Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 09:01 PM GMT

PDB ID	:	1N6M
Title	:	Rotation of the stalk/neck and one head in a new crystal structure of the kinesin motor protein, Ncd
Authors	:	Yun, M.; Bronner, C.E.; Park, C.-G.; Cha, S.-S.; Park, H.-W.; Endow, S.A.
Deposited on	:	2002-11-11
Resolution	:	2.50 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 (i) symbol.

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

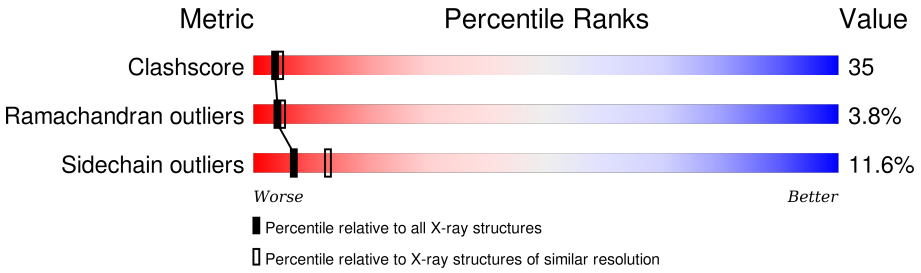
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	409	45% 38% 6% 11%
1	B	409	33% 39% 9% 18%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5886 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 Claret segregational protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2893	1804	507	561	21			
1	B	335	Total	C	N	O	S	0	0	0
			2656	1659	464	515	18			

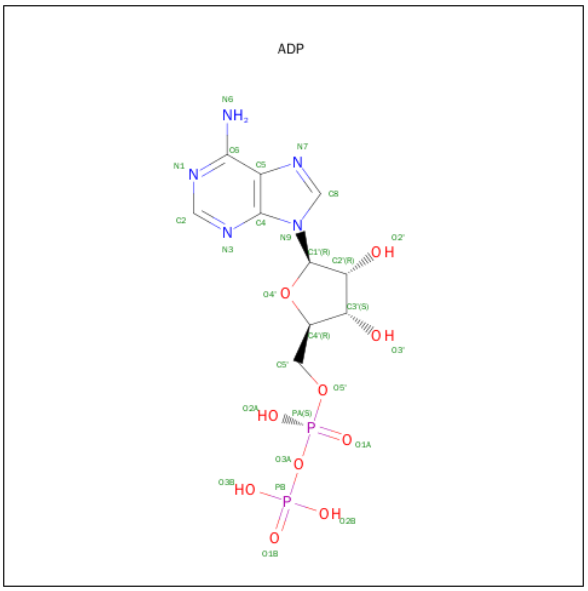
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	292	MET	-	INITIATING MET	UNP P20480
A	600	LYS	ASN	ENGINEERED	UNP P20480
B	292	MET	-	INITIATING MET	UNP P20480
B	600	LYS	ASN	ENGINEERED	UNP P20480

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	166	Total	O	0	0
			166	166		
4	B	115	Total	O	0	0
			115	115		

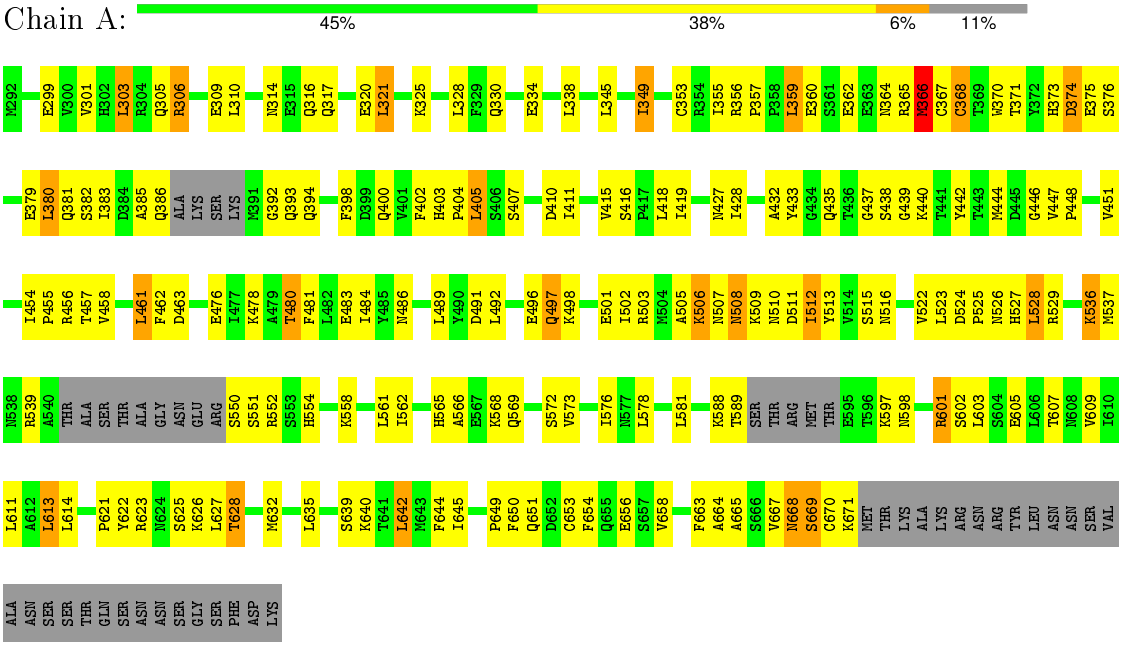
3 Residue-property plots (i)

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

Note EDS was not executed.

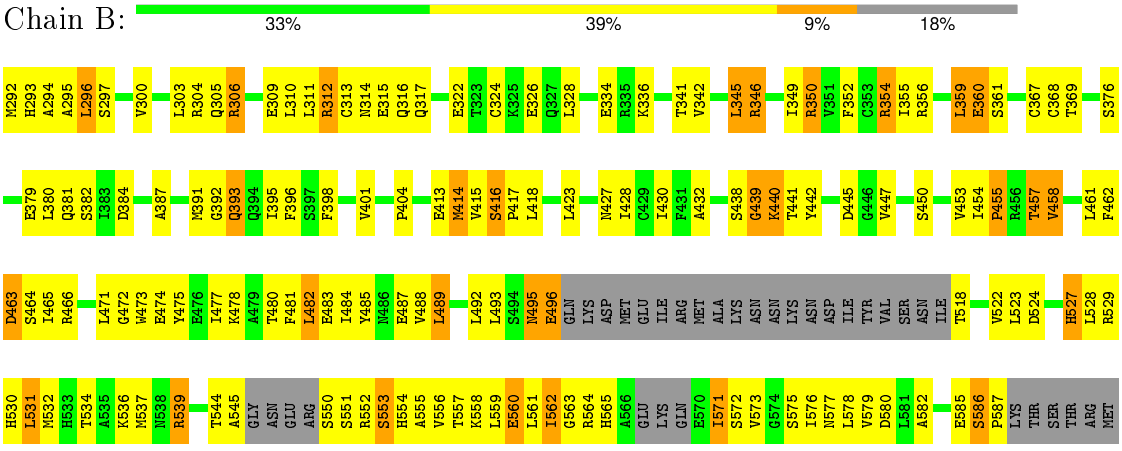
- Molecule 1: Claret segregational protein

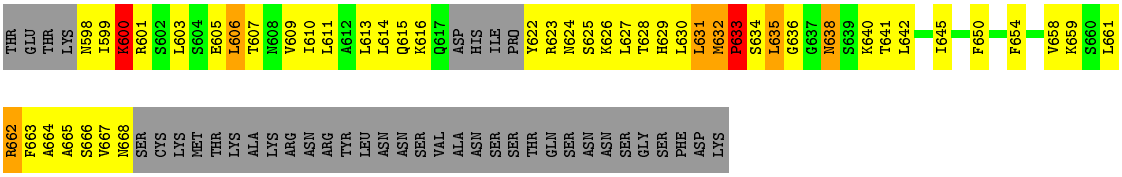
Chain A:



- Molecule 1: Claret segregational protein

Chain B:




THR GLU THR LYS N598 I599 R600 R601 S602 L603 S604 E605 L606 T607 N608 V609 I610 L611 A612 L613 L614 Q615 K616 Q617 ASP HIS ILE PRO Y622 R623 N624 S625 K626 L627 T628 H629 L630 L631 M632 P633 S634 L635 G636 G637 N638 S639 K640 T641 L642 I645 F650 F654 V658 K659 S660 L661
R662 F663 A664 A665 S666 V667 N668 SER CYS LYS MET THR LYS ALA LYS ARG ASN ARG TYR LEU ASN ASN SER VAL ALA ASN SER SER THR GLN SER ASN ASN SER GLY SER PHE ASP LYS

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	162.60Å 66.60Å 94.80Å 90.00° 98.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.260 , 0.302	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5886	wwPDB-VP
Average B, all atoms (Å^2)	61.0	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: MG, ADP

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	0.43	0/2939	0.65	0/3962
1	B	0.44	0/2697	0.70	2/3637 (0.1%)
All	All	0.44	0/5636	0.68	2/7599 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	495	ASN	N-CA-C	5.98	127.14	111.00
1	B	586	SER	C-N-CD	-5.24	109.07	120.60

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	2893	0	2860	166	0
1	B	2656	0	2621	234	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	12	6	0
3	B	27	0	12	6	0
4	A	166	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	115	0	0	1	0
All	All	5886	0	5505	394	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 (394) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:622:TYR:CD2	1:B:632:MET:HB2	1.97	0.99
1:A:359:LEU:HD13	1:A:359:LEU:H	1.22	0.98
1:A:497:GLN:HE21	1:A:498:LYS:N	1.63	0.95
1:A:382:SER:HB2	1:A:651:GLN:HE22	1.37	0.88
1:B:562:ILE:HG22	1:B:573:VAL:HG22	1.55	0.86
1:A:310:LEU:O	1:A:314:ASN:HB2	1.75	0.86
1:B:439:GLY:HA2	3:B:999:ADP:O1A	1.76	0.85
1:B:447:VAL:HG23	1:B:450:SER:H	1.42	0.84
1:B:376:SER:HB2	1:B:398:PHE:O	1.79	0.83
1:B:478:LYS:HD3	1:B:560:GLU:HB3	1.59	0.83
1:A:523:LEU:H	1:A:527:HIS:HD2	1.27	0.81
1:B:559:LEU:HB3	1:B:576:ILE:HB	1.63	0.80
1:B:600:LYS:HE3	1:B:600:LYS:HA	1.63	0.80
1:B:555:ALA:HB3	1:B:580:ASP:HB3	1.65	0.79
1:B:606:LEU:O	1:B:610:ILE:HG12	1.84	0.77
1:B:632:MET:N	1:B:633:PRO:HD3	2.00	0.77
1:B:654:PHE:O	1:B:658:VAL:HG23	1.84	0.76
1:A:356:ARG:HH12	1:A:359:LEU:HD12	1.51	0.76
1:B:472:GLY:O	1:B:565:HIS:ND1	2.18	0.75
1:B:633:PRO:CD	1:B:634:SER:H	2.00	0.75
1:B:633:PRO:HD2	1:B:634:SER:H	1.51	0.75
1:B:635:LEU:HD12	1:B:636:GLY:H	1.50	0.75
1:A:523:LEU:H	1:A:527:HIS:CD2	2.06	0.74
1:B:439:GLY:CA	3:B:999:ADP:O1A	2.35	0.73
1:B:544:THR:O	1:B:545:ALA:O	2.05	0.73
1:B:423:LEU:HD12	1:B:423:LEU:H	1.54	0.73
1:B:524:ASP:O	1:B:527:HIS:HB3	1.90	0.72
1:B:522:VAL:HG11	1:B:528:LEU:HB2	1.71	0.72
1:B:489:LEU:HG	1:B:626:LYS:HE2	1.71	0.72
1:A:301:VAL:O	1:A:305:GLN:HG3	1.90	0.72
1:A:454:ILE:HB	1:A:455:PRO:HD3	1.71	0.72
1:B:483:GLU:HG2	1:B:484:ILE:H	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:668:ASN:H	1:A:668:ASN:ND2	1.88	0.71
1:B:427:ASN:HD22	1:B:634:SER:HA	1.56	0.71
1:B:560:GLU:O	1:B:561:LEU:HD23	1.91	0.70
1:B:440:LYS:HZ3	1:B:440:LYS:HB2	1.56	0.70
1:B:544:THR:O	1:B:544:THR:CG2	2.40	0.70
1:A:359:LEU:HD23	1:A:362:GLU:H	1.57	0.69
1:B:342:VAL:O	1:B:346:ARG:HG2	1.92	0.69
1:B:454:ILE:HB	1:B:455:PRO:HD3	1.73	0.69
1:B:632:MET:N	1:B:633:PRO:CD	2.56	0.69
1:A:447:VAL:HG22	1:A:448:PRO:HD2	1.74	0.69
1:A:393:GLN:O	1:A:394:GLN:NE2	2.27	0.68
1:A:668:ASN:HD22	1:A:668:ASN:H	1.42	0.68
1:A:303:LEU:HD23	1:B:303:LEU:HB2	1.75	0.67
1:A:497:GLN:NE2	1:A:498:LYS:N	2.40	0.67
1:B:610:ILE:O	1:B:614:LEU:HG	1.94	0.67
1:B:635:LEU:HD12	1:B:636:GLY:N	2.10	0.66
1:B:312:ARG:NH1	1:B:316:GLN:HG3	2.10	0.66
1:A:435:GLN:OE1	1:A:656:GLU:HG2	1.95	0.66
1:B:445:ASP:OD2	1:B:454:ILE:HD12	1.96	0.66
1:B:480:THR:HB	1:B:558:LYS:HB2	1.76	0.66
1:A:511:ASP:OD1	1:A:512:ILE:N	2.28	0.66
1:B:607:THR:HG22	1:B:663:PHE:CE1	2.30	0.66
1:A:663:PHE:O	1:A:667:VAL:HG23	1.96	0.66
1:B:463:ASP:O	1:B:466:ARG:HG2	1.96	0.65
1:B:599:ILE:HG22	1:B:601:ARG:H	1.61	0.65
1:B:440:LYS:NZ	3:B:999:ADP:O2B	2.30	0.65
1:B:495:ASN:O	1:B:496:GLU:OE1	2.14	0.65
1:A:355:ILE:HG13	1:A:404:PRO:HD3	1.77	0.65
1:A:489:LEU:N	1:A:489:LEU:HD12	2.12	0.65
1:A:359:LEU:CD1	1:A:359:LEU:H	2.01	0.65
1:A:515:SER:O	1:A:516:ASN:HB2	1.94	0.65
1:A:357:PRO:HA	1:A:404:PRO:HB3	1.79	0.65
1:B:638:ASN:O	1:B:638:ASN:CG	2.35	0.65
1:A:523:LEU:N	1:A:527:HIS:HD2	1.95	0.64
1:B:292:MET:HG2	1:B:294:ALA:H	1.62	0.64
1:B:474:GLU:O	1:B:563:GLY:HA2	1.97	0.64
1:B:453:VAL:O	1:B:457:THR:HG23	1.98	0.64
1:A:303:LEU:HD23	1:B:303:LEU:CB	2.28	0.64
1:B:384:ASP:OD1	1:B:387:ALA:HB2	1.98	0.64
1:B:481:PHE:CZ	1:B:531:LEU:HD13	2.33	0.63
1:A:622:TYR:CG	1:A:632:MET:HG3	2.33	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:473:TRP:HE1	1:B:563:GLY:C	2.02	0.63
1:B:369:THR:OG1	1:B:381:GLN:HB2	1.99	0.63
1:A:427:ASN:C	1:A:428:ILE:HD12	2.19	0.63
1:A:497:GLN:HE21	1:A:497:GLN:C	2.01	0.62
1:A:668:ASN:N	1:A:668:ASN:HD22	1.96	0.62
1:B:458:VAL:HG13	1:B:528:LEU:HD23	1.80	0.62
1:B:631:LEU:C	1:B:633:PRO:CD	2.68	0.62
1:A:359:LEU:HD13	1:A:359:LEU:N	2.05	0.62
1:A:349:ILE:HD13	1:A:349:ILE:O	1.99	0.61
1:B:430:ILE:HD12	1:B:642:LEU:HD23	1.81	0.61
1:A:507:ASN:O	1:A:508:ASN:HB2	2.00	0.61
1:A:562:ILE:HA	1:A:573:VAL:HG12	1.83	0.61
1:B:544:THR:HG22	1:B:544:THR:O	2.00	0.61
1:A:613:LEU:HD21	1:A:635:LEU:O	2.01	0.61
1:B:440:LYS:NZ	1:B:440:LYS:HB2	2.15	0.61
1:B:623:ARG:HD3	1:B:629:HIS:CE1	2.36	0.60
1:A:444:MET:CE	1:A:578:LEU:HB3	2.32	0.60
1:A:524:ASP:HB2	1:A:525:PRO:HD2	1.83	0.60
1:A:357:PRO:CA	1:A:404:PRO:HB3	2.31	0.60
1:B:495:ASN:C	1:B:496:GLU:HG2	2.22	0.60
1:A:489:LEU:HD11	1:A:626:LYS:CG	2.32	0.60
1:A:489:LEU:HD11	1:A:626:LYS:HG2	1.83	0.60
1:B:600:LYS:CA	1:B:600:LYS:HE3	2.32	0.60
1:B:613:LEU:HD13	1:B:635:LEU:CD1	2.32	0.60
1:A:380:LEU:H	1:A:380:LEU:HD12	1.67	0.59
1:A:447:VAL:CG2	1:A:448:PRO:HD2	2.31	0.59
1:B:633:PRO:CD	1:B:634:SER:N	2.66	0.59
1:A:561:LEU:O	1:A:573:VAL:HA	2.03	0.59
1:A:411:ILE:N	1:A:411:ILE:HD12	2.16	0.59
1:B:605:GLU:O	1:B:609:VAL:HG23	2.03	0.59
1:B:605:GLU:CD	1:B:624:ASN:HB2	2.23	0.59
1:A:550:SER:C	1:A:552:ARG:H	2.05	0.58
1:A:418:LEU:HD12	1:A:642:LEU:HD22	1.85	0.58
1:A:360:GLU:CD	1:A:360:GLU:H	2.07	0.58
1:A:317:GLN:HB2	1:B:317:GLN:NE2	2.17	0.58
1:B:536:LYS:O	1:B:539:ARG:HB3	2.02	0.58
1:A:668:ASN:O	1:A:671:LYS:N	2.34	0.57
1:A:411:ILE:HG23	1:A:644:PHE:CE1	2.38	0.57
1:A:367:CYS:HA	1:A:650:PHE:HA	1.87	0.57
1:B:551:SER:OG	1:B:552:ARG:HD2	2.04	0.57
1:B:554:HIS:HE1	1:B:603:LEU:HD11	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:HIS:ND1	1:A:568:LYS:HB2	2.19	0.57
1:A:411:ILE:HG23	1:A:644:PHE:CZ	2.40	0.57
1:A:415:VAL:HG22	1:A:642:LEU:HD23	1.87	0.57
1:A:621:PRO:HA	4:A:6126:HOH:O	2.03	0.57
1:B:428:ILE:N	1:B:428:ILE:HD12	2.19	0.57
1:B:622:TYR:HB2	1:B:632:MET:SD	2.45	0.57
1:A:665:ALA:O	1:A:669:SER:HB3	2.04	0.57
1:B:349:ILE:HG21	1:B:614:LEU:CD2	2.34	0.56
1:B:306:ARG:NH2	1:B:310:LEU:HD11	2.19	0.56
1:B:454:ILE:HG12	1:B:578:LEU:HD13	1.87	0.56
1:A:435:GLN:HE22	1:A:653:CYS:HB3	1.70	0.56
1:B:631:LEU:C	1:B:633:PRO:HD2	2.26	0.56
1:A:359:LEU:CD2	1:A:362:GLU:H	2.17	0.56
1:B:457:THR:O	1:B:458:VAL:C	2.44	0.56
1:A:601:ARG:HG3	1:A:602:SER:N	2.21	0.56
1:B:427:ASN:HD22	1:B:634:SER:CB	2.19	0.56
1:A:625:SER:HB3	1:A:628:THR:OG1	2.05	0.56
1:B:609:VAL:O	1:B:613:LEU:HG	2.06	0.56
1:B:495:ASN:O	1:B:496:GLU:HG2	2.05	0.55
1:A:481:PHE:N	1:A:481:PHE:CD1	2.73	0.55
1:A:654:PHE:O	1:A:658:VAL:HG23	2.07	0.55
1:A:368:CYS:HB3	1:A:381:GLN:O	2.06	0.55
1:B:427:ASN:HD22	1:B:634:SER:CA	2.18	0.55
1:A:303:LEU:HB3	1:B:303:LEU:HD13	1.88	0.55
1:B:349:ILE:HD13	1:B:614:LEU:HD21	1.89	0.55
1:B:477:ILE:HG13	1:B:561:LEU:HD22	1.87	0.54
1:B:571:ILE:O	1:B:573:VAL:HG23	2.07	0.54
1:B:416:SER:HB3	1:B:417:PRO:HD3	1.89	0.54
1:B:562:ILE:HD13	1:B:562:ILE:N	2.22	0.54
1:B:293:HIS:HD2	1:B:296:LEU:HD22	1.71	0.54
1:A:439:GLY:CA	3:A:998:ADP:O1A	2.56	0.54
1:B:528:LEU:O	1:B:532:MET:HB2	2.07	0.54
1:B:633:PRO:C	1:B:635:LEU:H	2.11	0.54
1:B:352:PHE:CD1	1:B:642:LEU:HD11	2.42	0.54
1:B:605:GLU:HG3	1:B:628:THR:HG21	1.90	0.54
1:B:391:MET:O	1:B:393:GLN:N	2.39	0.54
1:B:658:VAL:O	1:B:662:ARG:HD3	2.08	0.54
1:B:482:LEU:HD23	1:B:482:LEU:N	2.22	0.54
1:A:400:GLN:HG2	1:A:402:PHE:CZ	2.43	0.54
1:A:359:LEU:HD22	1:A:359:LEU:O	2.07	0.53
1:B:531:LEU:HD12	1:B:532:MET:N	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:TRP:CZ2	1:A:649:PRO:HA	2.43	0.53
1:B:475:TYR:HA	1:B:562:ILE:O	2.09	0.53
1:A:614:LEU:HD11	1:A:667:VAL:HG13	1.90	0.53
1:B:552:ARG:HG3	1:B:600:LYS:HD3	1.91	0.53
1:B:622:TYR:CD2	1:B:632:MET:CB	2.84	0.53
1:B:623:ARG:HD3	1:B:629:HIS:ND1	2.24	0.53
1:B:559:LEU:CD2	1:B:561:LEU:HD21	2.39	0.52
1:A:668:ASN:N	1:A:668:ASN:ND2	2.53	0.52
1:B:293:HIS:O	1:B:296:LEU:HD23	2.09	0.52
1:A:328:LEU:HD13	1:B:328:LEU:HA	1.91	0.52
1:B:341:THR:O	1:B:345:LEU:HB2	2.09	0.52
1:A:439:GLY:N	3:A:998:ADP:O1A	2.43	0.52
1:B:480:THR:C	1:B:481:PHE:CD1	2.83	0.52
1:B:522:VAL:HG13	1:B:527:HIS:CD2	2.45	0.52
1:B:524:ASP:HB3	1:B:527:HIS:HB3	1.91	0.52
1:B:495:ASN:O	1:B:496:GLU:CG	2.57	0.52
1:B:560:GLU:C	1:B:561:LEU:HD23	2.30	0.52
1:B:614:LEU:CD1	1:B:667:VAL:HG13	2.40	0.52
1:A:605:GLU:O	1:A:609:VAL:HG23	2.10	0.52
1:B:658:VAL:HG12	1:B:662:ARG:HD3	1.91	0.52
1:B:613:LEU:HD21	1:B:622:TYR:OH	2.09	0.51
1:A:356:ARG:HD2	3:A:998:ADP:C8	2.45	0.51
1:B:631:LEU:HD12	1:B:631:LEU:H	1.75	0.51
1:B:529:ARG:HH11	1:B:529:ARG:HG2	1.75	0.51
1:B:599:ILE:HG23	1:B:601:ARG:HE	1.74	0.51
1:B:461:LEU:HD11	1:B:561:LEU:HD11	1.91	0.51
1:B:539:ARG:HB2	1:B:539:ARG:HH11	1.73	0.51
1:B:293:HIS:C	1:B:295:ALA:H	2.13	0.51
1:B:380:LEU:HD22	1:B:661:LEU:CD1	2.40	0.51
1:A:356:ARG:NE	1:A:437:GLY:O	2.40	0.51
1:A:394:GLN:OE1	1:A:658:VAL:HG13	2.10	0.51
1:A:503:ARG:O	1:A:513:TYR:HB3	2.10	0.51
1:A:321:LEU:CD1	1:A:325:LYS:HD3	2.40	0.51
1:A:491:ASP:OD1	1:A:492:LEU:N	2.44	0.51
1:B:550:SER:HA	1:B:553:SER:OG	2.10	0.51
1:B:293:HIS:C	1:B:295:ALA:N	2.63	0.51
1:A:383:ILE:O	1:A:383:ILE:HD12	2.11	0.51
1:A:428:ILE:HD12	1:A:428:ILE:N	2.26	0.51
1:A:489:LEU:CD1	1:A:489:LEU:N	2.73	0.51
1:B:623:ARG:HA	1:B:629:HIS:HB2	1.92	0.50
1:B:614:LEU:HD11	1:B:667:VAL:HG13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:THR:HG22	4:A:6154:HOH:O	2.10	0.50
1:A:439:GLY:HA2	3:A:998:ADP:O1A	2.12	0.50
1:B:663:PHE:O	1:B:666:SER:HB3	2.11	0.50
1:B:439:GLY:N	3:B:999:ADP:O1A	2.44	0.50
1:B:297:SER:O	1:B:300:VAL:HG12	2.12	0.50
1:A:503:ARG:HH21	1:A:513:TYR:HE1	1.60	0.50
1:B:428:ILE:HA	1:B:640:LYS:O	2.12	0.50
1:A:316:GLN:O	1:A:320:GLU:HG3	2.12	0.50
1:B:523:LEU:H	1:B:527:HIS:HD2	1.59	0.49
1:B:416:SER:CB	1:B:417:PRO:HD3	2.43	0.49
1:A:497:GLN:HE21	1:A:498:LYS:H	1.52	0.49
1:A:442:TYR:O	1:A:446:GLY:HA2	2.13	0.49
1:A:512:ILE:HG23	1:A:513:TYR:N	2.27	0.49
1:B:629:HIS:HD2	1:B:632:MET:SD	2.35	0.48
1:B:607:THR:O	1:B:611:LEU:HG	2.13	0.48
1:A:498:LYS:NZ	1:A:516:ASN:O	2.39	0.48
1:A:550:SER:O	1:A:552:ARG:N	2.43	0.48
1:A:403:HIS:HB2	1:A:404:PRO:HD2	1.95	0.48
1:A:480:THR:HG23	1:A:558:LYS:HB3	1.94	0.48
1:A:498:LYS:HD2	1:A:516:ASN:HB3	1.94	0.48
1:A:451:VAL:O	1:A:456:ARG:NE	2.41	0.48
1:A:484:ILE:HB	1:A:554:HIS:HB2	1.95	0.48
1:A:403:HIS:HB2	1:A:404:PRO:CD	2.43	0.48
1:B:369:THR:HG1	1:B:381:GLN:HB2	1.78	0.48
1:A:440:LYS:HB2	3:A:998:ADP:O3B	2.13	0.48
1:A:525:PRO:O	1:A:529:ARG:HG3	2.14	0.48
1:B:396:PHE:N	1:B:396:PHE:CD1	2.81	0.48
1:B:522:VAL:HA	1:B:527:HIS:NE2	2.28	0.47
1:B:428:ILE:CD1	1:B:428:ILE:N	2.77	0.47
1:A:451:VAL:HG23	1:A:456:ARG:CZ	2.44	0.47
1:B:447:VAL:O	1:B:450:SER:N	2.47	0.47
1:B:483:GLU:HG2	1:B:484:ILE:N	2.25	0.47
1:A:489:LEU:CD1	1:A:626:LYS:HG2	2.45	0.47
1:B:355:ILE:O	1:B:404:PRO:HA	2.14	0.47
1:B:662:ARG:HH11	1:B:662:ARG:HG2	1.79	0.47
1:A:373:HIS:NE2	1:A:379:GLU:OE2	2.42	0.47
1:A:486:ASN:OD1	1:A:598:ASN:ND2	2.48	0.47
1:B:603:LEU:O	1:B:607:THR:HG23	2.14	0.47
1:B:523:LEU:HD12	1:B:523:LEU:N	2.30	0.47
1:B:633:PRO:CG	1:B:634:SER:N	2.78	0.47
1:B:571:ILE:HG13	1:B:571:ILE:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:488:VAL:O	1:B:488:VAL:HG13	2.15	0.47
1:A:566:ALA:O	1:A:569:GLN:NE2	2.46	0.47
1:B:481:PHE:N	1:B:481:PHE:CD1	2.84	0.46
1:B:556:VAL:HG12	1:B:557:THR:N	2.30	0.46
1:A:667:VAL:O	1:A:670:CYS:HB2	2.15	0.46
1:B:349:ILE:HG21	1:B:614:LEU:HD22	1.97	0.46
1:B:393:GLN:HB2	1:B:393:GLN:HE21	1.60	0.46
1:A:365:ARG:O	1:A:366:MET:C	2.53	0.46
1:B:462:PHE:C	1:B:464:SER:H	2.19	0.46
1:A:310:LEU:O	1:A:314:ASN:CB	2.58	0.46
1:B:586:SER:HA	1:B:587:PRO:HD3	1.43	0.46
1:A:433:TYR:HB3	1:A:645:ILE:HD13	1.98	0.46
1:B:427:ASN:ND2	1:B:634:SER:OG	2.48	0.46
1:B:462:PHE:C	1:B:464:SER:N	2.68	0.46
1:A:306:ARG:HD2	4:A:6255:HOH:O	2.14	0.46
1:A:561:LEU:HD23	1:A:561:LEU:N	2.31	0.46
1:B:376:SER:HA	1:B:401:VAL:HG23	1.97	0.46
1:B:440:LYS:CB	1:B:440:LYS:NZ	2.76	0.46
1:A:371:THR:HG23	1:A:379:GLU:HB3	1.98	0.46
1:B:623:ARG:HD2	4:B:6291:HOH:O	2.16	0.45
1:B:575:SER:O	1:B:576:ILE:HG12	2.17	0.45
1:A:664:ALA:O	1:A:668:ASN:ND2	2.49	0.45
1:B:309:GLU:O	1:B:312:ARG:HG3	2.16	0.45
1:B:601:ARG:O	1:B:605:GLU:N	2.49	0.45
1:B:462:PHE:O	1:B:464:SER:N	2.49	0.45
1:B:524:ASP:HB3	1:B:527:HIS:CB	2.46	0.45
1:B:482:LEU:HD23	1:B:482:LEU:H	1.81	0.45
1:B:599:ILE:HG23	1:B:601:ARG:NE	2.31	0.45
1:A:356:ARG:HH12	1:A:359:LEU:CD1	2.24	0.45
1:B:352:PHE:CE1	1:B:642:LEU:HD11	2.51	0.45
1:A:508:ASN:OD1	1:A:510:ASN:HB2	2.16	0.45
1:A:370:TRP:HZ2	1:A:649:PRO:HA	1.81	0.45
1:B:665:ALA:HA	1:B:668:ASN:HB3	1.98	0.45
1:A:607:THR:O	1:A:611:LEU:HG	2.16	0.45
1:B:658:VAL:O	1:B:662:ARG:HB2	2.17	0.45
1:B:518:THR:O	1:B:518:THR:HG22	2.17	0.45
1:B:473:TRP:HA	1:B:565:HIS:HA	1.99	0.45
1:B:312:ARG:HG3	1:B:313:CYS:N	2.32	0.45
1:A:330:GLN:O	1:A:334:GLU:HG3	2.17	0.45
1:B:489:LEU:HD12	1:B:489:LEU:N	2.32	0.45
1:A:484:ILE:HD12	1:A:627:LEU:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:LYS:HD3	1:A:539:ARG:CZ	2.46	0.45
1:B:605:GLU:CD	1:B:624:ASN:HD22	2.21	0.44
1:A:306:ARG:NH2	1:A:310:LEU:HD11	2.31	0.44
1:B:613:LEU:HD22	1:B:636:GLY:HA2	1.98	0.44
1:A:622:TYR:CD2	1:A:632:MET:HG3	2.53	0.44
1:B:556:VAL:HG12	1:B:557:THR:H	1.81	0.44
1:B:629:HIS:HD2	1:B:632:MET:CE	2.30	0.44
1:A:668:ASN:C	1:A:670:CYS:N	2.71	0.44
1:B:418:LEU:O	1:B:428:ILE:HG12	2.17	0.44
1:B:396:PHE:HD2	1:B:665:ALA:HB2	1.82	0.44
1:B:463:ASP:C	1:B:466:ARG:HG2	2.37	0.44
1:A:522:VAL:CG2	1:A:528:LEU:HG	2.47	0.44
1:A:398:PHE:CZ	1:A:664:ALA:HB2	2.52	0.44
1:A:440:LYS:HB2	1:A:440:LYS:HZ3	1.82	0.44
1:A:394:GLN:OE1	1:A:658:VAL:CG1	2.65	0.44
1:B:293:HIS:O	1:B:296:LEU:CD2	2.65	0.44
1:B:645:ILE:HG13	1:B:664:ALA:HB2	2.00	0.44
1:B:489:LEU:CG	1:B:626:LYS:HE2	2.45	0.44
1:B:484:ILE:HG22	1:B:485:TYR:N	2.33	0.44
1:A:353:CYS:HB2	1:A:398:PHE:CZ	2.53	0.44
1:B:534:THR:O	1:B:537:MET:HG2	2.18	0.44
1:A:359:LEU:HD21	1:A:362:GLU:HB2	1.99	0.43
1:A:374:ASP:C	1:A:376:SER:H	2.22	0.43
1:B:524:ASP:CB	1:B:527:HIS:HB3	2.48	0.43
1:A:539:ARG:HG2	1:A:539:ARG:O	2.17	0.43
1:B:609:VAL:HG21	1:B:628:THR:HG21	2.00	0.43
1:A:383:ILE:C	1:A:383:ILE:HD12	2.38	0.43
1:A:665:ALA:HA	1:A:668:ASN:HD21	1.83	0.43
1:A:380:LEU:CD1	1:A:394:GLN:HB3	2.49	0.43
1:B:638:ASN:O	1:B:638:ASN:OD1	2.36	0.43
1:A:476:GLU:OE2	1:A:478:LYS:HE2	2.18	0.43
1:B:530:HIS:C	1:B:532:MET:N	2.71	0.43
1:B:352:PHE:CZ	1:B:414:MET:HG2	2.54	0.43
1:A:461:LEU:HD21	1:A:576:ILE:HD12	2.01	0.43
1:B:465:ILE:HD11	1:B:477:ILE:HD11	1.99	0.43
1:B:487:GLU:OE2	1:B:625:SER:HA	2.18	0.43
1:A:380:LEU:N	1:A:380:LEU:HD12	2.32	0.43
1:B:354:ARG:HG3	1:B:354:ARG:HH11	1.83	0.43
1:A:407:SER:O	1:A:410:ASP:HB2	2.19	0.43
1:B:635:LEU:CD1	1:B:636:GLY:H	2.27	0.43
1:A:386:GLN:OE1	1:A:386:GLN:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:629:HIS:CD2	1:B:632:MET:CE	3.01	0.43
1:A:444:MET:HE1	1:A:578:LEU:HB3	2.00	0.43
1:A:550:SER:C	1:A:552:ARG:N	2.72	0.43
1:A:356:ARG:HD3	1:A:438:SER:O	2.18	0.42
1:B:382:SER:N	1:B:654:PHE:CE2	2.87	0.42
1:A:403:HIS:HD2	1:A:405:LEU:HB2	1.84	0.42
1:B:415:VAL:O	1:B:416:SER:C	2.58	0.42
1:B:379:GLU:HB2	1:B:395:ILE:HB	2.01	0.42
1:B:442:TYR:N	3:B:999:ADP:O2A	2.47	0.42
1:A:357:PRO:HB3	1:A:404:PRO:HB3	2.01	0.42
1:B:522:VAL:HA	1:B:527:HIS:CD2	2.54	0.42
1:B:487:GLU:OE2	1:B:626:LYS:N	2.50	0.42
1:A:373:HIS:O	1:A:374:ASP:HB3	2.19	0.42
1:B:336:LYS:NZ	1:B:413:GLU:HG2	2.35	0.42
1:B:632:MET:O	1:B:632:MET:HG3	2.19	0.42
1:A:383:ILE:HG13	1:A:651:GLN:OE1	2.20	0.42
1:A:478:LYS:HD3	1:A:562:ILE:HD12	2.01	0.42
1:B:367:CYS:HA	1:B:650:PHE:HA	2.00	0.42
1:B:613:LEU:HD13	1:B:635:LEU:HD12	2.02	0.42
1:A:419:ILE:HD12	1:A:576:ILE:HD13	2.02	0.42
1:B:350:ARG:HH11	1:B:350:ARG:HG2	1.85	0.42
1:B:605:GLU:HG2	1:B:624:ASN:HB2	2.02	0.42
1:B:554:HIS:CD2	1:B:627:LEU:HD22	2.55	0.42
1:B:311:LEU:O	1:B:314:ASN:HB3	2.20	0.42
1:B:577:ASN:ND2	1:B:634:SER:HB3	2.34	0.42
1:B:442:TYR:CD1	1:B:442:TYR:C	2.93	0.42
1:B:493:LEU:CD1	1:B:518:THR:HG22	2.49	0.42
1:B:493:LEU:HB2	1:B:518:THR:HG21	2.01	0.42
1:A:497:GLN:NE2	1:A:498:LYS:H	2.13	0.42
1:B:432:ALA:HB1	1:B:440:LYS:HG2	2.01	0.42
1:B:665:ALA:O	1:B:668:ASN:HB3	2.20	0.42
1:A:437:GLY:H	3:A:998:ADP:PB	2.43	0.42
1:B:633:PRO:CG	1:B:634:SER:H	2.33	0.41
1:A:524:ASP:OD1	1:A:526:ASN:HB2	2.19	0.41
1:A:299:GLU:OE2	1:B:304:ARG:NH2	2.48	0.41
1:B:611:LEU:O	1:B:615:GLN:HG3	2.19	0.41
1:A:317:GLN:O	1:A:321:LEU:HB2	2.20	0.41
1:B:492:LEU:HD21	1:B:539:ARG:HD3	2.02	0.41
1:B:482:LEU:HD11	1:B:630:LEU:HD21	2.01	0.41
1:B:354:ARG:HG3	1:B:354:ARG:NH1	2.36	0.41
1:B:350:ARG:NH1	1:B:350:ARG:HG2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:577:ASN:O	1:B:579:VAL:HG23	2.20	0.41
1:A:457:THR:HG22	1:A:461:LEU:CD2	2.51	0.41
1:B:356:ARG:HG2	1:B:438:SER:O	2.19	0.41
1:A:359:LEU:HD23	1:A:362:GLU:N	2.27	0.41
1:A:505:ALA:O	1:A:506:LYS:C	2.59	0.41
1:A:458:VAL:HG13	1:A:462:PHE:CE2	2.56	0.41
1:B:465:ILE:HG23	1:B:475:TYR:HD2	1.86	0.41
1:B:552:ARG:HA	1:B:582:ALA:HB1	2.02	0.41
1:B:552:ARG:CG	1:B:600:LYS:HD3	2.51	0.41
1:A:432:ALA:O	1:A:440:LYS:HD2	2.20	0.41
1:B:600:LYS:N	1:B:600:LYS:HD2	2.36	0.41
1:B:522:VAL:HA	1:B:527:HIS:HE2	1.85	0.41
1:A:588:LYS:HB3	1:A:589:THR:H	1.64	0.41
1:B:442:TYR:O	1:B:442:TYR:CD1	2.74	0.41
1:B:447:VAL:HG23	1:B:450:SER:N	2.22	0.41
1:B:305:GLN:O	1:B:309:GLU:HB2	2.21	0.41
1:A:502:ILE:CG2	1:A:512:ILE:HG13	2.51	0.41
1:B:537:MET:C	1:B:539:ARG:N	2.74	0.41
1:A:480:THR:CG2	4:A:6154:HOH:O	2.69	0.41
1:A:355:ILE:HG12	1:A:355:ILE:H	1.76	0.41
1:A:411:ILE:HD12	1:A:411:ILE:H	1.84	0.41
1:B:462:PHE:O	1:B:465:ILE:N	2.55	0.40
1:A:483:GLU:HG2	1:A:484:ILE:N	2.37	0.40
1:B:605:GLU:CG	1:B:624:ASN:HB2	2.52	0.40
1:B:441:THR:HB	3:B:999:ADP:O2A	2.20	0.40
1:B:360:GLU:H	1:B:360:GLU:HG2	1.59	0.40
1:B:489:LEU:HD12	1:B:489:LEU:H	1.87	0.40
1:B:380:LEU:HD22	1:B:661:LEU:HD11	2.04	0.40
1:A:374:ASP:O	1:A:376:SER:N	2.55	0.40
1:B:662:ARG:HG2	1:B:662:ARG:NH1	2.36	0.40
1:A:360:GLU:N	1:A:360:GLU:CD	2.74	0.40
1:B:665:ALA:O	1:B:668:ASN:N	2.55	0.40
1:A:522:VAL:HG21	1:A:528:LEU:HG	2.03	0.40
1:B:359:LEU:C	1:B:361:SER:H	2.24	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	354/409 (87%)	309 (87%)	33 (9%)	12 (3%)	5 6
1	B	323/409 (79%)	263 (81%)	46 (14%)	14 (4%)	3 4
All	All	677/818 (83%)	572 (84%)	79 (12%)	26 (4%)	4 5

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	506	LYS
1	B	392	GLY
1	B	458	VAL
1	B	632	MET
1	B	633	PRO
1	A	385	ALA
1	A	508	ASN
1	A	551	SER
1	A	366	MET
1	A	368	CYS
1	A	374	ASP
1	B	360	GLU
1	B	572	SER
1	B	616	LYS
1	A	375	GLU
1	A	581	LEU
1	B	463	ASP
1	B	600	LYS
1	A	639	SER
1	A	597	LYS
1	B	457	THR
1	A	392	GLY
1	B	571	ILE
1	B	455	PRO
1	B	416	SER

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Mol	Chain	Res	Type
1	B	439	GLY

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	330/370 (89%)	296 (90%)	34 (10%)	9 17
1	B	301/370 (81%)	262 (87%)	39 (13%)	5 9
All	All	631/740 (85%)	558 (88%)	73 (12%)	7 13

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	303	LEU
1	A	306	ARG
1	A	309	GLU
1	A	321	LEU
1	A	338	LEU
1	A	345	LEU
1	A	349	ILE
1	A	359	LEU
1	A	364	ASN
1	A	366	MET
1	A	380	LEU
1	A	405	LEU
1	A	416	SER
1	A	461	LEU
1	A	463	ASP
1	A	480	THR
1	A	496	GLU
1	A	497	GLN
1	A	501	GLU
1	A	509	LYS
1	A	512	ILE
1	A	528	LEU

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Mol	Chain	Res	Type
1	A	536	LYS
1	A	537	MET
1	A	572	SER
1	A	601	ARG
1	A	603	LEU
1	A	613	LEU
1	A	623	ARG
1	A	628	THR
1	A	640	LYS
1	A	642	LEU
1	A	668	ASN
1	A	669	SER
1	B	296	LEU
1	B	306	ARG
1	B	312	ARG
1	B	315	GLU
1	B	322	GLU
1	B	324	CYS
1	B	326	GLU
1	B	334	GLU
1	B	345	LEU
1	B	346	ARG
1	B	350	ARG
1	B	354	ARG
1	B	359	LEU
1	B	368	CYS
1	B	393	GLN
1	B	414	MET
1	B	440	LYS
1	B	471	LEU
1	B	482	LEU
1	B	489	LEU
1	B	496	GLU
1	B	527	HIS
1	B	531	LEU
1	B	539	ARG
1	B	553	SER
1	B	560	GLU
1	B	562	ILE
1	B	564	ARG
1	B	585	GLU
1	B	598	ASN

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Mol	Chain	Res	Type
1	B	600	LYS
1	B	606	LEU
1	B	631	LEU
1	B	633	PRO
1	B	635	LEU
1	B	638	ASN
1	B	641	THR
1	B	659	LYS
1	B	662	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	317	GLN
1	A	364	ASN
1	A	393	GLN
1	A	403	HIS
1	A	486	ASN
1	A	495	ASN
1	A	497	GLN
1	A	510	ASN
1	A	516	ASN
1	A	527	HIS
1	A	598	ASN
1	A	608	ASN
1	A	629	HIS
1	A	668	ASN
1	B	293	HIS
1	B	305	GLN
1	B	317	GLN
1	B	327	GLN
1	B	348	ASN
1	B	364	ASN
1	B	393	GLN
1	B	403	HIS
1	B	427	ASN
1	B	486	ASN
1	B	495	ASN
1	B	527	HIS
1	B	577	ASN
1	B	598	ASN
1	B	615	GLN

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Mol	Chain	Res	Type
1	B	629	HIS
1	B	655	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	ADP	A	998	2	22,29,29	1.52	5 (22%)	27,45,45	2.78	4 (14%)
3	ADP	B	999	2	22,29,29	1.55	4 (18%)	27,45,45	2.86	5 (18%)

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	ADP	A	998	2	-	0/12/32/32	0/3/3/3
3	ADP	B	999	2	-	0/12/32/32	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	998	ADP	C5-N7	-2.45	1.31	1.39
3	B	999	ADP	PA-O1A	-2.13	1.43	1.51
3	A	998	ADP	PA-O2A	-2.05	1.46	1.54
3	A	998	ADP	PB-O2B	2.00	1.61	1.54
3	A	998	ADP	C2-N3	2.52	1.36	1.32
3	B	999	ADP	C2-N3	2.54	1.36	1.32
3	B	999	ADP	PB-O2B	2.62	1.64	1.54
3	A	998	ADP	O4'-C1'	3.98	1.46	1.41
3	B	999	ADP	O4'-C1'	4.13	1.46	1.41

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	999	ADP	N3-C2-N1	-13.46	118.59	128.89
3	A	998	ADP	N3-C2-N1	-12.88	119.03	128.89
3	B	999	ADP	C4-C5-N7	-2.91	106.81	109.48
3	A	998	ADP	PA-O3A-PB	-2.69	123.64	132.67
3	A	998	ADP	C4-C5-N7	-2.38	107.29	109.48
3	B	999	ADP	C2'-C1'-N9	-2.04	111.18	114.29
3	B	999	ADP	O4'-C1'-N9	2.27	112.85	108.10
3	B	999	ADP	O3A-PA-O5'	3.04	111.01	102.94
3	A	998	ADP	O3A-PA-O5'	4.13	113.91	102.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	998	ADP	6	0
3	B	999	ADP	6	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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.