Full wwPDB X-ray Structure Validation Report (i)

Feb 10, 2024 – 03:58 PM EST

PDB ID	:	2RIO
Title	:	Structure of the dual enzyme Ire1 reveals the basis for catalysis and regulation of non-conventional splicing
Authors	:	Lee, K.P.; Dey, M.; Neculai, D.; Cao, C.; Dever, T.E.; Sicheri, F.
Deposited on	:	2007-10-12
Resolution	:	2.40 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

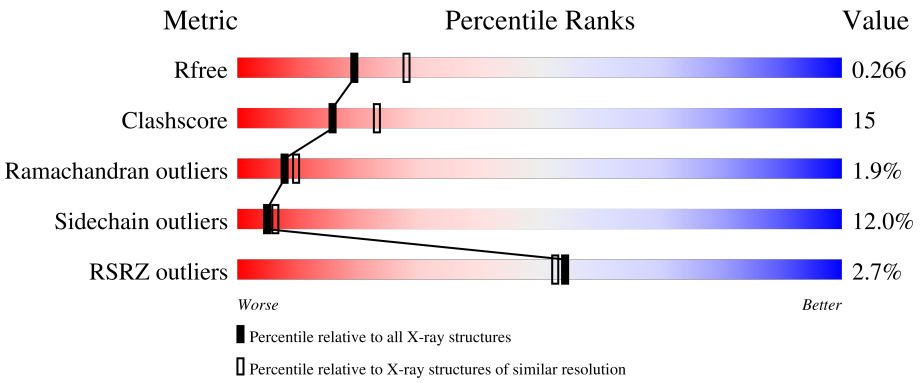
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	434	4% 60% 22% 7% • 9%
1	B	434	% 63% 21% • • 9%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6472 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 Serine/threonine-protein kinase/endoribonuclease IRE1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	396	3213	2058	547	590	18	0	0	0
1	B	395	3201	2051	544	588	18	0	0	0

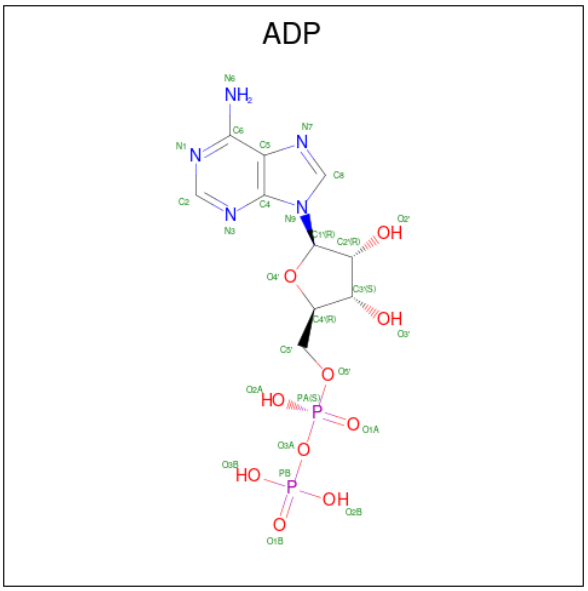
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is STRONTIUM ION (three-letter code: SR) (formula: Sr).

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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



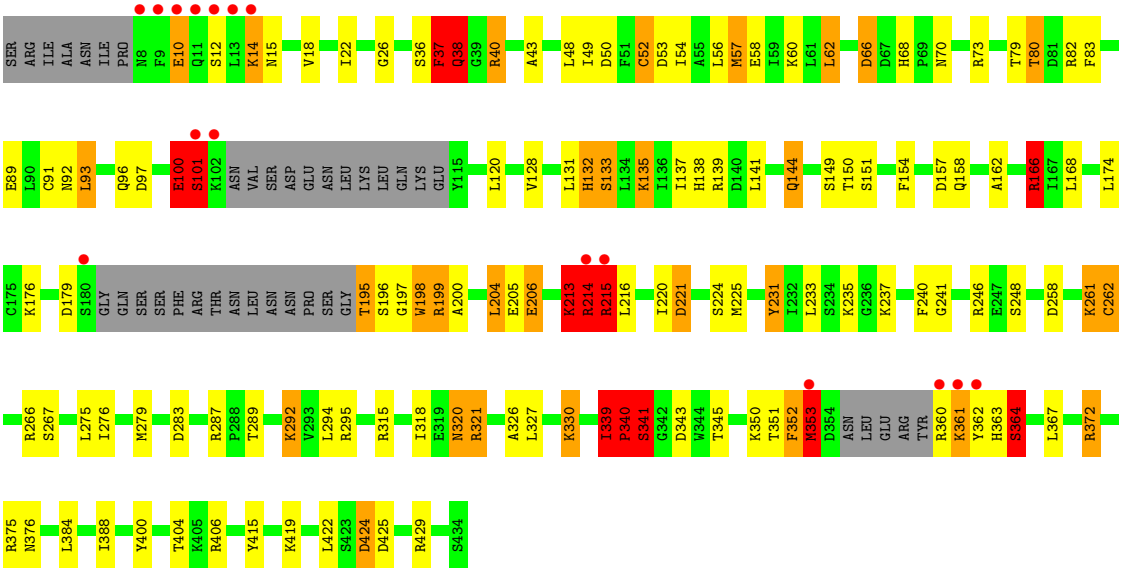
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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.

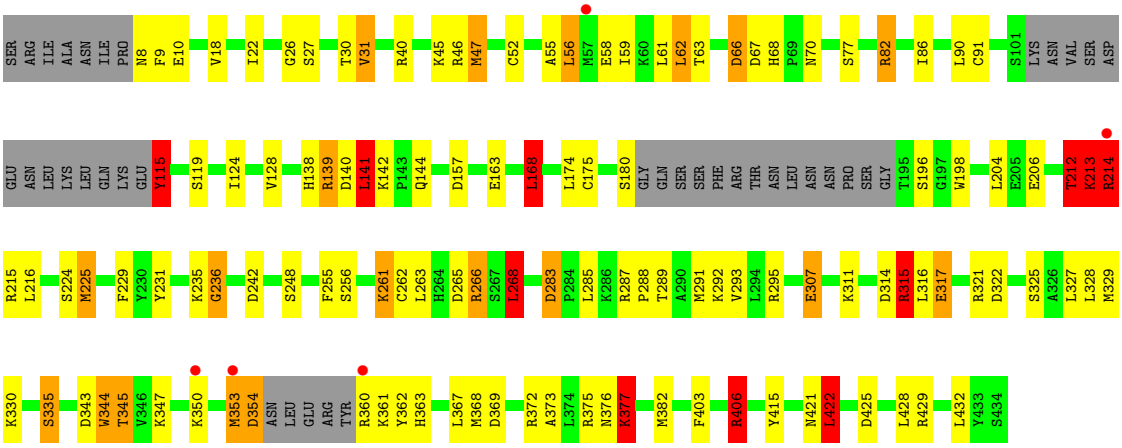
• Molecule 1: Serine/threonine-protein kinase/endoribonuclease IRE1

Chain A:



• Molecule 1: Serine/threonine-protein kinase/endoribonuclease IRE1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α, β, γ	130.31Å 130.31Å 175.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.83 – 2.40 19.83 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.83-2.40) 100.0 (19.83-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	82.65 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.222 , 0.266 0.224 , 0.266	Depositor DCC
R_{free} test set	3326 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	33.0	Xtriage
Anisotropy	0.003	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 12.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.162 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6472	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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, SR, 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	1.39	19/3279 (0.6%)	1.27	30/4417 (0.7%)
1	B	1.31	13/3267 (0.4%)	1.23	30/4402 (0.7%)
All	All	1.35	32/6546 (0.5%)	1.25	60/8819 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	11
1	B	0	1
All	All	0	12

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	283	ASP	CB-CG	11.80	1.76	1.51
1	B	262	CYS	CB-SG	-10.58	1.64	1.82
1	A	199	ARG	C-O	10.34	1.43	1.23
1	A	339	ILE	C-O	9.72	1.41	1.23
1	A	262	CYS	CB-SG	-9.66	1.65	1.82
1	A	206	GLU	CG-CD	7.65	1.63	1.51
1	A	339	ILE	CA-C	7.23	1.71	1.52
1	A	144	GLN	CG-CD	7.12	1.67	1.51
1	B	58	GLU	CG-CD	6.92	1.62	1.51
1	A	231	TYR	CD2-CE2	6.79	1.49	1.39
1	B	52	CYS	CB-SG	6.72	1.93	1.82
1	B	58	GLU	CB-CG	6.49	1.64	1.52
1	A	198	TRP	CE3-CZ3	6.25	1.49	1.38
1	B	415	TYR	CE2-CZ	6.12	1.46	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	58	GLU	CG-CD	6.10	1.61	1.51
1	B	163	GLU	CB-CG	6.08	1.63	1.52
1	B	236	GLY	N-CA	-6.01	1.37	1.46
1	A	241	GLY	CA-C	5.81	1.61	1.51
1	A	240	PHE	CE2-CZ	5.79	1.48	1.37
1	A	58	GLU	CB-CG	5.73	1.63	1.52
1	A	52	CYS	CB-SG	-5.72	1.72	1.81
1	A	240	PHE	CD1-CE1	5.55	1.50	1.39
1	A	400	TYR	CE2-CZ	5.54	1.45	1.38
1	B	403	PHE	CE2-CZ	5.43	1.47	1.37
1	A	221	ASP	CB-CG	-5.36	1.40	1.51
1	A	415	TYR	CD1-CE1	5.30	1.47	1.39
1	A	292	LYS	CD-CE	5.24	1.64	1.51
1	B	255	PHE	CE2-CZ	5.18	1.47	1.37
1	B	144	GLN	CG-CD	5.15	1.62	1.51
1	A	144	GLN	CB-CG	5.10	1.66	1.52
1	B	229	PHE	CD2-CE2	5.07	1.49	1.39
1	B	115	TYR	CD2-CE2	5.03	1.46	1.39

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	315	ARG	NE-CZ-NH2	-10.39	115.11	120.30
1	A	240	PHE	C-N-CA	9.53	142.32	122.30
1	A	240	PHE	CA-C-N	8.74	133.67	116.20
1	A	166	ARG	NE-CZ-NH2	-8.49	116.06	120.30
1	A	221	ASP	CB-CG-OD1	-8.28	110.85	118.30
1	A	221	ASP	CB-CA-C	8.07	126.55	110.40
1	A	66	ASP	CB-CG-OD1	8.05	125.54	118.30
1	A	40	ARG	NE-CZ-NH2	8.04	124.32	120.30
1	B	266	ARG	NE-CZ-NH2	-7.78	116.41	120.30
1	A	241	GLY	N-CA-C	-7.66	93.94	113.10
1	A	221	ASP	CB-CG-OD2	7.44	125.00	118.30
1	A	240	PHE	O-C-N	-7.34	110.71	123.20
1	B	406	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	A	287	ARG	NE-CZ-NH2	-7.11	116.75	120.30
1	A	133	SER	CA-CB-OG	-6.93	92.47	111.20
1	A	240	PHE	N-CA-C	6.86	129.51	111.00
1	B	291	MET	CG-SD-CE	6.82	111.11	100.20
1	B	47	MET	CA-CB-CG	6.79	124.85	113.30
1	B	235	LYS	CA-C-N	6.78	129.75	116.20
1	B	314	ASP	CB-CG-OD1	6.65	124.29	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	295	ARG	NE-CZ-NH2	-6.43	117.08	120.30
1	B	422	LEU	CA-CB-CG	6.43	130.09	115.30
1	B	406	ARG	CG-CD-NE	-6.26	98.64	111.80
1	A	315	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	B	168	LEU	CB-CG-CD1	6.15	121.46	111.00
1	B	46	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	B	139	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	A	372	ARG	NE-CZ-NH2	-6.11	117.24	120.30
1	A	315	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	B	295	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	B	235	LYS	C-N-CA	6.04	134.98	122.30
1	A	93	LEU	CA-CB-CG	6.02	129.15	115.30
1	B	66	ASP	CB-CG-OD2	5.99	123.69	118.30
1	B	295	ARG	CG-CD-NE	-5.98	99.24	111.80
1	B	175	CYS	CA-CB-SG	-5.86	103.45	114.00
1	B	315	ARG	NE-CZ-NH1	5.85	123.23	120.30
1	B	56	LEU	CB-CG-CD1	5.81	120.88	111.00
1	B	214	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	A	100	GLU	C-N-CA	5.76	136.10	121.70
1	B	235	LYS	O-C-N	-5.72	113.47	123.20
1	B	283	ASP	CB-CA-C	5.70	121.81	110.40
1	A	294	LEU	CB-CG-CD2	-5.68	101.34	111.00
1	A	199	ARG	N-CA-C	5.58	126.08	111.00
1	B	287	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	50	ASP	CB-CG-OD1	5.43	123.19	118.30
1	B	157	ASP	CB-CG-OD1	5.36	123.12	118.30
1	A	205	GLU	CA-CB-CG	5.31	125.09	113.40
1	A	283	ASP	CB-CG-OD1	5.31	123.08	118.30
1	A	38	GLN	CB-CA-C	5.30	121.01	110.40
1	A	339	ILE	CB-CA-C	5.24	122.09	111.60
1	B	268	LEU	CB-CG-CD2	5.19	119.82	111.00
1	A	166	ARG	CG-CD-NE	-5.19	100.91	111.80
1	A	406	ARG	CG-CD-NE	-5.12	101.06	111.80
1	B	141	LEU	CA-CB-CG	5.12	127.06	115.30
1	A	157	ASP	CB-CG-OD1	5.05	122.85	118.30
1	A	424	ASP	CB-CG-OD1	5.05	122.85	118.30
1	B	141	LEU	CB-CG-CD1	5.03	119.56	111.00
1	B	225	MET	CG-SD-CE	-5.03	92.15	100.20
1	B	168	LEU	CA-CB-CG	5.02	126.86	115.30
1	B	377	LYS	CD-CE-NZ	-5.00	100.19	111.70

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	GLU	Peptide
1	A	213	LYS	Peptide
1	A	214	ARG	Peptide
1	A	339	ILE	Peptide
1	A	340	PRO	Peptide
1	A	352	PHE	Peptide
1	A	353	MET	Peptide
1	A	361	LYS	Peptide
1	A	37	PHE	Mainchain,Peptide
1	A	38	GLN	Peptide
1	B	213	LYS	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3213	0	3239	107	0
1	B	3201	0	3222	86	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	27	0	12	2	0
4	B	27	0	12	1	0
All	All	6472	0	6485	191	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 15.

All (191) 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:283:ASP:CB	1:B:283:ASP:CG	1.76	1.50
1:A:128:VAL:HG21	1:A:225:MET:CE	1.71	1.21
1:B:124:ILE:HG22	1:B:225:MET:HE2	1.12	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:ARG:HB3	1:A:215:ARG:CB	1.90	1.01
1:A:214:ARG:HB3	1:A:215:ARG:HB2	1.39	1.00
1:A:214:ARG:NH2	1:A:216:LEU:HA	1.78	0.99
1:B:128:VAL:HG21	1:B:225:MET:HE3	1.43	0.99
1:A:100:GLU:HB3	1:A:101:SER:HB2	1.44	0.98
1:A:128:VAL:HG21	1:A:225:MET:HE3	1.47	0.96
1:A:80:THR:HG22	1:A:83:PHE:H	1.30	0.96
1:A:100:GLU:CB	1:A:101:SER:HB2	1.97	0.95
1:A:128:VAL:CG2	1:A:225:MET:CE	2.45	0.94
1:A:92:ASN:HD22	1:A:149:SER:HA	1.30	0.94
1:A:214:ARG:HH21	1:A:216:LEU:H	1.05	0.93
1:A:128:VAL:CG2	1:A:225:MET:HE1	1.99	0.93
1:B:330:LYS:HD3	1:B:421:ASN:OD1	1.71	0.91
1:A:320:ASN:HD22	1:A:321:ARG:H	1.18	0.90
1:B:360:ARG:HH22	1:B:372:ARG:HH21	1.20	0.90
1:A:214:ARG:HH22	1:A:216:LEU:HA	1.38	0.85
1:A:96:GLN:NE2	1:A:144:GLN:HE21	1.76	0.83
1:B:124:ILE:HG22	1:B:225:MET:CE	2.02	0.83
1:B:347:LYS:HB3	1:B:406:ARG:HG2	1.60	0.82
1:A:96:GLN:HE21	1:A:144:GLN:HE21	1.26	0.82
1:A:214:ARG:NH2	1:A:216:LEU:H	1.78	0.82
1:B:45:LYS:HG2	1:B:47:MET:CE	2.11	0.81
1:A:320:ASN:HD22	1:A:321:ARG:N	1.79	0.81
1:A:97:ASP:O	1:A:101:SER:HB3	1.81	0.80
1:B:353:MET:HA	1:B:354:ASP:C	2.02	0.80
1:A:100:GLU:HB3	1:A:101:SER:CB	2.12	0.79
1:A:289:THR:H	1:A:292:LYS:HE3	1.48	0.79
1:A:214:ARG:HH21	1:A:216:LEU:N	1.81	0.78
1:A:361:LYS:HB2	1:A:362:TYR:HA	1.66	0.78
1:A:73:ARG:HB3	1:A:89:GLU:HG2	1.67	0.77
1:A:214:ARG:CB	1:A:215:ARG:HB2	2.13	0.77
1:B:128:VAL:HG21	1:B:225:MET:CE	2.15	0.77
1:A:92:ASN:ND2	1:A:149:SER:HA	2.02	0.75
1:A:14:LYS:H	1:A:14:LYS:HD2	1.52	0.74
1:A:15:ASN:O	1:A:37:PHE:O	2.06	0.74
1:A:92:ASN:ND2	1:A:150:THR:H	1.86	0.73
1:B:68:HIS:HD2	1:B:70:ASN:H	1.36	0.72
1:B:361:LYS:HB2	1:B:362:TYR:HA	1.70	0.72
1:A:214:ARG:NH2	1:A:216:LEU:CA	2.51	0.71
1:A:235:LYS:HD2	1:A:261:LYS:HB2	1.71	0.71
1:A:276:ILE:HA	1:A:279:MET:CE	2.21	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:PHE:O	1:A:37:PHE:HD2	1.74	0.70
1:A:68:HIS:HD2	1:A:70:ASN:H	1.40	0.69
1:B:425:ASP:O	1:B:429:ARG:HG3	1.91	0.69
1:B:344:TRP:C	1:B:344:TRP:CD1	2.65	0.69
1:B:288:PRO:HB3	1:B:292:LYS:HD3	1.74	0.69
1:B:128:VAL:CG2	1:B:225:MET:CE	2.72	0.68
1:B:321:ARG:O	1:B:321:ARG:HG3	1.93	0.68
1:A:214:ARG:NH2	1:A:216:LEU:N	2.39	0.68
1:B:68:HIS:CD2	1:B:70:ASN:H	2.11	0.68
1:B:343:ASP:OD1	1:B:345:THR:HB	1.94	0.68
1:B:139:ARG:HD3	1:B:174:LEU:O	1.94	0.67
1:A:214:ARG:CB	1:A:215:ARG:CB	2.71	0.67
1:B:124:ILE:CG2	1:B:225:MET:HE2	2.08	0.66
1:A:97:ASP:O	1:A:101:SER:CB	2.44	0.66
1:A:363:HIS:O	1:A:364:SER:HB3	1.96	0.65
1:B:373:ALA:O	1:B:377:LYS:HG2	1.98	0.64
1:A:37:PHE:O	1:A:37:PHE:CD2	2.50	0.64
1:B:128:VAL:CG2	1:B:225:MET:HE3	2.23	0.64
1:A:80:THR:HG23	1:A:82:ARG:H	1.62	0.63
1:A:214:ARG:HH21	1:A:214:ARG:HB2	1.63	0.63
1:B:316:LEU:HD23	1:B:368:MET:HG3	1.80	0.63
1:A:195:THR:HG23	1:A:197:GLY:H	1.65	0.62
1:A:132:HIS:CE1	1:A:221:ASP:OD1	2.52	0.62
1:A:68:HIS:CD2	1:A:70:ASN:H	2.16	0.62
1:A:214:ARG:HB3	1:A:215:ARG:HB3	1.81	0.62
1:A:425:ASP:O	1:A:429:ARG:HG3	1.99	0.61
1:B:91:CYS:SG	1:B:168:LEU:HD23	2.41	0.61
1:A:276:ILE:HA	1:A:279:MET:HE2	1.81	0.61
1:B:55:ALA:O	1:B:59:ILE:HG12	2.02	0.60
1:A:37:PHE:CD2	1:A:37:PHE:C	2.75	0.60
1:A:276:ILE:HA	1:A:279:MET:HE3	1.84	0.59
1:A:318:ILE:HD13	1:B:382:MET:HG3	1.83	0.59
1:A:128:VAL:CB	1:A:225:MET:HE1	2.32	0.59
1:B:360:ARG:HH22	1:B:372:ARG:NH2	1.96	0.58
1:A:363:HIS:O	1:A:364:SER:CB	2.50	0.58
1:A:26:GLY:HA3	4:A:1101:ADP:O1B	2.04	0.57
1:A:96:GLN:HE21	1:A:144:GLN:NE2	2.00	0.57
1:A:214:ARG:HB2	1:A:216:LEU:H	1.69	0.57
1:B:128:VAL:HB	1:B:225:MET:HE1	1.85	0.57
1:A:214:ARG:CA	1:A:215:ARG:HB2	2.34	0.57
1:A:384:LEU:HD13	1:A:388:ILE:HG22	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:ASP:HB3	1:B:268:LEU:HD22	1.88	0.56
1:A:128:VAL:HB	1:A:225:MET:HE1	1.88	0.56
1:B:10:GLU:HG2	1:B:18:VAL:HB	1.88	0.55
1:A:139:ARG:NH1	1:A:176:LYS:HB2	2.22	0.55
1:B:361:LYS:HE2	1:B:372:ARG:HG2	1.88	0.55
1:A:327:LEU:HD11	1:A:422:LEU:HD21	1.89	0.54
1:A:139:ARG:NH1	1:A:174:LEU:O	2.41	0.54
1:A:80:THR:HG22	1:A:83:PHE:N	2.12	0.54
1:A:139:ARG:HH11	1:A:176:LYS:HB2	1.73	0.54
1:B:283:ASP:CG	1:B:283:ASP:CA	2.72	0.53
1:A:361:LYS:HB2	1:A:362:TYR:CG	2.44	0.53
1:B:316:LEU:HD23	1:B:368:MET:CG	2.39	0.53
1:B:321:ARG:NH1	1:B:363:HIS:ND1	2.56	0.53
1:B:360:ARG:HA	1:B:361:LYS:HG2	1.91	0.53
1:B:45:LYS:HG2	1:B:47:MET:HE1	1.90	0.53
1:B:128:VAL:CG2	1:B:225:MET:HE1	2.39	0.52
1:B:62:LEU:HD22	1:B:66:ASP:HB3	1.91	0.52
1:A:361:LYS:HB2	1:A:362:TYR:CA	2.38	0.52
1:B:213:LYS:HG2	1:B:214:ARG:HA	1.91	0.52
1:A:53:ASP:O	1:A:57:MET:SD	2.68	0.52
1:B:212:THR:N	1:B:213:LYS:HB2	2.24	0.52
1:B:361:LYS:NZ	1:B:368:MET:HE3	2.24	0.52
1:B:26:GLY:HA3	4:B:2101:ADP:O1B	2.10	0.51
1:B:22:ILE:HD13	1:B:31:VAL:HG21	1.91	0.51
1:A:100:GLU:HB2	1:A:101:SER:HB2	1.90	0.51
1:A:214:ARG:HH22	1:A:216:LEU:CA	2.17	0.51
1:A:91:CYS:SG	1:A:168:LEU:HD12	2.51	0.51
1:B:361:LYS:HE3	1:B:369:ASP:OD1	2.10	0.51
1:B:361:LYS:NZ	1:B:368:MET:CE	2.74	0.50
1:B:335:SER:HB2	1:B:367:LEU:H	1.76	0.50
1:B:375:ARG:HD3	1:B:375:ARG:C	2.31	0.50
1:A:166:ARG:NH2	1:B:67:ASP:OD2	2.41	0.50
1:A:326:ALA:O	1:A:330:LYS:HG2	2.11	0.50
1:A:198:TRP:O	1:A:224:SER:HB3	2.11	0.50
1:A:233:LEU:O	1:A:262:CYS:HB2	2.11	0.49
1:A:235:LYS:HD3	1:A:261:LYS:HE2	1.93	0.49
1:A:339:ILE:HB	1:A:340:PRO:O	2.12	0.49
1:B:45:LYS:HG2	1:B:47:MET:HE2	1.91	0.49
1:B:214:ARG:HE	1:B:216:LEU:HD11	1.76	0.49
1:B:360:ARG:NH2	1:B:372:ARG:HE	2.11	0.49
1:B:45:LYS:HE2	1:B:47:MET:CE	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:THR:CG2	1:A:82:ARG:H	2.26	0.48
1:A:138:HIS:O	1:A:139:ARG:HB2	2.14	0.48
1:A:141:LEU:CD1	1:A:221:ASP:HB2	2.44	0.47
1:B:330:LYS:HD3	1:B:421:ASN:CG	2.34	0.47
1:A:276:ILE:HG12	1:A:279:MET:CE	2.43	0.47
1:B:315:ARG:HD3	1:B:425:ASP:OD2	2.15	0.47
1:A:141:LEU:HD12	1:A:221:ASP:HB2	1.95	0.47
1:B:91:CYS:SG	1:B:168:LEU:CD2	3.03	0.47
1:B:196:SER:HB3	1:B:204:LEU:HD21	1.97	0.47
1:B:307:GLU:O	1:B:311:LYS:HG2	2.14	0.47
1:A:361:LYS:CB	1:A:362:TYR:HA	2.41	0.47
1:A:154:PHE:HB3	1:A:166:ARG:HG3	1.96	0.46
1:A:267:SER:HB3	1:A:404:THR:HG23	1.97	0.46
1:A:36:SER:HA	1:A:40:ARG:O	2.15	0.46
1:A:62:LEU:HD22	1:A:66:ASP:HB3	1.96	0.46
1:B:10:GLU:HG2	1:B:18:VAL:CB	2.46	0.46
1:A:375:ARG:HH11	1:A:376:ASN:ND2	2.12	0.46
1:A:350:LYS:O	1:A:353:MET:HG3	2.16	0.46
1:A:15:ASN:C	1:A:37:PHE:O	2.53	0.46
1:A:261:LYS:HB3	1:A:261:LYS:HE3	1.58	0.46
1:B:115:TYR:C	1:B:115:TYR:CD1	2.89	0.46
1:B:214:ARG:NE	1:B:216:LEU:HD11	2.31	0.46
1:B:372:ARG:HA	1:B:372:ARG:HD2	1.69	0.46
1:B:198:TRP:O	1:B:224:SER:HB3	2.16	0.46
1:A:128:VAL:CB	1:A:225:MET:CE	2.92	0.45
1:A:214:ARG:NH2	1:A:214:ARG:HB2	2.28	0.45
1:A:350:LYS:C	1:A:352:PHE:H	2.19	0.45
1:B:317:GLU:HG2	1:B:372:ARG:CZ	2.46	0.45
1:B:30:THR:HG21	1:B:45:LYS:HE3	1.98	0.45
1:B:59:ILE:HD11	1:B:86:ILE:HD13	1.98	0.45
1:B:138:HIS:CG	1:B:141:LEU:HD13	2.51	0.45
1:B:231:TYR:O	1:B:236:GLY:HA2	2.16	0.45
1:B:289:THR:O	1:B:293:VAL:HG23	2.16	0.45
1:A:54:ILE:HA	1:A:57:MET:SD	2.58	0.44
1:B:361:LYS:HZ2	1:B:368:MET:CE	2.31	0.44
1:A:135:LYS:HB2	1:A:135:LYS:HE2	1.72	0.44
1:A:158:GLN:HG2	1:A:162:ALA:HA	1.99	0.44
1:A:343:ASP:HA	1:A:364:SER:O	2.17	0.44
1:B:361:LYS:HB2	1:B:362:TYR:CA	2.42	0.43
1:A:340:PRO:N	1:A:341:SER:HB3	2.33	0.43
1:A:154:PHE:HB3	1:A:166:ARG:CG	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:LYS:O	1:B:353:MET:O	2.36	0.43
1:A:275:LEU:O	1:A:279:MET:HG3	2.19	0.42
1:B:288:PRO:CB	1:B:292:LYS:HD3	2.47	0.42
1:B:325:SER:O	1:B:329:MET:HG3	2.19	0.42
1:B:375:ARG:HH11	1:B:376:ASN:ND2	2.17	0.42
1:A:10:GLU:HB3	1:A:18:VAL:HB	2.01	0.42
1:B:377:LYS:HA	1:B:377:LYS:HD2	1.70	0.42
1:B:321:ARG:NH2	1:B:368:MET:HE1	2.34	0.42
1:B:327:LEU:HD11	1:B:422:LEU:HD22	2.01	0.42
1:B:128:VAL:CB	1:B:225:MET:HE1	2.48	0.42
1:A:43:ALA:HB3	4:A:1101:ADP:N6	2.35	0.42
1:B:82:ARG:HE	1:B:82:ARG:HB2	1.45	0.42
1:A:231:TYR:HA	1:A:237:LYS:O	2.20	0.42
1:B:47:MET:HE1	1:B:86:ILE:HD12	2.02	0.41
1:A:132:HIS:HE1	1:A:221:ASP:OD2	2.03	0.41
1:A:213:LYS:HB3	1:A:214:ARG:C	2.40	0.41
1:B:285:LEU:HD23	1:B:285:LEU:HA	1.90	0.41
1:B:215:ARG:HA	1:B:215:ARG:HD2	1.94	0.41
1:B:138:HIS:CE1	1:B:140:ASP:HB3	2.56	0.41
1:A:22:ILE:HD13	1:A:22:ILE:HG21	1.91	0.40
1:A:137:ILE:HG22	1:A:139:ARG:HG3	2.03	0.40
1:B:8:ASN:HB3	1:B:9:PHE:H	1.65	0.40
1:A:196:SER:HB3	1:A:204:LEU:HD11	2.03	0.40
1:B:328:LEU:HD23	1:B:328:LEU:HA	1.81	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	388/434 (89%)	353 (91%)	24 (6%)	11 (3%)	5 4

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	387/434 (89%)	365 (94%)	18 (5%)	4 (1%)	15 23
All	All	775/868 (89%)	718 (93%)	42 (5%)	15 (2%)	8 10

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	14	LYS
1	A	38	GLN
1	A	100	GLU
1	A	200	ALA
1	A	340	PRO
1	B	261	LYS
1	A	101	SER
1	A	215	ARG
1	A	364	SER
1	B	213	LYS
1	A	37	PHE
1	A	214	ARG
1	A	341	SER
1	B	212	THR
1	B	214	ARG

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	363/398 (91%)	315 (87%)	48 (13%)	4 4
1	B	361/398 (91%)	322 (89%)	39 (11%)	6 9
All	All	724/796 (91%)	637 (88%)	87 (12%)	5 6

All (87) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU

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Mol	Chain	Res	Type
1	A	12	SER
1	A	48	LEU
1	A	49	ILE
1	A	52	CYS
1	A	56	LEU
1	A	57	MET
1	A	60	LYS
1	A	62	LEU
1	A	79	THR
1	A	80	THR
1	A	93	LEU
1	A	100	GLU
1	A	101	SER
1	A	120	LEU
1	A	131	LEU
1	A	132	HIS
1	A	133	SER
1	A	135	LYS
1	A	151	SER
1	A	166	ARG
1	A	179	ASP
1	A	195	THR
1	A	199	ARG
1	A	204	LEU
1	A	206	GLU
1	A	213	LYS
1	A	214	ARG
1	A	215	ARG
1	A	220	ILE
1	A	246	ARG
1	A	248	SER
1	A	258	ASP
1	A	261	LYS
1	A	266	ARG
1	A	320	ASN
1	A	321	ARG
1	A	330	LYS
1	A	341	SER
1	A	345	THR
1	A	351	THR
1	A	353	MET
1	A	360	ARG

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Mol	Chain	Res	Type
1	A	364	SER
1	A	367	LEU
1	A	372	ARG
1	A	419	LYS
1	A	424	ASP
1	B	27	SER
1	B	31	VAL
1	B	40	ARG
1	B	56	LEU
1	B	61	LEU
1	B	62	LEU
1	B	63	THR
1	B	77	SER
1	B	82	ARG
1	B	90	LEU
1	B	115	TYR
1	B	119	SER
1	B	141	LEU
1	B	142	LYS
1	B	168	LEU
1	B	180	SER
1	B	206	GLU
1	B	212	THR
1	B	242	ASP
1	B	248	SER
1	B	256	SER
1	B	261	LYS
1	B	263	LEU
1	B	266	ARG
1	B	268	LEU
1	B	307	GLU
1	B	315	ARG
1	B	317	GLU
1	B	322	ASP
1	B	335	SER
1	B	344	TRP
1	B	345	THR
1	B	353	MET
1	B	354	ASP
1	B	377	LYS
1	B	406	ARG
1	B	422	LEU

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Mol	Chain	Res	Type
1	B	428	LEU
1	B	432	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	34	GLN
1	A	68	HIS
1	A	70	ASN
1	A	92	ASN
1	A	96	GLN
1	A	116	ASN
1	A	130	HIS
1	A	132	HIS
1	A	320	ASN
1	A	363	HIS
1	A	409	ASN
1	B	68	HIS
1	B	70	ASN
1	B	130	HIS
1	B	376	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, 4 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	$\#\|Z\| > 2$
4	ADP	B	2101	3,2	24,29,29	0.96	1 (4%)	29,45,45	1.51	4 (13%)
4	ADP	A	1101	3,2	24,29,29	0.96	1 (4%)	29,45,45	1.51	4 (13%)

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
4	ADP	B	2101	3,2	-	3/12/32/32	0/3/3/3
4	ADP	A	1101	3,2	-	3/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2101	ADP	C5-C4	2.50	1.47	1.40
4	A	1101	ADP	C5-C4	2.49	1.47	1.40

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	1101	ADP	PA-O3A-PB	-3.60	120.47	132.83
4	B	2101	ADP	PA-O3A-PB	-3.58	120.53	132.83
4	A	1101	ADP	C3'-C2'-C1'	3.51	106.26	100.98
4	B	2101	ADP	C3'-C2'-C1'	3.49	106.23	100.98
4	B	2101	ADP	N3-C2-N1	-3.16	123.73	128.68
4	A	1101	ADP	N3-C2-N1	-3.13	123.78	128.68
4	B	2101	ADP	C4-C5-N7	-2.72	106.56	109.40
4	A	1101	ADP	C4-C5-N7	-2.68	106.61	109.40

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1101	ADP	C5'-O5'-PA-O2A
4	A	1101	ADP	C5'-O5'-PA-O3A
4	B	2101	ADP	C5'-O5'-PA-O3A
4	A	1101	ADP	C5'-O5'-PA-O1A
4	B	2101	ADP	C5'-O5'-PA-O1A
4	B	2101	ADP	C5'-O5'-PA-O2A

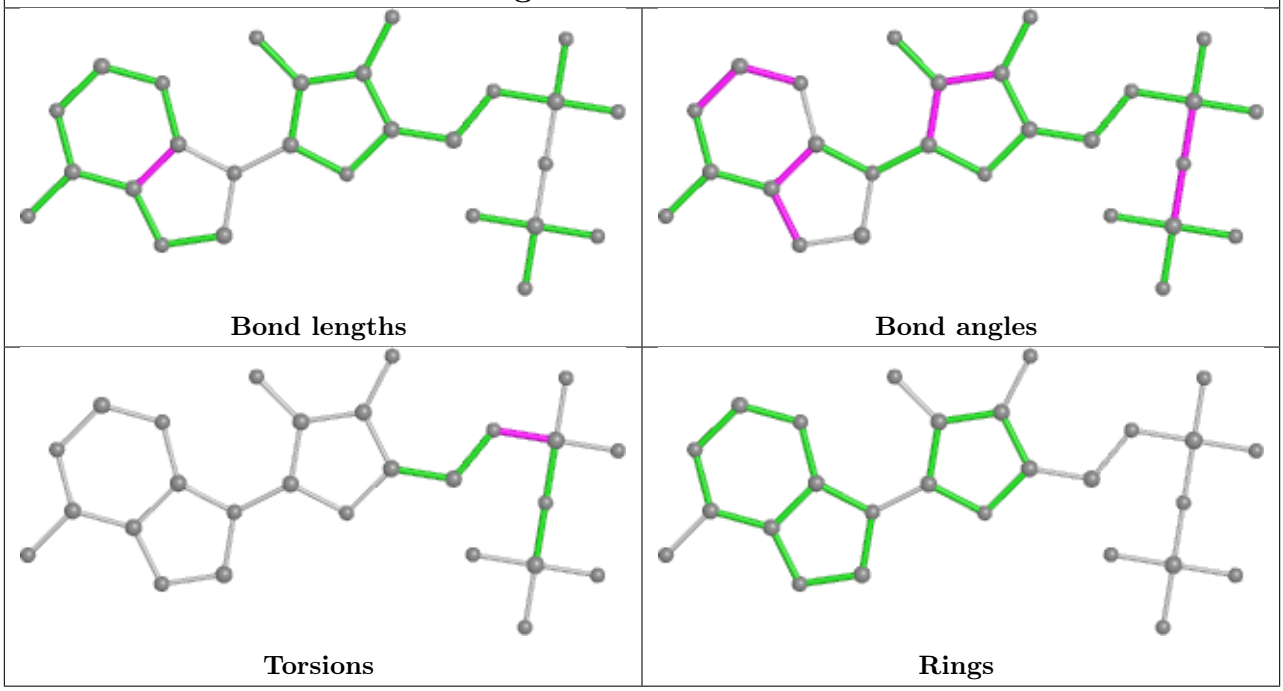
There are no ring outliers.

2 monomers are involved in 3 short contacts:

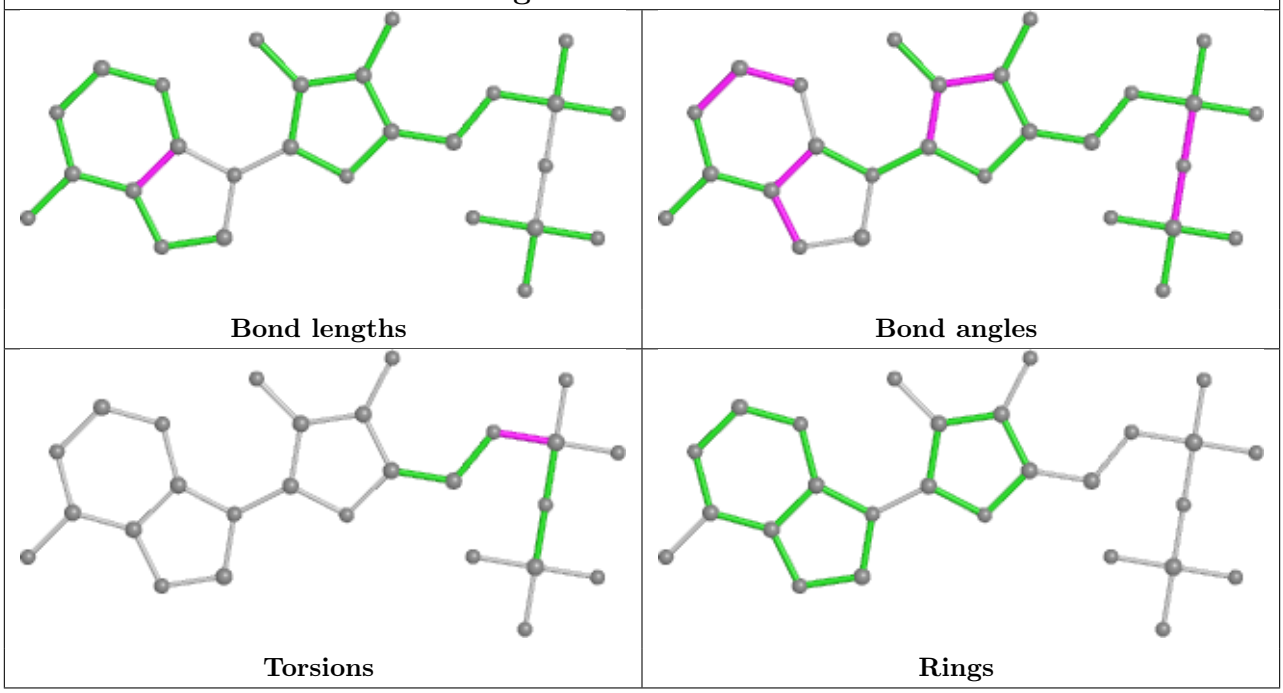
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2101	ADP	1	0
4	A	1101	ADP	2	0

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 2101



Ligand ADP A 1101



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	396/434 (91%)	0.12	16 (4%)	38	37	18, 31, 63, 105	0
1	B	395/434 (91%)	-0.06	5 (1%)	77	75	19, 32, 58, 78	0
All	All	791/868 (91%)	0.03	21 (2%)	54	52	18, 32, 61, 105	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	102	LYS	6.5
1	A	12	SER	6.0
1	A	8	ASN	5.9
1	A	215	ARG	5.8
1	A	9	PHE	5.8
1	A	11	GLN	5.1
1	A	360	ARG	3.9
1	B	360	ARG	3.7
1	A	13	LEU	3.5
1	A	353	MET	3.4
1	B	214	ARG	3.3
1	A	10	GLU	3.3
1	A	214	ARG	3.2
1	B	57	MET	2.8
1	B	353	MET	2.7
1	A	180	SER	2.6
1	B	350	LYS	2.4
1	A	361	LYS	2.3
1	A	101	SER	2.2
1	A	14	LYS	2.1
1	A	362	TYR	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)

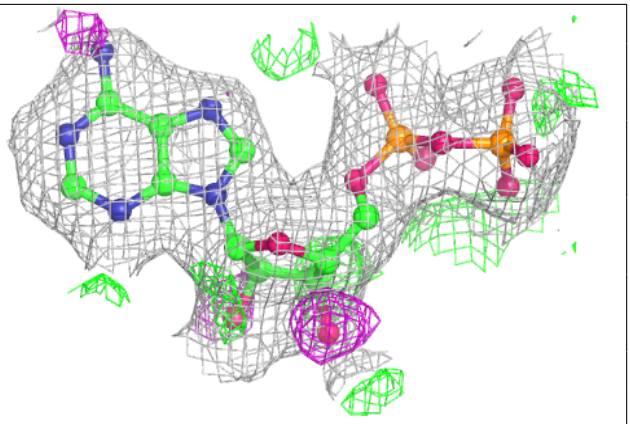
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

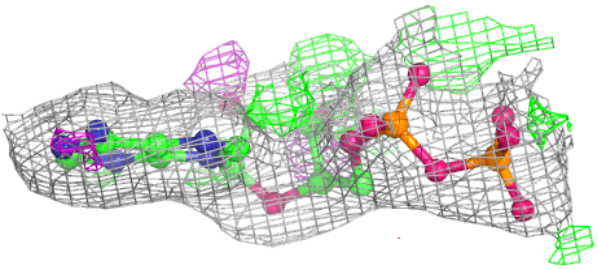
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MG	B	2102	1/1	0.81	0.23	26,26,26,26	0
2	MG	A	1102	1/1	0.84	0.20	28,28,28,28	0
4	ADP	A	1101	27/27	0.94	0.15	18,23,27,29	0
4	ADP	B	2101	27/27	0.94	0.18	18,20,25,26	0
3	SR	A	1103	1/1	0.99	0.19	34,34,34,34	0
3	SR	B	2103	1/1	0.99	0.20	37,37,37,37	0

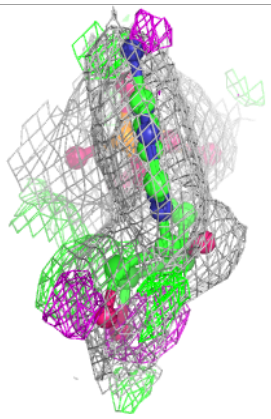
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 1101:

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

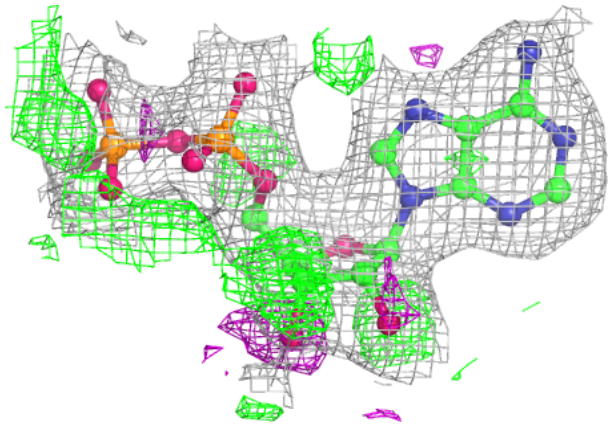


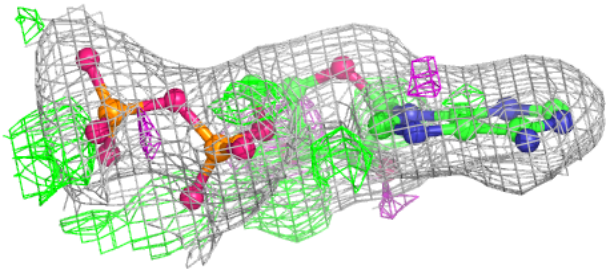


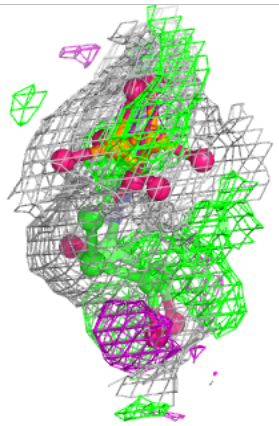


Electron density around ADP B 2101:

$2mF_o$-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.