



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:30 PM GMT

PDB ID : 1FVV
Title : THE STRUCTURE OF CDK2/CYCLIN A IN COMPLEX WITH AN OXINDOLE INHIBITOR
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Deposited on : 2000-09-20
Resolution : 2.80 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

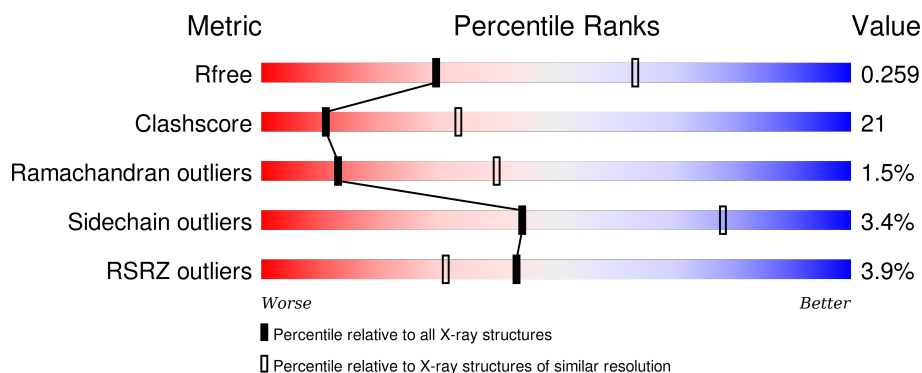
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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>3%</div> <div>65%</div> <div>32%</div> <div>.</div> </div>
1	C	298	<div> <div>5%</div> <div>57%</div> <div>39%</div> <div>.</div> </div>
2	B	260	<div> <div>3%</div> <div>55%</div> <div>41%</div> <div>.</div> </div>
2	D	260	<div> <div>5%</div> <div>65%</div> <div>31%</div> <div>.</div> </div>

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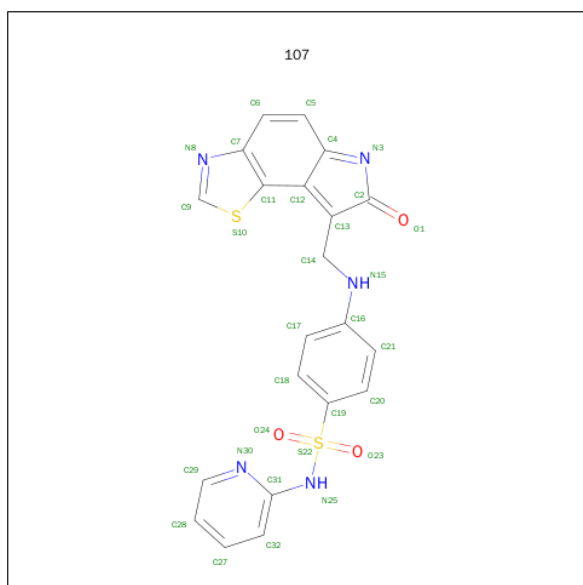
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 2 is a protein called CYCLIN A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	298	Total 2377	C 1546	N 405	O 418	S 8	0	0	0
1	C	298	Total 2381	C 1548	N 405	O 420	S 8	0	0	0

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

- Molecule 3 is 4-[(7-OXO-7H-THIAZOLO[5,4-E]INDOL-8-YLMETHYL)-AMINO]-N-PYRIDIN-2-YL-BENZENESULFONAMIDE (three-letter code: 107) (formula: $C_{21}H_{15}N_5O_3S_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			31	21	5	3	2		
3	C	1	Total	C	N	O	S	0	0
			31	21	5	3	2		

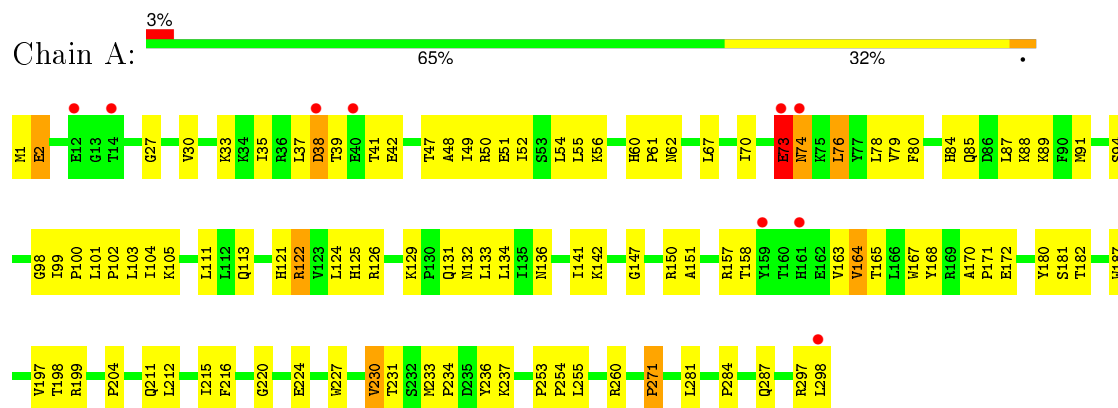
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total	O	0	0
			58	58		
4	B	14	Total	O	0	0
			14	14		
4	C	18	Total	O	0	0
			18	18		
4	D	39	Total	O	0	0
			39	39		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

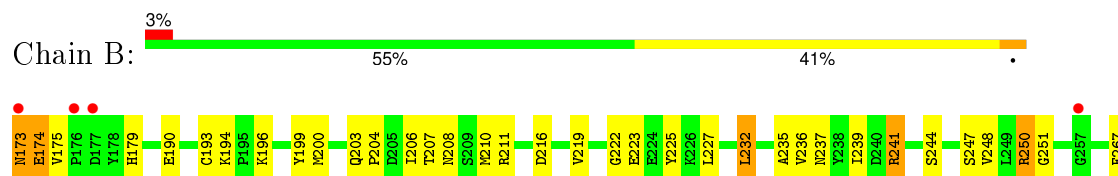
• Molecule 1: CYCLIN-DEPENDENT KINASE 2

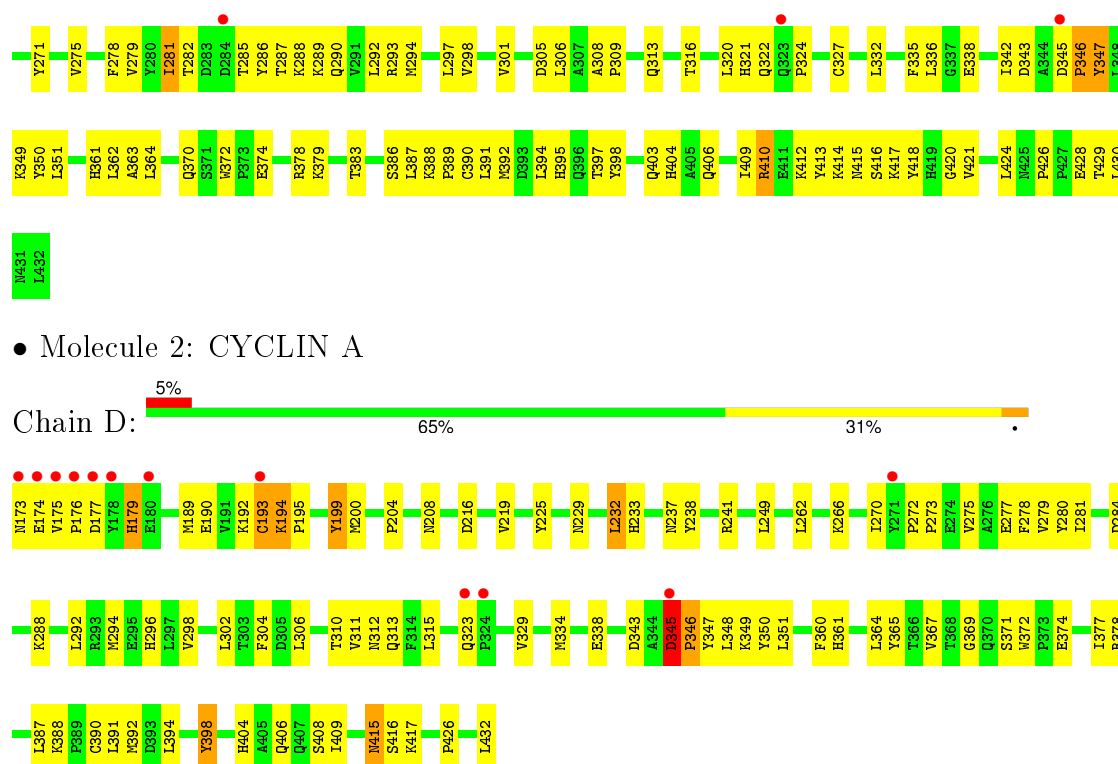


• Molecule 1: CYCLIN-DEPENDENT KINASE 2



• Molecule 2: CYCLIN A





• Molecule 2: CYCLIN A

4 Data and refinement statistics

Property	Value	Source
Space group	P 6 ₂ 2 2	Depositor
Cell constants a, b, c, α , β , γ	184.95Å 184.95Å 212.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.80 19.93 – 2.79	Depositor EDS
% Data completeness (in resolution range)	95.9 (20.00-2.80) 95.7 (19.93-2.79)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.03 (at 2.79Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.260 , 0.260 0.216 , 0.259	Depositor DCC
R_{free} test set	5137 reflections (10.09%)	DCC
Wilson B-factor (Å ²)	58.6	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 53318 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9151	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
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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	0.39	0/2438	0.63	0/3310
1	C	0.36	0/2442	0.62	1/3315 (0.0%)
2	B	0.40	1/2151 (0.0%)	0.60	0/2920
2	D	0.42	1/2151 (0.0%)	0.63	0/2920
All	All	0.39	2/9182 (0.0%)	0.62	1/12465 (0.0%)

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
2	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	193	CYS	CB-SG	6.68	1.93	1.82
2	D	193	CYS	CB-SG	5.27	1.91	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	154	VAL	N-CA-C	-6.12	94.48	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	347	TYR	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2377	0	2424	96	0
1	C	2381	0	2428	118	0
2	B	2101	0	2119	107	0
2	D	2101	0	2119	74	0
3	A	31	0	15	2	0
3	C	31	0	15	0	0
4	A	58	0	0	3	0
4	B	14	0	0	1	0
4	C	18	0	0	2	0
4	D	39	0	0	3	0
All	All	9151	0	9120	381	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 21.

All (381) 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:131:GLN:H	1:C:131:GLN:NE2	1.55	1.03
2:B:203:GLN:HG2	2:B:206:ILE:HG12	1.43	0.99
2:D:404:HIS:HD2	2:D:406:GLN:H	1.10	0.98
2:B:250:ARG:HB3	2:B:250:ARG:HH11	1.25	0.97
2:D:415:ASN:ND2	2:D:417:LYS:H	1.69	0.91
1:A:227:TRP:O	1:A:230:VAL:HG23	1.69	0.90
2:B:287:THR:HG22	2:B:289:LYS:H	1.39	0.88
2:D:345:ASP:HB2	2:D:346:PRO:HD3	1.56	0.86
2:D:415:ASN:HD22	2:D:417:LYS:H	1.24	0.86
2:B:345:ASP:HB3	2:B:346:PRO:HD3	1.59	0.84
2:D:219:VAL:HG22	2:D:232:LEU:HD11	1.56	0.84
2:B:207:THR:HG22	2:B:210:MET:HG3	1.58	0.83
1:C:131:GLN:HE21	1:C:131:GLN:H	1.22	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:289:LYS:O	2:B:293:ARG:HG3	1.81	0.81
2:D:415:ASN:HD22	2:D:416:SER:N	1.80	0.80
2:D:200:MET:HG2	2:D:208:ASN:ND2	1.97	0.80
1:A:61:PRO:O	1:A:142:LYS:HE2	1.82	0.79
1:A:172:GLU:HG2	1:A:271:PRO:HG3	1.65	0.79
2:D:404:HIS:CD2	2:D:406:GLN:H	1.98	0.78
1:C:2:GLU:H	1:C:2:GLU:CD	1.87	0.78
1:A:163:VAL:HB	4:A:555:HOH:O	1.83	0.77
2:B:417:LYS:HE3	2:B:418:TYR:CE1	2.20	0.77
2:D:345:ASP:CB	2:D:346:PRO:HD3	2.15	0.77
2:D:174:GLU:HG3	2:D:176:PRO:HD2	1.66	0.77
2:B:404:HIS:HE1	2:B:406:GLN:HE21	1.32	0.77
1:C:60:HIS:HD2	1:C:62:ASN:H	1.30	0.77
2:B:237:ASN:HD21	2:B:241:ARG:HE	1.31	0.76
2:B:250:ARG:HH11	2:B:250:ARG:CB	1.98	0.76
2:B:237:ASN:ND2	2:B:241:ARG:HE	1.84	0.75
2:B:428:GLU:HG3	2:B:429:THR:HG23	1.69	0.75
1:C:5:GLN:HB3	1:C:22:ARG:HG2	1.68	0.74
2:B:336:LEU:HD13	2:B:362:LEU:HD23	1.67	0.74
2:B:250:ARG:HB3	2:B:250:ARG:NH1	2.03	0.73
1:C:61:PRO:O	1:C:142:LYS:HE2	1.89	0.72
1:C:201:ALA:HB3	1:C:204:PRO:HG3	1.70	0.71
2:D:391:LEU:HD23	2:D:432:LEU:HD11	1.71	0.71
2:B:388:LYS:HB3	2:B:389:PRO:HD3	1.72	0.71
1:A:33:LYS:HB3	1:A:78:LEU:HB2	1.72	0.71
1:C:165:THR:HG22	4:C:607:HOH:O	1.89	0.70
2:D:175:VAL:HA	2:D:179:HIS:CD2	2.26	0.70
2:B:361:HIS:HD2	2:B:372:TRP:H	1.37	0.70
1:C:150:ARG:HD2	4:C:610:HOH:O	1.92	0.69
1:A:47:THR:HG22	1:A:147:GLY:O	1.92	0.69
2:B:410:ARG:NH1	2:B:410:ARG:HB3	2.07	0.69
1:A:30:VAL:HG21	1:A:79:VAL:HG12	1.74	0.69
2:B:203:GLN:HG2	2:B:206:ILE:CG1	2.21	0.69
2:B:332:LEU:HD23	2:B:363:ALA:HA	1.73	0.68
2:B:210:MET:HE1	2:B:250:ARG:HB2	1.74	0.68
1:A:74:ASN:H	1:A:74:ASN:HD22	1.41	0.68
2:D:262:LEU:HD11	2:D:266:LYS:HE3	1.73	0.68
1:C:136:ASN:HD21	1:C:138:GLU:HB2	1.58	0.68
1:C:136:ASN:ND2	1:C:138:GLU:HB2	2.09	0.68
1:A:51:GLU:O	1:A:55:LEU:HB2	1.94	0.68
1:A:52:ILE:O	1:A:56:LYS:HG3	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:241:PRO:HB2	1:C:243:TRP:CZ3	2.29	0.67
2:D:345:ASP:HB2	2:D:346:PRO:CD	2.25	0.67
2:B:207:THR:CG2	2:B:210:MET:HG3	2.25	0.67
2:B:287:THR:HB	2:B:290:GLN:HG3	1.76	0.66
1:C:38:ASP:HB3	1:C:42:GLU:HB3	1.77	0.66
1:A:182:THR:HG23	2:B:173:ASN:N	2.09	0.66
1:A:165:THR:HG21	1:A:168:TYR:HD1	1.59	0.66
1:C:52:ILE:HD11	1:C:78:LEU:HD21	1.78	0.66
1:A:60:HIS:CD2	1:A:62:ASN:H	2.14	0.66
1:C:74:ASN:HD22	1:C:75:LYS:CD	2.09	0.66
1:C:1:MET:HG2	1:C:70:ILE:HG21	1.78	0.66
1:C:60:HIS:CD2	1:C:62:ASN:H	2.14	0.64
1:C:252:VAL:HG21	1:C:255:LEU:HD12	1.79	0.64
2:B:404:HIS:CE1	2:B:406:GLN:HE21	2.16	0.63
1:C:200:ARG:HD2	1:C:201:ALA:H	1.63	0.63
1:C:40:GLU:O	2:D:288:LYS:HD3	1.98	0.63
2:D:204:PRO:HG2	4:D:49:HOH:O	1.98	0.63
1:C:74:ASN:HD22	1:C:75:LYS:HD2	1.65	0.62
2:D:374:GLU:HG3	2:D:378:ARG:NH1	2.15	0.62
1:A:60:HIS:HD2	1:A:62:ASN:H	1.47	0.61
1:C:34:LYS:HZ1	1:C:36:ARG:NH1	1.98	0.61
1:A:74:ASN:ND2	1:A:74:ASN:H	1.97	0.61
1:A:30:VAL:HG21	1:A:79:VAL:CG1	2.30	0.61
2:D:277:GLU:HA	2:D:280:TYR:CD1	2.36	0.61
1:C:131:GLN:HE21	1:C:131:GLN:N	1.95	0.61
1:A:231:THR:HA	1:A:236:TYR:CD2	2.35	0.61
1:C:126:ARG:HH12	1:C:159:TYR:HE1	1.48	0.61
2:D:361:HIS:CD2	2:D:391:LEU:HD21	2.35	0.60
1:C:203:PHE:CE2	1:C:215:ILE:HA	2.35	0.60
2:D:346:PRO:O	2:D:349:LYS:HG2	2.01	0.60
1:A:38:ASP:OD1	1:A:39:THR:HG23	2.01	0.60
2:B:282:THR:O	2:B:285:THR:HG23	2.01	0.60
2:D:415:ASN:C	2:D:415:ASN:HD22	2.03	0.60
2:B:374:GLU:HG3	2:B:378:ARG:NH2	2.17	0.60
1:A:164:VAL:HG12	1:A:165:THR:N	2.17	0.59
2:B:223:GLU:OE2	2:B:412:LYS:HE3	2.02	0.59
1:C:34:LYS:HE2	1:C:77:TYR:OH	2.02	0.59
1:C:36:ARG:O	1:C:37:LEU:HB2	2.02	0.59
2:B:219:VAL:HG22	2:B:232:LEU:HD11	1.85	0.59
2:B:225:TYR:HE2	2:B:281:ILE:HG21	1.68	0.58
1:A:85:GLN:NE2	1:A:89:LYS:HD3	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:HIS:O	1:A:122:ARG:HG3	2.03	0.58
2:D:229:ASN:HD22	2:D:334:MET:CE	2.16	0.58
1:C:224:GLU:OE2	1:C:231:THR:HB	2.03	0.58
1:C:74:ASN:ND2	1:C:75:LYS:HD2	2.19	0.57
2:D:345:ASP:CB	2:D:346:PRO:CD	2.83	0.57
2:B:361:HIS:HD2	2:B:372:TRP:N	2.02	0.57
1:C:101:LEU:N	1:C:102:PRO:HD2	2.19	0.57
1:A:94:SER:O	1:A:98:GLY:N	2.34	0.57
2:B:428:GLU:CG	2:B:429:THR:HG23	2.33	0.57
1:A:52:ILE:HD11	1:A:78:LEU:HD21	1.86	0.57
1:C:94:SER:HA	1:C:97:THR:HG23	1.85	0.56
2:D:388:LYS:O	2:D:392:MET:HG2	2.05	0.56
2:D:329:VAL:HG22	2:D:367:VAL:HB	1.88	0.56
2:D:312:ASN:OD1	2:D:334:MET:HE1	2.06	0.56
1:C:253:PRO:HB2	1:C:254:PRO:HD3	1.86	0.56
2:B:271:TYR:CD1	2:B:271:TYR:O	2.59	0.56
2:B:410:ARG:HH11	2:B:410:ARG:HB3	1.70	0.55
2:B:335:PHE:HB2	2:B:413:TYR:CD2	2.41	0.55
2:D:189:MET:CE	2:D:192:LYS:HD2	2.36	0.55
2:B:398:TYR:CE2	2:B:426:PRO:HB3	2.41	0.55
1:C:84:HIS:CD2	1:C:137:THR:HG23	2.42	0.55
2:B:173:ASN:O	2:B:175:VAL:HG22	2.07	0.55
1:A:74:ASN:N	1:A:74:ASN:HD22	1.99	0.55
2:D:278:PHE:O	2:D:281:ILE:HG12	2.07	0.55
1:C:202:LEU:HD13	1:C:203:PHE:CE1	2.42	0.55
1:C:33:LYS:HB3	1:C:78:LEU:HB2	1.88	0.54
2:D:225:TYR:HE1	2:D:277:GLU:HB3	1.72	0.54
2:B:410:ARG:HH11	2:B:410:ARG:CB	2.20	0.54
1:C:131:GLN:N	1:C:131:GLN:NE2	2.40	0.54
1:A:50:ARG:NE	1:A:150:ARG:NH1	2.55	0.54
2:D:190:GLU:HG3	2:D:351:LEU:HD22	1.89	0.54
1:A:129:LYS:HE2	1:A:132:ASN:ND2	2.23	0.54
2:B:222:GLY:HA2	2:B:227:LEU:HD12	1.90	0.54
2:D:350:TYR:CE2	2:D:390:CYS:HB2	2.43	0.54
1:C:2:GLU:OE2	1:C:2:GLU:N	2.41	0.54
2:D:237:ASN:OD1	2:D:241:ARG:HD3	2.08	0.54
1:A:87:LEU:O	1:A:91:MET:HG3	2.07	0.54
1:C:135:ILE:HA	1:C:140:ALA:O	2.07	0.54
1:C:111:LEU:HD21	1:C:141:ILE:HD13	1.89	0.54
1:C:60:HIS:CG	1:C:61:PRO:HD2	2.43	0.53
1:C:245:ARG:NH2	1:C:248:PHE:CE1	2.76	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:PRO:O	1:A:104:ILE:HG13	2.08	0.53
2:B:414:LYS:HA	2:B:420:GLY:HA2	1.89	0.53
2:B:343:ASP:O	2:B:346:PRO:HD2	2.09	0.53
1:C:94:SER:HA	1:C:97:THR:CG2	2.38	0.53
2:B:203:GLN:CG	2:B:206:ILE:HG12	2.27	0.53
1:C:167:TRP:CD1	1:C:204:PRO:HA	2.44	0.53
2:D:174:GLU:HG3	2:D:176:PRO:CD	2.36	0.53
1:C:34:LYS:HZ1	1:C:75:LYS:NZ	2.06	0.53
2:D:194:LYS:HG2	2:D:195:PRO:N	2.24	0.53
1:A:298:LEU:HD23	1:A:298:LEU:C	2.29	0.52
2:D:338:GLU:HB3	2:D:409:ILE:HD13	1.92	0.52
1:A:157:ARG:HG3	4:A:556:HOH:O	2.08	0.52
2:B:235:ALA:O	2:B:239:ILE:HG13	2.10	0.52
2:D:199:TYR:CE2	2:D:348:LEU:HD21	2.44	0.52
2:B:313:GLN:O	2:B:316:THR:HB	2.10	0.52
2:D:387:LEU:O	2:D:391:LEU:HB2	2.10	0.52
1:C:84:HIS:HD2	1:C:136:ASN:HA	1.73	0.52
2:B:175:VAL:HG23	2:B:175:VAL:O	2.10	0.52
2:D:398:TYR:CE1	2:D:426:PRO:HB3	2.45	0.52
2:B:216:ASP:HB2	2:B:406:GLN:HG2	1.92	0.52
2:B:196:LYS:HB2	2:B:199:TYR:HB3	1.91	0.52
2:B:374:GLU:HG3	2:B:378:ARG:HH22	1.74	0.52
1:C:152:PHE:C	1:C:154:VAL:H	2.13	0.52
1:A:129:LYS:HE3	1:A:131:GLN:HG2	1.92	0.51
1:C:63:ILE:O	1:C:64:VAL:C	2.47	0.51
2:B:210:MET:CE	2:B:250:ARG:HB2	2.40	0.51
2:B:350:TYR:CZ	2:B:390:CYS:HA	2.45	0.51
1:C:55:LEU:C	1:C:57:GLU:H	2.14	0.51
1:C:223:ASP:H	1:C:226:VAL:CG1	2.23	0.51
1:C:223:ASP:OD1	1:C:226:VAL:HG12	2.10	0.51
1:C:248:PHE:HB3	1:C:260:ARG:HD2	1.92	0.51
1:A:85:GLN:HA	3:A:501:107:C20	2.40	0.51
1:A:41:THR:HB	2:B:288:LYS:NZ	2.26	0.51
1:C:34:LYS:HZ1	1:C:75:LYS:HZ3	1.59	0.51
1:A:121:HIS:C	1:A:122:ARG:HG3	2.31	0.51
2:D:174:GLU:CG	2:D:176:PRO:HD2	2.38	0.50
1:C:34:LYS:NZ	1:C:75:LYS:NZ	2.58	0.50
2:B:297:LEU:O	2:B:301:VAL:HG23	2.10	0.50
1:A:27:GLY:HA3	2:D:249:LEU:HD12	1.94	0.50
1:C:127:ASP:HB2	1:C:148:LEU:HD12	1.92	0.50
1:C:57:GLU:OE1	1:C:122:ARG:NH2	2.38	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:346:PRO:HG2	2:D:347:TYR:CD1	2.46	0.50
1:C:34:LYS:NZ	1:C:75:LYS:HZ3	2.10	0.50
1:C:39:THR:HG22	1:C:39:THR:O	2.10	0.50
1:A:227:TRP:HB3	1:A:230:VAL:CG2	2.42	0.50
1:C:201:ALA:CB	1:C:204:PRO:HG3	2.42	0.50
2:B:200:MET:HG2	2:B:208:ASN:ND2	2.26	0.50
2:D:311:VAL:O	2:D:315:LEU:HB2	2.12	0.49
2:B:388:LYS:HE3	2:B:392:MET:HE3	1.94	0.49
1:C:109:PHE:HB2	1:C:286:PHE:CE1	2.47	0.49
1:A:60:HIS:CG	1:A:61:PRO:HD2	2.48	0.49
1:A:2:GLU:H	1:A:2:GLU:CD	2.14	0.49
2:B:322:GLN:HB3	2:B:324:PRO:O	2.12	0.49
1:A:227:TRP:CD2	1:A:230:VAL:HG22	2.47	0.49
1:C:60:HIS:CD2	1:C:61:PRO:HD2	2.48	0.49
1:A:216:PHE:O	1:A:220:GLY:N	2.44	0.49
1:A:297:ARG:O	1:A:298:LEU:C	2.50	0.49
2:B:203:GLN:OE1	2:B:248:VAL:N	2.43	0.49
2:B:190:GLU:O	2:B:194:LYS:HB3	2.12	0.49
1:A:284:PRO:HA	1:A:287:GLN:HG3	1.94	0.49
2:B:294:MET:O	2:B:298:VAL:HG23	2.12	0.49
2:B:346:PRO:O	2:B:349:LYS:HG2	2.13	0.48
1:C:134:LEU:O	1:C:141:ILE:HA	2.12	0.48
1:A:231:THR:HA	1:A:236:TYR:CE2	2.48	0.48
2:B:278:PHE:O	2:B:282:THR:HG23	2.13	0.48
1:C:87:LEU:O	1:C:91:MET:HG3	2.13	0.48
2:D:238:TYR:OH	2:D:306:LEU:HB3	2.13	0.48
1:A:88:LYS:HD2	1:A:131:GLN:HB3	1.95	0.48
1:C:99:ILE:HG22	1:C:104:ILE:HG13	1.96	0.48
2:B:237:ASN:O	2:B:241:ARG:HB2	2.14	0.48
2:B:388:LYS:O	2:B:392:MET:HG2	2.13	0.48
2:B:383:THR:N	2:B:386:SER:OG	2.46	0.48
1:C:165:THR:HG23	1:C:165:THR:O	2.13	0.48
1:C:152:PHE:CD1	1:C:152:PHE:N	2.82	0.48
2:D:189:MET:HE1	2:D:192:LYS:HD2	1.96	0.47
1:C:265:GLN:HB3	1:C:275:ILE:HB	1.95	0.47
1:C:188:SER:O	1:C:192:ILE:HG13	2.14	0.47
1:C:111:LEU:HD21	1:C:141:ILE:CD1	2.43	0.47
1:C:14:THR:HG23	1:C:15:TYR:N	2.29	0.47
1:A:181:SER:HB3	2:B:174:GLU:HG3	1.95	0.47
2:B:237:ASN:HD21	2:B:241:ARG:NE	2.07	0.47
1:C:105:LYS:HE2	1:C:285:PHE:CZ	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:THR:HA	1:A:180:TYR:CE2	2.50	0.47
1:A:111:LEU:CD2	1:A:133:LEU:HD22	2.45	0.47
2:D:371:SER:O	2:D:372:TRP:C	2.53	0.47
2:D:374:GLU:HA	2:D:377:ILE:HD12	1.97	0.47
1:C:84:HIS:CD2	1:C:137:THR:H	2.32	0.47
2:D:374:GLU:HG3	2:D:378:ARG:HH12	1.78	0.47
1:A:50:ARG:HD3	1:A:150:ARG:NH1	2.30	0.47
1:C:122:ARG:O	1:C:122:ARG:HD2	2.15	0.47
2:D:275:VAL:O	2:D:279:VAL:HG23	2.15	0.46
2:B:179:HIS:HD2	2:B:320:LEU:HD13	1.80	0.46
1:C:75:LYS:N	1:C:75:LYS:HD2	2.30	0.46
1:C:227:TRP:CE3	1:C:269:TYR:HB3	2.50	0.46
1:A:197:VAL:HG11	1:A:255:LEU:HG	1.97	0.46
1:A:30:VAL:HG23	1:A:67:LEU:HD12	1.97	0.46
1:C:70:ILE:O	1:C:70:ILE:HG22	2.14	0.46
2:B:275:VAL:HG21	2:B:292:LEU:HD11	1.98	0.46
2:B:404:HIS:CE1	2:B:406:GLN:HB2	2.51	0.46
1:C:34:LYS:NZ	1:C:36:ARG:NH1	2.62	0.46
1:C:100:PRO:C	1:C:102:PRO:HD2	2.36	0.46
1:C:7:VAL:HB	1:C:20:LYS:O	2.16	0.46
1:A:35:ILE:HD12	1:A:76:LEU:HD13	1.98	0.46
2:B:398:TYR:CD2	2:B:426:PRO:HB3	2.50	0.46
1:A:54:LEU:HD21	1:A:151:ALA:HB2	1.98	0.46
2:D:241:ARG:CZ	2:D:302:LEU:HD22	2.46	0.45
1:A:84:HIS:CD2	1:A:136:ASN:HA	2.51	0.45
2:B:287:THR:HG22	2:B:289:LYS:N	2.19	0.45
1:A:111:LEU:HD22	1:A:133:LEU:HD22	1.97	0.45
2:D:233:HIS:HD2	4:D:123:HOH:O	1.99	0.45
2:D:189:MET:HE2	2:D:192:LYS:HD2	1.97	0.45
2:B:241:ARG:NH2	2:B:305:ASP:O	2.45	0.45
1:C:200:ARG:HD2	1:C:201:ALA:N	2.30	0.45
1:A:181:SER:CB	2:B:174:GLU:HG3	2.47	0.45
2:B:203:GLN:HA	2:B:204:PRO:HD2	1.84	0.45
1:C:165:THR:HG23	1:C:168:TYR:HD1	1.81	0.45
1:A:1:MET:CE	1:A:70:ILE:HD12	2.47	0.45
2:B:216:ASP:O	2:B:219:VAL:HB	2.16	0.45
1:A:37:LEU:HD22	1:A:76:LEU:HB2	1.99	0.45
2:B:279:VAL:HG21	2:B:288:LYS:HA	1.99	0.45
1:A:35:ILE:HD13	1:A:48:ALA:HB2	1.99	0.45
1:C:113:GLN:HG3	1:C:281:LEU:HD21	1.98	0.45
1:A:73:GLU:HB2	1:A:74:ASN:H	1.56	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:LYS:HD2	1:C:131:GLN:OE1	2.17	0.44
2:B:345:ASP:HB3	2:B:346:PRO:CD	2.40	0.44
2:B:232:LEU:O	2:B:236:VAL:HG23	2.17	0.44
1:C:38:ASP:C	1:C:40:GLU:H	2.20	0.44
2:B:211:ARG:HG2	2:B:211:ARG:HH11	1.83	0.44
1:C:81:GLU:OE2	1:C:142:LYS:NZ	2.45	0.44
2:D:225:TYR:CE1	2:D:277:GLU:HB3	2.52	0.44
2:D:350:TYR:CD2	2:D:390:CYS:HB2	2.53	0.44
2:B:321:HIS:NE2	2:B:379:LYS:HD2	2.32	0.44
1:A:105:LYS:HB2	1:A:105:LYS:HE3	1.82	0.44
2:D:360:PHE:O	2:D:364:LEU:HB2	2.18	0.44
1:A:134:LEU:O	1:A:141:ILE:HA	2.17	0.44
1:C:198:THR:O	1:C:200:ARG:N	2.51	0.44
1:C:50:ARG:HE	1:C:150:ARG:NH1	2.15	0.44
1:A:150:ARG:NH1	2:B:267:PHE:O	2.51	0.44
1:A:101:LEU:N	1:A:102:PRO:CD	2.81	0.44
2:B:364:LEU:HG	2:B:370:GLN:HB2	1.99	0.44
1:C:69:VAL:CG1	1:C:76:LEU:HD21	2.48	0.44
2:B:417:LYS:HE3	2:B:418:TYR:HE1	1.77	0.44
1:C:121:HIS:C	1:C:122:ARG:HG3	2.38	0.44
1:A:212:LEU:HG	1:A:216:PHE:CZ	2.52	0.44
2:B:211:ARG:HG2	2:B:211:ARG:NH1	2.32	0.44
2:B:286:TYR:HA	2:B:290:GLN:OE1	2.17	0.44
1:A:164:VAL:CG1	1:A:165:THR:N	2.81	0.43
2:D:310:THR:OG1	2:D:313:GLN:HG3	2.17	0.43
1:A:253:PRO:HD2	1:A:254:PRO:HD3	2.00	0.43
2:B:421:VAL:HA	2:B:424:LEU:HG	2.00	0.43
1:C:223:ASP:H	1:C:226:VAL:HG12	1.83	0.43
1:A:198:THR:O	1:A:199:ARG:HB2	2.19	0.43
1:A:165:THR:HG21	1:A:168:TYR:CD1	2.46	0.43
2:D:225:TYR:CE2	2:D:281:ILE:CG2	3.01	0.43
2:D:334:MET:HE1	4:D:122:HOH:O	2.18	0.43
1:A:125:HIS:O	1:A:126:ARG:HB2	2.18	0.43
2:D:347:TYR:OH	2:D:394:LEU:HA	2.19	0.43
1:A:233:MET:HE2	1:A:233:MET:HB3	1.92	0.43
1:C:84:HIS:CD2	1:C:136:ASN:HA	2.53	0.43
1:C:7:VAL:HG11	1:C:20:LYS:HE3	2.00	0.43
1:A:167:TRP:CD1	1:A:204:PRO:HA	2.54	0.43
1:A:55:LEU:HD12	1:A:55:LEU:HA	1.78	0.43
2:B:190:GLU:HG3	2:B:351:LEU:HD22	2.00	0.43
1:C:106:SER:HB2	1:C:290:THR:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:203:GLN:OE1	2:B:247:SER:HA	2.19	0.43
2:B:415:ASN:OD1	2:B:416:SER:N	2.52	0.43
1:A:41:THR:HB	2:B:288:LYS:HZ3	1.83	0.42
1:C:122:ARG:O	1:C:151:ALA:HA	2.18	0.42
1:C:91:MET:HE1	1:C:195:GLU:HG2	2.01	0.42
2:D:275:VAL:HG21	2:D:292:LEU:HD21	2.00	0.42
1:C:257:GLU:HA	1:C:257:GLU:OE1	2.19	0.42
2:D:272:PRO:HA	2:D:273:PRO:HD3	1.89	0.42
1:A:62:ASN:HA	1:A:142:LYS:HG2	2.00	0.42
2:B:395:HIS:ND1	2:B:430:LEU:HG	2.34	0.42
2:D:343:ASP:OD1	2:D:404:HIS:HE1	2.01	0.42
1:C:46:SER:O	1:C:50:ARG:HG3	2.19	0.42
1:C:64:VAL:HG12	1:C:143:LEU:O	2.19	0.42
1:A:255:LEU:O	1:A:260:ARG:NH1	2.52	0.42
1:A:37:LEU:HB3	1:A:74:ASN:O	2.20	0.42
1:C:126:ARG:NH1	1:C:159:TYR:HE1	2.14	0.42
1:C:226:VAL:O	1:C:226:VAL:HG22	2.20	0.42
1:A:171:PRO:HD3	1:A:187:TRP:CZ2	2.54	0.42
1:A:181:SER:HA	2:B:173:ASN:HB3	2.02	0.42
2:B:387:LEU:O	2:B:391:LEU:HB2	2.20	0.42
2:B:308:ALA:HA	2:B:309:PRO:HD3	1.97	0.42
2:D:194:LYS:HE2	2:D:194:LYS:HB3	1.76	0.42
2:D:398:TYR:CD1	2:D:426:PRO:HB3	2.54	0.42
1:A:49:ILE:HG23	2:B:306:LEU:HD12	2.02	0.42
2:D:216:ASP:OD2	2:D:408:SER:HB2	2.20	0.42
1:A:30:VAL:CG2	1:A:80:PHE:O	2.68	0.42
2:D:229:ASN:HD22	2:D:334:MET:HE2	1.83	0.42
1:A:170:ALA:HB1	1:A:172:GLU:OE2	2.19	0.42
2:D:294:MET:O	2:D:298:VAL:HG23	2.20	0.42
1:C:88:LYS:HD2	1:C:131:GLN:HG3	2.02	0.42
1:C:150:ARG:NH2	2:D:270:ILE:HA	2.35	0.42
2:B:173:ASN:HD22	2:B:173:ASN:HA	1.69	0.42
1:C:60:HIS:HD2	1:C:62:ASN:N	2.08	0.41
2:B:388:LYS:CE	2:B:392:MET:HE3	2.50	0.41
1:A:165:THR:HG22	4:A:535:HOH:O	2.18	0.41
1:C:35:ILE:O	1:C:75:LYS:HB2	2.19	0.41
2:D:374:GLU:CG	2:D:378:ARG:HH12	2.32	0.41
1:A:42:GLU:OE1	2:B:275:VAL:HG23	2.20	0.41
1:C:203:PHE:HB2	1:C:211:GLN:HE22	1.85	0.41
2:B:287:THR:N	2:B:290:GLN:OE1	2.44	0.41
2:D:394:LEU:HD12	2:D:394:LEU:HA	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:LYS:HG3	1:C:82:PHE:CZ	2.55	0.41
1:C:18:VAL:HA	1:C:32:LEU:O	2.20	0.41
1:A:99:ILE:HG23	1:A:103:LEU:HD23	2.02	0.41
1:A:211:GLN:O	1:A:215:ILE:HG13	2.20	0.41
1:A:89:LYS:HD2	3:A:501:107:S22	2.60	0.41
1:C:237:LYS:HA	1:C:238:PRO:HD3	1.94	0.41
2:B:338:GLU:O	2:B:409:ILE:HD13	2.21	0.41
1:A:165:THR:HG23	1:A:168:TYR:H	1.86	0.41
1:A:124:LEU:CD2	1:A:182:THR:HA	2.50	0.41
2:B:196:LYS:HG2	2:B:244:SER:HB3	2.01	0.41
1:C:67:LEU:O	1:C:68:ASP:HB2	2.21	0.41
2:B:287:THR:HG22	2:B:288:LYS:N	2.36	0.41
2:B:342:ILE:HG22	2:B:343:ASP:OD1	2.21	0.41
2:D:175:VAL:N	2:D:176:PRO:CD	2.84	0.41
1:C:174:LEU:C	1:C:176:GLY:H	2.24	0.41
2:D:175:VAL:N	2:D:176:PRO:HD2	2.36	0.41
1:A:73:GLU:H	1:A:73:GLU:HG3	1.43	0.41
1:A:1:MET:HE2	1:A:70:ILE:HG21	2.02	0.41
1:C:17:VAL:CG2	1:C:18:VAL:N	2.83	0.41
1:C:71:HIS:NE2	2:D:296:HIS:CD2	2.89	0.41
1:A:113:GLN:HG3	1:A:281:LEU:HD21	2.02	0.41
2:B:293:ARG:HH22	1:C:2:GLU:HB2	1.85	0.41
1:A:50:ARG:CD	1:A:150:ARG:NH1	2.84	0.41
2:B:397:THR:HG23	4:B:56:HOH:O	2.21	0.41
1:A:101:LEU:HD23	1:A:104:ILE:HD12	2.02	0.40
1:C:91:MET:HE2	1:C:196:MET:HA	2.02	0.40
1:C:129:LYS:HA	1:C:192:ILE:HD11	2.03	0.40
2:D:365:TYR:O	2:D:369:GLY:HA2	2.21	0.40
1:A:224:GLU:OE2	1:A:231:THR:OG1	2.23	0.40
2:B:250:ARG:HG3	2:B:251:GLY:N	2.36	0.40
1:A:56:LYS:HB3	1:A:56:LYS:HE2	1.86	0.40
1:C:48:ALA:O	1:C:52:ILE:HG13	2.21	0.40
1:C:34:LYS:HG3	1:C:77:TYR:CE2	2.57	0.40
1:A:236:TYR:CG	1:A:237:LYS:N	2.88	0.40
2:D:415:ASN:HD22	2:D:417:LYS:N	2.04	0.40
1:C:165:THR:CG2	1:C:168:TYR:HD1	2.34	0.40
1:C:12:GLU:HA	1:C:17:VAL:HA	2.04	0.40
2:B:388:LYS:HG3	2:B:392:MET:HE3	2.03	0.40
2:B:196:LYS:CG	2:B:244:SER:HB3	2.51	0.40
2:B:347:TYR:OH	2:B:394:LEU:HA	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	296/298 (99%)	274 (93%)	17 (6%)	5 (2%)	11	36
1	C	296/298 (99%)	255 (86%)	35 (12%)	6 (2%)	9	30
2	B	258/260 (99%)	240 (93%)	15 (6%)	3 (1%)	16	47
2	D	258/260 (99%)	242 (94%)	13 (5%)	3 (1%)	16	47
All	All	1108/1116 (99%)	1011 (91%)	80 (7%)	17 (2%)	13	40

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	GLU
2	B	346	PRO
1	C	39	THR
2	D	346	PRO
1	C	199	ARG
1	C	297	ARG
1	C	145	ASP
2	B	327	CYS
2	D	304	PHE
1	C	234	PRO
1	A	234	PRO
2	D	345	ASP
1	A	271	PRO
2	B	281	ILE
1	A	230	VAL
1	C	64	VAL
1	A	164	VAL

5.3.2 Protein sidechains [i](#)

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	259/263 (98%)	253 (98%)	6 (2%)	58	88
1	C	260/263 (99%)	251 (96%)	9 (4%)	43	77
2	B	234/234 (100%)	227 (97%)	7 (3%)	48	82
2	D	234/234 (100%)	222 (95%)	12 (5%)	29	63
All	All	987/994 (99%)	953 (97%)	34 (3%)	44	78

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	38	ASP
1	A	73	GLU
1	A	74	ASN
1	A	76	LEU
1	A	122	ARG
2	B	173	ASN
2	B	174	GLU
2	B	232	LEU
2	B	241	ARG
2	B	250	ARG
2	B	403	GLN
2	B	410	ARG
1	C	47	THR
1	C	74	ASN
1	C	75	LYS
1	C	76	LEU
1	C	97	THR
1	C	122	ARG
1	C	131	GLN
1	C	152	PHE
1	C	202	LEU
2	D	173	ASN
2	D	177	ASP
2	D	179	HIS
2	D	193	CYS
2	D	194	LYS
2	D	199	TYR

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Mol	Chain	Res	Type
2	D	232	LEU
2	D	284	ASP
2	D	323	GLN
2	D	345	ASP
2	D	398	TYR
2	D	415	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	A	62	ASN
1	A	74	ASN
1	A	84	HIS
1	A	85	GLN
1	A	246	GLN
1	A	272	ASN
1	A	287	GLN
2	B	173	ASN
2	B	179	HIS
2	B	208	ASN
2	B	237	ASN
2	B	361	HIS
2	B	370	GLN
2	B	396	GLN
2	B	403	GLN
2	B	406	GLN
1	C	60	HIS
1	C	62	ASN
1	C	84	HIS
1	C	85	GLN
1	C	113	GLN
1	C	119	HIS
1	C	131	GLN
1	C	246	GLN
1	C	268	HIS
2	D	179	HIS
2	D	208	ASN
2	D	229	ASN
2	D	233	HIS
2	D	254	GLN
2	D	296	HIS

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Mol	Chain	Res	Type
2	D	322	GLN
2	D	404	HIS
2	D	406	GLN
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 carbohydrates 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	107	A	501	-	28,35,35	3.59	16 (57%)	34,51,51	1.67	7 (20%)
3	107	C	601	-	28,35,35	3.63	17 (60%)	34,51,51	1.91	12 (35%)

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	107	A	501	-	-	0/14/28/28	0/5/5/5
3	107	C	601	-	-	0/14/28/28	0/5/5/5

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	601	107	C19-S22	2.02	1.79	1.76
3	A	501	107	C6-C7	2.10	1.45	1.41
3	A	501	107	C14-N15	2.25	1.49	1.45
3	C	601	107	S22-N25	2.36	1.67	1.63
3	C	601	107	C6-C7	2.71	1.46	1.41
3	C	601	107	C7-C11	2.84	1.51	1.42
3	A	501	107	C7-C11	3.18	1.52	1.42
3	C	601	107	C28-C29	3.47	1.48	1.37
3	A	501	107	C28-C29	3.59	1.48	1.37
3	A	501	107	C27-C32	3.98	1.47	1.38
3	A	501	107	C28-C27	4.09	1.48	1.38
3	C	601	107	C28-C27	4.11	1.48	1.38
3	A	501	107	C18-C17	4.13	1.46	1.38
3	A	501	107	C29-N30	4.34	1.44	1.34
3	C	601	107	C32-C31	4.44	1.49	1.39
3	C	601	107	C27-C32	4.61	1.48	1.38
3	C	601	107	C18-C17	4.61	1.47	1.38
3	C	601	107	C29-N30	4.70	1.44	1.34
3	A	501	107	C32-C31	4.89	1.50	1.39
3	C	601	107	C21-C20	4.91	1.47	1.38
3	A	501	107	C6-C5	4.99	1.47	1.36
3	A	501	107	C21-C20	5.06	1.47	1.38
3	C	601	107	C21-C16	5.37	1.48	1.39
3	A	501	107	C31-N30	5.42	1.44	1.34
3	C	601	107	C31-N30	5.54	1.44	1.34
3	C	601	107	C18-C19	5.63	1.48	1.38
3	A	501	107	C20-C19	5.63	1.48	1.38
3	C	601	107	C6-C5	5.65	1.48	1.36
3	A	501	107	C21-C16	5.81	1.48	1.39
3	A	501	107	C18-C19	6.00	1.48	1.38
3	C	601	107	C17-C16	6.02	1.49	1.39
3	C	601	107	C20-C19	6.54	1.49	1.38
3	A	501	107	C17-C16	6.79	1.50	1.39

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	501	107	C12-C11-C7	-5.68	118.25	123.91
3	C	601	107	C12-C11-C7	-4.95	118.97	123.91
3	C	601	107	O23-S22-C19	-3.38	103.68	107.96
3	A	501	107	C21-C16-N15	-2.58	116.12	121.06
3	C	601	107	C21-C16-N15	-2.20	116.84	121.06
3	C	601	107	O24-S22-O23	-2.15	116.68	119.54
3	A	501	107	C18-C19-S22	-2.06	117.40	119.78
3	A	501	107	N25-C31-N30	2.16	116.48	114.27
3	C	601	107	C4-N3-C2	2.20	110.36	106.81
3	C	601	107	C14-N15-C16	2.21	127.12	122.68
3	C	601	107	C31-N25-S22	2.21	130.06	124.99
3	C	601	107	O24-S22-C19	2.37	110.96	107.96
3	A	501	107	C4-N3-C2	2.43	110.74	106.81
3	A	501	107	C14-N15-C16	2.46	127.63	122.68
3	C	601	107	C9-S10-C11	2.57	93.58	90.63
3	C	601	107	C14-C13-C2	2.96	127.30	122.53
3	A	501	107	C29-N30-C31	3.12	121.58	117.28
3	C	601	107	N25-C31-N30	3.32	117.67	114.27
3	C	601	107	C29-N30-C31	3.43	122.01	117.28

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	107	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	298/298 (100%)	-0.33	9 (3%) 54 41	31, 46, 83, 99	0
1	C	298/298 (100%)	-0.02	15 (5%) 32 21	38, 62, 87, 97	0
2	B	260/260 (100%)	-0.27	7 (2%) 58 45	34, 53, 75, 95	0
2	D	260/260 (100%)	-0.33	12 (4%) 36 25	29, 43, 73, 103	0
All	All	1116/1116 (100%)	-0.23	43 (3%) 43 31	29, 51, 84, 103	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	176	PRO	6.7
2	D	174	GLU	6.0
2	D	173	ASN	5.2
1	A	14	THR	4.4
2	D	177	ASP	4.4
1	C	40	GLU	4.2
2	D	175	VAL	4.2
1	A	298	LEU	4.2
2	D	323	GLN	3.8
1	C	96	LEU	3.5
1	A	38	ASP	3.4
1	C	14	THR	3.4
1	C	38	ASP	3.4
1	A	73	GLU	2.9
1	C	295	HIS	2.9
2	B	173	ASN	2.9
2	B	176	PRO	2.9
2	D	345	ASP	2.9
2	B	345	ASP	2.8
2	D	193	CYS	2.7
1	A	12	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	36	ARG	2.6
2	B	284	ASP	2.6
1	C	287	GLN	2.5
1	C	165	THR	2.5
2	B	323	GLN	2.5
1	C	250	LYS	2.4
1	C	288	ASP	2.4
1	C	293	VAL	2.4
1	C	13	GLY	2.3
1	A	159	TYR	2.2
1	A	74	ASN	2.2
2	D	324	PRO	2.2
1	C	161	HIS	2.2
2	B	177	ASP	2.1
1	A	40	GLU	2.1
1	C	39	THR	2.1
1	C	41	THR	2.1
1	A	161	HIS	2.1
2	D	178	TYR	2.1
2	D	271	TYR	2.1
2	B	257	GLY	2.0
2	D	180	GLU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	107	C	601	31/31	0.93	0.20	0.04	69,72,76,76	0
3	107	A	501	31/31	0.95	0.17	-0.29	39,48,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.