



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 04:55 am BST

PDB ID : 2C5N  
Title : Differential Binding Of Inhibitors To Active And Inactive Cdk2 Provides Insights For Drug Design  
Authors : Kontopidis, G.; McInnes, C.; Pandalaneni, S.R.; McNae, I.; Gibson, D.; Mezna, M.; Thomas, M.; Wood, G.; Wang, S.; Walkinshaw, M.D.; Fischer, P.M.  
Deposited on : 2005-10-30  
Resolution : 2.10 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](mailto: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.

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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

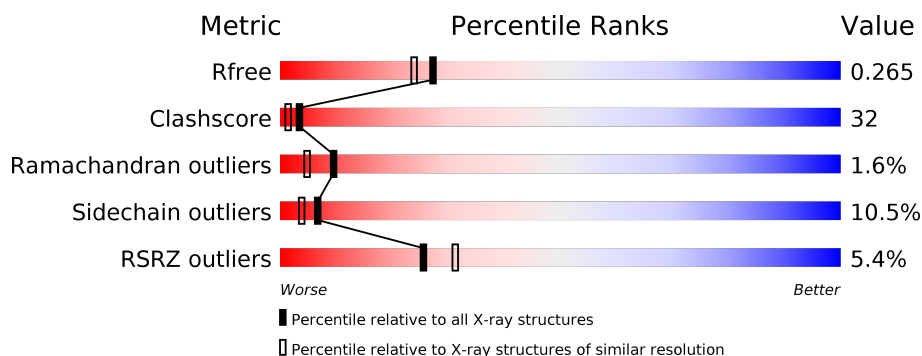
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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  $\geq 3$ , 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	298	<div> <div>6%</div> <div> <div></div> <div>64%</div> <div>26%</div> <div>7%</div> <div>..</div> </div> </div>
1	C	298	<div> <div>8%</div> <div> <div></div> <div>60%</div> <div>30%</div> <div>8%</div> <div>.</div> </div> </div>
2	B	259	<div> <div>5%</div> <div> <div></div> <div>63%</div> <div>28%</div> <div>8%</div> <div>.</div> </div> </div>
2	D	259	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>20%</div> <div>.</div> </div> </div>

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
3	CK8	A	1297[A]	-	-	X	-
3	CK8	A	1297[B]	-	-	X	-
3	CK8	C	1298	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9928 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.

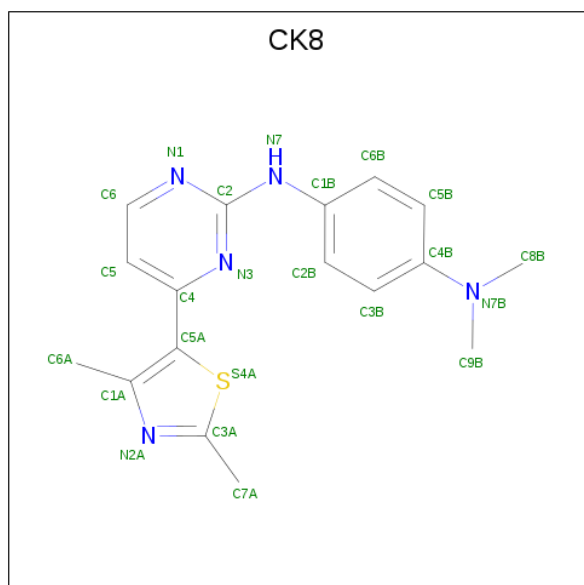
- Molecule 1 is a protein called CELL DIVISION PROTEIN KINASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2378	1547	403	420	8			
1	C	297	Total	C	N	O	S	0	0	1
			2379	1547	404	420	8			

- Molecule 2 is a protein called CYCLIN A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	258	Total	C	N	O	S	0	0	0
			2083	1350	339	383	11			
2	D	258	Total	C	N	O	S	0	0	0
			2084	1350	339	384	11			

- Molecule 3 is N-[4-(2,4-DIMETHYL-THIAZOL-5-YL)-PYRIMIDIN-2-YL]-N',N'-DIMETHYL-BENZENE-1,4-DIAMINE (three-letter code: CK8) (formula: C<sub>17</sub>H<sub>19</sub>N<sub>5</sub>S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	S	0	1
			46	34	10	2		
3	C	1	Total	C	N	S	0	0
			23	17	5	1		

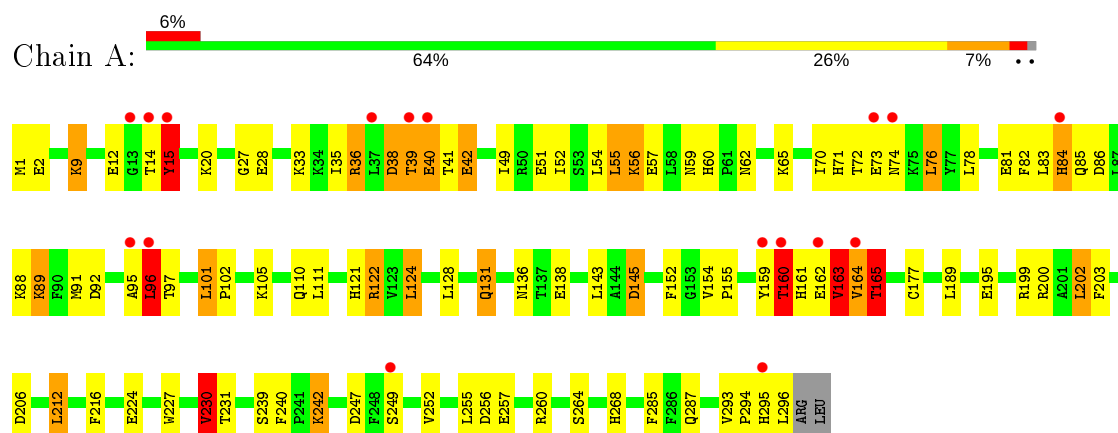
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	244	Total	O	0	0
			244	244		
4	B	188	Total	O	0	0
			188	188		
4	C	258	Total	O	0	0
			258	258		
4	D	245	Total	O	0	0
			245	245		

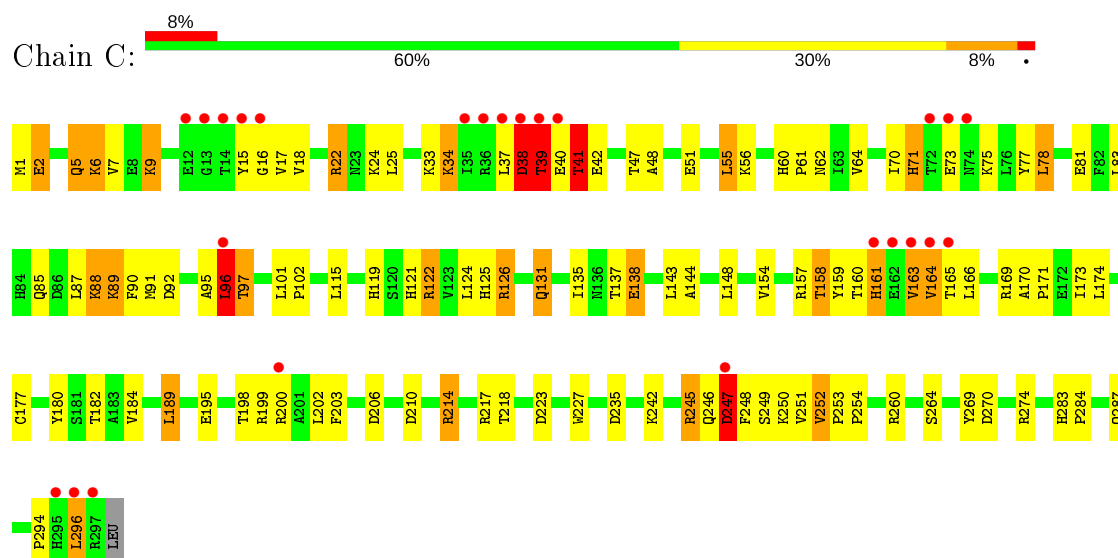
### 3 Residue-property plots [i](#)

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.

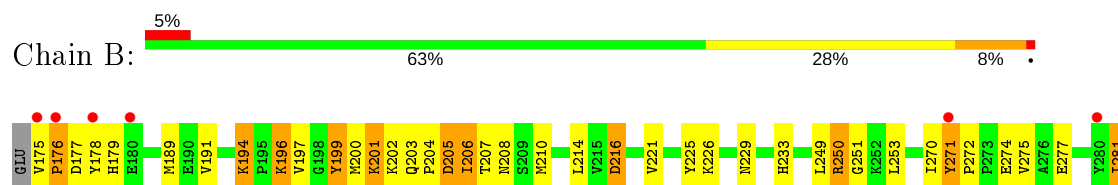
#### • Molecule 1: CELL DIVISION PROTEIN KINASE 2

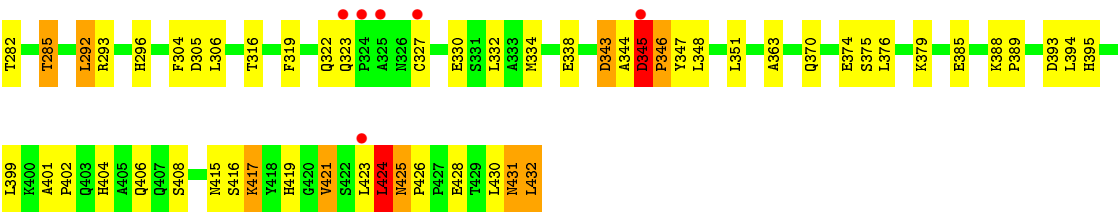


#### • Molecule 1: CELL DIVISION PROTEIN KINASE 2

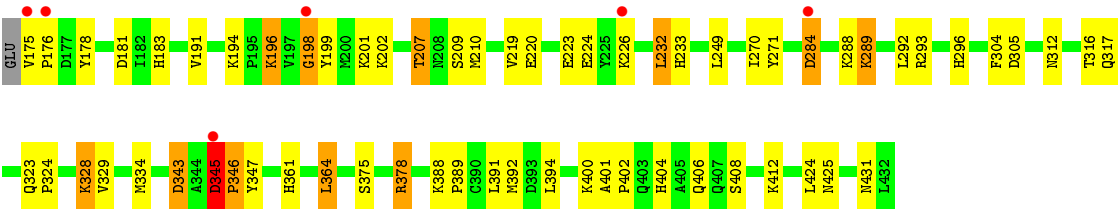
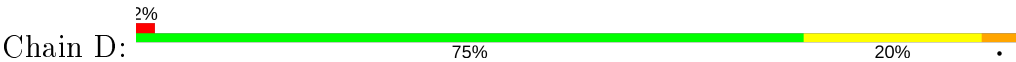


#### • Molecule 2: CYCLIN A2





● Molecule 2: CYCLIN A2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.54Å 113.14Å 157.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.10 14.99 – 2.10	Depositor EDS
% Data completeness (in resolution range)	91.7 (15.00-2.10) 91.7 (14.99-2.10)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.01 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, $R_{free}$	0.175 , 0.259 0.189 , 0.265	Depositor DCC
$R_{free}$ test set	2158 reflections (3.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.0	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 55.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9928	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CK8

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	0.56	0/2440	0.96	5/3313 (0.2%)
1	C	0.55	0/2441	0.96	6/3315 (0.2%)
2	B	0.54	0/2133	0.94	7/2896 (0.2%)
2	D	0.55	0/2134	0.89	3/2897 (0.1%)
All	All	0.55	0/9148	0.94	21/12421 (0.2%)

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	2
2	B	0	1
2	D	0	3
All	All	0	6

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	ASP	CB-CG-OD2	9.73	127.06	118.30
2	B	393	ASP	CB-CG-OD2	8.64	126.08	118.30
2	B	345	ASP	CB-CG-OD2	8.62	126.06	118.30
1	C	245	ARG	NE-CZ-NH2	-8.61	116.00	120.30
2	B	305	ASP	CB-CG-OD2	8.08	125.57	118.30
1	C	235	ASP	CB-CG-OD2	7.02	124.62	118.30
1	C	92	ASP	CB-CG-OD2	6.73	124.36	118.30
1	A	92	ASP	CB-CG-OD2	6.58	124.22	118.30
2	D	181	ASP	CB-CG-OD2	6.38	124.05	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	205	ASP	CB-CG-OD2	6.29	123.96	118.30
2	D	284	ASP	CB-CG-OD2	6.11	123.80	118.30
1	C	247	ASP	CB-CG-OD2	6.10	123.79	118.30
1	A	230	VAL	CB-CA-C	-5.86	100.26	111.40
1	C	274	ARG	NE-CZ-NH2	-5.82	117.39	120.30
2	B	345	ASP	N-CA-C	5.57	126.04	111.00
2	B	216	ASP	CB-CG-OD2	5.37	123.14	118.30
2	B	343	ASP	CB-CG-OD2	5.30	123.07	118.30
1	C	38	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	124	LEU	CB-CA-C	-5.18	100.35	110.20
2	D	343	ASP	CB-CG-OD2	5.17	122.95	118.30
1	A	206	ASP	CB-CG-OD2	5.16	122.94	118.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	15	TYR	Peptide
1	A	163	VAL	Peptide
2	B	345	ASP	Peptide
2	D	198	GLY	Peptide
2	D	345	ASP	Mainchain,Peptide

## 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	2378	0	2425	171	0
1	C	2379	0	2426	201	1
2	B	2083	0	2107	140	1
2	D	2084	0	2107	82	0
3	A	46	0	38	23	0
3	C	23	0	19	11	0
4	A	244	0	0	65	0
4	B	188	0	0	46	0
4	C	258	0	0	61	1
4	D	245	0	0	18	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	9928	0	9122	585	2

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 32.

All (585) 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:C:15:TYR:CD2	1:C:33:LYS:HE2	1.61	1.34
2:D:334:MET:HE1	4:D:2102:HOH:O	1.21	1.31
1:C:206:ASP:HB2	4:C:2186:HOH:O	1.28	1.31
2:B:425:ASN:HB2	4:B:2185:HOH:O	1.19	1.30
1:C:15:TYR:CE2	1:C:33:LYS:HE2	1.66	1.29
2:D:284:ASP:HB3	4:D:2139:HOH:O	1.17	1.28
1:C:177:CYS:HB2	4:C:2164:HOH:O	1.23	1.28
2:B:178:TYR:HB2	4:B:2033:HOH:O	1.14	1.27
2:B:281:ILE:HB	4:B:2097:HOH:O	1.29	1.26
1:C:71:HIS:CE1	2:D:296:HIS:NE2	2.03	1.25
2:D:345:ASP:OD2	2:D:346:PRO:HD3	1.36	1.25
1:A:131:GLN:HG2	4:A:2125:HOH:O	1.28	1.23
1:C:71:HIS:CE1	2:D:296:HIS:HE2	1.56	1.23
1:C:173:ILE:HG13	4:C:2144:HOH:O	1.05	1.20
1:C:64:VAL:HB	4:C:2070:HOH:O	1.04	1.19
1:A:268:HIS:CE1	4:A:2214:HOH:O	1.99	1.15
1:A:38:ASP:OD2	1:A:41:THR:HB	1.46	1.15
1:A:85:GLN:OE1	1:A:296:LEU:HG	1.45	1.14
2:B:327:CYS:HB3	4:B:2125:HOH:O	1.45	1.14
2:B:177:ASP:HB3	4:B:2034:HOH:O	1.48	1.14
1:C:64:VAL:HG23	1:C:143:LEU:O	1.45	1.14
1:C:163:VAL:HG12	4:C:2150:HOH:O	1.45	1.13
1:C:6:LYS:HB3	1:C:6:LYS:HZ2	0.99	1.12
1:C:64:VAL:HG21	1:C:144:ALA:HB2	1.33	1.11
1:C:248:PHE:CD1	4:C:2204:HOH:O	2.02	1.10
1:A:84:HIS:NE2	1:A:296:LEU:HD23	1.65	1.09
1:C:71:HIS:NE2	4:C:2074:HOH:O	1.84	1.09
1:C:248:PHE:CG	4:C:2204:HOH:O	2.01	1.08
1:C:71:HIS:CD2	4:C:2074:HOH:O	2.05	1.08
1:A:131:GLN:CG	4:A:2125:HOH:O	1.82	1.08
1:A:95:ALA:O	1:A:96:LEU:HB3	1.30	1.07
2:B:316:THR:HG22	4:B:2119:HOH:O	1.55	1.06
1:A:42:GLU:OE2	2:B:275:VAL:HG23	1.54	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:HIS:ND1	2:D:296:HIS:NE2	2.01	1.05
2:D:378:ARG:HG2	2:D:378:ARG:HH11	1.20	1.05
2:B:177:ASP:HB2	4:B:2035:HOH:O	1.57	1.04
1:A:177:CYS:HB2	4:A:2157:HOH:O	0.86	1.04
2:B:327:CYS:HB3	4:B:2018:HOH:O	1.57	1.03
1:A:84:HIS:CE1	1:A:296:LEU:HD23	1.93	1.03
1:C:126:ARG:NE	4:C:2131:HOH:O	1.91	1.03
1:C:6:LYS:NZ	1:C:6:LYS:HB3	1.73	1.02
1:A:95:ALA:O	1:A:96:LEU:CB	2.02	1.01
2:D:404:HIS:HD2	2:D:406:GLN:H	1.01	1.00
1:C:15:TYR:CE2	1:C:33:LYS:CE	2.44	1.00
2:D:334:MET:CE	4:D:2102:HOH:O	1.84	1.00
1:A:131:GLN:CD	4:A:2125:HOH:O	1.92	0.99
1:A:2:GLU:HG3	4:A:2005:HOH:O	1.60	0.98
1:A:84:HIS:CE1	1:A:296:LEU:CD2	2.47	0.98
1:C:131:GLN:HE21	1:C:131:GLN:H	1.05	0.98
2:D:345:ASP:OD2	2:D:346:PRO:CD	2.09	0.98
1:C:160:THR:HG23	4:C:2146:HOH:O	1.61	0.97
2:B:404:HIS:HD2	2:B:406:GLN:H	1.09	0.97
2:B:207:THR:HG23	2:B:210:MET:H	1.27	0.97
1:A:42:GLU:CB	4:A:2049:HOH:O	2.15	0.95
2:B:221:VAL:CG2	2:B:281:ILE:HD12	1.98	0.94
3:A:1297[B]:CK8:H6B	3:A:1297[B]:CK8:N3	1.81	0.94
1:C:39:THR:HG23	1:C:39:THR:O	1.66	0.93
1:C:71:HIS:CE1	2:D:296:HIS:CD2	2.56	0.93
1:C:126:ARG:CZ	4:C:2131:HOH:O	2.15	0.93
1:A:256:ASP:OD1	4:A:2200:HOH:O	1.87	0.93
2:B:175:VAL:O	2:B:177:ASP:N	2.01	0.93
1:A:36:ARG:NH1	4:A:2045:HOH:O	1.94	0.93
1:C:64:VAL:HG21	1:C:144:ALA:CB	2.00	0.92
2:D:296:HIS:CD2	4:D:2149:HOH:O	2.20	0.92
2:D:194:LYS:HE2	4:D:2070:HOH:O	1.68	0.92
2:B:419:HIS:O	4:B:2180:HOH:O	1.87	0.91
1:C:6:LYS:CB	1:C:6:LYS:HZ2	1.82	0.91
1:A:42:GLU:OE1	4:A:2053:HOH:O	1.86	0.91
1:A:56:LYS:HG3	4:B:2111:HOH:O	1.70	0.91
1:A:42:GLU:HB2	4:A:2049:HOH:O	1.70	0.90
2:B:206:ILE:HA	2:B:210:MET:CE	2.02	0.90
1:A:38:ASP:OD2	1:A:41:THR:CB	2.21	0.89
2:B:177:ASP:CB	4:B:2034:HOH:O	2.13	0.89
1:C:137:THR:OG1	1:C:138:GLU:OE2	1.91	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:214:ARG:CG	1:C:214:ARG:HH11	1.85	0.89
1:A:60:HIS:CD2	1:A:62:ASN:H	1.91	0.88
2:B:221:VAL:HG22	2:B:281:ILE:HD12	1.54	0.88
2:D:404:HIS:CD2	2:D:406:GLN:H	1.91	0.88
1:C:165:THR:HB	4:C:2155:HOH:O	1.72	0.88
1:A:240:PHE:HB2	4:A:2190:HOH:O	1.72	0.88
1:A:84:HIS:HB3	4:A:2093:HOH:O	1.72	0.88
1:C:165:THR:O	4:C:2154:HOH:O	1.91	0.88
1:C:163:VAL:CG1	1:C:163:VAL:O	2.21	0.86
1:C:88:LYS:HE3	4:C:2093:HOH:O	1.75	0.86
4:C:2140:HOH:O	2:D:316:THR:HB	1.76	0.85
1:C:159:TYR:CD1	2:D:270:ILE:HD13	2.10	0.85
3:A:1297[A]:CK8:C6A	3:A:1297[A]:CK8:H5	2.06	0.85
1:C:173:ILE:HD11	1:C:184:VAL:CG1	2.06	0.85
1:C:64:VAL:CG2	1:C:144:ALA:HB2	2.07	0.85
1:C:7:VAL:C	4:C:2009:HOH:O	2.15	0.85
2:B:345:ASP:HB2	2:B:346:PRO:HD3	1.58	0.85
1:A:91:MET:HE3	1:A:195:GLU:HG2	1.57	0.85
1:A:60:HIS:HD2	1:A:62:ASN:H	1.23	0.85
1:C:96:LEU:HD12	1:C:97:THR:N	1.92	0.84
1:C:85:GLN:NE2	1:C:89:LYS:HB3	1.91	0.84
1:C:38:ASP:O	1:C:40:GLU:N	2.11	0.84
2:B:207:THR:HG21	4:B:2011:HOH:O	1.76	0.84
1:C:160:THR:O	4:C:2147:HOH:O	1.96	0.84
2:D:198:GLY:O	2:D:201:LYS:NZ	2.09	0.84
2:B:408:SER:OG	4:B:2173:HOH:O	1.96	0.83
2:D:196:LYS:HG3	2:D:198:GLY:H	1.44	0.83
1:A:268:HIS:ND1	4:A:2214:HOH:O	2.03	0.83
1:C:6:LYS:CB	1:C:6:LYS:NZ	2.33	0.82
1:C:96:LEU:O	1:C:199:ARG:NH1	2.12	0.82
1:A:33:LYS:NZ	4:A:2040:HOH:O	2.11	0.82
2:B:425:ASN:ND2	4:B:2184:HOH:O	1.85	0.82
1:C:163:VAL:HG13	1:C:163:VAL:O	1.80	0.81
1:C:64:VAL:CG2	4:C:2070:HOH:O	2.09	0.81
2:D:378:ARG:CG	2:D:378:ARG:HH11	1.92	0.81
1:A:89:LYS:NZ	3:A:1297[B]:CK8:H8B1	1.95	0.80
2:B:293:ARG:CZ	4:B:2102:HOH:O	2.29	0.80
1:A:294:PRO:HB2	1:A:296:LEU:CD1	2.11	0.80
1:A:54:LEU:HD23	4:A:2065:HOH:O	1.81	0.80
2:B:327:CYS:CB	4:B:2018:HOH:O	2.20	0.80
1:C:60:HIS:HD2	1:C:62:ASN:H	1.29	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ALA:O	1:C:96:LEU:HG	1.81	0.80
2:B:282:THR:O	2:B:285:THR:HG23	1.82	0.80
1:A:40:GLU:HG3	1:A:40:GLU:O	1.80	0.79
1:C:15:TYR:CD2	1:C:33:LYS:CE	2.57	0.79
2:B:293:ARG:NE	4:B:2102:HOH:O	2.15	0.79
1:C:34:LYS:NZ	1:C:77:TYR:OH	2.13	0.79
1:A:256:ASP:N	4:A:2200:HOH:O	2.00	0.79
1:A:96:LEU:HD23	1:A:97:THR:HG23	1.65	0.79
3:A:1297[A]:CK8:H6A1	4:A:2060:HOH:O	1.81	0.79
1:A:81:GLU:OE2	4:A:2090:HOH:O	1.98	0.79
1:C:161:HIS:CE1	4:C:2148:HOH:O	2.35	0.79
1:A:36:ARG:HH11	1:A:36:ARG:HG2	1.48	0.78
2:B:207:THR:OG1	4:B:2056:HOH:O	2.00	0.78
1:A:91:MET:HE1	1:A:195:GLU:OE2	1.84	0.77
2:B:203:GLN:HB3	2:B:206:ILE:CD1	2.14	0.77
1:A:242:LYS:CB	4:A:2192:HOH:O	2.31	0.77
1:A:38:ASP:CG	1:A:41:THR:HB	2.03	0.77
1:C:173:ILE:HD11	1:C:184:VAL:HG11	1.65	0.77
1:A:57:GLU:CD	4:A:2066:HOH:O	2.22	0.77
1:C:40:GLU:O	2:D:288:LYS:NZ	2.18	0.76
2:D:404:HIS:HD2	2:D:406:GLN:N	1.82	0.76
1:C:159:TYR:CE1	2:D:270:ILE:HD13	2.21	0.76
1:A:81:GLU:CD	4:A:2090:HOH:O	2.22	0.76
2:B:177:ASP:CG	4:B:2034:HOH:O	2.22	0.76
2:B:196:LYS:N	2:B:196:LYS:HD2	2.00	0.75
1:C:39:THR:O	1:C:39:THR:CG2	2.35	0.75
1:A:12:GLU:OE1	1:A:36:ARG:NH2	2.20	0.75
1:C:37:LEU:O	1:C:39:THR:N	2.19	0.75
2:D:289:LYS:HD3	2:D:293:ARG:HD2	1.68	0.75
3:C:1298:CK8:C8B	4:C:2258:HOH:O	2.34	0.75
1:A:91:MET:CE	1:A:195:GLU:HG2	2.17	0.75
2:B:200:MET:O	4:B:2050:HOH:O	2.05	0.75
1:C:15:TYR:HE1	1:C:47:THR:C	1.90	0.75
1:A:85:GLN:OE1	1:A:296:LEU:CG	2.30	0.75
2:B:272:PRO:O	4:B:2092:HOH:O	2.05	0.75
2:D:378:ARG:HG2	2:D:378:ARG:NH1	1.91	0.75
4:A:2066:HOH:O	2:B:189:MET:HE2	1.86	0.74
2:B:282:THR:O	2:B:285:THR:CG2	2.36	0.74
1:A:257:GLU:H	1:A:257:GLU:CD	1.88	0.74
2:B:425:ASN:CB	4:B:2185:HOH:O	1.93	0.74
1:A:15:TYR:HB2	1:A:33:LYS:HE3	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:HIS:CD2	1:C:62:ASN:H	2.04	0.74
1:A:242:LYS:HB3	4:A:2192:HOH:O	1.87	0.74
1:C:15:TYR:CD1	1:C:47:THR:OG1	2.41	0.74
1:A:224:GLU:OE2	4:A:2184:HOH:O	2.06	0.73
1:A:227:TRP:O	1:A:230:VAL:HG22	1.88	0.73
1:C:251:VAL:HB	4:C:2089:HOH:O	1.87	0.73
1:A:260:ARG:CZ	4:A:2207:HOH:O	2.36	0.73
1:C:64:VAL:CG2	1:C:143:LEU:O	2.31	0.72
1:C:56:LYS:HD3	2:D:305:ASP:OD1	1.89	0.72
2:B:221:VAL:CG2	2:B:281:ILE:CD1	2.67	0.72
2:B:404:HIS:CD2	2:B:406:GLN:H	2.01	0.72
1:C:22:ARG:HD3	4:C:2022:HOH:O	1.88	0.72
2:B:277:GLU:O	2:B:281:ILE:HG23	1.90	0.72
2:D:431:ASN:ND2	4:D:2243:HOH:O	2.23	0.71
1:A:84:HIS:CE1	1:A:296:LEU:HD22	2.25	0.71
1:A:42:GLU:HB3	4:A:2049:HOH:O	1.86	0.71
2:B:415:ASN:OD1	2:B:416:SER:N	2.23	0.71
2:B:316:THR:CG2	4:B:2119:HOH:O	2.26	0.71
2:D:425:ASN:HB3	4:D:2238:HOH:O	1.90	0.71
2:B:345:ASP:HB2	2:B:346:PRO:CD	2.21	0.70
1:C:96:LEU:C	1:C:96:LEU:HD12	2.12	0.70
2:B:207:THR:CG2	2:B:210:MET:H	2.03	0.70
1:C:64:VAL:CB	4:C:2070:HOH:O	1.77	0.70
1:C:214:ARG:HG3	1:C:214:ARG:HH11	1.54	0.70
2:B:221:VAL:HG21	2:B:281:ILE:HD12	1.72	0.70
1:C:15:TYR:HD1	1:C:47:THR:OG1	1.72	0.70
1:C:96:LEU:HG	1:C:97:THR:H	1.55	0.70
1:C:170:ALA:HB3	1:C:173:ILE:CD1	2.22	0.69
1:A:202:LEU:HD13	1:A:203:PHE:CE2	2.28	0.69
2:B:179:HIS:NE2	2:B:379:LYS:HE2	2.07	0.69
2:B:327:CYS:SG	4:B:2018:HOH:O	2.50	0.69
2:D:207:THR:HG23	2:D:209:SER:H	1.56	0.69
1:A:36:ARG:HH11	1:A:36:ARG:CG	2.06	0.68
1:A:227:TRP:CD2	1:A:230:VAL:HG13	2.28	0.68
2:B:206:ILE:HA	2:B:210:MET:HE1	1.73	0.68
2:B:334:MET:HE2	4:B:2073:HOH:O	1.92	0.68
1:A:57:GLU:CG	4:A:2066:HOH:O	2.42	0.68
1:C:214:ARG:HG2	1:C:214:ARG:HH11	1.57	0.68
2:D:196:LYS:HB2	2:D:196:LYS:NZ	2.09	0.68
2:D:183:HIS:HB2	2:D:317:GLN:HE22	1.58	0.67
2:B:425:ASN:OD1	4:B:2185:HOH:O	2.12	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1298:CK8:H8B2	4:C:2258:HOH:O	1.94	0.67
1:A:1:MET:HE2	1:A:70:ILE:CD1	2.24	0.67
1:C:159:TYR:HE2	4:D:2027:HOH:O	1.77	0.67
1:A:49:ILE:HG23	2:B:306:LEU:HD12	1.76	0.67
2:D:296:HIS:CG	4:D:2149:HOH:O	2.41	0.67
2:B:203:GLN:HB3	2:B:206:ILE:HD13	1.77	0.67
1:C:5:GLN:HG3	1:C:5:GLN:O	1.94	0.67
2:B:428:GLU:OE1	4:B:2187:HOH:O	2.12	0.66
1:C:270:ASP:OD1	4:C:2232:HOH:O	2.12	0.66
1:C:88:LYS:CD	1:C:88:LYS:C	2.63	0.66
2:B:207:THR:HG23	2:B:210:MET:N	2.07	0.66
1:C:131:GLN:HE21	1:C:131:GLN:N	1.87	0.66
1:C:173:ILE:HD11	1:C:184:VAL:HG12	1.76	0.66
1:C:214:ARG:NH1	1:C:214:ARG:CG	2.55	0.66
1:C:22:ARG:CD	4:C:2022:HOH:O	2.42	0.66
1:A:138:GLU:CD	4:A:2131:HOH:O	2.33	0.66
1:A:1:MET:CE	1:A:70:ILE:CD1	2.73	0.66
2:B:221:VAL:HG21	2:B:281:ILE:CD1	2.25	0.66
2:B:415:ASN:OD1	2:B:417:LYS:N	2.27	0.66
2:B:425:ASN:CG	4:B:2185:HOH:O	2.24	0.66
1:C:131:GLN:H	1:C:131:GLN:NE2	1.85	0.66
1:C:91:MET:CE	1:C:195:GLU:HG2	2.25	0.66
4:C:2140:HOH:O	2:D:316:THR:CB	2.39	0.66
3:A:1297[A]:CK8:H5	3:A:1297[A]:CK8:H6A2	1.78	0.66
1:C:200:ARG:CD	4:C:2178:HOH:O	2.44	0.65
1:A:57:GLU:OE2	4:A:2066:HOH:O	2.14	0.65
2:D:400:LYS:HE3	4:D:2210:HOH:O	1.96	0.65
1:A:121:HIS:O	1:A:122:ARG:HG3	1.97	0.65
2:B:229:ASN:OD1	4:B:2072:HOH:O	2.15	0.65
1:A:242:LYS:CG	4:A:2192:HOH:O	2.45	0.64
1:C:88:LYS:CD	1:C:88:LYS:O	2.45	0.64
2:D:233:HIS:HD2	4:D:2157:HOH:O	1.81	0.64
1:C:33:LYS:HZ2	3:C:1298:CK8:H7A1	1.63	0.64
1:C:71:HIS:CE1	4:C:2074:HOH:O	2.40	0.64
3:C:1298:CK8:H6B	3:C:1298:CK8:H6A2	1.79	0.64
1:C:96:LEU:CD1	1:C:97:THR:N	2.60	0.64
2:B:194:LYS:HE2	4:B:2138:HOH:O	1.96	0.64
3:C:1298:CK8:H6B	3:C:1298:CK8:C6A	2.27	0.64
1:C:101:LEU:HD23	4:C:2108:HOH:O	1.97	0.63
1:A:42:GLU:OE2	2:B:275:VAL:CG2	2.40	0.63
1:C:251:VAL:CB	4:C:2089:HOH:O	2.44	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:260:ARG:NH1	4:C:2222:HOH:O	2.30	0.63
1:C:95:ALA:O	1:C:96:LEU:CG	2.46	0.63
1:A:163:VAL:O	1:A:163:VAL:HG13	1.97	0.63
2:B:388:LYS:N	2:B:389:PRO:CD	2.62	0.63
1:C:88:LYS:O	1:C:88:LYS:HD2	1.98	0.63
2:B:271:TYR:HD1	2:B:271:TYR:N	1.97	0.63
1:A:224:GLU:OE2	1:A:231:THR:OG1	2.11	0.62
1:A:88:LYS:HE3	4:A:2100:HOH:O	1.98	0.62
1:C:6:LYS:HG3	4:C:2008:HOH:O	1.99	0.62
2:B:233:HIS:HD2	4:B:2117:HOH:O	1.81	0.62
1:A:1:MET:CE	1:A:70:ILE:HD12	2.30	0.62
1:C:96:LEU:CA	1:C:199:ARG:HH11	2.13	0.62
1:C:38:ASP:N	1:C:38:ASP:OD1	2.33	0.62
1:A:294:PRO:HB2	1:A:296:LEU:HD13	1.80	0.62
1:A:88:LYS:CE	4:A:2100:HOH:O	2.48	0.62
1:C:245:ARG:NH2	4:C:2197:HOH:O	2.12	0.62
1:A:295:HIS:HE1	4:A:2243:HOH:O	1.82	0.61
1:C:159:TYR:CE2	4:D:2027:HOH:O	2.51	0.61
1:A:15:TYR:CD2	3:A:1297[A]:CK8:H7A2	2.34	0.61
1:A:89:LYS:HZ1	3:A:1297[B]:CK8:H8B1	1.65	0.61
1:A:73:GLU:OE2	1:C:2:GLU:HG2	2.01	0.61
1:C:248:PHE:HA	4:C:2204:HOH:O	2.00	0.61
1:A:89:LYS:HE3	3:A:1297[A]:CK8:H8B1	1.83	0.61
1:A:52:ILE:HD11	1:A:78:LEU:HD21	1.83	0.61
1:A:89:LYS:HE3	3:A:1297[B]:CK8:H9B1	1.83	0.61
1:C:166:LEU:HA	4:C:2154:HOH:O	1.99	0.61
1:C:96:LEU:C	1:C:199:ARG:NH1	2.55	0.60
4:C:2140:HOH:O	2:D:316:THR:CG2	2.49	0.60
1:A:28:GLU:OE2	4:A:2036:HOH:O	2.16	0.60
2:B:206:ILE:HA	2:B:210:MET:HE2	1.82	0.60
2:B:270:ILE:O	4:B:2091:HOH:O	2.16	0.60
2:B:203:GLN:HB3	2:B:206:ILE:HD11	1.81	0.60
1:A:1:MET:HE2	1:A:70:ILE:HD13	1.83	0.60
2:D:226:LYS:NZ	4:D:2095:HOH:O	2.34	0.60
1:C:15:TYR:HE2	1:C:33:LYS:CE	2.08	0.59
1:C:200:ARG:HD3	4:C:2178:HOH:O	2.01	0.59
3:A:1297[A]:CK8:H5	3:A:1297[A]:CK8:H6A3	1.82	0.59
1:C:95:ALA:O	1:C:96:LEU:CB	2.50	0.59
1:C:96:LEU:CA	1:C:199:ARG:NH1	2.66	0.59
1:A:57:GLU:HG2	4:A:2066:HOH:O	2.01	0.59
2:B:351:LEU:HG	4:B:2138:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:196:LYS:HZ2	2:D:196:LYS:HB2	1.68	0.59
1:C:115:LEU:HD12	1:C:189:LEU:HD22	1.84	0.59
2:D:361:HIS:CD2	2:D:391:LEU:HD21	2.38	0.59
2:B:370:GLN:CD	4:B:2145:HOH:O	2.41	0.58
1:A:138:GLU:OE1	4:A:2131:HOH:O	2.17	0.58
1:A:295:HIS:CE1	4:A:2243:HOH:O	2.54	0.58
1:A:35:ILE:HG13	1:A:76:LEU:HD13	1.84	0.58
1:C:121:HIS:O	1:C:122:ARG:HG3	2.03	0.58
1:C:158:THR:HG21	1:C:177:CYS:O	2.03	0.58
1:A:1:MET:CE	1:A:70:ILE:HD13	2.34	0.58
2:B:221:VAL:HG22	2:B:281:ILE:CD1	2.31	0.58
2:B:271:TYR:CD1	2:B:271:TYR:N	2.70	0.58
2:B:430:LEU:HB3	2:B:432:LEU:HD22	1.85	0.58
2:B:226:LYS:NZ	4:B:2068:HOH:O	2.34	0.58
1:C:165:THR:O	4:C:2151:HOH:O	2.17	0.58
1:A:164:VAL:O	1:A:165:THR:HB	2.03	0.58
2:B:395:HIS:NE2	2:B:399:LEU:HD11	2.18	0.58
1:C:101:LEU:CD2	4:C:2108:HOH:O	2.52	0.58
2:D:175:VAL:O	2:D:175:VAL:HG13	2.03	0.58
3:A:1297[A]:CK8:C6A	4:A:2060:HOH:O	2.47	0.57
2:B:203:GLN:OE1	2:B:206:ILE:HD11	2.03	0.57
1:C:138:GLU:N	1:C:138:GLU:CD	2.57	0.57
2:D:194:LYS:CE	4:D:2070:HOH:O	2.36	0.57
1:A:249:SER:HB2	4:A:2207:HOH:O	2.05	0.57
1:A:131:GLN:OE1	4:A:2125:HOH:O	2.10	0.57
1:A:159:TYR:N	4:A:2144:HOH:O	2.03	0.57
2:B:250:ARG:HG3	4:B:2085:HOH:O	2.05	0.56
1:A:84:HIS:NE2	1:A:296:LEU:CD2	2.52	0.56
1:A:159:TYR:HA	2:B:270:ILE:HD13	1.86	0.56
4:C:2074:HOH:O	2:D:296:HIS:CE1	2.58	0.56
1:A:86:ASP:OD2	3:A:1297[B]:CK8:H8B3	2.06	0.56
1:C:214:ARG:HG2	1:C:214:ARG:NH1	2.20	0.56
1:C:85:GLN:HE21	1:C:89:LYS:HB3	1.68	0.56
1:A:111:LEU:HD13	1:A:189:LEU:HD11	1.87	0.56
1:C:96:LEU:CG	1:C:97:THR:N	2.67	0.56
1:C:253:PRO:N	1:C:254:PRO:CD	2.69	0.56
2:D:207:THR:HG22	2:D:210:MET:HG3	1.88	0.56
1:A:51:GLU:O	1:A:55:LEU:HB2	2.05	0.56
2:D:312:ASN:OD1	2:D:334:MET:SD	2.64	0.56
2:B:225:TYR:HE1	2:B:281:ILE:HG21	1.70	0.55
1:C:91:MET:HE2	1:C:195:GLU:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:245:ARG:NH2	4:C:2196:HOH:O	2.32	0.55
2:D:270:ILE:HG22	2:D:271:TYR:CD2	2.40	0.55
1:A:159:TYR:HB2	4:A:2144:HOH:O	2.05	0.55
2:B:216:ASP:OD1	2:B:408:SER:HB2	2.07	0.55
1:C:15:TYR:CE1	1:C:47:THR:CB	2.89	0.55
1:A:89:LYS:CE	3:A:1297[B]:CK8:H8B1	2.35	0.55
1:A:54:LEU:HA	4:A:2065:HOH:O	2.06	0.55
2:B:430:LEU:O	2:B:431:ASN:HB2	2.06	0.55
1:C:71:HIS:NE2	2:D:304:PHE:CE2	2.75	0.55
1:C:96:LEU:CG	1:C:97:THR:H	2.20	0.55
1:A:159:TYR:CB	4:A:2144:HOH:O	2.55	0.55
1:A:177:CYS:CB	4:A:2157:HOH:O	1.73	0.55
1:A:249:SER:CB	4:A:2207:HOH:O	2.54	0.55
2:B:197:VAL:N	4:B:2049:HOH:O	2.39	0.55
2:B:205:ASP:OD2	2:B:250:ARG:HD2	2.06	0.54
2:B:206:ILE:CA	2:B:210:MET:CE	2.81	0.54
3:A:1297[A]:CK8:C5	3:A:1297[A]:CK8:C6A	2.81	0.54
1:A:260:ARG:NE	4:A:2207:HOH:O	2.40	0.54
1:C:91:MET:HE1	1:C:195:GLU:HG2	1.87	0.54
1:C:96:LEU:HA	1:C:199:ARG:NH1	2.23	0.54
2:D:347:TYR:OH	2:D:394:LEU:HA	2.07	0.54
2:B:395:HIS:CE1	2:B:399:LEU:HD11	2.43	0.54
1:C:85:GLN:HA	3:C:1298:CK8:C3B	2.38	0.54
1:C:163:VAL:HG12	1:C:163:VAL:O	2.06	0.54
1:A:96:LEU:O	1:A:199:ARG:NH1	2.41	0.54
2:D:284:ASP:CB	4:D:2139:HOH:O	2.03	0.54
3:A:1297[B]:CK8:H6B	3:A:1297[B]:CK8:C6A	2.38	0.53
1:A:71:HIS:CE1	2:B:296:HIS:CE1	2.95	0.53
1:A:164:VAL:HG13	4:A:2146:HOH:O	2.07	0.53
1:A:38:ASP:OD1	1:A:41:THR:HB	2.08	0.53
1:A:86:ASP:OD2	1:A:89:LYS:HE2	2.08	0.53
1:C:249:SER:N	4:C:2205:HOH:O	2.41	0.53
1:C:88:LYS:C	1:C:88:LYS:HD3	2.27	0.53
2:B:203:GLN:CD	2:B:206:ILE:HD11	2.29	0.53
2:B:345:ASP:OD1	2:B:345:ASP:N	2.39	0.53
1:A:15:TYR:HD1	1:A:15:TYR:H	1.52	0.53
1:C:101:LEU:HB2	1:C:102:PRO:HD3	1.91	0.53
1:C:6:LYS:CG	4:C:2008:HOH:O	2.56	0.53
1:A:162:GLU:HG3	1:A:164:VAL:HB	1.91	0.53
1:A:96:LEU:C	1:A:199:ARG:NH1	2.61	0.53
1:C:15:TYR:OH	1:C:51:GLU:OE1	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:271:TYR:HB3	4:B:2091:HOH:O	2.09	0.52
1:C:126:ARG:NH2	4:C:2131:HOH:O	2.30	0.52
2:D:198:GLY:CA	2:D:201:LYS:HZ1	2.22	0.52
1:A:227:TRP:CE3	1:A:230:VAL:HG13	2.44	0.52
1:C:1:MET:HE2	1:C:70:ILE:HD13	1.91	0.52
1:C:15:TYR:HD2	1:C:33:LYS:HE2	1.53	0.52
1:A:65:LYS:NZ	4:A:2069:HOH:O	2.42	0.52
1:C:125:HIS:O	1:C:126:ARG:HB2	2.08	0.52
2:D:388:LYS:N	2:D:389:PRO:HD2	2.25	0.52
3:A:1297[B]:CK8:N3	3:A:1297[B]:CK8:C6B	2.51	0.52
1:A:91:MET:HE1	1:A:195:GLU:CD	2.29	0.52
1:C:60:HIS:HE1	4:C:2063:HOH:O	1.92	0.52
1:A:60:HIS:HD2	1:A:62:ASN:N	2.01	0.52
2:B:206:ILE:HB	2:B:210:MET:HE3	1.92	0.52
2:B:370:GLN:OE1	4:B:2145:HOH:O	2.18	0.52
1:C:15:TYR:CE1	1:C:47:THR:HB	2.45	0.51
1:C:166:LEU:HD22	1:C:169:ARG:NH1	2.25	0.51
1:C:166:LEU:CD2	1:C:169:ARG:NH1	2.72	0.51
2:B:179:HIS:NE2	2:B:379:LYS:CE	2.72	0.51
2:D:400:LYS:HE2	4:D:2212:HOH:O	2.09	0.51
1:A:38:ASP:OD2	1:A:41:THR:CG2	2.58	0.51
1:A:84:HIS:CD2	1:A:296:LEU:HD23	2.42	0.51
2:B:351:LEU:N	4:B:2138:HOH:O	2.43	0.51
1:C:159:TYR:CD1	2:D:270:ILE:CD1	2.89	0.51
2:D:431:ASN:CG	4:D:2243:HOH:O	2.48	0.51
2:B:221:VAL:CG1	2:B:281:ILE:HD13	2.41	0.51
2:D:334:MET:HE3	4:D:2169:HOH:O	2.11	0.50
2:B:206:ILE:H	2:B:206:ILE:HD13	1.75	0.50
1:C:138:GLU:H	1:C:138:GLU:CD	2.15	0.50
1:C:260:ARG:HD3	4:C:2222:HOH:O	2.12	0.50
1:C:51:GLU:O	1:C:55:LEU:HB2	2.11	0.50
1:A:101:LEU:N	1:A:102:PRO:CD	2.74	0.50
2:B:388:LYS:N	2:B:389:PRO:HD2	2.27	0.50
2:D:207:THR:HG22	2:D:210:MET:H	1.75	0.50
1:A:124:LEU:HG	1:A:152:PHE:CD1	2.46	0.50
1:A:91:MET:CE	4:A:2166:HOH:O	2.58	0.50
1:C:88:LYS:CG	4:C:2093:HOH:O	2.60	0.50
1:C:88:LYS:O	1:C:88:LYS:HD3	2.10	0.50
2:D:220:GLU:HG2	2:D:408:SER:OG	2.12	0.50
1:A:212:LEU:HD22	1:A:216:PHE:CE1	2.47	0.50
1:A:36:ARG:CZ	4:A:2045:HOH:O	2.49	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:287:GLN:NE2	4:C:2247:HOH:O	2.43	0.50
2:B:319:PHE:CE2	2:B:330:GLU:HG2	2.47	0.49
1:A:121:HIS:C	1:A:122:ARG:HG3	2.33	0.49
2:B:225:TYR:HE1	2:B:281:ILE:CG2	2.25	0.49
1:C:101:LEU:N	1:C:102:PRO:CD	2.75	0.49
1:C:24:LYS:NZ	4:C:2030:HOH:O	2.44	0.49
2:D:220:GLU:CG	2:D:408:SER:OG	2.60	0.49
1:A:84:HIS:HE2	1:A:296:LEU:HB3	1.77	0.49
1:A:88:LYS:HD3	1:A:131:GLN:OE1	2.13	0.48
1:A:91:MET:HE2	4:A:2166:HOH:O	2.13	0.48
1:A:164:VAL:HG13	1:A:164:VAL:O	2.12	0.48
1:C:294:PRO:HG2	1:C:296:LEU:HD22	1.96	0.48
2:D:343:ASP:O	2:D:346:PRO:HD2	2.13	0.48
1:A:145:ASP:OD2	4:A:2133:HOH:O	2.20	0.48
1:A:165:THR:HG22	4:A:2146:HOH:O	2.13	0.48
1:A:42:GLU:OE2	1:A:42:GLU:HA	2.09	0.48
1:A:294:PRO:CB	1:A:296:LEU:HD13	2.44	0.48
1:C:283:HIS:CG	1:C:284:PRO:HD2	2.49	0.48
1:C:71:HIS:HE1	2:D:296:HIS:CD2	2.28	0.48
1:A:161:HIS:CD2	4:A:2145:HOH:O	2.66	0.48
1:A:294:PRO:CG	1:A:296:LEU:HD13	2.44	0.48
1:A:59:ASN:ND2	4:A:2069:HOH:O	2.36	0.48
1:C:126:ARG:NH2	1:C:148:LEU:HD22	2.29	0.48
1:A:159:TYR:O	1:A:160:THR:C	2.51	0.48
2:B:225:TYR:CE1	2:B:281:ILE:HG21	2.47	0.48
1:A:27:GLY:HA3	2:D:249:LEU:HD22	1.95	0.47
2:B:206:ILE:N	2:B:206:ILE:HD13	2.28	0.47
1:C:170:ALA:HB3	1:C:173:ILE:HD12	1.94	0.47
2:B:203:GLN:CB	2:B:206:ILE:HD11	2.44	0.47
2:B:250:ARG:HG2	2:B:251:GLY:N	2.28	0.47
1:A:202:LEU:HD13	1:A:203:PHE:CZ	2.48	0.47
2:B:175:VAL:HG13	4:B:2035:HOH:O	2.13	0.47
2:B:345:ASP:CB	2:B:346:PRO:HD3	2.37	0.47
3:C:1298:CK8:N3	3:C:1298:CK8:C6A	2.77	0.47
1:C:15:TYR:OH	1:C:51:GLU:HB3	2.15	0.47
2:B:425:ASN:HA	2:B:426:PRO:HD3	1.76	0.47
2:B:250:ARG:HB2	4:B:2083:HOH:O	2.14	0.47
2:B:271:TYR:HD1	2:B:271:TYR:H	1.63	0.47
2:B:201:LYS:CE	2:B:202:LYS:HG3	2.44	0.47
2:B:343:ASP:OD1	2:B:404:HIS:HE1	1.98	0.47
1:C:137:THR:HG22	1:C:296:LEU:HD23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:HIS:HB3	4:A:2092:HOH:O	2.13	0.47
1:C:170:ALA:HB3	1:C:173:ILE:HD13	1.96	0.47
1:A:2:GLU:HG2	1:C:73:GLU:HG3	1.96	0.47
2:D:323:GLN:HA	2:D:324:PRO:HA	1.72	0.47
1:A:83:LEU:HD23	1:A:136:ASN:HB3	1.96	0.47
1:A:9:LYS:HG3	1:A:9:LYS:O	2.11	0.47
1:A:15:TYR:CE2	3:A:1297[A]:CK8:H7A2	2.50	0.46
2:B:179:HIS:NE2	2:B:379:LYS:NZ	2.62	0.46
2:B:199:TYR:C	2:B:199:TYR:CD1	2.87	0.46
2:B:319:PHE:CD2	2:B:330:GLU:HG2	2.49	0.46
1:A:20:LYS:HE2	1:A:82:PHE:CZ	2.50	0.46
1:A:83:LEU:O	3:A:1297[A]:CK8:H2B	2.16	0.46
1:A:89:LYS:CE	3:A:1297[A]:CK8:H8B1	2.43	0.46
1:A:128:LEU:CD2	1:A:143:LEU:HD22	2.46	0.46
2:B:203:GLN:CG	2:B:206:ILE:HD11	2.46	0.46
1:C:51:GLU:HG3	1:C:55:LEU:HD22	1.98	0.46
1:A:15:TYR:HB2	1:A:33:LYS:CE	2.43	0.46
1:C:81:GLU:O	3:C:1298:CK8:H6	2.15	0.46
1:C:165:THR:HG21	4:C:2152:HOH:O	2.14	0.46
1:C:9:LYS:O	1:C:9:LYS:HE3	2.16	0.46
2:B:175:VAL:HG23	2:B:176:PRO:HD2	1.98	0.46
1:C:157:ARG:NE	4:C:2142:HOH:O	2.47	0.46
2:D:329:VAL:HG11	2:D:364:LEU:HD13	1.96	0.46
2:B:200:MET:HG2	2:B:208:ASN:ND2	2.31	0.46
1:C:245:ARG:NE	4:C:2197:HOH:O	2.45	0.46
2:D:388:LYS:N	2:D:389:PRO:CD	2.79	0.46
2:D:401:ALA:N	2:D:402:PRO:CD	2.79	0.46
1:A:39:THR:HA	4:A:2051:HOH:O	2.15	0.45
1:A:85:GLN:HA	3:A:1297[A]:CK8:C3B	2.47	0.45
1:C:15:TYR:HE1	1:C:47:THR:CB	2.28	0.45
2:B:385:GLU:HG2	4:B:2029:HOH:O	2.16	0.45
3:C:1298:CK8:H6B	3:C:1298:CK8:H6A3	1.98	0.45
2:B:196:LYS:N	2:B:196:LYS:CD	2.77	0.45
2:D:223:GLU:OE1	2:D:412:LYS:HE3	2.16	0.45
1:C:246:GLN:O	1:C:247:ASP:C	2.55	0.45
1:C:41:THR:HB	1:C:42:GLU:H	1.35	0.45
1:C:71:HIS:ND1	2:D:296:HIS:CE1	2.80	0.45
1:A:154:VAL:HG13	1:A:155:PRO:HD2	1.99	0.45
1:A:36:ARG:NH2	4:A:2045:HOH:O	2.50	0.45
2:B:203:GLN:OE1	2:B:206:ILE:CD1	2.65	0.45
1:C:33:LYS:NZ	3:C:1298:CK8:H7A1	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:198:GLY:CA	2:D:201:LYS:NZ	2.80	0.45
1:C:85:GLN:HA	3:C:1298:CK8:C2B	2.47	0.45
1:C:198:THR:O	1:C:199:ARG:HB2	2.17	0.45
1:A:1:MET:N	4:A:2003:HOH:O	2.47	0.44
1:A:257:GLU:CD	1:A:257:GLU:N	2.64	0.44
1:A:105:LYS:HE2	1:A:285:PHE:CZ	2.53	0.44
2:B:332:LEU:HD23	2:B:363:ALA:HA	1.98	0.44
1:C:17:VAL:CG1	1:C:18:VAL:N	2.79	0.44
2:D:312:ASN:OD1	2:D:334:MET:CE	2.65	0.44
2:B:421:VAL:O	2:B:424:LEU:HB2	2.17	0.44
1:C:242:LYS:HD2	1:C:242:LYS:HA	1.66	0.44
1:C:296:LEU:HA	1:C:296:LEU:HD12	1.76	0.44
1:C:223:ASP:C	1:C:223:ASP:OD1	2.56	0.44
2:D:176:PRO:HB3	2:D:178:TYR:CE1	2.52	0.44
1:A:15:TYR:CE2	3:A:1297[A]:CK8:C7A	3.01	0.44
1:A:177:CYS:SG	4:A:2157:HOH:O	2.53	0.44
1:C:71:HIS:CD2	2:D:304:PHE:HE2	2.35	0.44
1:A:294:PRO:HG2	1:A:296:LEU:HD13	1.99	0.44
1:C:252:VAL:N	1:C:253:PRO:HD3	2.32	0.44
2:D:183:HIS:CB	2:D:317:GLN:HE22	2.29	0.44
1:A:145:ASP:HB2	3:A:1297[B]:CK8:H7A2	1.99	0.44
1:C:251:VAL:CG1	4:C:2089:HOH:O	2.66	0.44
1:A:128:LEU:HD21	1:A:143:LEU:HD22	2.00	0.43
2:B:221:VAL:HG13	2:B:281:ILE:HD13	1.99	0.43
2:B:430:LEU:O	2:B:431:ASN:CB	2.66	0.43
1:A:84:HIS:CD2	1:A:84:HIS:C	2.90	0.43
1:C:37:LEU:O	1:C:38:ASP:C	2.56	0.43
2:D:207:THR:CG2	2:D:210:MET:H	2.31	0.43
1:C:1:MET:CE	1:C:70:ILE:HD13	2.48	0.43
2:D:219:VAL:HG22	2:D:232:LEU:HD11	2.00	0.43
1:C:199:ARG:NH2	4:C:2100:HOH:O	2.51	0.43
1:C:15:TYR:CE1	1:C:47:THR:C	2.81	0.43
2:D:207:THR:HG23	2:D:209:SER:N	2.29	0.43
1:A:51:GLU:HG3	1:A:55:LEU:HD22	2.00	0.43
2:B:316:THR:CB	4:B:2119:HOH:O	2.62	0.43
1:C:15:TYR:HE1	1:C:48:ALA:N	2.17	0.43
1:A:131:GLN:HG2	1:A:131:GLN:H	1.26	0.43
1:C:71:HIS:NE2	2:D:304:PHE:HE2	2.16	0.43
2:B:430:LEU:HA	2:B:430:LEU:HD23	1.75	0.42
1:A:84:HIS:O	1:A:84:HIS:CD2	2.72	0.42
1:A:85:GLN:OE1	1:A:296:LEU:CD2	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:THR:O	2:B:296:HIS:HE1	2.02	0.42
1:C:22:ARG:HD2	4:C:2022:HOH:O	2.14	0.42
1:C:218:THR:O	1:C:246:GLN:HG2	2.18	0.42
4:C:2074:HOH:O	2:D:296:HIS:CD2	2.72	0.42
2:B:175:VAL:CG2	2:B:176:PRO:HD2	2.50	0.42
2:B:293:ARG:CD	4:B:2102:HOH:O	2.64	0.42
2:B:344:ALA:HB1	2:B:348:LEU:HD22	2.01	0.42
2:B:430:LEU:CB	2:B:432:LEU:HD22	2.48	0.42
1:A:89:LYS:CE	3:A:1297[A]:CK8:C8B	2.97	0.42
1:C:189:LEU:HD12	1:C:189:LEU:HA	1.93	0.42
1:C:202:LEU:HD23	1:C:203:PHE:CE2	2.55	0.42
1:C:15:TYR:CE1	1:C:48:ALA:N	2.88	0.42
2:B:376:LEU:HA	2:B:376:LEU:HD23	1.86	0.42
1:C:78:LEU:N	1:C:78:LEU:HD23	2.34	0.42
2:D:198:GLY:HA2	2:D:201:LYS:HZ1	1.83	0.42
2:D:323:GLN:O	2:D:323:GLN:HG2	2.19	0.42
1:A:111:LEU:CD1	1:A:189:LEU:HD11	2.49	0.42
1:C:251:VAL:HG11	4:C:2089:HOH:O	2.20	0.42
2:D:388:LYS:O	2:D:392:MET:HG2	2.19	0.42
2:B:401:ALA:HB3	2:B:402:PRO:HD3	2.02	0.42
1:C:25:LEU:HA	1:C:25:LEU:HD23	1.79	0.42
2:B:201:LYS:HE3	2:B:202:LYS:HG3	2.01	0.41
2:B:225:TYR:CE1	2:B:281:ILE:CG2	3.03	0.41
1:C:171:PRO:HA	1:C:174:LEU:HD12	2.02	0.41
1:A:256:ASP:CG	4:A:2200:HOH:O	2.48	0.41
1:C:15:TYR:HE2	1:C:33:LYS:CD	2.33	0.41
2:D:328:LYS:HG2	2:D:328:LYS:HZ3	1.62	0.41
2:B:175:VAL:C	2:B:177:ASP:H	2.18	0.41
2:B:202:LYS:O	2:B:204:PRO:HD3	2.20	0.41
2:B:347:TYR:OH	2:B:394:LEU:HA	2.21	0.41
1:A:2:GLU:HG2	1:C:73:GLU:CG	2.50	0.41
2:B:214:LEU:HD22	2:B:253:LEU:HG	2.02	0.41
2:B:205:ASP:O	2:B:210:MET:HE1	2.21	0.41
1:C:126:ARG:NH1	1:C:180:TYR:OH	2.53	0.41
1:C:90:PHE:HB2	1:C:135:ILE:HD11	2.02	0.41
1:C:96:LEU:HA	1:C:199:ARG:HH11	1.80	0.41
1:C:60:HIS:CG	1:C:61:PRO:HD2	2.56	0.41
1:A:62:ASN:ND2	1:A:110:GLN:HB3	2.36	0.41
1:C:15:TYR:CE2	1:C:33:LYS:NZ	2.88	0.41
1:A:36:ARG:NH1	1:A:36:ARG:CG	2.71	0.41
2:B:191:VAL:O	2:B:191:VAL:HG12	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:206:ILE:CB	2:B:210:MET:HE3	2.51	0.41
1:C:214:ARG:HG2	4:C:2187:HOH:O	2.19	0.41
1:C:96:LEU:HG	1:C:97:THR:N	2.26	0.41
1:A:89:LYS:NZ	4:A:2097:HOH:O	2.54	0.41
2:B:292:LEU:HD12	2:B:292:LEU:HA	1.88	0.41
2:D:183:HIS:HB2	2:D:317:GLN:NE2	2.32	0.41
2:D:191:VAL:HG12	2:D:191:VAL:O	2.21	0.41
2:D:202:LYS:HB3	2:D:202:LYS:HZ2	1.86	0.41
2:B:338:GLU:HG3	4:B:2129:HOH:O	2.21	0.40
1:C:15:TYR:HB2	1:C:16:GLY:H	1.35	0.40
2:B:201:LYS:HE2	2:B:202:LYS:HG3	2.03	0.40
1:A:91:MET:HE3	4:A:2167:HOH:O	2.21	0.40
2:B:348:LEU:HD12	2:B:348:LEU:HA	1.98	0.40
2:B:431:ASN:HD22	2:B:431:ASN:HA	1.64	0.40
1:C:227:TRP:CE3	1:C:269:TYR:HB3	2.56	0.40
1:A:84:HIS:CE1	1:A:296:LEU:CB	3.05	0.40
1:C:119:HIS:CE1	1:C:182:THR:HB	2.56	0.40
1:C:15:TYR:HD1	1:C:47:THR:HG1	1.40	0.40
1:C:6:LYS:HE2	4:C:2083:HOH:O	2.21	0.40

All (2) 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)
4:C:2211:HOH:O	4:D:2115:HOH:O[1_655]	2.13	0.07
2:B:431:ASN:ND2	1:C:210:ASP:OD1[2_664]	2.19	0.01

## 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	294/298 (99%)	281 (96%)	9 (3%)	4 (1%)	11 6

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	295/298 (99%)	280 (95%)	10 (3%)	5 (2%)	9	4
2	B	256/259 (99%)	245 (96%)	4 (2%)	7 (3%)	5	1
2	D	256/259 (99%)	250 (98%)	4 (2%)	2 (1%)	19	15
All	All	1101/1114 (99%)	1056 (96%)	27 (2%)	18 (2%)	9	5

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	LEU
2	B	176	PRO
1	C	38	ASP
1	C	39	THR
1	C	41	THR
2	D	346	PRO
1	A	160	THR
1	A	165	THR
2	B	346	PRO
2	B	421	VAL
2	B	424	LEU
1	C	96	LEU
1	A	145	ASP
2	B	417	LYS
2	B	304	PHE
2	B	345	ASP
1	C	164	VAL
2	D	345	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	261/263 (99%)	228 (87%)	33 (13%)	4	2
1	C	261/263 (99%)	224 (86%)	37 (14%)	3	1
2	B	232/233 (100%)	211 (91%)	21 (9%)	9	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	232/233 (100%)	219 (94%)	13 (6%)	21	18
All	All	986/992 (99%)	882 (90%)	104 (10%)	7	4

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	14	THR
1	A	15	TYR
1	A	36	ARG
1	A	38	ASP
1	A	39	THR
1	A	40	GLU
1	A	42	GLU
1	A	55	LEU
1	A	56	LYS
1	A	74	ASN
1	A	76	LEU
1	A	84	HIS
1	A	89	LYS
1	A	96	LEU
1	A	101	LEU
1	A	122	ARG
1	A	131	GLN
1	A	160	THR
1	A	163	VAL
1	A	164	VAL
1	A	165	THR
1	A	200	ARG
1	A	202	LEU
1	A	212	LEU
1	A	230	VAL
1	A	239	SER
1	A	242	LYS
1	A	252	VAL
1	A	255	LEU
1	A	264	SER
1	A	287	GLN
1	A	293	VAL
2	B	194	LYS
2	B	196	LYS
2	B	199	TYR

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Mol	Chain	Res	Type
2	B	201	LYS
2	B	206	ILE
2	B	249	LEU
2	B	250	ARG
2	B	271	TYR
2	B	274	GLU
2	B	281	ILE
2	B	285	THR
2	B	292	LEU
2	B	322	GLN
2	B	323	GLN
2	B	374	GLU
2	B	375	SER
2	B	423	LEU
2	B	424	LEU
2	B	425	ASN
2	B	431	ASN
2	B	432	LEU
1	C	2	GLU
1	C	5	GLN
1	C	6	LYS
1	C	9	LYS
1	C	22	ARG
1	C	34	LYS
1	C	38	ASP
1	C	39	THR
1	C	41	THR
1	C	55	LEU
1	C	71	HIS
1	C	75	LYS
1	C	78	LEU
1	C	83	LEU
1	C	87	LEU
1	C	88	LYS
1	C	89	LYS
1	C	96	LEU
1	C	97	THR
1	C	122	ARG
1	C	124	LEU
1	C	126	ARG
1	C	131	GLN
1	C	138	GLU

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Mol	Chain	Res	Type
1	C	154	VAL
1	C	158	THR
1	C	161	HIS
1	C	163	VAL
1	C	164	VAL
1	C	189	LEU
1	C	214	ARG
1	C	217	ARG
1	C	247	ASP
1	C	250	LYS
1	C	252	VAL
1	C	264	SER
1	C	296	LEU
2	D	196	LYS
2	D	199	TYR
2	D	207	THR
2	D	224	GLU
2	D	232	LEU
2	D	289	LYS
2	D	292	LEU
2	D	328	LYS
2	D	345	ASP
2	D	364	LEU
2	D	375	SER
2	D	378	ARG
2	D	424	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	84	HIS
2	B	208	ASN
2	B	233	HIS
2	B	323	GLN
2	B	404	HIS
2	B	419	HIS
2	B	425	ASN
2	B	431	ASN
1	C	59	ASN
1	C	60	HIS
1	C	84	HIS

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Mol	Chain	Res	Type
1	C	85	GLN
1	C	113	GLN
1	C	131	GLN
1	C	287	GLN
2	D	179	HIS
2	D	233	HIS
2	D	317	GLN
2	D	322	GLN
2	D	370	GLN
2	D	404	HIS
2	D	419	HIS

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

3 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
3	CK8	C	1298	-	21,25,25	1.36	3 (14%)	26,35,35	2.55	9 (34%)
3	CK8	A	1297[B]	-	21,25,25	1.22	2 (9%)	26,35,35	2.14	8 (30%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	CK8	A	1297[A]	-	21,25,25	1.18	1 (4%)	26,35,35	2.37	11 (42%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CK8	C	1298	-	-	0/8/12/12	0/3/3/3
3	CK8	A	1297[B]	-	-	0/8/12/12	0/3/3/3
3	CK8	A	1297[A]	-	-	0/8/12/12	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1298	CK8	C4-C5A	-4.54	1.38	1.49
3	A	1297[B]	CK8	C4-C5A	-4.44	1.39	1.49
3	A	1297[A]	CK8	C4-C5A	-4.13	1.39	1.49
3	C	1298	CK8	C1B-N7	-2.49	1.35	1.40
3	C	1298	CK8	C6A-C1A	-2.45	1.46	1.50
3	A	1297[B]	CK8	C1B-N7	-2.14	1.36	1.40

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1297[A]	CK8	C4-N3-C2	7.14	122.57	116.69
3	C	1298	CK8	N1-C2-N3	-6.22	120.65	126.55
3	C	1298	CK8	C6-N1-C2	5.90	120.68	115.45
3	C	1298	CK8	C4-N3-C2	5.65	121.34	116.69
3	A	1297[B]	CK8	C4-N3-C2	5.15	120.94	116.69
3	A	1297[B]	CK8	N1-C2-N3	-4.68	122.11	126.55
3	A	1297[A]	CK8	N1-C2-N3	-4.51	122.28	126.55
3	A	1297[B]	CK8	C6-N1-C2	4.48	119.42	115.45
3	A	1297[B]	CK8	C5-C6-N1	-3.69	119.38	123.96
3	C	1298	CK8	C5-C6-N1	-3.52	119.58	123.96
3	A	1297[B]	CK8	C5A-C4-N3	3.52	120.70	116.59
3	A	1297[A]	CK8	C5-C4-N3	-3.47	117.47	121.97
3	A	1297[A]	CK8	C5-C4-C5A	3.05	124.50	120.66
3	C	1298	CK8	N7-C2-N3	3.02	127.17	116.92
3	C	1298	CK8	C7A-C3A-S4A	2.71	123.75	120.12
3	C	1298	CK8	C7A-C3A-N2A	-2.68	116.75	123.80
3	A	1297[A]	CK8	C7A-C3A-S4A	2.53	123.51	120.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1297[B]	CK8	C5-C4-N3	-2.49	118.75	121.97
3	A	1297[A]	CK8	C5-C6-N1	-2.43	120.94	123.96
3	A	1297[A]	CK8	C3B-C4B-N7B	-2.43	118.35	121.63
3	C	1298	CK8	C5B-C4B-N7B	-2.35	118.45	121.63
3	A	1297[A]	CK8	C7A-C3A-N2A	-2.34	117.65	123.80
3	A	1297[B]	CK8	C6-C5-C4	2.30	119.39	117.22
3	A	1297[A]	CK8	C6-C5-C4	2.28	119.37	117.22
3	C	1298	CK8	C5-C4-N3	-2.27	119.03	121.97
3	A	1297[B]	CK8	C5B-C4B-N7B	-2.08	118.82	121.63
3	A	1297[A]	CK8	C6-N1-C2	2.03	117.25	115.45
3	A	1297[A]	CK8	C3B-C2B-C1B	2.02	122.63	120.30

There are no chirality outliers.

There are no torsion outliers.

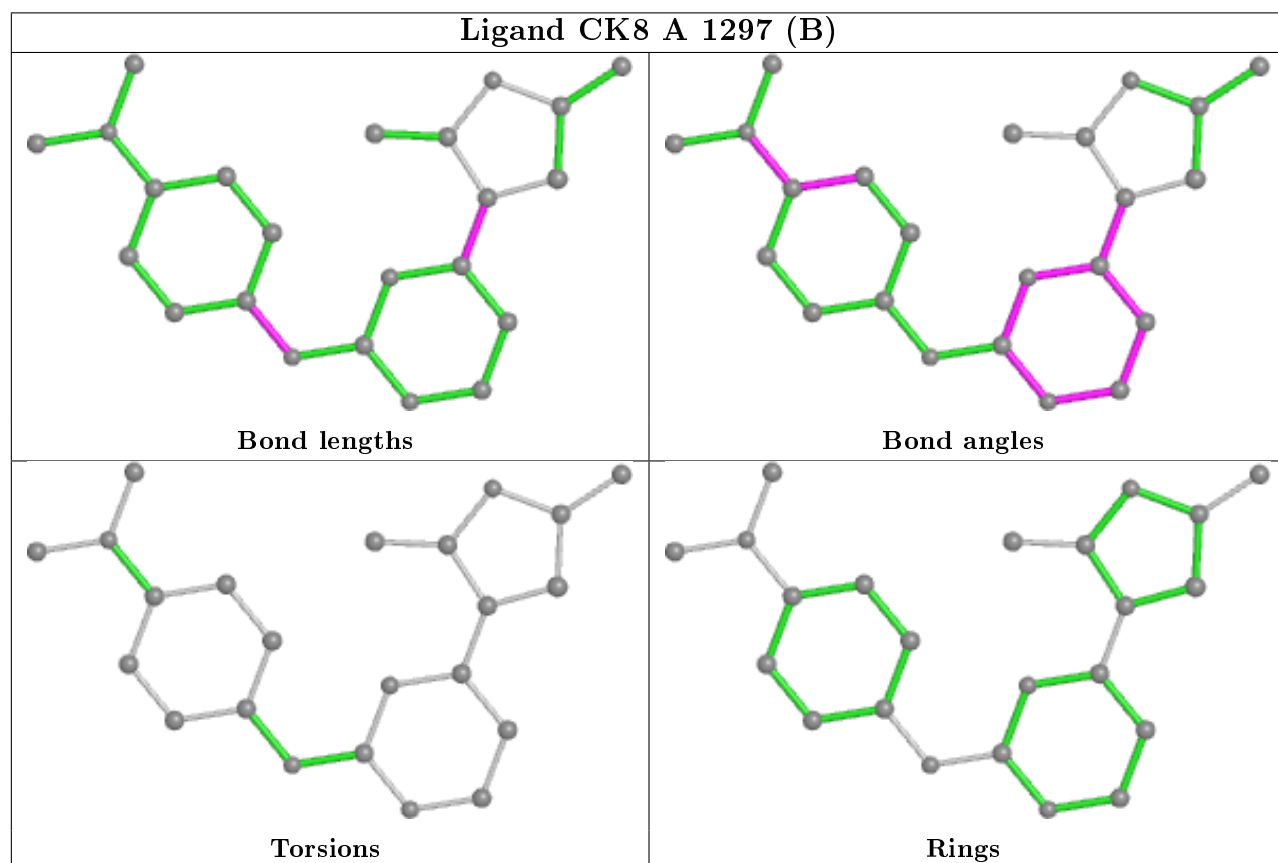
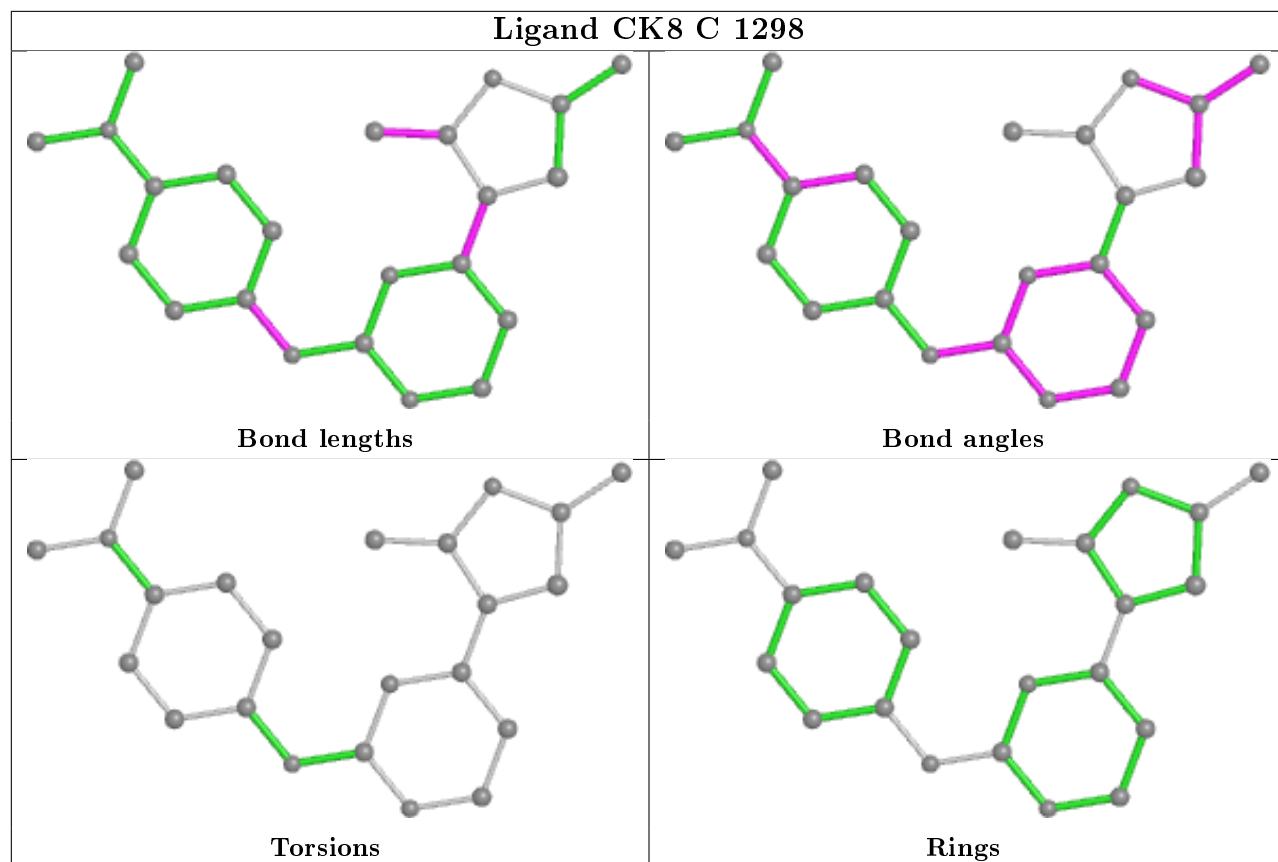
There are no ring outliers.

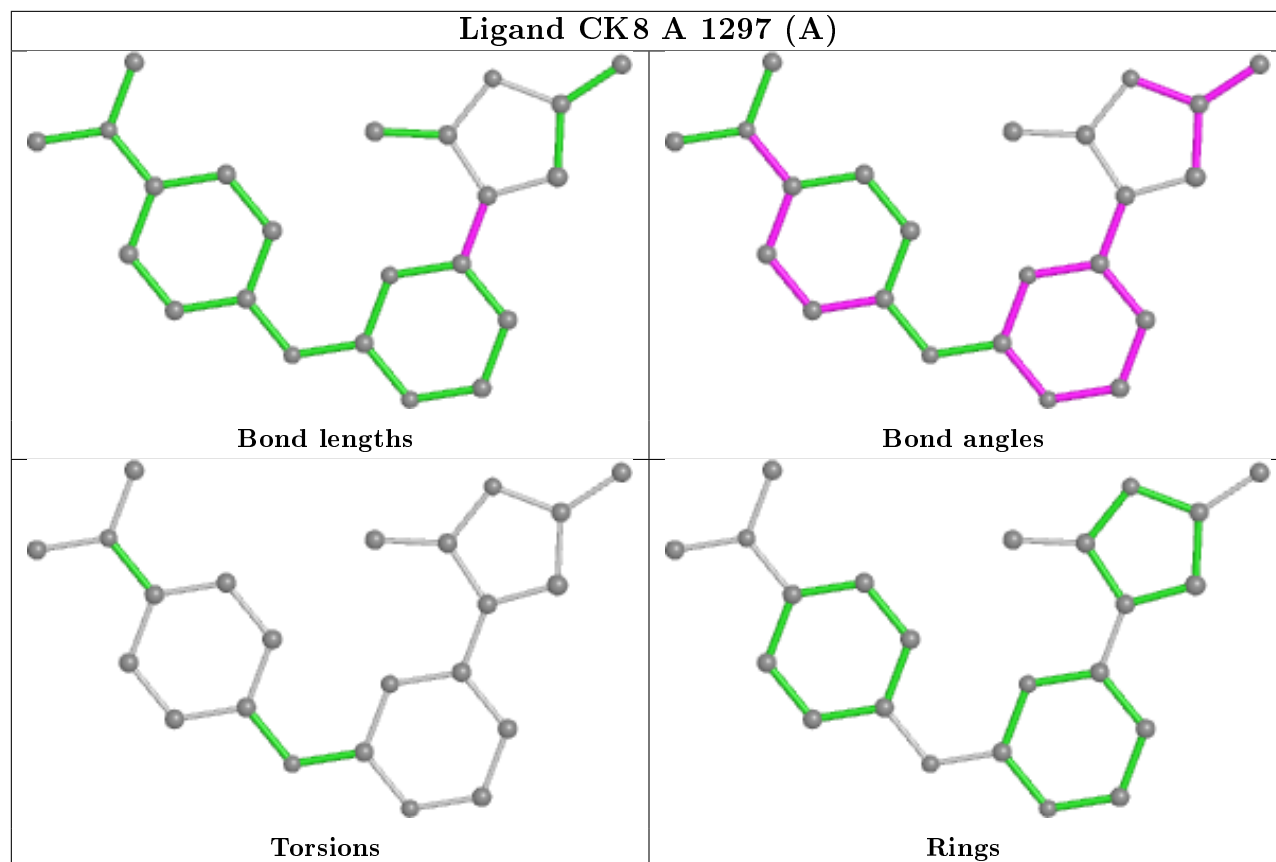
3 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1298	CK8	11	0
3	A	1297[B]	CK8	9	0
3	A	1297[A]	CK8	14	0

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







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	296/298 (99%)	-0.03	17 (5%)	23 29	13, 24, 51, 75	1 (0%)
1	C	297/298 (99%)	0.16	25 (8%)	11 14	13, 26, 61, 83	0
2	B	258/259 (99%)	0.05	12 (4%)	31 37	15, 29, 51, 79	0
2	D	258/259 (99%)	-0.11	6 (2%)	60 65	14, 26, 45, 59	0
All	All	1109/1114 (99%)	0.02	60 (5%)	25 31	13, 26, 53, 83	1 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	175	VAL	9.7
1	C	15	TYR	8.8
1	C	39	THR	8.4
1	C	161	HIS	8.2
1	C	13	GLY	8.1
1	C	37	LEU	7.5
2	B	271	TYR	6.8
1	C	163	VAL	6.5
1	C	14	THR	6.4
1	C	40	GLU	4.8
2	D	198	GLY	4.6
1	C	38	ASP	4.3
2	B	176	PRO	4.0
1	C	36	ARG	4.0
2	B	324	PRO	3.7
1	C	165	THR	3.7
1	A	14	THR	3.7
1	A	74	ASN	3.6
1	C	96	LEU	3.5
1	C	247	ASP	3.5
2	D	175	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	39	THR	3.3
1	A	95	ALA	3.3
1	C	73	GLU	3.0
1	C	12	GLU	3.0
2	B	280	TYR	2.9
1	C	16	GLY	2.8
2	B	323	GLN	2.8
1	C	162	GLU	2.8
2	B	178	TYR	2.8
1	A	164	VAL	2.7
1	C	72	THR	2.7
1	A	73	GLU	2.6
1	A	37	LEU	2.6
1	A	84	HIS	2.6
2	D	284	ASP	2.6
1	C	164	VAL	2.5
1	C	74	ASN	2.5
1	C	297	ARG	2.5
2	D	345	ASP	2.5
1	A	40	GLU	2.5
1	C	296	LEU	2.4
1	A	249	SER	2.4
1	A	96	LEU	2.3
1	A	13	GLY	2.3
2	B	345	ASP	2.3
1	C	200	ARG	2.3
2	B	327	CYS	2.3
1	A	295	HIS	2.2
1	A	162	GLU	2.2
1	A	160	THR	2.2
1	A	159	TYR	2.2
1	C	295	HIS	2.2
1	A	15	TYR	2.1
2	D	226	LYS	2.1
2	B	423	LEU	2.1
1	C	35	ILE	2.1
2	D	176	PRO	2.1
2	B	180	GLU	2.0
2	B	325	ALA	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

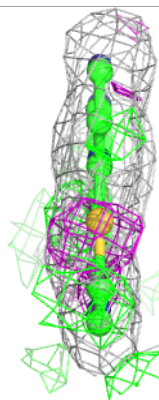
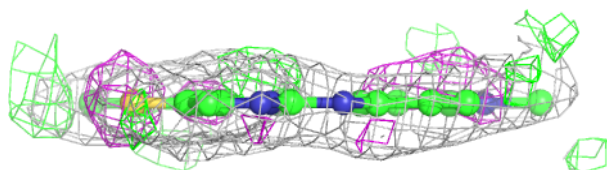
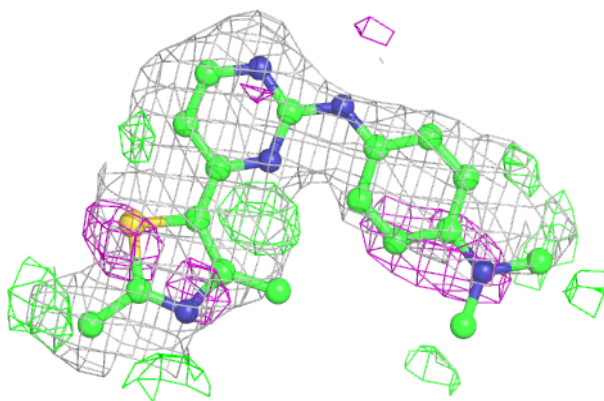
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	CK8	A	1297[B]	23/23	0.82	0.24	2,35,57,57	23
3	CK8	A	1297[A]	23/23	0.82	0.24	16,39,59,64	23
3	CK8	C	1298	23/23	0.91	0.13	17,36,50,52	0

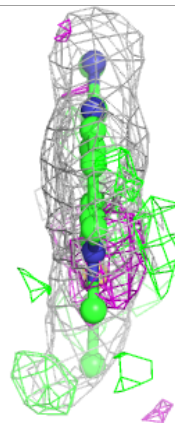
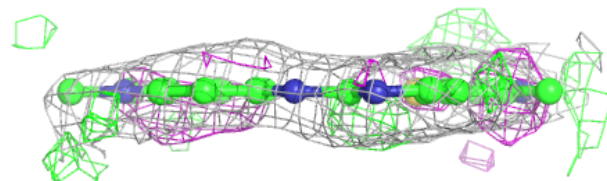
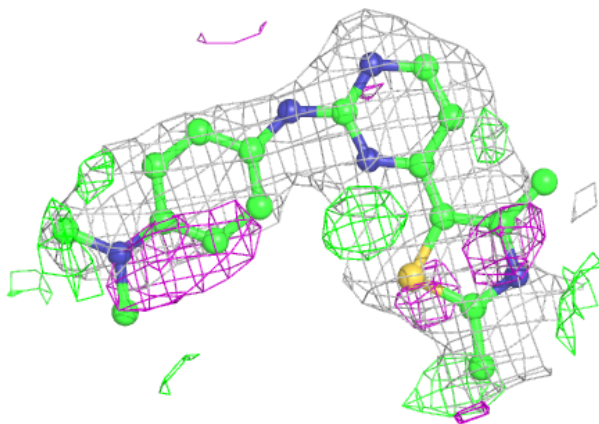
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

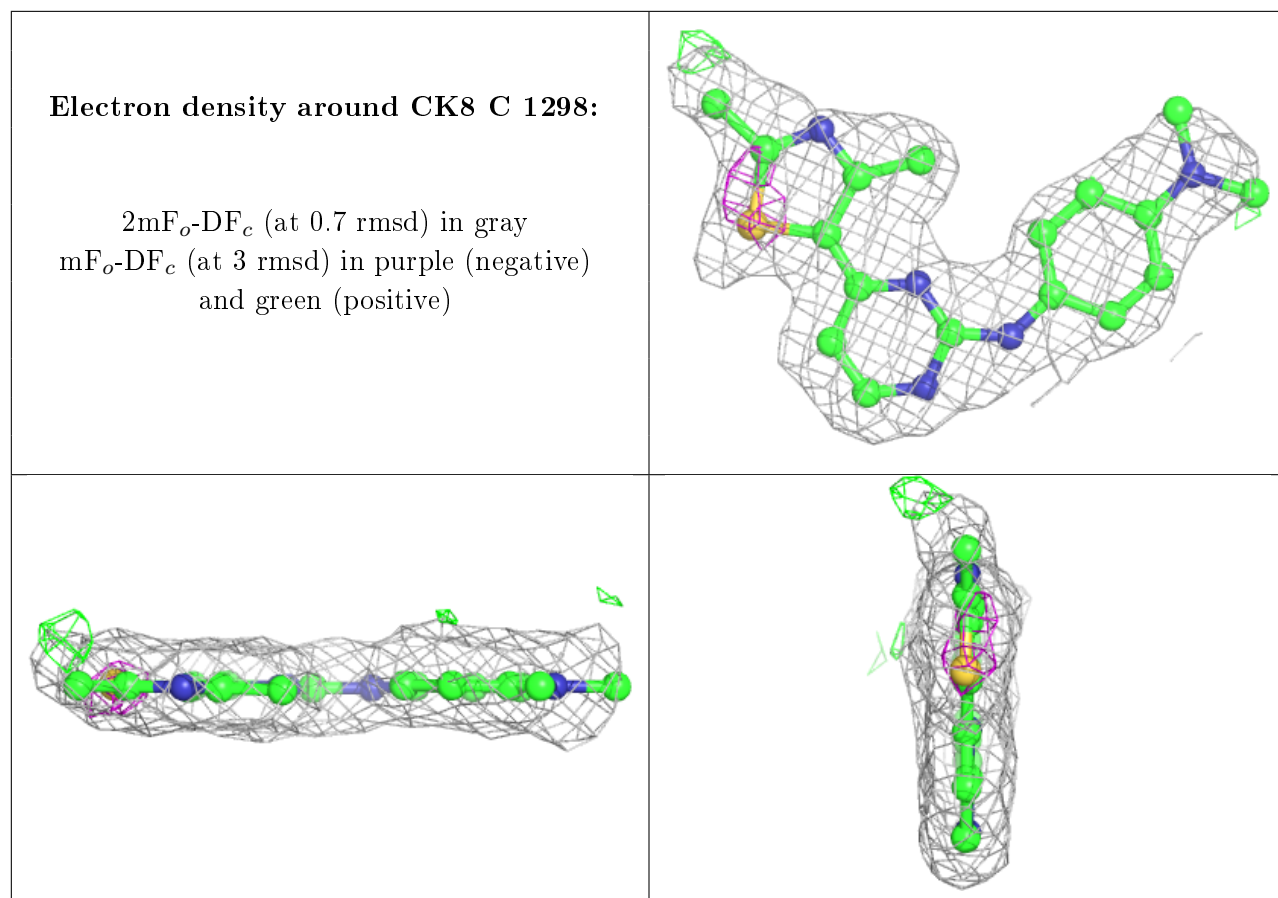
**Electron density around CK8 A 1297 (B):**

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

**Electron density around CK8 A 1297 (A):**

$2mF_o-DF_c$  (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.