Full wwPDB X-ray Structure Validation Report (i)

Jun 21, 2021 – 12:03 PM BST

PDB ID	:	6Y1I
Title	:	Human Eg5 motor domain mutant R234C
Authors	:	Garcia-Saez, I.; Skoufias, D.A.
Deposited on	:	2020-02-12
Resolution	:	3.00 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 (i) symbol.

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

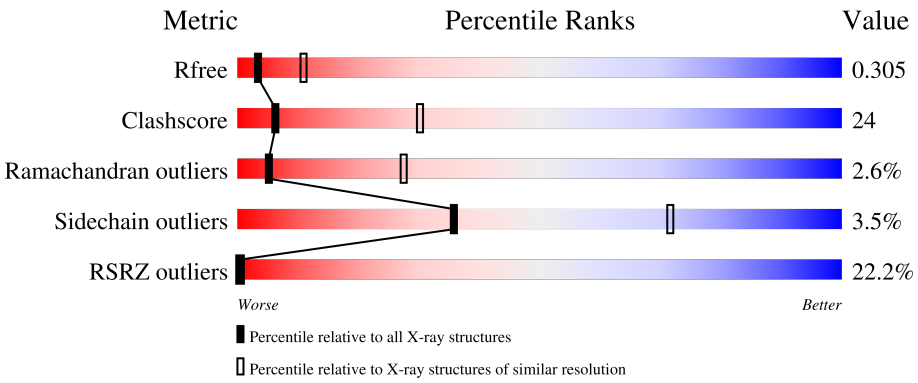
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.20
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.20

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 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 <=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	368	2%, 52%, 32%, •, 11%
1	B	368	5%, 51%, 32%, •, 13%
1	C	368	51%, 53%, 33%, •, 11%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7727 atoms, of which 36 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.

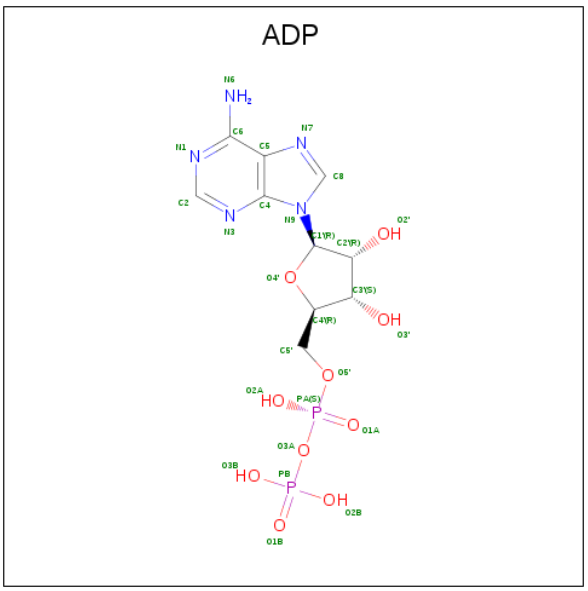
- Molecule 1 is a protein called Kinesin-like protein KIF11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	326	2552	1601	445	495	11	0	0	0
1	B	319	2503	1573	434	485	11	0	0	0
1	C	326	2552	1601	445	495	11	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	234	CYS	ARG	engineered mutation	UNP P52732
B	234	CYS	ARG	engineered mutation	UNP P52732
C	234	CYS	ARG	engineered mutation	UNP P52732

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	A	1	39	10	12	5	10	2	0	0
2	B	1	39	10	12	5	10	2	0	0
2	C	1	39	10	12	5	10	2	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

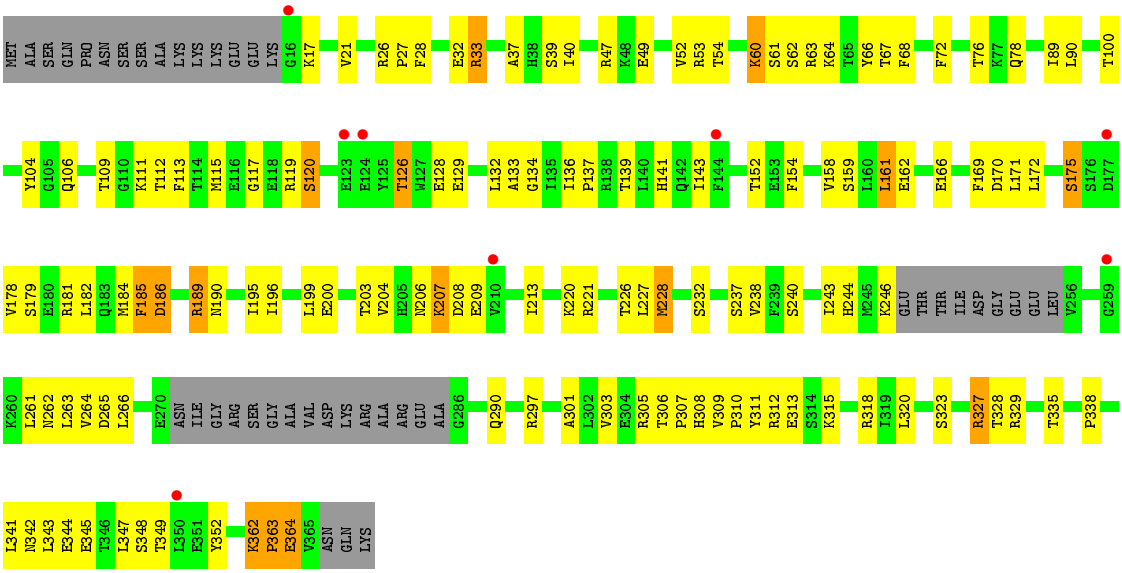
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	1	1	1	0	0
3	B	1	1	1	0	0
3	C	1	1	1	0	0

3 Residue-property plots (i)

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 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.

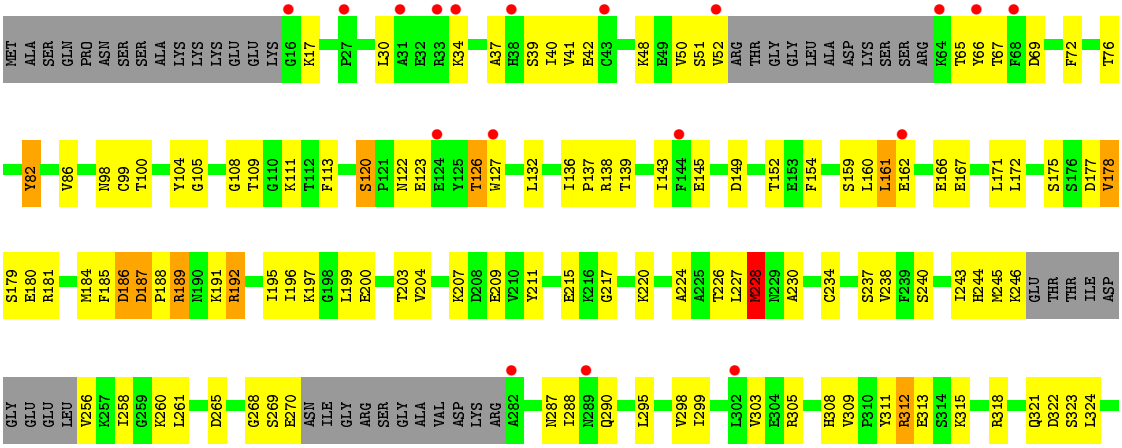
- Molecule 1: Kinesin-like protein KIF11

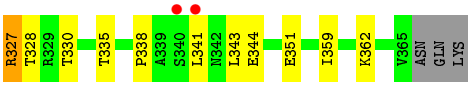
Chain A:



- Molecule 1: Kinesin-like protein KIF11

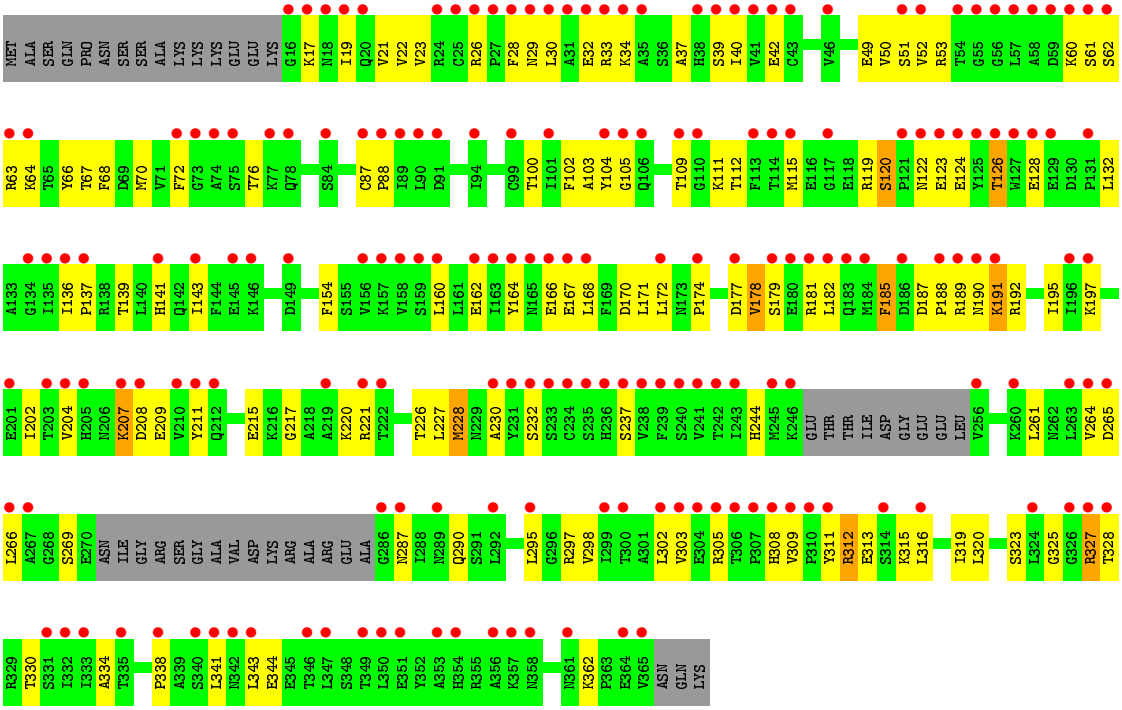
Chain B:





- Molecule 1: Kinesin-like protein KIF11

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	87.43Å 87.60Å 95.06Å 90.00° 116.82° 90.00°	Depositor
Resolution (Å)	47.70 – 3.00 47.70 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (47.70-3.00) 97.0 (47.70-3.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.75 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.272 , 0.306 0.274 , 0.305	Depositor DCC
R_{free} test set	1269 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	77.9	Xtriage
Anisotropy	0.855	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 104.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.227 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7727	wwPDB-VP
Average B, all atoms (Å^2)	194.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.40% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.29	0/2590	0.49	0/3500
1	B	0.26	0/2540	0.45	0/3433
1	C	0.24	0/2590	0.43	0/3500
All	All	0.26	0/7720	0.46	0/10433

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2552	0	2582	148	0
1	B	2503	0	2528	128	0
1	C	2552	0	2582	112	0
2	A	27	12	12	2	0
2	B	27	12	12	3	0
2	C	27	12	12	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
All	All	7691	36	7728	374	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 24.

All (374) 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:160:LEU:HB3	1:B:172:LEU:HD21	1.15	1.12
1:A:203:THR:HB	1:B:175:SER:HB2	1.32	1.08
1:B:160:LEU:HB3	1:B:172:LEU:CD2	1.92	0.99
1:A:40:ILE:HD12	1:A:343:LEU:HD23	1.45	0.99
1:C:37:ALA:HB2	1:C:341:LEU:HD13	1.48	0.94
1:C:40:ILE:HD12	1:C:343:LEU:HD23	1.46	0.94
1:B:177:ASP:HA	1:B:220:LYS:HE3	1.51	0.90
1:B:67:THR:HG22	1:B:359:ILE:HD11	1.51	0.90
1:B:82:TYR:CD1	1:B:86:VAL:HB	2.08	0.89
1:B:111:LYS:HE2	2:B:401:ADP:O2B	1.74	0.88
1:C:112:THR:HB	2:C:401:ADP:O1A	1.73	0.87
1:C:160:LEU:HB3	1:C:172:LEU:HD13	1.54	0.86
1:A:60:LYS:HD3	1:A:61:SER:H	1.41	0.86
1:A:362:LYS:H	1:A:362:LYS:HD3	1.42	0.85
1:A:26:ARG:HG3	1:A:27:PRO:HD2	1.59	0.85
1:A:309:VAL:HG13	1:A:311:TYR:CE2	2.12	0.84
1:A:162:GLU:CB	1:A:171:LEU:HD11	2.07	0.84
1:A:60:LYS:HD3	1:A:61:SER:N	1.92	0.83
1:A:120:SER:HB3	1:A:132:LEU:HD12	1.57	0.83
1:A:162:GLU:HB2	1:A:171:LEU:HD11	1.60	0.82
1:B:17:LYS:HE3	1:B:303:VAL:HA	1.60	0.82
1:B:160:LEU:CB	1:B:172:LEU:HD21	2.05	0.80
1:A:309:VAL:CG1	1:A:311:TYR:CE2	2.66	0.79
1:B:298:VAL:HG13	1:B:309:VAL:HG21	1.65	0.79
1:B:105:GLY:O	1:B:111:LYS:NZ	2.16	0.78
1:A:362:LYS:HB2	1:A:363:PRO:HD2	1.65	0.77
1:B:98:ASN:OD1	1:B:323:SER:HA	1.84	0.77
1:B:167:GLU:OE1	1:B:181:ARG:HD3	1.85	0.77
1:A:189:ARG:NH1	1:A:190:ASN:HB2	2.01	0.76
1:A:89:ILE:HD13	1:A:329:ARG:HB2	1.67	0.75
1:B:256:VAL:HG21	1:C:190:ASN:HA	1.69	0.75
1:A:17:LYS:HE3	1:A:303:VAL:HA	1.69	0.75
1:C:178:VAL:HG12	1:C:179:SER:H	1.52	0.75
1:C:141:HIS:HD2	1:C:207:LYS:HD3	1.52	0.74
1:A:53:ARG:HA	1:A:62:SER:HB2	1.68	0.74
1:B:40:ILE:CD1	1:B:343:LEU:HD23	2.17	0.74
1:B:40:ILE:HD12	1:B:343:LEU:HD23	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:PHE:HB3	1:A:76:THR:HG21	1.71	0.73
1:B:41:VAL:CG2	1:B:338:PRO:HA	2.18	0.73
1:C:120:SER:HB3	1:C:132:LEU:HD12	1.71	0.73
1:A:169:PHE:CD2	1:A:179:SER:HB2	2.24	0.73
1:C:49:GLU:HG2	1:C:67:THR:HG22	1.69	0.72
1:A:185:PHE:HZ	1:B:195:ILE:HD13	1.54	0.72
1:C:170:ASP:HB2	1:C:182:LEU:HD11	1.70	0.72
1:C:17:LYS:HE3	1:C:303:VAL:HA	1.70	0.72
1:C:192:ARG:NH1	1:C:325:GLY:HA3	2.04	0.72
1:A:109:THR:HB	1:A:335:THR:HG22	1.69	0.72
1:A:169:PHE:CE2	1:A:179:SER:HB2	2.25	0.71
1:A:209:GLU:O	1:A:213:ILE:HD13	1.89	0.71
1:A:115:MET:CE	1:A:263:LEU:HB3	2.20	0.71
1:C:111:LYS:HE3	1:C:266:LEU:O	1.91	0.71
1:B:41:VAL:HG21	1:B:338:PRO:HA	1.72	0.71
1:A:54:THR:HG23	1:A:62:SER:HB3	1.74	0.70
1:B:120:SER:HB3	1:B:132:LEU:HD12	1.73	0.70
1:B:178:VAL:HG12	1:B:179:SER:H	1.57	0.70
1:B:166:GLU:HB2	1:B:287:ASN:OD1	1.91	0.70
1:B:189:ARG:HG2	1:C:185:PHE:CE1	2.27	0.70
1:C:177:ASP:OD2	1:C:220:LYS:HE3	1.92	0.69
1:A:184:MET:CE	1:A:318:ARG:HD3	2.22	0.69
1:A:26:ARG:CG	1:A:27:PRO:HD2	2.23	0.68
1:A:349:THR:HA	1:A:352:TYR:HD2	1.58	0.68
1:C:49:GLU:HG2	1:C:67:THR:CG2	2.23	0.68
1:A:310:PRO:HB2	1:A:313:GLU:HG3	1.74	0.68
1:C:204:VAL:HG13	1:C:209:GLU:HB3	1.75	0.68
1:A:40:ILE:CD1	1:A:343:LEU:HD23	2.22	0.67
1:C:40:ILE:CD1	1:C:343:LEU:HD23	2.22	0.67
1:B:226:THR:OG1	1:B:227:LEU:HD22	1.95	0.67
1:A:344:GLU:N	1:A:344:GLU:OE1	2.27	0.66
1:A:49:GLU:HG2	1:A:67:THR:CG2	2.25	0.66
1:C:139:THR:HG21	1:C:261:LEU:HD23	1.75	0.66
1:C:178:VAL:HG12	1:C:179:SER:N	2.10	0.66
1:A:301:ALA:HA	1:A:306:THR:CG2	2.26	0.66
1:B:100:THR:HG21	1:B:324:LEU:HD21	1.76	0.66
1:B:166:GLU:HG2	1:B:315:LYS:HG3	1.78	0.65
1:C:298:VAL:HA	1:C:309:VAL:CG2	2.25	0.65
1:A:89:ILE:HD13	1:A:329:ARG:CB	2.26	0.65
1:A:185:PHE:CZ	1:B:195:ILE:HD13	2.32	0.65
1:A:184:MET:HE3	1:A:318:ARG:HD3	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:THR:HG1	1:C:323:SER:HG	1.44	0.65
1:C:237:SER:HB3	1:C:265:ASP:HB3	1.78	0.65
1:B:40:ILE:O	1:B:52:VAL:HA	1.97	0.65
1:A:49:GLU:HG2	1:A:67:THR:HG22	1.79	0.64
1:B:67:THR:HG22	1:B:359:ILE:CD1	2.27	0.64
1:A:175:SER:HB2	1:B:203:THR:HB	1.79	0.64
1:B:177:ASP:HA	1:B:220:LYS:CE	2.25	0.64
1:B:344:GLU:OE1	1:B:344:GLU:N	2.29	0.64
1:C:103:ALA:O	1:C:111:LYS:HD2	1.98	0.64
1:B:143:ILE:HD13	1:B:243:ILE:HD11	1.79	0.63
1:A:237:SER:HB3	1:A:265:ASP:HB3	1.81	0.63
1:B:109:THR:HG1	1:B:335:THR:HG1	1.45	0.63
1:B:237:SER:HB3	1:B:265:ASP:HB3	1.81	0.62
1:A:226:THR:OG1	1:A:227:LEU:HD22	1.98	0.62
1:C:39:SER:HA	1:C:338:PRO:O	1.99	0.62
1:B:184:MET:CE	1:B:318:ARG:HD3	2.29	0.62
1:A:362:LYS:H	1:A:362:LYS:CD	2.11	0.62
1:A:204:VAL:HG13	1:A:209:GLU:HB3	1.80	0.62
1:C:166:GLU:HG2	1:C:315:LYS:HG3	1.82	0.61
1:A:189:ARG:HD2	1:A:189:ARG:O	1.99	0.61
1:B:298:VAL:HA	1:B:309:VAL:HG23	1.83	0.61
1:C:40:ILE:O	1:C:52:VAL:HA	2.00	0.61
1:B:39:SER:HA	1:B:338:PRO:O	2.00	0.61
1:C:226:THR:OG1	1:C:227:LEU:HD22	2.00	0.61
1:A:78:GLN:NE2	1:A:113:PHE:O	2.29	0.60
1:B:162:GLU:CB	1:B:171:LEU:HD11	2.30	0.60
1:A:221:ARG:HD3	1:A:232:SER:HB3	1.83	0.60
1:C:312:ARG:HD3	1:C:312:ARG:N	2.16	0.60
1:B:37:ALA:HB2	1:B:341:LEU:HD13	1.82	0.60
1:B:312:ARG:HD3	1:B:312:ARG:N	2.16	0.60
1:C:189:ARG:HG3	1:C:190:ASN:N	2.16	0.60
1:A:137:PRO:O	1:A:141:HIS:ND1	2.23	0.60
1:C:167:GLU:HB3	1:C:181:ARG:HG3	1.83	0.59
1:A:106:GLN:HE22	1:A:342:ASN:HB3	1.66	0.59
1:C:221:ARG:HD3	1:C:232:SER:HB3	1.85	0.59
1:A:196:ILE:HG22	1:A:199:LEU:HB2	1.85	0.59
1:B:327:ARG:HA	1:B:327:ARG:NE	2.18	0.59
1:C:162:GLU:HB2	1:C:171:LEU:HD11	1.84	0.59
1:A:39:SER:HA	1:A:338:PRO:O	2.03	0.58
1:A:308:HIS:O	1:A:308:HIS:CG	2.55	0.58
1:A:323:SER:O	1:A:328:THR:HG21	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:THR:HG21	1:C:362:LYS:NZ	2.18	0.58
1:C:344:GLU:OE1	1:C:344:GLU:N	2.34	0.58
1:A:115:MET:HE1	1:A:263:LEU:HB3	1.84	0.58
1:A:143:ILE:HD12	1:A:243:ILE:CD1	2.33	0.58
1:B:171:LEU:O	1:B:217:GLY:HA2	2.03	0.58
1:C:168:LEU:HD13	1:C:181:ARG:NH2	2.18	0.58
1:B:139:THR:HG21	1:B:261:LEU:HD23	1.85	0.58
1:C:295:LEU:HD22	1:C:316:LEU:CD2	2.33	0.58
1:B:246:LYS:HG3	1:B:256:VAL:HG22	1.85	0.58
1:C:136:ILE:HB	1:C:137:PRO:HD3	1.84	0.58
1:B:82:TYR:HD1	1:B:86:VAL:HB	1.67	0.58
1:C:181:ARG:HD3	1:C:182:LEU:O	2.04	0.58
1:C:141:HIS:CD2	1:C:207:LYS:HD3	2.35	0.58
1:C:168:LEU:HD21	1:C:319:ILE:HD11	1.85	0.57
1:A:162:GLU:HB3	1:A:171:LEU:HD11	1.86	0.57
1:A:171:LEU:HA	1:A:220:LYS:HG2	1.87	0.57
1:B:234:CYS:SG	1:B:288:ILE:HD11	2.45	0.57
1:B:177:ASP:CA	1:B:220:LYS:HE3	2.29	0.57
1:C:295:LEU:HD22	1:C:316:LEU:HD21	1.87	0.56
1:A:111:LYS:NZ	2:A:401:ADP:O2B	2.30	0.56
1:C:119:ARG:O	1:C:120:SER:OG	2.20	0.56
1:C:323:SER:O	1:C:328:THR:HG21	2.04	0.56
1:A:60:LYS:CD	1:A:61:SER:N	2.67	0.56
1:B:82:TYR:CE1	1:B:86:VAL:HB	2.41	0.56
1:B:228:MET:O	1:B:228:MET:HG2	2.05	0.56
1:B:98:ASN:HA	1:B:260:LYS:O	2.05	0.55
1:B:98:ASN:HB2	1:B:260:LYS:HB3	1.88	0.55
1:C:327:ARG:NE	1:C:327:ARG:HA	2.21	0.55
1:A:170:ASP:HB2	1:A:182:LEU:HD11	1.88	0.55
1:A:327:ARG:NE	1:A:327:ARG:HA	2.21	0.55
1:B:187:ASP:OD1	1:B:189:ARG:HB3	2.06	0.55
1:C:211:TYR:O	1:C:215:GLU:HG3	2.07	0.55
1:B:30:LEU:HG	1:B:34:LYS:HD2	1.89	0.55
1:B:298:VAL:HG13	1:B:309:VAL:CG2	2.37	0.54
1:B:309:VAL:HG13	1:B:311:TYR:CE2	2.41	0.54
1:A:364:GLU:O	1:A:364:GLU:HG3	2.08	0.54
1:A:143:ILE:HD12	1:A:243:ILE:HD11	1.89	0.54
1:B:82:TYR:CE2	1:B:138:ARG:HB3	2.42	0.54
1:B:184:MET:HE1	1:B:318:ARG:HD3	1.90	0.54
1:A:228:MET:HG2	1:A:228:MET:O	2.07	0.54
1:C:29:ASN:O	1:C:33:ARG:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:187:ASP:CB	1:C:195:ILE:HD12	2.38	0.53
1:C:181:ARG:HH22	1:C:315:LYS:HD2	1.73	0.53
1:A:115:MET:HE3	1:A:263:LEU:HB3	1.90	0.53
1:A:309:VAL:HG11	1:A:311:TYR:CE2	2.41	0.53
1:C:171:LEU:O	1:C:217:GLY:HA2	2.09	0.53
1:B:98:ASN:ND2	1:B:322:ASP:O	2.41	0.53
1:C:72:PHE:HB3	1:C:76:THR:HG21	1.90	0.53
1:A:195:ILE:HD11	1:B:197:LYS:CD	2.39	0.53
1:A:28:PHE:HB2	1:A:33:ARG:HH12	1.73	0.53
1:B:188:PRO:O	1:B:189:ARG:HB2	2.08	0.52
1:B:298:VAL:HA	1:B:309:VAL:CG2	2.39	0.52
1:A:306:THR:OG1	1:A:307:PRO:HD2	2.08	0.52
1:B:145:GLU:O	1:B:149:ASP:HB2	2.09	0.52
1:A:90:LEU:HD11	1:A:143:ILE:HD11	1.91	0.52
1:C:298:VAL:HA	1:C:309:VAL:HG23	1.91	0.52
1:C:167:GLU:HA	1:C:181:ARG:HE	1.73	0.52
1:A:40:ILE:O	1:A:52:VAL:HA	2.10	0.52
1:B:211:TYR:O	1:B:215:GLU:HG3	2.10	0.52
1:A:175:SER:CB	1:B:203:THR:HB	2.41	0.51
1:A:90:LEU:HD11	1:A:143:ILE:CD1	2.41	0.51
1:A:106:GLN:NE2	1:A:345:GLU:HB2	2.26	0.51
1:A:308:HIS:O	1:A:308:HIS:ND1	2.42	0.51
1:B:82:TYR:HD1	1:B:82:TYR:O	1.93	0.51
1:B:42:GLU:O	1:B:50:VAL:HA	2.11	0.51
1:A:204:VAL:CG1	1:A:209:GLU:HB3	2.40	0.51
1:B:179:SER:HB2	1:B:228:MET:CE	2.41	0.51
1:A:185:PHE:CZ	1:B:195:ILE:CD1	2.94	0.51
1:B:100:THR:HG23	1:B:330:THR:HG23	1.93	0.51
1:B:245:MET:O	1:B:256:VAL:HA	2.11	0.51
1:C:139:THR:O	1:C:143:ILE:HG13	2.11	0.51
1:C:308:HIS:O	1:C:308:HIS:ND1	2.44	0.51
1:A:139:THR:HG21	1:A:261:LEU:HD23	1.92	0.51
1:A:301:ALA:HB1	1:A:306:THR:HG23	1.93	0.51
1:B:108:GLY:H	2:B:401:ADP:PB	2.33	0.50
1:B:256:VAL:HG21	1:C:190:ASN:CA	2.38	0.50
1:A:196:ILE:CG2	1:A:199:LEU:HB2	2.41	0.50
1:B:162:GLU:HB2	1:B:171:LEU:HD11	1.92	0.50
1:B:104:TYR:CZ	1:B:269:SER:HB3	2.47	0.50
1:A:52:VAL:HG12	1:A:53:ARG:N	2.27	0.50
1:B:111:LYS:HZ1	1:B:268:GLY:HA2	1.77	0.50
1:A:17:LYS:HD2	1:A:303:VAL:HG13	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:52:VAL:HG12	1:C:53:ARG:N	2.28	0.49
1:A:301:ALA:HA	1:A:306:THR:HG22	1.94	0.49
1:A:362:LYS:HD3	1:A:362:LYS:N	2.21	0.49
1:B:186:ASP:O	1:B:188:PRO:HD3	2.11	0.49
1:B:72:PHE:HB3	1:B:76:THR:HG21	1.94	0.49
1:B:136:ILE:HB	1:B:137:PRO:HD3	1.93	0.49
1:A:166:GLU:HG2	1:A:315:LYS:HG3	1.95	0.49
1:C:154:PHE:HA	1:C:244:HIS:O	2.13	0.49
1:C:187:ASP:HB2	1:C:195:ILE:HD12	1.94	0.49
1:B:111:LYS:HE2	2:B:401:ADP:PB	2.52	0.48
1:C:309:VAL:HG13	1:C:311:TYR:CE2	2.48	0.48
1:A:47:ARG:O	1:A:362:LYS:NZ	2.41	0.48
1:B:204:VAL:HG13	1:B:209:GLU:HB3	1.95	0.48
1:A:264:VAL:HG21	1:A:320:LEU:HD11	1.95	0.48
1:B:41:VAL:HG23	1:B:338:PRO:HA	1.95	0.48
1:C:17:LYS:NZ	1:C:17:LYS:HB3	2.28	0.48
1:A:227:LEU:O	1:A:228:MET:HB2	2.13	0.48
1:A:17:LYS:CD	1:A:303:VAL:HG13	2.44	0.48
1:A:309:VAL:HG13	1:A:309:VAL:O	2.13	0.48
1:C:17:LYS:CD	1:C:303:VAL:HG13	2.43	0.48
1:A:112:THR:HB	2:A:401:ADP:O2A	2.13	0.48
1:A:141:HIS:HD2	1:A:207:LYS:HD3	1.77	0.48
1:A:169:PHE:HD2	1:A:179:SER:HB2	1.77	0.48
1:B:17:LYS:HB3	1:B:17:LYS:NZ	2.28	0.48
1:B:17:LYS:HD2	1:B:303:VAL:HG13	1.95	0.48
1:C:115:MET:O	1:C:136:ILE:HG13	2.14	0.48
1:A:17:LYS:NZ	1:A:17:LYS:HB3	2.29	0.48
1:A:37:ALA:HB2	1:A:341:LEU:HD13	1.96	0.47
1:A:309:VAL:HG13	1:A:311:TYR:CD2	2.48	0.47
1:A:28:PHE:HB2	1:A:33:ARG:NH1	2.29	0.47
1:A:100:THR:HA	1:A:262:ASN:O	2.14	0.47
1:A:348:SER:O	1:A:352:TYR:CD2	2.66	0.47
1:C:102:PHE:CE1	1:C:295:LEU:HD21	2.49	0.47
1:A:169:PHE:HE2	1:A:179:SER:HB2	1.78	0.47
1:A:170:ASP:O	1:A:220:LYS:NZ	2.48	0.47
1:C:30:LEU:HG	1:C:34:LYS:HD2	1.96	0.47
1:C:308:HIS:O	1:C:308:HIS:CG	2.66	0.47
1:A:26:ARG:NH2	1:A:32:GLU:OE1	2.37	0.47
1:A:154:PHE:HA	1:A:244:HIS:O	2.14	0.47
1:B:82:TYR:CD2	1:B:138:ARG:HB3	2.49	0.47
1:B:327:ARG:HA	1:B:327:ARG:CZ	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:ARG:HH21	1:A:310:PRO:HG3	1.79	0.47
1:B:270:GLU:OE1	1:B:270:GLU:N	2.46	0.47
1:B:256:VAL:HB	1:C:190:ASN:N	2.30	0.47
1:A:139:THR:O	1:A:143:ILE:HG12	2.15	0.46
1:A:349:THR:HA	1:A:352:TYR:CD2	2.44	0.46
1:B:309:VAL:CG1	1:B:311:TYR:CE2	2.98	0.46
1:C:60:LYS:O	1:C:62:SER:N	2.49	0.46
1:C:190:ASN:O	1:C:191:LYS:O	2.33	0.46
1:A:189:ARG:HD2	1:A:189:ARG:C	2.33	0.46
1:A:26:ARG:HH22	1:A:32:GLU:CD	2.16	0.46
1:B:178:VAL:HG12	1:B:179:SER:N	2.28	0.46
1:A:301:ALA:CA	1:A:306:THR:CG2	2.92	0.46
1:B:159:SER:OG	1:B:240:SER:HB2	2.15	0.46
1:C:19:ILE:HD13	1:C:330:THR:HB	1.98	0.46
1:A:161:LEU:CD1	1:A:238:VAL:HB	2.46	0.46
1:B:312:ARG:HD3	1:B:312:ARG:H	1.79	0.46
1:C:227:LEU:O	1:C:228:MET:HB2	2.16	0.46
1:A:158:VAL:HA	1:A:240:SER:O	2.15	0.46
1:B:98:ASN:OD1	1:B:323:SER:CA	2.59	0.46
1:B:154:PHE:HA	1:B:244:HIS:O	2.15	0.46
1:B:191:LYS:O	1:B:192:ARG:HB2	2.15	0.46
1:A:126:THR:HB	1:A:128:GLU:OE1	2.16	0.46
1:A:141:HIS:CD2	1:A:207:LYS:HD3	2.51	0.46
1:A:195:ILE:HD11	1:B:197:LYS:HD2	1.98	0.46
1:A:213:ILE:HD12	1:A:213:ILE:N	2.31	0.46
1:A:159:SER:OG	1:A:240:SER:HB2	2.17	0.45
1:A:21:VAL:O	1:A:68:PHE:HB3	2.16	0.45
1:B:200:GLU:HA	1:B:200:GLU:OE1	2.16	0.45
1:C:264:VAL:HG21	1:C:320:LEU:HD11	1.97	0.45
1:B:113:PHE:CD2	1:B:113:PHE:C	2.90	0.45
1:B:152:THR:O	1:B:154:PHE:HD2	1.99	0.45
1:A:301:ALA:HA	1:A:306:THR:HG21	1.97	0.45
1:B:161:LEU:HD11	1:B:238:VAL:HB	1.99	0.45
1:B:167:GLU:OE1	1:B:181:ARG:NH1	2.40	0.45
1:C:22:VAL:CG1	1:C:70:MET:HB2	2.47	0.45
1:C:137:PRO:O	1:C:141:HIS:ND1	2.38	0.45
1:A:161:LEU:HD11	1:A:238:VAL:HB	1.98	0.45
1:C:177:ASP:CG	1:C:220:LYS:HE3	2.36	0.45
1:A:28:PHE:CB	1:A:33:ARG:NH1	2.80	0.45
1:A:26:ARG:HD3	1:A:109:THR:O	2.17	0.45
1:A:129:GLU:OE2	1:A:141:HIS:CG	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:LEU:HD12	1:A:171:LEU:N	2.32	0.45
1:B:51:SER:HB3	1:B:65:THR:OG1	2.17	0.45
1:B:258:ILE:CD1	1:C:188:PRO:HB3	2.47	0.45
1:A:49:GLU:HA	1:A:66:TYR:O	2.18	0.44
1:B:179:SER:O	1:B:179:SER:OG	2.26	0.44
1:B:187:ASP:OD2	1:B:195:ILE:HD11	2.17	0.44
1:C:21:VAL:O	1:C:68:PHE:HB3	2.17	0.44
1:C:290:GLN:NE2	1:C:313:GLU:O	2.50	0.44
1:B:161:LEU:CD1	1:B:238:VAL:HB	2.47	0.44
1:B:234:CYS:O	1:B:288:ILE:HD11	2.17	0.44
1:C:168:LEU:H	1:C:181:ARG:NE	2.14	0.44
1:A:72:PHE:HB3	1:A:76:THR:CG2	2.44	0.44
1:A:327:ARG:HA	1:A:327:ARG:CZ	2.48	0.44
1:A:26:ARG:O	1:A:338:PRO:HG3	2.17	0.44
1:A:343:LEU:O	1:A:347:LEU:HD12	2.17	0.44
1:B:122:ASN:O	1:B:123:GLU:HB2	2.18	0.44
1:B:189:ARG:HA	1:C:185:PHE:HZ	1.83	0.44
1:A:133:ALA:HB1	1:A:137:PRO:HB2	1.98	0.44
1:A:290:GLN:NE2	1:A:313:GLU:O	2.49	0.44
1:B:162:GLU:HB3	1:B:171:LEU:HD11	1.99	0.44
1:C:228:MET:O	1:C:228:MET:HG2	2.17	0.44
1:A:117:GLY:HA3	1:A:134:GLY:N	2.33	0.43
1:C:172:LEU:HD23	1:C:202:ILE:HG12	1.99	0.43
1:A:119:ARG:HA	1:A:119:ARG:HD3	1.68	0.43
1:A:172:LEU:HD12	1:A:172:LEU:HA	1.86	0.43
1:A:181:ARG:O	1:A:181:ARG:HG2	2.18	0.43
1:C:67:THR:HG21	1:C:362:LYS:HZ1	1.83	0.43
1:C:298:VAL:O	1:C:302:LEU:HG	2.18	0.43
1:C:23:VAL:HG22	1:C:334:ALA:HB3	2.01	0.43
1:C:197:LYS:NZ	1:C:197:LYS:HB2	2.33	0.43
1:A:200:GLU:OE1	1:A:200:GLU:HA	2.18	0.43
1:C:28:PHE:HB2	1:C:33:ARG:NH2	2.34	0.43
1:C:187:ASP:HB3	1:C:195:ILE:HD12	1.99	0.43
1:B:196:ILE:HG22	1:B:199:LEU:HB2	2.01	0.43
1:C:42:GLU:O	1:C:50:VAL:HA	2.19	0.43
1:C:126:THR:HB	1:C:128:GLU:OE1	2.18	0.43
1:C:166:GLU:HG2	1:C:315:LYS:CG	2.47	0.43
1:C:111:LYS:HB2	1:C:111:LYS:HE2	1.80	0.43
1:A:206:ASN:OD1	1:A:208:ASP:HB2	2.18	0.42
1:C:17:LYS:HD2	1:C:303:VAL:HG13	1.99	0.42
1:B:166:GLU:HB2	1:B:287:ASN:CG	2.39	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:28:PHE:HB2	1:C:33:ARG:CZ	2.50	0.42
1:A:186:ASP:OD1	1:A:186:ASP:N	2.53	0.42
1:A:309:VAL:CG1	1:A:311:TYR:HE2	2.26	0.42
1:B:99:CYS:HG	1:B:328:THR:HG1	1.54	0.42
1:C:164:TYR:CE2	1:C:230:ALA:HA	2.55	0.42
1:A:185:PHE:CE1	1:A:195:ILE:CG2	3.03	0.42
1:B:172:LEU:N	1:B:172:LEU:HD22	2.35	0.42
1:C:362:LYS:HE2	1:C:362:LYS:HB2	1.87	0.42
1:A:134:GLY:C	1:A:137:PRO:HD2	2.39	0.42
1:A:152:THR:HG23	1:A:246:LYS:O	2.19	0.42
1:A:221:ARG:CD	1:A:232:SER:HB3	2.50	0.42
1:B:99:CYS:SG	1:B:328:THR:OG1	2.57	0.42
1:A:136:ILE:HB	1:A:137:PRO:HD3	2.02	0.42
1:B:171:LEU:N	1:B:171:LEU:HD12	2.35	0.42
1:C:170:ASP:O	1:C:220:LYS:HE2	2.20	0.41
1:C:327:ARG:HA	1:C:327:ARG:CZ	2.50	0.41
1:C:207:LYS:HD2	1:C:208:ASP:N	2.35	0.41
1:B:227:LEU:O	1:B:228:MET:HB2	2.20	0.41
1:B:295:LEU:O	1:B:299:ILE:HG12	2.19	0.41
1:C:42:GLU:HB3	1:C:51:SER:OG	2.20	0.41
1:B:290:GLN:NE2	1:B:313:GLU:O	2.53	0.41
1:C:87:CYS:HB2	1:C:88:PRO:HD3	2.01	0.41
1:A:119:ARG:O	1:A:120:SER:OG	2.29	0.41
1:B:66:TYR:OH	1:B:351:GLU:OE2	2.30	0.41
1:C:312:ARG:HD3	1:C:312:ARG:H	1.82	0.41
1:A:72:PHE:CD1	1:A:76:THR:HG21	2.55	0.41
1:A:143:ILE:HD12	1:A:243:ILE:HD13	2.00	0.41
1:B:311:TYR:CD2	1:B:321:GLN:HG3	2.56	0.41
1:C:112:THR:CB	2:C:401:ADP:O1A	2.58	0.41
1:A:67:THR:CG2	1:A:362:LYS:HZ2	2.34	0.41
1:B:126:THR:HB	1:B:127:TRP:H	1.67	0.41
1:C:26:ARG:HD3	1:C:109:THR:O	2.21	0.41
1:C:28:PHE:CB	1:C:33:ARG:CZ	2.99	0.41
1:A:47:ARG:HE	1:A:49:GLU:CD	2.24	0.41
1:A:111:LYS:HE3	1:A:266:LEU:O	2.20	0.41
1:A:343:LEU:O	1:A:343:LEU:HD22	2.20	0.41
1:A:161:LEU:HD23	1:A:196:ILE:HD13	2.03	0.41
1:B:48:LYS:HD3	1:B:69:ASP:O	2.21	0.40
1:B:234:CYS:HG	1:B:288:ILE:HD11	1.84	0.40
1:B:323:SER:O	1:B:328:THR:HG21	2.22	0.40
1:C:49:GLU:HA	1:C:66:TYR:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:104:TYR:CZ	1:C:269:SER:HB3	2.56	0.40
1:C:362:LYS:HZ3	1:C:362:LYS:H	1.68	0.40
1:C:105:GLY:O	1:C:111:LYS:NZ	2.51	0.40
1:B:362:LYS:HE2	1:B:362:LYS:HB2	1.89	0.40
1:C:26:ARG:NH2	1:C:32:GLU:OE2	2.51	0.40
1:C:122:ASN:O	1:C:123:GLU:HB2	2.22	0.40
1:B:82:TYR:CD1	1:B:82:TYR:C	2.94	0.40
1:B:224:ALA:O	1:B:230:ALA:HB3	2.21	0.40
1:C:123:GLU:O	1:C:124:GLU:C	2.60	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	320/368 (87%)	290 (91%)	22 (7%)	8 (2%)	5 28
1	B	311/368 (84%)	283 (91%)	20 (6%)	8 (3%)	5 27
1	C	320/368 (87%)	283 (88%)	28 (9%)	9 (3%)	5 25
All	All	951/1104 (86%)	856 (90%)	70 (7%)	25 (3%)	5 27

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	THR
1	A	228	MET
1	A	363	PRO
1	B	126	THR
1	B	192	ARG
1	B	228	MET
1	C	61	SER
1	C	126	THR

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Mol	Chain	Res	Type
1	C	191	LYS
1	C	228	MET
1	A	178	VAL
1	A	189	ARG
1	B	180	GLU
1	B	189	ARG
1	C	178	VAL
1	A	63	ARG
1	A	120	SER
1	C	63	ARG
1	A	64	LYS
1	B	120	SER
1	B	308	HIS
1	C	120	SER
1	C	174	PRO
1	B	178	VAL
1	C	64	LYS

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	288/322 (89%)	275 (96%)	13 (4%)	27	64
1	B	282/322 (88%)	272 (96%)	10 (4%)	36	71
1	C	288/322 (89%)	281 (98%)	7 (2%)	49	79
All	All	858/966 (89%)	828 (96%)	30 (4%)	36	71

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ARG
1	A	60	LYS
1	A	104	TYR
1	A	161	LEU
1	A	175	SER

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Mol	Chain	Res	Type
1	A	185	PHE
1	A	186	ASP
1	A	207	LYS
1	A	305	ARG
1	A	312	ARG
1	A	327	ARG
1	A	362	LYS
1	A	364	GLU
1	B	82	TYR
1	B	161	LEU
1	B	185	PHE
1	B	186	ASP
1	B	187	ASP
1	B	207	LYS
1	B	228	MET
1	B	305	ARG
1	B	312	ARG
1	B	327	ARG
1	C	185	PHE
1	C	207	LYS
1	C	287	ASN
1	C	297	ARG
1	C	305	ARG
1	C	312	ARG
1	C	327	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	342	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	$\#\|Z\| > 2$
2	ADP	B	401	3	24,29,29	0.99	2 (8%)	29,45,45	1.70	7 (24%)
2	ADP	A	401	3	24,29,29	1.09	2 (8%)	29,45,45	1.54	4 (13%)
2	ADP	C	401	3	24,29,29	0.96	1 (4%)	29,45,45	1.87	9 (31%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	B	401	3	-	1/12/32/32	0/3/3/3
2	ADP	A	401	3	-	4/12/32/32	0/3/3/3
2	ADP	C	401	3	-	4/12/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	ADP	O4'-C1'	3.34	1.45	1.41
2	B	401	ADP	O4'-C1'	2.50	1.44	1.41
2	A	401	ADP	C5-C4	2.42	1.47	1.40
2	B	401	ADP	C5-C4	2.38	1.47	1.40
2	C	401	ADP	C5-C4	2.33	1.47	1.40

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	401	ADP	C3'-C2'-C1'	4.69	108.04	100.98
2	A	401	ADP	PA-O3A-PB	-4.22	118.35	132.83
2	C	401	ADP	PA-O3A-PB	-4.17	118.52	132.83
2	B	401	ADP	PA-O3A-PB	-4.13	118.67	132.83
2	B	401	ADP	N3-C2-N1	-3.94	122.52	128.68
2	C	401	ADP	N3-C2-N1	-3.87	122.63	128.68
2	A	401	ADP	N3-C2-N1	-3.81	122.73	128.68
2	C	401	ADP	O3'-C3'-C4'	-2.90	102.67	111.05
2	B	401	ADP	C3'-C2'-C1'	2.84	105.25	100.98
2	C	401	ADP	C4-C5-N7	-2.70	106.59	109.40
2	B	401	ADP	C4-C5-N7	-2.61	106.67	109.40
2	B	401	ADP	O3A-PB-O1B	-2.58	96.87	111.19
2	A	401	ADP	C4-C5-N7	-2.45	106.84	109.40
2	C	401	ADP	O4'-C4'-C3'	2.37	109.81	105.11
2	C	401	ADP	C1'-N9-C4	-2.29	122.61	126.64
2	B	401	ADP	C2-N1-C6	2.27	122.64	118.75
2	A	401	ADP	C2-N1-C6	2.25	122.61	118.75
2	C	401	ADP	O2'-C2'-C3'	-2.24	104.58	111.82
2	C	401	ADP	C2-N1-C6	2.21	122.53	118.75
2	B	401	ADP	C1'-N9-C4	-2.06	123.02	126.64

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ADP	C5'-O5'-PA-O3A
2	C	401	ADP	C5'-O5'-PA-O2A
2	C	401	ADP	C5'-O5'-PA-O3A
2	A	401	ADP	C3'-C4'-C5'-O5'
2	C	401	ADP	C3'-C4'-C5'-O5'
2	A	401	ADP	O4'-C4'-C5'-O5'
2	A	401	ADP	C5'-O5'-PA-O1A
2	C	401	ADP	O4'-C4'-C5'-O5'
2	B	401	ADP	C5'-O5'-PA-O1A

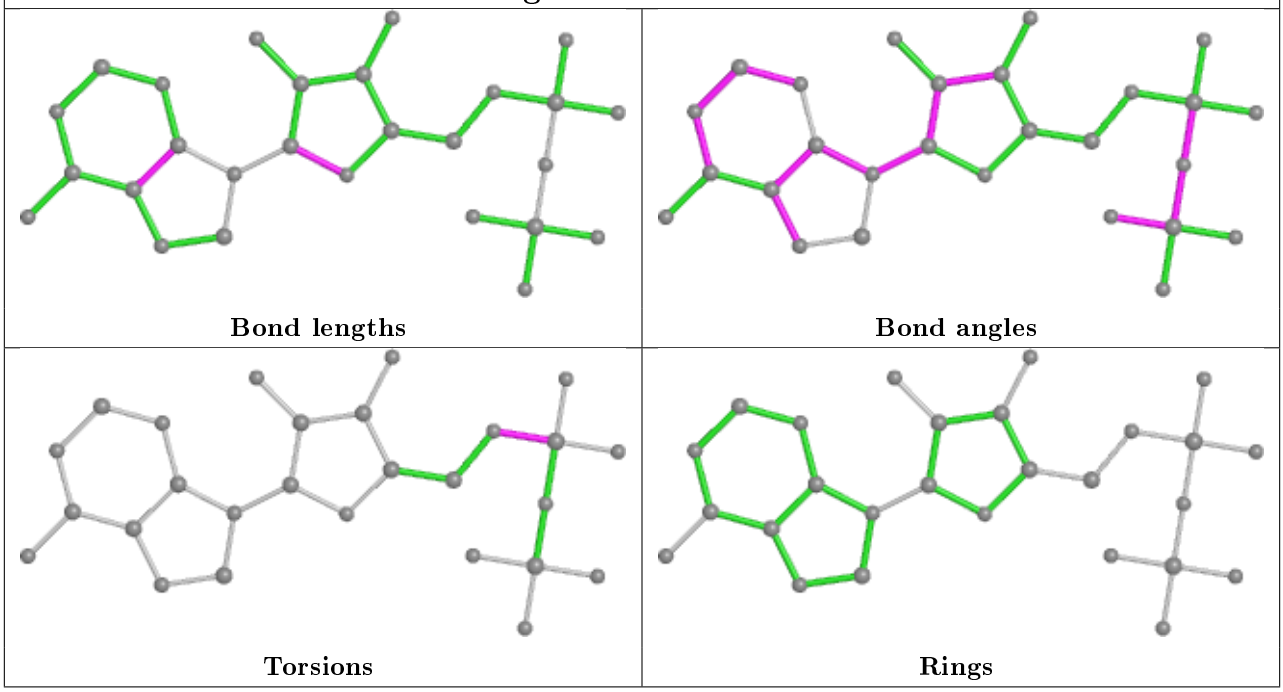
There are no ring outliers.

3 monomers are involved in 7 short contacts:

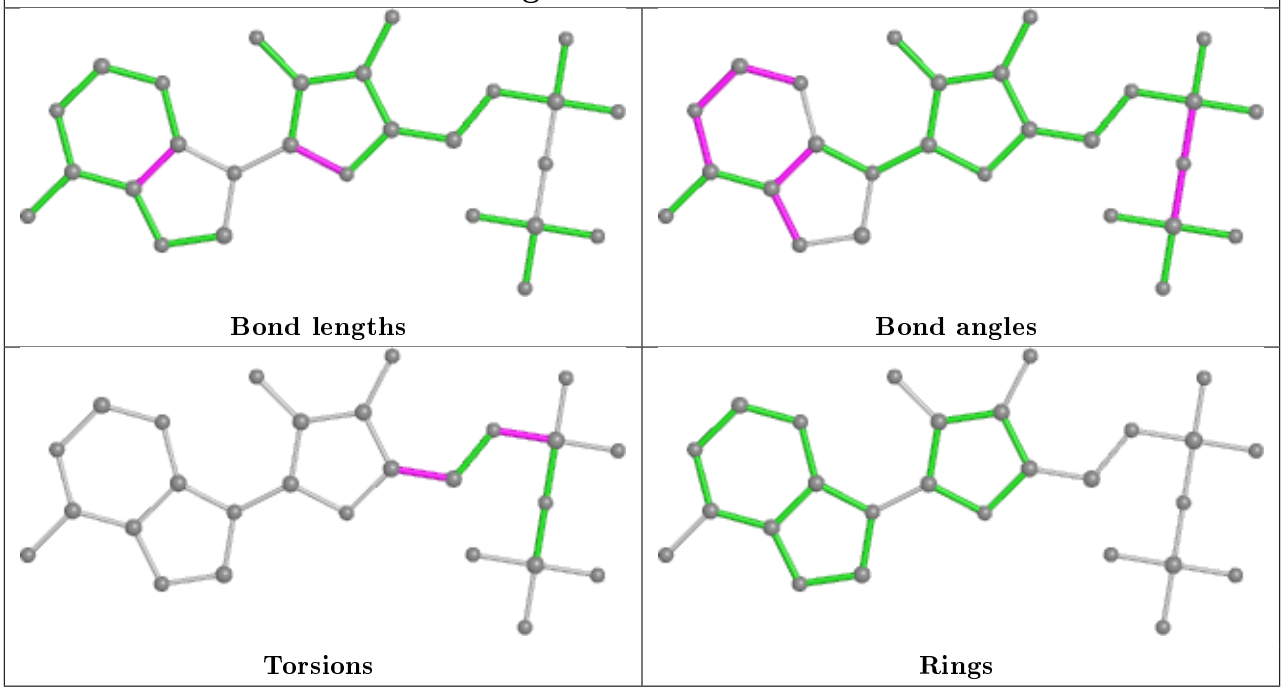
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	ADP	3	0
2	A	401	ADP	2	0
2	C	401	ADP	2	0

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 then 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.

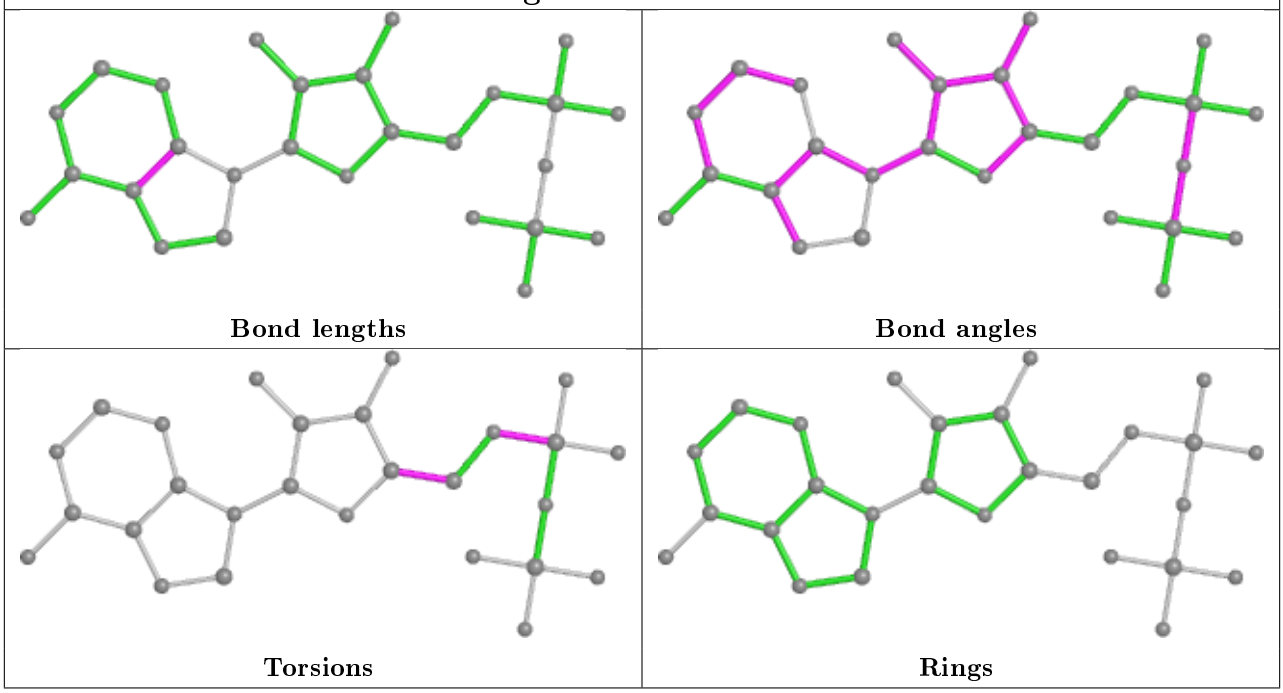
Ligand ADP B 401



Ligand ADP A 401



Ligand ADP C 401



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	326/368 (88%)	0.27	8 (2%)	57	29	50, 107, 191, 304	0
1	B	319/368 (86%)	0.41	20 (6%)	20	6	67, 121, 212, 280	0
1	C	326/368 (88%)	3.31	188 (57%)	0	0	279, 338, 366, 409	0
All	All	971/1104 (87%)	1.33	216 (22%)	0	0	50, 148, 356, 409	0

All (216) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	74	ALA	20.7
1	C	56	GLY	18.6
1	C	127	TRP	14.3
1	C	364	GLU	13.9
1	C	25	CYS	13.1
1	C	236	HIS	12.6
1	C	233	SER	12.1
1	C	18	ASN	12.0
1	C	114	THR	11.6
1	C	90	LEU	11.6
1	C	58	ALA	11.5
1	C	232	SER	10.8
1	C	303	VAL	10.7
1	C	365	VAL	10.7
1	C	172	LEU	10.3
1	C	179	SER	10.2
1	C	135	ILE	9.7
1	C	286	GLY	9.6
1	C	57	LEU	9.5
1	C	180	GLU	9.3
1	C	16	GLY	9.1
1	C	356	ALA	9.0
1	C	113	PHE	9.0

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Mol	Chain	Res	Type	RSRZ
1	C	182	LEU	8.8
1	C	126	THR	8.6
1	C	239	PHE	8.6
1	C	110	GLY	8.6
1	C	246	LYS	8.2
1	C	243	ILE	8.0
1	C	61	SER	8.0
1	C	299	ILE	7.9
1	C	30	LEU	7.9
1	C	353	ALA	7.8
1	C	289	ASN	7.7
1	C	51	SER	7.7
1	B	16	GLY	7.6
1	C	75	SER	7.5
1	C	266	LEU	7.5
1	C	190	ASN	7.4
1	C	106	GLN	7.3
1	C	41	VAL	7.2
1	C	181	ARG	7.2
1	C	160	LEU	6.7
1	C	310	PRO	6.6
1	C	168	LEU	6.6
1	C	186	ASP	6.5
1	C	256	VAL	6.5
1	C	357	LYS	6.4
1	C	307	PRO	6.4
1	C	29	ASN	6.2
1	C	42	GLU	6.2
1	C	241	VAL	6.2
1	C	234	CYS	6.1
1	C	332	ILE	6.1
1	C	73	GLY	6.0
1	C	235	SER	5.9
1	C	231	TYR	5.9
1	C	39	SER	5.9
1	C	264	VAL	5.8
1	C	105	GLY	5.8
1	C	302	LEU	5.8
1	C	211	TYR	5.8
1	C	91	ASP	5.8
1	B	38	HIS	5.7
1	A	177	ASP	5.7

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Mol	Chain	Res	Type	RSRZ
1	C	52	VAL	5.6
1	C	146	LYS	5.6
1	C	335	THR	5.6
1	B	127	TRP	5.5
1	C	149	ASP	5.5
1	C	237	SER	5.5
1	C	55	GLY	5.5
1	C	134	GLY	5.3
1	C	210	VAL	5.3
1	C	124	GLU	5.2
1	C	267	ALA	5.1
1	C	159	SER	5.1
1	C	163	ILE	5.1
1	C	287	ASN	5.0
1	C	104	TYR	4.9
1	C	122	ASN	4.7
1	C	333	ILE	4.7
1	C	94	ILE	4.7
1	C	316	LEU	4.6
1	C	347	LEU	4.6
1	C	328	THR	4.6
1	C	230	ALA	4.6
1	C	338	PRO	4.6
1	C	131	PRO	4.5
1	C	191	LYS	4.4
1	C	87	CYS	4.4
1	C	309	VAL	4.4
1	A	259	GLY	4.4
1	C	89	ILE	4.3
1	C	204	VAL	4.3
1	C	59	ASP	4.3
1	B	124	GLU	4.2
1	C	26	ARG	4.2
1	C	156	VAL	4.2
1	C	343	LEU	4.2
1	A	16	GLY	4.1
1	C	165	ASN	4.1
1	C	117	GLY	4.0
1	C	221	ARG	4.0
1	C	63	ARG	3.9
1	C	141	HIS	3.9
1	C	164	TYR	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	295	LEU	3.8
1	B	289	ASN	3.8
1	C	121	PRO	3.8
1	B	33	ARG	3.8
1	C	145	GLU	3.7
1	C	64	LYS	3.7
1	C	341	LEU	3.7
1	C	54	THR	3.7
1	C	166	GLU	3.7
1	C	324	LEU	3.7
1	C	300	THR	3.7
1	C	143	ILE	3.6
1	C	101	ILE	3.6
1	C	99	CYS	3.6
1	C	27	PRO	3.6
1	C	43	CYS	3.6
1	C	35	ALA	3.6
1	C	238	VAL	3.6
1	C	242	THR	3.5
1	C	33	ARG	3.5
1	C	84	SER	3.5
1	C	88	PRO	3.5
1	C	351	GLU	3.4
1	C	77	LYS	3.4
1	B	302	LEU	3.4
1	C	354	HIS	3.3
1	C	24	ARG	3.2
1	C	342	ASN	3.2
1	C	201	GLU	3.2
1	C	115	MET	3.1
1	C	212	GLN	3.1
1	B	27	PRO	3.1
1	C	305	ARG	3.1
1	C	31	ALA	3.1
1	C	60	LYS	3.1
1	C	62	SER	3.1
1	C	136	ILE	3.0
1	C	222	THR	3.0
1	B	68	PHE	3.0
1	C	183	GLN	3.0
1	C	260	LYS	3.0
1	C	19	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	311	TYR	3.0
1	C	128	GLU	2.9
1	C	196	ILE	2.9
1	C	20	GLN	2.9
1	B	66	TYR	2.9
1	B	34	LYS	2.8
1	B	52	VAL	2.8
1	C	326	GLY	2.8
1	C	203	THR	2.8
1	C	17	LYS	2.8
1	C	162	GLU	2.8
1	C	314	SER	2.7
1	A	123	GLU	2.7
1	B	282	ALA	2.7
1	C	137	PRO	2.7
1	C	349	THR	2.7
1	C	306	THR	2.7
1	B	340	SER	2.6
1	C	28	PHE	2.6
1	C	245	MET	2.6
1	C	219	ALA	2.6
1	C	129	GLU	2.6
1	B	31	ALA	2.5
1	C	207	LYS	2.5
1	C	184	MET	2.5
1	C	358	ASN	2.5
1	B	341	LEU	2.5
1	C	327	ARG	2.4
1	C	197	LYS	2.4
1	A	350	LEU	2.4
1	C	346	THR	2.4
1	C	263	LEU	2.4
1	C	308	HIS	2.4
1	C	32	GLU	2.4
1	C	188	PRO	2.4
1	B	144	PHE	2.4
1	C	123	GLU	2.4
1	C	78	GLN	2.4
1	C	350	LEU	2.4
1	A	124	GLU	2.4
1	B	162	GLU	2.4
1	C	340	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	208	ASP	2.4
1	A	210	VAL	2.3
1	C	189	ARG	2.3
1	C	34	LYS	2.3
1	C	167	GLU	2.3
1	C	361	ASN	2.3
1	C	174	PRO	2.2
1	C	46	VAL	2.2
1	C	125	TYR	2.2
1	C	177	ASP	2.2
1	C	38	HIS	2.2
1	C	240	SER	2.1
1	C	205	HIS	2.1
1	C	304	GLU	2.1
1	C	158	VAL	2.1
1	C	292	LEU	2.1
1	B	43	CYS	2.1
1	C	331	SER	2.1
1	C	40	ILE	2.1
1	B	64	LYS	2.1
1	C	265	ASP	2.0
1	C	109	THR	2.0
1	A	144	PHE	2.0
1	C	72	PHE	2.0
1	C	157	LYS	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 monosaccharides in this entry.

6.4 Ligands (i)

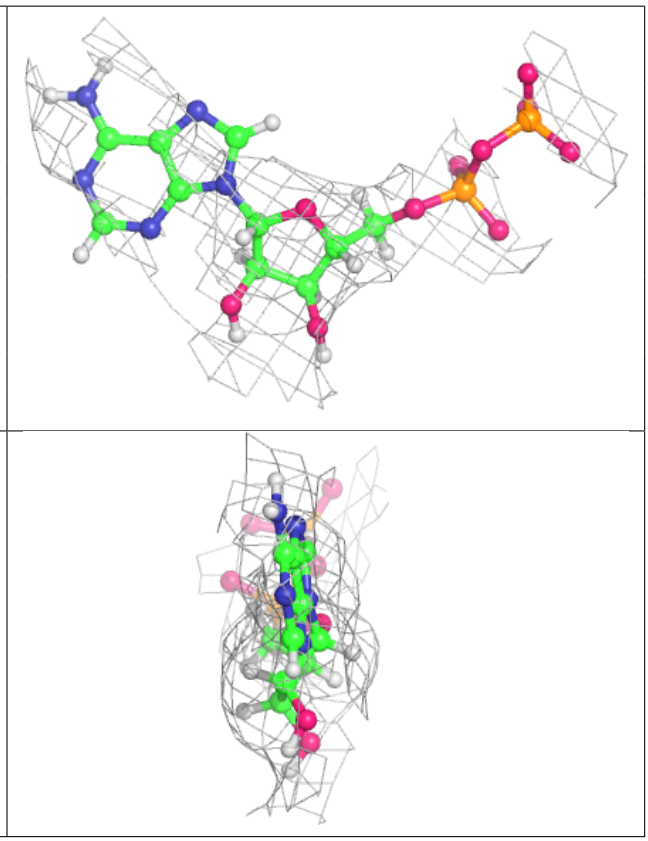
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MG	C	402	1/1	0.51	0.26	331,331,331,331	0
2	ADP	C	401	27/27	0.53	0.20	291,351,424,448	0
2	ADP	B	401	27/27	0.90	0.21	54,154,191,206	0
3	MG	B	402	1/1	0.93	0.19	69,69,69,69	0
2	ADP	A	401	27/27	0.94	0.21	55,116,153,176	0
3	MG	A	402	1/1	0.95	0.35	57,57,57,57	0

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

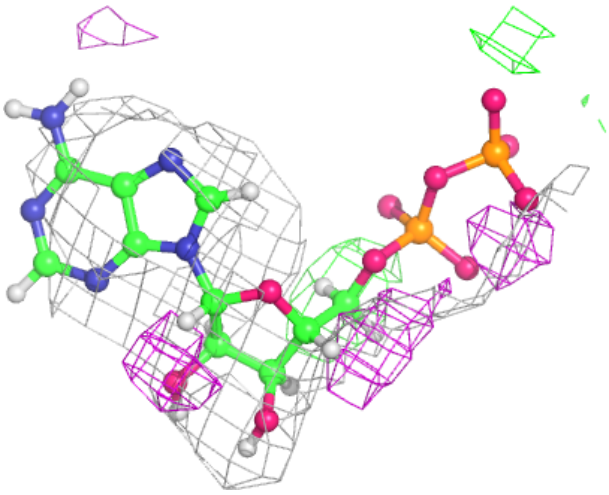
Electron density around ADP C 401:

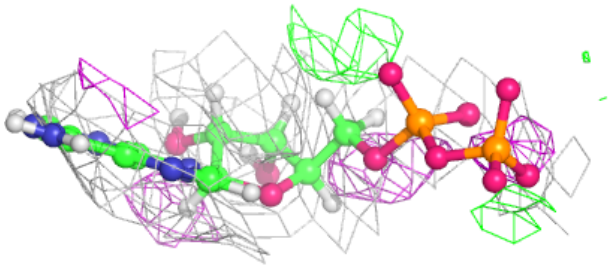
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

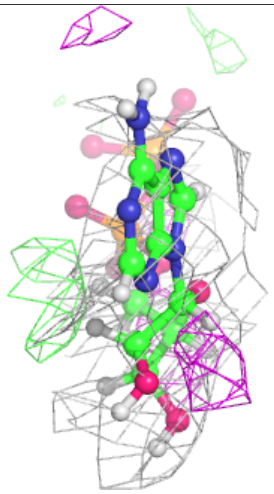


Electron density around ADP B 401:

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

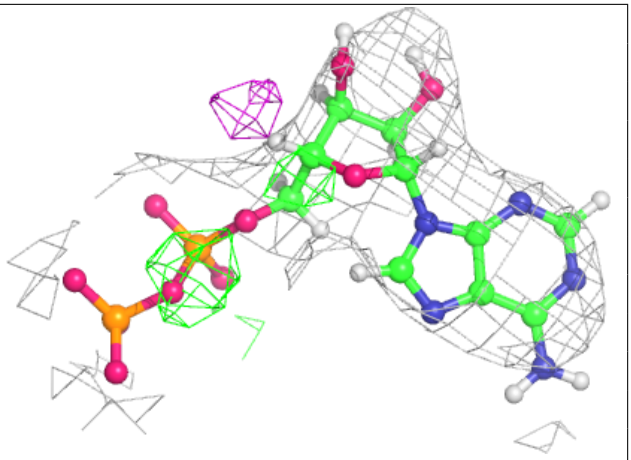


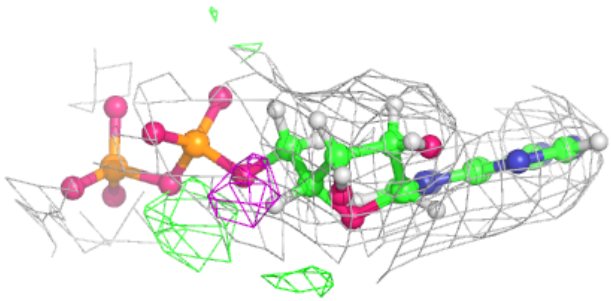


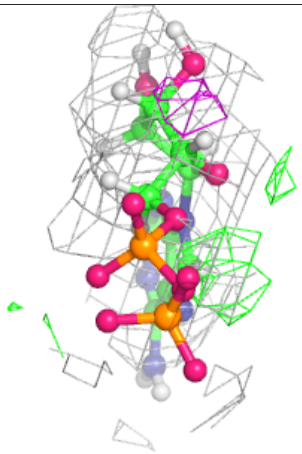


Electron density around ADP A 401:

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







6.5 Other polymers (i)

There are no such residues in this entry.