Full wwPDB X-ray Structure Validation Report (i)

Dec 17, 2023 – 08:15 am GMT

PDB ID	:	4CEI
Title	:	Crystal structure of ADPNP-bound AddAB with a forked DNA substrate
Authors	:	Krajewski, W.W.; Wilkinson, M.; Fu, X.; Cronin, N.B.; Wigley, D.
Deposited on	:	2013-11-11
Resolution	:	2.80 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
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:

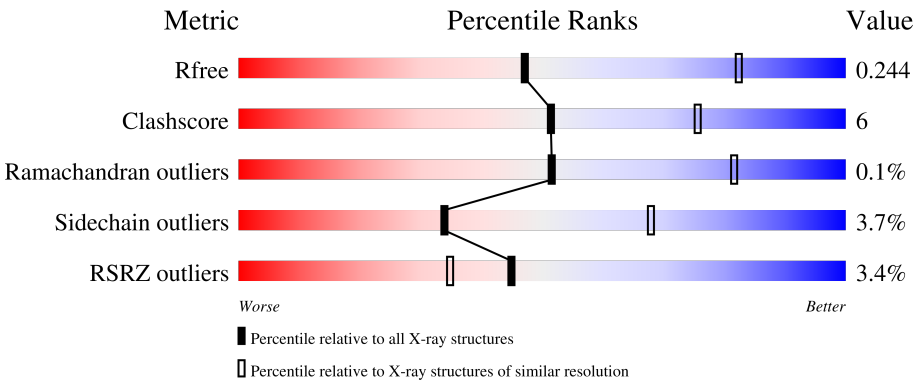
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
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 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	1232	5% 77% 17% • 5%
2	B	1166	% 83% 16% ••
3	X	65	12% 40% 17% 43%

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 19688 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 ATP-DEPENDENT HELICASE/NUCLEASE SUBUNIT A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1170	9468	6040	1611	1790	27	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	780	GLY	ALA	variant	UNP P23478
A	1172	ALA	ASP	engineered mutation	UNP P23478

- Molecule 2 is a protein called ATP-DEPENDENT HELICASE/DEOXYRIBONUCLEASE SUBUNIT B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	1152	9383	5964	1605	1770	44	0	1	0

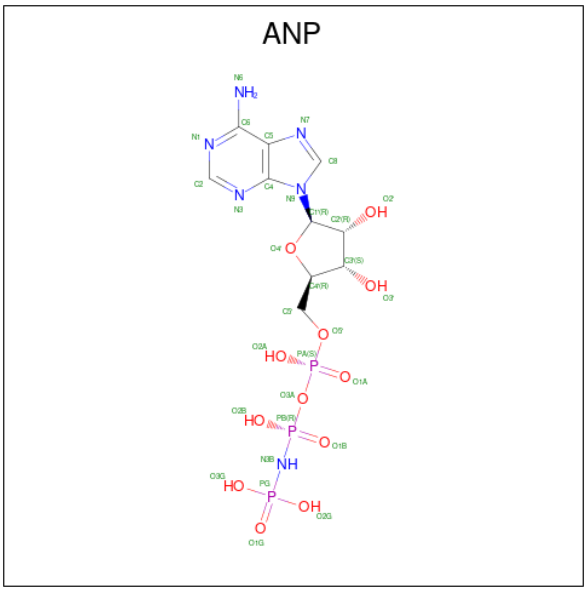
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	843	ASP	GLU	variant	UNP P23477
B	844	GLU	GLN	variant	UNP P23477
B	961	ALA	ASP	engineered mutation	UNP P23477

- Molecule 3 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	X	37	757	363	129	228	37	0	0	0

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).

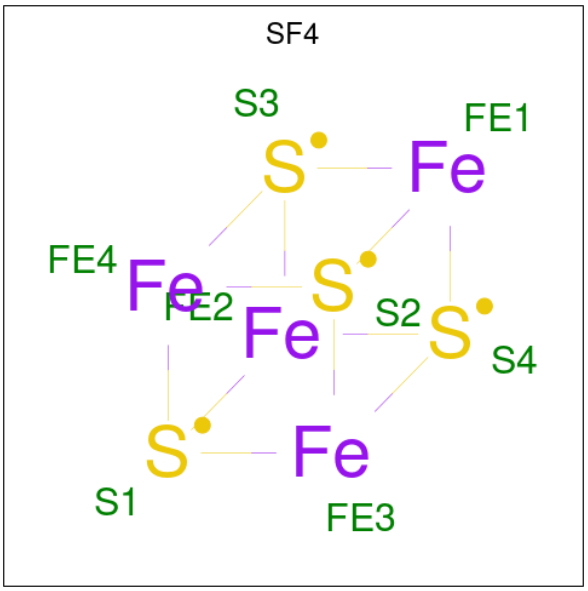


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	31	10	6	12	3	0	0
			Total	C	N	O	P		
4	B	1	31	10	6	12	3	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 6 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Fe	S		
6	B	1	8	4	4	0	0

- Molecule 7 is water.

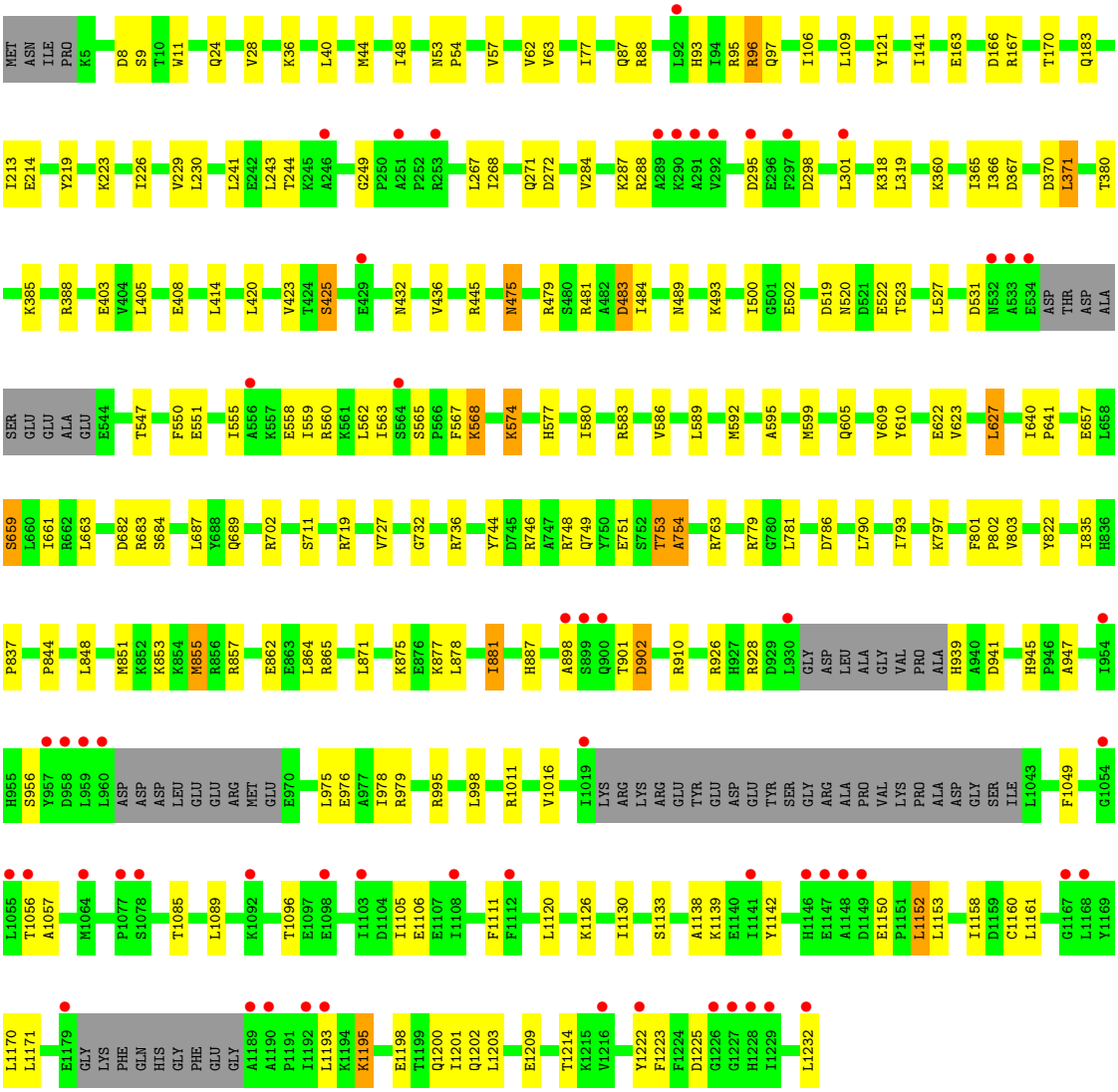
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	4	4	4	0	0
			Total	O		
7	B	4	4	4	0	0

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.

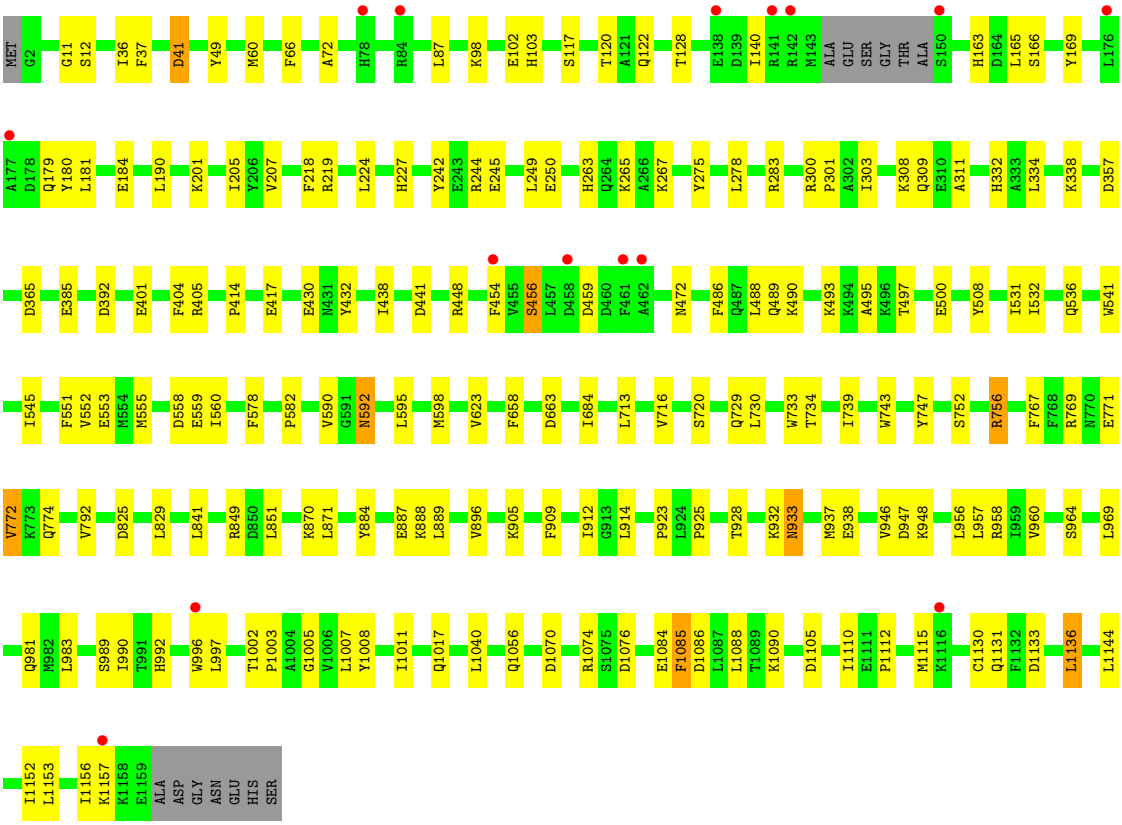
• Molecule 1: ATP-DEPENDENT HELICASE/NUCLEASE SUBUNIT A

Chain A:



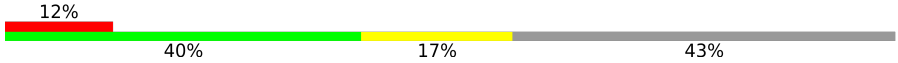
• Molecule 2: ATP-DEPENDENT HELICASE/DEOXYRIBONUCLEASE SUBUNIT B

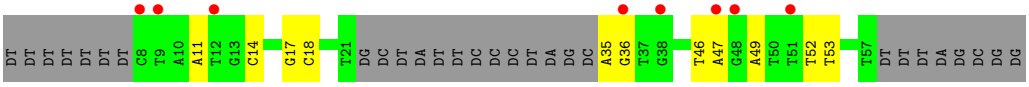
Chain B:



- Molecule 3: DNA

Chain X:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	75.94Å 151.34Å 124.56Å 90.00° 95.95° 90.00°	Depositor
Resolution (Å)	29.95 – 2.80 50.45 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.95-2.80) 99.8 (50.45-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.85 (at 2.81Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.195 , 0.240 0.199 , 0.244	Depositor DCC
R_{free} test set	3459 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	49.8	Xtriage
Anisotropy	0.341	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 39.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	19688	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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: SF4, MG, ANP

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	0.24	0/9657	0.42	0/13034
2	B	0.24	0/9577	0.41	0/12904
3	X	0.60	0/845	1.10	2/1300 (0.2%)
All	All	0.26	0/20079	0.47	2/27238 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	X	49	DA	C3'-C2'-C1'	-5.14	96.33	102.50
3	X	11	DA	C1'-O4'-C4'	-5.05	105.05	110.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	574	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	9468	0	9441	127	1
2	B	9383	0	9329	100	1
3	X	757	0	423	6	0
4	A	31	0	13	0	0
4	B	31	0	13	2	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	8	0	0	1	0
7	A	4	0	0	0	0
7	B	4	0	0	0	0
All	All	19688	0	19219	217	1

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 6.

All (217) 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:A:380:THR:HG21	1:A:388:ARG:HH11	1.46	0.79
2:B:912:ILE:HD11	2:B:948:LYS:HB3	1.67	0.76
2:B:244:ARG:HE	2:B:245:GLU:H	1.34	0.74
1:A:28:VAL:HB	1:A:436:VAL:HG12	1.70	0.72
2:B:896:VAL:HG21	2:B:1011:ILE:HG23	1.72	0.71
1:A:902:ASP:N	1:A:902:ASP:OD1	2.24	0.70
1:A:360:LYS:NZ	1:A:367:ASP:OD1	2.21	0.70
2:B:12:SER:HB3	2:B:278:LEU:HB3	1.73	0.68
1:A:719:ARG:NH1	2:B:365:ASP:OD1	2.25	0.68
2:B:947:ASP:HB2	2:B:958:ARG:HG2	1.76	0.67
2:B:392:ASP:OD2	2:B:405:ARG:NH2	2.29	0.66
1:A:1198:GLU:OE1	1:A:1202:GLN:NE2	2.29	0.65
2:B:283:ARG:HD3	4:B:2161:ANP:H4'	1.78	0.65
1:A:683:ARG:O	1:A:689:GLN:NE2	2.29	0.65
2:B:592:ASN:ND2	2:B:595:LEU:H	1.97	0.63
1:A:1138:ALA:HA	1:A:1152:LEU:HD11	1.80	0.62
1:A:502:GLU:OE1	1:A:865:ARG:NH2	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:871:LEU:HD22	1:A:878:LEU:HD21	1.82	0.60
2:B:180:TYR:OH	2:B:553:GLU:OE1	2.19	0.60
2:B:459:ASP:O	2:B:870:LYS:NZ	2.35	0.60
2:B:263:HIS:NE2	2:B:275:TYR:OH	2.30	0.59
2:B:456:SER:HB2	2:B:459:ASP:HB2	1.83	0.59
1:A:555:ILE:HG12	1:A:881:ILE:HD12	1.85	0.59
2:B:909:PHE:HB3	2:B:947:ASP:HB3	1.83	0.58
2:B:1112:PRO:HG3	6:B:2160:SF4:S4	2.43	0.58
2:B:334:LEU:HD23	2:B:338:LYS:HG3	1.85	0.58
1:A:166:ASP:OD1	2:B:884:TYR:OH	2.20	0.57
1:A:95:ARG:HD2	2:B:249:LEU:HD13	1.87	0.56
2:B:140:ILE:HG22	2:B:166:SER:HB3	1.86	0.56
2:B:925:PRO:HG2	2:B:992:HIS:CE1	2.40	0.56
1:A:567:PHE:HB3	1:A:580:ILE:HB	1.87	0.56
2:B:441:ASP:OD1	2:B:441:ASP:N	2.35	0.56
2:B:414:PRO:HB2	2:B:417:GLU:HG3	1.88	0.56
1:A:170:THR:O	2:B:888:LYS:NZ	2.37	0.55
1:A:481:ARG:HG2	1:A:483:ASP:H	1.71	0.55
1:A:1150:GLU:HB3	2:B:49:TYR:OH	2.05	0.55
1:A:288:ARG:HH12	3:X:17:DG:H21	1.52	0.55
1:A:939:HIS:CE1	1:A:941:ASP:HB3	2.41	0.55
3:X:35:DA:H2"	3:X:36:DG:C8	2.42	0.55
3:X:46:DT:H2"	3:X:47:DA:C8	2.41	0.55
1:A:268:ILE:HA	1:A:271:GLN:HG3	1.88	0.55
1:A:599:MET:HG2	1:A:609:VAL:HG11	1.89	0.55
2:B:495:ALA:HB1	2:B:500:GLU:HG3	1.89	0.54
1:A:748:ARG:NH1	2:B:357:ASP:OD2	2.41	0.54
1:A:167:ARG:NH1	1:A:835:ILE:O	2.40	0.54
1:A:298:ASP:HB3	1:A:301:LEU:HB2	1.90	0.54
2:B:990:ILE:HG13	2:B:1003:PRO:HG3	1.90	0.54
1:A:523:THR:HA	1:A:877:LYS:HD3	1.89	0.53
2:B:488:LEU:HD13	2:B:508:TYR:HB2	1.91	0.53
2:B:928:THR:HG22	2:B:938:GLU:HG3	1.90	0.53
1:A:54:PRO:HB2	1:A:93:HIS:CE1	2.44	0.52
1:A:520:ASN:HB2	1:A:523:THR:HG23	1.91	0.52
1:A:551:GLU:O	1:A:555:ILE:HG13	2.10	0.52
1:A:366:ILE:HB	1:A:370:ASP:HB2	1.91	0.52
1:A:622:GLU:HG2	1:A:736:ARG:HB3	1.91	0.52
1:A:141:ILE:HD12	1:A:365:ILE:HD11	1.90	0.52
1:A:403:GLU:HA	1:A:432:ASN:HB2	1.91	0.52
1:A:214:GLU:HA	1:A:219:TYR:CD2	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:849:ARG:HG2	2:B:905:LYS:HG2	1.92	0.52
1:A:558:GLU:O	1:A:562:LEU:HG	2.10	0.51
1:A:1133:SER:HB3	1:A:1153:LEU:HD11	1.91	0.51
2:B:36:ILE:HG23	2:B:66:PHE:HD2	1.75	0.51
2:B:829:LEU:HD22	2:B:889:LEU:HD13	1.93	0.51
2:B:658:PHE:HZ	2:B:684:ILE:HD12	1.76	0.51
1:A:230:LEU:HD12	1:A:271:GLN:HG2	1.93	0.51
1:A:1011:ARG:HD3	2:B:582:PRO:HB2	1.93	0.51
2:B:448:ARG:NH2	2:B:472:ASN:OD1	2.44	0.51
2:B:536:GLN:HB2	2:B:578:PHE:CE1	2.46	0.51
1:A:425:SER:OG	1:A:432:ASN:ND2	2.44	0.51
1:A:531:ASP:O	1:A:956:SER:OG	2.19	0.51
1:A:711:SER:N	1:A:751:GLU:OE2	2.32	0.51
2:B:825:ASP:HB3	2:B:871:LEU:HD11	1.92	0.51
1:A:1016:VAL:HG11	1:A:1200:GLN:HG3	1.92	0.50
2:B:98:LYS:HD2	2:B:555:MET:HG3	1.93	0.50
2:B:165:LEU:HD11	2:B:623:VAL:HG11	1.93	0.50
1:A:1120:LEU:O	1:A:1126:LYS:HD3	2.10	0.50
1:A:48:ILE:HD13	1:A:57:VAL:HG22	1.94	0.50
2:B:489:GLN:HG2	2:B:493:LYS:HE3	1.94	0.50
2:B:72:ALA:HA	2:B:190:LEU:HD13	1.93	0.49
2:B:960:VAL:HG22	2:B:1007:LEU:HB2	1.95	0.49
1:A:267:LEU:HD11	1:A:284:VAL:HG21	1.93	0.49
1:A:567:PHE:HA	1:A:568:LYS:HE3	1.95	0.49
1:A:500:ILE:HB	1:A:910:ARG:NH1	2.28	0.48
1:A:1111:PHE:HB2	1:A:1222:TYR:CZ	2.48	0.48
2:B:730:LEU:O	2:B:734:THR:HG23	2.13	0.48
2:B:432:TYR:CE2	2:B:438:ILE:HD11	2.48	0.48
1:A:445:ARG:NH1	1:A:862:GLU:OE1	2.46	0.48
1:A:163:GLU:OE1	1:A:167:ARG:NH2	2.38	0.48
2:B:303:ILE:O	2:B:303:ILE:HG13	2.14	0.48
2:B:1115:MET:HB2	2:B:1144:LEU:HB2	1.95	0.48
1:A:763:ARG:NE	2:B:401:GLU:OE1	2.46	0.48
1:A:802:PRO:HA	1:A:875:LYS:HB2	1.95	0.48
2:B:983:LEU:HD13	2:B:1088:LEU:HB3	1.96	0.47
1:A:1160:CYS:HB3	1:A:1171:LEU:HB3	1.96	0.47
1:A:901:THR:HB	1:A:939:HIS:HD2	1.79	0.47
2:B:531:ILE:HD12	2:B:532:ILE:HG13	1.96	0.47
1:A:975:LEU:HD22	2:B:767:PHE:HE1	1.80	0.47
1:A:183:GLN:NE2	1:A:822:TYR:OH	2.49	0.46
1:A:560:ARG:NH1	1:A:605:GLN:HB3	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1105:ILE:HD12	1:A:1105:ILE:H	1.81	0.46
1:A:241:LEU:O	1:A:244:THR:OG1	2.21	0.46
1:A:1195:LYS:NZ	1:A:1198:GLU:HG3	2.30	0.46
1:A:77:ILE:HD13	1:A:106:ILE:HD13	1.97	0.46
2:B:933:ASN:O	2:B:933:ASN:ND2	2.49	0.46
2:B:912:ILE:HD13	2:B:997:LEU:HD11	1.97	0.46
2:B:1074:ARG:HG2	3:X:14:DC:H5"	1.98	0.46
1:A:484:ILE:HD12	1:A:878:LEU:HB2	1.96	0.46
1:A:746:ARG:HA	1:A:749:GLN:HG2	1.97	0.46
2:B:128:THR:HG23	2:B:184:GLU:HG3	1.98	0.46
2:B:497:THR:HG22	2:B:559:GLU:HG2	1.96	0.46
2:B:713:LEU:O	2:B:716:VAL:HG22	2.15	0.45
1:A:226:ILE:HG12	1:A:319:LEU:HD11	1.98	0.45
1:A:531:ASP:OD2	1:A:887:HIS:N	2.49	0.45
1:A:803:VAL:HG22	1:A:877:LYS:HB2	1.98	0.45
2:B:102:GLU:HG3	2:B:103:HIS:ND1	2.31	0.45
2:B:969:LEU:HB2	2:B:1008:TYR:HE1	1.82	0.45
1:A:623:VAL:O	1:A:627:LEU:HB2	2.17	0.45
1:A:1201:ILE:HG13	1:A:1232:LEU:HD12	1.98	0.45
2:B:117:SER:O	2:B:120:THR:HG22	2.17	0.45
1:A:53:ASN:OD1	1:A:53:ASN:N	2.50	0.45
1:A:489:ASN:O	1:A:493:LYS:HG3	2.17	0.45
1:A:93:HIS:O	1:A:97:GLN:HG2	2.16	0.45
1:A:547:THR:HA	1:A:550:PHE:HD2	1.81	0.45
1:A:1085:THR:O	1:A:1089:LEU:HG	2.17	0.45
1:A:1139:LYS:HA	1:A:1142:TYR:O	2.17	0.45
1:A:371:LEU:HD12	1:A:371:LEU:HA	1.89	0.45
1:A:627:LEU:HA	1:A:627:LEU:HD12	1.82	0.45
1:A:657:GLU:HB3	1:A:687:LEU:HD13	1.99	0.44
1:A:978:ILE:HA	2:B:747:TYR:OH	2.16	0.44
1:A:243:LEU:O	1:A:249:GLY:HA3	2.17	0.44
2:B:792:VAL:HG11	2:B:981:GLN:HA	1.98	0.44
1:A:229:VAL:HG21	1:A:319:LEU:HD13	1.99	0.44
2:B:11:GLY:H	4:B:2161:ANP:HNB1	1.64	0.44
2:B:592:ASN:HD21	2:B:595:LEU:H	1.63	0.44
1:A:9:SER:HB3	1:A:11:TRP:H	1.83	0.44
1:A:732:GLY:O	1:A:736:ARG:HG3	2.17	0.44
2:B:716:VAL:O	2:B:716:VAL:HG23	2.17	0.44
1:A:851:MET:O	1:A:855:MET:HB2	2.17	0.44
2:B:772:VAL:HG11	2:B:1110:ILE:HG13	2.00	0.44
1:A:475:ASN:HD22	1:A:475:ASN:HA	1.57	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:753:THR:HG22	1:A:754:ALA:N	2.32	0.44
2:B:1084:GLU:HG2	2:B:1153:LEU:HD13	1.99	0.44
1:A:975:LEU:HD23	1:A:975:LEU:HA	1.88	0.43
2:B:385:GLU:OE2	2:B:405:ARG:HD2	2.17	0.43
1:A:682:ASP:OD2	1:A:684:SER:OG	2.27	0.43
2:B:912:ILE:HG21	2:B:997:LEU:HD21	1.99	0.43
1:A:555:ILE:O	1:A:559:ILE:HG13	2.19	0.43
1:A:622:GLU:CD	1:A:622:GLU:H	2.21	0.43
2:B:486:PHE:O	2:B:490:LYS:HG2	2.19	0.43
1:A:40:LEU:HG	1:A:405:LEU:HD23	2.01	0.43
1:A:595:ALA:O	1:A:599:MET:HG3	2.18	0.43
1:A:1161:LEU:HD23	1:A:1170:LEU:HA	2.00	0.43
2:B:946:VAL:HG21	2:B:996:TRP:HH2	1.84	0.43
1:A:44:MET:SD	1:A:62:VAL:HG21	2.59	0.43
1:A:975:LEU:HD22	2:B:767:PHE:CE1	2.53	0.43
2:B:957:LEU:HD23	2:B:989:SER:HB3	2.00	0.43
1:A:519:ASP:OD1	1:A:520:ASN:N	2.52	0.43
1:A:657:GLU:O	1:A:661:ILE:HG13	2.19	0.43
2:B:729:GLN:HG3	2:B:743:TRP:CD1	2.54	0.43
1:A:36:LYS:HB3	1:A:436:VAL:HB	2.00	0.43
2:B:1005:GLY:HA2	2:B:1085:PHE:CZ	2.54	0.43
2:B:311:ALA:HB1	2:B:663:ASP:O	2.18	0.43
2:B:774:GLN:NE2	2:B:1105:ASP:O	2.52	0.43
2:B:1040:LEU:HD12	2:B:1040:LEU:HA	1.83	0.43
2:B:560:ILE:HA	2:B:560:ILE:HD12	1.76	0.43
1:A:583:ARG:HB2	1:A:786:ASP:OD1	2.19	0.42
1:A:610:TYR:CE1	1:A:779:ARG:HD2	2.53	0.42
2:B:267:LYS:HE3	2:B:267:LYS:HB2	1.87	0.42
2:B:308:LYS:HG2	2:B:309:GLN:N	2.34	0.42
2:B:201:LYS:HA	2:B:227:HIS:O	2.19	0.42
1:A:844:PRO:HB3	1:A:848:LEU:HD23	2.00	0.42
1:A:864:LEU:HD12	1:A:864:LEU:HA	1.90	0.42
1:A:1138:ALA:HA	1:A:1152:LEU:CD1	2.48	0.42
2:B:541:TRP:O	2:B:545:ILE:HG12	2.18	0.42
1:A:408:GLU:OE1	1:A:797:LYS:HE2	2.20	0.42
1:A:1056:THR:OG1	1:A:1057:ALA:N	2.52	0.42
2:B:1130:CYS:SG	2:B:1131:GLN:N	2.92	0.42
1:A:711:SER:HB2	1:A:744:TYR:CE1	2.55	0.42
2:B:733:TRP:HB2	2:B:739:ILE:HD13	2.02	0.42
1:A:95:ARG:HG3	1:A:96:ARG:N	2.34	0.42
2:B:1152:ILE:O	2:B:1156:ILE:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:586:VAL:HG21	1:A:801:PHE:CD1	2.56	0.41
1:A:589:LEU:HD11	1:A:790:LEU:HD22	2.02	0.41
1:A:945:HIS:C	1:A:947:ALA:H	2.22	0.41
2:B:163:HIS:O	2:B:166:SER:OG	2.33	0.41
1:A:1193:LEU:HD13	1:A:1223:PHE:CZ	2.55	0.41
1:A:385:LYS:HE2	1:A:385:LYS:HB3	1.81	0.41
2:B:205:ILE:HD13	2:B:224:LEU:HD13	2.02	0.41
2:B:300:ARG:HA	2:B:301:PRO:HA	1.94	0.41
1:A:121:TYR:OH	2:B:122:GLN:NE2	2.53	0.41
1:A:976:GLU:CD	1:A:979:ARG:HH21	2.23	0.41
1:A:1049:PHE:CD2	2:B:552:VAL:HG21	2.55	0.41
1:A:1209:GLU:HA	1:A:1214:THR:O	2.20	0.41
1:A:287:LYS:HD2	1:A:287:LYS:N	2.36	0.41
1:A:898:ALA:HB1	1:A:926:ARG:HG3	2.02	0.41
2:B:404:PHE:CZ	2:B:430:GLU:HA	2.55	0.41
2:B:87:LEU:HD23	2:B:87:LEU:HA	1.88	0.41
1:A:223:LYS:NZ	1:A:272:ASP:OD1	2.50	0.41
3:X:52:DT:H2"	3:X:53:DT:OP2	2.20	0.41
1:A:63:VAL:HG11	1:A:109:LEU:HD13	2.01	0.41
1:A:589:LEU:O	1:A:793:ILE:HG13	2.21	0.41
1:A:998:LEU:HD23	1:A:998:LEU:HA	1.91	0.41
2:B:41:ASP:OD1	2:B:41:ASP:N	2.52	0.41
1:A:659:SER:O	1:A:663:LEU:HG	2.21	0.41
1:A:995:ARG:NH1	2:B:752:SER:OG	2.54	0.41
2:B:956:LEU:HD23	2:B:1002:THR:HB	2.03	0.41
1:A:166:ASP:HB3	1:A:837:PRO:HB3	2.02	0.40
2:B:1133:ASP:HB3	2:B:1136:LEU:HD22	2.03	0.40
3:X:17:DG:H2"	3:X:18:DC:C6	2.56	0.40
1:A:420:LEU:HA	1:A:423:VAL:HG22	2.03	0.40
1:A:1130:ILE:HB	1:A:1158:ILE:HB	2.03	0.40
2:B:308:LYS:HE2	2:B:308:LYS:HB3	1.87	0.40
2:B:769:ARG:O	2:B:771:GLU:N	2.52	0.40
1:A:213:ILE:HG22	1:A:219:TYR:CD1	2.56	0.40
1:A:640:ILE:HB	1:A:641:PRO:HD3	2.03	0.40
2:B:841:LEU:HD21	2:B:851:LEU:HD21	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:901:THR:OG1	2:B:756:ARG:NH1[2_545]	2.18	0.02

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	1158/1232 (94%)	1120 (97%)	36 (3%)	2 (0%)	47	78
2	B	1149/1166 (98%)	1112 (97%)	36 (3%)	1 (0%)	51	81
All	All	2307/2398 (96%)	2232 (97%)	72 (3%)	3 (0%)	51	81

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	754	ALA
1	A	592	MET
2	B	923	PRO

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	1013/1062 (95%)	975 (96%)	38 (4%)	33	67
2	B	1021/1029 (99%)	984 (96%)	37 (4%)	35	69
All	All	2034/2091 (97%)	1959 (96%)	75 (4%)	34	68

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASP
1	A	24	GLN
1	A	87	GLN

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Mol	Chain	Res	Type
1	A	88	ARG
1	A	96	ARG
1	A	295	ASP
1	A	318	LYS
1	A	371	LEU
1	A	414	LEU
1	A	425	SER
1	A	475	ASN
1	A	479	ARG
1	A	483	ASP
1	A	522	GLU
1	A	527	LEU
1	A	563	ILE
1	A	565	SER
1	A	568	LYS
1	A	574	LYS
1	A	577	HIS
1	A	627	LEU
1	A	659	SER
1	A	702	ARG
1	A	727	VAL
1	A	753	THR
1	A	781	LEU
1	A	853	LYS
1	A	855	MET
1	A	857	ARG
1	A	881	ILE
1	A	902	ASP
1	A	928	ARG
1	A	1096	THR
1	A	1106	GLU
1	A	1152	LEU
1	A	1195	LYS
1	A	1203	LEU
1	A	1225	ASP
2	B	37	PHE
2	B	41	ASP
2	B	60	MET
2	B	169	TYR
2	B	179	GLN
2	B	181	LEU
2	B	207	VAL

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Mol	Chain	Res	Type
2	B	218	PHE
2	B	219	ARG
2	B	242	TYR
2	B	250	GLU
2	B	265	LYS
2	B	332	HIS
2	B	454	PHE
2	B	456	SER
2	B	551	PHE
2	B	558	ASP
2	B	590	VAL
2	B	592	ASN
2	B	598	MET
2	B	720	SER
2	B	756	ARG
2	B	772	VAL
2	B	887	GLU
2	B	914	LEU
2	B	932	LYS
2	B	933	ASN
2	B	937	MET
2	B	964	SER
2	B	1017	GLN
2	B	1056	GLN
2	B	1076	ASP
2	B	1085	PHE
2	B	1086	ASP
2	B	1090	LYS
2	B	1136	LEU
2	B	1157	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	93	HIS
1	A	183	GLN
1	A	475	ASN
1	A	579	ASN
1	A	1069	GLN
2	B	122	GLN
2	B	592	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 5 ligands modelled in this entry, 2 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	$\#\|Z\| > 2$
4	ANP	B	2161	5	29,33,33	1.09	4 (13%)	31,52,52	1.15	5 (16%)
6	SF4	B	2160	2	0,12,12	-	-	-		
4	ANP	A	2233	5	29,33,33	1.09	4 (13%)	31,52,52	1.13	4 (12%)

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	ANP	B	2161	5	-	6/14/38/38	0/3/3/3
6	SF4	B	2160	2	-	-	0/6/5/5
4	ANP	A	2233	5	-	3/14/38/38	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	2161	ANP	PG-O1G	2.67	1.50	1.46
4	A	2233	ANP	PG-O1G	2.57	1.50	1.46
4	A	2233	ANP	PG-N3B	2.42	1.69	1.63
4	A	2233	ANP	PB-O3A	-2.39	1.56	1.59
4	B	2161	ANP	PB-O1B	2.36	1.49	1.46
4	B	2161	ANP	PG-N3B	2.34	1.69	1.63
4	B	2161	ANP	PB-O3A	-2.12	1.56	1.59
4	A	2233	ANP	PB-O1B	2.09	1.49	1.46

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	2233	ANP	PB-O3A-PA	-3.07	121.81	132.62
4	B	2161	ANP	PB-O3A-PA	-2.91	122.36	132.62
4	A	2233	ANP	O3G-PG-O1G	-2.63	106.84	113.45
4	A	2233	ANP	C4-C5-N7	2.61	112.12	109.40
4	A	2233	ANP	C5-C6-N6	2.33	123.89	120.35
4	B	2161	ANP	O1G-PG-N3B	-2.31	108.38	111.77
4	B	2161	ANP	C5-C6-N6	2.27	123.80	120.35
4	B	2161	ANP	C4-C5-N7	2.16	111.65	109.40
4	B	2161	ANP	O3G-PG-O1G	-2.10	108.18	113.45

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2233	ANP	PB-N3B-PG-O1G
4	A	2233	ANP	PG-N3B-PB-O1B
4	A	2233	ANP	PG-N3B-PB-O3A
4	B	2161	ANP	PG-N3B-PB-O1B
4	B	2161	ANP	C5'-O5'-PA-O2A
4	B	2161	ANP	O4'-C4'-C5'-O5'
4	B	2161	ANP	C3'-C4'-C5'-O5'
4	B	2161	ANP	C5'-O5'-PA-O3A
4	B	2161	ANP	C5'-O5'-PA-O1A

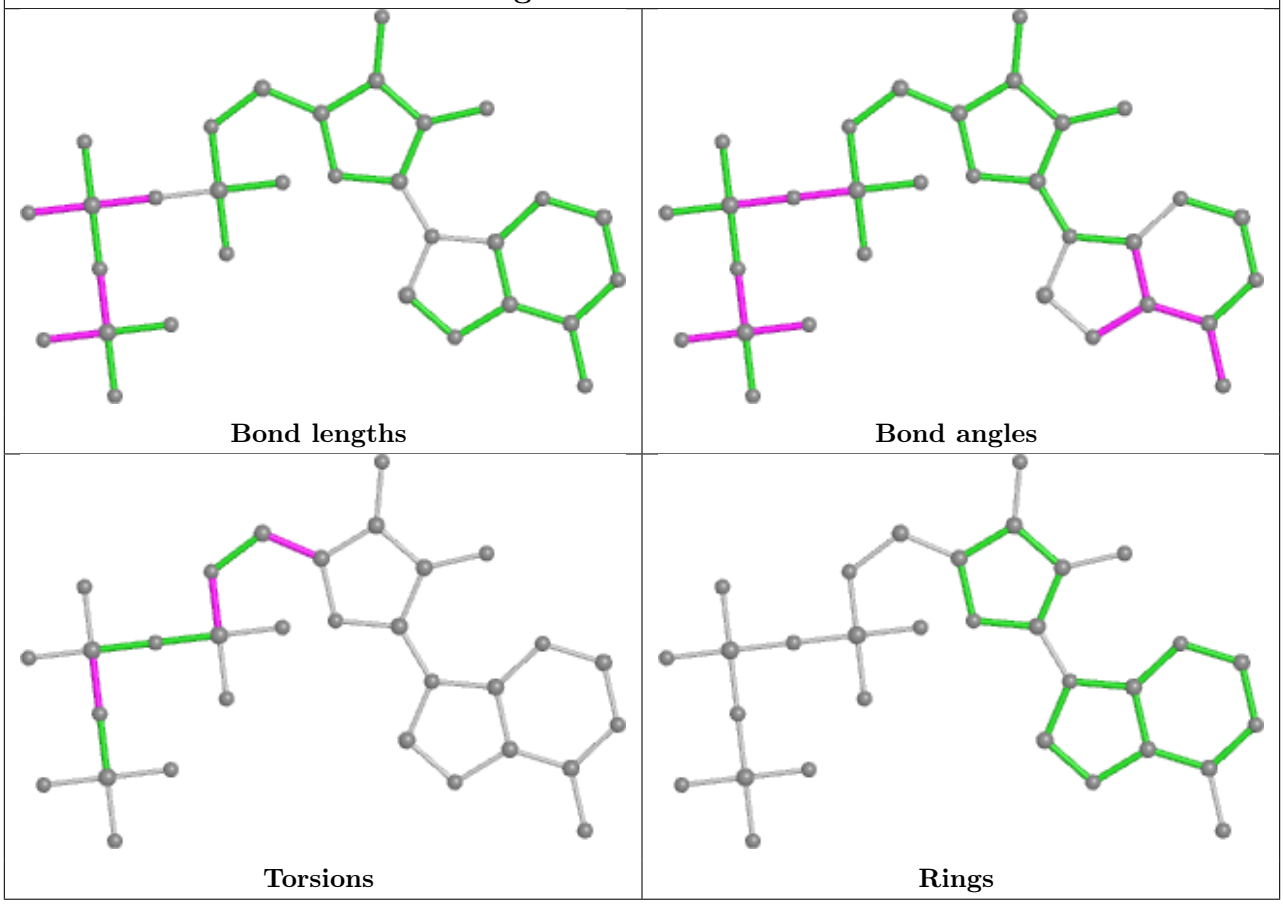
There are no ring outliers.

2 monomers are involved in 3 short contacts:

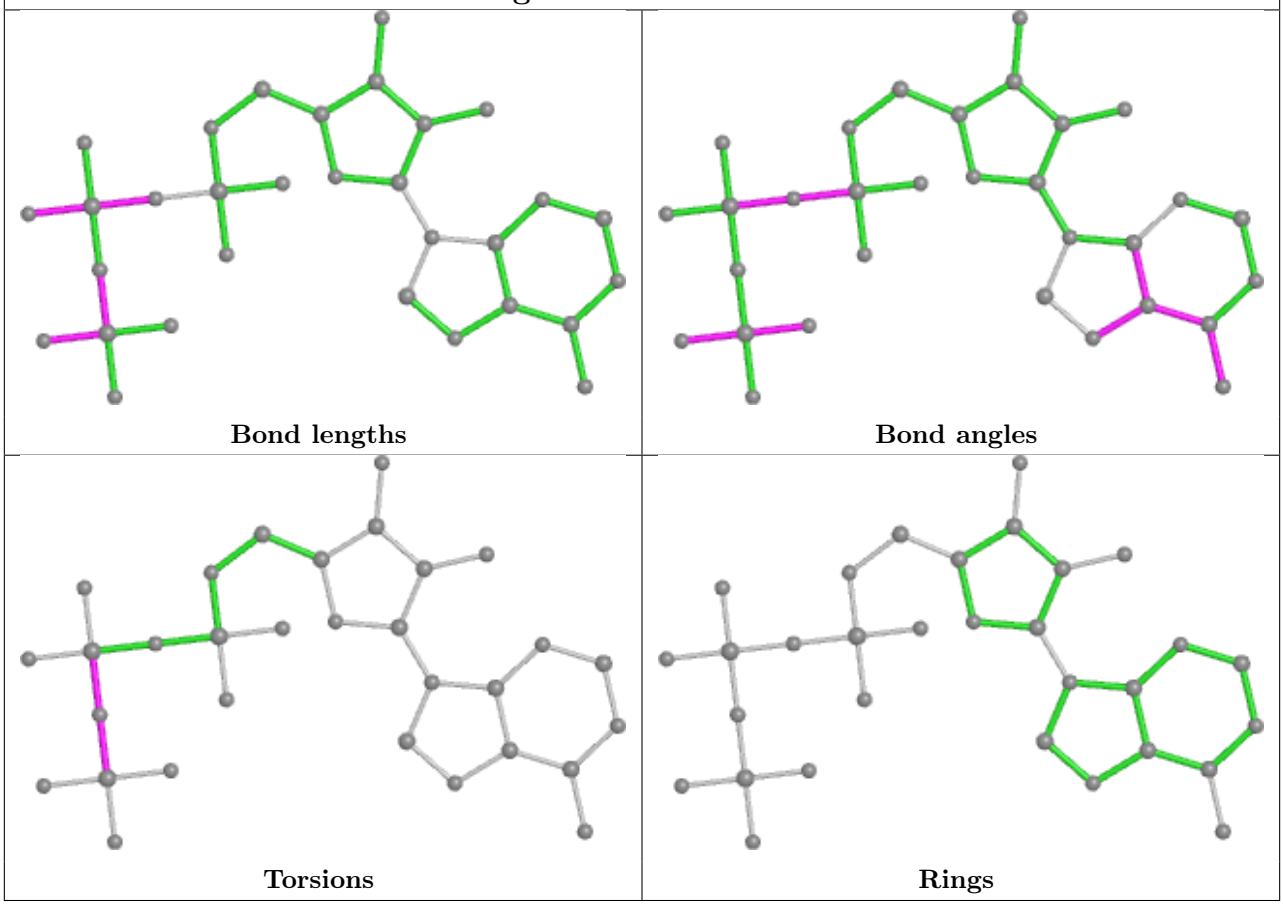
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2161	ANP	2	0
6	B	2160	SF4	1	0

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 ANP B 2161



Ligand ANP A 2233



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	1170/1232 (94%)	-0.01	57 (4%)	29	20	18, 48, 103, 129	0
2	B	1152/1166 (98%)	-0.18	15 (1%)	77	72	20, 48, 82, 122	0
3	X	37/65 (56%)	1.26	8 (21%)	0	0	49, 122, 140, 153	0
All	All	2359/2463 (95%)	-0.07	80 (3%)	45	35	18, 48, 100, 153	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	X	48	DG	4.8
1	A	1148	ALA	4.8
2	B	461	PHE	4.5
3	X	47	DA	4.5
2	B	150	SER	4.3
1	A	1077	PRO	4.2
3	X	8	DC	3.9
1	A	92	LEU	3.9
1	A	1112	PHE	3.8
1	A	1193	LEU	3.7
1	A	1108	ILE	3.7
1	A	960	LEU	3.7
1	A	1147	GLU	3.6
1	A	1179	GLU	3.6
1	A	1216	VAL	3.5
2	B	141	ARG	3.5
2	B	138	GLU	3.4
1	A	1228	HIS	3.3
1	A	1222	TYR	3.2
1	A	1226	GLY	3.2
1	A	1055	LEU	3.1
1	A	1227	GLY	3.1
1	A	954	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	1054	GLY	3.0
1	A	556	ALA	2.9
1	A	301	LEU	2.9
1	A	1229	ILE	2.9
2	B	458	ASP	2.9
1	A	251	ALA	2.8
2	B	996	TRP	2.7
1	A	898	ALA	2.7
3	X	9	DT	2.7
1	A	930	LEU	2.7
2	B	78	HIS	2.7
1	A	899	SER	2.6
1	A	1232	LEU	2.6
1	A	290	LYS	2.6
1	A	1192	ILE	2.6
1	A	1056	THR	2.6
1	A	1098	GLU	2.6
1	A	1146	HIS	2.6
1	A	295	ASP	2.6
3	X	38	DG	2.6
1	A	1167	GLY	2.6
1	A	429	GLU	2.5
1	A	957	TYR	2.5
3	X	51	DT	2.5
1	A	1141	ILE	2.5
1	A	959	LEU	2.4
1	A	1019	ILE	2.4
1	A	289	ALA	2.4
1	A	534	GLU	2.4
2	B	1157	LYS	2.4
1	A	246	ALA	2.4
1	A	900	GLN	2.4
1	A	1103	ILE	2.4
1	A	564	SER	2.4
1	A	533	ALA	2.4
1	A	1168	LEU	2.3
1	A	532	ASN	2.3
2	B	177	ALA	2.3
1	A	1064	MET	2.2
2	B	454	PHE	2.2
1	A	291	ALA	2.2
3	X	36	DG	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	462	ALA	2.2
1	A	1190	ALA	2.2
2	B	84	ARG	2.2
1	A	958	ASP	2.1
2	B	142	ARG	2.1
1	A	1092	LYS	2.1
2	B	1116	LYS	2.1
1	A	253	ARG	2.1
1	A	1149	ASP	2.0
1	A	1189	ALA	2.0
3	X	12	DT	2.0
1	A	297	PHE	2.0
1	A	292	VAL	2.0
2	B	176	LEU	2.0
1	A	1078	SER	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.

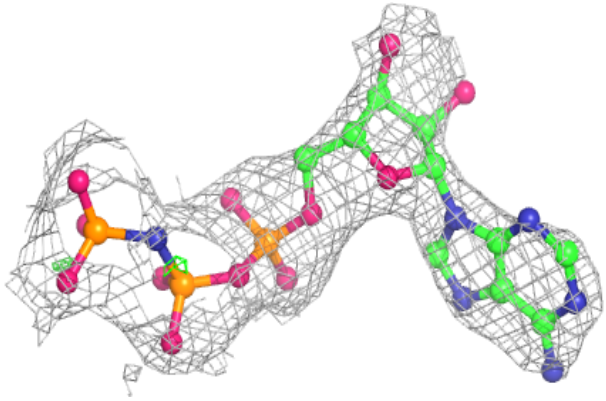
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	MG	B	2162	1/1	0.73	0.14	65,65,65,65	0
4	ANP	B	2161	31/31	0.94	0.17	48,78,95,96	0
4	ANP	A	2233	31/31	0.96	0.14	37,50,61,67	0
5	MG	A	2234	1/1	0.97	0.10	51,51,51,51	0
6	SF4	B	2160	8/8	0.99	0.09	24,36,47,51	0

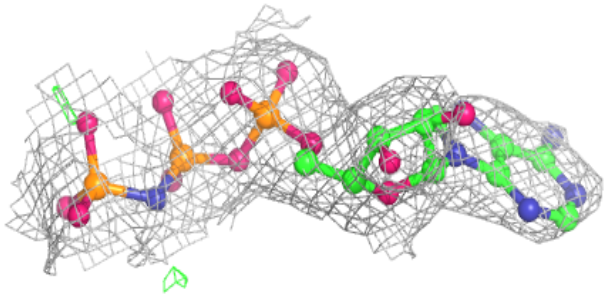
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

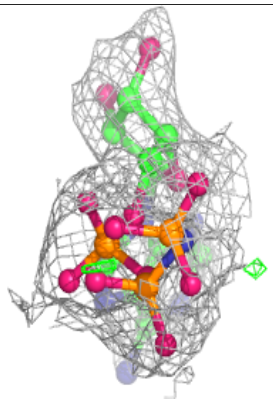
as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ANP B 2161:

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

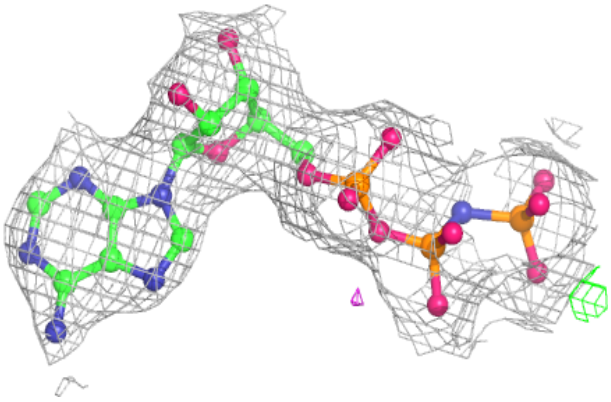


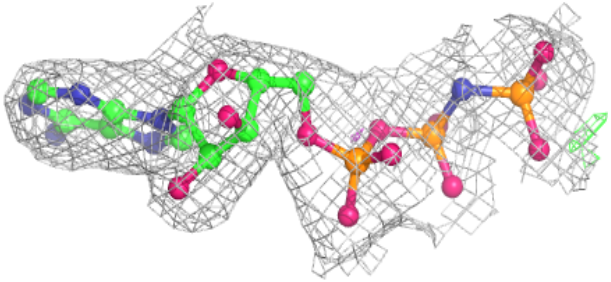


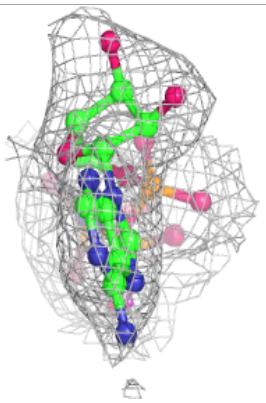


Electron density around ANP A 2233:

$2mF_o$-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.