



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:14 AM UTC

PDB ID : 4LL0 / pdb_00004ll0
Title : EGFR L858R/T790M in complex with PD168393
Authors : Yun, C.H.; Eck, M.J.
Deposited on : 2013-07-09
Resolution : 4.00 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

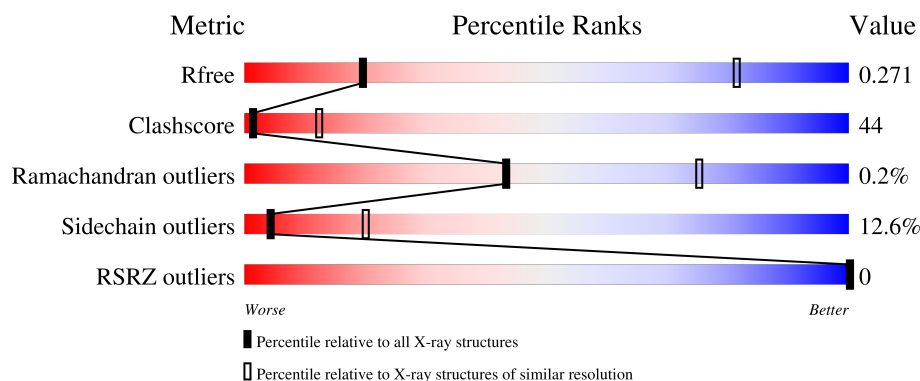
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

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	331	
1	B	331	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	YUN	A	1101	-	-	X	-
2	YUN	B	1101	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4552 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.

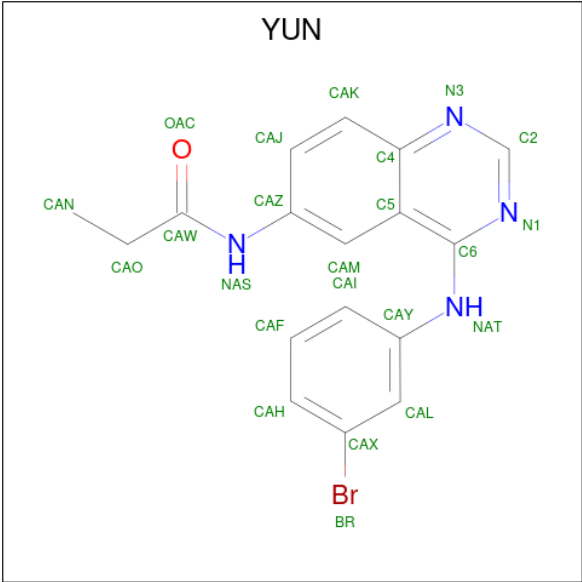
- Molecule 1 is a protein called Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			2287	1471	383	418	15			
1	B	281	Total	C	N	O	S	0	0	0
			2219	1434	371	399	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	692	GLY	-	expression tag	UNP P00533
A	693	SER	-	expression tag	UNP P00533
A	790	MET	THR	engineered mutation	UNP P00533
A	858	ARG	LEU	engineered mutation	UNP P00533
B	692	GLY	-	expression tag	UNP P00533
B	693	SER	-	expression tag	UNP P00533
B	790	MET	THR	engineered mutation	UNP P00533
B	858	ARG	LEU	engineered mutation	UNP P00533

- Molecule 2 is N-{4-[(3-bromophenyl)amino]quinazolin-6-yl}propanamide (CCD ID: YUN) (formula: C₁₇H₁₅BrN₄O).

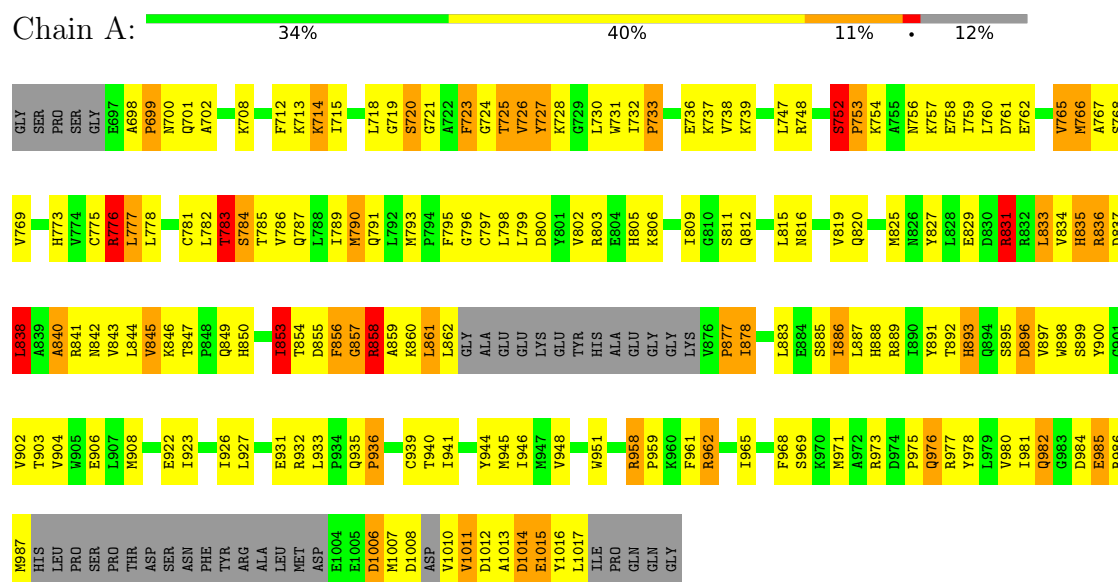


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	Br	C	N	O	0	0
			23	1	17	4	1		
2	B	1	Total	Br	C	N	O	0	0
			23	1	17	4	1		

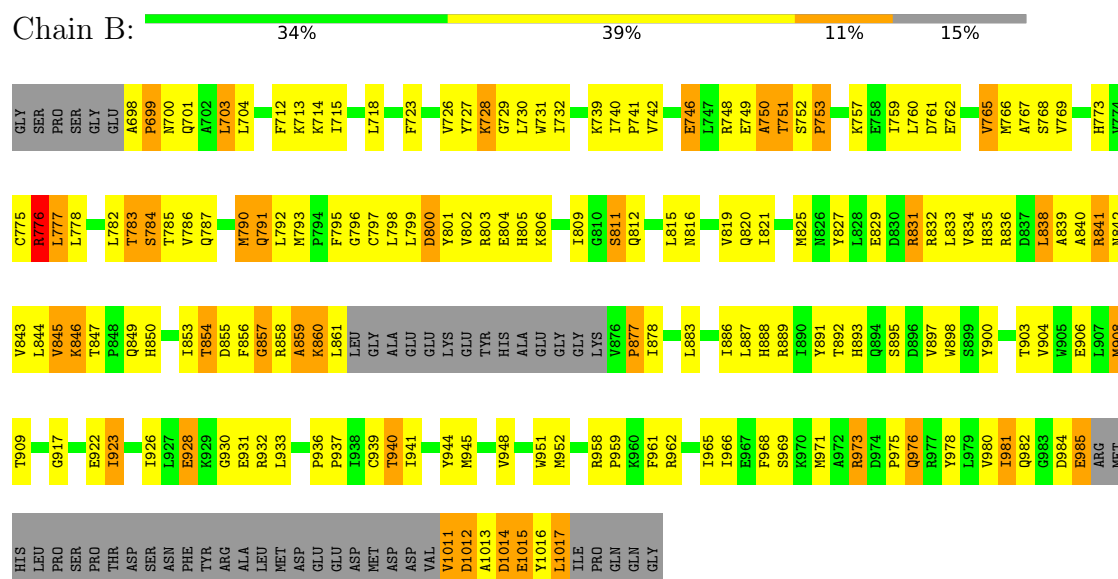
3 Residue-property plots

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: Epidermal growth factor receptor



• Molecule 1: Epidermal growth factor receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.64Å 72.71Å 167.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.60 – 4.00 49.60 – 4.00	Depositor EDS
% Data completeness (in resolution range)	95.8 (49.60-4.00) 95.7 (49.60-4.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 4.00Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.238 , 0.257 0.228 , 0.271	Depositor DCC
R_{free} test set	334 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	200.5	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 256.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4552	wwPDB-VP
Average B, all atoms (Å ²)	257.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% 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: YUN

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	0.92	3/2334 (0.1%)	1.45	47/3163 (1.5%)
1	B	0.87	0/2267	1.39	39/3073 (1.3%)
All	All	0.89	3/4601 (0.1%)	1.42	86/6236 (1.4%)

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	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	858	ARG	N-CA	-8.65	1.35	1.46
1	A	861	LEU	CA-C	-6.43	1.44	1.52
1	A	835	HIS	CG-ND1	-5.92	1.31	1.38

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	857	GLY	N-CA-C	-13.97	95.78	112.83
1	A	858	ARG	N-CA-C	11.65	123.72	111.14
1	A	783	THR	N-CA-C	-11.37	94.66	110.35
1	A	857	GLY	CA-C-N	-9.83	107.07	120.44
1	A	857	GLY	C-N-CA	-9.83	107.07	120.44
1	B	783	THR	N-CA-C	-9.57	97.14	110.35
1	A	896	ASP	N-CA-C	-8.86	100.79	111.69
1	B	765	VAL	N-CA-C	8.85	118.84	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	940	THR	N-CA-C	-8.61	98.11	110.59
1	A	861	LEU	N-CA-C	8.45	123.04	108.76
1	B	757	LYS	N-CA-C	-8.01	102.77	112.54
1	B	750	ALA	N-CA-C	-7.80	99.24	110.24
1	B	857	GLY	N-CA-C	-7.79	102.92	113.27
1	A	831	ARG	NE-CZ-NH2	7.69	126.12	119.20
1	B	940	THR	N-CA-C	-7.68	99.45	110.59
1	B	748	ARG	N-CA-C	-7.67	99.11	110.23
1	A	831	ARG	NE-CZ-NH1	-7.61	113.89	121.50
1	B	728	LYS	N-CA-C	-7.50	96.68	108.90
1	B	776	ARG	NE-CZ-NH1	-7.46	114.04	121.50
1	B	776	ARG	NE-CZ-NH2	7.37	125.83	119.20
1	A	892	THR	N-CA-C	-7.37	98.72	109.18
1	B	768	SER	N-CA-C	7.26	121.74	113.02
1	A	765	VAL	N-CA-C	7.20	117.99	110.72
1	B	854	THR	N-CA-C	7.17	124.53	112.99
1	A	728	LYS	N-CA-C	-7.14	97.27	108.90
1	B	1014	ASP	N-CA-C	-7.04	103.61	111.28
1	A	878	ILE	N-CA-C	6.98	117.77	110.72
1	A	768	SER	N-CA-C	6.93	121.34	113.02
1	A	936	PRO	N-CA-C	-6.92	102.25	110.70
1	A	1014	ASP	N-CA-C	6.88	118.78	111.28
1	B	892	THR	N-CA-C	-6.84	99.47	109.18
1	A	853	ILE	CB-CA-C	-6.42	102.41	111.28
1	B	791	GLN	N-CA-C	-6.31	101.77	110.35
1	A	776	ARG	NE-CZ-NH2	-6.28	113.55	119.20
1	B	784	SER	N-CA-C	-6.18	104.64	111.82
1	A	962	ARG	CD-NE-CZ	6.14	132.99	124.40
1	A	721	GLY	N-CA-C	-6.10	100.24	111.04
1	A	720	SER	N-CA-C	6.07	119.14	108.75
1	B	936	PRO	N-CA-C	-6.05	104.66	110.47
1	A	753	PRO	N-CA-C	-6.04	105.13	113.53
1	B	878	ILE	N-CA-C	5.94	116.72	110.72
1	A	841	ARG	NE-CZ-NH2	5.92	124.53	119.20
1	A	962	ARG	NE-CZ-NH2	-5.91	113.88	119.20
1	A	831	ARG	CD-NE-CZ	5.88	132.63	124.40
1	B	753	PRO	N-CA-C	5.88	122.51	113.75
1	A	886	ILE	N-CA-C	5.83	116.48	110.36
1	A	841	ARG	NE-CZ-NH1	-5.75	115.75	121.50
1	A	893	HIS	N-CA-C	-5.74	105.16	111.82
1	A	836	ARG	N-CA-CB	-5.72	107.65	114.17
1	B	773	HIS	N-CA-C	5.64	120.16	112.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	840	ALA	N-CA-C	5.64	117.88	111.11
1	B	859	ALA	N-CA-C	-5.62	105.04	112.24
1	A	757	LYS	N-CA-C	-5.62	105.69	112.54
1	A	897	VAL	CB-CA-C	-5.60	104.70	112.04
1	B	843	VAL	N-CA-C	-5.60	100.38	108.89
1	A	766	MET	N-CA-C	-5.59	105.28	111.71
1	B	917	GLY	CA-C-N	-5.58	118.57	122.59
1	B	917	GLY	C-N-CA	-5.58	118.57	122.59
1	B	928	GLU	N-CA-C	-5.55	105.13	111.07
1	B	798	LEU	N-CA-C	-5.54	105.33	111.71
1	B	895	SER	N-CA-C	-5.52	106.04	112.89
1	A	798	LEU	N-CA-C	-5.51	105.43	111.82
1	B	923	ILE	N-CA-C	5.51	116.24	110.62
1	A	843	VAL	N-CA-C	-5.48	100.56	108.89
1	B	751	THR	N-CA-C	5.46	119.51	112.41
1	A	958	ARG	CA-C-N	5.46	125.22	119.76
1	A	958	ARG	C-N-CA	5.46	125.22	119.76
1	A	838	LEU	N-CA-C	5.44	118.11	109.96
1	A	773	HIS	N-CA-C	5.42	119.87	112.88
1	A	985	GLU	N-CA-C	5.42	118.30	110.28
1	A	784	SER	N-CA-C	-5.29	105.69	111.82
1	B	746	GLU	N-CA-C	-5.27	101.94	110.32
1	B	846	LYS	N-CA-C	-5.24	105.48	111.14
1	B	909	THR	N-CA-C	-5.23	105.63	113.89
1	B	804	GLU	N-CA-C	5.20	117.85	111.82
1	B	841	ARG	NE-CZ-NH2	-5.17	114.55	119.20
1	A	980	VAL	N-CA-C	5.15	114.16	106.85
1	A	752	SER	N-CA-C	-5.07	102.05	109.50
1	B	831	ARG	N-CA-C	-5.05	106.96	113.02
1	B	897	VAL	CB-CA-C	-5.04	104.69	112.05
1	A	727	TYR	CA-C-N	-5.04	115.88	122.99
1	A	727	TYR	C-N-CA	-5.04	115.88	122.99
1	B	841	ARG	CD-NE-CZ	5.04	131.46	124.40
1	B	1015	GLU	N-CA-C	-5.03	105.80	111.28
1	A	733	PRO	N-CA-C	-5.02	103.14	111.11
1	B	800	ASP	N-CA-C	5.01	116.55	111.14

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	856	PHE	Peptide

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	2287	0	2278	220	0
1	B	2219	0	2241	184	0
2	A	23	0	14	12	0
2	B	23	0	14	7	0
All	All	4552	0	4547	404	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 44.

All (404) 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:836:ARG:CB	1:B:860:LYS:HE2	1.60	1.32
1:B:860:LYS:HE3	1:B:891:TYR:CD2	1.69	1.27
1:B:836:ARG:CB	1:B:860:LYS:CE	2.25	1.14
1:A:982:GLN:HA	1:A:982:GLN:OE1	1.44	1.13
1:A:860:LYS:NZ	1:A:891:TYR:HD2	1.46	1.12
1:B:698:ALA:CB	1:B:699:PRO:CD	2.29	1.10
1:A:833:LEU:HD23	1:A:834:VAL:N	1.66	1.10
1:B:698:ALA:HB1	1:B:699:PRO:HD3	1.26	1.08
1:B:698:ALA:CB	1:B:699:PRO:HD3	1.88	1.00
1:A:724:GLY:O	1:A:725:THR:HG22	1.60	1.00
1:B:1012:ASP:HB2	1:B:1015:GLU:HG3	1.45	0.98
1:A:860:LYS:NZ	1:A:891:TYR:CD2	2.23	0.98
1:B:860:LYS:CE	1:B:891:TYR:CD2	2.49	0.96
1:A:718:LEU:HB2	1:A:726:VAL:CG1	1.95	0.96
1:A:853:ILE:CD1	1:A:853:ILE:N	2.30	0.95
1:A:853:ILE:N	1:A:853:ILE:HD12	1.80	0.94
1:A:726:VAL:CG1	1:A:727:TYR:N	2.30	0.94
1:B:833:LEU:HG	1:B:861:LEU:HD13	1.48	0.93
1:A:1011:VAL:CG1	1:A:1015:GLU:HG2	1.98	0.93
1:A:708:LYS:HB3	1:B:931:GLU:HB2	1.51	0.92
2:A:1101:YUN:N1	2:A:1101:YUN:H1	1.84	0.91
2:B:1101:YUN:H1	2:B:1101:YUN:N1	1.84	0.91
1:A:1011:VAL:HG11	1:A:1015:GLU:HG2	1.50	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:LEU:CB	1:A:726:VAL:HG11	2.01	0.90
1:B:795:PHE:HB2	1:B:845:VAL:HB	1.54	0.90
1:B:698:ALA:HB3	1:B:699:PRO:CD	2.02	0.89
1:A:739:LYS:O	1:A:1010:VAL:HA	1.73	0.89
1:A:753:PRO:CD	1:A:754:LYS:H	1.86	0.88
1:B:1016:TYR:O	1:B:1017:LEU:HD23	1.74	0.88
1:A:833:LEU:HD23	1:A:833:LEU:C	1.96	0.87
1:B:698:ALA:HB1	1:B:699:PRO:CD	1.98	0.87
1:B:698:ALA:HB3	1:B:699:PRO:HD2	1.56	0.86
2:A:1101:YUN:H6	2:A:1101:YUN:OAC	1.77	0.85
2:B:1101:YUN:OAC	2:B:1101:YUN:H6	1.76	0.85
1:B:923:ILE:HA	1:B:926:ILE:HD12	1.59	0.85
1:A:726:VAL:O	1:A:727:TYR:CD1	2.29	0.85
1:A:795:PHE:HB2	1:A:845:VAL:HB	1.59	0.84
1:A:923:ILE:HA	1:A:926:ILE:HD12	1.59	0.84
1:A:712:PHE:HB3	1:A:731:TRP:HA	1.59	0.84
1:A:854:THR:HG23	1:A:854:THR:O	1.79	0.82
1:A:718:LEU:HB2	1:A:726:VAL:HG12	1.62	0.81
1:B:712:PHE:HB3	1:B:731:TRP:HA	1.60	0.81
1:B:793:MET:H	2:B:1101:YUN:H8	1.46	0.81
1:A:718:LEU:HB2	1:A:726:VAL:HG11	1.59	0.81
1:A:724:GLY:C	1:A:725:THR:CG2	2.54	0.80
1:A:1015:GLU:CA	1:A:1015:GLU:OE2	2.30	0.80
1:B:797:CYS:H	2:B:1101:YUN:H15	1.45	0.80
1:A:726:VAL:HG12	1:A:727:TYR:N	1.96	0.79
1:A:723:PHE:N	1:A:723:PHE:CD1	2.44	0.79
1:B:1016:TYR:O	1:B:1017:LEU:CD2	2.30	0.79
1:B:1016:TYR:O	1:B:1017:LEU:CG	2.30	0.79
1:A:835:HIS:CE1	1:A:855:ASP:O	2.35	0.79
1:A:797:CYS:H	2:A:1101:YUN:H15	1.47	0.79
1:A:753:PRO:HD2	1:A:754:LYS:H	1.46	0.78
1:A:936:PRO:HG2	1:A:939:CYS:SG	2.22	0.78
1:A:877:PRO:HG2	1:A:877:PRO:O	1.81	0.78
1:A:833:LEU:C	1:A:833:LEU:CD2	2.56	0.78
1:B:833:LEU:HD23	1:B:834:VAL:N	1.97	0.78
1:B:856:PHE:HD1	1:B:859:ALA:CB	1.97	0.77
1:B:883:LEU:HD12	1:B:886:ILE:HD12	1.66	0.77
1:A:723:PHE:HB2	1:A:748:ARG:CB	2.15	0.77
1:A:860:LYS:HZ1	1:A:891:TYR:HD2	0.80	0.77
1:A:753:PRO:CG	1:A:754:LYS:N	2.48	0.77
1:B:877:PRO:O	1:B:877:PRO:HG2	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1012:ASP:HB2	1:B:1015:GLU:CG	2.16	0.76
1:A:853:ILE:CD1	1:A:853:ILE:H	1.97	0.75
1:A:718:LEU:HD22	2:A:1101:YUN:CAO	2.16	0.75
1:A:723:PHE:H	1:A:723:PHE:HD1	1.34	0.74
1:B:984:ASP:OD1	1:B:984:ASP:C	2.30	0.74
1:A:853:ILE:H	1:A:853:ILE:HD13	1.52	0.74
1:A:726:VAL:HG13	1:A:727:TYR:N	2.01	0.74
1:A:793:MET:H	2:A:1101:YUN:H8	1.51	0.74
1:B:836:ARG:CB	1:B:860:LYS:NZ	2.51	0.74
1:A:718:LEU:CB	1:A:726:VAL:CG1	2.64	0.74
2:B:1101:YUN:N1	2:B:1101:YUN:CAL	2.47	0.73
1:B:836:ARG:CB	1:B:860:LYS:HZ3	2.03	0.72
1:A:753:PRO:HG2	1:A:754:LYS:N	2.05	0.72
1:B:1016:TYR:O	1:B:1017:LEU:HG	1.89	0.72
1:B:860:LYS:HE3	1:B:891:TYR:HD2	1.49	0.72
1:A:715:ILE:HD11	1:A:730:LEU:HG	1.72	0.71
1:B:715:ILE:HD11	1:B:730:LEU:HG	1.73	0.71
1:A:724:GLY:O	1:A:725:THR:CG2	2.39	0.71
1:A:753:PRO:CD	1:A:754:LYS:N	2.52	0.71
1:A:883:LEU:HD12	1:A:886:ILE:HD12	1.70	0.71
1:A:724:GLY:C	1:A:725:THR:HG22	2.16	0.71
1:A:753:PRO:CG	1:A:754:LYS:H	2.02	0.71
1:A:718:LEU:HB3	1:A:726:VAL:HG11	1.72	0.71
1:B:802:VAL:O	1:B:806:LYS:HB3	1.91	0.70
1:A:1016:TYR:O	1:A:1017:LEU:C	2.34	0.70
2:A:1101:YUN:N1	2:A:1101:YUN:CAL	2.47	0.70
1:A:1015:GLU:OE2	1:A:1015:GLU:HA	1.90	0.69
1:B:819:VAL:HG22	1:B:968:PHE:HB3	1.74	0.69
1:A:856:PHE:O	1:A:857:GLY:C	2.35	0.69
1:A:976:GLN:NE2	1:A:984:ASP:OD2	2.26	0.69
1:A:738:VAL:CG2	1:A:1011:VAL:HG23	2.24	0.68
1:A:1011:VAL:HG11	1:A:1015:GLU:CG	2.21	0.68
1:A:736:GLU:CD	1:A:1016:TYR:HH	2.02	0.68
1:B:803:ARG:O	1:B:806:LYS:HG2	1.93	0.68
1:A:819:VAL:HG22	1:A:968:PHE:HB3	1.76	0.68
1:A:877:PRO:O	1:A:877:PRO:CG	2.42	0.68
1:A:835:HIS:O	1:A:836:ARG:CB	2.43	0.67
1:A:783:THR:OG1	1:A:787:GLN:NE2	2.27	0.67
1:A:856:PHE:HB3	1:A:859:ALA:HB2	1.76	0.66
1:B:778:LEU:HD11	1:B:791:GLN:HG2	1.76	0.66
1:A:982:GLN:OE1	1:A:982:GLN:CA	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:833:LEU:HD23	1:A:834:VAL:H	1.58	0.66
1:A:802:VAL:O	1:A:806:LYS:HB3	1.96	0.66
1:A:968:PHE:HA	1:A:971:MET:HE3	1.78	0.66
1:A:803:ARG:O	1:A:806:LYS:HG2	1.96	0.65
1:A:985:GLU:OE1	1:A:985:GLU:N	2.30	0.65
1:A:726:VAL:HG13	1:A:727:TYR:H	1.61	0.65
1:A:856:PHE:O	1:A:859:ALA:N	2.30	0.65
1:A:791:GLN:CG	1:A:1013:ALA:HB2	2.27	0.65
1:A:1015:GLU:OE2	1:A:1015:GLU:N	2.30	0.65
1:B:858:ARG:HD2	1:B:860:LYS:HZ2	1.61	0.64
1:B:877:PRO:O	1:B:877:PRO:CG	2.45	0.64
1:B:749:GLU:O	1:B:750:ALA:C	2.41	0.64
1:A:777:LEU:HA	1:A:790:MET:HG3	1.80	0.64
1:B:835:HIS:CE1	1:B:855:ASP:O	2.52	0.63
1:B:856:PHE:HB3	1:B:859:ALA:HB2	1.80	0.63
1:A:815:LEU:O	1:A:819:VAL:HG23	1.99	0.63
1:A:790:MET:HE1	1:A:854:THR:OG1	1.98	0.63
1:B:831:ARG:C	1:B:832:ARG:HG2	2.24	0.63
1:A:791:GLN:CD	1:A:1013:ALA:HB3	2.23	0.62
1:A:805:HIS:O	1:A:809:ILE:HG13	2.00	0.62
1:B:860:LYS:CE	1:B:891:TYR:HD2	2.09	0.62
1:B:968:PHE:HA	1:B:971:MET:HE3	1.82	0.62
1:B:856:PHE:HD1	1:B:859:ALA:HB2	1.64	0.62
1:A:791:GLN:CD	1:A:1013:ALA:CB	2.74	0.61
1:A:791:GLN:HB2	1:A:1013:ALA:HB2	1.82	0.61
1:A:833:LEU:CD2	1:A:834:VAL:N	2.54	0.61
1:A:933:LEU:HB2	1:A:951:TRP:CH2	2.35	0.61
1:A:1011:VAL:HG12	1:A:1015:GLU:HG2	1.81	0.61
1:A:855:ASP:C	1:A:856:PHE:HD2	2.08	0.61
1:B:856:PHE:CD1	1:B:859:ALA:HB2	2.35	0.61
1:A:738:VAL:HG21	1:A:1011:VAL:HG23	1.83	0.61
1:B:712:PHE:CB	1:B:731:TRP:HA	2.31	0.61
1:A:856:PHE:O	1:A:859:ALA:HB2	2.01	0.60
1:A:898:TRP:CD1	1:A:898:TRP:C	2.79	0.60
1:A:844:LEU:CD1	1:A:854:THR:HG21	2.31	0.60
1:A:1014:ASP:OD2	1:A:1014:ASP:N	2.33	0.60
1:B:858:ARG:O	1:B:859:ALA:HB3	2.01	0.60
1:B:903:THR:O	1:B:906:GLU:HB2	2.02	0.59
1:A:835:HIS:NE2	1:A:855:ASP:O	2.35	0.59
1:A:856:PHE:O	1:A:858:ARG:N	2.36	0.59
1:B:815:LEU:O	1:B:819:VAL:HG23	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:819:VAL:HG22	1:B:968:PHE:CB	2.33	0.59
1:B:751:THR:HG23	1:B:751:THR:O	2.02	0.59
1:B:856:PHE:HD1	1:B:859:ALA:HB1	1.65	0.59
1:B:718:LEU:HD11	1:B:728:LYS:HB2	1.85	0.59
1:B:829:GLU:HA	1:B:893:HIS:CE1	2.38	0.58
1:A:723:PHE:CD1	1:A:723:PHE:O	2.56	0.58
1:B:699:PRO:CG	1:B:699:PRO:O	2.51	0.58
1:B:777:LEU:HA	1:B:790:MET:HG3	1.85	0.58
1:A:908:MET:HG2	1:A:939:CYS:SG	2.43	0.58
1:B:844:LEU:CD1	1:B:854:THR:HG21	2.33	0.58
1:A:753:PRO:HG2	1:A:754:LYS:H	1.65	0.58
1:B:860:LYS:NZ	1:B:891:TYR:CE2	2.72	0.58
1:A:718:LEU:HD22	2:A:1101:YUN:H12	1.86	0.58
1:B:898:TRP:CD1	1:B:898:TRP:C	2.81	0.58
1:B:805:HIS:O	1:B:809:ILE:HG13	2.02	0.58
1:B:933:LEU:HB2	1:B:951:TRP:CH2	2.38	0.57
1:A:712:PHE:CB	1:A:731:TRP:HA	2.32	0.57
1:A:931:GLU:O	1:A:932:ARG:HG2	2.05	0.57
1:B:699:PRO:CD	1:B:699:PRO:O	2.52	0.56
1:A:753:PRO:O	1:A:754:LYS:C	2.48	0.56
1:A:923:ILE:HD13	1:A:926:ILE:HD12	1.86	0.56
1:B:833:LEU:HG	1:B:861:LEU:CD1	2.30	0.56
1:A:778:LEU:HD11	1:A:791:GLN:HB2	1.87	0.56
1:A:736:GLU:O	1:A:738:VAL:HG12	2.06	0.56
1:A:853:ILE:HG22	1:A:854:THR:N	2.20	0.56
1:A:922:GLU:O	1:A:926:ILE:HG13	2.06	0.56
1:B:746:GLU:HA	1:B:787:GLN:HG2	1.87	0.56
1:A:724:GLY:C	1:A:725:THR:HG23	2.30	0.56
1:B:833:LEU:HD23	1:B:833:LEU:C	2.30	0.55
1:B:762:GLU:OE1	1:B:766:MET:HE3	2.06	0.55
1:B:795:PHE:CE1	1:B:847:THR:HA	2.40	0.55
1:A:726:VAL:O	1:A:727:TYR:CG	2.60	0.55
1:A:829:GLU:HA	1:A:893:HIS:CE1	2.42	0.55
1:B:776:ARG:O	1:B:776:ARG:HG3	2.05	0.55
1:B:976:GLN:NE2	1:B:984:ASP:OD2	2.40	0.55
2:B:1101:YUN:OAC	2:B:1101:YUN:CAM	2.46	0.55
1:A:736:GLU:CD	1:A:1016:TYR:OH	2.49	0.55
1:B:860:LYS:CD	1:B:891:TYR:HD2	2.20	0.55
1:A:819:VAL:HG22	1:A:968:PHE:CB	2.37	0.55
1:B:769:VAL:HB	1:B:827:TYR:HE2	1.70	0.54
1:B:846:LYS:HD3	1:B:850:HIS:ND1	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:LEU:CD2	2:A:1101:YUN:H11	2.36	0.54
1:A:723:PHE:N	1:A:723:PHE:HD1	1.95	0.54
1:A:714:LYS:HD3	1:A:727:TYR:CD2	2.42	0.54
1:A:775:CYS:SG	1:A:854:THR:HB	2.47	0.54
2:A:1101:YUN:OAC	2:A:1101:YUN:CAM	2.46	0.54
1:B:847:THR:C	1:B:849:GLN:H	2.15	0.54
1:B:703:LEU:HD23	1:B:703:LEU:H	1.73	0.54
1:B:1012:ASP:CB	1:B:1015:GLU:HB2	2.38	0.54
1:A:718:LEU:HD13	2:A:1101:YUN:CAW	2.38	0.54
1:A:766:MET:HB3	1:A:777:LEU:HB2	1.89	0.54
1:B:699:PRO:O	1:B:699:PRO:HG2	2.06	0.54
1:B:944:TYR:O	1:B:948:VAL:HG23	2.07	0.54
1:A:795:PHE:CE1	1:A:847:THR:HA	2.43	0.53
1:B:969:SER:O	1:B:973:ARG:HG2	2.09	0.53
1:A:791:GLN:CB	1:A:1013:ALA:HB2	2.39	0.53
1:B:726:VAL:HG12	1:B:727:TYR:N	2.23	0.53
1:A:847:THR:C	1:A:849:GLN:H	2.16	0.53
1:A:903:THR:O	1:A:906:GLU:HB2	2.09	0.53
1:A:723:PHE:CB	1:A:748:ARG:CB	2.86	0.52
1:B:858:ARG:HB3	1:B:860:LYS:HG3	1.90	0.52
1:B:860:LYS:CE	1:B:891:TYR:CE2	2.91	0.52
1:A:723:PHE:HD1	1:A:723:PHE:O	1.93	0.52
1:A:1011:VAL:CG1	1:A:1015:GLU:CG	2.79	0.52
1:A:718:LEU:CD2	2:A:1101:YUN:CAO	2.87	0.52
1:B:887:LEU:HB3	1:B:888:HIS:CE1	2.44	0.52
1:A:844:LEU:HD11	1:A:854:THR:HG21	1.91	0.52
1:B:700:ASN:O	1:B:700:ASN:CG	2.52	0.52
1:B:767:ALA:HB2	1:B:777:LEU:HD23	1.91	0.52
1:B:1012:ASP:CB	1:B:1015:GLU:HG3	2.31	0.52
1:B:829:GLU:HG3	1:B:893:HIS:CD2	2.45	0.52
1:B:923:ILE:HD13	1:B:926:ILE:HD12	1.92	0.52
1:A:944:TYR:O	1:A:948:VAL:HG23	2.10	0.52
1:B:856:PHE:O	1:B:857:GLY:C	2.52	0.52
1:B:856:PHE:CB	1:B:859:ALA:HB2	2.40	0.52
1:B:766:MET:HB3	1:B:777:LEU:HB2	1.91	0.51
1:A:1007:MET:O	1:A:1008:ASP:C	2.54	0.51
1:A:702:ALA:O	1:B:941:ILE:HB	2.11	0.51
1:A:782:LEU:O	1:B:930:GLY:HA2	2.10	0.51
1:B:858:ARG:HD2	1:B:860:LYS:NZ	2.25	0.51
1:B:1012:ASP:HB3	1:B:1015:GLU:H	1.75	0.51
1:A:837:ASP:OD2	1:A:858:ARG:HD3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:856:PHE:C	1:A:858:ARG:N	2.57	0.51
1:A:847:THR:C	1:A:849:GLN:N	2.67	0.51
1:A:767:ALA:HB2	1:A:777:LEU:HD23	1.93	0.51
1:B:740:ILE:HG21	1:B:1013:ALA:HB2	1.93	0.51
1:B:835:HIS:NE2	1:B:855:ASP:O	2.43	0.51
1:A:860:LYS:NZ	1:A:860:LYS:HB2	2.27	0.50
1:B:833:LEU:HA	1:B:861:LEU:HD12	1.93	0.50
1:B:860:LYS:HD3	1:B:891:TYR:HD2	1.77	0.50
1:A:782:LEU:HD21	1:B:952:MET:CE	2.41	0.50
1:B:984:ASP:OD1	1:B:984:ASP:O	2.29	0.50
1:B:1012:ASP:HB2	1:B:1015:GLU:HB2	1.93	0.50
1:A:782:LEU:HD21	1:B:952:MET:HE2	1.92	0.50
1:A:986:ARG:O	1:A:987:MET:C	2.54	0.50
1:A:736:GLU:OE2	1:A:1016:TYR:OH	2.29	0.50
1:A:1014:ASP:O	1:A:1015:GLU:C	2.54	0.50
1:B:700:ASN:O	1:B:700:ASN:OD1	2.29	0.50
1:B:701:GLN:O	1:B:701:GLN:OE1	2.30	0.50
1:A:1012:ASP:OD1	1:A:1014:ASP:OD2	2.30	0.50
1:B:829:GLU:HG3	1:B:893:HIS:CG	2.46	0.50
1:B:860:LYS:HE3	1:B:891:TYR:CE2	2.39	0.50
1:A:856:PHE:O	1:A:859:ALA:CB	2.60	0.50
1:B:847:THR:C	1:B:849:GLN:N	2.66	0.50
1:A:1011:VAL:HG11	1:A:1015:GLU:CB	2.42	0.50
1:B:775:CYS:SG	1:B:854:THR:HB	2.52	0.50
1:A:698:ALA:HB1	1:A:699:PRO:HD2	1.93	0.49
1:A:1008:ASP:HB2	1:A:1010:VAL:HG23	1.93	0.49
1:B:1012:ASP:O	1:B:1015:GLU:HB2	2.11	0.49
1:A:736:GLU:OE1	1:A:1016:TYR:OH	2.30	0.49
1:B:847:THR:O	1:B:849:GLN:N	2.46	0.49
1:A:829:GLU:HG3	1:A:893:HIS:CD2	2.48	0.49
1:A:986:ARG:O	1:A:987:MET:O	2.30	0.49
1:B:855:ASP:OD2	1:B:856:PHE:N	2.45	0.49
1:A:778:LEU:HG	1:A:790:MET:HA	1.93	0.49
1:A:931:GLU:C	1:A:932:ARG:HG2	2.37	0.49
1:A:887:LEU:HB3	1:A:888:HIS:CE1	2.48	0.49
1:B:796:GLY:O	1:B:845:VAL:HG23	2.12	0.49
1:A:769:VAL:HB	1:A:827:TYR:HE2	1.78	0.49
1:A:825:MET:HE2	1:A:838:LEU:HD22	1.95	0.48
1:A:701:GLN:CD	1:B:980:VAL:HG21	2.38	0.48
1:A:714:LYS:HD3	1:A:727:TYR:HD2	1.79	0.48
1:B:728:LYS:HE2	1:B:792:LEU:HD21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1012:ASP:HB2	1:B:1015:GLU:CB	2.43	0.48
1:B:700:ASN:OD1	1:B:700:ASN:C	2.55	0.48
1:B:908:MET:HG2	1:B:939:CYS:SG	2.53	0.48
1:B:790:MET:HE1	1:B:854:THR:OG1	2.13	0.48
1:A:747:LEU:CB	1:A:759:ILE:HD12	2.43	0.48
1:A:840:ALA:C	1:A:842:ASN:H	2.21	0.48
1:B:825:MET:HE2	1:B:838:LEU:HD22	1.95	0.48
1:B:883:LEU:HD12	1:B:883:LEU:HA	1.68	0.48
1:A:762:GLU:OE1	1:A:766:MET:HE3	2.13	0.47
1:B:740:ILE:CG2	1:B:1013:ALA:HB2	2.44	0.47
1:B:704:LEU:HD12	1:B:704:LEU:HA	1.16	0.47
1:B:799:LEU:O	1:B:802:VAL:HG22	2.15	0.47
1:A:969:SER:O	1:A:973:ARG:HG2	2.13	0.47
1:B:778:LEU:HD11	1:B:791:GLN:CG	2.44	0.47
1:A:797:CYS:HB2	1:A:800:ASP:OD1	2.14	0.47
1:A:837:ASP:HB2	1:A:858:ARG:CG	2.43	0.47
1:B:797:CYS:HB2	1:B:800:ASP:OD1	2.13	0.47
1:A:776:ARG:O	1:A:776:ARG:HG3	2.15	0.47
1:B:985:GLU:OE1	1:B:985:GLU:O	2.33	0.47
1:A:827:TYR:O	1:A:831:ARG:HG3	2.15	0.46
1:A:941:ILE:HG12	1:A:945:MET:HE2	1.96	0.46
1:A:853:ILE:CG2	1:A:854:THR:N	2.78	0.46
1:B:922:GLU:O	1:B:926:ILE:HG13	2.15	0.46
1:A:898:TRP:CD1	1:A:898:TRP:O	2.68	0.46
1:B:844:LEU:HD11	1:B:854:THR:HG21	1.97	0.46
1:A:756:ASN:HA	1:A:759:ILE:HG22	1.96	0.46
1:A:791:GLN:HB2	1:A:1013:ALA:CB	2.45	0.46
1:A:846:LYS:HD3	1:A:850:HIS:ND1	2.30	0.46
1:B:752:SER:HB2	1:B:753:PRO:CD	2.45	0.46
1:A:799:LEU:O	1:A:802:VAL:HG22	2.15	0.46
1:A:854:THR:O	1:A:854:THR:CG2	2.48	0.46
1:A:971:MET:HA	1:A:978:TYR:CE1	2.50	0.46
1:B:840:ALA:C	1:B:842:ASN:H	2.23	0.46
1:A:961:PHE:O	1:A:965:ILE:HG13	2.16	0.46
1:A:738:VAL:HG23	1:A:1011:VAL:HG23	1.98	0.46
1:B:856:PHE:O	1:B:858:ARG:C	2.59	0.46
1:A:855:ASP:C	1:A:856:PHE:CD2	2.91	0.46
1:A:837:ASP:HB2	1:A:858:ARG:HG2	1.98	0.45
1:B:769:VAL:HG23	1:B:769:VAL:O	2.16	0.45
1:A:789:ILE:HG22	1:A:790:MET:N	2.30	0.45
1:B:931:GLU:C	1:B:932:ARG:HG2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:985:GLU:OE1	1:B:985:GLU:CA	2.63	0.45
1:A:885:SER:HB3	1:A:891:TYR:HE1	1.81	0.45
1:A:984:ASP:HB3	1:A:985:GLU:OE1	2.16	0.45
1:B:854:THR:OG1	1:B:854:THR:O	2.26	0.45
1:B:860:LYS:HG3	1:B:860:LYS:H	1.55	0.45
1:B:1016:TYR:C	1:B:1017:LEU:HD23	2.41	0.45
1:A:829:GLU:HG3	1:A:893:HIS:CG	2.52	0.45
1:B:833:LEU:C	1:B:833:LEU:CD2	2.90	0.45
1:A:736:GLU:O	1:A:738:VAL:N	2.50	0.44
1:A:760:LEU:HD23	1:A:760:LEU:HA	1.79	0.44
1:B:858:ARG:HB3	1:B:860:LYS:CG	2.48	0.44
1:A:719:GLY:O	1:A:720:SER:OG	2.30	0.44
1:A:789:ILE:HD12	1:A:789:ILE:N	2.32	0.44
1:A:812:GLN:CD	1:A:975:PRO:HG3	2.43	0.44
1:B:742:VAL:O	1:B:792:LEU:HB2	2.17	0.44
1:B:1013:ALA:O	1:B:1016:TYR:N	2.49	0.44
1:A:718:LEU:HD21	2:A:1101:YUN:H11	2.00	0.44
1:A:986:ARG:H	1:A:986:ARG:HG3	1.40	0.44
1:A:761:ASP:O	1:A:765:VAL:HG23	2.17	0.44
1:A:923:ILE:HD13	1:A:926:ILE:CD1	2.47	0.44
1:A:733:PRO:HG2	1:A:1016:TYR:HE2	1.82	0.44
1:A:847:THR:O	1:A:849:GLN:N	2.51	0.44
1:A:860:LYS:HE2	1:A:891:TYR:HE2	1.82	0.43
1:A:888:HIS:O	1:A:889:ARG:C	2.57	0.43
1:B:739:LYS:O	1:B:1011:VAL:HG22	2.18	0.43
1:B:928:GLU:C	1:B:930:GLY:N	2.76	0.43
1:A:708:LYS:HD3	1:B:931:GLU:OE2	2.18	0.43
1:A:753:PRO:O	1:A:756:ASN:N	2.51	0.43
1:A:796:GLY:O	1:A:845:VAL:HG23	2.17	0.43
1:B:838:LEU:O	1:B:839:ALA:HB2	2.19	0.43
1:B:941:ILE:HG12	1:B:945:MET:HE2	2.01	0.43
1:B:816:ASN:O	1:B:820:GLN:HG3	2.18	0.43
1:B:790:MET:O	1:B:791:GLN:C	2.60	0.43
1:A:752:SER:O	1:A:753:PRO:C	2.62	0.43
1:A:899:SER:O	1:A:900:TYR:C	2.60	0.43
1:B:856:PHE:O	1:B:859:ALA:HB2	2.19	0.43
1:A:789:ILE:CG2	1:A:790:MET:N	2.82	0.43
1:B:1012:ASP:CB	1:B:1015:GLU:CG	2.91	0.43
1:A:883:LEU:HD12	1:A:883:LEU:HA	1.72	0.43
1:A:958:ARG:HA	1:A:959:PRO:HD3	1.85	0.43
1:B:898:TRP:CD1	1:B:898:TRP:O	2.72	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:747:LEU:CB	1:A:759:ILE:CD1	2.97	0.42
1:B:726:VAL:CG1	1:B:727:TYR:N	2.83	0.42
1:B:759:ILE:HG23	1:B:760:LEU:N	2.33	0.42
1:A:736:GLU:O	1:A:737:LYS:C	2.62	0.42
1:B:844:LEU:HD12	1:B:854:THR:HG21	1.99	0.42
1:B:761:ASP:O	1:B:765:VAL:HG23	2.20	0.42
1:B:888:HIS:O	1:B:889:ARG:C	2.62	0.42
1:A:816:ASN:O	1:A:820:GLN:HG3	2.20	0.42
1:B:812:GLN:NE2	1:B:975:PRO:HG3	2.34	0.42
1:B:793:MET:N	2:B:1101:YUN:H8	2.25	0.42
1:B:931:GLU:O	1:B:932:ARG:HG2	2.19	0.42
1:A:791:GLN:HG3	1:A:1013:ALA:HB2	2.00	0.42
1:A:726:VAL:HG12	1:A:727:TYR:CA	2.50	0.42
1:B:971:MET:HA	1:B:978:TYR:CE1	2.55	0.42
1:B:937:PRO:C	1:B:939:CYS:H	2.28	0.42
1:B:778:LEU:HG	1:B:790:MET:HA	2.02	0.41
1:B:861:LEU:O	1:B:861:LEU:HG	2.19	0.41
1:B:900:TYR:O	1:B:904:VAL:HG23	2.20	0.41
1:B:856:PHE:O	1:B:858:ARG:N	2.53	0.41
1:A:754:LYS:HE2	1:A:758:GLU:OE2	2.20	0.41
1:A:781:CYS:C	1:A:782:LEU:HD12	2.45	0.41
1:A:977:ARG:O	1:A:977:ARG:HG2	2.20	0.41
1:B:958:ARG:HA	1:B:959:PRO:HD3	1.89	0.41
1:B:961:PHE:O	1:B:965:ILE:HG13	2.20	0.41
1:A:791:GLN:CD	1:A:1013:ALA:HB2	2.44	0.41
1:A:902:VAL:HG13	1:A:933:LEU:HD11	2.03	0.41
1:A:718:LEU:HA	1:A:718:LEU:HD23	1.51	0.41
1:A:861:LEU:O	1:A:862:LEU:CB	2.65	0.41
1:B:740:ILE:N	1:B:740:ILE:HD12	2.36	0.41
1:B:811:SER:N	1:B:981:ILE:HD11	2.36	0.41
1:B:821:ILE:HG23	1:B:853:ILE:HD11	2.03	0.41
1:B:928:GLU:O	1:B:930:GLY:N	2.53	0.41
1:B:760:LEU:HA	1:B:760:LEU:HD23	1.74	0.41
1:A:700:ASN:O	1:B:940:THR:HB	2.21	0.41
1:B:797:CYS:HA	1:B:844:LEU:HD23	2.01	0.41
1:B:858:ARG:C	1:B:860:LYS:N	2.75	0.41
1:B:928:GLU:C	1:B:930:GLY:H	2.29	0.41
1:A:769:VAL:O	1:A:769:VAL:HG23	2.21	0.41
1:A:896:ASP:O	1:A:899:SER:N	2.54	0.41
1:B:718:LEU:HA	1:B:718:LEU:HD23	1.78	0.41
1:B:856:PHE:O	1:B:859:ALA:CB	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:946:ILE:HD13	1:A:971:MET:HE1	2.03	0.41
1:A:1011:VAL:CG1	1:A:1015:GLU:HB2	2.51	0.41
1:B:740:ILE:HA	1:B:741:PRO:HD3	1.94	0.41
1:A:878:ILE:HA	1:A:878:ILE:HD13	1.80	0.40
1:B:715:ILE:HG13	1:B:729:GLY:HA2	2.02	0.40
1:A:927:LEU:CD2	1:A:932:ARG:HD3	2.51	0.40
1:A:935:GLN:HA	1:A:936:PRO:HD3	1.87	0.40
1:A:1011:VAL:CG1	1:A:1015:GLU:CB	2.99	0.40
1:A:900:TYR:O	1:A:904:VAL:HG23	2.21	0.40
1:B:801:TYR:CD1	1:B:801:TYR:C	2.99	0.40
1:B:805:HIS:ND1	1:B:805:HIS:N	2.70	0.40
1:B:962:ARG:O	1:B:966:ILE:HD13	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	283/331 (86%)	261 (92%)	21 (7%)	1 (0%)	30	65
1	B	275/331 (83%)	256 (93%)	19 (7%)	0	100	100
All	All	558/662 (84%)	517 (93%)	40 (7%)	1 (0%)	43	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1006	ASP

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	245/290 (84%)	214 (87%)	31 (13%)	4	20
1	B	240/290 (83%)	210 (88%)	30 (12%)	4	20
All	All	485/580 (84%)	424 (87%)	61 (13%)	4	20

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	699	PRO
1	A	713	LYS
1	A	714	LYS
1	A	723	PHE
1	A	725	THR
1	A	726	VAL
1	A	732	ILE
1	A	752	SER
1	A	776	ARG
1	A	777	LEU
1	A	783	THR
1	A	784	SER
1	A	785	THR
1	A	786	VAL
1	A	790	MET
1	A	811	SER
1	A	831	ARG
1	A	833	LEU
1	A	838	LEU
1	A	845	VAL
1	A	853	ILE
1	A	858	ARG
1	A	877	PRO
1	A	895	SER
1	A	962	ARG
1	A	976	GLN
1	A	981	ILE

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Mol	Chain	Res	Type
1	A	982	GLN
1	A	1006	ASP
1	A	1011	VAL
1	A	1015	GLU
1	B	699	PRO
1	B	703	LEU
1	B	713	LYS
1	B	714	LYS
1	B	723	PHE
1	B	732	ILE
1	B	776	ARG
1	B	777	LEU
1	B	782	LEU
1	B	783	THR
1	B	784	SER
1	B	785	THR
1	B	786	VAL
1	B	790	MET
1	B	811	SER
1	B	838	LEU
1	B	841	ARG
1	B	845	VAL
1	B	860	LYS
1	B	877	PRO
1	B	908	MET
1	B	973	ARG
1	B	976	GLN
1	B	981	ILE
1	B	982	GLN
1	B	985	GLU
1	B	1011	VAL
1	B	1012	ASP
1	B	1014	ASP
1	B	1017	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	773	HIS
1	A	808	ASN
1	A	893	HIS
1	A	976	GLN

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Mol	Chain	Res	Type
1	B	701	GLN
1	B	756	ASN
1	B	773	HIS
1	B	808	ASN
1	B	893	HIS
1	B	976	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](#)

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	YUN	A	1101	1	25,25,25	1.64	3 (12%)	33,34,34	2.31	9 (27%)
2	YUN	B	1101	1	25,25,25	1.61	3 (12%)	33,34,34	2.32	10 (30%)

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	YUN	A	1101	1	-	0/10/10/10	0/3/3/3
2	YUN	B	1101	1	-	0/10/10/10	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1101	YUN	C6-C5	-4.46	1.39	1.44
2	B	1101	YUN	CAZ-NAS	-4.34	1.32	1.41
2	A	1101	YUN	CAZ-NAS	-4.25	1.33	1.41
2	B	1101	YUN	C6-C5	-4.08	1.39	1.44
2	B	1101	YUN	CAY-NAT	-3.34	1.33	1.40
2	A	1101	YUN	CAY-NAT	-3.17	1.33	1.40

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1101	YUN	C2-N1-C6	6.36	121.54	116.60
2	A	1101	YUN	C2-N1-C6	6.16	121.38	116.60
2	A	1101	YUN	N3-C2-N1	-5.44	121.28	128.67
2	B	1101	YUN	N3-C2-N1	-5.40	121.34	128.67
2	A	1101	YUN	C2-N3-C4	4.77	120.85	115.43
2	B	1101	YUN	C2-N3-C4	4.60	120.66	115.43
2	A	1101	YUN	CAM-C5-C6	-4.40	121.27	124.84
2	B	1101	YUN	CAM-C5-C6	-4.29	121.35	124.84
2	A	1101	YUN	C6-C5-C4	3.90	118.16	115.86
2	B	1101	YUN	C6-C5-C4	3.82	118.12	115.86
2	B	1101	YUN	C5-C6-N1	-3.55	118.72	121.35
2	A	1101	YUN	CAZ-NAS-CAW	-3.54	121.26	127.52
2	A	1101	YUN	C5-C6-N1	-3.39	118.84	121.35
2	B	1101	YUN	CAZ-NAS-CAW	-3.33	121.62	127.52
2	A	1101	YUN	C5-C4-N3	-3.19	119.44	122.82
2	B	1101	YUN	C5-C4-N3	-3.06	119.57	122.82
2	B	1101	YUN	CAY-CAL-CAX	2.35	120.96	118.66
2	B	1101	YUN	CAY-NAT-C6	-2.24	122.48	128.22
2	A	1101	YUN	CAY-NAT-C6	-2.16	122.68	128.22

There are no chirality outliers.

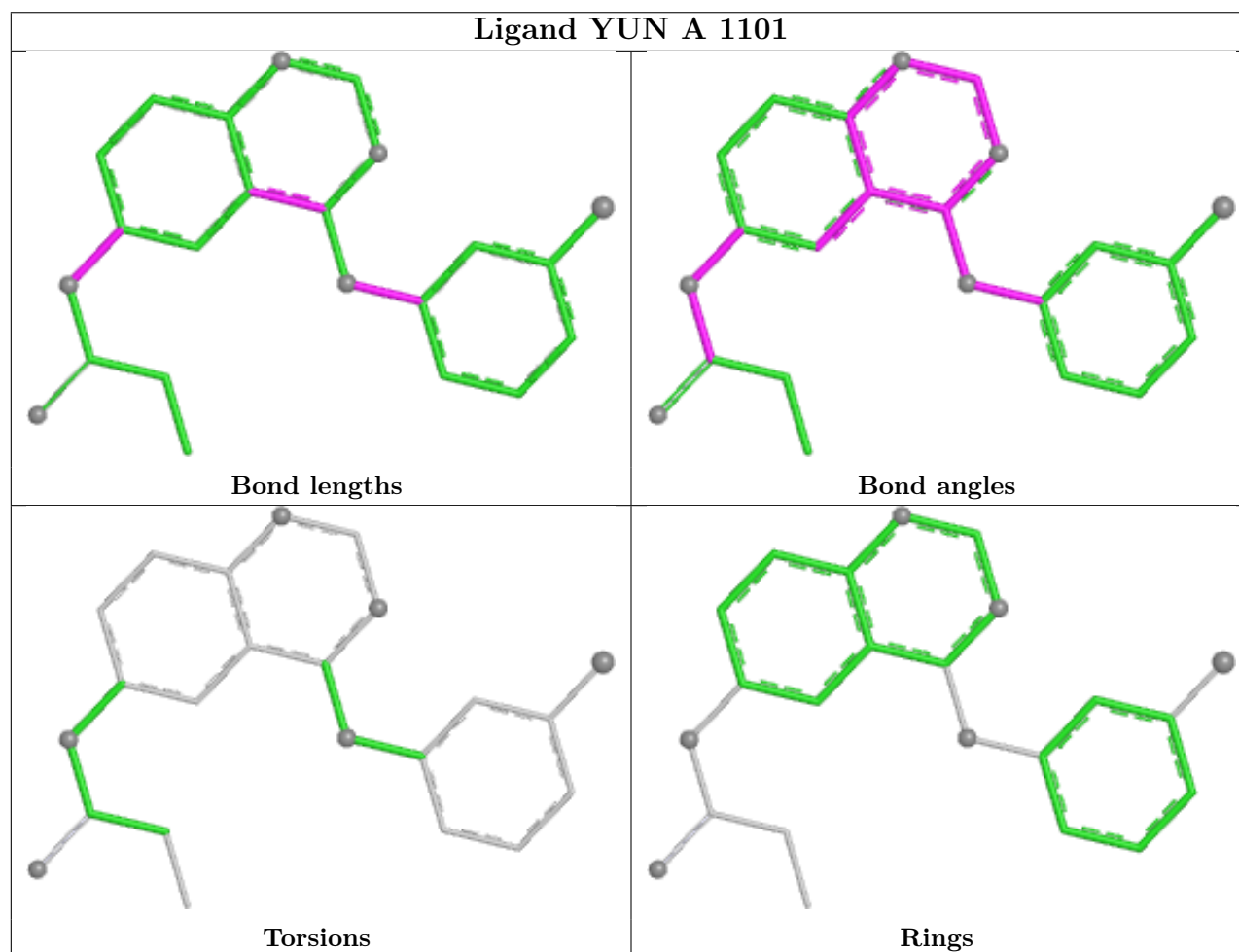
There are no torsion outliers.

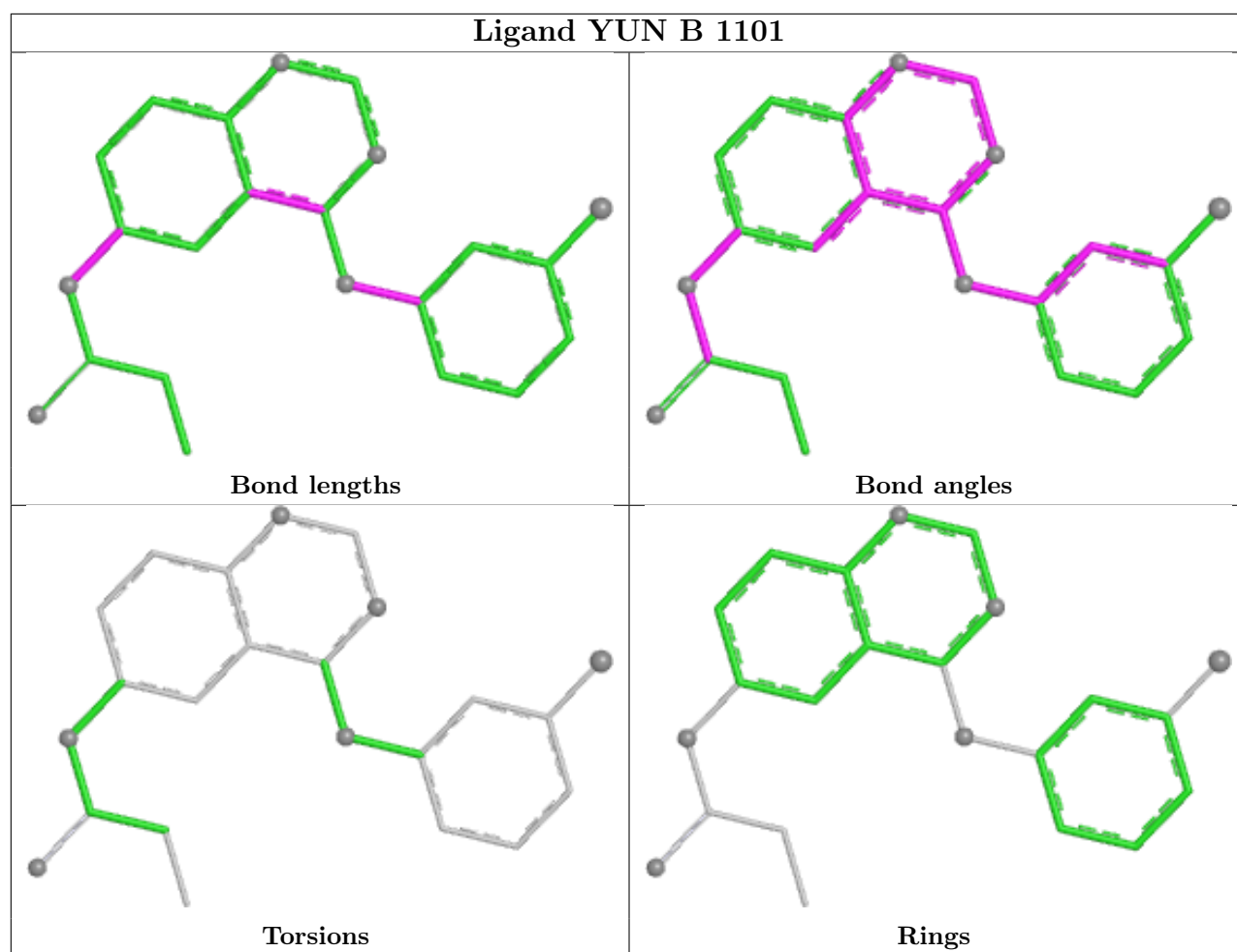
There are no ring outliers.

2 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1101	YUN	12	0
2	B	1101	YUN	7	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	291/331 (87%)	-0.97	0 100 100	159, 256, 310, 335	1 (0%)
1	B	281/331 (84%)	-1.01	0 100 100	161, 255, 301, 327	1 (0%)
All	All	572/662 (86%)	-0.99	0 100 100	159, 256, 307, 335	2 (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](#)

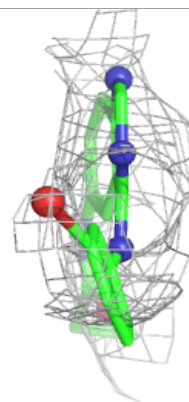
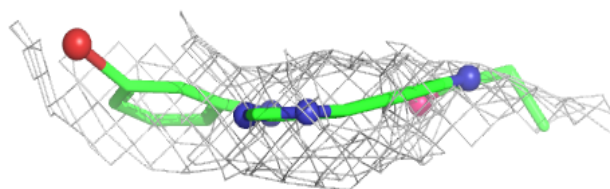
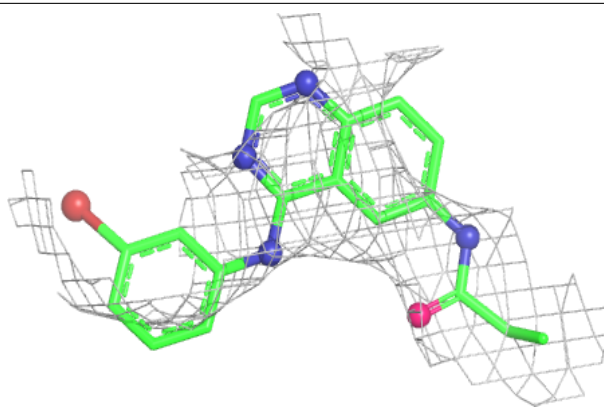
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	YUN	B	1101	23/23	0.91	0.06	211,256,286,297	0
2	YUN	A	1101	23/23	0.92	0.05	208,256,288,291	0

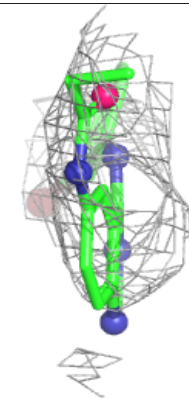
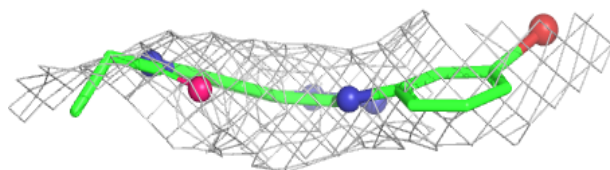
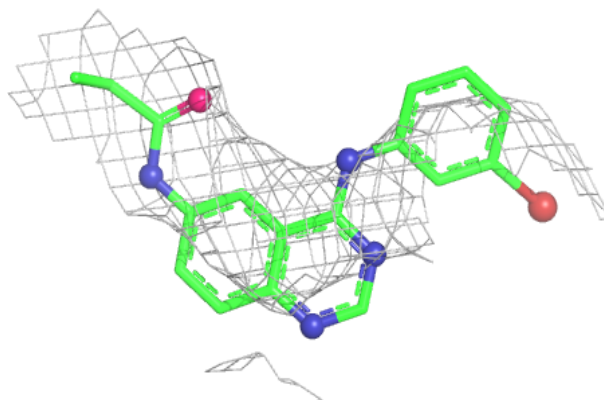
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around YUN B 1101:

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

**Electron density around YUN A 1101:**

$2mF_o-DF_c$ (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.