:B:358:ARG:HA	2:B:362:THR:HA	1.80	0.62
1:A:97:PRO:HA	1:A:100:LEU:CD1	2.29	0.62
1:A:345:PRO:O	1:A:346:PHE:CD2	2.52	0.62
1:A:390:LYS:HB3	1:A:417:VAL:HG21	1.80	0.62
1:A:542:ILE:CG2	1:A:543:GLY:N	2.62	0.62
2:B:315:HIS:CG	2:B:316:GLY:N	2.68	0.62
2:B:132:ILE:HD12	2:B:133:PRO:CD	2.29	0.62
2:B:132:ILE:HD12	2:B:132:ILE:C	2.25	0.61
2:B:244:ILE:H	2:B:244:ILE:CD1	2.12	0.61
2:B:316:GLY:O	2:B:318:TYR:N	2.34	0.60
1:A:31:ILE:HD13	1:A:133:PRO:O	2.01	0.60
1:A:276:VAL:HG12	1:A:276:VAL:O	2.01	0.60
2:B:104:LYS:HG3	2:B:237:ASP:OD2	2.02	0.60
1:A:282:LEU:HD21	1:A:296:THR:HG23	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:GLY:C	1:A:114:ALA:H	2.09	0.59
1:A:130:PHE:CZ	1:A:144:TYR:HB2	2.38	0.59
1:A:552:VAL:HG12	1:A:552:VAL:O	2.02	0.59
2:B:308:GLU:O	2:B:310:LEU:N	2.35	0.59
2:B:320:ASP:OD1	2:B:322:SER:HB3	2.03	0.59
1:A:266:TRP:O	1:A:269:GLN:HG2	2.03	0.59
2:B:12:LEU:O	2:B:14:PRO:HD3	2.03	0.59
1:A:174:GLN:C	1:A:176:PRO:HD3	2.28	0.58
2:B:371:ALA:O	2:B:375:ILE:HG13	2.03	0.58
1:A:37:ILE:CG2	1:A:41:MET:HE3	2.34	0.58
1:A:498:ASP:HB2	1:A:538:ALA:HA	1.86	0.58
2:B:244:ILE:HG12	2:B:244:ILE:O	2.04	0.58
1:A:138:GLU:O	1:A:140:PRO:HD3	2.04	0.57
1:A:239:TRP:CZ2	1:A:270:ILE:HD12	2.39	0.57
1:A:23:GLN:O	1:A:23:GLN:HG3	2.02	0.57
2:B:422:LEU:O	2:B:426:TRP:CZ3	2.57	0.57
2:B:209:LEU:HD22	2:B:214:LEU:HD12	1.87	0.57
2:B:423:VAL:HG13	2:B:423:VAL:O	2.03	0.57
2:B:132:ILE:HG23	2:B:142:ILE:HB	1.87	0.57
1:A:556:ILE:HD12	1:A:556:ILE:N	2.19	0.57
2:B:328:GLU:HG2	2:B:390:LYS:HD2	1.86	0.57
2:B:178:ILE:HD11	2:B:201:LYS:HD3	1.85	0.56
1:A:543:GLY:HA2	2:B:284:ARG:O	2.04	0.56
2:B:234:LEU:HD11	2:B:377:THR:HG21	1.85	0.56
2:B:330:GLN:HE22	2:B:340:GLN:HE22	1.51	0.56
1:A:5:ILE:HD11	1:A:167:ILE:HD11	1.87	0.56
1:A:10:VAL:HG12	1:A:11:LYS:N	2.20	0.56
1:A:411:ILE:HG13	1:A:411:ILE:O	2.05	0.56
1:A:108:VAL:HG13	1:A:223:LYS:CB	2.36	0.56
1:A:454:LYS:HD2	1:A:552:VAL:HG13	1.87	0.56
2:B:306:ASN:O	2:B:308:GLU:N	2.38	0.56
1:A:254:VAL:HG13	1:A:283:LEU:CD2	2.33	0.56
2:B:56:TYR:HE1	2:B:126:LYS:HE3	1.71	0.56
1:A:27:THR:C	1:A:29:GLU:N	2.64	0.56
1:A:177:ASP:OD1	1:A:193:LEU:HD21	2.06	0.56
2:B:87:PHE:O	2:B:88:TRP:HD1	1.87	0.56
2:B:306:ASN:C	2:B:308:GLU:N	2.62	0.56
1:A:151:GLN:O	1:A:151:GLN:HG2	2.04	0.55
2:B:60:VAL:HG11	2:B:130:PHE:CD2	2.41	0.55
2:B:109:LEU:HD22	2:B:216:THR:HG21	1.88	0.55
2:B:293:ILE:HG12	2:B:293:ILE:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:257:ILE:HD13	2:B:283:LEU:HD13	1.87	0.55
1:A:27:THR:C	1:A:29:GLU:H	2.14	0.55
1:A:33:ALA:HB1	1:A:71:TRP:CG	2.41	0.55
2:B:64:LYS:HE3	2:B:71:TRP:CH2	2.42	0.55
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.89	0.55
1:A:229:TRP:CH2	3:A:561:T50:H14	2.42	0.55
2:B:132:ILE:HD12	2:B:133:PRO:N	2.22	0.55
2:B:308:GLU:C	2:B:310:LEU:N	2.64	0.55
1:A:137:ASN:O	1:A:139:THR:N	2.37	0.55
2:B:363:ASN:ND2	2:B:364:ASP:N	2.54	0.54
2:B:79:GLU:OE2	2:B:83:ARG:NH1	2.40	0.54
2:B:139:THR:HB	2:B:140:PRO:HD2	1.89	0.54
1:A:199:ARG:HH11	1:A:220:LYS:HE2	1.73	0.54
1:A:362:THR:CG2	1:A:367:GLN:HG3	2.37	0.54
1:A:228:LEU:HB3	1:A:242:GLN:NE2	2.23	0.54
2:B:423:VAL:O	2:B:423:VAL:CG1	2.56	0.54
1:A:170:PRO:O	1:A:174:GLN:HG3	2.07	0.54
1:A:215:THR:C	1:A:217:PRO:HD3	2.33	0.54
1:A:544:GLY:HA2	1:A:548:VAL:HG23	1.88	0.54
1:A:51:GLY:N	1:A:52:PRO:CD	2.62	0.54
1:A:543:GLY:HA3	2:B:285:GLY:O	2.07	0.54
1:A:324:ASP:O	1:A:343:GLN:HG2	2.08	0.53
1:A:60:VAL:CG2	1:A:130:PHE:HB2	2.37	0.53
2:B:306:ASN:C	2:B:308:GLU:H	2.17	0.53
2:B:197:GLN:O	2:B:200:THR:HB	2.09	0.53
1:A:334:GLN:CD	1:A:512:LYS:HG3	2.33	0.53
1:A:556:ILE:HD12	1:A:556:ILE:H	1.74	0.53
1:A:345:PRO:O	1:A:346:PHE:CG	2.62	0.53
1:A:366:LYS:HE3	1:A:370:GLU:OE2	2.08	0.53
1:A:270:ILE:HD11	1:A:316:GLY:N	2.25	0.52
1:A:57:ASN:OD1	1:A:131:THR:HB	2.09	0.52
1:A:77:PHE:O	1:A:79:GLU:N	2.42	0.52
2:B:423:VAL:O	2:B:427:TYR:HE1	1.92	0.52
1:A:454:LYS:CG	1:A:552:VAL:HG13	2.40	0.52
2:B:319:TYR:O	2:B:321:PRO:HD3	2.10	0.52
1:A:88:TRP:O	1:A:89:GLU:C	2.52	0.52
2:B:241:VAL:HG23	2:B:241:VAL:O	2.09	0.52
2:B:103:LYS:HE2	2:B:103:LYS:HA	1.92	0.52
2:B:173:LYS:O	2:B:176:PRO:HD3	2.10	0.51
1:A:328:GLU:CG	1:A:390:LYS:HB2	2.40	0.51
1:A:108:VAL:CG1	1:A:223:LYS:CB	2.88	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:295:LEU:HB3	2:B:299:ALA:HB3	1.93	0.51
1:A:115:TYR:CE1	1:A:156:SER:HB3	2.44	0.51
1:A:219:LYS:O	1:A:220:LYS:HB3	2.10	0.51
1:A:290:THR:O	1:A:291:GLU:C	2.54	0.51
2:B:308:GLU:C	2:B:310:LEU:H	2.18	0.51
1:A:5:ILE:HG22	1:A:6:GLU:O	2.11	0.51
1:A:13:LYS:HB2	1:A:16:MET:HG3	1.92	0.51
1:A:135:ILE:HG13	1:A:138:GLU:OE1	2.10	0.51
2:B:5:ILE:HG22	2:B:6:GLU:O	2.11	0.51
1:A:448:ARG:HG2	1:A:448:ARG:HH11	1.77	0.50
1:A:458:VAL:HG22	1:A:464:GLN:HG2	1.93	0.50
1:A:97:PRO:HG2	1:A:232:TYR:CD2	2.47	0.50
1:A:215:THR:O	1:A:217:PRO:HD3	2.11	0.50
1:A:500:GLN:HE21	2:B:422:LEU:CD1	2.24	0.50
1:A:194:GLU:O	1:A:196:GLY:N	2.45	0.50
2:B:281:LYS:HA	2:B:284:ARG:NH1	2.26	0.50
1:A:457:TYR:CD1	1:A:457:TYR:C	2.88	0.50
2:B:244:ILE:HD13	2:B:244:ILE:N	2.26	0.50
2:B:64:LYS:HE3	2:B:71:TRP:CZ2	2.47	0.50
1:A:30:LYS:HD2	1:A:62:ALA:HB3	1.92	0.50
1:A:111:VAL:O	1:A:111:VAL:HG13	2.12	0.50
1:A:277:ARG:HB2	1:A:336:GLN:CD	2.37	0.50
1:A:290:THR:O	1:A:290:THR:HG22	2.11	0.50
1:A:330:GLN:NE2	1:A:340:GLN:OE1	2.34	0.50
1:A:112:GLY:O	1:A:114:ALA:N	2.34	0.50
1:A:66:LYS:C	1:A:68:SER:N	2.70	0.50
1:A:420:PRO:HA	1:A:421:PRO:C	2.37	0.50
2:B:183:TYR:CD2	2:B:380:ILE:HD13	2.47	0.50
2:B:242:GLN:N	2:B:243:PRO:CD	2.75	0.49
1:A:112:GLY:C	1:A:114:ALA:N	2.70	0.49
1:A:60:VAL:HG23	1:A:130:PHE:HB2	1.94	0.49
1:A:398:TRP:CD1	1:A:398:TRP:C	2.91	0.49
2:B:234:LEU:HD11	2:B:377:THR:CG2	2.42	0.49
2:B:395:LYS:O	2:B:399:GLU:CG	2.61	0.49
1:A:143:ARG:HH11	1:A:143:ARG:HG2	1.75	0.49
2:B:65:LYS:HG3	2:B:72:ARG:HB2	1.94	0.49
1:A:50:ILE:HD11	1:A:54:ASN:HD22	1.78	0.48
2:B:157:PRO:HG3	2:B:184:MET:HA	1.95	0.48
2:B:169:GLU:HB2	2:B:170:PRO:HD3	1.96	0.48
2:B:403:THR:HG23	2:B:403:THR:O	2.12	0.48
1:A:273:GLY:O	1:A:274:ILE:C	2.56	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:LYS:HB2	1:A:192:ASP:HA	1.95	0.48
2:B:178:ILE:CD1	2:B:201:LYS:HD3	2.44	0.48
1:A:354:TYR:OH	1:A:370:GLU:OE1	2.31	0.48
1:A:433:PRO:HB3	2:B:289:LEU:CD1	2.44	0.48
2:B:2:ILE:O	2:B:4:PRO:HD3	2.14	0.48
1:A:103:LYS:HE3	1:A:179:VAL:HG11	1.96	0.47
1:A:276:VAL:HG12	1:A:280:SER:HB2	1.95	0.47
2:B:115:TYR:O	2:B:117:SER:N	2.47	0.47
1:A:369:THR:HG23	1:A:398:TRP:HH2	1.79	0.47
2:B:49:LYS:HG3	2:B:144:TYR:CE1	2.49	0.47
1:A:49:LYS:O	1:A:50:ILE:CG2	2.62	0.47
2:B:27:THR:HG22	2:B:30:LYS:HG3	1.97	0.47
1:A:53:GLU:C	1:A:55:PRO:HD3	2.39	0.47
1:A:131:THR:HB	1:A:143:ARG:HH11	1.79	0.47
2:B:77:PHE:O	2:B:78:ARG:C	2.58	0.47
2:B:101:LYS:O	2:B:236:PRO:HB2	2.15	0.47
1:A:135:ILE:H	1:A:135:ILE:HG12	1.48	0.47
1:A:326:ILE:O	1:A:341:ILE:HA	2.15	0.47
2:B:235:HIS:C	2:B:237:ASP:H	2.23	0.47
1:A:49:LYS:HG3	1:A:144:TYR:CE2	2.50	0.47
1:A:390:LYS:HB3	1:A:417:VAL:CG2	2.44	0.47
2:B:88:TRP:CZ3	2:B:92:LEU:HD12	2.49	0.47
2:B:266:TRP:CZ3	2:B:426:TRP:HB3	2.50	0.47
1:A:434:ILE:HD11	1:A:530:LYS:HB3	1.97	0.47
2:B:88:TRP:HZ3	2:B:159:ILE:HG13	1.80	0.47
1:A:94:ILE:HG21	1:A:230:MET:HE2	1.96	0.46
1:A:358:ARG:HH12	1:A:366:LYS:HE2	1.80	0.46
2:B:423:VAL:O	2:B:427:TYR:CE1	2.68	0.46
1:A:314:VAL:HG12	1:A:315:HIS:N	2.29	0.46
1:A:344:GLU:HB3	1:A:345:PRO:HD2	1.97	0.46
1:A:472:THR:OG1	1:A:476:LYS:HB2	2.16	0.46
1:A:177:ASP:OD2	1:A:177:ASP:N	2.48	0.46
1:A:278:GLN:OE1	1:A:298:GLU:HB3	2.14	0.46
1:A:483:TYR:CZ	1:A:520:GLN:HG2	2.49	0.46
1:A:226:PRO:HB3	1:A:235:HIS:CE1	2.50	0.46
2:B:330:GLN:HG2	2:B:338:THR:OG1	2.16	0.46
1:A:219:LYS:O	1:A:220:LYS:CB	2.63	0.46
1:A:220:LYS:O	1:A:221:HIS:CB	2.63	0.46
1:A:433:PRO:HB3	2:B:289:LEU:HD13	1.98	0.46
1:A:5:ILE:HG22	1:A:6:GLU:N	2.31	0.46
2:B:85:GLN:O	2:B:86:ASP:C	2.56	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:392:PRO:HA	2:B:417:VAL:HG22	1.96	0.46
1:A:12:LEU:O	1:A:13:LYS:C	2.58	0.46
1:A:111:VAL:CG1	1:A:185:ASP:O	2.62	0.46
1:A:451:LYS:O	1:A:471:ASN:N	2.49	0.46
2:B:115:TYR:C	2:B:117:SER:H	2.24	0.46
2:B:175:ASN:OD1	2:B:201:LYS:HE2	2.15	0.46
2:B:214:LEU:N	2:B:214:LEU:HD23	2.31	0.46
1:A:497:THR:HG22	1:A:499:SER:H	1.81	0.45
1:A:500:GLN:HE21	2:B:422:LEU:HD13	1.81	0.45
1:A:49:LYS:O	1:A:50:ILE:HG23	2.16	0.45
2:B:124:PHE:O	2:B:125:ARG:C	2.59	0.45
1:A:147:ASN:C	1:A:148:VAL:HG13	2.41	0.45
1:A:411:ILE:O	1:A:412:PRO:O	2.35	0.45
1:A:441:TYR:HB3	1:A:548:VAL:HG21	1.98	0.45
2:B:194:GLU:O	2:B:196:GLY:N	2.49	0.45
1:A:134:SER:HB3	1:A:139:THR:HA	1.99	0.45
2:B:198:HIS:C	2:B:200:THR:N	2.72	0.45
2:B:407:GLN:HE21	2:B:407:GLN:HB3	1.59	0.45
1:A:50:ILE:CD1	1:A:54:ASN:HD22	2.30	0.45
2:B:395:LYS:O	2:B:399:GLU:HG2	2.17	0.45
1:A:171:PHE:CD2	1:A:171:PHE:C	2.94	0.45
2:B:238:LYS:HG2	2:B:238:LYS:O	2.16	0.45
2:B:245:VAL:HG12	2:B:246:LEU:N	2.31	0.45
2:B:332:GLN:HB3	2:B:333:GLY:H	1.51	0.45
1:A:111:VAL:O	1:A:185:ASP:HB2	2.17	0.45
1:A:137:ASN:C	1:A:139:THR:N	2.75	0.45
1:A:10:VAL:HG21	1:A:153:TRP:HH2	1.82	0.45
1:A:291:GLU:O	1:A:293:ILE:N	2.50	0.45
2:B:270:ILE:O	2:B:272:PRO:HD3	2.17	0.45
1:A:194:GLU:O	1:A:195:ILE:C	2.60	0.45
2:B:232:TYR:CE2	2:B:234:LEU:HD21	2.52	0.45
2:B:279:LEU:C	2:B:281:LYS:H	2.26	0.44
1:A:382:ILE:O	2:B:136:ASN:HB2	2.16	0.44
2:B:206:ARG:CZ	2:B:217:PRO:O	2.66	0.44
2:B:369:THR:HG23	2:B:406:TRP:CE3	2.52	0.44
1:A:142:ILE:HG23	1:A:142:ILE:O	2.17	0.44
1:A:175:ASN:N	1:A:176:PRO:HD3	2.33	0.44
2:B:87:PHE:C	2:B:88:TRP:CD1	2.96	0.44
2:B:258:GLN:HG3	2:B:283:LEU:HD21	1.98	0.44
1:A:10:VAL:CG1	1:A:11:LYS:N	2.81	0.44
1:A:52:PRO:HG2	1:A:53:GLU:N	2.31	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:PRO:O	1:A:219:LYS:N	2.51	0.44
2:B:257:ILE:HG12	2:B:283:LEU:HD22	1.99	0.44
1:A:77:PHE:O	1:A:80:LEU:N	2.50	0.44
1:A:405:TYR:CE2	1:A:407:GLN:HB2	2.53	0.44
1:A:27:THR:OG1	1:A:29:GLU:HB3	2.16	0.44
1:A:134:SER:CB	1:A:139:THR:HA	2.48	0.44
1:A:358:ARG:O	1:A:359:GLY:C	2.59	0.44
2:B:31:ILE:HG23	2:B:132:ILE:CD1	2.29	0.44
2:B:76:ASP:OD1	2:B:76:ASP:C	2.61	0.44
1:A:108:VAL:HG12	1:A:227:PHE:HE1	1.82	0.43
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.53	0.43
2:B:241:VAL:CG1	2:B:351:THR:H	2.30	0.43
2:B:242:GLN:H	2:B:243:PRO:HD3	1.81	0.43
1:A:362:THR:CG2	1:A:363:ASN:N	2.79	0.43
2:B:203:GLU:O	2:B:203:GLU:CG	2.65	0.43
1:A:545:ASN:HA	1:A:549:ASP:CG	2.42	0.43
2:B:125:ARG:HD3	2:B:147:ASN:HA	2.00	0.43
2:B:315:HIS:ND1	2:B:316:GLY:N	2.66	0.43
1:A:501:TYR:CE1	1:A:505:ILE:HD11	2.53	0.43
2:B:87:PHE:C	2:B:88:TRP:HD1	2.25	0.43
2:B:395:LYS:O	2:B:399:GLU:HG3	2.18	0.43
1:A:169:GLU:O	1:A:173:LYS:HB2	2.18	0.43
1:A:334:GLN:NE2	1:A:512:LYS:HG3	2.33	0.43
2:B:34:LEU:HD23	2:B:34:LEU:HA	1.81	0.43
1:A:515:SER:O	1:A:519:ASN:ND2	2.50	0.43
2:B:376:THR:HG21	2:B:410:TRP:CZ3	2.53	0.43
1:A:64:LYS:HD3	1:A:71:TRP:CZ2	2.53	0.43
1:A:130:PHE:CD1	1:A:130:PHE:N	2.86	0.43
1:A:224:GLU:H	1:A:224:GLU:HG3	1.55	0.43
1:A:310:LEU:HD12	1:A:310:LEU:HA	1.88	0.43
1:A:335:GLY:O	1:A:355:ALA:HA	2.18	0.43
2:B:323:LYS:NZ	2:B:344:GLU:OE2	2.49	0.43
2:B:422:LEU:HD23	2:B:422:LEU:HA	1.70	0.43
1:A:121:ASP:O	1:A:122:GLU:C	2.62	0.43
1:A:443:ASP:HB3	1:A:548:VAL:HG12	2.00	0.43
2:B:23:GLN:OE1	2:B:60:VAL:HG22	2.18	0.43
2:B:85:GLN:CG	2:B:154:LYS:HB2	2.45	0.43
2:B:139:THR:CB	2:B:140:PRO:HD2	2.46	0.43
1:A:417:VAL:CG1	1:A:419:THR:HG23	2.33	0.43
1:A:444:GLY:O	1:A:445:ALA:HB2	2.19	0.43
1:A:377:THR:O	1:A:381:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LYS:CE	1:A:130:PHE:CE2	3.03	0.42
1:A:114:ALA:HB1	1:A:160:PHE:CZ	2.54	0.42
2:B:85:GLN:C	2:B:87:PHE:N	2.76	0.42
1:A:406:TRP:CZ3	1:A:407:GLN:HG2	2.54	0.42
1:A:408:ALA:HB1	2:B:364:ASP:HB3	2.01	0.42
2:B:53:GLU:OE1	2:B:53:GLU:N	2.41	0.42
1:A:178:ILE:HD11	1:A:201:LYS:HB3	2.02	0.42
1:A:238:LYS:HD2	1:A:315:HIS:CG	2.54	0.42
2:B:266:TRP:CG	2:B:426:TRP:CE3	3.07	0.42
1:A:24:TRP:O	1:A:25:PRO:C	2.60	0.42
1:A:131:THR:HA	1:A:143:ARG:HG2	2.01	0.42
1:A:3:SER:C	1:A:5:ILE:H	2.28	0.42
2:B:193:LEU:HD11	2:B:201:LYS:HD2	2.02	0.42
2:B:216:THR:HA	2:B:217:PRO:HD3	1.89	0.42
2:B:423:VAL:HG13	2:B:427:TYR:OH	2.19	0.42
1:A:21:VAL:HG12	1:A:22:LYS:N	2.35	0.42
1:A:115:TYR:N	1:A:115:TYR:CD2	2.88	0.42
2:B:149:LEU:HA	2:B:150:PRO:HD3	1.75	0.42
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.55	0.42
1:A:540:LYS:O	1:A:542:ILE:HG13	2.19	0.42
2:B:237:ASP:C	2:B:239:TRP:H	2.27	0.42
1:A:9:PRO:O	1:A:9:PRO:HG2	2.20	0.42
2:B:115:TYR:HE1	2:B:185:ASP:OD1	2.02	0.42
2:B:194:GLU:O	2:B:197:GLN:N	2.50	0.41
1:A:13:LYS:HD2	1:A:16:MET:HE3	2.01	0.41
1:A:199:ARG:HH12	1:A:220:LYS:HE2	1.79	0.41
1:A:398:TRP:CH2	1:A:411:ILE:HD13	2.55	0.41
2:B:28:GLU:CB	2:B:135:ILE:HD11	2.50	0.41
2:B:426:TRP:O	2:B:427:TYR:CB	2.69	0.41
1:A:308:GLU:O	1:A:311:LYS:CB	2.68	0.41
1:A:417:VAL:CG1	1:A:419:THR:CG2	2.97	0.41
2:B:277:ARG:NH2	2:B:302:GLU:OE1	2.53	0.41
1:A:503:LEU:HD11	1:A:507:GLN:OE1	2.20	0.41
1:A:179:VAL:HG22	3:A:561:T50:H20	2.02	0.41
1:A:276:VAL:O	1:A:276:VAL:CG1	2.68	0.41
2:B:85:GLN:CB	2:B:154:LYS:HB2	2.50	0.41
2:B:215:THR:CG2	2:B:216:THR:H	2.28	0.41
2:B:312:GLU:HA	2:B:313:PRO:HD2	1.81	0.41
1:A:398:TRP:CZ2	1:A:411:ILE:HD13	2.55	0.41
1:A:446:ALA:HB2	1:A:472:THR:HG23	2.02	0.41
1:A:478:GLU:CD	1:A:499:SER:HB2	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:GLN:N	1:A:510:PRO:CD	2.83	0.41
2:B:210:LEU:HD12	2:B:214:LEU:O	2.20	0.41
1:A:20:LYS:HE2	1:A:55:PRO:O	2.21	0.41
1:A:270:ILE:HD11	1:A:316:GLY:CA	2.50	0.41
2:B:2:ILE:O	2:B:4:PRO:CD	2.69	0.41
2:B:308:GLU:O	2:B:309:ILE:C	2.64	0.41
2:B:369:THR:HG23	2:B:406:TRP:HE3	1.86	0.41
1:A:49:LYS:CG	1:A:144:TYR:CE2	3.05	0.40
1:A:194:GLU:C	1:A:196:GLY:N	2.79	0.40
1:A:424:LYS:HE3	1:A:426:TRP:CZ2	2.56	0.40
1:A:320:ASP:HA	1:A:321:PRO:HD3	1.79	0.40
1:A:398:TRP:CH2	1:A:411:ILE:HG21	2.57	0.40
1:A:254:VAL:O	1:A:258:GLN:HG3	2.21	0.40
2:B:354:TYR:OH	2:B:370:GLU:HB3	2.21	0.40
1:A:131:THR:CB	1:A:143:ARG:HG2	2.51	0.40
2:B:207:GLN:O	2:B:211:ARG:HB2	2.21	0.40
2:B:266:TRP:HB2	2:B:426:TRP:CZ3	2.56	0.40
2:B:306:ASN:O	2:B:307:ARG:C	2.64	0.40
1:A:376:THR:HG23	1:A:386:THR:HG22	2.02	0.40
2:B:75:VAL:HG11	2:B:77:PHE:CZ	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	556/560 (99%)	480 (86%)	60 (11%)	16 (3%)	3	5
2	B	425/430 (99%)	347 (82%)	59 (14%)	19 (4%)	2	2
All	All	981/990 (99%)	827 (84%)	119 (12%)	35 (4%)	2	4

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	412	PRO
1	A	538	ALA
2	B	225	PRO
2	B	295	LEU
2	B	316	GLY
2	B	317	VAL
2	B	357	MET
1	A	78	ARG
1	A	218	ASP
1	A	345	PRO
1	A	358	ARG
1	A	541	GLY
2	B	85	GLN
2	B	93	GLY
2	B	116	PHE
2	B	307	ARG
2	B	309	ILE
2	B	359	GLY
1	A	90	VAL
1	A	113	ASP
1	A	114	ALA
1	A	135	ILE
1	A	270	ILE
2	B	238	LYS
2	B	14	PRO
2	B	362	THR
1	A	195	ILE
2	B	126	LYS
2	B	313	PRO
1	A	67	ASP
1	A	292	VAL
1	A	445	ALA
2	B	195	ILE
2	B	224	GLU
2	B	236	PRO

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	472/500 (94%)	441 (93%)	31 (7%)	15	26
2	B	368/392 (94%)	340 (92%)	28 (8%)	12	21
All	All	840/892 (94%)	781 (93%)	59 (7%)	14	24

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	MET
1	A	22	LYS
1	A	23	GLN
1	A	40	GLU
1	A	75	VAL
1	A	89	GLU
1	A	90	VAL
1	A	110	ASP
1	A	122	GLU
1	A	130	PHE
1	A	135	ILE
1	A	178	ILE
1	A	185	ASP
1	A	224	GLU
1	A	241	VAL
1	A	245	VAL
1	A	274	ILE
1	A	278	GLN
1	A	286	THR
1	A	310	LEU
1	A	340	GLN
1	A	345	PRO
1	A	362	THR
1	A	387	PRO
1	A	396	GLU
1	A	402	TRP
1	A	431	LYS
1	A	497	THR
1	A	520	GLN
1	A	547	GLN
1	A	548	VAL
2	B	2	ILE
2	B	14	PRO
2	B	27	THR
2	B	35	VAL

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Mol	Chain	Res	Type
2	B	55	PRO
2	B	69	THR
2	B	111	VAL
2	B	132	ILE
2	B	174	GLN
2	B	193	LEU
2	B	214	LEU
2	B	244	ILE
2	B	248	GLU
2	B	252	TRP
2	B	257	ILE
2	B	274	ILE
2	B	283	LEU
2	B	284	ARG
2	B	293	ILE
2	B	314	VAL
2	B	315	HIS
2	B	330	GLN
2	B	332	GLN
2	B	334	GLN
2	B	399	GLU
2	B	403	THR
2	B	423	VAL
2	B	427	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	GLN
1	A	182	GLN
1	A	197	GLN
1	A	207	GLN
1	A	334	GLN
1	A	394	GLN
1	A	428	GLN
1	A	471	ASN
1	A	475	GLN
1	A	500	GLN
2	B	91	GLN
2	B	147	ASN
2	B	161	GLN
2	B	174	GLN

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Mol	Chain	Res	Type
2	B	235	HIS
2	B	258	GLN
2	B	330	GLN
2	B	343	GLN
2	B	363	ASN
2	B	373	GLN
2	B	407	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	$\# Z > 2$
3	T50	A	561	-	30,30,30	2.58	17 (56%)	43,44,44	1.88	10 (23%)

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	T50	A	561	-	-	0/8/8/8	0/4/4/4

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	561	T50	C6-C5	4.89	1.47	1.38
3	A	561	T50	C2-N1	-4.46	1.33	1.38
3	A	561	T50	C7-C8	3.83	1.45	1.39
3	A	561	T50	C20-C19	3.57	1.45	1.38
3	A	561	T50	C6-C7	3.54	1.45	1.38
3	A	561	T50	C20-C21	3.44	1.44	1.38
3	A	561	T50	C17-C2	3.30	1.53	1.48
3	A	561	T50	C21-C22	3.05	1.44	1.37
3	A	561	T50	C5-C4	2.89	1.45	1.39
3	A	561	T50	C17-C18	2.74	1.43	1.39
3	A	561	T50	C11-C12	2.58	1.42	1.38
3	A	561	T50	C13-C12	2.56	1.43	1.37
3	A	561	T50	C14-C13	2.51	1.43	1.38
3	A	561	T50	C4A-C4	2.48	1.55	1.51
3	A	561	T50	C9-C4	2.42	1.44	1.40
3	A	561	T50	C15-C16	2.37	1.42	1.37
3	A	561	T50	C14-C15	2.23	1.42	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	561	T50	C9-N3-C2	5.85	109.98	104.29
3	A	561	T50	C8-C9-N3	-4.42	104.95	110.31
3	A	561	T50	C17-C2-N3	-3.84	120.53	123.95
3	A	561	T50	C22-C17-C2	-3.79	119.39	122.33
3	A	561	T50	C16-C11-C12	3.52	119.28	114.54
3	A	561	T50	C8-N1-C2	3.25	108.32	106.25
3	A	561	T50	N1-C2-N3	-2.23	110.73	112.96
3	A	561	T50	C22-C17-C18	2.19	118.70	114.93
3	A	561	T50	C10-C11-C16	-2.18	117.63	122.50
3	A	561	T50	F3-C12-C11	2.08	120.23	117.64

There are no chirality outliers.

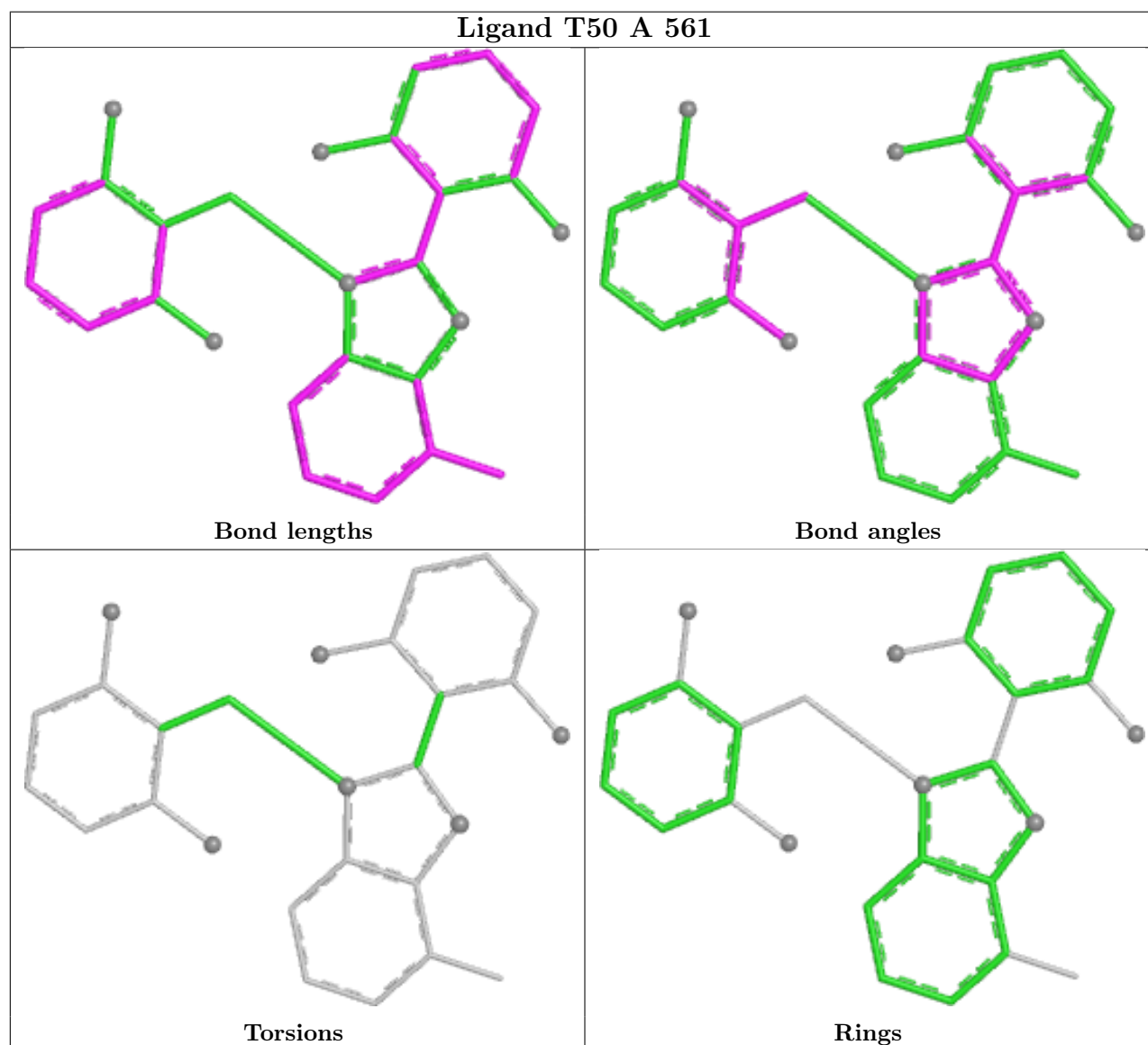
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	561	T50	2	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	558/560 (99%)	0.15	15 (2%)	56 49	47, 87, 121, 121	0
2	B	427/430 (99%)	0.10	27 (6%)	26 19	41, 74, 121, 121	0
All	All	985/990 (99%)	0.13	42 (4%)	40 32	41, 83, 121, 121	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	552	VAL	6.9
1	A	551	LEU	5.7
2	B	231	GLY	4.7
1	A	543	GLY	4.2
2	B	222	GLN	4.0
2	B	244	ILE	3.7
2	B	2	ILE	3.6
1	A	288	ALA	3.3
2	B	232	TYR	3.1
2	B	243	PRO	2.9
2	B	4	PRO	2.9
1	A	287	LYS	2.9
2	B	423	VAL	2.8
1	A	544	GLY	2.7
2	B	419	THR	2.7
2	B	221	HIS	2.7
2	B	427	TYR	2.7
1	A	51	GLY	2.6
2	B	1	PRO	2.6
2	B	294	PRO	2.6
1	A	225	PRO	2.6
2	B	357	MET	2.5
2	B	426	TRP	2.4
2	B	293	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	240	THR	2.3
1	A	556	ILE	2.3
1	A	360	ALA	2.3
2	B	226	PRO	2.2
2	B	299	ALA	2.2
2	B	225	PRO	2.2
1	A	295	LEU	2.2
1	A	539	HIS	2.2
2	B	87	PHE	2.2
2	B	285	GLY	2.1
1	A	136	ASN	2.1
2	B	292	VAL	2.1
2	B	271	TYR	2.1
1	A	256	ASP	2.0
2	B	266	TRP	2.0
2	B	422	LEU	2.0
1	A	69	THR	2.0
2	B	295	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](#)

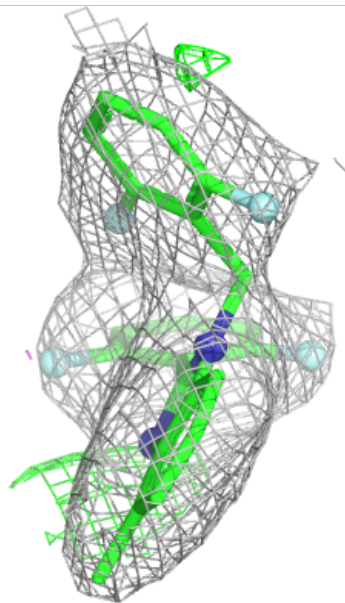
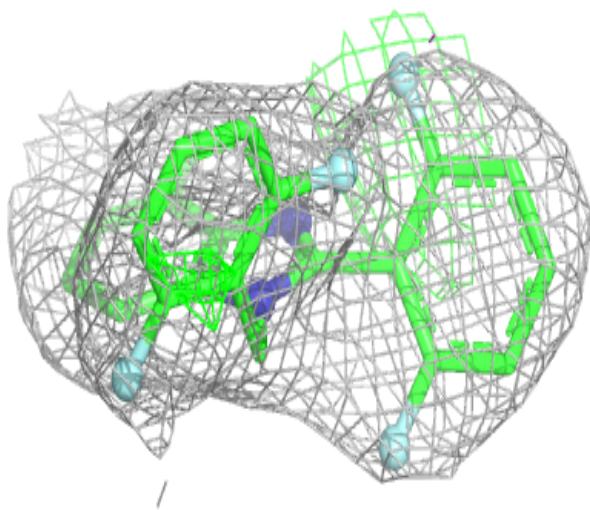
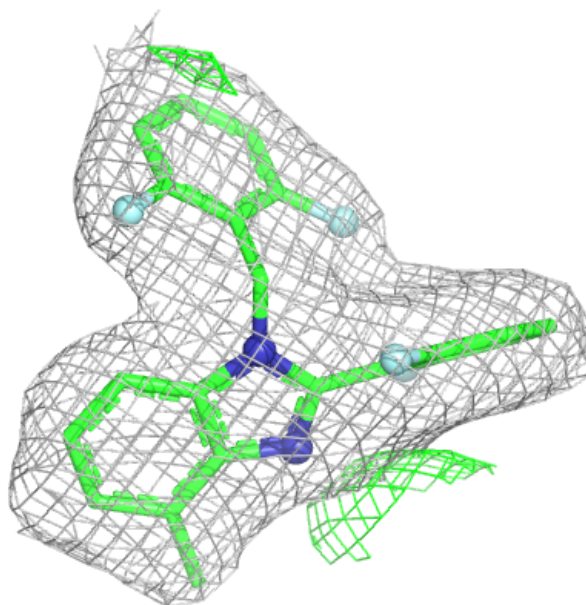
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	T50	A	561	27/27	0.94	0.07	47,53,66,72	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 T50 A 561:

$2mF_o-DF_c$ (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.