



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:44 AM UTC

PDB ID : 3VPC / pdb_00003vpc
Title : ArgX from Sulfolobus tokodaii complexed with ADP
Authors : Tomita, T.; Horie, A.; Kuzuyama, T.; Nishiyama, M.
Deposited on : 2012-02-29
Resolution : 1.87 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

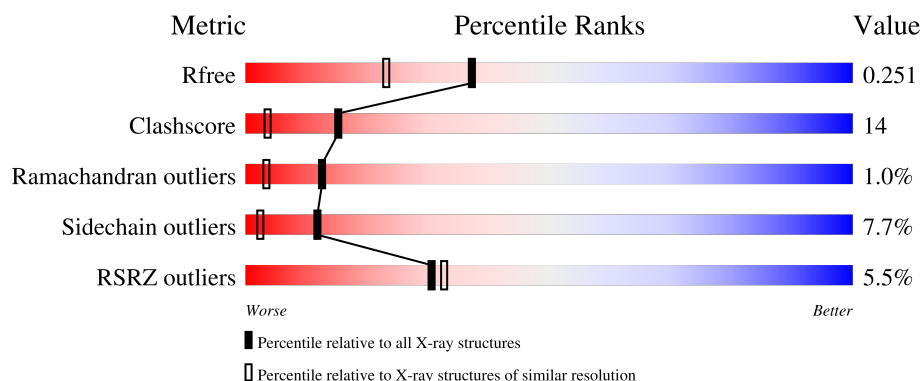
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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}	180053	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	282	8% 76% 18% 5%
1	B	282	4% 77% 20%
1	C	282	5% 77% 16% 5%
1	D	282	6% 75% 20% 5%

2 Entry composition [i](#)

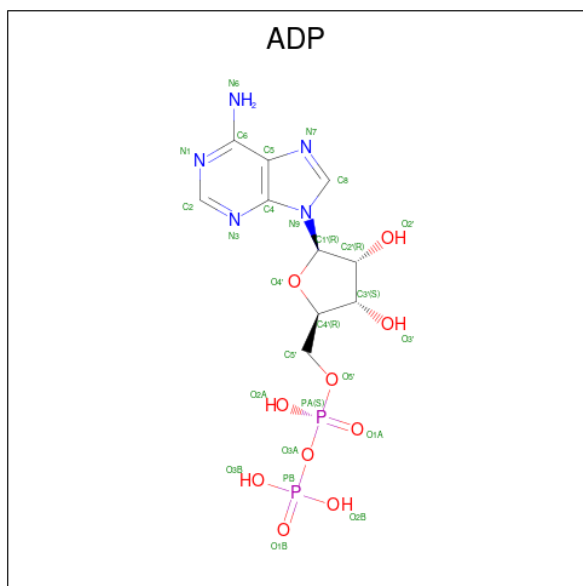
There are 3 unique types of molecules in this entry. The entry contains 10021 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Putative acetylornithine deacetylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	282	Total	C	N	O	S	0	5	0
			2246	1440	382	417	7			
1	B	282	Total	C	N	O	S	0	8	0
			2260	1450	382	421	7			
1	C	275	Total	C	N	O	S	0	7	0
			2213	1420	375	411	7			
1	D	282	Total	C	N	O	S	0	3	0
			2238	1435	379	417	7			

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



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

Continued on next page...

Continued from previous page...

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

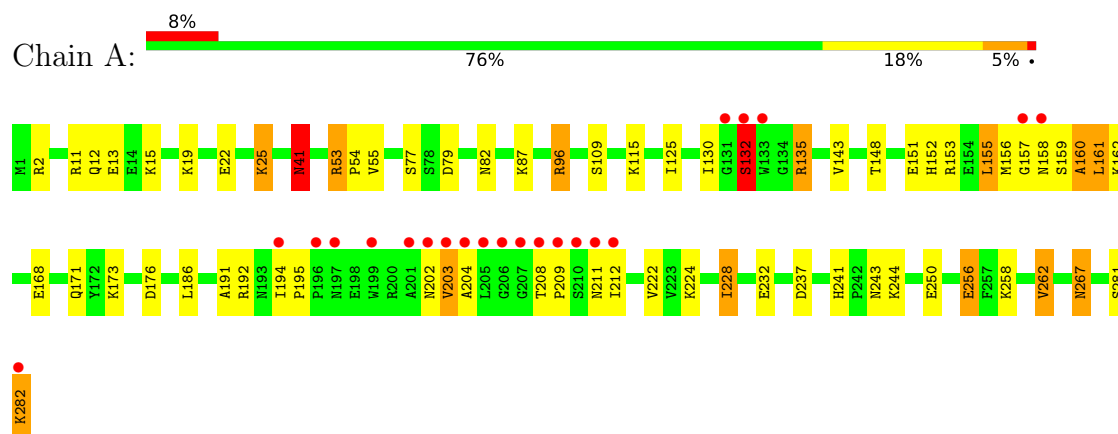
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	247	Total	O	0	0
			247	247		
3	B	270	Total	O	0	0
			270	270		
3	C	221	Total	O	0	0
			221	221		
3	D	218	Total	O	0	0
			218	218		

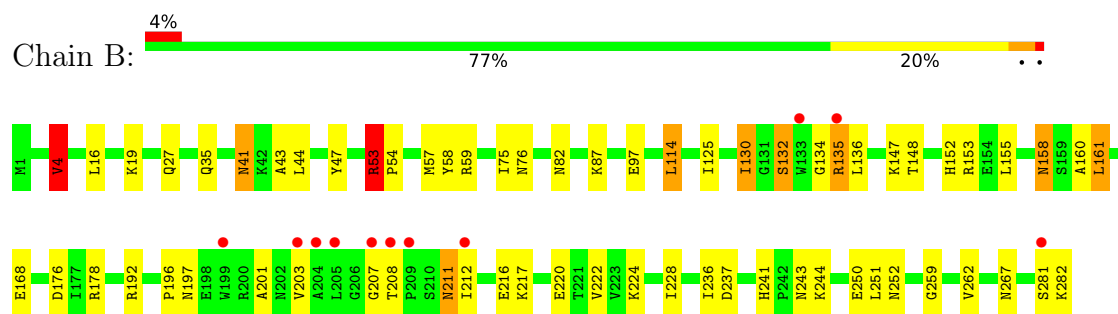
3 Residue-property plots [i](#)

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.

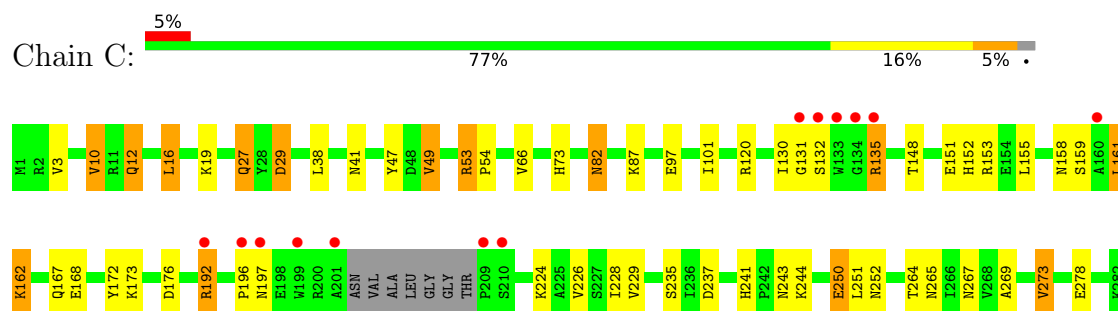
• Molecule 1: Putative acetylornithine deacetylase



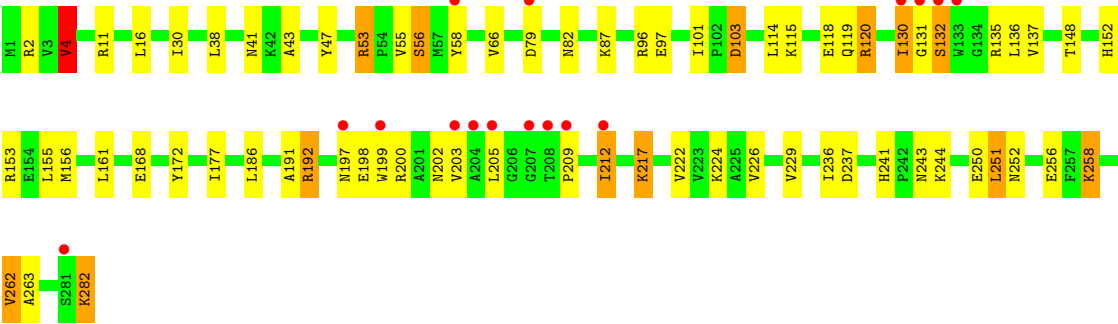
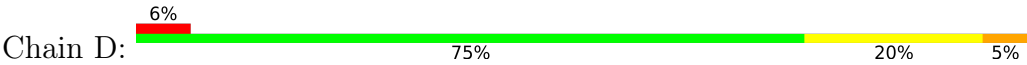
• Molecule 1: Putative acetylornithine deacetylase



• Molecule 1: Putative acetylornithine deacetylase



• Molecule 1: Putative acetylornithine deacetylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	63.27Å 67.84Å 80.03Å 87.77° 75.29° 67.03°	Depositor
Resolution (Å)	34.96 – 1.87 34.96 – 1.87	Depositor EDS
% Data completeness (in resolution range)	95.6 (34.96-1.87) 96.0 (34.96-1.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 1.87Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.198 , 0.253 0.197 , 0.251	Depositor DCC
R_{free} test set	4915 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10021	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.68	0/2301	0.91	4/3117 (0.1%)
1	B	0.69	0/2321	0.90	4/3146 (0.1%)
1	C	0.68	1/2273 (0.0%)	0.88	3/3076 (0.1%)
1	D	0.70	0/2287	0.97	5/3099 (0.2%)
All	All	0.69	1/9182 (0.0%)	0.92	16/12438 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	10	VAL	CA-CB	5.36	1.59	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	53	ARG	CA-C-N	-7.24	112.31	119.76
1	D	53	ARG	C-N-CA	-7.24	112.31	119.76
1	B	4	VAL	CB-CA-C	-6.93	100.27	110.81
1	D	4	VAL	CB-CA-C	-6.42	101.27	110.83
1	A	41	ASN	N-CA-C	5.86	116.02	108.34
1	C	101	ILE	CA-C-N	-5.81	113.78	119.76
1	C	101	ILE	C-N-CA	-5.81	113.78	119.76
1	B	53	ARG	CA-C-N	-5.77	113.84	119.78
1	B	53	ARG	C-N-CA	-5.77	113.84	119.78
1	A	195	PRO	CA-C-N	5.38	126.56	119.84
1	A	195	PRO	C-N-CA	5.38	126.56	119.84
1	D	101	ILE	CA-C-N	-5.17	114.59	119.76
1	D	101	ILE	C-N-CA	-5.17	114.59	119.76
1	A	160	ALA	N-CA-C	-5.16	106.96	113.20
1	B	134	GLY	N-CA-C	-5.12	108.55	115.36
1	C	159	SER	N-CA-C	5.06	116.88	111.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2246	0	2316	75	0
1	B	2260	0	2331	71	0
1	C	2213	0	2282	66	0
1	D	2238	0	2299	60	0
2	A	27	0	12	2	0
2	B	27	0	12	2	0
2	C	27	0	12	2	0
2	D	27	0	12	0	0
3	A	247	0	0	9	0
3	B	270	0	0	7	0
3	C	221	0	0	7	0
3	D	218	0	0	7	0
All	All	10021	0	9276	249	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 14.

All (249) 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:135:ARG:HD3	1:B:136:LEU:H	1.10	1.16
1:C:135:ARG:HG2	1:C:135:ARG:HH21	1.06	1.13
1:B:135:ARG:CD	1:B:136:LEU:H	1.67	1.06
1:A:96[A]:ARG:HH11	1:A:96[A]:ARG:HG2	0.89	1.06
1:C:192[B]:ARG:HH21	1:C:192[B]:ARG:HG2	0.90	1.03
1:B:135:ARG:HD3	1:B:136:LEU:N	1.73	1.02
1:A:132:SER:HB3	1:A:135:ARG:HB2	1.43	0.98
1:A:153:ARG:HE	1:D:152:HIS:HE1	1.09	0.96
1:B:87:LYS:NZ	1:B:250:GLU:HG2	1.80	0.96
1:B:152:HIS:HE1	1:C:153:ARG:HE	1.07	0.96
1:C:192[B]:ARG:HG2	1:C:192[B]:ARG:NH2	1.66	0.96

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96[A]:ARG:HG2	1:A:96[A]:ARG:NH1	1.67	0.96
1:C:132:SER:O	1:C:135:ARG:HB2	1.66	0.95
1:A:96[A]:ARG:HH11	1:A:96[A]:ARG:CG	1.80	0.94
1:B:87:LYS:NZ	1:B:250:GLU:CG	2.32	0.92
1:C:192[B]:ARG:HH21	1:C:192[B]:ARG:CG	1.81	0.92
1:B:41:ASN:C	1:B:41:ASN:HD22	1.79	0.91
1:B:87:LYS:HZ3	1:B:250:GLU:CD	1.79	0.90
1:A:152:HIS:HE1	1:D:153:ARG:HE	1.13	0.89
1:B:153:ARG:HE	1:C:152:HIS:HE1	1.16	0.89
1:A:153:ARG:HE	1:D:152:HIS:CE1	1.93	0.85
1:A:132:SER:HB3	1:A:135:ARG:CB	2.07	0.85
1:C:176[A]:ASP:OD2	1:C:192[A]:ARG:NH1	2.09	0.84
1:C:224:LYS:O	1:C:228:ILE:HD13	1.76	0.84
1:C:135:ARG:HG2	1:C:135:ARG:NH2	1.86	0.84
1:B:241:HIS:HD2	1:B:243:ASN:H	1.25	0.83
1:D:198:GLU:HG2	1:D:199:TRP:H	1.43	0.82
1:A:132:SER:O	1:A:135:ARG:HB2	1.79	0.81
1:B:87:LYS:NZ	1:B:250:GLU:CD	2.38	0.81
1:D:87:LYS:HZ3	1:D:250:GLU:HG2	1.44	0.81
1:A:87:LYS:HZ3	1:A:250:GLU:HG2	1.46	0.81
1:C:97:GLU:HG2	1:C:228:ILE:HG12	1.62	0.81
1:A:162:LYS:NZ	1:B:59:ARG:HH11	1.78	0.80
1:A:87:LYS:NZ	1:A:250:GLU:HG2	1.99	0.78
1:A:115:LYS:HE3	1:B:35:GLN:O	1.84	0.78
1:B:241:HIS:CD2	1:B:243:ASN:H	2.01	0.77
1:C:135:ARG:HH21	1:C:135:ARG:CG	1.94	0.76
1:B:224:LYS:HD3	3:B:776:HOH:O	1.86	0.76
1:B:152:HIS:CE1	1:C:153:ARG:HE	1.99	0.75
1:B:87:LYS:HZ2	1:B:250:GLU:HG2	1.52	0.75
1:D:252:ASN:HB2	3:D:608:HOH:O	1.88	0.74
1:A:153:ARG:NE	1:D:152:HIS:HE1	1.85	0.73
1:A:152:HIS:CE1	1:D:153:ARG:HE	2.03	0.73
1:A:12:GLN:HA	1:A:15:LYS:HE2	1.72	0.72
1:C:87:LYS:HZ3	1:C:250:GLU:HG2	1.55	0.71
1:D:87:LYS:NZ	1:D:250:GLU:HG2	2.06	0.70
1:D:132:SER:OG	1:D:135:ARG:HB2	1.91	0.70
1:A:168:GLU:OE1	1:A:241:HIS:HE1	1.74	0.70
1:A:241:HIS:CD2	1:A:244:LYS:H	2.11	0.69
1:A:241:HIS:HD2	1:A:244:LYS:H	1.37	0.69
1:D:130:ILE:O	3:D:695:HOH:O	2.10	0.68
1:B:135:ARG:HE	1:C:151:GLU:HG3	1.57	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:SER:O	1:A:282:LYS:HB3	1.93	0.68
1:B:87:LYS:HZ3	1:B:250:GLU:CG	2.00	0.67
1:A:151:GLU:OE2	1:D:135:ARG:HA	1.95	0.66
1:C:132:SER:HB2	3:C:605:HOH:O	1.95	0.66
1:B:41:ASN:C	1:B:41:ASN:ND2	2.48	0.66
1:C:73:HIS:HE1	3:C:729:HOH:O	1.77	0.66
1:D:41:ASN:ND2	1:D:43:ALA:HB3	2.11	0.66
1:C:38:LEU:HB3	1:C:66:VAL:HG11	1.78	0.65
1:C:168:GLU:OE1	1:C:241:HIS:HE1	1.80	0.65
1:B:241:HIS:CD2	1:B:244:LYS:H	2.15	0.65
1:B:252:ASN:HB2	3:B:791:HOH:O	1.97	0.65
2:B:500:ADP:O2B	3:B:620:HOH:O	2.15	0.64
1:B:41:ASN:HD21	1:B:44:LEU:H	1.46	0.64
1:A:162:LYS:HZ2	1:B:59:ARG:NH1	1.96	0.64
1:B:135:ARG:HD2	1:B:136:LEU:H	1.61	0.63
1:C:12:GLN:HE21	1:C:12:GLN:HA	1.62	0.63
1:A:176[B]:ASP:OD2	2:A:500:ADP:O3'	2.16	0.62
1:A:87:LYS:NZ	1:A:250:GLU:CG	2.62	0.62
1:D:87:LYS:NZ	1:D:250:GLU:CG	2.63	0.62
1:A:191:ALA:HB2	1:A:212:ILE:HD12	1.82	0.62
1:B:148:THR:O	1:B:152:HIS:HD2	1.81	0.62
1:D:237:ASP:HB2	1:D:250:GLU:HB2	1.81	0.62
1:D:241:HIS:HD2	1:D:243:ASN:H	1.50	0.60
1:A:96[A]:ARG:NH1	1:A:96[A]:ARG:CG	2.47	0.60
1:D:241:HIS:CD2	1:D:243:ASN:H	2.20	0.60
1:B:135:ARG:NE	1:C:151:GLU:HG3	2.16	0.60
1:D:87:LYS:HE3	3:D:715:HOH:O	2.02	0.59
1:A:162:LYS:NZ	1:B:59:ARG:NH1	2.46	0.59
1:D:87:LYS:HZ3	1:D:250:GLU:CG	2.16	0.59
1:A:115:LYS:CE	1:B:35:GLN:O	2.50	0.59
1:C:241:HIS:HD2	1:C:243[B]:ASN:H	1.50	0.59
1:A:132:SER:O	1:A:135:ARG:HD3	2.03	0.59
1:A:237:ASP:HB2	1:A:250:GLU:HB2	1.84	0.59
1:B:132:SER:OG	1:B:135:ARG:O	2.20	0.58
1:C:241:HIS:HD2	1:C:243[A]:ASN:H	1.50	0.58
1:D:192:ARG:HH21	1:D:203:VAL:HG23	1.66	0.58
1:D:217:LYS:HE2	1:D:217:LYS:H	1.67	0.58
1:B:135:ARG:HG3	1:C:151:GLU:HB3	1.85	0.58
1:A:258:LYS:O	1:A:262:VAL:HG13	2.03	0.58
1:D:258:LYS:O	1:D:262:VAL:HG13	2.04	0.58
1:C:241:HIS:CD2	1:C:244:LYS:H	2.22	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ARG:HE	1:C:152:HIS:CE1	2.08	0.57
1:A:25:LYS:HD3	1:A:25:LYS:N	2.21	0.56
1:A:202:ASN:HB2	1:A:203:VAL:HA	1.87	0.56
1:C:241:HIS:HD2	1:C:244:LYS:H	1.54	0.56
3:C:650:HOH:O	1:D:79:ASP:HB3	2.06	0.55
1:C:172:TYR:CD1	1:C:173:LYS:HD2	2.42	0.55
1:A:96[A]:ARG:CZ	3:A:711:HOH:O	2.53	0.55
1:D:209:PRO:HB2	1:D:263:ALA:HB2	1.87	0.55
1:A:250:GLU:OE1	3:A:823:HOH:O	2.18	0.55
1:D:229:VAL:HG21	1:D:251:LEU:HD22	1.89	0.55
1:D:168:GLU:OE1	1:D:241:HIS:HE1	1.90	0.55
1:C:12:GLN:HG3	1:C:16:LEU:HD22	1.89	0.55
1:A:87:LYS:HZ3	1:A:250:GLU:CG	2.17	0.55
1:B:241:HIS:HD2	1:B:244:LYS:H	1.55	0.55
1:D:41:ASN:HD21	1:D:43:ALA:HB3	1.71	0.55
1:B:211:ASN:ND2	1:B:212:ILE:H	2.04	0.55
1:C:158:ASN:HB3	1:C:161:LEU:HD22	1.88	0.55
1:B:241:HIS:HD2	1:B:243:ASN:N	2.02	0.54
1:B:135:ARG:HG3	1:C:151:GLU:OE1	2.07	0.54
1:C:87:LYS:HZ3	1:C:250:GLU:CG	2.20	0.54
1:B:87:LYS:HZ3	1:B:250:GLU:HG2	1.60	0.54
1:B:135:ARG:CD	1:B:136:LEU:N	2.47	0.54
1:C:29[A]:ASP:OD1	1:C:47:TYR:CE1	2.61	0.53
1:C:237:ASP:HB2	1:C:250:GLU:HB2	1.90	0.53
1:A:162:LYS:HZ3	1:B:59:ARG:HH11	1.56	0.53
1:D:222:VAL:HG13	1:D:236:ILE:HD13	1.90	0.53
1:A:13:GLU:OE1	1:A:256:GLU:HG3	2.09	0.53
1:B:152:HIS:HE1	1:C:153:ARG:NE	1.91	0.52
1:D:56:SER:OG	1:D:58:TYR:CD1	2.62	0.52
1:C:267:ASN:ND2	3:C:682:HOH:O	2.42	0.52
1:A:155:LEU:HD22	1:D:136:LEU:HD11	1.92	0.52
1:D:282:LYS:HA	3:D:751:HOH:O	2.10	0.52
1:A:162:LYS:HD2	1:B:58:TYR:OH	2.10	0.52
1:A:41:ASN:C	1:A:41:ASN:HD22	2.17	0.51
1:C:252:ASN:HB2	3:C:629:HOH:O	2.09	0.51
1:D:241:HIS:HD2	1:D:244:LYS:H	1.57	0.51
1:D:250:GLU:OE1	3:D:760:HOH:O	2.19	0.51
1:A:109:SER:HA	1:A:162:LYS:NZ	2.26	0.51
1:C:73:HIS:CE1	3:C:729:HOH:O	2.57	0.51
1:C:241:HIS:CD2	1:C:243[A]:ASN:H	2.28	0.51
1:C:241:HIS:CD2	1:C:243[B]:ASN:H	2.28	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:LYS:HE3	3:B:681:HOH:O	2.11	0.51
1:A:176[A]:ASP:HB3	1:A:192:ARG:HD3	1.93	0.51
1:C:97:GLU:HG2	1:C:228:ILE:CG1	2.38	0.50
1:B:216:GLU:OE1	1:B:220:GLU:OE2	2.29	0.50
1:B:97:GLU:HG3	1:B:228:ILE:HG12	1.93	0.50
1:D:103:ASP:HB3	1:D:120:ARG:HH22	1.77	0.50
1:D:258:LYS:HE2	3:D:777:HOH:O	2.10	0.50
1:D:132:SER:HB3	1:D:135:ARG:O	2.12	0.50
1:A:202:ASN:N	1:A:203:VAL:HB	2.27	0.49
1:B:168:GLU:OE1	1:B:241:HIS:HE1	1.94	0.49
1:A:256:GLU:HG3	3:A:772:HOH:O	2.12	0.49
1:C:176[B]:ASP:OD2	2:C:500:ADP:O3'	2.25	0.49
1:B:41:ASN:OD1	1:B:43:ALA:HB3	2.13	0.49
1:C:130:ILE:HG22	1:C:131:GLY:N	2.27	0.49
1:A:53:ARG:N	1:A:54:PRO:CD	2.76	0.49
1:C:148:THR:O	1:C:152:HIS:HD2	1.96	0.48
1:C:87:LYS:NZ	1:C:250:GLU:HG2	2.26	0.48
1:D:4:VAL:HG13	1:D:47:TYR:CZ	2.48	0.48
1:D:198:GLU:HG2	1:D:200:ARG:H	1.77	0.48
1:A:109:SER:HA	1:A:162:LYS:HZ1	1.77	0.48
1:A:77:SER:OG	1:A:79:ASP:OD1	2.29	0.48
1:A:232:GLU:OE1	3:A:679:HOH:O	2.20	0.48
1:D:202:ASN:HB2	1:D:205:LEU:HD23	1.95	0.48
1:C:53:ARG:N	1:C:54:PRO:CD	2.76	0.48
1:A:267:ASN:C	1:A:267:ASN:HD22	2.22	0.48
1:B:158:ASN:O	1:B:161:LEU:HB2	2.14	0.48
1:B:222:VAL:HG13	1:B:236:ILE:HG21	1.96	0.48
1:A:241:HIS:CD2	1:A:243:ASN:H	2.32	0.47
1:D:177:ILE:HG12	1:D:212:ILE:HD12	1.95	0.47
1:A:53:ARG:H	1:A:54:PRO:HD3	1.79	0.47
1:C:229:VAL:HG21	1:C:251:LEU:HD22	1.96	0.47
1:A:11:ARG:CZ	1:A:55:VAL:HG22	2.45	0.47
1:B:158:ASN:HD22	1:B:160:ALA:H	1.62	0.47
1:A:156:MET:N	1:A:157:GLY:HA2	2.29	0.47
1:B:224:LYS:O	1:B:228:ILE:HD12	2.14	0.47
1:A:87:LYS:HZ1	1:A:250:GLU:CG	2.28	0.47
1:A:241:HIS:HD2	1:A:243:ASN:H	1.63	0.47
1:D:241:HIS:CD2	1:D:244:LYS:H	2.32	0.46
1:B:57:MET:HE1	1:B:130:ILE:HG13	1.97	0.46
1:D:4:VAL:HG22	1:D:47:TYR:CG	2.51	0.46
1:D:156:MET:SD	1:D:161:LEU:HD23	2.55	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:130:ILE:HD11	1:D:161:LEU:HD11	1.98	0.46
1:A:53:ARG:N	1:A:54:PRO:HD3	2.31	0.46
1:D:38:LEU:HB3	1:D:66:VAL:HG11	1.97	0.46
1:B:196:PRO:HA	1:B:197:ASN:HA	1.64	0.45
1:B:207:GLY:HA2	1:B:208:THR:HA	1.61	0.45
1:D:186:LEU:HD13	1:D:222:VAL:HG11	1.98	0.45
1:C:87:LYS:NZ	1:C:250:GLU:CG	2.79	0.45
1:C:167:GLN:NE2	2:C:500:ADP:HN61	2.15	0.45
1:B:237:ASP:HB2	1:B:250:GLU:HB2	1.99	0.45
1:D:11:ARG:CZ	1:D:55:VAL:HG22	2.46	0.45
1:C:131:GLY:HA3	1:C:132:SER:HB2	1.98	0.44
1:D:87:LYS:NZ	1:D:250:GLU:CD	2.75	0.44
1:D:137:VAL:O	1:D:199:TRP:HH2	2.00	0.44
1:B:259:GLY:HA2	1:B:262:VAL:HG22	2.00	0.44
1:D:226:VAL:CG1	1:D:236:ILE:HD11	2.47	0.44
1:D:130:ILE:HD12	1:D:131:GLY:H	1.82	0.44
1:D:226:VAL:HG11	1:D:236:ILE:HD11	1.99	0.44
1:C:192[B]:ARG:NH2	1:C:192[B]:ARG:CG	2.51	0.44
1:D:11:ARG:NH1	1:D:256[A]:GLU:OE2	2.51	0.44
1:C:264:THR:O	1:C:265:ASN:HB2	2.16	0.44
1:C:269:ALA:O	1:C:273:VAL:HG12	2.18	0.44
1:A:143:VAL:HG23	3:A:635:HOH:O	2.17	0.44
1:A:171:GLN:NE2	3:A:748:HOH:O	2.50	0.44
1:B:178:ARG:HD3	3:B:777:HOH:O	2.18	0.43
1:D:115:LYS:O	1:D:119:GLN:HG3	2.18	0.43
1:C:53:ARG:H	1:C:54:PRO:HD3	1.83	0.43
1:A:258:LYS:NZ	3:A:772:HOH:O	2.51	0.43
1:B:259:GLY:O	1:B:262:VAL:HG22	2.18	0.43
1:B:125:ILE:HG13	2:B:500:ADP:C6	2.52	0.43
1:D:96:ARG:NH2	3:D:748:HOH:O	2.51	0.43
1:B:53:ARG:N	1:B:54:PRO:CD	2.81	0.43
1:B:267:ASN:ND2	3:B:648:HOH:O	2.48	0.43
1:B:4:VAL:HG13	1:B:47:TYR:CZ	2.54	0.43
1:B:176[A]:ASP:OD2	1:B:201:ALA:HB3	2.19	0.43
1:C:3:VAL:HG22	1:C:49:VAL:HG12	2.01	0.43
1:C:130:ILE:CG2	1:C:131:GLY:N	2.81	0.43
1:C:235:SER:OG	1:C:252:ASN:HB3	2.19	0.43
1:D:198:GLU:CG	1:D:199:TRP:H	2.16	0.43
1:A:281:SER:O	1:A:282:LYS:CB	2.65	0.42
1:A:256:GLU:CG	3:A:772:HOH:O	2.66	0.42
1:B:241:HIS:CE1	1:B:244:LYS:HE2	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:THR:O	1:A:152:HIS:HD2	2.02	0.42
1:A:203:VAL:HG13	1:A:203:VAL:O	2.18	0.42
1:B:135:ARG:HD2	1:C:151:GLU:HB3	2.02	0.42
1:C:120:ARG:HD3	3:C:653:HOH:O	2.20	0.42
1:A:87:LYS:NZ	1:A:250:GLU:CD	2.78	0.42
1:A:132:SER:HB3	1:A:135:ARG:HB3	1.96	0.42
1:A:224:LYS:O	1:A:228:ILE:HG12	2.20	0.42
1:B:87:LYS:HZ1	1:B:250:GLU:CD	2.25	0.42
1:C:131:GLY:HA3	1:C:132:SER:CB	2.48	0.42
1:A:160:ALA:HB2	3:B:651:HOH:O	2.20	0.42
1:A:159:SER:C	1:A:161:LEU:H	2.27	0.42
1:C:135:ARG:NH2	1:C:135:ARG:CG	2.64	0.42
1:B:75:ILE:HA	1:B:76:ASN:HA	1.83	0.41
1:C:27:GLN:HE21	1:C:27:GLN:HB2	1.66	0.41
1:B:176[A]:ASP:CG	1:B:201:ALA:HB3	2.45	0.41
1:B:114:LEU:HD11	1:B:147:LYS:HE2	2.01	0.41
1:D:217:LYS:HE2	1:D:217:LYS:N	2.33	0.41
1:B:281:SER:O	1:B:282:LYS:HB2	2.21	0.41
1:D:148:THR:O	1:D:152:HIS:HD2	2.02	0.41
1:A:208:THR:HA	1:A:209:PRO:HD3	1.80	0.41
1:B:87:LYS:HZ1	1:B:250:GLU:CG	2.29	0.41
1:A:194:ILE:HD11	3:A:756:HOH:O	2.20	0.41
1:D:191:ALA:HB2	1:D:212:ILE:HD11	2.03	0.41
1:B:212:ILE:O	1:B:212:ILE:HG13	2.21	0.41
1:D:114:LEU:O	1:D:118:GLU:HG2	2.19	0.41
1:C:82:ASN:HD22	1:C:82:ASN:HA	1.73	0.41
1:C:267:ASN:C	1:C:267:ASN:HD22	2.28	0.40
1:A:186:LEU:HD13	1:A:222:VAL:HG11	2.03	0.40
1:A:41:ASN:C	1:A:41:ASN:ND2	2.79	0.40
1:A:87:LYS:HG2	1:A:250:GLU:OE2	2.20	0.40
1:C:196:PRO:HA	1:C:197:ASN:HA	1.61	0.40
1:A:125:ILE:HG13	2:A:500:ADP:C6	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	285/282 (101%)	272 (95%)	9 (3%)	4 (1%)	9	1
1	B	288/282 (102%)	280 (97%)	5 (2%)	3 (1%)	12	3
1	C	278/282 (99%)	270 (97%)	7 (2%)	1 (0%)	30	18
1	D	283/282 (100%)	274 (97%)	6 (2%)	3 (1%)	11	3
All	All	1134/1128 (100%)	1096 (97%)	27 (2%)	11 (1%)	12	3

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	SER
1	A	203	VAL
1	A	204	ALA
1	B	132	SER
1	D	132	SER
1	D	172	TYR
1	B	203	VAL
1	A	53	ARG
1	C	53	ARG
1	B	53	ARG
1	D	53	ARG

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	247/242 (102%)	225 (91%)	22 (9%)	9	1
1	B	250/242 (103%)	233 (93%)	17 (7%)	14	3
1	C	245/242 (101%)	224 (91%)	21 (9%)	10	2
1	D	245/242 (101%)	224 (91%)	21 (9%)	10	2
All	All	987/968 (102%)	906 (92%)	81 (8%)	12	2

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	19[A]	LYS
1	A	19[B]	LYS
1	A	22	GLU
1	A	25	LYS
1	A	41	ASN
1	A	82	ASN
1	A	96[A]	ARG
1	A	96[B]	ARG
1	A	130	ILE
1	A	132	SER
1	A	135	ARG
1	A	155	LEU
1	A	158	ASN
1	A	161	LEU
1	A	173	LYS
1	A	211	ASN
1	A	228	ILE
1	A	256	GLU
1	A	262	VAL
1	A	267	ASN
1	A	282	LYS
1	B	4	VAL
1	B	16[A]	LEU
1	B	16[B]	LEU
1	B	19	LYS
1	B	27	GLN
1	B	41	ASN
1	B	82	ASN
1	B	114	LEU
1	B	130	ILE
1	B	135	ARG
1	B	155	LEU
1	B	158	ASN
1	B	161	LEU
1	B	192	ARG
1	B	211	ASN
1	B	217	LYS
1	B	251	LEU
1	C	10	VAL
1	C	12	GLN
1	C	16	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	19	LYS
1	C	27	GLN
1	C	29[A]	ASP
1	C	29[B]	ASP
1	C	41	ASN
1	C	49	VAL
1	C	82	ASN
1	C	135	ARG
1	C	155	LEU
1	C	161	LEU
1	C	162[A]	LYS
1	C	162[B]	LYS
1	C	192[A]	ARG
1	C	192[B]	ARG
1	C	226	VAL
1	C	250	GLU
1	C	273	VAL
1	C	278	GLU
1	D	2	ARG
1	D	4	VAL
1	D	16	LEU
1	D	30	ILE
1	D	56	SER
1	D	82	ASN
1	D	97	GLU
1	D	103	ASP
1	D	120	ARG
1	D	130	ILE
1	D	155	LEU
1	D	192	ARG
1	D	197	ASN
1	D	212	ILE
1	D	217	LYS
1	D	224[A]	LYS
1	D	224[B]	LYS
1	D	251	LEU
1	D	258	LYS
1	D	262	VAL
1	D	282	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	41	ASN
1	A	82	ASN
1	A	119	GLN
1	A	152	HIS
1	A	197	ASN
1	A	211	ASN
1	A	241	HIS
1	A	267	ASN
1	A	270	GLN
1	A	279	ASN
1	B	12	GLN
1	B	24	ASN
1	B	27	GLN
1	B	41	ASN
1	B	82	ASN
1	B	119	GLN
1	B	152	HIS
1	B	158	ASN
1	B	193	ASN
1	B	211	ASN
1	B	230	HIS
1	B	241	HIS
1	B	267	ASN
1	B	270	GLN
1	B	279	ASN
1	C	12	GLN
1	C	27	GLN
1	C	35	GLN
1	C	82	ASN
1	C	119	GLN
1	C	152	HIS
1	C	167	GLN
1	C	197	ASN
1	C	211	ASN
1	C	241	HIS
1	C	267	ASN
1	C	279	ASN
1	D	24	ASN
1	D	27	GLN
1	D	82	ASN
1	D	152	HIS
1	D	193	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	241	HIS
1	D	267	ASN
1	D	270	GLN
1	D	279	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ADP	D	500	-	28,29,29	1.43	5 (17%)	43,45,45	1.76	9 (20%)
2	ADP	C	500	-	28,29,29	1.39	4 (14%)	43,45,45	1.86	10 (23%)
2	ADP	B	500	-	28,29,29	1.39	5 (17%)	43,45,45	1.81	9 (20%)
2	ADP	A	500	-	28,29,29	1.44	5 (17%)	43,45,45	1.83	8 (18%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	D	500	-	-	1/16/32/32	0/3/3/3
2	ADP	C	500	-	-	2/16/32/32	0/3/3/3
2	ADP	B	500	-	-	1/16/32/32	0/3/3/3
2	ADP	A	500	-	-	2/16/32/32	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	ADP	C5-C4	4.70	1.47	1.39
2	D	500	ADP	C5-C4	4.65	1.47	1.39
2	C	500	ADP	C5-C4	4.64	1.47	1.39
2	B	500	ADP	C5-C4	4.15	1.46	1.39
2	A	500	ADP	C5-C6	2.79	1.48	1.41
2	B	500	ADP	C5-C6	2.53	1.48	1.41
2	D	500	ADP	C5-C6	2.48	1.47	1.41
2	D	500	ADP	C8-N7	2.45	1.36	1.31
2	D	500	ADP	C4-N9	-2.40	1.32	1.37
2	B	500	ADP	C8-N7	2.38	1.36	1.31
2	C	500	ADP	C5-C6	2.33	1.47	1.41
2	C	500	ADP	C4-N9	-2.27	1.33	1.37
2	B	500	ADP	C5-N7	-2.25	1.35	1.39
2	A	500	ADP	C4-N9	-2.25	1.33	1.37
2	D	500	ADP	C5-N7	-2.21	1.35	1.39
2	C	500	ADP	C8-N7	2.16	1.35	1.31
2	A	500	ADP	C5-N7	-2.11	1.35	1.39
2	A	500	ADP	C8-N7	2.04	1.35	1.31
2	B	500	ADP	C4-N9	-2.03	1.33	1.37

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	ADP	C5-C4-N3	-5.52	119.11	126.72
2	C	500	ADP	C5-C4-N3	-5.24	119.50	126.72
2	C	500	ADP	N3-C4-N9	4.83	135.38	127.17
2	B	500	ADP	C5-C4-N3	-4.79	120.12	126.72
2	A	500	ADP	N3-C4-N9	4.72	135.20	127.17
2	D	500	ADP	C5-C4-N3	-4.67	120.28	126.72
2	B	500	ADP	N3-C4-N9	4.51	134.84	127.17
2	D	500	ADP	N3-C4-N9	4.40	134.65	127.17
2	B	500	ADP	N3-C2-N1	-4.00	122.52	128.58
2	C	500	ADP	C4-N9-C8	3.92	109.86	105.74
2	D	500	ADP	C4-N9-C8	3.82	109.75	105.74
2	B	500	ADP	C4-N9-C8	3.78	109.70	105.74

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	ADP	N3-C2-N1	-3.63	123.08	128.58
2	C	500	ADP	C2-N3-C4	3.60	120.61	111.83
2	B	500	ADP	C2-N3-C4	3.56	120.52	111.83
2	A	500	ADP	C2-N3-C4	3.53	120.45	111.83
2	A	500	ADP	C4-N9-C8	3.50	109.42	105.74
2	D	500	ADP	N3-C2-N1	-3.16	123.80	128.58
2	A	500	ADP	C4-C5-N7	-3.13	107.00	110.58
2	D	500	ADP	C2-N3-C4	3.00	119.15	111.83
2	A	500	ADP	N3-C2-N1	-2.98	124.07	128.58
2	C	500	ADP	C4-C5-N7	-2.73	107.47	110.58
2	A	500	ADP	C5-N7-C8	2.62	107.57	103.45
2	A	500	ADP	N9-C8-N7	-2.54	110.33	113.94
2	B	500	ADP	N9-C8-N7	-2.34	110.62	113.94
2	D	500	ADP	C4-C5-N7	-2.30	107.95	110.58
2	D	500	ADP	C2-N1-C6	2.26	122.45	118.73
2	C	500	ADP	N9-C8-N7	-2.23	110.77	113.94
2	D	500	ADP	C3'-C2'-C1'	2.20	105.62	101.46
2	B	500	ADP	C4-C5-N7	-2.17	108.10	110.58
2	C	500	ADP	C5-N7-C8	2.12	106.78	103.45
2	C	500	ADP	O3'-C3'-C4'	-2.07	105.15	111.08
2	B	500	ADP	C2-N1-C6	2.06	122.12	118.73
2	B	500	ADP	C6-C5-N7	2.03	136.01	132.09
2	D	500	ADP	N9-C8-N7	-2.03	111.05	113.94
2	C	500	ADP	O2'-C2'-C1'	-2.02	103.15	110.10

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	ADP	PB-O3A-PA-O2A
2	B	500	ADP	PB-O3A-PA-O2A
2	C	500	ADP	PB-O3A-PA-O2A
2	D	500	ADP	PB-O3A-PA-O2A
2	A	500	ADP	PB-O3A-PA-O1A
2	C	500	ADP	PB-O3A-PA-O1A

There are no ring outliers.

3 monomers are involved in 6 short contacts:

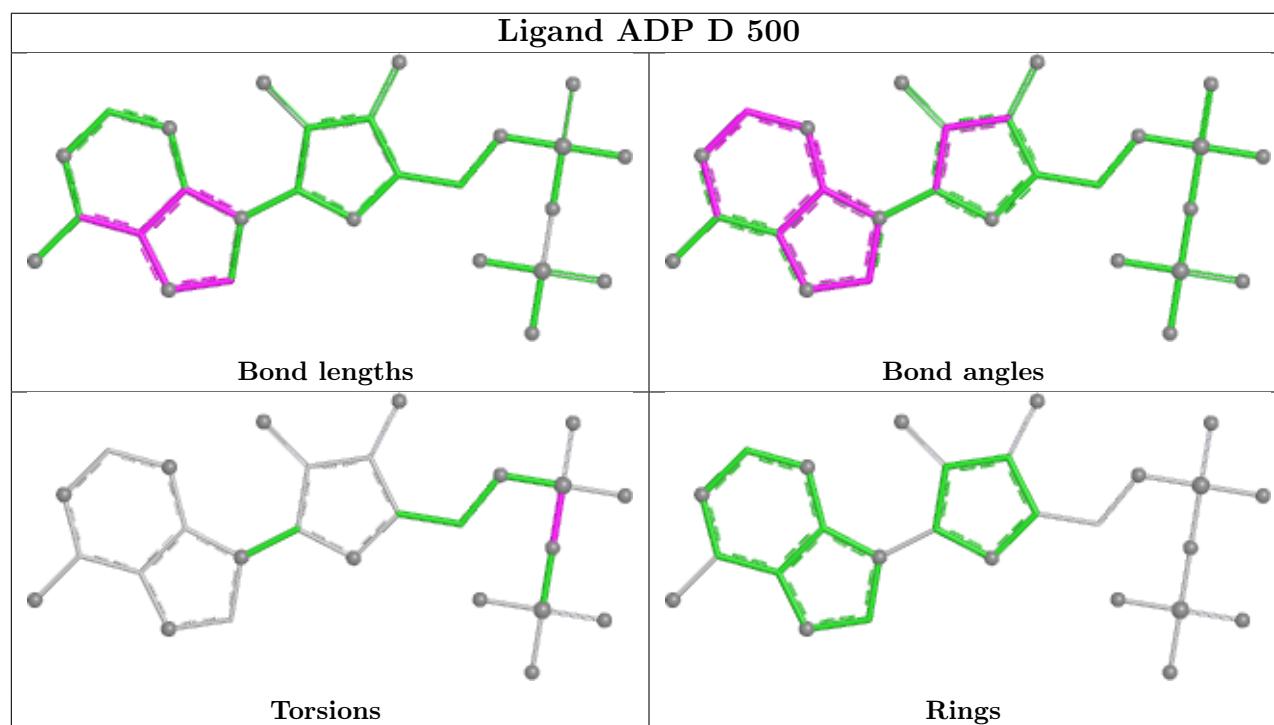
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	500	ADP	2	0

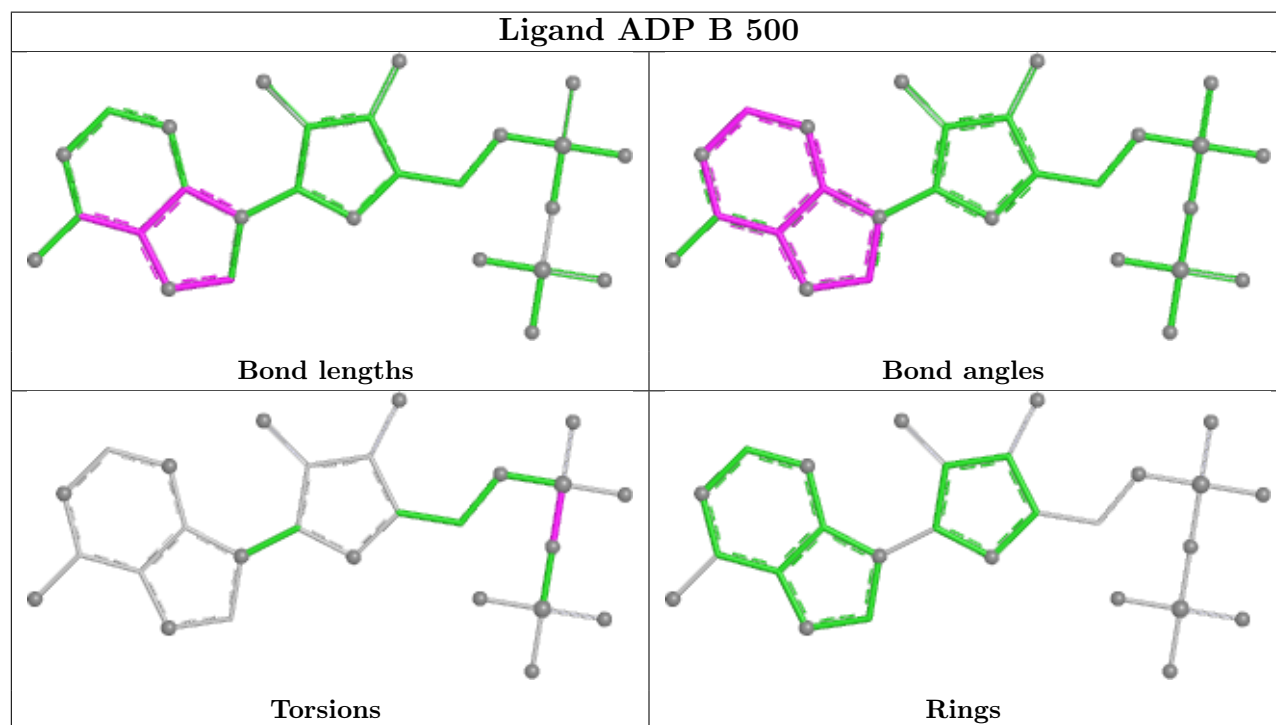
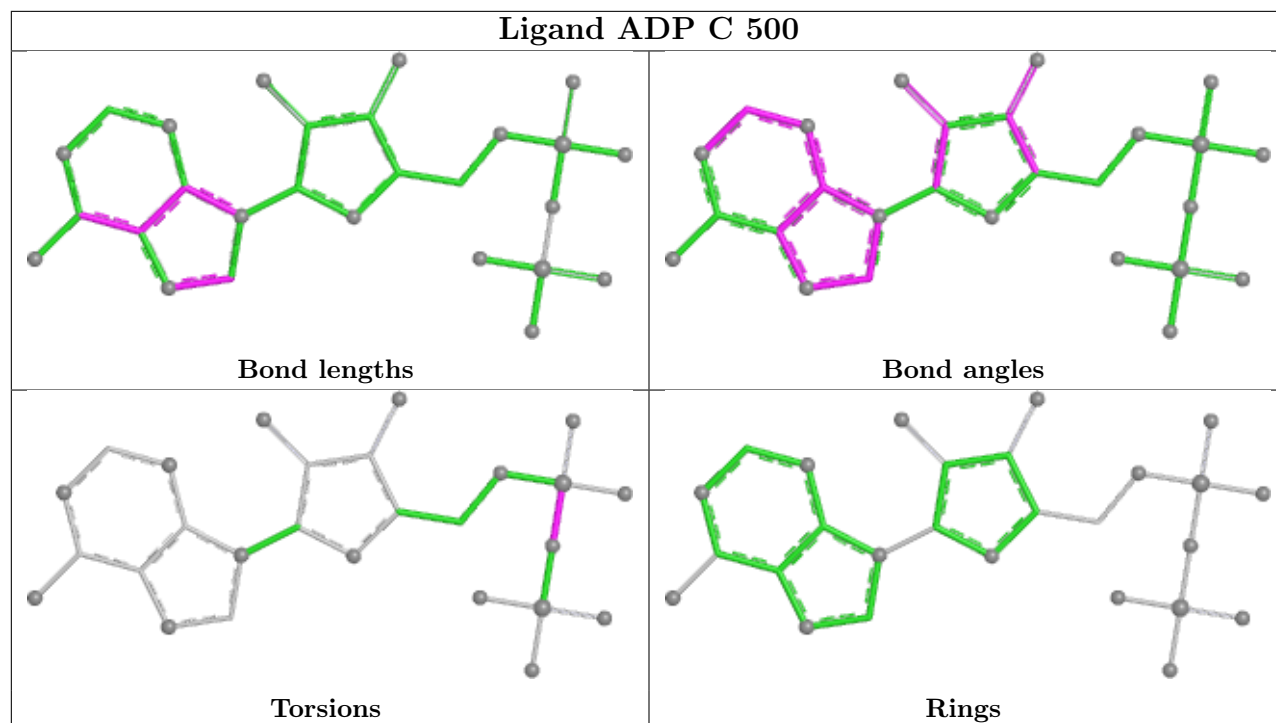
Continued on next page...

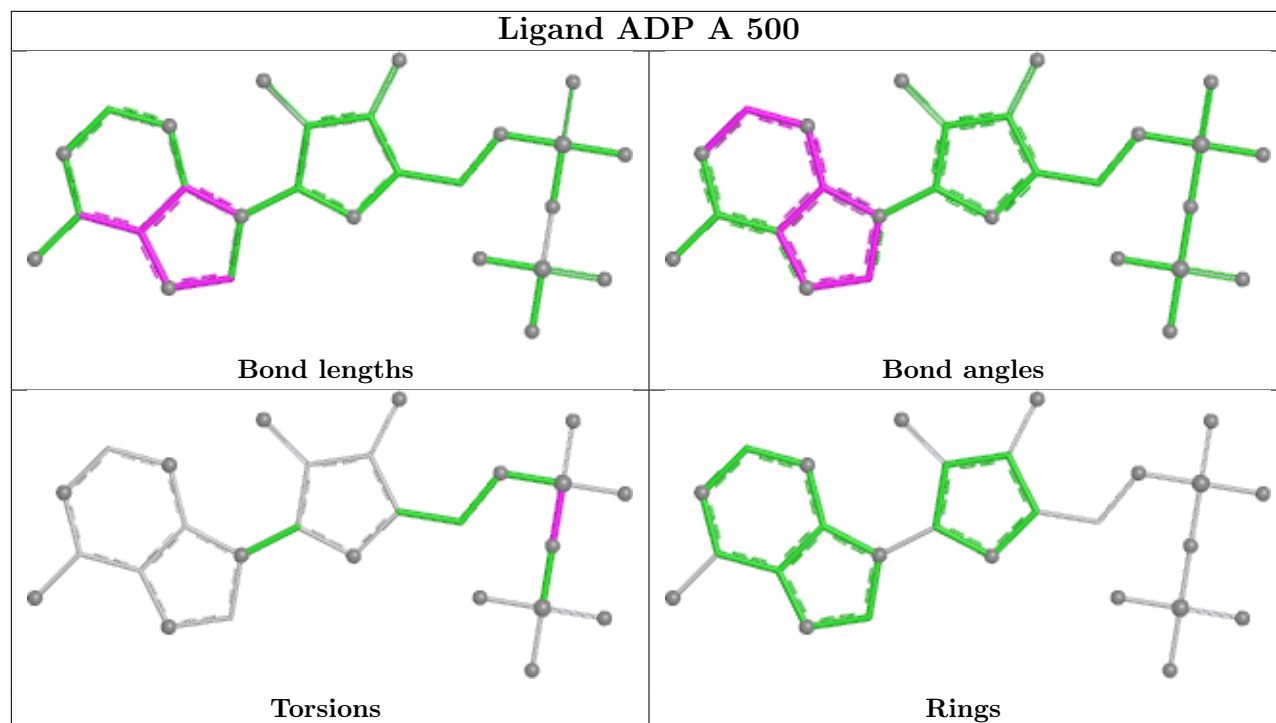
Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	ADP	2	0
2	A	500	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 than 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.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	282/282 (100%)	0.12	22 (7%)	19 21	10, 22, 50, 63	5 (1%)
1	B	282/282 (100%)	0.07	11 (3%)	43 47	9, 22, 43, 60	8 (2%)
1	C	275/282 (97%)	0.22	13 (4%)	36 38	12, 25, 46, 55	7 (2%)
1	D	282/282 (100%)	0.21	16 (5%)	29 31	13, 25, 46, 58	3 (1%)
All	All	1121/1128 (99%)	0.16	62 (5%)	30 33	9, 24, 47, 63	23 (2%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	207	GLY	7.0
1	A	205	LEU	5.6
1	C	133	TRP	5.4
1	B	208	THR	4.9
1	C	131	GLY	4.7
1	D	204	ALA	4.5
1	A	199	TRP	4.4
1	A	204	ALA	4.2
1	D	203	VAL	4.0
1	B	133	TRP	3.9
1	D	133	TRP	3.9
1	B	203	VAL	3.9
1	A	206	GLY	3.7
1	B	204	ALA	3.6
1	D	205	LEU	3.6
1	A	201	ALA	3.5
1	C	209	PRO	3.4
1	D	132	SER	3.4
1	A	133	TRP	3.4
1	B	199	TRP	3.2
1	C	196	PRO	3.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	207	GLY	3.2
1	A	208	THR	3.1
1	A	131	GLY	3.1
1	B	209	PRO	3.0
1	A	209	PRO	3.0
1	A	203	VAL	2.8
1	C	132	SER	2.8
1	A	158	ASN	2.8
1	A	157	GLY	2.8
1	D	212	ILE	2.8
1	D	130	ILE	2.7
1	C	160	ALA	2.7
1	B	205	LEU	2.6
1	D	199	TRP	2.6
1	B	212	ILE	2.6
1	D	58	TYR	2.5
1	D	79	ASP	2.5
1	A	212	ILE	2.5
1	C	201	ALA	2.5
1	A	211	ASN	2.4
1	A	197	ASN	2.4
1	C	134	GLY	2.4
1	D	207	GLY	2.4
1	B	281	SER	2.4
1	B	135	ARG	2.3
1	D	208	THR	2.3
1	D	197	ASN	2.3
1	D	209	PRO	2.3
1	C	199	TRP	2.2
1	D	131	GLY	2.2
1	C	210	SER	2.2
1	A	202	ASN	2.2
1	A	132	SER	2.2
1	A	196	PRO	2.2
1	A	210	SER	2.1
1	A	194	ILE	2.1
1	C	135	ARG	2.1
1	C	192[A]	ARG	2.1
1	D	281	SER	2.1
1	C	197	ASN	2.1
1	A	282	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

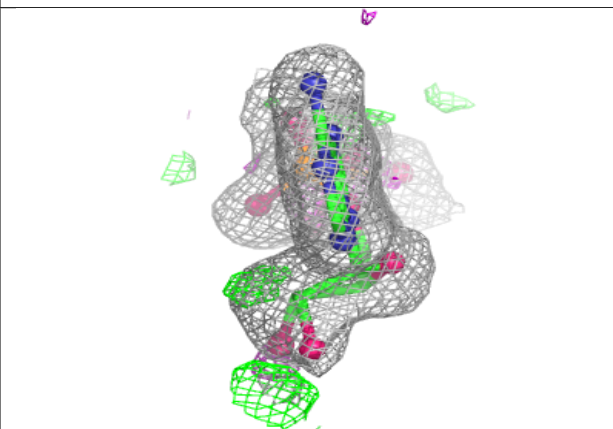
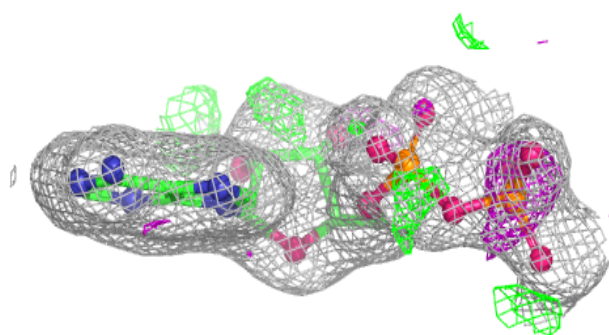
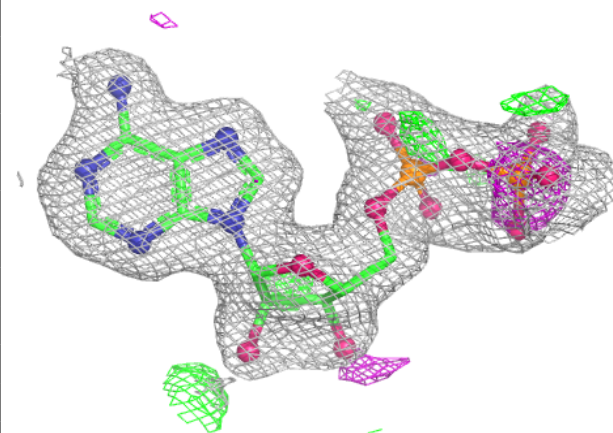
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ADP	C	500	27/27	0.92	0.09	18,25,40,42	0
2	ADP	A	500	27/27	0.93	0.07	17,22,36,38	0
2	ADP	D	500	27/27	0.95	0.07	15,22,38,39	0
2	ADP	B	500	27/27	0.96	0.06	15,20,33,33	0

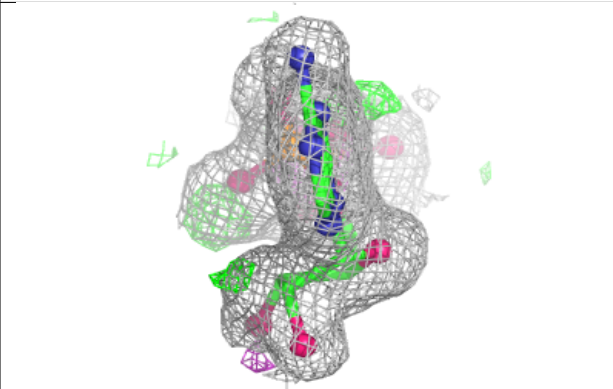
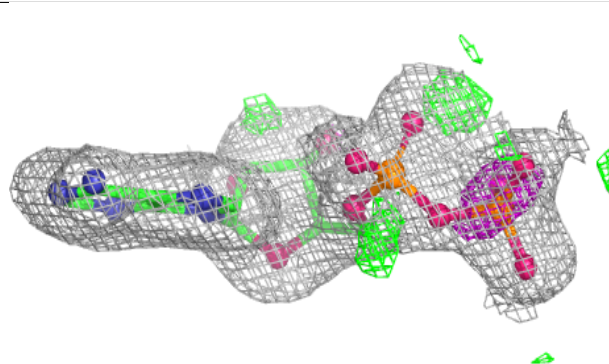
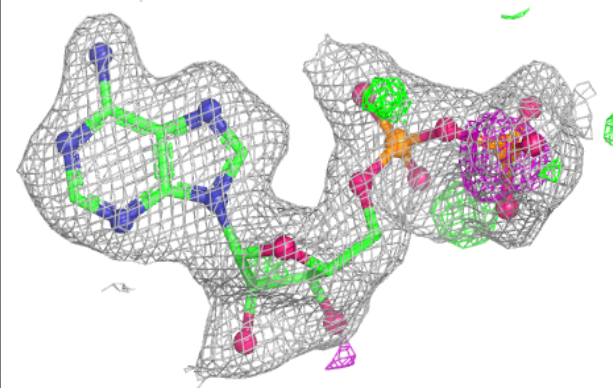
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP C 500:

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

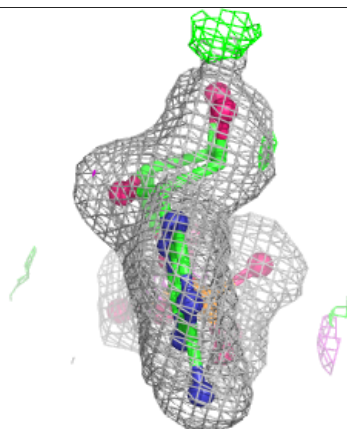
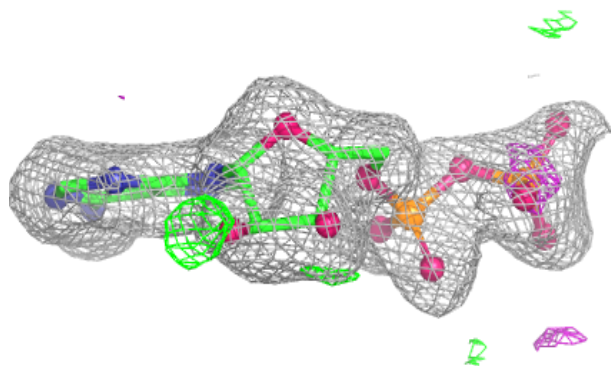
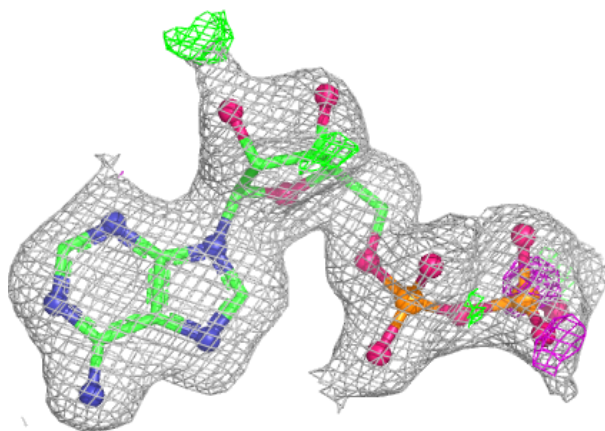
**Electron density around ADP A 500:**

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

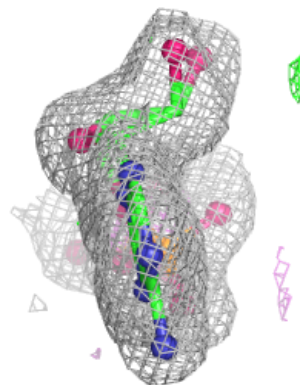
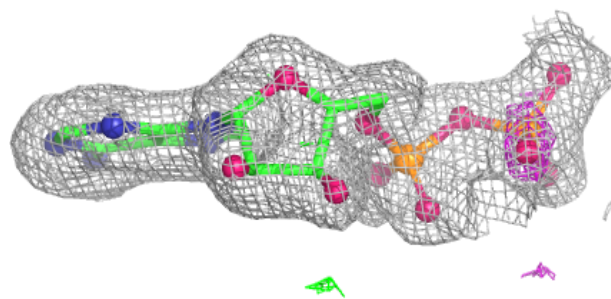
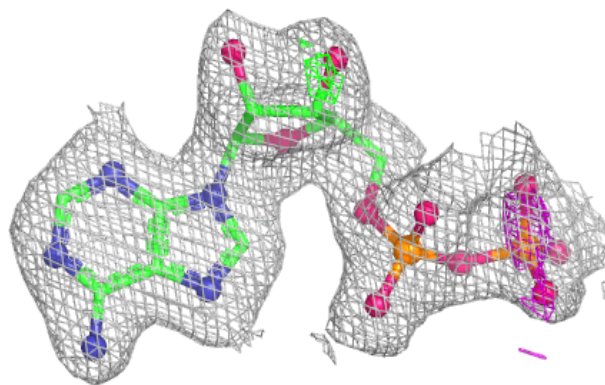


Electron density around ADP D 500:

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

**Electron density around ADP B 500:**

$2mF_o-DF_c$ (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.