



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:38 AM UTC

PDB ID : 3CKP / pdb_00003ckp
Title : Crystal structure of BACE-1 in complex with inhibitor
Authors : Min, K.
Deposited on : 2008-03-16
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

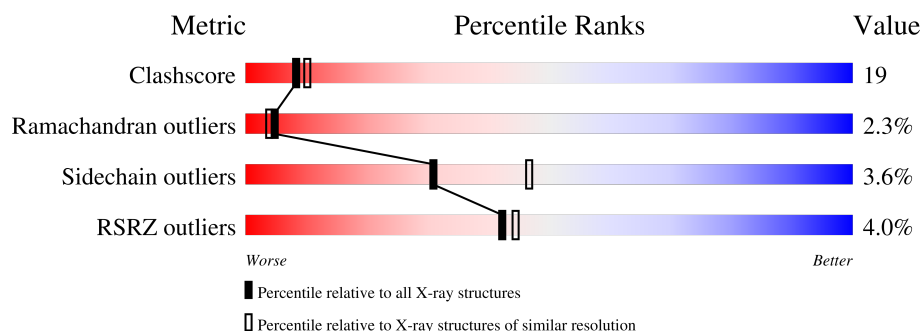
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	412	2% 59% 30% • 8%
1	B	412	4% 59% 29% • 9%
1	C	412	5% 45% 39% 6% • 8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9210 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 Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			2974	1904	496	560	14			
1	B	376	Total	C	N	O	S	0	0	0
			2956	1894	492	556	14			
1	C	377	Total	C	N	O	S	0	0	0
			2967	1900	494	559	14			

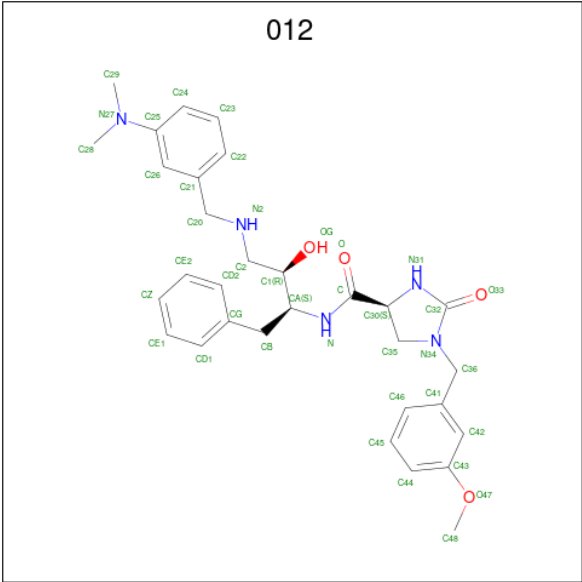
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LYS	ARG	engineered mutation	UNP P56817
A	-5	LYS	ARG	engineered mutation	UNP P56817
B	-6	LYS	ARG	engineered mutation	UNP P56817
B	-5	LYS	ARG	engineered mutation	UNP P56817
C	-6	LYS	ARG	engineered mutation	UNP P56817
C	-5	LYS	ARG	engineered mutation	UNP P56817

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Cl	0	0
			5	5		
2	B	2	Total	Cl	0	0
			2	2		
2	C	2	Total	Cl	0	0
			2	2		

- Molecule 3 is (4S)-N-[(1S,2R)-1-benzyl-3-{[3-(dimethylamino)benzyl]amino}-2-hydroxypropyl]-1-(3-methoxybenzyl)-2-oxoimidazolidine-4-carboxamide (CCD ID: 012) (formula: C₃₁H₃₉N₅O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			40	31	5	4		
3	B	1	Total	C	N	O	0	0
			40	31	5	4		
3	C	1	Total	C	N	O	0	0
			40	31	5	4		

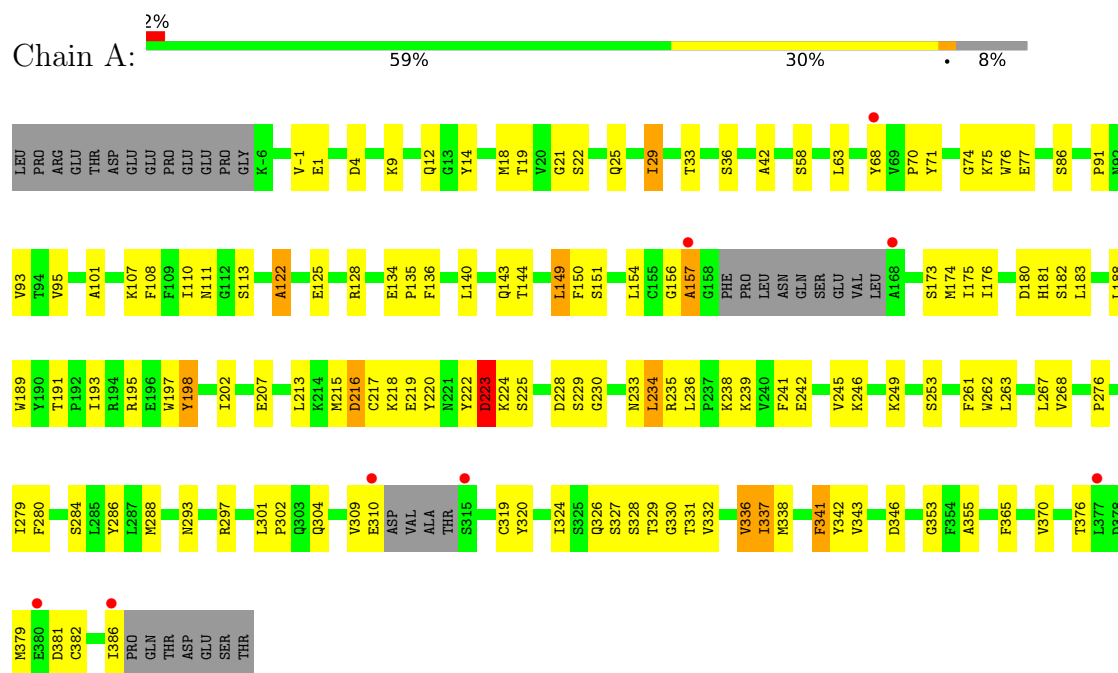
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	81	Total	O	0	0
			81	81		
4	B	64	Total	O	0	0
			64	64		
4	C	39	Total	O	0	0
			39	39		

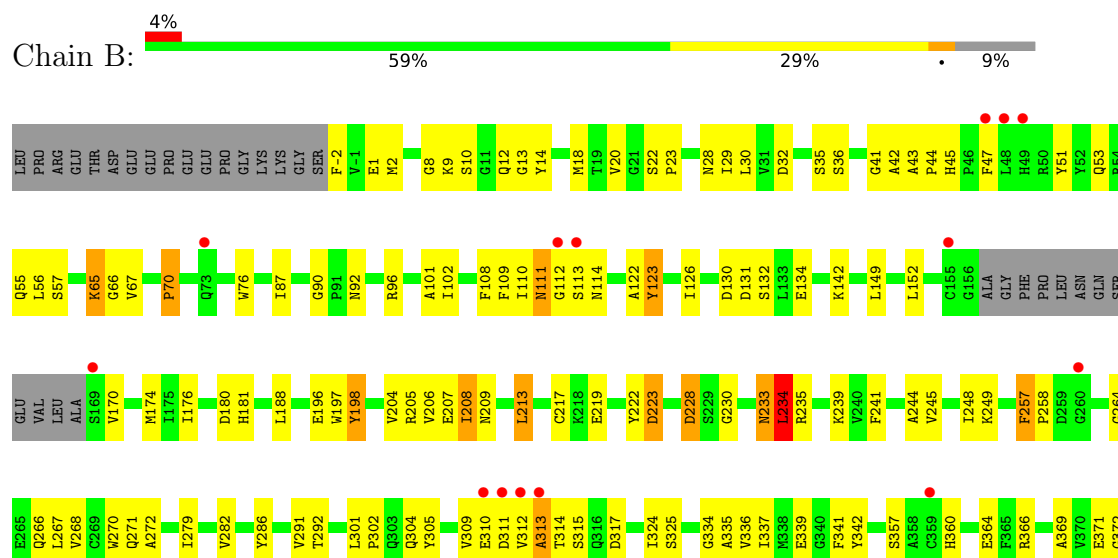
3 Residue-property plots

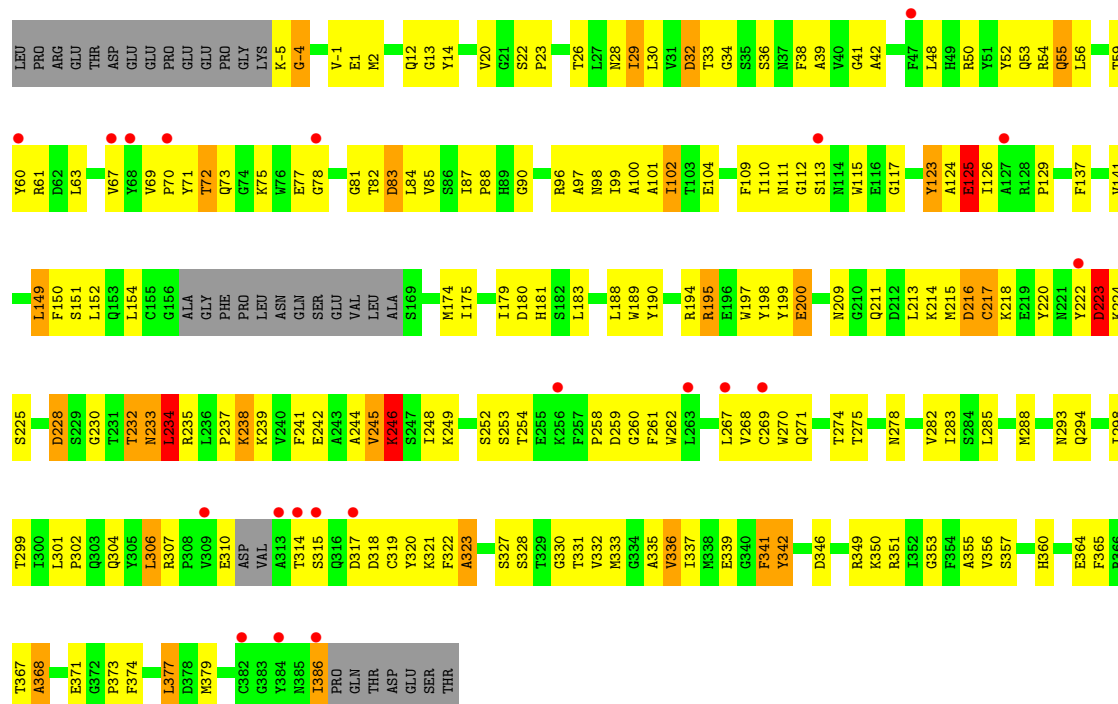
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Beta-secretase 1



• Molecule 1: Beta-secretase 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	224.06Å 105.65Å 63.70Å 90.00° 101.47° 90.00°	Depositor
Resolution (Å)	19.97 – 2.30 19.97 – 2.30	Depositor EDS
% Data completeness (in resolution range)	87.3 (19.97-2.30) 87.1 (19.97-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.245 , 0.304 0.258 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.761	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9210	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: CL, 012

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	0.54	0/3048	1.04	17/4138 (0.4%)
1	B	0.50	0/3031	1.06	16/4119 (0.4%)
1	C	0.50	0/3041	1.04	21/4129 (0.5%)
All	All	0.52	0/9120	1.05	54/12386 (0.4%)

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
1	C	0	1

There are no bond length outliers.

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	198	TYR	N-CA-C	-10.15	95.43	110.46
1	C	198	TYR	N-CA-C	-9.69	95.74	109.69
1	A	198	TYR	N-CA-C	-9.53	95.78	110.17
1	C	223	ASP	N-CA-C	-8.24	93.24	110.80
1	B	213	LEU	N-CA-C	-8.11	102.01	112.23
1	B	233	ASN	N-CA-C	8.01	122.20	110.59
1	A	233	ASN	N-CA-C	7.82	121.15	110.35
1	A	342	TYR	N-CA-C	-7.59	96.36	108.73
1	C	341	PHE	N-CA-C	7.25	119.08	108.86
1	A	216	ASP	N-CA-C	-7.17	100.12	110.24
1	B	204	VAL	N-CA-C	6.79	119.42	112.90
1	C	214	LYS	N-CA-C	6.62	118.66	110.91
1	B	342	TYR	N-CA-C	-6.52	98.60	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	96	ARG	N-CA-C	-6.42	99.18	109.96
1	C	365	PHE	N-CA-C	6.39	121.21	113.41
1	B	123	TYR	N-CA-C	6.38	118.25	110.41
1	A	213	LEU	N-CA-C	-6.24	104.37	112.23
1	C	102	ILE	N-CA-C	6.23	116.21	107.37
1	B	369	ALA	N-CA-C	6.20	118.49	109.07
1	A	193	ILE	N-CA-C	-6.12	98.39	107.51
1	C	228	ASP	N-CA-C	6.09	118.34	107.80
1	C	246	LYS	N-CA-C	-6.06	104.37	110.97
1	B	234	LEU	N-CA-C	-6.05	97.86	108.20
1	C	234	LEU	N-CA-C	-6.03	98.89	108.73
1	B	208	ILE	N-CA-C	-5.99	98.85	107.78
1	A	136	PHE	N-CA-C	5.83	117.31	111.07
1	A	343	VAL	N-CA-C	5.79	115.94	107.37
1	C	278	ASN	N-CA-C	5.75	118.32	111.71
1	A	365	PHE	N-CA-C	5.67	120.33	112.90
1	B	341	PHE	N-CA-C	5.65	118.28	109.52
1	C	216	ASP	N-CA-C	-5.64	101.30	109.59
1	A	223	ASP	N-CA-C	-5.60	98.88	110.80
1	A	341	PHE	N-CA-C	5.57	117.87	109.41
1	A	29	ILE	N-CA-C	5.55	115.89	108.11
1	C	282	VAL	N-CA-C	-5.53	101.77	109.45
1	C	323	ALA	N-CA-C	5.48	119.59	113.02
1	C	328	SER	N-CA-C	-5.42	106.67	113.28
1	A	280	PHE	CA-C-N	5.41	125.40	120.21
1	A	280	PHE	C-N-CA	5.41	125.40	120.21
1	B	228	ASP	N-CA-C	5.35	117.06	107.80
1	A	246	LYS	N-CA-C	-5.32	105.38	111.07
1	C	123	TYR	N-CA-C	5.27	116.89	110.41
1	B	181	HIS	N-CA-C	5.23	119.14	112.34
1	C	29	ILE	N-CA-C	5.17	115.34	108.11
1	C	364	GLU	N-CA-C	5.15	116.90	111.28
1	C	342	TYR	N-CA-C	-5.12	99.07	108.02
1	C	32	ASP	N-CA-C	5.11	116.64	107.80
1	C	38	PHE	N-CA-C	-5.07	100.47	108.73
1	B	23	PRO	CA-C-N	5.05	125.05	119.90
1	B	23	PRO	C-N-CA	5.05	125.05	119.90
1	A	21	GLY	N-CA-C	5.03	120.26	112.81
1	B	13	GLY	N-CA-C	5.03	120.65	112.61
1	A	382	CYS	N-CA-C	-5.02	107.01	113.23
1	C	368	ALA	N-CA-C	-5.02	103.86	110.43

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	52	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	2974	0	2889	94	0
1	B	2956	0	2865	106	0
1	C	2967	0	2886	147	0
2	A	5	0	0	1	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	40	0	39	5	0
3	B	40	0	39	1	0
3	C	40	0	39	6	0
4	A	81	0	0	3	0
4	B	64	0	0	2	0
4	C	39	0	0	2	0
All	All	9210	0	8757	342	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 19.

All (342) 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:B:53:GLN:HB3	1:B:55:GLN:HE22	1.27	0.96
1:C:55:GLN:H	1:C:55:GLN:HE21	1.11	0.95
1:B:234:LEU:HB2	1:B:336:VAL:HG21	1.47	0.94
1:C:55:GLN:HG2	1:C:56:LEU:HD12	1.53	0.91
1:A:4:ASP:H	1:A:173:SER:HB3	1.39	0.88
1:B:271:GLN:CD	1:B:271:GLN:H	1.81	0.88
1:C:225:SER:OG	1:C:331:THR:HB	1.75	0.87
1:B:53:GLN:HB3	1:B:55:GLN:NE2	1.91	0.85
1:C:246:LYS:HA	1:C:246:LYS:HE2	1.56	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:110:ILE:HB	1:C:113:SER:HB3	1.59	0.84
1:C:20:VAL:HG12	1:C:85:VAL:HG22	1.59	0.84
1:B:9:LYS:HD3	1:B:12:GLN:OE1	1.81	0.81
1:A:301:LEU:H	1:A:304:GLN:NE2	1.80	0.80
1:A:301:LEU:H	1:A:304:GLN:HE21	1.25	0.80
1:A:68:TYR:HB2	1:A:77:GLU:HG2	1.67	0.76
1:A:180:ASP:OD2	1:A:182:SER:HB3	1.86	0.76
1:C:125:GLU:HG2	1:C:197:TRP:HB2	1.68	0.75
1:B:55:GLN:HE21	1:B:56:LEU:HG	1.51	0.75
1:C:188:LEU:HD23	1:C:355:ALA:HB2	1.67	0.75
1:C:194:ARG:HB3	1:C:200:GLU:HG2	1.69	0.75
1:B:241:PHE:O	1:B:245:VAL:HG23	1.88	0.74
1:A:238:LYS:O	1:A:242:GLU:HG3	1.87	0.74
1:B:310:GLU:HG3	1:B:311:ASP:N	2.01	0.73
1:B:234:LEU:HB2	1:B:336:VAL:CG2	2.18	0.73
1:B:301:LEU:H	1:B:304:GLN:NE2	1.86	0.73
1:A:110:ILE:HB	1:A:113:SER:HB3	1.71	0.72
1:C:82:THR:O	1:C:83:ASP:HB2	1.89	0.72
1:C:41:GLY:HA2	1:C:102:ILE:HB	1.71	0.72
1:C:55:GLN:H	1:C:55:GLN:NE2	1.86	0.72
1:C:314:THR:HG22	1:C:315:SER:N	2.07	0.70
1:A:241:PHE:O	1:A:245:VAL:HG23	1.90	0.70
1:B:279:ILE:HG22	1:C:211:GLN:HE21	1.56	0.69
1:B:245:VAL:HG12	1:B:249:LYS:HE2	1.74	0.69
1:C:-5:LYS:HG2	1:C:-4:GLY:H	1.57	0.68
1:B:2:MET:HG2	1:B:90:GLY:HA2	1.77	0.67
1:B:310:GLU:HG3	1:B:311:ASP:H	1.60	0.67
1:B:311:ASP:C	1:B:313:ALA:H	2.01	0.67
1:C:149:LEU:HD23	1:C:149:LEU:C	2.20	0.66
1:C:55:GLN:HE21	1:C:55:GLN:N	1.90	0.66
1:C:225:SER:HA	1:C:331:THR:O	1.95	0.66
1:C:111:ASN:HD22	1:C:112:GLY:N	1.94	0.66
1:C:261:PHE:CE1	1:C:322:PHE:HB2	2.32	0.65
1:B:219:GLU:OE1	1:B:239:LYS:HD2	1.96	0.64
1:B:335:ALA:O	1:B:339:GLU:HG3	1.98	0.64
1:C:20:VAL:CG1	1:C:85:VAL:HG22	2.28	0.64
1:A:149:LEU:HD23	1:A:149:LEU:C	2.23	0.64
1:B:385:ASN:O	1:B:386:ILE:HG22	1.98	0.64
1:C:314:THR:HG22	1:C:315:SER:H	1.63	0.64
1:C:-5:LYS:HG2	1:C:-4:GLY:N	2.13	0.63
1:C:222:TYR:O	1:C:223:ASP:HB2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:TYR:CB	1:A:77:GLU:HG2	2.29	0.63
1:C:302:PRO:O	1:C:306:LEU:HB2	1.98	0.63
1:A:276:PRO:O	1:A:279:ILE:HG12	1.98	0.63
1:C:285:LEU:HD12	1:C:333:MET:HE3	1.80	0.63
1:A:74:GLY:HA2	1:A:107:LYS:HB2	1.80	0.62
1:B:241:PHE:HE1	1:B:324:ILE:O	1.81	0.62
1:B:309:VAL:HG23	1:B:309:VAL:O	1.98	0.62
1:A:267:LEU:HD13	1:A:319:CYS:HB3	1.80	0.62
1:C:234:LEU:HB2	1:C:336:VAL:HG21	1.81	0.62
1:B:149:LEU:C	1:B:149:LEU:HD23	2.25	0.61
1:B:386:ILE:HD13	1:B:386:ILE:O	2.00	0.61
1:C:97:ALA:O	1:C:99:ILE:HD12	2.01	0.61
1:A:293:ASN:HA	1:A:376:THR:O	2.01	0.60
1:C:234:LEU:HB2	1:C:336:VAL:CG2	2.31	0.60
1:B:311:ASP:C	1:B:313:ALA:N	2.58	0.60
1:B:14:TYR:CE2	1:B:170:VAL:HB	2.37	0.60
1:B:301:LEU:HB3	1:B:302:PRO:HD2	1.83	0.60
1:A:238:LYS:HG3	1:A:326:GLN:NE2	2.16	0.59
1:B:111:ASN:C	1:B:111:ASN:HD22	2.10	0.59
1:C:137:PHE:O	1:C:141:VAL:HG23	2.03	0.59
1:C:241:PHE:O	1:C:245:VAL:HG23	2.03	0.59
1:A:151:SER:OG	1:A:175:ILE:HB	2.04	0.58
1:C:301:LEU:H	1:C:304:GLN:HE21	1.52	0.58
1:A:218:LYS:HE3	4:A:640:HOH:O	2.04	0.58
1:C:238:LYS:HE2	1:C:242:GLU:OE2	2.03	0.58
1:C:271:GLN:HA	1:C:271:GLN:HE21	1.66	0.58
1:C:285:LEU:CD1	1:C:333:MET:HE3	2.33	0.57
1:A:4:ASP:N	1:A:173:SER:HB3	2.16	0.57
1:C:14:TYR:CG	1:C:154:LEU:HD22	2.40	0.57
1:C:298:ILE:HB	1:C:341:PHE:CZ	2.40	0.57
1:A:14:TYR:CG	1:A:154:LEU:HD22	2.40	0.56
1:B:197:TRP:CD1	1:B:198:TYR:H	2.23	0.56
1:B:206:VAL:HG11	1:B:213:LEU:HD22	1.87	0.56
1:B:267:LEU:HD22	1:B:309:VAL:HG21	1.87	0.56
1:C:110:ILE:HD11	3:C:503:012:H45	1.87	0.56
1:B:197:TRP:CG	1:B:198:TYR:H	2.23	0.56
1:B:205:ARG:HB3	1:B:286:TYR:HB2	1.87	0.56
1:B:65:LYS:HD3	1:B:66:GLY:N	2.21	0.56
1:C:34:GLY:HA3	1:C:228:ASP:OD1	2.05	0.56
1:C:72:THR:HG22	1:C:73:GLN:N	2.20	0.56
1:A:183:LEU:HA	4:A:678:HOH:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179:ILE:HG23	1:C:342:TYR:HE2	1.71	0.56
1:B:14:TYR:O	1:B:30:LEU:HD22	2.06	0.55
1:B:205:ARG:NE	1:B:207:GLU:OE2	2.37	0.55
1:C:83:ASP:O	1:C:96:ARG:HA	2.06	0.55
1:B:228:ASP:OD1	1:B:230:GLY:N	2.34	0.55
1:C:30:LEU:HD21	3:C:503:012:H48A	1.89	0.55
1:A:286:TYR:CZ	1:A:297:ARG:HD3	2.41	0.55
1:C:77:GLU:HB3	1:C:104:GLU:HB2	1.88	0.55
1:B:314:THR:O	1:B:315:SER:HB3	2.07	0.55
1:B:12:GLN:HE21	1:B:110:ILE:HG21	1.72	0.55
1:C:55:GLN:HG2	1:C:56:LEU:CD1	2.34	0.55
1:C:222:TYR:O	1:C:223:ASP:CB	2.54	0.54
1:A:42:ALA:CB	1:A:101:ALA:HB1	2.37	0.54
1:C:14:TYR:CD2	1:C:154:LEU:HD22	2.42	0.54
1:C:53:GLN:HB3	1:C:56:LEU:HD13	1.88	0.54
1:A:180:ASP:HB3	1:A:183:LEU:HG	1.89	0.54
1:B:310:GLU:CG	1:B:311:ASP:H	2.18	0.54
1:C:333:MET:HE2	1:C:337:ILE:HG21	1.89	0.54
1:C:367:THR:HG23	1:C:368:ALA:O	2.08	0.54
1:A:238:LYS:CG	1:A:326:GLN:NE2	2.70	0.54
1:C:33:THR:OG1	1:C:228:ASP:HA	2.08	0.54
1:B:244:ALA:O	1:B:248:ILE:HG13	2.07	0.54
1:C:29:ILE:HD12	1:C:117:GLY:HA3	1.89	0.54
1:B:55:GLN:NE2	1:B:56:LEU:HG	2.20	0.53
1:A:207:GLU:HB2	1:A:284:SER:HB2	1.91	0.53
1:C:268:VAL:O	1:C:319:CYS:HA	2.07	0.53
1:C:217:CYS:HA	1:C:220:TYR:CD1	2.43	0.53
1:C:71:TYR:O	1:C:72:THR:C	2.52	0.53
1:C:314:THR:CG2	1:C:315:SER:N	2.72	0.53
1:C:12:GLN:HG2	1:C:110:ILE:HD13	1.91	0.52
1:C:87:ILE:N	4:C:642:HOH:O	2.40	0.52
1:B:271:GLN:H	1:B:271:GLN:NE2	2.07	0.52
1:B:264:GLY:HA2	4:B:615:HOH:O	2.09	0.52
1:C:237:PRO:HD3	1:C:331:THR:OG1	2.09	0.52
1:C:30:LEU:HD11	3:C:503:012:H48A	1.90	0.52
1:A:149:LEU:HD23	1:A:150:PHE:N	2.25	0.52
1:C:97:ALA:O	1:C:99:ILE:CD1	2.58	0.52
1:C:48:LEU:HD21	1:C:109:PHE:CD2	2.45	0.51
1:C:149:LEU:HD23	1:C:150:PHE:N	2.24	0.51
1:A:217:CYS:HA	1:A:220:TYR:CD1	2.44	0.51
1:A:235:ARG:HB3	1:A:327:SER:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:ARG:HB3	1:B:286:TYR:CG	2.46	0.51
1:B:301:LEU:H	1:B:304:GLN:HE21	1.57	0.51
1:B:42:ALA:CB	1:B:101:ALA:HB1	2.40	0.51
1:C:349:ARG:NH1	1:C:351:ARG:NH2	2.59	0.51
1:B:130:ASP:OD2	1:B:132:SER:HB2	2.10	0.51
1:C:314:THR:CG2	1:C:315:SER:H	2.22	0.51
1:A:225:SER:OG	1:A:331:THR:HB	2.11	0.51
1:C:70:PRO:HA	1:C:75:LYS:HB3	1.92	0.51
1:C:230:GLY:O	3:C:503:012:H48B	2.11	0.51
1:C:151:SER:OG	1:C:175:ILE:HB	2.11	0.51
1:C:307:ARG:HH22	1:C:321:LYS:HE3	1.76	0.51
1:C:195:ARG:HB3	1:C:197:TRP:CD1	2.46	0.50
1:A:234:LEU:HD13	1:A:234:LEU:O	2.11	0.50
1:B:-2:PHE:HB3	1:B:2:MET:HE3	1.93	0.50
1:B:357:SER:HB3	1:B:360:HIS:HB3	1.93	0.50
1:C:386:ILE:HD13	1:C:386:ILE:C	2.36	0.50
1:A:219:GLU:HG3	1:A:239:LYS:HD3	1.93	0.50
1:B:111:ASN:C	1:B:111:ASN:ND2	2.69	0.50
1:C:235:ARG:HB2	1:C:332:VAL:HB	1.93	0.50
1:A:241:PHE:HE1	1:A:324:ILE:O	1.94	0.50
1:B:112:GLY:O	1:B:113:SER:C	2.54	0.50
1:B:180:ASP:C	1:B:180:ASP:OD2	2.55	0.50
1:B:230:GLY:O	3:B:502:012:H48B	2.12	0.50
1:A:36:SER:OG	1:A:122:ALA:HB3	2.12	0.49
1:A:143:GLN:HG3	1:C:1:GLU:OE1	2.11	0.49
1:B:208:ILE:O	1:B:209:ASN:C	2.55	0.49
1:C:288:MET:HE2	1:C:379:MET:HB3	1.94	0.49
1:A:341:PHE:HB3	1:A:355:ALA:O	2.12	0.49
1:B:47:PHE:CD1	1:B:109:PHE:O	2.65	0.49
1:C:36:SER:OG	1:C:123:TYR:O	2.24	0.49
1:B:43:ALA:HB1	1:B:44:PRO:HD2	1.93	0.49
1:C:215:MET:HE3	1:C:239:LYS:NZ	2.27	0.49
1:A:95:VAL:HG11	1:A:140:LEU:HA	1.94	0.49
1:A:228:ASP:OD1	1:A:230:GLY:N	2.42	0.49
1:B:42:ALA:HB2	1:B:101:ALA:HB1	1.94	0.49
1:C:113:SER:HB2	1:C:115:TRP:NE1	2.28	0.49
1:C:274:THR:O	1:C:275:THR:C	2.53	0.49
1:B:271:GLN:CD	1:B:271:GLN:N	2.60	0.49
1:C:213:LEU:HB3	1:C:220:TYR:OH	2.13	0.49
1:A:218:LYS:HG3	1:A:381:ASP:O	2.12	0.49
1:C:32:ASP:OD1	1:C:230:GLY:HA3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:N	1:A:336:VAL:HG11	2.28	0.48
1:A:188:LEU:HD23	1:A:355:ALA:HB2	1.95	0.48
1:A:197:TRP:CG	1:A:198:TYR:H	2.31	0.48
1:A:235:ARG:HB2	1:A:332:VAL:HB	1.95	0.48
1:B:174:MET:SD	1:B:176:ILE:HD11	2.53	0.48
1:A:234:LEU:HB2	1:A:336:VAL:CG1	2.43	0.48
1:A:93:VAL:HG21	1:A:144:THR:HG21	1.95	0.48
1:A:261:PHE:CD1	1:A:268:VAL:HG23	2.48	0.48
1:C:301:LEU:HB3	1:C:302:PRO:HD2	1.95	0.48
1:C:26:THR:HG22	1:C:50:ARG:NH1	2.29	0.48
1:B:28:ASN:O	1:B:29:ILE:HD13	2.14	0.48
1:B:205:ARG:HB3	1:B:286:TYR:CB	2.44	0.47
1:C:59:THR:HG21	1:C:84:LEU:HD12	1.96	0.47
1:A:110:ILE:HD11	3:A:501:012:H45	1.96	0.47
1:C:224:LYS:O	1:C:331:THR:N	2.47	0.47
1:C:22:SER:HA	1:C:23:PRO:C	2.39	0.47
1:A:337:ILE:HG22	1:A:338:MET:N	2.29	0.47
1:B:257:PHE:CB	1:B:258:PRO:HD2	2.44	0.47
1:A:267:LEU:C	1:A:267:LEU:HD12	2.39	0.47
1:A:297:ARG:O	1:A:370:VAL:HA	2.15	0.47
1:B:209:ASN:ND2	1:C:209:ASN:HB3	2.29	0.47
1:B:282:VAL:HG13	1:B:366:ARG:NH1	2.30	0.47
1:B:47:PHE:HD1	1:B:109:PHE:O	1.98	0.47
1:B:268:VAL:HG13	1:B:270:TRP:CZ3	2.49	0.47
1:C:238:LYS:O	1:C:242:GLU:HG3	2.15	0.47
1:A:14:TYR:CZ	1:A:154:LEU:HB3	2.50	0.46
1:A:238:LYS:HA	1:A:326:GLN:HG3	1.96	0.46
1:A:128:ARG:NH2	3:A:501:012:H28B	2.29	0.46
1:B:305:TYR:HB2	1:B:324:ILE:HG13	1.96	0.46
1:C:200:GLU:HA	1:C:225:SER:O	2.16	0.46
1:B:1:GLU:HG3	1:B:2:MET:HG3	1.98	0.46
1:C:216:ASP:OD2	1:C:218:LYS:HB2	2.16	0.46
1:C:244:ALA:O	1:C:248:ILE:HG13	2.15	0.46
1:A:202:ILE:HB	1:A:288:MET:HE3	1.96	0.46
1:B:1:GLU:CG	1:B:2:MET:H	2.29	0.46
1:B:18:MET:HG2	1:B:87:ILE:HG12	1.96	0.46
1:B:257:PHE:N	1:B:257:PHE:CD1	2.84	0.46
1:A:222:TYR:O	1:A:223:ASP:CB	2.63	0.46
1:B:197:TRP:CG	1:B:198:TYR:N	2.84	0.46
1:A:18:MET:SD	1:A:29:ILE:HG13	2.55	0.46
1:B:207:GLU:O	1:B:208:ILE:HD13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:111:ASN:HD22	1:C:111:ASN:C	2.21	0.46
1:A:-1:VAL:HG12	1:A:1:GLU:OE1	2.16	0.45
1:A:293:ASN:HA	1:A:379:MET:HE2	1.98	0.45
1:B:67:VAL:HG13	1:B:76:TRP:HZ3	1.81	0.45
1:C:293:ASN:HA	1:C:379:MET:HE2	1.98	0.45
1:C:216:ASP:O	1:C:218:LYS:N	2.49	0.45
1:C:252:SER:O	1:C:254:THR:N	2.50	0.45
1:C:283:ILE:O	1:C:299:THR:HA	2.16	0.45
1:A:71:TYR:CG	3:A:501:012:HD2	2.52	0.45
1:C:69:VAL:O	1:C:75:LYS:HB2	2.17	0.45
1:C:189:TRP:O	1:C:353:GLY:HA2	2.16	0.45
1:B:41:GLY:HA2	1:B:102:ILE:HB	1.97	0.45
1:C:110:ILE:CD1	3:C:503:012:H45	2.47	0.45
1:A:249:LYS:HD3	1:A:262:TRP:CD1	2.52	0.45
1:A:174:MET:HE3	1:A:176:ILE:HD11	1.97	0.45
1:B:12:GLN:NE2	1:B:110:ILE:HG21	2.31	0.45
1:B:336:VAL:CG2	1:B:337:ILE:N	2.80	0.45
1:A:261:PHE:C	1:A:263:LEU:H	2.24	0.45
1:C:20:VAL:HG12	1:C:85:VAL:HG13	1.99	0.45
1:A:230:GLY:O	3:A:501:012:H48B	2.16	0.45
1:C:152:LEU:HD23	1:C:174:MET:HA	1.98	0.45
1:A:19:THR:OG1	1:A:86:SER:HB2	2.17	0.44
1:A:319:CYS:C	1:A:320:TYR:CD1	2.95	0.44
1:B:22:SER:O	1:B:57:SER:HA	2.17	0.44
1:B:152:LEU:HD23	1:B:174:MET:HA	1.98	0.44
1:B:228:ASP:OD1	1:B:228:ASP:C	2.60	0.44
1:C:233:ASN:HB3	1:C:323:ALA:O	2.17	0.44
1:C:59:THR:HG21	1:C:84:LEU:CD1	2.47	0.44
1:C:180:ASP:HB3	1:C:183:LEU:HD12	1.99	0.44
1:B:257:PHE:HB3	1:B:258:PRO:HD2	1.99	0.44
1:A:110:ILE:HG22	1:A:111:ASN:N	2.33	0.44
1:C:42:ALA:CB	1:C:101:ALA:HB1	2.48	0.44
1:A:301:LEU:N	1:A:304:GLN:HE21	2.05	0.44
1:B:9:LYS:O	1:B:10:SER:C	2.60	0.44
1:C:269:CYS:HA	1:C:318:ASP:O	2.18	0.44
1:B:45:HIS:HE1	1:B:47:PHE:CD1	2.36	0.44
1:B:108:PHE:O	1:B:110:ILE:HD12	2.18	0.44
1:B:233:ASN:HA	1:B:336:VAL:HG13	1.99	0.44
1:B:335:ALA:O	1:B:336:VAL:C	2.60	0.44
1:C:13:GLY:HA3	1:C:30:LEU:HD11	1.99	0.44
1:C:190:TYR:CG	1:C:351:ARG:HD2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:271:GLN:HA	1:C:271:GLN:NE2	2.31	0.43
1:A:63:LEU:O	1:C:181:HIS:HD2	2.00	0.43
1:C:356:VAL:HG23	4:C:612:HOH:O	2.17	0.43
1:A:346:ASP:OD1	1:A:346:ASP:C	2.60	0.43
1:A:75:LYS:HD3	2:A:607:CL:CL	2.54	0.43
1:C:327:SER:OG	1:C:330:GLY:O	2.36	0.43
1:A:191:THR:HA	1:A:288:MET:O	2.19	0.43
1:A:261:PHE:C	1:A:263:LEU:N	2.76	0.43
1:C:28:ASN:N	1:C:50:ARG:HH21	2.16	0.43
1:C:232:THR:HB	1:C:233:ASN:H	1.61	0.43
1:A:197:TRP:CD1	1:A:198:TYR:H	2.36	0.43
1:B:222:TYR:O	1:B:223:ASP:CB	2.66	0.43
1:B:271:GLN:O	1:B:272:ALA:C	2.62	0.43
1:A:22:SER:HB2	1:A:58:SER:OG	2.17	0.43
1:B:291:VAL:HG12	1:B:292:THR:N	2.34	0.43
1:C:26:THR:HG22	1:C:50:ARG:HH12	1.83	0.43
1:C:267:LEU:HB2	1:C:320:TYR:O	2.19	0.43
1:C:293:ASN:OD1	1:C:377:LEU:HA	2.19	0.43
1:A:9:LYS:HB2	1:A:12:GLN:OE1	2.19	0.43
1:A:71:TYR:CD1	3:A:501:012:HD2	2.54	0.43
1:B:266:GLN:HG3	4:B:635:HOH:O	2.19	0.42
1:A:235:ARG:HG2	1:A:235:ARG:HH21	1.82	0.42
1:B:134:GLU:OE1	1:B:142:LYS:HE3	2.19	0.42
1:B:314:THR:O	1:B:315:SER:CB	2.66	0.42
1:C:188:LEU:HD23	1:C:355:ALA:CB	2.44	0.42
1:B:51:TYR:CD1	1:B:51:TYR:N	2.88	0.42
1:C:63:LEU:HG	1:C:81:GLY:HA2	2.00	0.42
1:C:124:ALA:O	1:C:126:ILE:N	2.52	0.42
1:C:249:LYS:HG3	1:C:262:TRP:CE2	2.54	0.42
1:C:270:TRP:O	1:C:271:GLN:C	2.63	0.42
1:A:134:GLU:HA	1:A:135:PRO:HD3	1.91	0.42
1:B:32:ASP:OD1	1:B:230:GLY:HA3	2.19	0.42
1:A:234:LEU:HB2	1:A:336:VAL:HG13	2.01	0.42
1:B:92:ASN:OD1	1:B:92:ASN:O	2.38	0.42
1:C:71:TYR:CG	3:C:503:012:HD2	2.55	0.42
1:C:357:SER:HB3	1:C:360:HIS:HB3	2.01	0.42
1:C:267:LEU:C	1:C:267:LEU:HD12	2.44	0.42
1:A:224:LYS:O	1:A:330:GLY:HA3	2.20	0.42
1:B:8:GLY:C	1:B:170:VAL:HG22	2.45	0.42
1:B:235:ARG:HA	1:B:325:SER:O	2.20	0.42
1:A:215:MET:O	1:A:216:ASP:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:HD13	1:A:236:LEU:HG	2.02	0.42
1:A:320:TYR:CD1	1:A:320:TYR:N	2.88	0.42
1:C:61:ARG:HB2	1:C:82:THR:OG1	2.20	0.42
1:C:216:ASP:C	1:C:218:LYS:H	2.28	0.42
1:C:232:THR:O	1:C:335:ALA:N	2.51	0.42
1:B:123:TYR:CZ	1:B:196:GLU:HG2	2.55	0.42
1:C:67:VAL:HG23	1:C:129:PRO:HG3	2.02	0.42
1:C:237:PRO:O	1:C:238:LYS:C	2.62	0.42
1:C:294:GLN:HG2	1:C:373:PRO:HB2	2.02	0.42
1:C:258:PRO:O	1:C:260:GLY:N	2.54	0.41
1:A:267:LEU:HB2	1:A:320:TYR:O	2.20	0.41
1:C:28:ASN:O	1:C:29:ILE:HD13	2.19	0.41
1:A:14:TYR:CD1	1:A:154:LEU:HD22	2.55	0.41
1:A:70:PRO:HD2	1:A:128:ARG:NH1	2.35	0.41
1:A:19:THR:HA	1:A:25:GLN:O	2.20	0.41
1:B:336:VAL:HG23	1:B:337:ILE:N	2.36	0.41
1:B:364:GLU:OE2	1:C:374:PHE:HB3	2.20	0.41
1:B:372:GLY:HA2	1:B:374:PHE:CE1	2.56	0.41
1:C:149:LEU:C	1:C:149:LEU:CD2	2.89	0.41
1:B:35:SER:O	1:B:122:ALA:HB3	2.21	0.41
1:C:216:ASP:C	1:C:218:LYS:N	2.76	0.41
1:A:76:TRP:HD1	1:A:108:PHE:CE1	2.39	0.41
1:A:157:ALA:HA	4:A:668:HOH:O	2.21	0.41
1:C:335:ALA:O	1:C:339:GLU:HG3	2.21	0.41
1:A:91:PRO:HD3	1:A:176:ILE:HB	2.03	0.41
1:B:36:SER:OG	1:B:126:ILE:HG13	2.21	0.41
1:B:76:TRP:HB2	1:B:102:ILE:HG23	2.02	0.41
1:C:98:ASN:C	1:C:99:ILE:HD12	2.46	0.41
1:A:189:TRP:O	1:A:353:GLY:HA2	2.21	0.41
1:C:54:ARG:HD2	1:C:60:TYR:CZ	2.56	0.41
1:A:202:ILE:HD12	1:A:288:MET:HE1	2.04	0.40
1:C:39:ALA:HB2	1:C:100:ALA:HB3	2.03	0.40
1:B:270:TRP:O	1:B:317:ASP:HB3	2.21	0.40
1:B:309:VAL:O	1:B:309:VAL:CG2	2.66	0.40
1:C:2:MET:HG2	1:C:90:GLY:HA2	2.04	0.40
1:C:123:TYR:HE1	1:C:199:TYR:CE2	2.40	0.40
1:C:77:GLU:N	1:C:104:GLU:O	2.44	0.40
1:C:310:GLU:C	1:C:310:GLU:CD	2.90	0.40
1:A:301:LEU:O	1:A:302:PRO:C	2.63	0.40
1:A:309:VAL:O	1:A:310:GLU:C	2.64	0.40
1:B:42:ALA:O	1:B:51:TYR:HD2	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:252:SER:C	1:C:254:THR:H	2.30	0.40
1:C:288:MET:HE2	1:C:379:MET:CB	2.51	0.40
1:A:33:THR:HG23	1:A:229:SER:OG	2.21	0.40
1:C:78:GLY:HA3	1:C:101:ALA:O	2.22	0.40
1:C:125:GLU:HG2	1:C:197:TRP:CB	2.43	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	373/412 (90%)	341 (91%)	27 (7%)	5 (1%)	9	10
1	B	372/412 (90%)	329 (88%)	35 (9%)	8 (2%)	5	4
1	C	371/412 (90%)	324 (87%)	34 (9%)	13 (4%)	3	1
All	All	1116/1236 (90%)	994 (89%)	96 (9%)	26 (2%)	5	4

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	ASP
1	B	223	ASP
1	C	125	GLU
1	C	223	ASP
1	C	233	ASN
1	B	312	VAL
1	C	72	THR
1	C	238	LYS
1	C	253	SER
1	C	259	ASP
1	B	111	ASN
1	B	313	ALA

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Mol	Chain	Res	Type
1	B	334	GLY
1	C	83	ASP
1	C	217	CYS
1	C	232	THR
1	A	253	SER
1	C	350	LYS
1	A	122	ALA
1	A	157	ALA
1	B	131	ASP
1	B	217	CYS
1	C	-4	GLY
1	A	156	GLY
1	C	346	ASP
1	B	70	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	321/353 (91%)	311 (97%)	10 (3%)	35	52
1	B	319/353 (90%)	310 (97%)	9 (3%)	38	56
1	C	322/353 (91%)	306 (95%)	16 (5%)	22	33
All	All	962/1059 (91%)	927 (96%)	35 (4%)	31	47

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	GLU
1	A	149	LEU
1	A	181	HIS
1	A	195	ARG
1	A	234	LEU
1	A	328	SER
1	A	329	THR
1	A	336	VAL

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Mol	Chain	Res	Type
1	A	337	ILE
1	A	386	ILE
1	B	20	VAL
1	B	65	LYS
1	B	70	PRO
1	B	114	ASN
1	B	188	LEU
1	B	234	LEU
1	B	257	PHE
1	B	371	GLU
1	B	386	ILE
1	C	-1	VAL
1	C	55	GLN
1	C	88	PRO
1	C	125	GLU
1	C	149	LEU
1	C	195	ARG
1	C	200	GLU
1	C	234	LEU
1	C	245	VAL
1	C	246	LYS
1	C	306	LEU
1	C	317	ASP
1	C	336	VAL
1	C	371	GLU
1	C	377	LEU
1	C	386	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	89	HIS
1	A	111	ASN
1	A	304	GLN
1	A	326	GLN
1	A	362	HIS
1	B	28	ASN
1	B	53	GLN
1	B	55	GLN
1	B	73	GLN
1	B	92	ASN

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Mol	Chain	Res	Type
1	B	111	ASN
1	B	304	GLN
1	B	316	GLN
1	C	28	ASN
1	C	55	GLN
1	C	73	GLN
1	C	111	ASN
1	C	181	HIS
1	C	211	GLN
1	C	266	GLN
1	C	271	GLN
1	C	294	GLN
1	C	304	GLN
1	C	326	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 9 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	012	C	503	-	42,43,43	1.59	9 (21%)	53,58,58	2.07	5 (9%)
3	012	B	502	-	42,43,43	1.48	6 (14%)	53,58,58	2.14	6 (11%)
3	012	A	501	-	42,43,43	1.53	5 (11%)	53,58,58	2.07	7 (13%)

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	012	C	503	-	-	9/32/44/44	0/4/4/4
3	012	B	502	-	-	10/32/44/44	0/4/4/4
3	012	A	501	-	-	3/32/44/44	0/4/4/4

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	503	012	C32-N31	4.28	1.41	1.35
3	B	502	012	C32-N31	3.94	1.40	1.35
3	A	501	012	C35-C30	-3.52	1.52	1.55
3	A	501	012	C32-N31	3.41	1.40	1.35
3	C	503	012	C36-N34	3.06	1.51	1.46
3	B	502	012	C36-N34	3.06	1.51	1.46
3	C	503	012	C35-C30	-2.53	1.53	1.55
3	A	501	012	C24-C25	2.44	1.43	1.39
3	A	501	012	C44-C43	2.24	1.42	1.38
3	A	501	012	C36-N34	2.18	1.50	1.46
3	C	503	012	C44-C43	2.15	1.42	1.38
3	C	503	012	C24-C25	2.15	1.43	1.39
3	C	503	012	C42-C41	2.13	1.42	1.39
3	B	502	012	C42-C43	2.11	1.42	1.39
3	C	503	012	CD1-CG	2.11	1.43	1.38
3	B	502	012	C45-C46	2.06	1.42	1.38
3	C	503	012	C22-C21	2.05	1.43	1.38
3	B	502	012	C23-C22	2.05	1.42	1.38
3	B	502	012	CE1-CD1	2.03	1.42	1.38
3	C	503	012	C42-C43	2.01	1.42	1.39

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	012	C35-N34-C32	-9.80	98.98	111.51
3	C	503	012	N31-C32-N34	9.36	117.45	108.32
3	A	501	012	N31-C32-N34	9.20	117.30	108.32
3	B	502	012	N31-C32-N34	8.70	116.81	108.32
3	C	503	012	C35-N34-C32	-8.68	100.41	111.51
3	A	501	012	C35-N34-C32	-8.56	100.56	111.51
3	B	502	012	C41-C36-N34	4.96	120.94	113.15
3	C	503	012	C41-C36-N34	4.42	120.10	113.15
3	A	501	012	C41-C36-N34	4.22	119.78	113.15
3	A	501	012	O33-C32-N31	-3.80	119.94	126.84
3	C	503	012	O33-C32-N31	-3.75	120.03	126.84
3	B	502	012	O33-C32-N31	-3.58	120.34	126.84
3	A	501	012	C36-N34-C35	-2.85	117.61	121.76
3	C	503	012	C36-N34-C35	-2.51	118.11	121.76
3	A	501	012	O-C-N	2.37	127.20	122.96
3	B	502	012	C36-N34-C35	-2.32	118.38	121.76
3	B	502	012	CB-CA-N	-2.12	107.07	110.08
3	A	501	012	C30-C-N	-2.10	112.15	116.63

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	012	OG-C1-C2-N2
3	C	503	012	OG-C1-C2-N2
3	B	502	012	C24-C25-N27-C29
3	B	502	012	C26-C25-N27-C28
3	B	502	012	C24-C25-N27-C28
3	B	502	012	C26-C25-N27-C29
3	A	501	012	C44-C43-O47-C48
3	B	502	012	C42-C43-O47-C48
3	C	503	012	C44-C43-O47-C48
3	C	503	012	C24-C25-N27-C29
3	C	503	012	C26-C25-N27-C28
3	A	501	012	C42-C43-O47-C48
3	C	503	012	C42-C43-O47-C48
3	B	502	012	C44-C43-O47-C48
3	C	503	012	C24-C25-N27-C28
3	C	503	012	C26-C25-N27-C29
3	B	502	012	CA-C1-C2-N2
3	C	503	012	CA-C1-C2-N2
3	B	502	012	C41-C36-N34-C35
3	C	503	012	C41-C36-N34-C35

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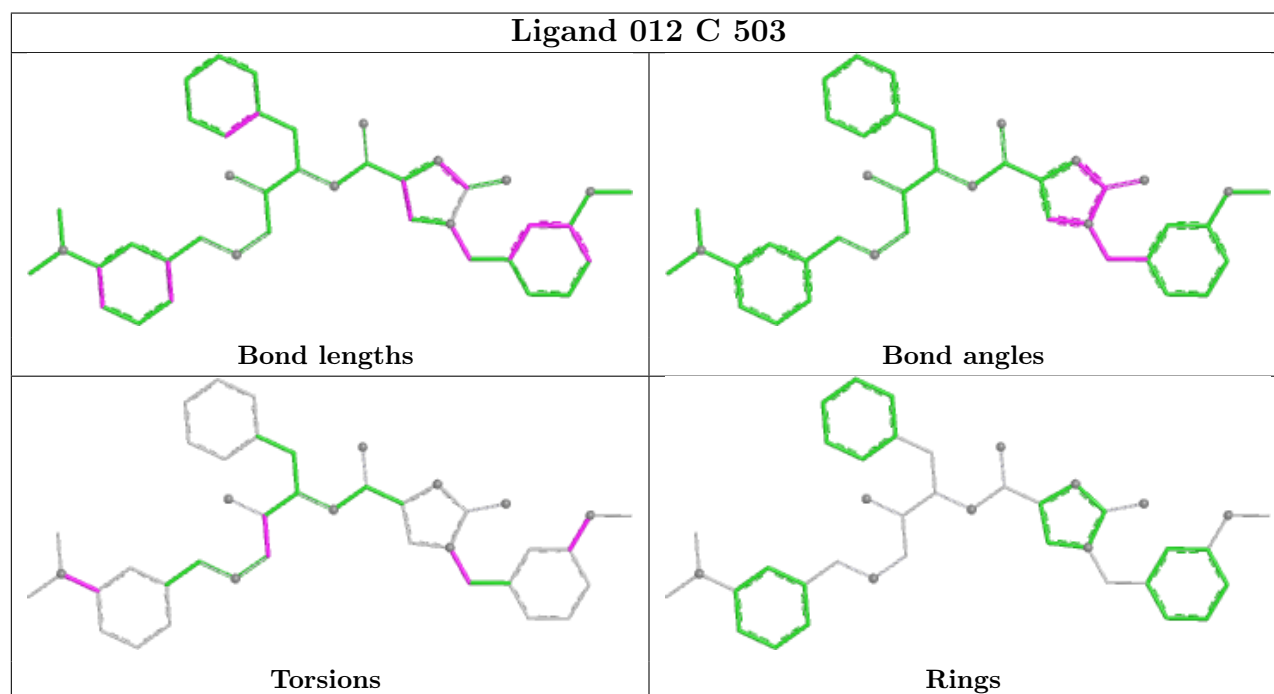
Mol	Chain	Res	Type	Atoms
3	B	502	012	O-C-C30-N31
3	A	501	012	C41-C36-N34-C35

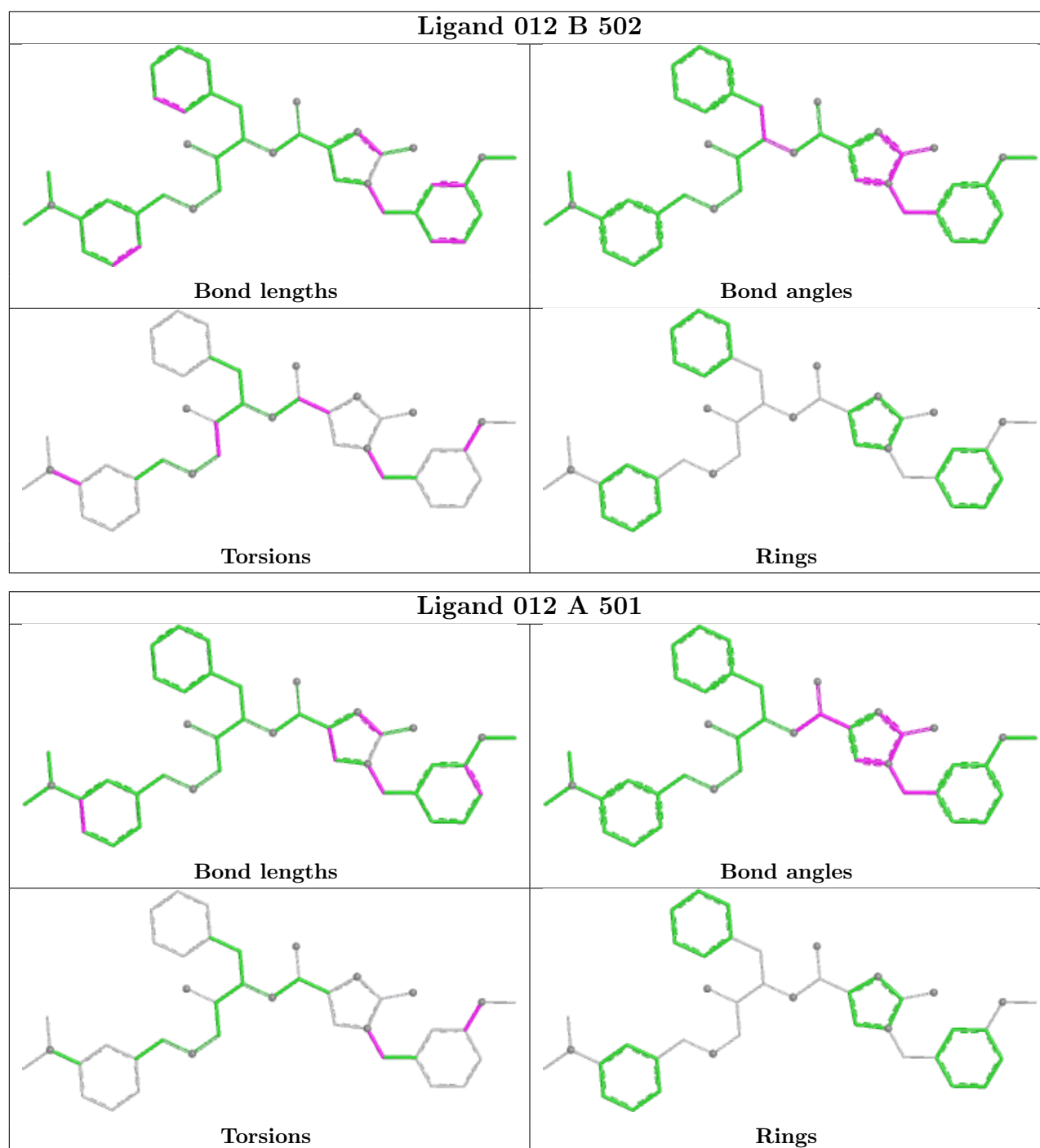
There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	503	012	6	0
3	B	502	012	1	0
3	A	501	012	5	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	379/412 (91%)	0.11	8 (2%) 63 65	18, 35, 61, 91	0
1	B	376/412 (91%)	0.34	16 (4%) 40 41	22, 40, 68, 89	0
1	C	377/412 (91%)	0.69	21 (5%) 30 32	25, 50, 74, 105	0
All	All	1132/1236 (91%)	0.38	45 (3%) 42 44	18, 41, 69, 105	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	386	ILE	4.7
1	A	168	ALA	4.4
1	B	311	ASP	4.3
1	B	312	VAL	4.2
1	C	386	ILE	4.2
1	C	384	TYR	3.8
1	B	113	SER	3.7
1	A	386	ILE	3.7
1	A	157	ALA	3.6
1	B	313	ALA	3.3
1	B	112	GLY	3.3
1	C	313	ALA	3.2
1	C	68	TYR	3.1
1	A	315	SER	3.0
1	C	309	VAL	2.9
1	C	127	ALA	2.8
1	B	359	CYS	2.8
1	C	317	ASP	2.7
1	C	67	VAL	2.5
1	B	384	TYR	2.5
1	C	263	LEU	2.5
1	B	260	GLY	2.5
1	C	222	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	269	CYS	2.4
1	B	48	LEU	2.3
1	C	314	THR	2.3
1	A	68	TYR	2.2
1	B	310	GLU	2.2
1	C	78	GLY	2.2
1	B	169	SER	2.2
1	A	377	LEU	2.2
1	C	267	LEU	2.2
1	C	47	PHE	2.2
1	A	380	GLU	2.2
1	C	113	SER	2.2
1	A	310	GLU	2.1
1	C	256	LYS	2.1
1	C	60	TYR	2.1
1	B	47	PHE	2.1
1	C	315	SER	2.1
1	B	73	GLN	2.1
1	C	70	PRO	2.1
1	B	49	HIS	2.1
1	B	155	CYS	2.0
1	C	382	CYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	012	C	503	40/40	0.89	0.12	40,45,58,58	0

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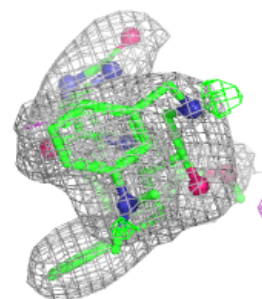
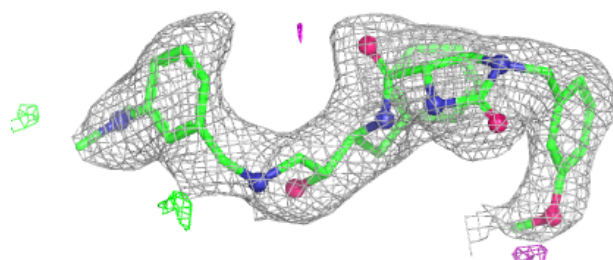
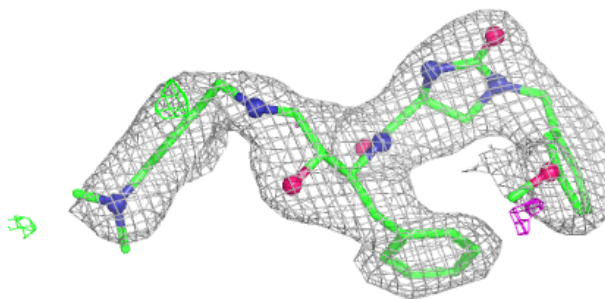
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	012	B	502	40/40	0.90	0.11	36,43,50,50	0
3	012	A	501	40/40	0.93	0.09	30,37,45,45	0
2	CL	A	608	1/1	0.94	0.05	39,39,39,39	0
2	CL	B	602	1/1	0.95	0.12	23,23,23,23	0
2	CL	C	609	1/1	0.97	0.07	43,43,43,43	0
2	CL	C	604	1/1	0.98	0.08	32,32,32,32	0
2	CL	A	607	1/1	0.98	0.06	34,34,34,34	0
2	CL	A	601	1/1	0.98	0.16	12,12,12,12	0
2	CL	A	605	1/1	0.98	0.11	32,32,32,32	0
2	CL	B	606	1/1	0.98	0.06	36,36,36,36	0
2	CL	A	603	1/1	0.99	0.09	21,21,21,21	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.

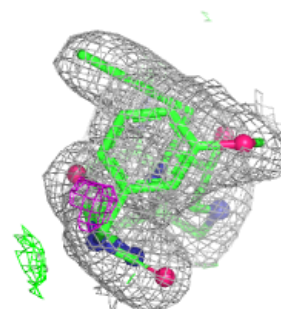
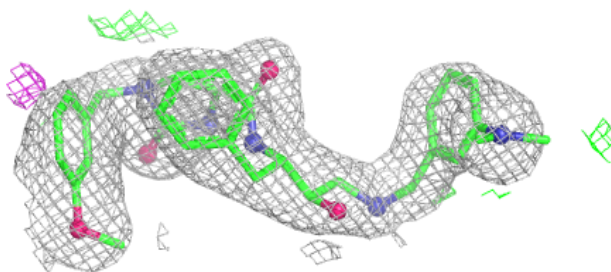
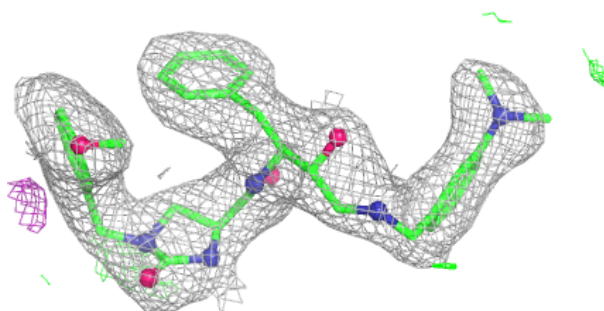
Electron density around 012 C 503:

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

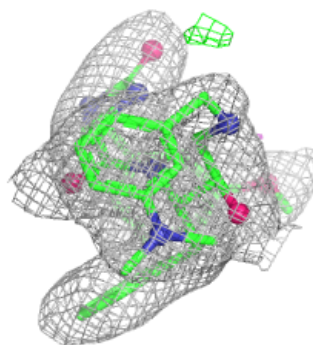
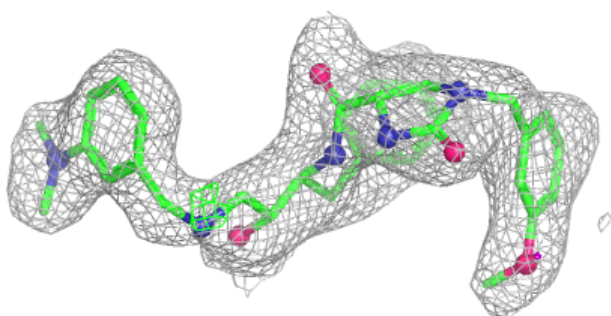
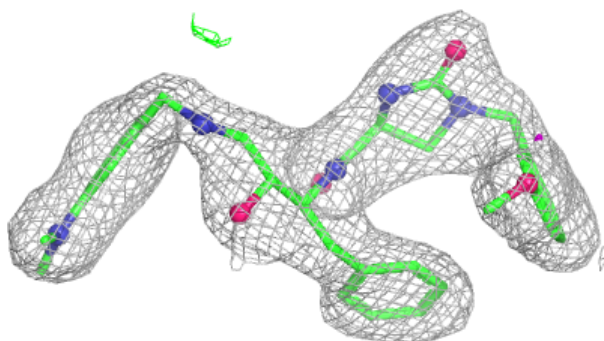


Electron density around 012 B 502:

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

**Electron density around 012 A 501:**

$2mF_o-DF_c$ (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.