



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:15 PM UTC

PDB ID : 1FIN / pdb_00001fin
Title : CYCLIN A-CYCLIN-DEPENDENT KINASE 2 COMPLEX
Authors : Jeffrey, P.D.; Russo, A.A.; Pavletich, N.P.
Deposited on : 1996-07-14
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

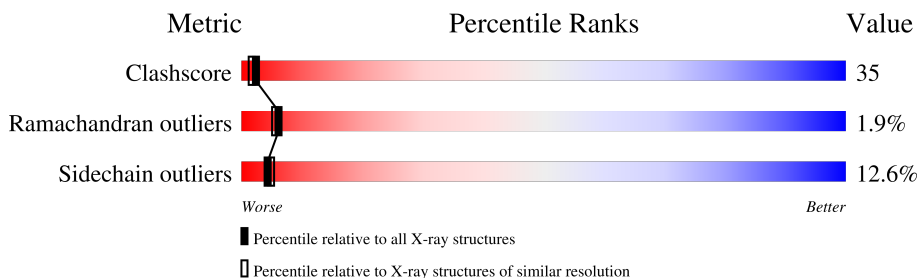
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	298	51% 37% 10% .
1	C	298	36% 51% 12% .
2	B	260	47% 42% 10% .
2	D	260	57% 34% 8% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9476 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.

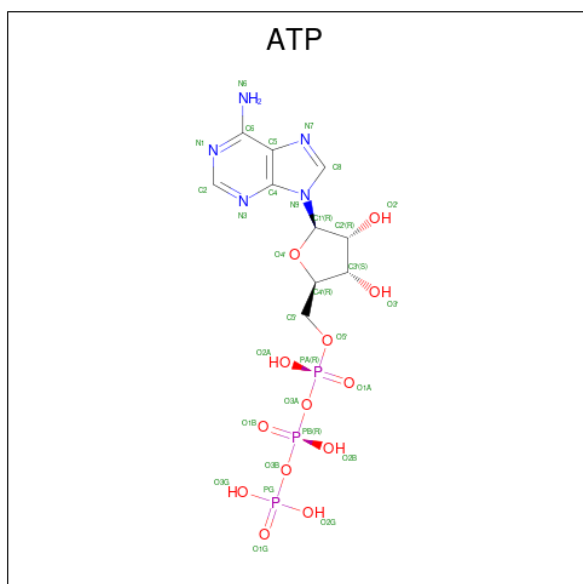
- Molecule 1 is a protein called CYCLIN-DEPENDENT KINASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total	C	N	O	S	0	0	0
			2398	1559	408	423	8			
1	C	298	Total	C	N	O	S	0	0	0
			2398	1559	408	423	8			

- Molecule 2 is a protein called CYCLIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	260	Total	C	N	O	S	0	0	0
			2101	1359	342	389	11			
2	D	260	Total	C	N	O	S	0	0	0
			2101	1359	342	389	11			

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
3	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 4 is water.

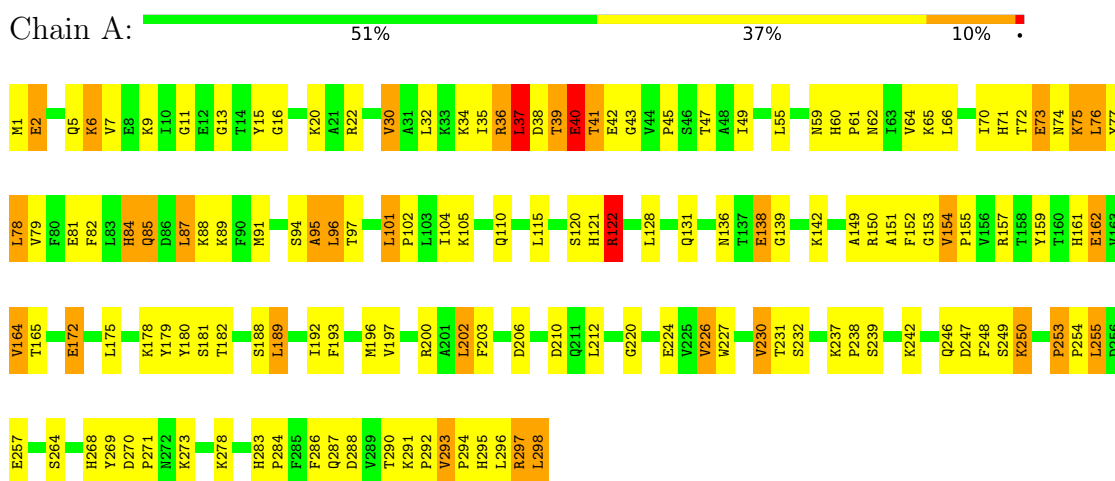
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	136	Total	O	0	0
			136	136		
4	B	96	Total	O	0	0
			96	96		
4	C	58	Total	O	0	0
			58	58		
4	D	126	Total	O	0	0
			126	126		

3 Residue-property plots

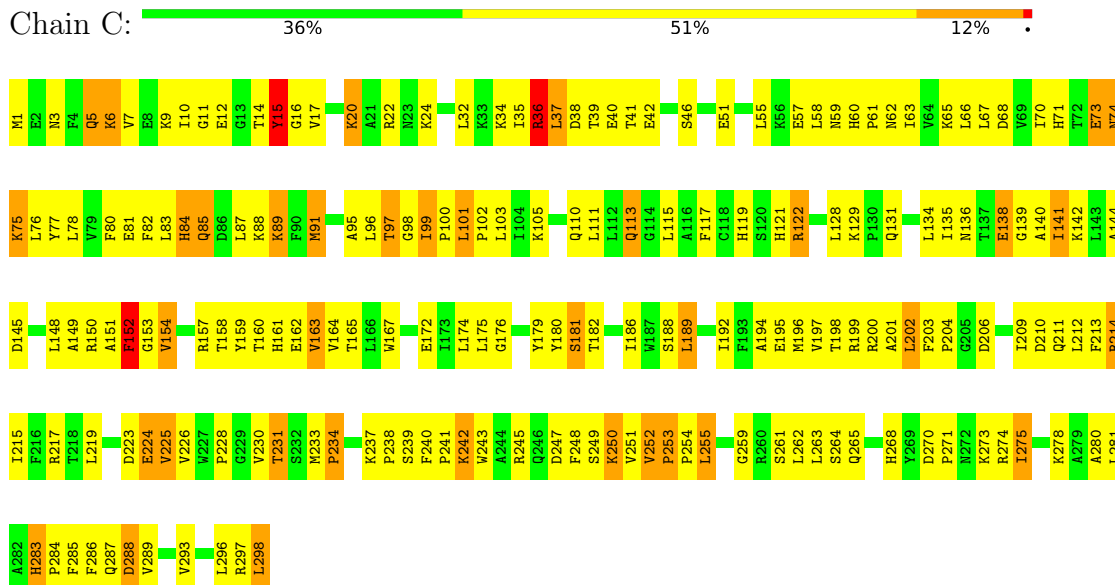
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

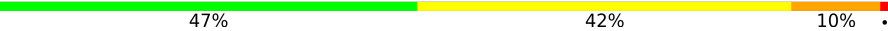
• Molecule 1: CYCLIN-DEPENDENT KINASE 2

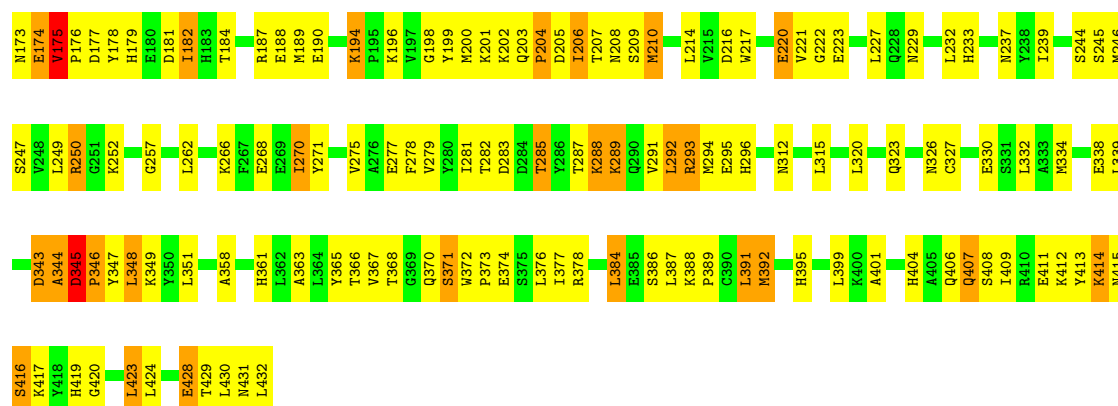


• Molecule 1: CYCLIN-DEPENDENT KINASE 2



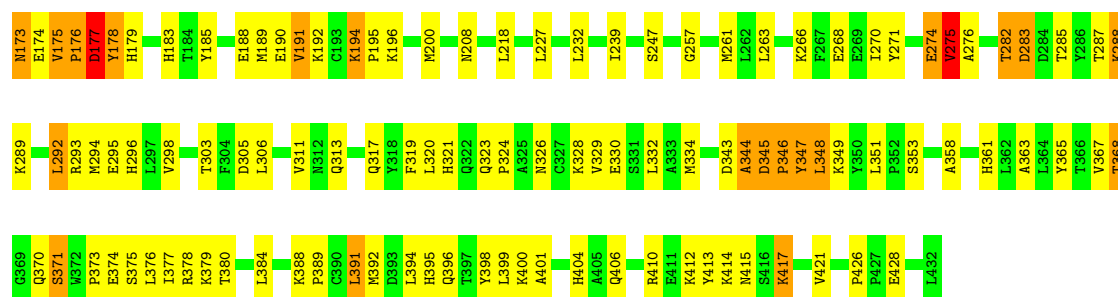
• Molecule 2: CYCLIN A

Chain B:  47% 42% 10%



• Molecule 2: CYCLIN A

Chain D:  57% 34% 8%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	185.10Å 185.10Å 214.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.30	Depositor
% Data completeness (in resolution range)	90.2 (6.00-2.30)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.208 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9476	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.81	0/2460	1.21	20/3338 (0.6%)
1	C	0.75	0/2460	1.21	18/3338 (0.5%)
2	B	0.79	0/2151	1.22	18/2920 (0.6%)
2	D	0.80	0/2151	1.16	16/2920 (0.5%)
All	All	0.79	0/9222	1.20	72/12516 (0.6%)

There are no bond length outliers.

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	347	TYR	N-CA-C	11.70	124.03	111.28
2	B	344	ALA	N-CA-C	-10.04	100.16	111.71
2	D	344	ALA	N-CA-C	-9.87	100.34	112.38
1	C	84	HIS	N-CA-C	9.65	124.39	112.23
2	B	174	GLU	N-CA-C	9.45	130.92	110.80
2	D	347	TYR	N-CA-C	8.80	121.83	111.71
2	B	367	VAL	N-CA-C	8.26	118.83	110.82
1	C	73	GLU	N-CA-C	-7.97	101.88	111.69
1	A	95	ALA	N-CA-C	-7.92	99.58	110.35
1	A	164	VAL	CB-CA-C	-7.82	101.20	111.13
2	D	414	LYS	N-CA-C	-7.79	103.80	113.38
1	C	15	TYR	N-CA-C	7.76	122.13	113.21
1	A	253	PRO	N-CA-C	7.72	120.12	110.70
1	C	20	LYS	N-CA-C	-7.53	96.97	109.24
1	C	36	ARG	N-CA-C	7.46	119.50	111.36
2	D	194	LYS	N-CA-C	7.37	120.47	110.31
2	B	268	GLU	N-CA-C	7.32	120.69	111.69
1	A	81	GLU	N-CA-C	-7.24	100.94	110.43
1	A	84	HIS	N-CA-C	7.17	118.74	111.07
1	A	153	GLY	N-CA-C	-7.11	106.50	115.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	175	VAL	N-CA-C	-7.10	103.80	112.67
1	C	249	SER	N-CA-C	-6.85	103.88	111.82
1	C	99	ILE	N-CA-C	-6.79	102.79	108.63
1	A	20	LYS	N-CA-C	-6.72	99.64	110.32
1	C	283	HIS	N-CA-C	6.55	117.13	109.60
2	D	268	GLU	N-CA-C	6.55	121.56	112.45
1	A	11	GLY	N-CA-C	6.52	120.08	110.63
2	B	371	SER	N-CA-C	6.48	119.72	109.81
1	A	122	ARG	N-CA-C	6.44	120.30	111.54
2	D	421	VAL	N-CA-C	6.40	119.05	111.05
2	D	275	VAL	CB-CA-C	-6.38	103.53	112.14
2	D	178	TYR	N-CA-C	6.34	121.02	113.28
2	D	428	GLU	N-CA-C	-6.31	104.50	111.82
2	D	177	ASP	N-CA-C	6.22	120.90	113.12
2	B	206	ILE	N-CA-C	6.09	117.10	110.21
2	D	306	LEU	N-CA-C	6.07	120.76	113.23
2	B	220	GLU	N-CA-C	-6.07	104.36	110.97
2	B	283	ASP	N-CA-C	-5.99	104.58	112.24
2	B	239	ILE	N-CA-C	5.98	116.15	110.53
1	A	248	PHE	N-CA-C	-5.97	105.26	112.54
1	A	7	VAL	CB-CA-C	-5.84	104.27	111.92
1	A	36	ARG	N-CA-C	5.83	117.71	111.36
2	B	414	LYS	N-CA-C	-5.76	104.66	112.26
2	B	252	LYS	N-CA-C	-5.76	106.12	114.12
1	C	238	PRO	N-CA-C	-5.75	106.96	114.03
1	C	296	LEU	N-CA-C	5.74	119.45	110.32
2	B	345	ASP	C-N-CD	-5.74	101.48	125.00
2	B	343	ASP	N-CA-C	5.71	118.94	107.41
1	C	165	THR	N-CA-C	-5.66	102.02	110.23
1	C	97	THR	N-CA-C	-5.66	98.74	110.80
1	C	214	ARG	N-CA-C	-5.66	105.11	111.28
1	A	249	SER	N-CA-C	-5.57	105.36	111.82
2	B	294	MET	N-CA-C	-5.57	105.36	111.82
2	D	282	THR	N-CA-C	-5.56	106.26	113.16
1	A	66	LEU	N-CA-C	-5.51	100.20	108.96
1	A	180	TYR	N-CA-C	-5.40	103.56	110.53
1	A	220	GLY	N-CA-C	-5.37	106.28	112.29
1	C	253	PRO	N-CA-C	5.33	117.21	110.70
2	B	401	ALA	N-CA-C	5.27	120.77	113.45
2	D	311	VAL	N-CA-C	-5.24	105.28	110.62
1	C	154	VAL	N-CA-C	-5.22	97.60	108.88
1	A	165	THR	N-CA-C	-5.14	103.25	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	152	PHE	N-CA-CB	5.12	119.14	110.49
2	B	175	VAL	C-N-CD	-5.11	104.07	125.00
1	A	78	LEU	N-CA-C	-5.10	100.85	109.07
2	D	175	VAL	CB-CA-C	-5.09	108.87	113.70
1	C	149	ALA	N-CA-C	5.07	117.39	110.24
2	D	368	THR	CB-CA-C	-5.07	102.16	110.72
1	A	85	GLN	N-CA-C	5.03	115.17	108.38
1	A	30	VAL	N-CA-CB	-5.02	104.03	112.06
1	C	153	GLY	N-CA-C	-5.01	101.30	113.18
2	B	428	GLU	N-CA-C	-5.00	106.77	112.92

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2398	0	2450	141	0
1	C	2398	0	2450	255	0
2	B	2101	0	2119	153	0
2	D	2101	0	2119	133	0
3	A	31	0	12	2	0
3	C	31	0	12	7	0
4	A	136	0	0	11	0
4	B	96	0	0	12	0
4	C	58	0	0	7	0
4	D	126	0	0	4	0
All	All	9476	0	9162	645	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 35.

All (645) 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:34:LYS:HZ1	1:C:75:LYS:NZ	1.45	1.13
1:C:278:LYS:HE3	2:D:177:ASP:HB3	1.32	1.12
2:B:210:MET:HE1	2:B:250:ARG:HB2	1.31	1.10
1:C:34:LYS:NZ	1:C:75:LYS:HZ2	1.48	1.09
2:D:374:GLU:HA	2:D:377:ILE:HD12	1.15	1.08
2:B:207:THR:HG22	2:B:210:MET:HG3	1.15	1.08
1:A:247:ASP:HB2	1:A:250:LYS:HE3	1.16	1.07
2:B:414:LYS:HE2	2:B:423:LEU:HD21	1.38	1.01
1:A:38:ASP:HB3	1:A:42:GLU:HB3	1.40	1.00
1:C:11:GLY:HA3	3:C:299:ATP:H5'2	1.39	1.00
2:B:332:LEU:HD23	2:B:363:ALA:HA	1.41	0.98
1:C:7:VAL:HG11	1:C:20:LYS:HD3	1.46	0.98
1:C:96:LEU:N	1:C:199:ARG:HE	1.61	0.98
2:B:404:HIS:HD2	2:B:406:GLN:H	0.99	0.97
1:A:73:GLU:HG3	1:A:74:ASN:H	1.30	0.96
2:D:404:HIS:HD2	2:D:406:GLN:H	1.01	0.96
1:A:122:ARG:HB3	2:B:182:ILE:HD13	1.47	0.96
2:D:289:LYS:HG3	2:D:293:ARG:HH21	1.28	0.94
2:B:207:THR:HG22	2:B:210:MET:CG	1.97	0.94
1:C:96:LEU:H	1:C:199:ARG:HE	1.11	0.94
1:C:60:HIS:HD2	1:C:62:ASN:H	1.14	0.91
2:B:250:ARG:HH11	2:B:250:ARG:HB3	1.36	0.91
2:D:185:TYR:CE1	2:D:189:MET:HE3	2.08	0.88
2:D:374:GLU:CA	2:D:377:ILE:HD12	2.02	0.88
2:B:404:HIS:CD2	2:B:406:GLN:H	1.90	0.88
1:C:7:VAL:CG1	1:C:20:LYS:HD3	2.03	0.87
2:D:398:TYR:CE2	2:D:426:PRO:HB3	2.08	0.87
1:A:60:HIS:HD2	1:A:62:ASN:H	1.21	0.87
1:C:1:MET:CE	1:C:70:ILE:HD13	2.04	0.87
1:C:96:LEU:HA	1:C:199:ARG:HH21	1.39	0.87
1:C:85:GLN:NE2	1:C:297:ARG:NH1	2.24	0.86
1:A:247:ASP:CB	1:A:250:LYS:HE3	2.04	0.86
2:B:374:GLU:HG3	2:B:378:ARG:NH1	1.90	0.86
1:A:227:TRP:O	1:A:230:VAL:HG22	1.76	0.85
1:C:1:MET:HE3	1:C:70:ILE:HD13	1.59	0.85
1:C:182:THR:OG1	2:D:174:GLU:HB2	1.77	0.85
1:C:99:ILE:HG23	1:C:103:LEU:HD23	1.56	0.85
2:D:404:HIS:CD2	2:D:406:GLN:H	1.92	0.84
1:C:60:HIS:CD2	1:C:62:ASN:H	1.94	0.84
1:C:11:GLY:HA3	3:C:299:ATP:C5'	2.06	0.84
2:D:343:ASP:HB3	2:D:345:ASP:HB2	1.59	0.84
2:B:270:ILE:HG22	2:B:271:TYR:HD1	1.43	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:VAL:HB	4:A:398:HOH:O	1.79	0.82
2:B:282:THR:O	2:B:285:THR:HG23	1.80	0.82
1:A:172:GLU:HG2	1:A:271:PRO:HG3	1.59	0.82
1:A:202:LEU:HD13	1:A:203:PHE:CE2	2.13	0.82
2:D:188:GLU:HG2	2:D:189:MET:HE2	1.60	0.81
2:D:191:VAL:HA	2:D:194:LYS:NZ	1.95	0.81
1:C:250:LYS:HD2	1:C:250:LYS:N	1.95	0.80
1:C:96:LEU:N	1:C:199:ARG:NE	2.29	0.80
1:C:34:LYS:HZ3	1:C:75:LYS:HG3	1.46	0.79
2:B:201:LYS:HZ3	2:B:202:LYS:HE2	1.48	0.79
1:C:35:ILE:O	1:C:75:LYS:HB2	1.83	0.79
2:B:201:LYS:HG3	2:B:202:LYS:N	1.96	0.78
2:B:196:LYS:HB2	2:B:244:SER:HB3	1.66	0.78
1:A:101:LEU:HD12	1:A:104:ILE:HD12	1.66	0.77
1:C:241:PRO:HB2	1:C:243:TRP:CZ3	2.20	0.77
1:C:34:LYS:HZ1	1:C:75:LYS:HZ2	0.79	0.77
1:C:159:TYR:CD2	1:C:162:GLU:HB2	2.20	0.77
1:A:6:LYS:HB2	1:A:6:LYS:HZ3	1.50	0.77
2:B:250:ARG:HH11	2:B:250:ARG:CB	1.97	0.77
1:C:11:GLY:CA	3:C:299:ATP:H5'2	2.15	0.76
1:C:84:HIS:HD2	1:C:136:ASN:HA	1.49	0.76
1:A:247:ASP:HB2	1:A:250:LYS:CE	2.09	0.76
1:A:131:GLN:HG2	4:A:349:HOH:O	1.84	0.76
2:D:388:LYS:O	2:D:392:MET:HG2	1.85	0.76
1:A:1:MET:HE1	1:A:70:ILE:CD1	2.16	0.76
1:C:96:LEU:HA	1:C:199:ARG:NH2	2.01	0.76
1:C:159:TYR:HD2	1:C:162:GLU:HB2	1.49	0.76
2:D:200:MET:HE3	2:D:208:ASN:HD22	1.51	0.75
1:A:157:ARG:HG3	1:A:179:TYR:CE1	2.20	0.75
2:B:217:TRP:O	2:B:221:VAL:HG23	1.86	0.75
1:A:15:TYR:CD1	1:A:35:ILE:HG12	2.21	0.74
2:B:332:LEU:HD23	2:B:363:ALA:CA	2.17	0.74
2:B:345:ASP:HB2	2:B:346:PRO:CD	2.17	0.74
2:B:200:MET:SD	2:B:206:ILE:HD12	2.28	0.74
2:D:175:VAL:N	2:D:176:PRO:HD2	2.02	0.74
1:C:253:PRO:HB2	1:C:254:PRO:HD3	1.70	0.74
2:B:332:LEU:CD2	2:B:363:ALA:HA	2.17	0.74
2:D:179:HIS:NE2	2:D:320:LEU:HD12	2.02	0.74
1:C:111:LEU:HD21	1:C:141:ILE:HG12	1.68	0.74
2:D:289:LYS:HG3	2:D:293:ARG:NH2	2.02	0.74
1:C:159:TYR:HD2	1:C:162:GLU:CB	2.01	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:9:LYS:NZ	1:C:17:VAL:HG11	2.03	0.73
1:A:60:HIS:CD2	1:A:62:ASN:H	2.04	0.73
2:B:179:HIS:CE1	2:B:320:LEU:HD12	2.24	0.73
2:D:334:MET:HE1	4:D:541:HOH:O	1.86	0.73
1:C:22:ARG:HH12	1:C:24:LYS:HA	1.52	0.73
2:B:210:MET:HE1	2:B:250:ARG:CB	2.16	0.73
1:C:34:LYS:NZ	1:C:75:LYS:HG3	2.03	0.73
1:A:40:GLU:HG3	1:A:41:THR:OG1	1.87	0.73
2:B:203:GLN:HG2	2:B:206:ILE:HG13	1.68	0.73
1:C:60:HIS:CG	1:C:61:PRO:HD2	2.23	0.73
1:C:152:PHE:CE2	2:D:173:ASN:HA	2.23	0.72
2:B:388:LYS:HG3	2:B:432:LEU:HD13	1.72	0.72
2:D:415:ASN:HD22	2:D:417:LYS:H	1.38	0.72
1:C:138:GLU:O	1:C:140:ALA:N	2.22	0.72
1:A:15:TYR:CE1	1:A:35:ILE:HD11	2.24	0.72
1:A:162:GLU:HB2	4:A:391:HOH:O	1.88	0.72
1:A:40:GLU:OE2	1:A:41:THR:HG23	1.90	0.71
2:B:345:ASP:HB2	2:B:346:PRO:HD3	1.73	0.71
1:A:73:GLU:CG	1:A:74:ASN:H	1.94	0.71
2:B:201:LYS:HG3	2:B:202:LYS:HG3	1.71	0.71
1:C:101:LEU:N	1:C:102:PRO:HD2	2.05	0.71
2:D:174:GLU:HG2	2:D:177:ASP:H	1.55	0.71
2:D:287:THR:OG1	2:D:289:LYS:HG2	1.89	0.71
1:C:247:ASP:HB3	1:C:250:LYS:HD3	1.73	0.71
2:D:319:PHE:HE2	2:D:334:MET:HE2	1.54	0.71
1:C:159:TYR:CZ	1:C:161:HIS:HB2	2.26	0.71
1:A:1:MET:HE3	1:A:77:TYR:CD1	2.25	0.71
2:B:312:ASN:OD1	2:B:334:MET:HE1	1.91	0.71
2:B:404:HIS:HD2	2:B:406:GLN:N	1.82	0.71
2:B:334:MET:HG3	4:B:527:HOH:O	1.90	0.71
1:A:1:MET:HE1	1:A:70:ILE:HD13	1.73	0.70
1:C:159:TYR:HD2	1:C:162:GLU:CG	2.04	0.70
1:C:5:GLN:OE1	1:C:22:ARG:NE	2.22	0.70
1:C:255:LEU:HG	1:C:259:GLY:HA3	1.74	0.70
2:B:388:LYS:HB3	2:B:389:PRO:HD3	1.72	0.70
1:C:59:ASN:OD1	1:C:65:LYS:HE2	1.92	0.70
1:A:38:ASP:CB	1:A:42:GLU:HB3	2.21	0.69
1:C:57:GLU:OE2	1:C:122:ARG:NH2	2.24	0.69
1:A:157:ARG:HG3	1:A:179:TYR:HE1	1.57	0.69
2:D:345:ASP:HB2	2:D:346:PRO:HD2	1.74	0.69
1:A:72:THR:OG1	1:A:75:LYS:HE3	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:74:ASN:HA	4:C:335:HOH:O	1.92	0.69
1:A:84:HIS:HB3	1:A:298:LEU:CD2	2.23	0.69
1:C:9:LYS:HZ2	1:C:17:VAL:HG11	1.57	0.69
2:D:174:GLU:OE2	2:D:176:PRO:HB2	1.92	0.69
2:B:374:GLU:HA	2:B:377:ILE:HD12	1.73	0.69
1:C:158:THR:HA	1:C:180:TYR:CE2	2.28	0.69
2:D:191:VAL:HA	2:D:194:LYS:HZ1	1.56	0.68
1:A:71:HIS:HA	1:A:76:LEU:HD23	1.75	0.68
1:C:22:ARG:NH1	1:C:24:LYS:HA	2.09	0.68
2:D:404:HIS:HD2	2:D:406:GLN:N	1.84	0.68
2:D:345:ASP:HB2	2:D:346:PRO:CD	2.24	0.68
1:C:195:GLU:HB2	4:C:337:HOH:O	1.93	0.67
1:C:270:ASP:CB	1:C:273:LYS:HE2	2.23	0.67
2:D:174:GLU:HA	4:D:502:HOH:O	1.92	0.67
1:A:36:ARG:O	1:A:37:LEU:HB2	1.94	0.67
2:B:262:LEU:HD11	2:B:266:LYS:HE3	1.75	0.67
1:C:213:PHE:CE2	1:C:241:PRO:HD2	2.30	0.67
1:C:128:LEU:HD13	1:C:189:LEU:HD13	1.77	0.67
1:A:89:LYS:HD2	1:A:89:LYS:N	2.10	0.67
1:C:89:LYS:HD3	1:C:89:LYS:N	2.10	0.67
2:D:412:LYS:HE2	2:D:413:TYR:CE1	2.29	0.66
1:C:95:ALA:HA	1:C:199:ARG:HD2	1.77	0.66
1:A:88:LYS:HD3	4:A:433:HOH:O	1.94	0.66
1:A:39:THR:HG23	4:B:513:HOH:O	1.95	0.66
1:C:15:TYR:CE1	1:C:35:ILE:HD11	2.31	0.66
1:A:84:HIS:HB3	1:A:298:LEU:HD21	1.78	0.66
1:C:200:ARG:HD2	1:C:201:ALA:H	1.60	0.66
2:B:201:LYS:HE2	2:B:202:LYS:CG	2.25	0.65
1:C:38:ASP:HB3	1:C:42:GLU:HB3	1.78	0.65
1:C:157:ARG:HH21	2:D:270:ILE:HG13	1.62	0.65
1:C:100:PRO:C	1:C:102:PRO:HD2	2.22	0.65
1:C:159:TYR:HB3	1:C:162:GLU:HB2	1.78	0.65
2:D:191:VAL:HG22	2:D:194:LYS:NZ	2.12	0.65
2:B:201:LYS:HE2	2:B:202:LYS:HG3	1.79	0.65
2:B:207:THR:HG21	4:B:455:HOH:O	1.95	0.64
2:B:270:ILE:HG22	2:B:271:TYR:CD1	2.31	0.64
2:D:345:ASP:CB	2:D:346:PRO:CD	2.76	0.64
1:C:172:GLU:HG2	1:C:271:PRO:HG3	1.80	0.64
2:D:174:GLU:CG	2:D:176:PRO:HB2	2.27	0.64
1:C:85:GLN:HE21	1:C:297:ARG:NH1	1.95	0.64
1:C:181:SER:HA	2:D:173:ASN:N	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:138:GLU:HA	1:C:293:VAL:HG13	1.79	0.64
2:D:174:GLU:HG3	2:D:176:PRO:CG	2.27	0.64
1:A:15:TYR:HD1	1:A:35:ILE:HG12	1.60	0.64
2:B:293:ARG:NH1	1:C:3:ASN:OD1	2.31	0.64
1:C:55:LEU:CD1	1:C:80:PHE:HE1	2.12	0.63
2:D:174:GLU:HG3	2:D:176:PRO:HB2	1.79	0.63
2:B:179:HIS:NE2	2:B:320:LEU:HD12	2.13	0.63
1:C:40:GLU:O	2:D:288:LYS:HD2	1.99	0.63
1:A:59:ASN:HD21	1:A:65:LYS:HD2	1.64	0.63
4:A:330:HOH:O	2:B:296:HIS:HB3	1.98	0.63
2:D:176:PRO:HG2	4:D:502:HOH:O	1.98	0.63
1:A:6:LYS:HB2	1:A:6:LYS:NZ	2.13	0.63
2:B:250:ARG:HB3	2:B:250:ARG:NH1	2.11	0.63
1:C:270:ASP:HB3	1:C:273:LYS:HE2	1.79	0.63
2:D:395:HIS:CE1	2:D:399:LEU:HD11	2.34	0.63
1:C:60:HIS:CD2	1:C:61:PRO:HD2	2.34	0.62
1:C:98:GLY:HA2	1:C:199:ARG:HD3	1.81	0.62
1:C:248:PHE:HA	1:C:251:VAL:HB	1.81	0.62
1:A:1:MET:CE	1:A:70:ILE:HG21	2.29	0.62
1:A:159:TYR:CZ	1:A:161:HIS:HB2	2.35	0.62
2:B:343:ASP:HB3	2:B:346:PRO:HD2	1.81	0.62
1:C:61:PRO:O	1:C:142:LYS:HD3	1.99	0.62
2:D:396:GLN:O	2:D:400:LYS:HG3	1.99	0.62
1:A:182:THR:H	2:B:173:ASN:N	1.97	0.62
2:B:203:GLN:HG3	4:B:433:HOH:O	2.00	0.62
1:C:213:PHE:HD2	1:C:243:TRP:CH2	2.18	0.61
1:C:241:PRO:HB2	1:C:243:TRP:CH2	2.35	0.61
1:C:283:HIS:CG	1:C:284:PRO:HD2	2.34	0.61
2:D:174:GLU:HG3	2:D:176:PRO:CB	2.30	0.61
2:B:374:GLU:O	2:B:378:ARG:HG3	2.00	0.61
1:C:152:PHE:CD2	2:D:173:ASN:HA	2.35	0.61
1:A:128:LEU:HD13	1:A:189:LEU:HD13	1.82	0.61
1:C:297:ARG:O	1:C:298:LEU:HB2	1.99	0.61
3:A:299:ATP:H3'	4:A:429:HOH:O	2.00	0.61
2:B:278:PHE:O	2:B:282:THR:HG23	2.00	0.61
1:A:1:MET:CE	1:A:70:ILE:HD13	2.30	0.61
1:A:41:THR:HG23	2:B:288:LYS:NZ	2.16	0.60
1:C:60:HIS:HD2	1:C:62:ASN:N	1.94	0.60
1:C:242:LYS:HB2	1:C:242:LYS:NZ	2.16	0.60
2:B:344:ALA:HB1	2:B:348:LEU:HD22	1.81	0.60
1:C:96:LEU:H	1:C:199:ARG:NE	1.90	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:222:GLY:HA2	2:B:227:LEU:HD12	1.82	0.60
1:C:297:ARG:HH11	1:C:298:LEU:HB2	1.66	0.60
1:A:71:HIS:HE1	2:B:296:HIS:CE1	2.20	0.59
2:B:198:GLY:O	2:B:201:LYS:HG2	2.01	0.59
2:B:201:LYS:NZ	2:B:202:LYS:HE2	2.16	0.59
1:C:247:ASP:OD2	1:C:250:LYS:HD3	2.01	0.59
2:D:200:MET:HE3	2:D:208:ASN:ND2	2.17	0.59
2:D:398:TYR:CD2	2:D:426:PRO:HB3	2.36	0.59
1:C:122:ARG:O	1:C:151:ALA:HA	2.02	0.59
1:C:215:ILE:HG23	1:C:219:LEU:HD12	1.82	0.59
1:C:46:SER:HB2	2:D:266:LYS:O	2.02	0.59
2:B:345:ASP:CB	2:B:346:PRO:HD3	2.31	0.59
2:D:183:HIS:CE1	2:D:380:THR:HG22	2.38	0.59
1:C:223:ASP:OD1	1:C:225:VAL:HG12	2.02	0.58
2:D:326:ASN:O	2:D:330:GLU:HG3	2.03	0.58
1:A:1:MET:HE2	1:A:70:ILE:HG21	1.83	0.58
1:C:213:PHE:CD2	1:C:243:TRP:CH2	2.91	0.58
2:D:374:GLU:HA	2:D:377:ILE:CD1	2.10	0.58
1:C:247:ASP:O	1:C:250:LYS:HB2	2.04	0.58
2:B:332:LEU:HB3	2:B:363:ALA:HB1	1.86	0.58
1:C:172:GLU:OE2	1:C:274:ARG:NH1	2.36	0.58
2:D:274:GLU:OE1	2:D:276:ALA:N	2.30	0.58
2:B:275:VAL:O	2:B:279:VAL:HG23	2.04	0.58
2:B:287:THR:HG22	2:B:288:LYS:N	2.18	0.58
1:C:288:ASP:OD1	1:C:288:ASP:N	2.36	0.58
1:C:134:LEU:HD12	1:C:144:ALA:HB2	1.85	0.58
2:B:374:GLU:HG3	2:B:378:ARG:HH11	1.68	0.57
1:C:159:TYR:CD2	1:C:162:GLU:CB	2.83	0.57
2:D:282:THR:O	2:D:285:THR:HG23	2.03	0.57
1:C:84:HIS:HD2	1:C:136:ASN:CA	2.15	0.57
1:C:157:ARG:HD3	1:C:157:ARG:N	2.19	0.57
1:A:202:LEU:HD13	1:A:203:PHE:CZ	2.40	0.57
1:C:3:ASN:HA	1:C:24:LYS:HE3	1.86	0.57
1:C:34:LYS:CE	1:C:75:LYS:HZ2	2.16	0.57
2:B:189:MET:HG3	4:B:498:HOH:O	2.04	0.57
2:B:206:ILE:HG22	2:B:250:ARG:HA	1.87	0.57
1:A:71:HIS:CE1	2:B:296:HIS:ND1	2.73	0.57
1:C:210:ASP:O	1:C:214:ARG:HG3	2.04	0.57
1:A:73:GLU:CG	1:A:75:LYS:HG2	2.35	0.56
1:C:12:GLU:CD	1:C:17:VAL:HG12	2.30	0.56
1:A:89:LYS:HD2	1:A:89:LYS:H	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:ASP:HB2	4:A:390:HOH:O	2.03	0.56
2:B:175:VAL:O	2:B:177:ASP:N	2.38	0.56
2:B:245:SER:C	2:B:246:MET:HG2	2.29	0.56
1:C:268:HIS:HD2	1:C:270:ASP:N	2.02	0.56
2:D:332:LEU:HD23	2:D:363:ALA:HA	1.85	0.56
1:C:95:ALA:HA	1:C:199:ARG:CD	2.35	0.56
2:B:351:LEU:HD12	4:B:494:HOH:O	2.05	0.56
1:C:157:ARG:HG3	1:C:179:TYR:HE1	1.70	0.56
1:C:188:SER:O	1:C:192:ILE:HG13	2.06	0.56
2:B:207:THR:CG2	2:B:210:MET:HG3	2.10	0.56
1:C:88:LYS:HA	1:C:91:MET:HE2	1.87	0.56
1:C:150:ARG:HD2	4:C:315:HOH:O	2.06	0.56
1:C:201:ALA:HB3	1:C:204:PRO:HG3	1.88	0.56
1:A:1:MET:HE1	1:A:70:ILE:HB	1.88	0.56
1:C:36:ARG:O	1:C:37:LEU:HB2	2.05	0.56
2:B:395:HIS:CE1	2:B:399:LEU:HD11	2.42	0.55
1:C:159:TYR:CE1	2:D:271:TYR:OH	2.58	0.55
2:D:227:LEU:HD12	2:D:261:MET:HE3	1.88	0.55
2:B:345:ASP:CB	2:B:346:PRO:CD	2.80	0.55
1:C:223:ASP:OD1	1:C:226:VAL:HG12	2.06	0.55
2:D:191:VAL:HA	2:D:194:LYS:CE	2.36	0.55
1:A:40:GLU:O	2:B:288:LYS:NZ	2.35	0.55
2:D:319:PHE:CE2	2:D:334:MET:HE2	2.37	0.55
2:B:223:GLU:HA	2:B:223:GLU:OE1	2.06	0.55
2:D:194:LYS:HZ1	2:D:351:LEU:HD21	1.72	0.55
2:B:208:ASN:HB3	2:B:345:ASP:OD1	2.07	0.55
2:B:216:ASP:OD1	2:B:408:SER:HB2	2.07	0.55
2:B:411:GLU:O	2:B:414:LYS:HG2	2.07	0.55
1:C:16:GLY:HA3	1:C:34:LYS:O	2.07	0.54
1:C:51:GLU:O	1:C:55:LEU:HB2	2.08	0.54
1:C:100:PRO:HB2	1:C:102:PRO:HG2	1.88	0.54
1:C:253:PRO:CB	1:C:254:PRO:HD3	2.35	0.54
1:C:297:ARG:NH1	1:C:298:LEU:HB2	2.22	0.54
1:A:61:PRO:O	1:A:142:LYS:HE2	2.07	0.54
1:C:157:ARG:CD	1:C:157:ARG:H	2.20	0.54
1:A:71:HIS:CE1	2:B:296:HIS:CE1	2.96	0.54
2:B:414:LYS:HA	2:B:420:GLY:HA2	1.89	0.54
1:C:135:ILE:CG2	1:C:141:ILE:HD12	2.38	0.54
1:C:36:ARG:HD3	1:C:75:LYS:HZ3	1.73	0.54
2:B:365:TYR:OH	2:B:431:ASN:ND2	2.41	0.54
1:C:113:GLN:HE22	1:C:281:LEU:HD22	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:159:TYR:CE2	1:C:161:HIS:HB2	2.42	0.53
2:D:294:MET:O	2:D:298:VAL:HG23	2.08	0.53
1:C:209:ILE:HG22	4:C:342:HOH:O	2.07	0.53
1:A:162:GLU:HG3	4:A:360:HOH:O	2.08	0.53
1:A:91:MET:HE3	1:A:196:MET:HG2	1.91	0.53
1:A:269:TYR:O	1:A:271:PRO:HD3	2.08	0.53
2:B:327:CYS:SG	2:B:419:HIS:NE2	2.79	0.53
2:B:345:ASP:HB2	2:B:346:PRO:HD2	1.88	0.53
2:D:175:VAL:N	2:D:176:PRO:CD	2.69	0.53
2:B:223:GLU:OE2	2:B:412:LYS:HG3	2.09	0.53
2:D:401:ALA:HB1	2:D:410:ARG:HD2	1.91	0.53
2:B:175:VAL:O	2:B:175:VAL:HG12	2.06	0.53
2:D:345:ASP:CB	2:D:346:PRO:HD3	2.39	0.53
1:C:213:PHE:HE2	1:C:241:PRO:HD2	1.72	0.53
1:A:253:PRO:HB2	1:A:254:PRO:HD3	1.91	0.52
1:C:160:THR:O	1:C:163:VAL:HG22	2.09	0.52
2:B:206:ILE:CG2	2:B:250:ARG:HA	2.39	0.52
1:C:60:HIS:HB3	1:C:63:ILE:HG13	1.90	0.52
1:C:71:HIS:CE1	2:D:296:HIS:CD2	2.97	0.52
1:C:157:ARG:N	1:C:157:ARG:CD	2.72	0.52
1:C:102:PRO:HG2	4:C:351:HOH:O	2.10	0.52
1:C:245:ARG:NH2	1:C:248:PHE:CE1	2.77	0.52
2:D:344:ALA:HB1	2:D:348:LEU:HD22	1.91	0.52
1:C:242:LYS:HB2	1:C:242:LYS:HZ2	1.73	0.52
1:C:297:ARG:HH12	1:C:298:LEU:HD23	1.74	0.52
2:B:412:LYS:HD3	2:B:413:TYR:CE1	2.45	0.52
1:C:12:GLU:HA	1:C:17:VAL:HA	1.91	0.52
1:A:1:MET:N	4:A:367:HOH:O	2.42	0.52
1:A:15:TYR:CE1	1:A:35:ILE:CD1	2.92	0.52
1:C:209:ILE:HD11	1:C:213:PHE:HE1	1.74	0.52
2:D:174:GLU:HG3	2:D:176:PRO:HG2	1.91	0.52
2:B:201:LYS:HG3	2:B:202:LYS:H	1.74	0.52
2:D:358:ALA:HB1	2:D:391:LEU:HD13	1.92	0.52
1:C:9:LYS:HZ2	1:C:17:VAL:CG1	2.22	0.51
1:A:283:HIS:CG	1:A:284:PRO:HD2	2.45	0.51
1:C:159:TYR:OH	1:C:161:HIS:HB2	2.10	0.51
1:C:268:HIS:CD2	1:C:270:ASP:N	2.78	0.51
2:D:321:HIS:HD2	2:D:375:SER:HB2	1.74	0.51
1:C:202:LEU:HD13	1:C:203:PHE:CE2	2.45	0.51
2:D:188:GLU:CG	2:D:189:MET:HE2	2.37	0.51
1:A:73:GLU:CG	1:A:74:ASN:N	2.66	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:179:HIS:CE1	2:B:320:LEU:CD1	2.93	0.51
2:D:367:VAL:HG12	2:D:368:THR:HG23	1.92	0.51
2:D:178:TYR:CD1	2:D:178:TYR:N	2.78	0.51
1:C:250:LYS:O	1:C:253:PRO:HD3	2.11	0.51
1:A:40:GLU:HG3	1:A:41:THR:N	2.26	0.51
1:C:175:LEU:HD21	1:C:212:LEU:HD21	1.91	0.51
1:C:110:GLN:O	1:C:113:GLN:HB2	2.11	0.51
1:C:224:GLU:OE1	1:C:231:THR:HB	2.11	0.51
1:C:65:LYS:HB3	1:C:81:GLU:HG2	1.93	0.51
1:C:157:ARG:HG3	1:C:179:TYR:CE1	2.46	0.51
1:A:290:THR:C	1:A:292:PRO:HD3	2.36	0.50
2:B:282:THR:O	2:B:285:THR:CG2	2.57	0.50
2:B:339:LEU:HD23	2:B:409:ILE:HG21	1.92	0.50
1:C:158:THR:HA	1:C:180:TYR:HE2	1.73	0.50
1:A:41:THR:CG2	2:B:288:LYS:NZ	2.74	0.50
1:C:217:ARG:HG2	1:C:243:TRP:CE2	2.46	0.50
2:D:190:GLU:OE2	2:D:353:SER:OG	2.28	0.50
2:B:292:LEU:O	2:B:295:GLU:HB3	2.11	0.50
1:C:194:ALA:O	1:C:198:THR:HG23	2.11	0.50
2:B:370:GLN:CD	2:B:371:SER:H	2.19	0.50
1:C:152:PHE:N	1:C:152:PHE:CD1	2.79	0.50
1:C:247:ASP:O	1:C:250:LYS:N	2.44	0.50
1:C:9:LYS:HD2	1:C:17:VAL:HB	1.92	0.50
1:C:96:LEU:CA	1:C:199:ARG:HE	2.25	0.50
2:B:204:PRO:HG2	4:B:433:HOH:O	2.12	0.50
1:C:34:LYS:NZ	1:C:75:LYS:NZ	2.27	0.50
1:C:242:LYS:NZ	1:C:242:LYS:CB	2.75	0.49
2:D:174:GLU:CD	2:D:176:PRO:HB2	2.37	0.49
1:C:85:GLN:NE2	1:C:298:LEU:CB	2.75	0.49
1:C:278:LYS:HE3	2:D:177:ASP:CB	2.23	0.49
1:C:283:HIS:ND1	1:C:284:PRO:HD2	2.27	0.49
2:B:412:LYS:HD3	2:B:413:TYR:HE1	1.78	0.49
1:C:60:HIS:CD2	1:C:61:PRO:CD	2.95	0.49
1:C:203:PHE:CB	1:C:211:GLN:HE22	2.25	0.49
1:A:149:ALA:O	1:A:150:ARG:HG2	2.13	0.49
1:A:294:PRO:HB2	1:A:296:LEU:HD23	1.94	0.49
1:C:15:TYR:O	1:C:35:ILE:HG12	2.13	0.49
1:A:91:MET:HE3	1:A:196:MET:CG	2.42	0.49
1:C:84:HIS:CD2	1:C:136:ASN:HA	2.38	0.49
2:D:175:VAL:HB	2:D:176:PRO:CD	2.42	0.49
2:D:358:ALA:CB	2:D:391:LEU:HD13	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:GLU:HG3	1:A:75:LYS:HG2	1.95	0.49
1:A:95:ALA:O	1:A:96:LEU:CB	2.61	0.49
1:C:200:ARG:HD2	1:C:201:ALA:N	2.26	0.49
2:B:207:THR:HG23	2:B:210:MET:H	1.77	0.49
2:B:373:PRO:HG2	2:B:376:LEU:HB2	1.94	0.49
1:C:74:ASN:HB2	1:C:75:LYS:HE2	1.94	0.49
1:C:85:GLN:NE2	1:C:298:LEU:HB3	2.28	0.49
1:A:122:ARG:O	1:A:151:ALA:HA	2.12	0.48
2:D:174:GLU:CG	2:D:177:ASP:H	2.23	0.48
2:D:313:GLN:O	2:D:317:GLN:HG2	2.13	0.48
1:A:224:GLU:OE2	1:A:231:THR:OG1	2.21	0.48
1:C:101:LEU:N	1:C:102:PRO:CD	2.75	0.48
2:D:274:GLU:CD	2:D:276:ALA:H	2.20	0.48
2:D:282:THR:O	2:D:283:ASP:HB2	2.12	0.48
1:A:38:ASP:CG	1:A:41:THR:H	2.22	0.48
1:C:40:GLU:O	2:D:288:LYS:HG2	2.13	0.48
1:C:75:LYS:H	1:C:75:LYS:HD2	1.79	0.48
1:C:159:TYR:CZ	2:D:271:TYR:CE1	3.02	0.48
2:B:200:MET:HE3	2:B:208:ASN:ND2	2.29	0.48
1:C:268:HIS:CD2	1:C:270:ASP:H	2.31	0.48
1:C:75:LYS:HD2	1:C:75:LYS:N	2.28	0.48
2:D:292:LEU:HD12	2:D:292:LEU:HA	1.55	0.48
1:A:36:ARG:O	1:A:74:ASN:O	2.31	0.48
1:C:36:ARG:HD3	1:C:75:LYS:NZ	2.29	0.48
1:C:39:THR:HG22	1:C:39:THR:O	2.13	0.48
1:A:227:TRP:CE3	1:A:230:VAL:HG13	2.49	0.48
2:D:275:VAL:HG11	2:D:292:LEU:HD13	1.96	0.48
1:A:41:THR:HG23	2:B:288:LYS:HZ1	1.78	0.48
1:C:167:TRP:CZ3	1:C:201:ALA:HB2	2.49	0.48
1:C:77:TYR:C	1:C:78:LEU:HD23	2.38	0.48
1:C:223:ASP:H	1:C:226:VAL:HG12	1.79	0.48
1:C:1:MET:HG2	1:C:70:ILE:HG21	1.96	0.47
1:A:120:SER:HB2	2:B:181:ASP:HB3	1.96	0.47
2:B:194:LYS:HE3	4:B:473:HOH:O	2.14	0.47
2:D:175:VAL:HA	2:D:179:HIS:HB2	1.96	0.47
2:D:185:TYR:O	2:D:189:MET:HG2	2.14	0.47
2:D:288:LYS:HB2	2:D:288:LYS:HE3	1.54	0.47
2:D:415:ASN:ND2	4:D:444:HOH:O	2.46	0.47
3:A:299:ATP:O2B	3:A:299:ATP:H5'2	2.14	0.47
4:C:344:HOH:O	2:D:189:MET:HB3	2.13	0.47
1:A:159:TYR:HB2	2:B:270:ILE:HG13	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:185:TYR:HE1	2:D:189:MET:HE3	1.70	0.47
1:C:40:GLU:O	1:C:40:GLU:HG3	2.14	0.47
2:D:179:HIS:CE1	2:D:379:LYS:NZ	2.82	0.47
1:A:291:LYS:N	1:A:292:PRO:HD3	2.29	0.47
2:B:407:GLN:HA	4:B:492:HOH:O	2.14	0.47
1:C:82:PHE:CD1	1:C:82:PHE:C	2.92	0.47
1:C:223:ASP:C	1:C:225:VAL:H	2.21	0.47
2:D:398:TYR:CZ	2:D:426:PRO:HB3	2.50	0.47
1:A:38:ASP:C	1:A:40:GLU:H	2.23	0.47
2:B:392:MET:HA	2:B:392:MET:HE3	1.97	0.47
2:B:428:GLU:C	2:B:429:THR:HG23	2.40	0.47
1:C:172:GLU:OE2	1:C:274:ARG:NH2	2.44	0.47
2:D:303:THR:OG1	2:D:305:ASP:OD2	2.32	0.47
2:D:343:ASP:C	2:D:345:ASP:N	2.66	0.47
1:A:2:GLU:H	1:A:2:GLU:CD	2.23	0.47
1:A:206:ASP:N	1:A:210:ASP:OD2	2.45	0.47
1:C:36:ARG:HE	1:C:36:ARG:HB2	1.51	0.47
1:A:5:GLN:HG2	1:A:22:ARG:HH21	1.81	0.46
2:D:191:VAL:HA	2:D:194:LYS:HE3	1.97	0.46
1:A:49:ILE:CD1	2:B:266:LYS:HD2	2.45	0.46
1:A:159:TYR:CD2	1:A:162:GLU:HG3	2.50	0.46
1:C:206:ASP:N	1:C:210:ASP:OD2	2.44	0.46
2:B:205:ASP:OD2	2:B:250:ARG:HG2	2.15	0.46
2:B:287:THR:HG22	2:B:288:LYS:H	1.79	0.46
1:C:230:VAL:HG13	1:C:231:THR:N	2.31	0.46
1:A:72:THR:OG1	1:A:75:LYS:HG3	2.15	0.46
2:B:395:HIS:HB2	2:B:430:LEU:HD11	1.98	0.46
2:D:329:VAL:HG23	2:D:367:VAL:HG11	1.96	0.46
2:D:373:PRO:HG2	2:D:376:LEU:HB2	1.97	0.46
1:A:1:MET:HE1	1:A:70:ILE:CB	2.44	0.46
1:C:95:ALA:HA	1:C:199:ARG:CG	2.46	0.46
1:A:15:TYR:CD1	1:A:15:TYR:C	2.94	0.46
1:C:159:TYR:CD2	1:C:162:GLU:CG	2.92	0.46
1:A:43:GLY:O	1:A:45:PRO:HD3	2.16	0.46
1:A:226:VAL:HG12	1:A:227:TRP:N	2.31	0.46
1:A:288:ASP:OD1	1:A:288:ASP:N	2.49	0.46
2:B:289:LYS:HG3	2:B:293:ARG:HD2	1.98	0.46
1:C:11:GLY:HA3	3:C:299:ATP:C4'	2.45	0.46
2:B:326:ASN:O	2:B:330:GLU:HG3	2.15	0.46
1:C:159:TYR:CD2	1:C:162:GLU:HG3	2.51	0.46
2:D:374:GLU:HG3	2:D:378:ARG:NH1	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:220:GLU:O	2:B:223:GLU:HB2	2.16	0.46
2:B:415:ASN:OD1	2:B:417:LYS:N	2.33	0.46
1:C:157:ARG:HD3	1:C:157:ARG:H	1.78	0.46
2:D:191:VAL:HG22	2:D:194:LYS:HZ1	1.81	0.46
2:B:210:MET:CE	2:B:250:ARG:HB2	2.23	0.45
1:C:197:VAL:HG11	1:C:255:LEU:HD13	1.98	0.45
2:D:346:PRO:O	2:D:349:LYS:HG2	2.16	0.45
2:D:417:LYS:HB3	2:D:417:LYS:HE3	1.58	0.45
1:A:60:HIS:HD2	1:A:62:ASN:N	2.01	0.45
1:A:71:HIS:NE2	2:B:296:HIS:ND1	2.64	0.45
1:A:152:PHE:CD2	2:B:173:ASN:ND2	2.84	0.45
1:C:1:MET:HE2	1:C:70:ILE:HD13	1.92	0.45
1:C:113:GLN:NE2	1:C:281:LEU:CD2	2.79	0.45
1:C:159:TYR:CD1	1:C:160:THR:N	2.85	0.45
2:D:388:LYS:HB3	2:D:389:PRO:HD3	1.97	0.45
1:A:13:GLY:N	1:A:16:GLY:O	2.50	0.45
1:A:32:LEU:CD2	1:A:79:VAL:HG22	2.47	0.45
2:B:201:LYS:HE2	2:B:202:LYS:CD	2.47	0.45
1:C:34:LYS:CE	1:C:75:LYS:NZ	2.79	0.45
1:C:121:HIS:C	1:C:122:ARG:HG3	2.42	0.45
1:C:159:TYR:HE1	2:D:271:TYR:OH	2.00	0.45
1:A:227:TRP:CD2	1:A:230:VAL:HG13	2.52	0.45
2:B:372:TRP:CZ3	2:B:376:LEU:HD13	2.51	0.45
1:C:41:THR:HG22	2:D:288:LYS:HD2	1.98	0.45
1:C:113:GLN:HE22	1:C:281:LEU:CD2	2.30	0.45
1:C:117:PHE:CE1	1:C:121:HIS:CE1	3.05	0.45
1:C:225:VAL:CG1	1:C:226:VAL:N	2.79	0.45
2:D:179:HIS:NE2	2:D:320:LEU:CD1	2.75	0.45
1:A:37:LEU:HD23	1:A:37:LEU:HA	1.69	0.45
1:A:84:HIS:HB3	1:A:298:LEU:HD22	1.97	0.45
1:A:84:HIS:ND1	1:A:136:ASN:HA	2.31	0.45
1:A:85:GLN:HE21	1:A:85:GLN:HB2	1.50	0.45
1:A:159:TYR:CD2	1:A:162:GLU:CG	2.99	0.45
1:C:275:ILE:HD13	1:C:280:ALA:CA	2.47	0.45
1:A:286:PHE:C	1:A:288:ASP:H	2.25	0.44
1:C:15:TYR:CE1	1:C:35:ILE:CD1	3.00	0.44
1:C:233:MET:HA	1:C:234:PRO:HD3	1.76	0.44
1:C:297:ARG:NH1	1:C:298:LEU:HD23	2.32	0.44
2:B:414:LYS:HB2	2:B:423:LEU:HD11	1.98	0.44
1:C:6:LYS:CD	1:C:6:LYS:H	2.30	0.44
1:C:226:VAL:O	1:C:226:VAL:HG22	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:TYR:HB3	4:A:391:HOH:O	2.18	0.44
2:B:277:GLU:O	2:B:281:ILE:HG12	2.16	0.44
2:B:327:CYS:HG	2:B:419:HIS:HE2	1.55	0.44
2:B:430:LEU:C	2:B:432:LEU:N	2.73	0.44
1:A:2:GLU:HG2	1:C:73:GLU:OE1	2.18	0.44
1:A:121:HIS:C	1:A:122:ARG:HG3	2.43	0.44
1:C:203:PHE:HB2	1:C:211:GLN:HE22	1.83	0.44
1:C:268:HIS:HB3	1:C:274:ARG:HA	2.00	0.44
2:D:189:MET:HE1	2:D:192:LYS:NZ	2.32	0.44
2:D:263:LEU:HD21	2:D:295:GLU:HG3	1.98	0.44
2:D:328:LYS:HB3	2:D:328:LYS:HE2	1.72	0.44
1:A:74:ASN:N	1:A:74:ASN:OD1	2.51	0.44
1:A:159:TYR:CE2	1:A:161:HIS:HB2	2.52	0.44
2:B:201:LYS:NZ	2:B:202:LYS:CE	2.80	0.44
1:C:11:GLY:HA3	3:C:299:ATP:H4'	2.00	0.44
1:C:158:THR:HG22	1:C:180:TYR:CE2	2.53	0.44
1:C:6:LYS:H	1:C:6:LYS:HD2	1.82	0.44
2:D:179:HIS:CE1	2:D:320:LEU:HD12	2.53	0.44
1:A:82:PHE:CD1	1:A:82:PHE:C	2.96	0.44
1:A:273:LYS:HA	1:A:273:LYS:HD2	1.79	0.44
2:B:338:GLU:OE2	2:B:412:LYS:NZ	2.43	0.44
1:C:225:VAL:HG12	1:C:226:VAL:N	2.33	0.44
1:A:49:ILE:HD12	2:B:266:LYS:HD2	2.00	0.44
1:A:72:THR:HG1	1:A:75:LYS:HG3	1.83	0.44
2:B:233:HIS:HE1	4:B:515:HOH:O	1.99	0.44
2:B:361:HIS:CE1	2:B:384:LEU:HD21	2.53	0.44
1:C:213:PHE:HE2	1:C:241:PRO:CD	2.30	0.44
1:C:55:LEU:HD23	1:C:55:LEU:HA	1.83	0.43
1:C:135:ILE:HG23	1:C:141:ILE:HD12	1.99	0.43
2:B:203:GLN:HG2	2:B:206:ILE:CG1	2.44	0.43
1:A:255:LEU:HD12	1:A:255:LEU:HA	1.80	0.43
2:B:200:MET:HE3	2:B:208:ASN:HD22	1.83	0.43
2:B:370:GLN:HG3	4:B:479:HOH:O	2.18	0.43
2:D:415:ASN:HD22	2:D:417:LYS:HB3	1.83	0.43
1:A:101:LEU:N	1:A:102:PRO:CD	2.82	0.43
1:A:138:GLU:C	1:A:293:VAL:HG12	2.43	0.43
2:B:387:LEU:O	2:B:391:LEU:HB2	2.19	0.43
1:C:105:LYS:HE2	1:C:285:PHE:O	2.18	0.43
1:C:135:ILE:HA	1:C:140:ALA:O	2.18	0.43
1:C:159:TYR:CD2	1:C:162:GLU:N	2.86	0.43
2:D:263:LEU:HD21	2:D:295:GLU:CG	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ASN:ND2	1:A:110:GLN:HB3	2.33	0.43
1:A:268:HIS:CE1	4:A:390:HOH:O	2.71	0.43
1:C:1:MET:HG3	1:C:77:TYR:CE1	2.54	0.43
1:C:55:LEU:HD13	1:C:66:LEU:HB2	2.01	0.43
1:C:262:LEU:HD21	1:C:286:PHE:CZ	2.53	0.43
1:A:212:LEU:HA	1:A:212:LEU:HD23	1.79	0.43
1:A:286:PHE:O	1:A:288:ASP:N	2.52	0.43
1:C:223:ASP:OD1	1:C:226:VAL:N	2.45	0.43
1:C:247:ASP:CB	1:C:250:LYS:HD3	2.47	0.43
2:D:194:LYS:HA	2:D:195:PRO:HD3	1.79	0.43
2:B:187:ARG:O	2:B:190:GLU:HG2	2.18	0.43
2:B:187:ARG:HD2	2:B:190:GLU:OE2	2.18	0.43
2:B:270:ILE:CG2	2:B:271:TYR:HD1	2.20	0.43
1:C:85:GLN:HE21	1:C:298:LEU:HB3	1.82	0.43
1:C:182:THR:N	2:D:173:ASN:O	2.40	0.43
1:A:159:TYR:HD2	1:A:162:GLU:HG3	1.83	0.43
1:A:178:LYS:HD3	1:A:179:TYR:CZ	2.54	0.43
1:C:237:LYS:HE3	1:C:240:PHE:CE1	2.54	0.43
1:A:95:ALA:O	1:A:96:LEU:HB3	2.19	0.43
1:A:152:PHE:C	1:A:154:VAL:H	2.25	0.43
2:B:229:ASN:HD22	2:B:334:MET:CE	2.32	0.43
2:B:287:THR:O	2:B:291:VAL:HG23	2.17	0.43
2:B:358:ALA:HB1	2:B:391:LEU:HD13	2.00	0.43
1:C:15:TYR:CD1	1:C:15:TYR:C	2.96	0.43
1:C:250:LYS:N	1:C:250:LYS:CD	2.74	0.43
1:C:252:VAL:HG21	1:C:263:LEU:HD23	2.01	0.43
2:D:175:VAL:HB	2:D:176:PRO:HD3	2.00	0.43
1:A:188:SER:O	1:A:192:ILE:HG13	2.19	0.43
2:B:184:THR:O	2:B:188:GLU:HG3	2.19	0.43
2:B:245:SER:O	2:B:246:MET:HG2	2.18	0.43
2:B:370:GLN:CG	2:B:371:SER:N	2.82	0.43
1:C:12:GLU:N	3:C:299:ATP:H5'2	2.34	0.43
1:A:157:ARG:NH1	1:A:179:TYR:HE1	2.17	0.42
1:A:193:PHE:O	1:A:197:VAL:HG23	2.19	0.42
1:C:15:TYR:C	1:C:15:TYR:HD1	2.27	0.42
2:D:368:THR:C	2:D:370:GLN:H	2.25	0.42
1:A:270:ASP:HB3	1:A:273:LYS:HB2	2.00	0.42
2:B:199:TYR:CD1	2:B:199:TYR:C	2.97	0.42
1:C:141:ILE:C	1:C:142:LYS:HG2	2.44	0.42
2:D:347:TYR:C	2:D:349:LYS:N	2.76	0.42
2:D:384:LEU:HD12	2:D:384:LEU:HA	1.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:159:TYR:CZ	2:D:271:TYR:HE1	2.37	0.42
2:D:239:ILE:HD11	2:D:257:GLY:HA2	2.00	0.42
1:A:105:LYS:HB2	1:A:105:LYS:HE3	1.76	0.42
1:A:181:SER:HA	2:B:173:ASN:N	2.34	0.42
1:A:237:LYS:HA	1:A:238:PRO:HD3	1.88	0.42
1:C:67:LEU:O	1:C:68:ASP:HB2	2.19	0.42
1:C:95:ALA:CA	1:C:199:ARG:HD2	2.48	0.42
1:C:141:ILE:O	1:C:142:LYS:HG2	2.19	0.42
1:C:223:ASP:C	1:C:225:VAL:N	2.77	0.42
2:D:376:LEU:HD23	2:D:376:LEU:HA	1.87	0.42
1:C:89:LYS:HB3	1:C:297:ARG:HH21	1.85	0.42
1:C:213:PHE:CD2	1:C:243:TRP:HH2	2.36	0.42
1:C:255:LEU:HD23	1:C:259:GLY:C	2.44	0.42
2:D:348:LEU:HA	2:D:348:LEU:HD12	1.71	0.42
2:D:412:LYS:HE2	2:D:413:TYR:HE1	1.82	0.42
1:C:297:ARG:NH1	1:C:298:LEU:CD2	2.82	0.42
2:D:191:VAL:O	2:D:191:VAL:HG13	2.19	0.42
2:D:343:ASP:HB3	2:D:345:ASP:CB	2.40	0.42
2:D:361:HIS:HE1	2:D:371:SER:HB3	1.83	0.42
2:B:270:ILE:HG22	2:B:271:TYR:N	2.33	0.42
2:B:201:LYS:CE	2:B:202:LYS:HG3	2.46	0.42
2:B:370:GLN:HG3	2:B:371:SER:N	2.34	0.42
2:D:347:TYR:C	2:D:349:LYS:H	2.27	0.42
1:A:175:LEU:HD23	1:A:212:LEU:HD11	2.02	0.42
1:C:55:LEU:CD1	1:C:80:PHE:CE1	2.99	0.42
1:C:134:LEU:HD12	1:C:144:ALA:CB	2.49	0.42
1:C:284:PRO:HA	1:C:287:GLN:HG3	2.02	0.42
1:A:182:THR:N	2:B:173:ASN:N	2.67	0.41
2:B:207:THR:CG2	2:B:210:MET:H	2.33	0.41
2:B:233:HIS:HD2	4:B:446:HOH:O	2.02	0.41
1:C:203:PHE:HB3	1:C:211:GLN:NE2	2.35	0.41
2:B:270:ILE:HD12	2:B:270:ILE:HA	1.74	0.41
1:C:275:ILE:CD1	1:C:280:ALA:HA	2.50	0.41
1:A:55:LEU:HD12	1:A:55:LEU:HA	1.82	0.41
1:C:41:THR:HG22	2:D:288:LYS:CD	2.51	0.41
2:B:315:LEU:HA	2:B:315:LEU:HD23	1.85	0.41
2:B:366:THR:HG22	2:B:424:LEU:HD21	2.03	0.41
1:A:34:LYS:HE2	1:A:34:LYS:HB3	1.88	0.41
1:C:138:GLU:O	1:C:138:GLU:HG3	2.19	0.41
2:D:323:GLN:HA	2:D:324:PRO:HA	1.63	0.41
2:B:415:ASN:CG	2:B:416:SER:N	2.79	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:175:VAL:O	2:D:176:PRO:C	2.63	0.41
2:B:178:TYR:O	2:B:182:ILE:HG13	2.21	0.41
2:B:214:LEU:HD11	2:B:257:GLY:HA3	2.03	0.41
2:B:372:TRP:HA	2:B:373:PRO:HD2	1.80	0.41
2:B:404:HIS:O	2:B:407:GLN:NE2	2.54	0.41
1:C:129:LYS:HE3	1:C:131:GLN:CG	2.50	0.41
1:C:253:PRO:N	1:C:254:PRO:CD	2.84	0.41
1:A:40:GLU:O	2:B:288:LYS:CD	2.69	0.41
1:A:87:LEU:O	1:A:91:MET:HB2	2.21	0.41
1:C:85:GLN:HE21	1:C:297:ARG:HH12	1.66	0.41
1:C:140:ALA:N	4:C:354:HOH:O	2.54	0.41
1:A:49:ILE:HD12	2:B:266:LYS:CD	2.51	0.41
1:A:152:PHE:C	1:A:154:VAL:N	2.78	0.41
1:C:174:LEU:C	1:C:176:GLY:H	2.29	0.41
1:C:261:SER:OG	1:C:283:HIS:NE2	2.52	0.41
2:D:358:ALA:HB1	2:D:391:LEU:CD1	2.51	0.41
1:A:295:HIS:O	1:A:297:ARG:NH2	2.54	0.40
1:C:119:HIS:CD2	1:C:182:THR:HB	2.56	0.40
1:C:261:SER:O	1:C:265:GLN:HG3	2.21	0.40
1:C:83:LEU:O	3:C:299:ATP:H2	2.03	0.40
1:C:268:HIS:CD2	1:C:268:HIS:C	3.00	0.40
2:D:394:LEU:HD12	2:D:394:LEU:HA	1.89	0.40
2:D:415:ASN:ND2	2:D:417:LYS:HB3	2.35	0.40
1:A:139:GLY:HA2	1:A:294:PRO:HD3	2.03	0.40
2:B:200:MET:SD	2:B:206:ILE:CD1	3.05	0.40
2:D:218:LEU:HD23	2:D:218:LEU:HA	1.90	0.40
2:D:365:TYR:HA	2:D:370:GLN:O	2.21	0.40
1:C:115:LEU:HD22	1:C:189:LEU:HD22	2.04	0.40
2:D:191:VAL:CG2	2:D:194:LYS:HZ1	2.35	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	296/298 (99%)	270 (91%)	20 (7%)	6 (2%)	6	5
1	C	296/298 (99%)	267 (90%)	22 (7%)	7 (2%)	4	4
2	B	258/260 (99%)	238 (92%)	15 (6%)	5 (2%)	6	5
2	D	258/260 (99%)	247 (96%)	8 (3%)	3 (1%)	10	12
All	All	1108/1116 (99%)	1022 (92%)	65 (6%)	21 (2%)	6	5

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	GLU
1	A	96	LEU
2	B	174	GLU
2	B	176	PRO
2	B	346	PRO
1	C	139	GLY
2	D	346	PRO
1	A	287	GLN
2	B	204	PRO
1	C	14	THR
1	A	40	GLU
1	A	155	PRO
1	C	97	THR
1	A	37	LEU
1	C	37	LEU
2	D	345	ASP
1	C	228	PRO
1	C	289	VAL
2	D	176	PRO
2	B	345	ASP
1	C	234	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	263/263 (100%)	223 (85%)	40 (15%)	3	3
1	C	263/263 (100%)	221 (84%)	42 (16%)	2	2
2	B	234/234 (100%)	206 (88%)	28 (12%)	5	5
2	D	234/234 (100%)	219 (94%)	15 (6%)	16	23
All	All	994/994 (100%)	869 (87%)	125 (13%)	4	5

All (125) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	6	LYS
1	A	9	LYS
1	A	30	VAL
1	A	37	LEU
1	A	39	THR
1	A	40	GLU
1	A	41	THR
1	A	47	THR
1	A	75	LYS
1	A	76	LEU
1	A	78	LEU
1	A	87	LEU
1	A	94	SER
1	A	97	THR
1	A	101	LEU
1	A	115	LEU
1	A	122	ARG
1	A	138	GLU
1	A	154	VAL
1	A	162	GLU
1	A	164	VAL
1	A	172	GLU
1	A	189	LEU
1	A	200	ARG
1	A	202	LEU
1	A	226	VAL
1	A	230	VAL
1	A	232	SER
1	A	239	SER
1	A	242	LYS
1	A	246	GLN
1	A	250	LYS

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Mol	Chain	Res	Type
1	A	255	LEU
1	A	257	GLU
1	A	264	SER
1	A	278	LYS
1	A	293	VAL
1	A	297	ARG
1	A	298	LEU
2	B	175	VAL
2	B	182	ILE
2	B	194	LYS
2	B	209	SER
2	B	210	MET
2	B	232	LEU
2	B	237	ASN
2	B	247	SER
2	B	249	LEU
2	B	250	ARG
2	B	270	ILE
2	B	285	THR
2	B	288	LYS
2	B	289	LYS
2	B	292	LEU
2	B	293	ARG
2	B	323	GLN
2	B	345	ASP
2	B	348	LEU
2	B	349	LYS
2	B	368	THR
2	B	384	LEU
2	B	386	SER
2	B	391	LEU
2	B	392	MET
2	B	407	GLN
2	B	416	SER
2	B	423	LEU
1	C	5	GLN
1	C	6	LYS
1	C	10	ILE
1	C	15	TYR
1	C	32	LEU
1	C	36	ARG
1	C	58	LEU

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Mol	Chain	Res	Type
1	C	74	ASN
1	C	75	LYS
1	C	76	LEU
1	C	85	GLN
1	C	87	LEU
1	C	89	LYS
1	C	91	MET
1	C	101	LEU
1	C	113	GLN
1	C	122	ARG
1	C	138	GLU
1	C	141	ILE
1	C	145	ASP
1	C	148	LEU
1	C	152	PHE
1	C	154	VAL
1	C	163	VAL
1	C	164	VAL
1	C	181	SER
1	C	186	ILE
1	C	189	LEU
1	C	196	MET
1	C	202	LEU
1	C	224	GLU
1	C	225	VAL
1	C	231	THR
1	C	239	SER
1	C	242	LYS
1	C	250	LYS
1	C	252	VAL
1	C	255	LEU
1	C	264	SER
1	C	275	ILE
1	C	288	ASP
1	C	298	LEU
2	D	173	ASN
2	D	177	ASP
2	D	191	VAL
2	D	196	LYS
2	D	232	LEU
2	D	247	SER
2	D	274	GLU

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Mol	Chain	Res	Type
2	D	275	VAL
2	D	283	ASP
2	D	288	LYS
2	D	292	LEU
2	D	348	LEU
2	D	371	SER
2	D	391	LEU
2	D	417	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	ASN
1	A	60	HIS
1	A	71	HIS
1	A	85	GLN
1	A	268	HIS
2	B	173	ASN
2	B	179	HIS
2	B	208	ASN
2	B	313	GLN
2	B	317	GLN
2	B	323	GLN
2	B	404	HIS
2	B	431	ASN
1	C	60	HIS
1	C	71	HIS
1	C	84	HIS
1	C	85	GLN
1	C	113	GLN
1	C	119	HIS
1	C	161	HIS
1	C	211	GLN
2	D	173	ASN
2	D	179	HIS
2	D	183	HIS
2	D	208	ASN
2	D	229	ASN
2	D	254	GLN
2	D	296	HIS
2	D	317	GLN
2	D	321	HIS

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Mol	Chain	Res	Type
2	D	404	HIS
2	D	415	ASN
2	D	431	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ATP	A	299	-	32,33,33	1.88	3 (9%)	48,52,52	2.50	10 (20%)
3	ATP	C	299	-	32,33,33	1.67	3 (9%)	48,52,52	2.03	10 (20%)

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	ATP	A	299	-	-	9/22/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ATP	C	299	-	-	7/22/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	299	ATP	PB-O3B	7.05	1.67	1.59
3	C	299	ATP	PB-O3B	5.67	1.65	1.59
3	A	299	ATP	PA-O3A	5.43	1.65	1.59
3	C	299	ATP	PA-O3A	5.02	1.64	1.59
3	A	299	ATP	PB-O3A	4.77	1.64	1.59
3	C	299	ATP	PB-O3A	4.32	1.64	1.59

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	299	ATP	O2A-PA-O3A	-10.36	79.28	107.27
3	A	299	ATP	O2A-PA-O3A	-9.88	80.56	107.27
3	A	299	ATP	O3B-PB-O1B	-6.40	91.44	110.70
3	A	299	ATP	O3A-PA-O1A	6.15	129.19	110.70
3	A	299	ATP	O2B-PB-O3B	5.24	121.43	107.27
3	A	299	ATP	O4'-C1'-N9	4.69	117.10	108.09
3	A	299	ATP	O4'-C4'-C5'	3.65	121.02	109.33
3	A	299	ATP	C2'-C1'-N9	-3.42	104.81	113.30
3	C	299	ATP	O3G-PG-O1G	3.11	122.96	110.83
3	C	299	ATP	O4'-C1'-N9	2.91	113.67	108.09
3	C	299	ATP	O3A-PA-O1A	2.81	119.17	110.70
3	A	299	ATP	O2B-PB-O1B	2.66	124.84	112.44
3	A	299	ATP	O5'-C5'-C4'	2.63	117.96	108.99
3	C	299	ATP	O2B-PB-O1B	2.26	122.98	112.44
3	C	299	ATP	C4-N9-C1'	2.25	131.90	126.63
3	C	299	ATP	O2A-PA-O1A	2.24	122.87	112.44
3	C	299	ATP	O3A-PB-O1B	-2.12	104.32	110.70
3	C	299	ATP	O4'-C4'-C3'	-2.12	100.95	105.15
3	A	299	ATP	C4'-O4'-C1'	-2.08	104.88	109.47
3	C	299	ATP	O3B-PB-O1B	-2.06	104.51	110.70

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	299	ATP	PB-O3B-PG-O2G

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Mol	Chain	Res	Type	Atoms
3	A	299	ATP	PB-O3B-PG-O3G
3	A	299	ATP	C5'-O5'-PA-O2A
3	A	299	ATP	C5'-O5'-PA-O3A
3	C	299	ATP	PB-O3B-PG-O3G
3	C	299	ATP	C5'-O5'-PA-O1A
3	C	299	ATP	C5'-O5'-PA-O3A
3	C	299	ATP	O4'-C4'-C5'-O5'
3	A	299	ATP	O4'-C4'-C5'-O5'
3	C	299	ATP	C3'-C4'-C5'-O5'
3	A	299	ATP	C3'-C4'-C5'-O5'
3	A	299	ATP	C4'-C5'-O5'-PA
3	A	299	ATP	C5'-O5'-PA-O1A
3	C	299	ATP	C4'-C5'-O5'-PA
3	C	299	ATP	PB-O3B-PG-O2G
3	A	299	ATP	PG-O3B-PB-O2B

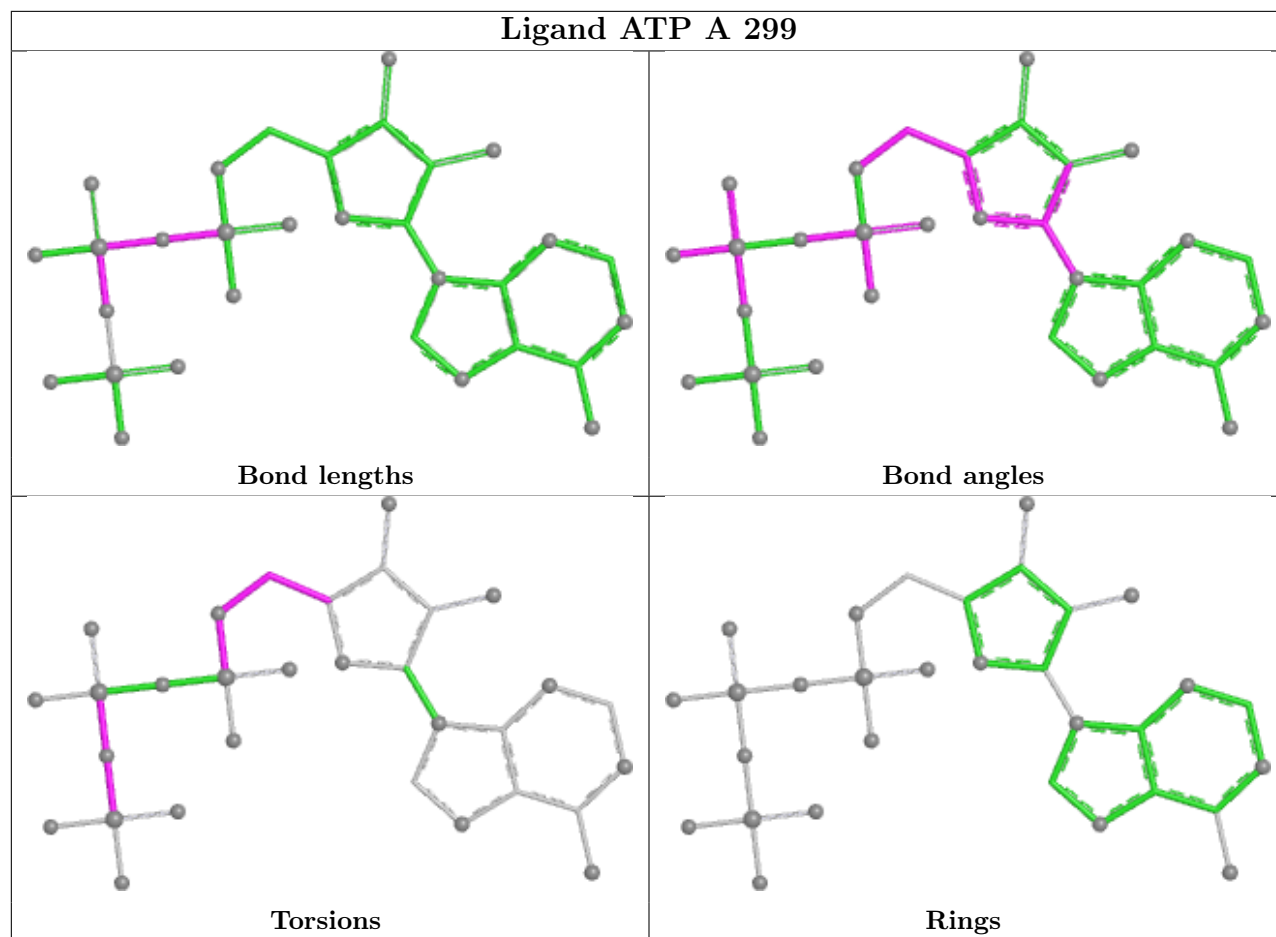
There are no ring outliers.

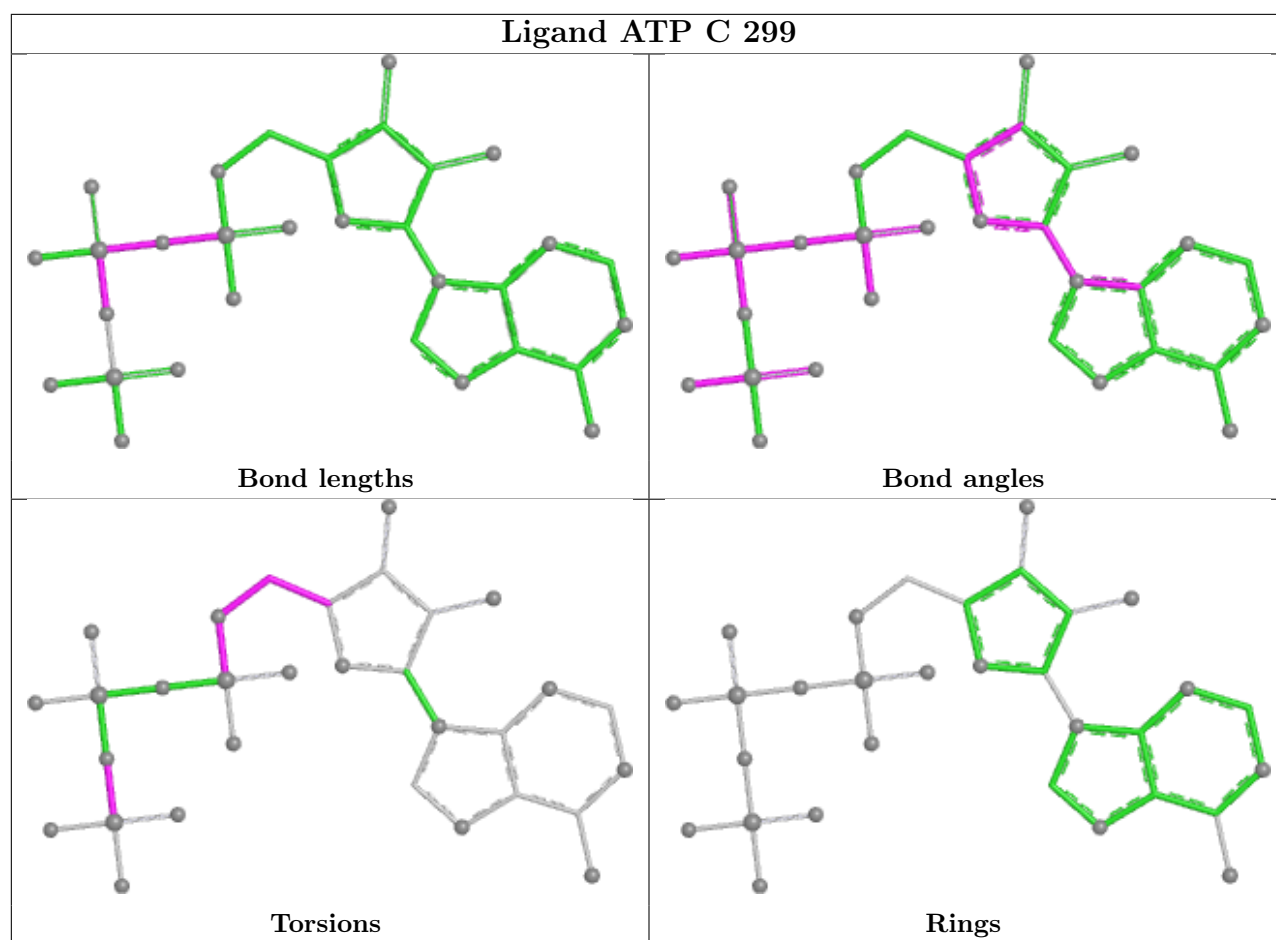
2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	299	ATP	2	0
3	C	299	ATP	7	0

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

Ligand ATP A 299





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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.