



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 12:49 am BST

PDB ID : 3DS6
Title : P38 complex with a phthalazine inhibitor
Authors : Herberich, B.; Syed, R.; Li, V.; Grosfeld, D.
Deposited on : 2008-07-11
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

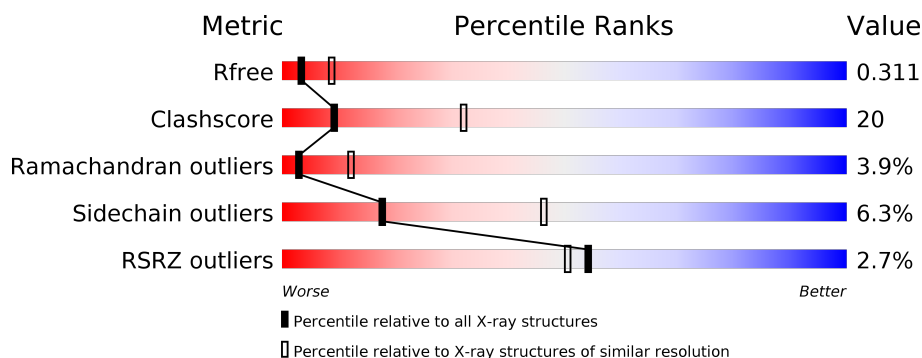
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on 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	366	
1	B	366	
1	C	366	
1	D	366	

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	A17	A	361	-	X	-	-
2	A17	B	361	-	X	-	-
2	A17	D	361	-	X	-	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11384 atoms, of which 92 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 Mitogen-activated protein kinase 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	343	Total	C	N	O	S	0	0	0
			2769	1774	475	508	12			
1	B	343	Total	C	N	O	S	0	0	0
			2769	1774	475	508	12			
1	C	349	Total	C	N	O	S	0	0	0
			2817	1803	481	520	13			
1	D	349	Total	C	N	O	S	0	0	0
			2817	1803	481	520	13			

There are 24 discrepancies between the modelled and reference sequences:

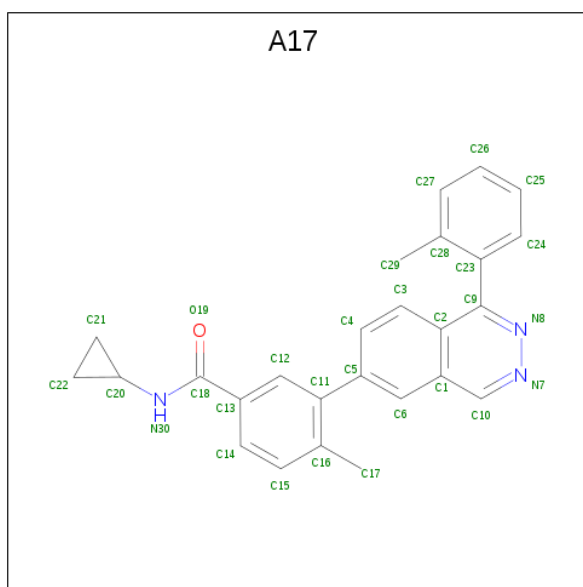
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	EXPRESSION TAG	UNP Q16539
A	-4	SER	-	EXPRESSION TAG	UNP Q16539
A	-3	HIS	-	EXPRESSION TAG	UNP Q16539
A	-2	MET	-	EXPRESSION TAG	UNP Q16539
A	-1	LEU	-	EXPRESSION TAG	UNP Q16539
A	0	GLU	-	EXPRESSION TAG	UNP Q16539
B	-5	GLY	-	EXPRESSION TAG	UNP Q16539
B	-4	SER	-	EXPRESSION TAG	UNP Q16539
B	-3	HIS	-	EXPRESSION TAG	UNP Q16539
B	-2	MET	-	EXPRESSION TAG	UNP Q16539
B	-1	LEU	-	EXPRESSION TAG	UNP Q16539
B	0	GLU	-	EXPRESSION TAG	UNP Q16539
C	-5	GLY	-	EXPRESSION TAG	UNP Q16539
C	-4	SER	-	EXPRESSION TAG	UNP Q16539
C	-3	HIS	-	EXPRESSION TAG	UNP Q16539
C	-2	MET	-	EXPRESSION TAG	UNP Q16539
C	-1	LEU	-	EXPRESSION TAG	UNP Q16539
C	0	GLU	-	EXPRESSION TAG	UNP Q16539
D	-5	GLY	-	EXPRESSION TAG	UNP Q16539
D	-4	SER	-	EXPRESSION TAG	UNP Q16539
D	-3	HIS	-	EXPRESSION TAG	UNP Q16539

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	MET	-	EXPRESSION TAG	UNP Q16539
D	-1	LEU	-	EXPRESSION TAG	UNP Q16539
D	0	GLU	-	EXPRESSION TAG	UNP Q16539

- Molecule 2 is N-cyclopropyl-4-methyl-3-[1-(2-methylphenyl)phthalazin-6-yl]benzamide (three-letter code: A17) (formula: $\text{C}_{26}\text{H}_{23}\text{N}_3\text{O}$).

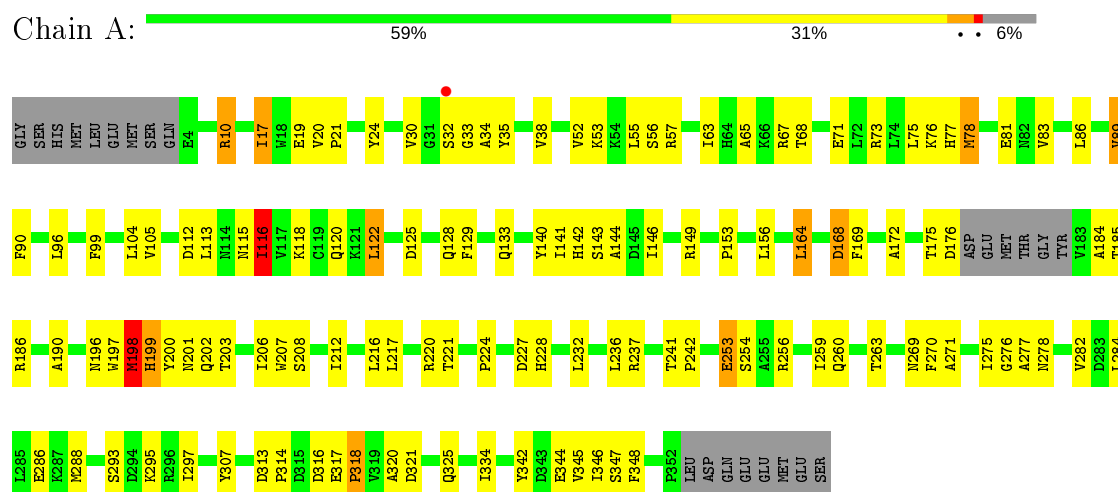


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 53	C 26	H 23	N 3	O 1	23	0
2	B	1	Total 53	C 26	H 23	N 3	O 1	23	0
2	C	1	Total 53	C 26	H 23	N 3	O 1	23	0
2	D	1	Total 53	C 26	H 23	N 3	O 1	23	0

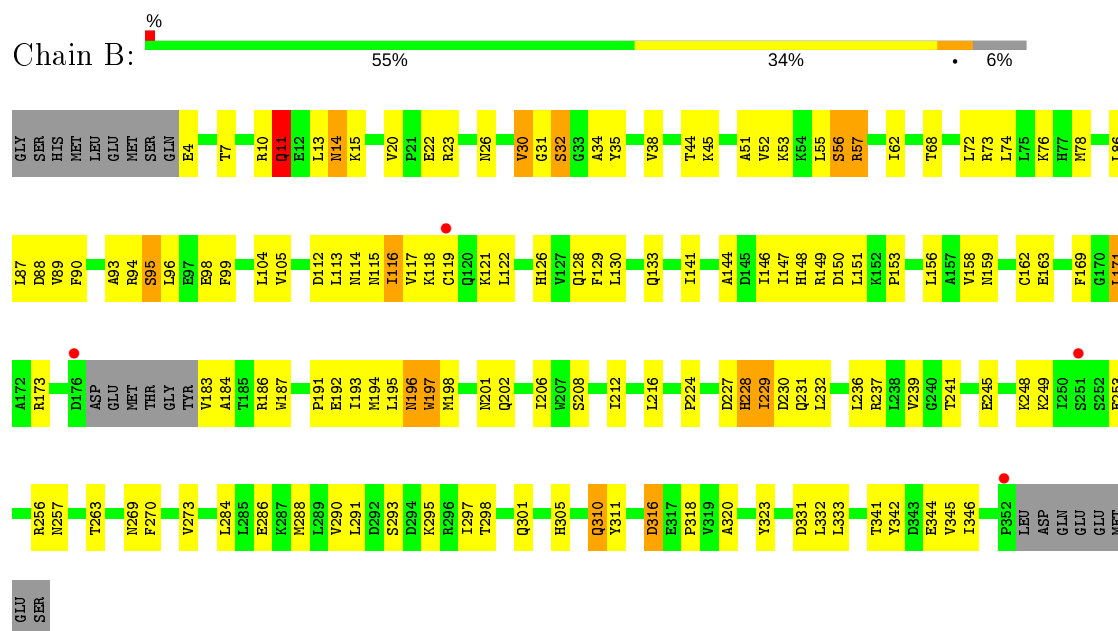
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

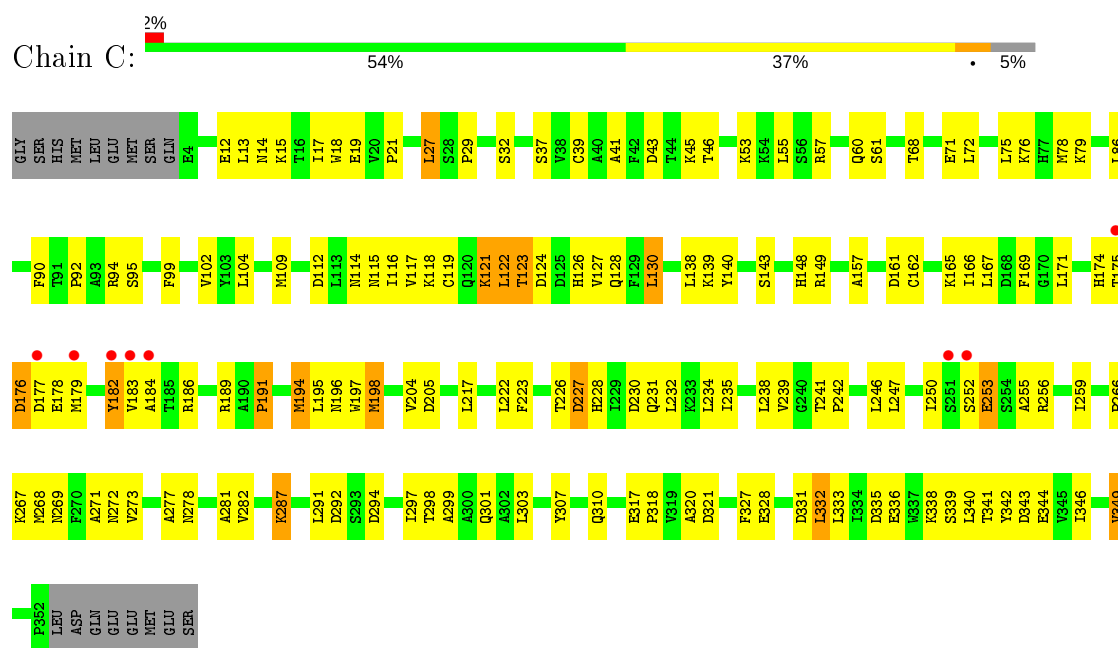
- Molecule 1: Mitogen-activated protein kinase 14



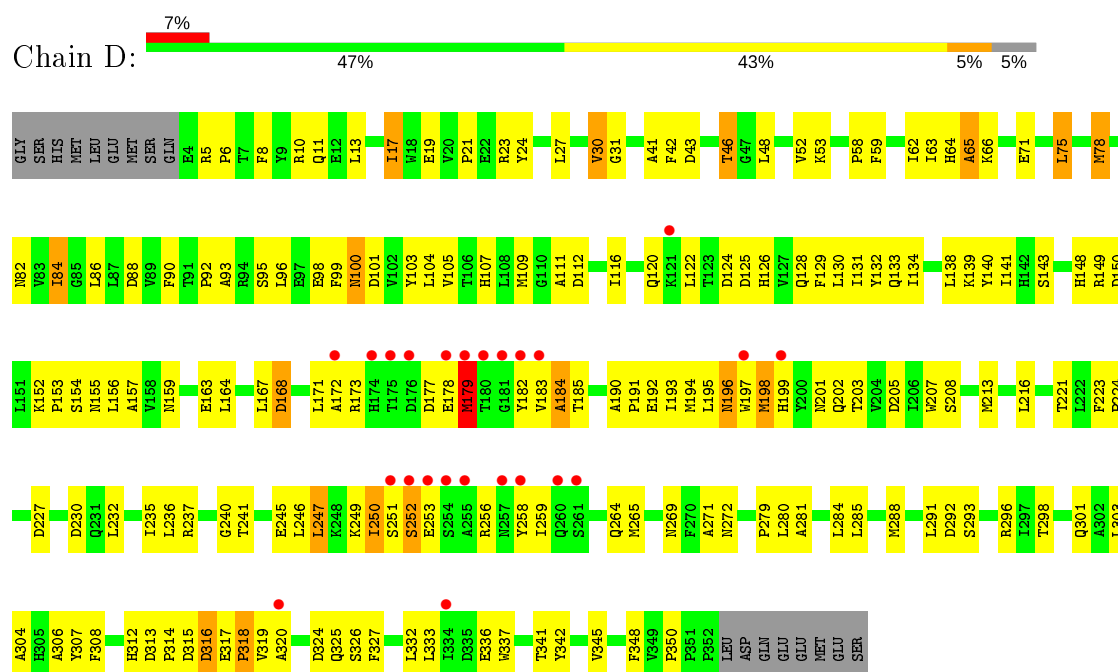
- Molecule 1: Mitogen-activated protein kinase 14



- Molecule 1: Mitogen-activated protein kinase 14



• Molecule 1: Mitogen-activated protein kinase 14



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.95Å 87.88Å 224.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.80 – 2.90 45.80 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.9 (45.80-2.90) 75.2 (45.80-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.228 , 0.319 0.224 , 0.311	Depositor DCC
R_{free} test set	3365 reflections (8.13%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtriage
Anisotropy	0.728	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.7	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	0.92	EDS
Total number of atoms	11384	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: A17

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	0.50	0/2833	0.73	1/3846 (0.0%)
1	B	0.45	0/2833	0.70	0/3846
1	C	0.48	0/2883	0.73	0/3915
1	D	0.40	0/2883	0.62	0/3915
All	All	0.46	0/11432	0.70	1/15522 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2769	0	2767	100	0
1	B	2769	0	2767	109	1
1	C	2817	0	2806	119	0
1	D	2817	0	2806	132	1
2	A	30	23	23	4	0
2	B	30	23	23	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	30	23	23	3	0
2	D	30	23	23	3	0
All	All	11292	92	11238	448	1

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 20.

All (448) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:LEU:HB3	1:A:86:LEU:HD13	1.47	0.95
1:B:86:LEU:HD21	1:B:89:VAL:HG22	1.50	0.93
1:D:316:ASP:O	1:D:318:PRO:HD3	1.70	0.92
1:C:140:TYR:CE1	1:C:320:ALA:HB2	2.08	0.89
1:A:76:LYS:HE2	1:A:344:GLU:HG3	1.54	0.89
1:A:34:ALA:HB1	1:C:15:LYS:HG2	1.55	0.86
1:C:246:LEU:HD11	1:C:292:ASP:HB2	1.57	0.85
1:D:53:LYS:HB2	2:D:361:A17:H17B	1.57	0.84
1:D:140:TYR:CE2	1:D:320:ALA:HB2	2.13	0.84
1:A:185:THR:HG21	1:C:12:GLU:OE2	1.77	0.83
1:A:78:MET:HG2	1:A:83:VAL:HG11	1.61	0.83
1:C:195:LEU:HD13	1:C:255:ALA:HB2	1.62	0.81
1:C:195:LEU:HB3	1:C:197:TRP:HE1	1.43	0.81
1:C:127:VAL:HG11	1:C:217:LEU:HD21	1.63	0.81
1:C:27:LEU:HD22	1:C:41:ALA:HB2	1.63	0.80
1:A:86:LEU:HD11	1:A:104:LEU:HD22	1.62	0.79
1:B:86:LEU:HD21	1:B:89:VAL:CG2	2.14	0.78
1:A:17:ILE:HD12	1:C:182:TYR:O	1.84	0.77
1:D:96:LEU:HD13	1:D:342:TYR:CD2	2.18	0.77
1:B:76:LYS:HE3	1:B:344:GLU:O	1.86	0.75
1:B:331:ASP:O	1:B:332:LEU:HD23	1.86	0.75
1:D:8:PHE:CE2	1:D:21:PRO:HD3	2.22	0.74
1:C:45:LYS:HG3	1:C:45:LYS:O	1.87	0.74
1:D:332:LEU:HB3	1:D:336:GLU:OE2	1.89	0.73
1:B:57:ARG:O	1:B:57:ARG:HG2	1.87	0.72
1:A:346:ILE:HD12	1:A:346:ILE:C	2.09	0.72
1:D:27:LEU:HD23	1:D:41:ALA:HB2	1.71	0.72
1:C:179:MET:HG3	1:C:189:ARG:HH12	1.56	0.71
1:D:131:ILE:HG21	1:D:213:MET:HE3	1.72	0.71
1:B:147:ILE:CG2	1:B:149:ARG:HG3	2.19	0.71
1:A:200:TYR:HD1	1:A:200:TYR:O	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:ASP:HA	1:D:348:PHE:CE2	2.26	0.70
1:D:298:THR:OG1	1:D:301:GLN:HG3	1.90	0.70
1:C:223:PHE:CE1	1:C:235:ILE:HA	2.26	0.70
1:D:150:ASP:HB3	1:D:171:LEU:HD12	1.74	0.69
1:B:146:ILE:HD11	1:B:323:TYR:CD2	2.27	0.69
1:C:76:LYS:HE2	1:C:344:GLU:OE2	1.92	0.69
1:B:341:THR:O	1:B:345:VAL:HG23	1.93	0.69
1:C:195:LEU:HB3	1:C:197:TRP:NE1	2.07	0.69
1:D:155:ASN:OD1	1:D:168:ASP:HB2	1.93	0.69
1:B:153:PRO:HA	1:B:156:LEU:HD12	1.73	0.69
1:B:147:ILE:HG22	1:B:149:ARG:HG3	1.74	0.68
1:D:62:ILE:O	1:D:66:LYS:HG2	1.94	0.68
1:A:256:ARG:O	1:A:259:ILE:HG22	1.93	0.68
1:B:126:HIS:O	1:B:130:LEU:HD12	1.93	0.68
1:C:195:LEU:O	1:C:197:TRP:HD1	1.77	0.67
1:B:113:LEU:O	1:B:117:VAL:HG23	1.95	0.67
1:A:57:ARG:HG2	1:C:177:ASP:HB3	1.76	0.67
1:C:117:VAL:HG12	1:C:117:VAL:O	1.94	0.66
1:D:88:ASP:HA	1:D:348:PHE:HE2	1.61	0.66
1:D:58:PRO:HG2	1:D:59:PHE:CD2	2.30	0.66
1:D:96:LEU:HD13	1:D:342:TYR:CG	2.31	0.65
1:B:30:VAL:HG22	1:B:38:VAL:O	1.96	0.65
1:C:148:HIS:O	1:C:149:ARG:HB2	1.97	0.65
1:D:227:ASP:OD1	1:D:230:ASP:HB2	1.96	0.65
1:B:245:GLU:HG2	1:B:249:LYS:HE3	1.78	0.65
1:D:280:LEU:HD21	1:D:306:ALA:HB1	1.79	0.64
1:A:203:THR:OG1	1:A:293:SER:HB2	1.98	0.64
1:D:65:ALA:HB1	1:D:337:TRP:HB2	1.78	0.64
1:A:115:ASN:OD1	1:A:118:LYS:HD2	1.97	0.64
1:B:53:LYS:HB2	2:B:361:A17:H17B	1.80	0.64
1:B:87:LEU:HD23	1:B:87:LEU:O	1.97	0.63
1:A:34:ALA:HB2	1:C:14:ASN:HA	1.79	0.63
1:B:13:LEU:O	1:B:14:ASN:HB2	1.98	0.63
1:C:343:ASP:O	1:C:346:ILE:HG13	1.99	0.62
1:B:56:SER:C	1:B:57:ARG:HD3	2.19	0.62
1:D:131:ILE:HG21	1:D:213:MET:CE	2.30	0.62
1:D:48:LEU:HD13	1:D:107:HIS:HE1	1.65	0.62
1:B:239:VAL:HG21	1:B:291:LEU:HG	1.81	0.61
1:A:76:LYS:HG2	1:A:86:LEU:HD22	1.81	0.61
1:A:75:LEU:CB	1:A:86:LEU:HD13	2.26	0.61
1:C:112:ASP:OD1	1:C:115:ASN:HB2	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:183:VAL:O	1:D:184:ALA:HB3	1.99	0.61
1:C:278:ASN:O	1:C:282:VAL:HG23	2.01	0.61
1:C:127:VAL:HG11	1:C:217:LEU:CD2	2.31	0.61
1:D:152:LYS:HE2	1:D:185:THR:OG1	1.98	0.61
1:A:17:ILE:HG23	1:A:17:ILE:O	2.00	0.61
1:A:241:THR:HB	1:A:242:PRO:HD2	1.82	0.61
1:D:19:GLU:OE1	1:D:92:PRO:HB3	2.01	0.61
1:A:270:PHE:CD2	1:A:286:GLU:HG2	2.36	0.60
1:C:124:ASP:OD2	1:C:278:ASN:HB2	2.01	0.60
1:B:116:ILE:HA	1:B:119:CYS:HB3	1.83	0.60
1:C:18:TRP:HZ2	1:C:37:SER:HB3	1.66	0.60
1:B:56:SER:O	1:B:57:ARG:HD3	2.01	0.60
1:D:116:ILE:O	1:D:120:GLN:HB3	2.01	0.60
1:B:113:LEU:O	1:B:116:ILE:HG22	2.01	0.60
1:A:57:ARG:HA	1:C:177:ASP:OD1	2.01	0.60
1:D:93:ALA:HB2	1:D:99:PHE:HA	1.82	0.60
1:C:191:PRO:O	1:C:194:MET:HB2	2.01	0.60
1:B:31:GLY:O	1:B:32:SER:O	2.19	0.60
1:D:100:ASN:O	1:D:101:ASP:OD2	2.20	0.60
1:A:277:ALA:O	1:A:278:ASN:C	2.41	0.59
1:B:342:TYR:CZ	1:B:346:ILE:HD11	2.37	0.59
1:C:124:ASP:O	1:C:128:GLN:HG3	2.02	0.59
1:C:169:PHE:O	1:C:171:LEU:HG	2.02	0.59
1:D:246:LEU:HD13	1:D:292:ASP:OD2	2.03	0.59
1:A:156:LEU:HD22	1:A:164:LEU:HD11	1.84	0.59
1:C:269:ASN:HB3	1:C:272:ASN:HD22	1.67	0.59
1:D:48:LEU:HD13	1:D:107:HIS:CE1	2.37	0.59
1:D:53:LYS:CB	2:D:361:A17:H17B	2.29	0.59
1:A:232:LEU:O	1:A:236:LEU:HG	2.03	0.59
1:A:200:TYR:CD1	1:A:200:TYR:O	2.55	0.59
1:D:71:GLU:O	1:D:75:LEU:HD12	2.03	0.59
1:C:269:ASN:OD1	1:C:271:ALA:HB3	2.02	0.58
1:D:62:ILE:O	1:D:65:ALA:HB3	2.03	0.58
1:B:298:THR:OG1	1:B:301:GLN:HG3	2.02	0.58
1:D:245:GLU:O	1:D:249:LYS:HG3	2.03	0.58
1:B:148:HIS:O	1:B:149:ARG:HB2	2.03	0.58
1:A:142:HIS:HB3	1:A:202:GLN:NE2	2.17	0.58
1:C:140:TYR:CZ	1:C:320:ALA:HB2	2.37	0.58
1:D:195:LEU:HD12	1:D:232:LEU:HD22	1.83	0.58
1:B:151:LEU:HD11	1:B:156:LEU:HD21	1.84	0.58
1:C:255:ALA:O	1:C:259:ILE:HG12	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:269:ASN:HB3	1:C:272:ASN:ND2	2.18	0.58
1:A:71:GLU:HG3	2:A:361:A17:H21A	1.86	0.57
1:D:109:MET:HB2	1:D:157:ALA:HB1	1.87	0.56
1:C:250:ILE:HG21	1:C:255:ALA:HB3	1.87	0.56
1:D:139:LYS:HE3	1:D:319:VAL:HG12	1.88	0.56
1:D:24:TYR:OH	1:D:88:ASP:OD2	2.21	0.56
1:A:38:VAL:HG22	1:A:53:LYS:HB2	1.88	0.55
1:D:221:THR:HG21	1:D:224:PRO:HB3	1.88	0.55
1:A:129:PHE:O	1:A:133:GLN:HG3	2.07	0.55
1:A:153:PRO:HA	1:A:156:LEU:HD12	1.88	0.55
1:A:83:VAL:HG13	1:A:169:PHE:CE2	2.41	0.55
1:D:128:GLN:HG2	1:D:307:TYR:CZ	2.41	0.55
1:D:112:ASP:OD1	1:D:154:SER:HA	2.06	0.55
1:A:17:ILE:HB	1:C:184:ALA:N	2.21	0.55
1:B:94:ARG:O	1:B:95:SER:CB	2.55	0.55
1:C:297:ILE:CG1	1:C:301:GLN:HG3	2.36	0.55
1:A:253:GLU:HG3	1:A:254:SER:H	1.71	0.55
1:C:239:VAL:O	1:C:267:LYS:HB2	2.07	0.55
1:C:19:GLU:HG2	1:C:92:PRO:HB3	1.89	0.55
1:D:246:LEU:HD13	1:D:292:ASP:CG	2.27	0.55
1:A:256:ARG:O	1:A:260:GLN:HG2	2.07	0.55
1:B:245:GLU:O	1:B:248:LYS:HB3	2.06	0.55
1:C:241:THR:HG23	1:C:242:PRO:HD2	1.88	0.55
1:D:178:GLU:HA	1:D:183:VAL:HG11	1.88	0.55
1:D:17:ILE:HD13	1:D:17:ILE:C	2.27	0.55
1:D:84:ILE:O	1:D:84:ILE:HG23	2.07	0.55
1:D:341:THR:O	1:D:345:VAL:HG23	2.07	0.54
1:A:185:THR:HG22	1:C:17:ILE:HG13	1.89	0.54
1:D:253:GLU:HA	1:D:256:ARG:HB3	1.88	0.54
1:A:208:SER:O	1:A:212:ILE:HG13	2.07	0.54
1:C:109:MET:SD	1:C:165:LYS:HD2	2.48	0.54
1:D:342:TYR:O	1:D:345:VAL:HB	2.08	0.54
1:C:99:PHE:O	1:C:338:LYS:HE3	2.08	0.53
1:A:284:LEU:HG	1:A:288:MET:HE1	1.89	0.53
1:D:93:ALA:HB2	1:D:98:GLU:O	2.09	0.53
1:B:94:ARG:O	1:B:95:SER:HB3	2.08	0.53
1:D:159:ASN:OD1	1:D:163:GLU:HB2	2.07	0.53
1:B:208:SER:O	1:B:212:ILE:HG13	2.09	0.53
1:B:7:THR:O	1:B:22:GLU:HG2	2.09	0.53
1:D:190:ALA:HA	1:D:207:TRP:CD1	2.44	0.53
1:B:194:MET:HG2	1:B:194:MET:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:LEU:HD22	1:B:241:THR:HA	1.92	0.52
1:D:269:ASN:HD21	1:D:271:ALA:HB3	1.74	0.52
1:D:148:HIS:O	1:D:149:ARG:HB2	2.08	0.52
1:B:122:LEU:HD13	1:B:216:LEU:O	2.09	0.52
1:D:284:LEU:O	1:D:288:MET:HG3	2.08	0.52
1:D:304:ALA:HB2	1:D:312:HIS:CE1	2.44	0.52
1:C:253:GLU:H	1:C:256:ARG:HH12	1.57	0.52
1:D:221:THR:CG2	1:D:224:PRO:HB3	2.39	0.52
1:A:141:ILE:HG22	1:A:146:ILE:O	2.09	0.52
1:D:193:ILE:O	1:D:193:ILE:HG12	2.09	0.52
1:A:284:LEU:HG	1:A:288:MET:CE	2.40	0.52
1:B:78:MET:HG3	1:B:169:PHE:CZ	2.45	0.52
1:D:130:LEU:O	1:D:134:ILE:HG13	2.09	0.52
1:A:52:VAL:HG22	1:A:105:VAL:HG22	1.91	0.52
1:D:333:LEU:HB2	1:D:336:GLU:HG3	1.92	0.52
1:B:144:ALA:HB2	1:B:320:ALA:HB3	1.91	0.51
1:C:161:ASP:O	1:C:162:CYS:HB2	2.09	0.51
1:B:90:PHE:C	1:B:90:PHE:CD1	2.84	0.51
1:D:237:ARG:HG2	1:D:265:MET:SD	2.50	0.51
1:B:269:ASN:O	1:B:273:VAL:HG23	2.10	0.51
1:A:168:ASP:HA	2:A:361:A17:H20	1.93	0.51
1:A:33:GLY:HA2	1:C:32:SER:O	2.10	0.51
1:A:75:LEU:HD13	1:A:86:LEU:CD1	2.41	0.51
1:B:56:SER:O	1:B:57:ARG:C	2.49	0.51
1:C:195:LEU:O	1:C:197:TRP:CD1	2.62	0.50
1:D:124:ASP:O	1:D:128:GLN:HG3	2.11	0.50
1:A:113:LEU:O	1:A:116:ILE:HB	2.12	0.50
1:B:227:ASP:O	1:B:228:HIS:C	2.49	0.50
1:A:17:ILE:HB	1:C:184:ALA:H	1.76	0.50
1:C:186:ARG:CZ	1:C:227:ASP:HA	2.42	0.50
1:C:43:ASP:HB3	1:C:46:THR:OG1	2.11	0.50
1:B:183:VAL:O	1:B:184:ALA:HB3	2.12	0.50
1:C:222:LEU:HD23	1:C:223:PHE:CE2	2.46	0.50
1:B:146:ILE:HD11	1:B:323:TYR:CE2	2.46	0.50
1:B:93:ALA:HB2	1:B:98:GLU:O	2.11	0.50
1:C:45:LYS:O	1:C:45:LYS:CG	2.60	0.50
1:B:270:PHE:CD2	1:B:286:GLU:HG2	2.46	0.50
1:B:297:ILE:HD12	1:B:305:HIS:CE1	2.46	0.50
1:C:342:TYR:CE2	1:C:346:ILE:HG21	2.47	0.50
1:B:147:ILE:HD11	1:B:202:GLN:HA	1.94	0.50
1:B:284:LEU:HG	1:B:288:MET:HE2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:LEU:HD22	1:C:29:PRO:HD3	1.93	0.50
1:A:221:THR:HG22	1:A:224:PRO:HD3	1.93	0.49
1:A:17:ILE:HD11	1:A:19:GLU:OE1	2.12	0.49
1:B:297:ILE:HD12	1:B:305:HIS:HE1	1.77	0.49
1:D:52:VAL:HA	1:D:104:LEU:O	2.13	0.49
1:A:190:ALA:HA	1:A:207:TRP:CD1	2.48	0.49
1:B:113:LEU:HB3	1:B:153:PRO:HB3	1.95	0.49
1:B:51:ALA:O	1:B:105:VAL:HA	2.11	0.49
1:C:227:ASP:O	1:C:228:HIS:C	2.50	0.49
1:D:203:THR:HB	1:D:296:ARG:HD2	1.93	0.49
1:D:308:PHE:O	1:D:312:HIS:HB3	2.12	0.49
1:A:275:ILE:HG13	1:A:275:ILE:O	2.13	0.49
1:A:35:TYR:HB3	2:A:361:A17:H4	1.94	0.49
1:D:129:PHE:O	1:D:132:TYR:HB3	2.13	0.49
1:D:112:ASP:HA	1:D:156:LEU:O	2.12	0.49
1:D:250:ILE:HD11	1:D:259:ILE:HD12	1.94	0.49
1:A:17:ILE:CG2	1:C:183:VAL:HA	2.43	0.48
1:D:13:LEU:CD1	1:D:27:LEU:HD13	2.43	0.48
1:C:195:LEU:CB	1:C:197:TRP:HE1	2.21	0.48
1:C:186:ARG:NH2	1:C:227:ASP:HA	2.28	0.48
1:C:99:PHE:CD2	1:C:342:TYR:HD1	2.31	0.48
1:C:223:PHE:HB3	1:C:231:GLN:HE22	1.78	0.48
1:C:53:LYS:HB2	2:C:361:A17:H17B	1.96	0.48
1:D:348:PHE:O	1:D:350:PRO:HD3	2.13	0.48
1:D:27:LEU:CD2	1:D:41:ALA:HB2	2.42	0.48
1:C:78:MET:HG3	1:C:169:PHE:CZ	2.48	0.48
1:A:232:LEU:HD21	1:A:259:ILE:CD1	2.44	0.48
1:B:171:LEU:HD13	1:B:171:LEU:O	2.13	0.48
1:C:75:LEU:HB3	1:C:86:LEU:HG	1.96	0.48
1:D:131:ILE:HG13	1:D:213:MET:CE	2.43	0.48
1:C:269:ASN:O	1:C:273:VAL:HG23	2.13	0.48
1:D:179:MET:N	1:D:183:VAL:HB	2.28	0.48
1:C:246:LEU:HD21	1:C:292:ASP:CG	2.33	0.48
1:D:5:ARG:HG2	1:D:6:PRO:HD2	1.94	0.48
1:D:75:LEU:HB3	1:D:86:LEU:HB2	1.95	0.48
1:A:201:ASN:OD1	1:A:203:THR:HG23	2.13	0.48
1:B:121:LYS:HD2	1:B:121:LYS:HA	1.72	0.48
1:C:186:ARG:NH1	1:C:227:ASP:HA	2.28	0.48
1:A:128:GLN:HG2	1:A:307:TYR:CZ	2.49	0.47
1:B:72:LEU:O	1:B:76:LYS:HB2	2.14	0.47
1:C:55:LEU:HD12	1:C:102:VAL:HB	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:252:SER:H	1:D:256:ARG:HD2	1.78	0.47
1:D:88:ASP:O	1:D:105:VAL:HB	2.13	0.47
1:B:55:LEU:HD13	1:B:68:THR:HG23	1.96	0.47
1:C:281:ALA:HB2	1:C:307:TYR:CE1	2.49	0.47
1:D:152:LYS:HB2	1:D:153:PRO:HD2	1.97	0.47
1:A:186:ARG:NH1	1:A:227:ASP:HA	2.29	0.47
1:B:290:VAL:HG11	1:B:295:LYS:HB2	1.97	0.47
1:D:201:ASN:CG	1:D:202:GLN:H	2.17	0.47
1:D:247:LEU:O	1:D:250:ILE:HG12	2.15	0.47
1:A:76:LYS:HD2	1:A:348:PHE:HD1	1.80	0.47
1:B:114:ASN:HA	1:B:117:VAL:HG23	1.97	0.47
1:D:324:ASP:O	1:D:327:PHE:HE1	1.98	0.47
1:C:55:LEU:HD13	1:C:68:THR:HG23	1.97	0.47
1:D:269:ASN:HD22	1:D:272:ASN:CG	2.18	0.47
1:A:112:ASP:HA	1:A:156:LEU:O	2.14	0.47
1:C:226:THR:O	1:C:227:ASP:HB3	2.14	0.47
1:C:72:LEU:HD23	1:C:341:THR:HG23	1.96	0.46
1:A:89:VAL:HG21	1:A:345:VAL:HG22	1.96	0.46
1:C:99:PHE:CG	1:C:342:TYR:HD1	2.34	0.46
1:D:198:MET:H	1:D:198:MET:HE3	1.81	0.46
1:A:55:LEU:HD13	1:A:68:THR:HG23	1.96	0.46
1:C:223:PHE:HZ	1:C:238:LEU:HD23	1.80	0.46
1:A:142:HIS:HB3	1:A:202:GLN:HE22	1.79	0.46
1:B:206:ILE:HG13	1:B:297:ILE:O	2.16	0.46
1:C:112:ASP:OD2	1:C:114:ASN:HB3	2.16	0.46
1:D:63:ILE:H	1:D:63:ILE:HD12	1.80	0.46
1:B:201:ASN:HB2	1:B:293:SER:HB3	1.97	0.46
1:D:303:LEU:HD13	1:D:317:GLU:OE1	2.16	0.46
1:A:20:VAL:HG12	1:A:90:PHE:HZ	1.81	0.46
1:C:317:GLU:N	1:C:318:PRO:HD3	2.30	0.46
1:A:317:GLU:N	1:A:318:PRO:HD3	2.31	0.46
1:A:76:LYS:HG2	1:A:86:LEU:CD2	2.46	0.46
1:B:195:LEU:HD11	1:B:232:LEU:HD22	1.98	0.46
1:D:122:LEU:CD1	1:D:216:LEU:HD22	2.46	0.46
1:A:113:LEU:HD22	1:A:156:LEU:HD13	1.98	0.45
1:C:297:ILE:HG12	1:C:301:GLN:HG3	1.98	0.45
1:C:297:ILE:HG13	1:C:301:GLN:HG3	1.98	0.45
1:A:65:ALA:HB2	1:A:334:ILE:HD12	1.99	0.45
1:A:346:ILE:C	1:A:346:ILE:CD1	2.77	0.45
1:A:56:SER:O	1:A:57:ARG:C	2.54	0.45
1:C:298:THR:OG1	1:C:301:GLN:HG2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:147:ILE:HD11	1:B:202:GLN:HG3	1.97	0.45
1:C:14:ASN:O	1:C:15:LYS:HB2	2.16	0.45
1:D:140:TYR:CD2	1:D:320:ALA:HB2	2.49	0.45
1:B:32:SER:C	1:B:34:ALA:H	2.20	0.45
1:C:253:GLU:H	1:C:256:ARG:NH1	2.12	0.45
1:A:269:ASN:HD21	1:A:271:ALA:HB3	1.82	0.45
1:B:194:MET:HE2	1:B:195:LEU:HG	1.99	0.45
1:A:89:VAL:CG2	1:A:345:VAL:HG22	2.46	0.45
1:B:129:PHE:O	1:B:133:GLN:HG3	2.16	0.45
1:C:277:ALA:O	1:C:278:ASN:C	2.54	0.45
1:D:152:LYS:HB2	1:D:153:PRO:CD	2.46	0.45
1:D:111:ALA:O	1:D:157:ALA:HA	2.15	0.45
1:D:13:LEU:HD11	1:D:27:LEU:HB3	1.98	0.45
1:A:20:VAL:HG12	1:A:90:PHE:CZ	2.51	0.45
1:C:124:ASP:HA	1:C:127:VAL:HG23	1.98	0.45
1:D:291:LEU:HA	1:D:291:LEU:HD12	1.78	0.45
1:B:32:SER:O	1:B:34:ALA:N	2.49	0.45
1:C:327:PHE:CE2	1:C:328:GLU:HG2	2.52	0.45
1:D:223:PHE:CE1	1:D:235:ILE:HA	2.52	0.45
1:A:198:MET:SD	1:A:199:HIS:O	2.75	0.45
1:B:195:LEU:O	1:B:196:ASN:C	2.54	0.45
1:B:229:ILE:O	1:B:230:ASP:C	2.54	0.45
1:B:316:ASP:C	1:B:318:PRO:HD3	2.37	0.45
1:B:99:PHE:CE1	1:B:342:TYR:HB2	2.51	0.45
1:C:223:PHE:HB3	1:C:231:GLN:NE2	2.32	0.45
1:D:78:MET:HE3	1:D:141:ILE:HG12	1.99	0.45
1:D:269:ASN:ND2	1:D:271:ALA:HB3	2.32	0.45
1:D:281:ALA:HB2	1:D:307:TYR:CE1	2.52	0.44
1:A:115:ASN:HA	1:A:118:LYS:HE3	1.98	0.44
1:B:112:ASP:HA	1:B:156:LEU:O	2.17	0.44
1:C:138:LEU:HD23	1:C:138:LEU:HA	1.71	0.44
1:C:287:LYS:HG2	1:C:287:LYS:H	1.50	0.44
1:D:63:ILE:N	1:D:63:ILE:HD12	2.32	0.44
1:C:123:THR:HG23	1:C:126:HIS:CE1	2.53	0.44
1:C:166:ILE:C	1:C:167:LEU:HD23	2.37	0.44
1:A:55:LEU:HA	1:A:55:LEU:HD23	1.69	0.44
1:A:346:ILE:HD12	1:A:347:SER:N	2.31	0.44
1:C:27:LEU:HD22	1:C:41:ALA:CB	2.42	0.44
1:D:232:LEU:O	1:D:236:LEU:HG	2.18	0.44
1:A:57:ARG:HG2	1:C:177:ASP:CB	2.46	0.44
1:B:38:VAL:HG22	1:B:53:LYS:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:301:GLN:HE21	1:C:301:GLN:HB3	1.58	0.44
1:C:349:VAL:O	1:C:349:VAL:HG12	2.18	0.44
1:D:316:ASP:C	1:D:318:PRO:HD3	2.33	0.44
1:B:297:ILE:CD1	1:B:305:HIS:CE1	3.01	0.44
1:D:280:LEU:HD21	1:D:306:ALA:CB	2.46	0.44
1:D:11:GLN:HG3	1:D:13:LEU:HG	2.00	0.44
1:A:206:ILE:HG13	1:A:297:ILE:O	2.18	0.43
1:B:112:ASP:OD1	1:B:112:ASP:C	2.56	0.43
1:D:183:VAL:O	1:D:184:ALA:CB	2.65	0.43
1:D:23:ARG:HG2	1:D:43:ASP:OD1	2.18	0.43
1:B:20:VAL:HG12	1:B:90:PHE:HZ	1.83	0.43
1:D:197:TRP:CE3	1:D:250:ILE:HG21	2.53	0.43
1:D:280:LEU:HD12	1:D:280:LEU:HA	1.63	0.43
1:D:178:GLU:OE2	1:D:178:GLU:HA	2.18	0.43
1:B:62:ILE:HD11	1:B:333:LEU:HD23	2.00	0.43
1:C:109:MET:HB2	1:C:157:ALA:HB1	1.99	0.43
1:B:141:ILE:HG23	1:B:146:ILE:HB	2.01	0.43
1:B:187:TRP:CD1	1:B:224:PRO:HA	2.53	0.43
1:C:78:MET:C	1:C:79:LYS:HG3	2.39	0.43
1:D:58:PRO:O	1:D:64:HIS:HB3	2.18	0.43
1:A:96:LEU:HD13	1:A:342:TYR:CD2	2.54	0.43
1:C:195:LEU:HB3	1:C:197:TRP:CD1	2.54	0.43
1:A:10:ARG:HH11	1:A:10:ARG:HG2	1.84	0.43
1:C:230:ASP:O	1:C:234:LEU:HG	2.18	0.43
1:D:90:PHE:CE1	1:D:103:TYR:CG	3.07	0.43
1:D:308:PHE:O	1:D:312:HIS:CB	2.67	0.43
1:B:195:LEU:HD23	1:B:195:LEU:HA	1.88	0.43
1:C:179:MET:HA	1:C:183:VAL:HG21	2.00	0.43
1:A:21:PRO:HG2	1:A:24:TYR:CD2	2.54	0.43
1:B:316:ASP:CG	1:B:316:ASP:O	2.58	0.43
1:C:281:ALA:HB2	1:C:307:TYR:CZ	2.53	0.43
1:C:340:LEU:HA	1:C:340:LEU:HD23	1.66	0.43
1:A:217:LEU:HA	1:A:217:LEU:HD23	1.79	0.42
1:C:122:LEU:HA	1:C:122:LEU:HD12	1.79	0.42
1:C:175:THR:O	1:C:176:ASP:HB2	2.19	0.42
1:D:195:LEU:O	1:D:196:ASN:C	2.58	0.42
1:D:65:ALA:HB1	1:D:337:TRP:CB	2.47	0.42
1:A:116:ILE:HA	1:A:116:ILE:HD13	1.58	0.42
1:B:310:GLN:HG3	1:B:310:GLN:O	2.19	0.42
1:C:204:VAL:HG23	1:C:205:ASP:N	2.34	0.42
1:D:173:ARG:HD3	1:D:173:ARG:HA	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:325:GLN:HG3	1:D:325:GLN:O	2.19	0.42
1:C:332:LEU:HD22	1:C:336:GLU:OE1	2.19	0.42
1:B:73:ARG:O	1:B:74:LEU:C	2.56	0.42
1:D:30:VAL:HB	2:D:361:A17:H25	2.01	0.42
1:A:316:ASP:C	1:A:318:PRO:HD3	2.40	0.42
1:A:38:VAL:HG22	1:A:53:LYS:CB	2.48	0.42
1:D:313:ASP:C	1:D:315:ASP:H	2.22	0.42
1:C:242:PRO:CG	1:C:247:LEU:HD13	2.49	0.42
1:D:140:TYR:O	1:D:143:SER:HB3	2.20	0.42
1:D:281:ALA:HB2	1:D:307:TYR:CZ	2.55	0.42
1:D:82:ASN:HD22	1:D:82:ASN:HA	1.62	0.42
1:A:140:TYR:O	1:A:143:SER:OG	2.29	0.42
1:B:159:ASN:OD1	1:B:159:ASN:C	2.57	0.42
1:A:122:LEU:CD1	1:A:216:LEU:HD22	2.50	0.42
1:A:83:VAL:CG1	1:A:169:PHE:CE2	3.02	0.42
1:B:144:ALA:HB2	1:B:320:ALA:CB	2.49	0.42
1:B:158:VAL:HA	1:B:163:GLU:O	2.20	0.42
1:B:192:GLU:HG2	1:B:193:ILE:N	2.35	0.42
1:C:223:PHE:CB	1:C:231:GLN:HE22	2.32	0.42
1:B:115:ASN:O	1:B:116:ILE:HB	2.18	0.41
1:C:157:ALA:HB2	1:C:167:LEU:HD11	2.02	0.41
1:B:171:LEU:CD1	1:B:171:LEU:O	2.69	0.41
1:B:23:ARG:O	1:B:44:THR:HG23	2.20	0.41
1:B:23:ARG:CZ	1:B:45:LYS:HD2	2.50	0.41
1:A:259:ILE:HG23	1:A:260:GLN:N	2.34	0.41
1:A:282:VAL:O	1:A:286:GLU:HG3	2.20	0.41
1:A:295:LYS:HB2	1:A:295:LYS:HE3	1.87	0.41
1:A:30:VAL:HG23	1:A:32:SER:H	1.85	0.41
1:A:73:ARG:HD2	1:A:73:ARG:HA	1.81	0.41
1:B:192:GLU:HG2	1:B:193:ILE:H	1.84	0.41
1:C:268:MET:HE1	1:C:273:VAL:HG22	2.01	0.41
1:D:251:SER:O	1:D:252:SER:HB2	2.19	0.41
1:D:325:GLN:C	1:D:327:PHE:H	2.23	0.41
1:A:20:VAL:HA	1:A:90:PHE:CE1	2.55	0.41
1:C:124:ASP:HA	1:C:127:VAL:CG2	2.51	0.41
1:C:75:LEU:HD11	2:C:361:A17:H14	2.02	0.41
1:D:131:ILE:HG13	1:D:213:MET:HE2	2.01	0.41
1:D:109:MET:CB	1:D:157:ALA:HB1	2.50	0.41
1:A:197:TRP:O	1:A:198:MET:O	2.39	0.41
1:A:77:HIS:HD2	1:A:78:MET:CE	2.33	0.41
1:B:291:LEU:HD23	1:B:291:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:ARG:O	1:B:11:GLN:CB	2.68	0.41
1:B:197:TRP:HD1	1:B:198:MET:N	2.19	0.41
1:D:246:LEU:HD22	1:D:292:ASP:HB2	2.01	0.41
1:B:195:LEU:O	1:B:197:TRP:N	2.54	0.41
1:B:228:HIS:O	1:B:231:GLN:HB3	2.20	0.41
1:C:246:LEU:HD11	1:C:292:ASP:CB	2.41	0.41
1:C:232:LEU:HD21	1:C:259:ILE:HD11	2.02	0.41
1:C:71:GLU:HG3	2:C:361:A17:H21A	2.01	0.41
1:D:191:PRO:O	1:D:194:MET:HB3	2.21	0.41
1:B:128:GLN:HG2	1:B:311:TYR:HE2	1.84	0.41
1:B:52:VAL:HA	1:B:104:LEU:O	2.21	0.41
1:D:8:PHE:HB3	1:D:19:GLU:HG2	2.02	0.41
1:D:93:ALA:HB1	1:D:95:SER:O	2.21	0.41
1:C:138:LEU:O	1:C:139:LYS:C	2.59	0.41
1:C:299:ALA:O	1:C:303:LEU:HG	2.20	0.41
1:D:236:LEU:O	1:D:240:GLY:HA2	2.21	0.41
1:D:43:ASP:OD2	1:D:46:THR:OG1	2.33	0.41
1:A:313:ASP:HA	1:A:314:PRO:HD2	1.77	0.41
1:A:53:LYS:HB2	2:A:361:A17:H17B	2.01	0.41
1:C:130:LEU:HD12	1:C:130:LEU:HA	1.80	0.41
1:C:226:THR:H	1:C:230:ASP:HB3	1.86	0.41
1:A:325:GLN:HB2	1:A:325:GLN:HE21	1.69	0.41
1:A:63:ILE:O	1:A:67:ARG:HB2	2.21	0.41
1:B:253:GLU:HA	1:B:256:ARG:NH1	2.35	0.41
1:C:53:LYS:HB3	1:C:104:LEU:HB2	2.03	0.41
1:D:134:ILE:O	1:D:138:LEU:HB2	2.21	0.41
1:D:84:ILE:HB	1:D:167:LEU:CD2	2.51	0.41
1:A:10:ARG:HG3	1:A:19:GLU:HG3	2.03	0.40
1:A:78:MET:HA	1:A:78:MET:HE3	2.02	0.40
1:B:88:ASP:O	1:B:105:VAL:HG23	2.21	0.40
1:D:10:ARG:HD2	1:D:17:ILE:HD11	2.02	0.40
1:D:133:GLN:HB2	1:D:164:LEU:HD22	2.04	0.40
1:B:113:LEU:HA	1:B:113:LEU:HD12	1.86	0.40
1:B:128:GLN:HG2	1:B:311:TYR:CE2	2.56	0.40
1:B:196:ASN:OD1	1:B:196:ASN:O	2.40	0.40
1:B:195:LEU:CD1	1:B:232:LEU:HD22	2.51	0.40
1:D:205:ASP:O	1:D:208:SER:HB2	2.22	0.40
1:D:327:PHE:CD1	1:D:327:PHE:N	2.89	0.40
1:D:23:ARG:CG	1:D:43:ASP:OD1	2.69	0.40
1:B:342:TYR:CE2	1:B:346:ILE:HD11	2.56	0.40
1:B:55:LEU:HA	1:B:55:LEU:HD23	1.92	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:PRO:HD3	1:C:90:PHE:CE1	2.56	0.40
1:A:144:ALA:HB2	1:A:320:ALA:HB3	2.04	0.40
1:B:86:LEU:HD12	1:B:86:LEU:HA	1.84	0.40
1:C:27:LEU:HD13	1:C:39:CYS:HB2	2.03	0.40
1:B:14:ASN:HB3	1:B:15:LYS:H	1.68	0.40
1:B:194:MET:SD	1:B:231:GLN:HG2	2.61	0.40
1:B:35:TYR:O	1:B:53:LYS:HE3	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:VAL:O	1:D:31:GLY:O[2_554]	2.13	0.07

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	339/366 (93%)	293 (86%)	35 (10%)	11 (3%)	4	16
1	B	339/366 (93%)	288 (85%)	39 (12%)	12 (4%)	3	14
1	C	347/366 (95%)	287 (83%)	46 (13%)	14 (4%)	3	11
1	D	347/366 (95%)	267 (77%)	63 (18%)	17 (5%)	2	8
All	All	1372/1464 (94%)	1135 (83%)	183 (13%)	54 (4%)	3	12

All (54) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	PHE
1	A	172	ALA
1	A	198	MET

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Mol	Chain	Res	Type
1	B	11	GLN
1	B	32	SER
1	B	95	SER
1	B	116	ILE
1	C	118	LYS
1	C	174	HIS
1	C	176	ASP
1	D	252	SER
1	D	318	PRO
1	A	120	GLN
1	A	196	ASN
1	B	30	VAL
1	B	173	ARG
1	B	196	ASN
1	C	178	GLU
1	C	252	SER
1	C	331	ASP
1	D	172	ALA
1	D	179	MET
1	D	196	ASN
1	A	122	LEU
1	B	14	ASN
1	B	150	ASP
1	B	228	HIS
1	C	196	ASN
1	C	198	MET
1	C	227	ASP
1	C	253	GLU
1	D	84	ILE
1	D	293	SER
1	D	316	ASP
1	D	326	SER
1	A	184	ALA
1	A	253	GLU
1	C	122	LEU
1	C	332	LEU
1	D	65	ALA
1	D	100	ASN
1	D	184	ALA
1	A	276	GLY
1	B	96	LEU
1	B	229	ILE

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Mol	Chain	Res	Type
1	D	46	THR
1	D	199	HIS
1	C	121	LYS
1	A	116	ILE
1	A	318	PRO
1	C	116	ILE
1	D	250	ILE
1	D	30	VAL
1	D	314	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	304/325 (94%)	286 (94%)	18 (6%)	19	49
1	B	304/325 (94%)	288 (95%)	16 (5%)	22	54
1	C	309/325 (95%)	284 (92%)	25 (8%)	11	33
1	D	309/325 (95%)	291 (94%)	18 (6%)	20	50
All	All	1226/1300 (94%)	1149 (94%)	77 (6%)	18	46

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ARG
1	A	17	ILE
1	A	78	MET
1	A	81	GLU
1	A	89	VAL
1	A	116	ILE
1	A	125	ASP
1	A	149	ARG
1	A	168	ASP
1	A	175	THR
1	A	176	ASP
1	A	198	MET

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Mol	Chain	Res	Type
1	A	199	HIS
1	A	220	ARG
1	A	228	HIS
1	A	237	ARG
1	A	263	THR
1	A	321	ASP
1	B	4	GLU
1	B	11	GLN
1	B	26	ASN
1	B	56	SER
1	B	57	ARG
1	B	118	LYS
1	B	162	CYS
1	B	171	LEU
1	B	186	ARG
1	B	191	PRO
1	B	197	TRP
1	B	237	ARG
1	B	257	ASN
1	B	263	THR
1	B	310	GLN
1	B	316	ASP
1	C	27	LEU
1	C	57	ARG
1	C	60	GLN
1	C	61	SER
1	C	94	ARG
1	C	95	SER
1	C	119	CYS
1	C	121	LYS
1	C	123	THR
1	C	130	LEU
1	C	143	SER
1	C	182	TYR
1	C	191	PRO
1	C	194	MET
1	C	198	MET
1	C	266	PRO
1	C	287	LYS
1	C	291	LEU
1	C	294	ASP
1	C	310	GLN

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Mol	Chain	Res	Type
1	C	321	ASP
1	C	333	LEU
1	C	335	ASP
1	C	339	SER
1	C	349	VAL
1	D	17	ILE
1	D	42	PHE
1	D	75	LEU
1	D	78	MET
1	D	125	ASP
1	D	126	HIS
1	D	168	ASP
1	D	177	ASP
1	D	179	MET
1	D	182	TYR
1	D	192	GLU
1	D	198	MET
1	D	241	THR
1	D	247	LEU
1	D	258	TYR
1	D	264	GLN
1	D	279	PRO
1	D	285	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	64	HIS
1	A	77	HIS
1	A	120	GLN
1	A	128	GLN
1	A	174	HIS
1	A	202	GLN
1	A	264	GLN
1	A	269	ASN
1	A	301	GLN
1	A	325	GLN
1	B	60	GLN
1	B	77	HIS
1	B	120	GLN
1	B	196	ASN

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Mol	Chain	Res	Type
1	B	202	GLN
1	B	257	ASN
1	B	301	GLN
1	B	310	GLN
1	C	14	ASN
1	C	60	GLN
1	C	64	HIS
1	C	114	ASN
1	C	120	GLN
1	C	126	HIS
1	C	128	GLN
1	C	202	GLN
1	C	231	GLN
1	C	257	ASN
1	C	272	ASN
1	C	301	GLN
1	C	310	GLN
1	D	11	GLN
1	D	14	ASN
1	D	64	HIS
1	D	107	HIS
1	D	128	GLN
1	D	202	GLN
1	D	269	ASN
1	D	301	GLN
1	D	310	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 carbohydrates in this entry.

5.6 Ligand geometry

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	A17	A	361	-	34,34,34	3.47	24 (70%)	48,49,49	2.62	19 (39%)
2	A17	D	361	-	34,34,34	3.87	25 (73%)	48,49,49	2.34	20 (41%)
2	A17	B	361	-	34,34,34	3.72	24 (70%)	48,49,49	2.16	17 (35%)
2	A17	C	361	-	34,34,34	3.78	23 (67%)	48,49,49	2.65	15 (31%)

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	A17	A	361	-	-	2/16/18/18	0/5/5/5
2	A17	D	361	-	-	5/16/18/18	0/5/5/5
2	A17	B	361	-	-	2/16/18/18	0/5/5/5
2	A17	C	361	-	-	1/16/18/18	0/5/5/5

All (96) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	361	A17	N7-N8	14.38	1.53	1.35
2	D	361	A17	N7-N8	13.90	1.52	1.35
2	B	361	A17	N7-N8	12.53	1.50	1.35
2	A	361	A17	N7-N8	10.56	1.48	1.35
2	B	361	A17	C9-N8	5.95	1.42	1.32
2	A	361	A17	C9-N8	5.81	1.42	1.32
2	D	361	A17	C9-N8	5.73	1.42	1.32
2	B	361	A17	C10-N7	5.47	1.42	1.31
2	C	361	A17	C9-N8	5.37	1.41	1.32
2	A	361	A17	C10-N7	5.20	1.41	1.31
2	D	361	A17	C10-N7	5.18	1.41	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	361	A17	C3-C4	5.01	1.47	1.36
2	B	361	A17	C3-C4	4.97	1.47	1.36
2	C	361	A17	C10-N7	4.96	1.41	1.31
2	A	361	A17	C3-C4	4.62	1.46	1.36
2	C	361	A17	C14-C15	4.52	1.47	1.38
2	A	361	A17	C14-C15	4.40	1.46	1.38
2	C	361	A17	C3-C4	4.34	1.45	1.36
2	B	361	A17	C23-C9	-4.21	1.44	1.49
2	D	361	A17	C12-C13	4.19	1.45	1.39
2	A	361	A17	C12-C13	4.09	1.45	1.39
2	D	361	A17	C14-C15	4.09	1.46	1.38
2	A	361	A17	C24-C23	4.04	1.46	1.40
2	D	361	A17	C24-C23	3.92	1.46	1.40
2	B	361	A17	C12-C13	3.89	1.45	1.39
2	D	361	A17	C14-C13	3.79	1.45	1.39
2	B	361	A17	C14-C15	3.79	1.45	1.38
2	B	361	A17	C4-C5	3.77	1.47	1.39
2	A	361	A17	C14-C13	3.75	1.45	1.39
2	C	361	A17	C12-C13	3.75	1.45	1.39
2	D	361	A17	C12-C11	3.74	1.46	1.39
2	C	361	A17	C23-C9	-3.72	1.44	1.49
2	B	361	A17	C12-C11	3.61	1.45	1.39
2	C	361	A17	C14-C13	3.59	1.45	1.39
2	D	361	A17	C4-C5	3.54	1.46	1.39
2	A	361	A17	C12-C11	3.49	1.45	1.39
2	D	361	A17	C26-C27	3.48	1.46	1.38
2	B	361	A17	C14-C13	3.46	1.45	1.39
2	C	361	A17	C12-C11	3.40	1.45	1.39
2	D	361	A17	C23-C28	3.39	1.46	1.40
2	A	361	A17	C15-C16	3.37	1.47	1.39
2	A	361	A17	C25-C24	3.36	1.46	1.38
2	A	361	A17	C11-C16	3.35	1.46	1.40
2	C	361	A17	C24-C23	3.35	1.45	1.40
2	D	361	A17	C25-C24	3.34	1.46	1.38
2	B	361	A17	C25-C24	3.30	1.45	1.38
2	A	361	A17	C23-C28	3.30	1.46	1.40
2	B	361	A17	C26-C27	3.29	1.45	1.38
2	B	361	A17	C11-C16	3.29	1.46	1.40
2	B	361	A17	C24-C23	3.29	1.45	1.40
2	C	361	A17	C25-C24	3.27	1.45	1.38
2	C	361	A17	C26-C27	3.24	1.45	1.38
2	C	361	A17	C15-C16	3.22	1.46	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	361	A17	C6-C5	3.20	1.46	1.38
2	D	361	A17	C6-C5	3.19	1.46	1.38
2	C	361	A17	C27-C28	3.17	1.46	1.39
2	C	361	A17	C11-C5	-3.16	1.43	1.49
2	B	361	A17	C27-C28	3.12	1.46	1.39
2	D	361	A17	C23-C9	-3.12	1.45	1.49
2	D	361	A17	C11-C16	3.12	1.46	1.40
2	D	361	A17	C27-C28	3.11	1.46	1.39
2	A	361	A17	C26-C27	3.09	1.45	1.38
2	D	361	A17	C26-C25	3.02	1.46	1.38
2	A	361	A17	C27-C28	3.01	1.46	1.39
2	B	361	A17	C26-C25	2.99	1.45	1.38
2	D	361	A17	C15-C16	2.96	1.46	1.39
2	B	361	A17	C15-C16	2.93	1.46	1.39
2	B	361	A17	C17-C16	-2.87	1.45	1.51
2	C	361	A17	C26-C25	2.86	1.45	1.38
2	C	361	A17	C6-C5	2.85	1.45	1.38
2	B	361	A17	C23-C28	2.81	1.45	1.40
2	C	361	A17	C11-C16	2.74	1.45	1.40
2	C	361	A17	C23-C28	2.73	1.45	1.40
2	C	361	A17	C4-C5	2.70	1.45	1.39
2	A	361	A17	C4-C5	2.69	1.45	1.39
2	A	361	A17	C26-C25	2.69	1.45	1.38
2	A	361	A17	C11-C5	-2.67	1.44	1.49
2	A	361	A17	C6-C5	2.60	1.44	1.38
2	D	361	A17	C17-C16	-2.60	1.45	1.51
2	B	361	A17	C20-N30	-2.56	1.41	1.46
2	C	361	A17	C17-C16	-2.48	1.46	1.51
2	A	361	A17	C23-C9	-2.45	1.46	1.49
2	A	361	A17	C3-C2	2.45	1.47	1.42
2	A	361	A17	C20-N30	-2.44	1.41	1.46
2	D	361	A17	C20-N30	-2.41	1.41	1.46
2	C	361	A17	C29-C28	-2.40	1.46	1.51
2	D	361	A17	C29-C28	-2.39	1.46	1.51
2	B	361	A17	C2-C1	2.36	1.47	1.43
2	D	361	A17	C2-C1	2.34	1.47	1.43
2	A	361	A17	C29-C28	-2.34	1.46	1.51
2	D	361	A17	C3-C2	2.34	1.47	1.42
2	A	361	A17	C17-C16	-2.33	1.46	1.51
2	B	361	A17	C3-C2	2.28	1.46	1.42
2	B	361	A17	C29-C28	-2.25	1.46	1.51
2	D	361	A17	C11-C5	-2.25	1.45	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	361	A17	C20-N30	-2.18	1.42	1.46

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	361	A17	C1-C10-N7	-9.31	114.82	124.56
2	D	361	A17	C1-C10-N7	-7.45	116.77	124.56
2	A	361	A17	C1-C10-N7	-6.66	117.59	124.56
2	A	361	A17	C22-C20-N30	-6.55	109.22	118.61
2	C	361	A17	C9-C2-C1	6.20	121.06	117.60
2	C	361	A17	C10-C1-C2	5.82	123.37	117.78
2	A	361	A17	C10-C1-C2	5.48	123.04	117.78
2	A	361	A17	C20-N30-C18	5.27	129.84	122.55
2	D	361	A17	C24-C23-C9	-5.03	113.91	120.14
2	B	361	A17	C15-C14-C13	-5.02	114.94	120.78
2	C	361	A17	C23-C9-C2	4.62	128.66	122.23
2	D	361	A17	C10-C1-C2	4.55	122.14	117.78
2	C	361	A17	C2-C9-N8	-4.50	117.67	123.00
2	B	361	A17	C1-C10-N7	-4.35	120.01	124.56
2	B	361	A17	C23-C9-N8	-4.33	109.54	116.62
2	A	361	A17	C15-C14-C13	-4.21	115.88	120.78
2	B	361	A17	C9-N8-N7	-4.15	115.87	120.41
2	B	361	A17	C14-C13-C12	4.14	124.13	119.24
2	A	361	A17	C4-C5-C11	4.13	127.59	120.91
2	D	361	A17	C3-C2-C1	-4.04	112.67	117.89
2	A	361	A17	C3-C2-C9	3.90	128.76	123.57
2	A	361	A17	C21-C20-N30	-3.87	113.06	118.61
2	C	361	A17	C3-C2-C1	-3.80	112.97	117.89
2	B	361	A17	C24-C23-C9	-3.77	115.46	120.14
2	B	361	A17	C17-C16-C15	-3.68	113.13	120.31
2	D	361	A17	C23-C9-C2	3.57	127.20	122.23
2	A	361	A17	C4-C3-C2	3.56	126.08	121.13
2	A	361	A17	C3-C2-C1	-3.48	113.39	117.89
2	C	361	A17	C6-C1-C10	-3.40	112.65	121.75
2	C	361	A17	C6-C1-C2	3.40	123.72	119.25
2	D	361	A17	C9-C2-C1	3.33	119.46	117.60
2	C	361	A17	C4-C3-C2	3.28	125.69	121.13
2	D	361	A17	C15-C14-C13	-3.25	116.99	120.78
2	D	361	A17	C3-C2-C9	3.23	127.87	123.57
2	C	361	A17	C12-C11-C5	-3.20	112.62	118.58
2	B	361	A17	C6-C1-C2	3.19	123.45	119.25
2	A	361	A17	C24-C23-C9	-3.14	116.25	120.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	361	A17	C4-C3-C2	3.10	125.44	121.13
2	B	361	A17	C3-C2-C9	3.07	127.65	123.57
2	C	361	A17	C10-N7-N8	2.96	123.63	119.23
2	B	361	A17	C23-C9-C2	2.92	126.30	122.23
2	C	361	A17	C22-C20-N30	2.92	122.80	118.61
2	A	361	A17	C6-C5-C11	-2.90	116.25	120.72
2	D	361	A17	C10-N7-N8	2.78	123.37	119.23
2	D	361	A17	C29-C28-C27	-2.75	114.94	120.31
2	D	361	A17	C4-C5-C11	2.71	125.30	120.91
2	D	361	A17	C6-C1-C2	2.70	122.80	119.25
2	D	361	A17	C23-C9-N8	-2.69	112.22	116.62
2	A	361	A17	C12-C13-C18	-2.69	111.52	120.44
2	B	361	A17	C10-C1-C2	2.61	120.28	117.78
2	A	361	A17	C25-C26-C27	-2.61	116.21	120.19
2	B	361	A17	C3-C2-C1	-2.58	114.55	117.89
2	B	361	A17	C21-C20-N30	-2.55	114.96	118.61
2	D	361	A17	C6-C1-C10	-2.50	115.06	121.75
2	C	361	A17	C12-C11-C16	2.47	122.67	119.02
2	D	361	A17	C9-N8-N7	-2.46	117.72	120.41
2	B	361	A17	C25-C26-C27	-2.46	116.44	120.19
2	B	361	A17	C20-N30-C18	-2.45	119.15	122.55
2	A	361	A17	C2-C9-N8	-2.40	120.16	123.00
2	C	361	A17	C21-C20-N30	-2.36	115.22	118.61
2	C	361	A17	C15-C16-C11	-2.36	115.00	118.44
2	A	361	A17	C12-C11-C5	-2.32	114.26	118.58
2	A	361	A17	C5-C6-C1	2.31	125.54	121.53
2	B	361	A17	C6-C1-C10	-2.12	116.08	121.75
2	D	361	A17	C17-C16-C15	-2.12	116.18	120.31
2	D	361	A17	C25-C26-C27	-2.10	117.00	120.19
2	B	361	A17	C17-C16-C11	2.09	125.89	122.33
2	D	361	A17	C2-C9-N8	-2.09	120.53	123.00
2	A	361	A17	C5-C11-C16	2.09	126.73	122.48
2	A	361	A17	C6-C1-C10	-2.03	116.32	121.75
2	D	361	A17	C29-C28-C23	2.03	125.78	122.33

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	361	A17	C16-C11-C5-C4
2	D	361	A17	C16-C11-C5-C6
2	D	361	A17	C28-C23-C9-N8

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Mol	Chain	Res	Type	Atoms
2	B	361	A17	C28-C23-C9-N8
2	C	361	A17	C28-C23-C9-N8
2	D	361	A17	C12-C11-C5-C4
2	D	361	A17	C12-C11-C5-C6
2	A	361	A17	C12-C11-C5-C4
2	A	361	A17	C16-C11-C5-C4
2	B	361	A17	C16-C11-C5-C4

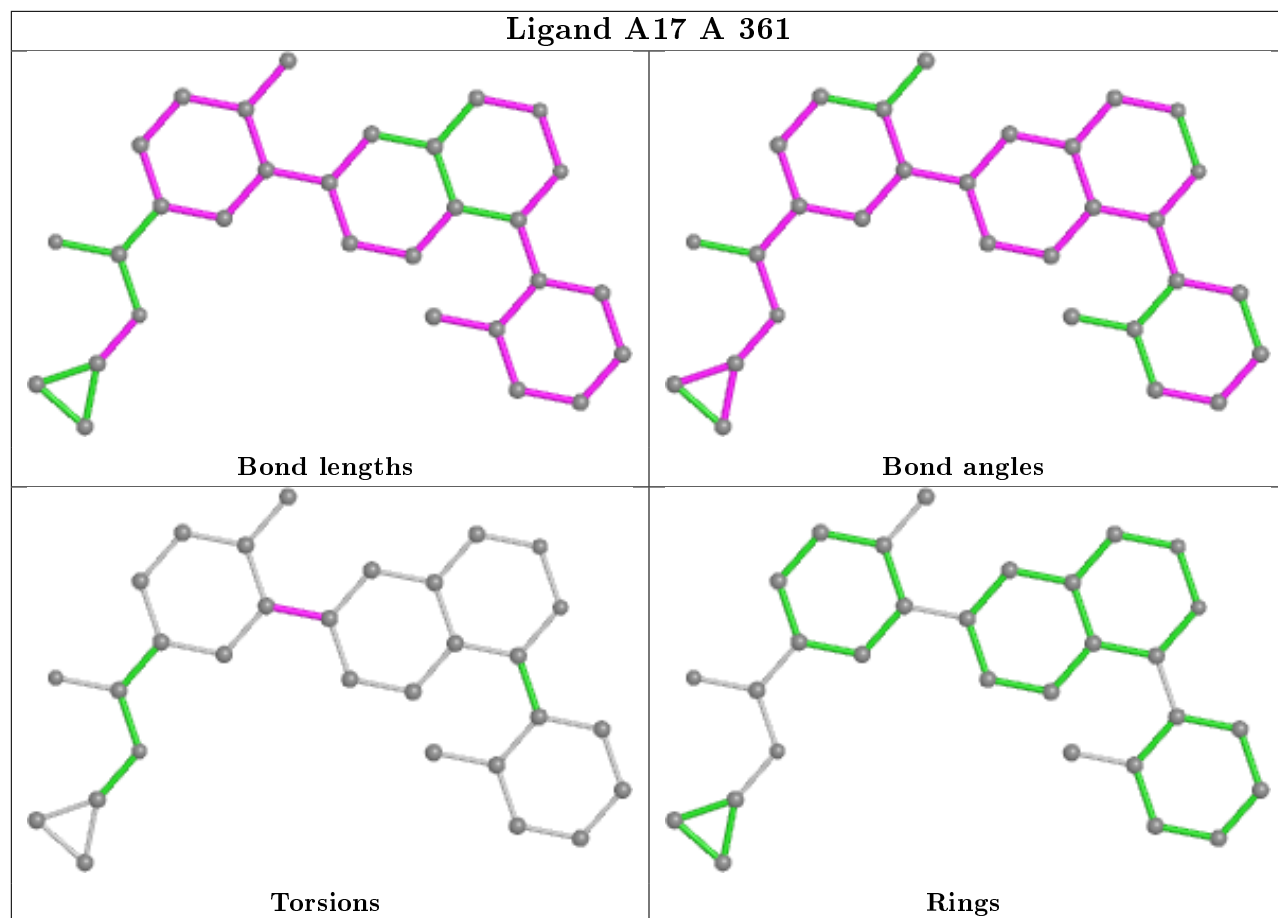
There are no ring outliers.

4 monomers are involved in 11 short contacts:

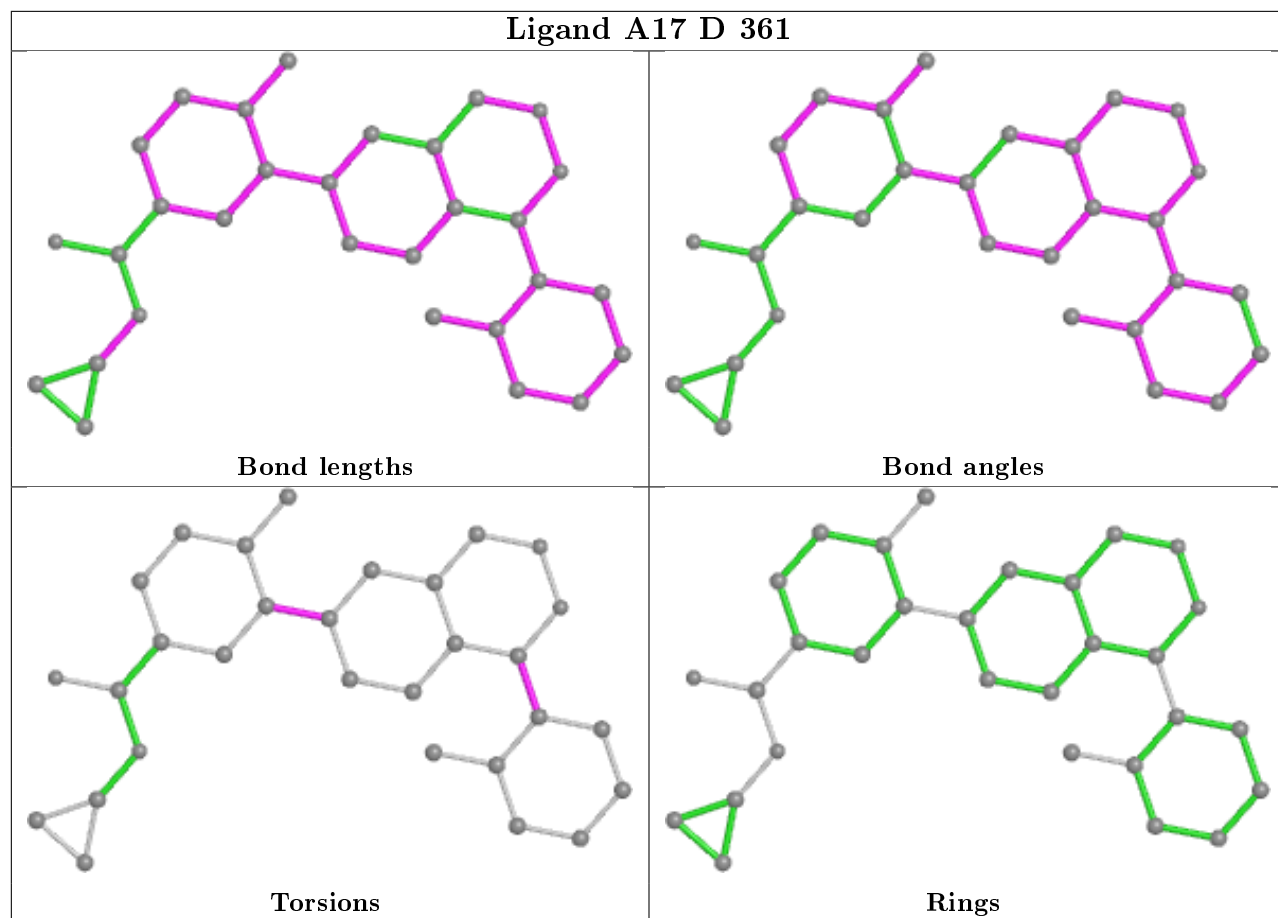
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	361	A17	4	0
2	D	361	A17	3	0
2	B	361	A17	1	0
2	C	361	A17	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

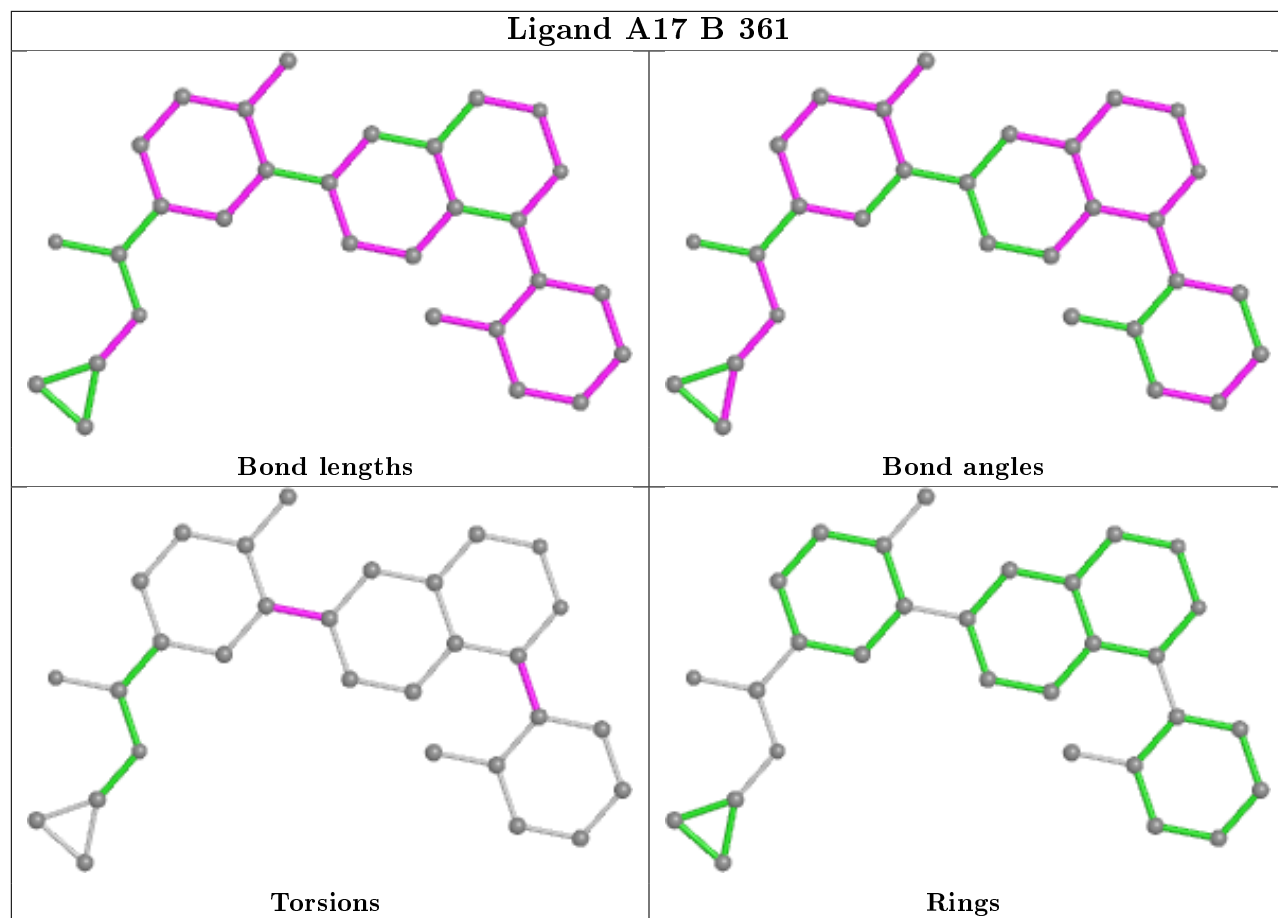
Ligand A17 A 361

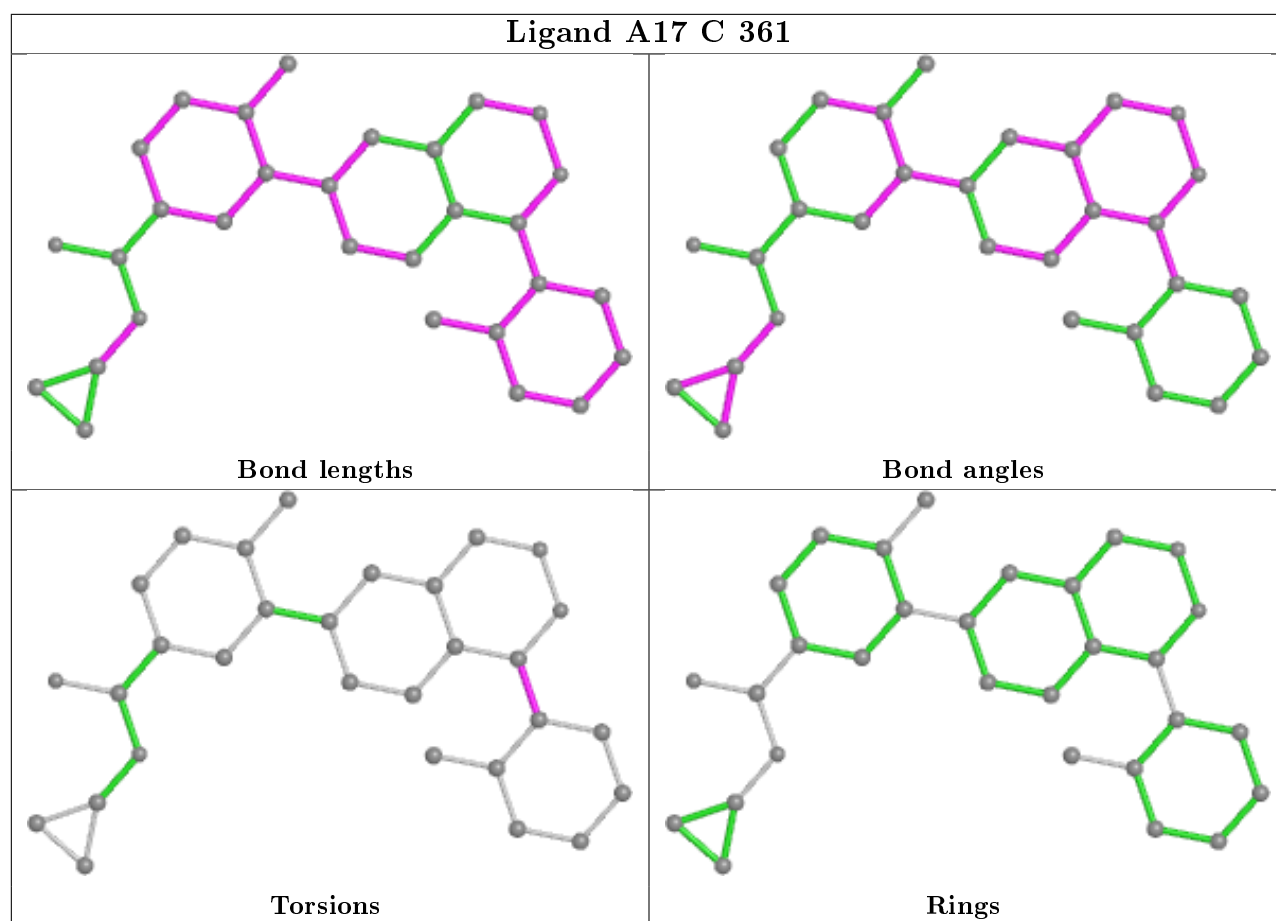


Ligand A17 D 361



Ligand A17 B 361





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	343/366 (93%)	-0.46	1 (0%) 94 94	18, 35, 69, 95	0
1	B	343/366 (93%)	-0.35	4 (1%) 79 79	16, 43, 74, 89	0
1	C	349/366 (95%)	-0.21	8 (2%) 60 58	17, 43, 83, 104	0
1	D	349/366 (95%)	0.13	24 (6%) 16 13	31, 64, 95, 110	0
All	All	1384/1464 (94%)	-0.22	37 (2%) 54 50	16, 46, 83, 110	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	181	GLY	7.8
1	D	174	HIS	7.7
1	D	180	THR	7.4
1	C	179	MET	7.2
1	C	183	VAL	6.5
1	D	197	TRP	5.1
1	C	184	ALA	4.6
1	C	182	TYR	4.6
1	D	182	TYR	4.4
1	D	255	ALA	4.1
1	B	119	CYS	4.0
1	C	251	SER	3.7
1	D	175	THR	3.6
1	D	261	SER	3.5
1	C	175	THR	3.5
1	D	178	GLU	3.5
1	D	254	SER	3.4
1	B	352	PRO	3.3
1	D	260	GLN	3.2
1	B	251	SER	3.2
1	D	320	ALA	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	172	ALA	3.0
1	D	252	SER	2.8
1	D	258	TYR	2.6
1	D	253	GLU	2.6
1	D	334	ILE	2.5
1	C	177	ASP	2.5
1	D	179	MET	2.4
1	D	199	HIS	2.4
1	C	252	SER	2.4
1	D	176	ASP	2.3
1	D	251	SER	2.2
1	D	257	ASN	2.2
1	D	183	VAL	2.1
1	A	32	SER	2.1
1	B	176	ASP	2.1
1	D	121	LYS	2.1

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 carbohydrates in this entry.

6.4 Ligands [i](#)

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

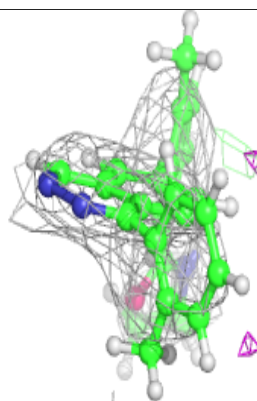
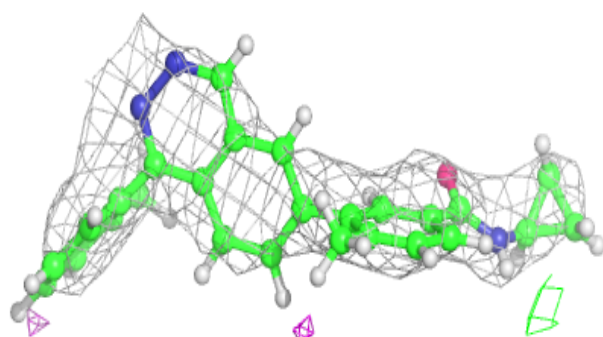
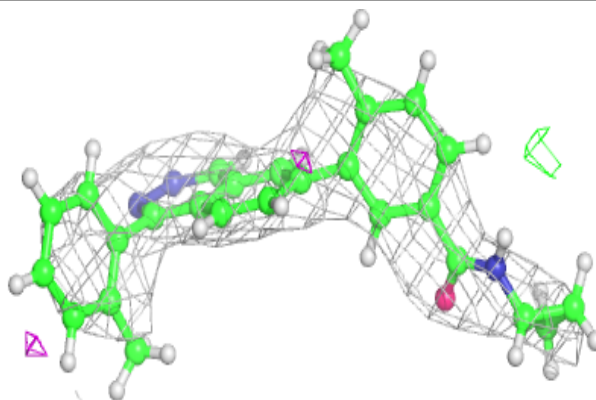
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	A17	D	361	30/30	0.91	0.29	85,87,89,89	23
2	A17	B	361	30/30	0.95	0.17	16,21,34,35	23
2	A17	A	361	30/30	0.96	0.16	21,24,32,32	23
2	A17	C	361	30/30	0.96	0.15	14,18,21,21	23

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.

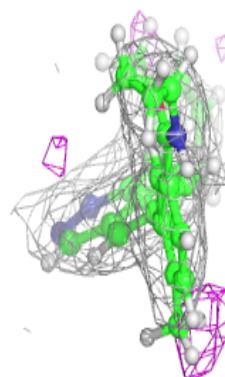
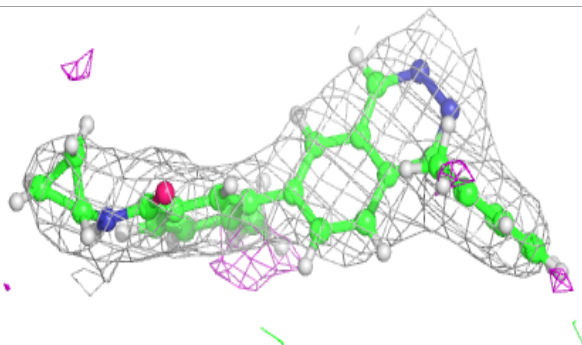
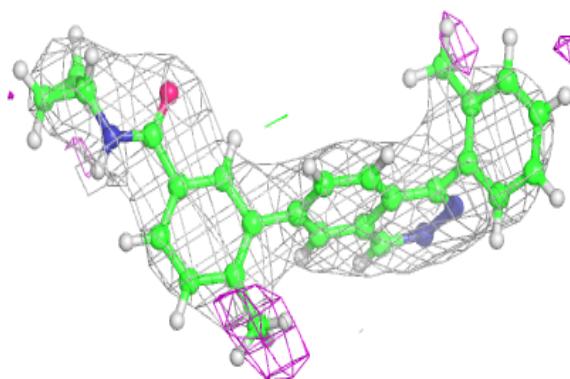
Electron density around A17 D 361:

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



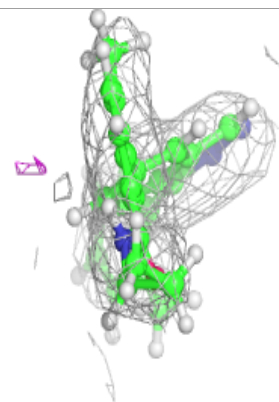
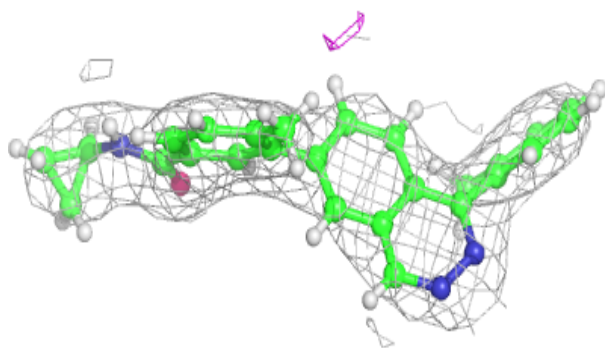
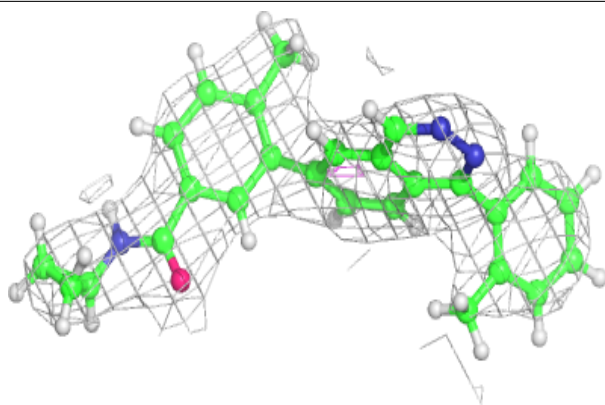
Electron density around A17 B 361:

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

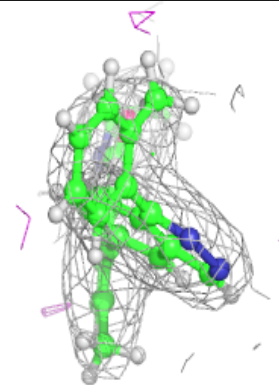
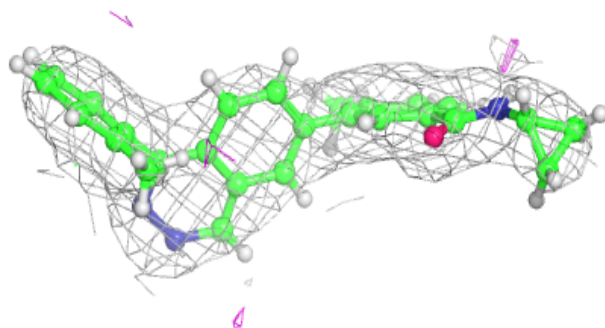
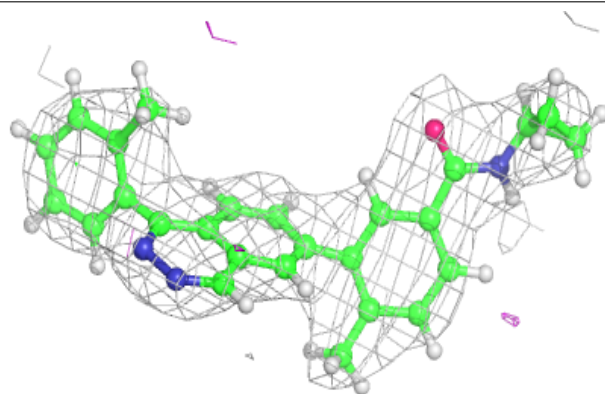


Electron density around A17 A 361:

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

**Electron density around A17 C 361:**

$2mF_o-DF_c$ (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.