



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:17 AM UTC

PDB ID : 2YZN / pdb_00002yzn
Title : Crystal structure of D-alanine:D-Alanine Ligase with AMPPNP from *Thermus thermophilus* HB8.
Authors : Kitamuta, Y.; Yokoyama, S.; Kuramitsu, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-05-06
Resolution : 2.60 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

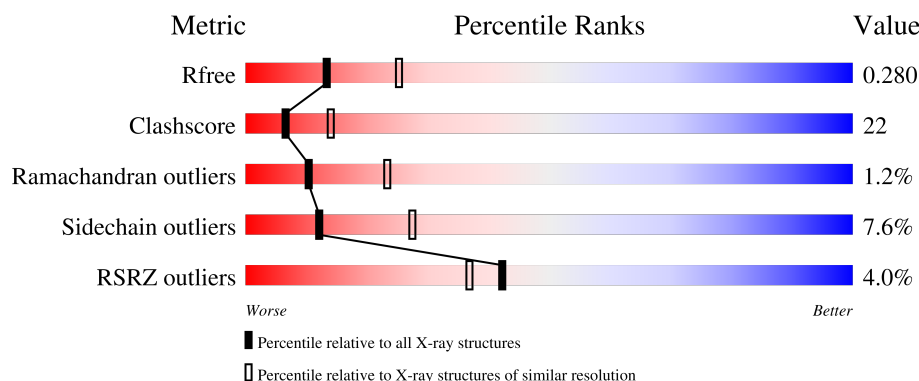
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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.

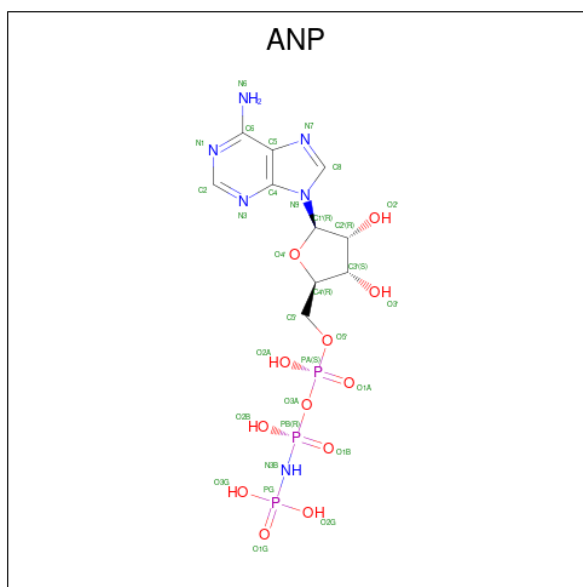
Mol	Chain	Length	Quality of chain
1	A	319	
1	B	319	
1	C	319	

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 D-alanine-D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total 2443	C 1588	N 404	O 446	S 5	0	0	0
1	B	313	Total 2396	C 1556	N 399	O 436	S 5	0	0	0
1	C	309	Total 2359	C 1529	N 395	O 430	S 5	0	0	0

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $\text{C}_{10}\text{H}_{17}\text{N}_6\text{O}_{12}\text{P}_3$).



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

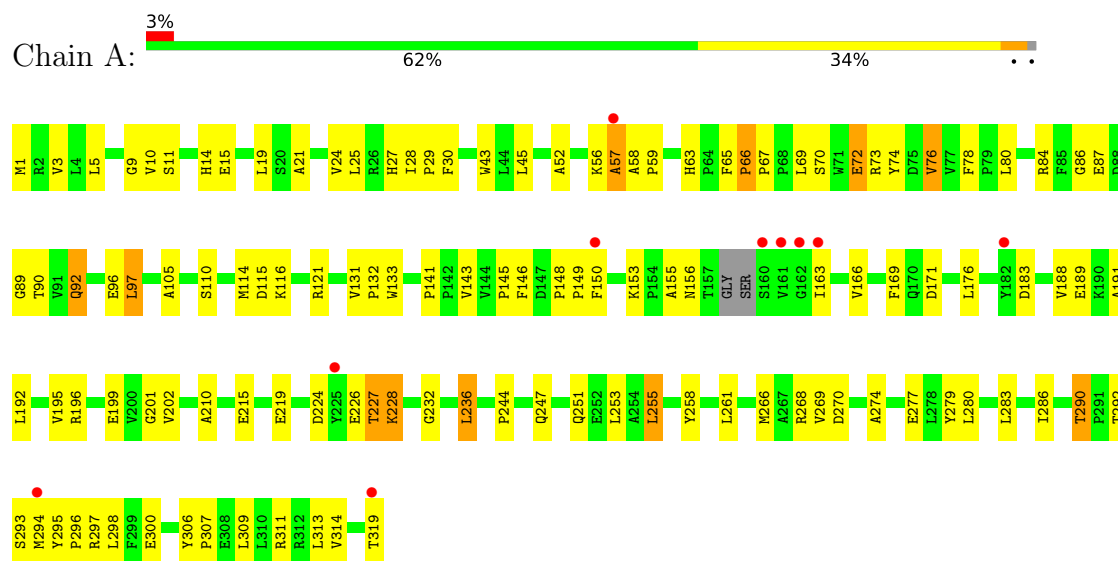
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	24	Total 24	O 24	0	0
3	B	22	Total 22	O 22	0	0
3	C	35	Total 35	O 35	0	0

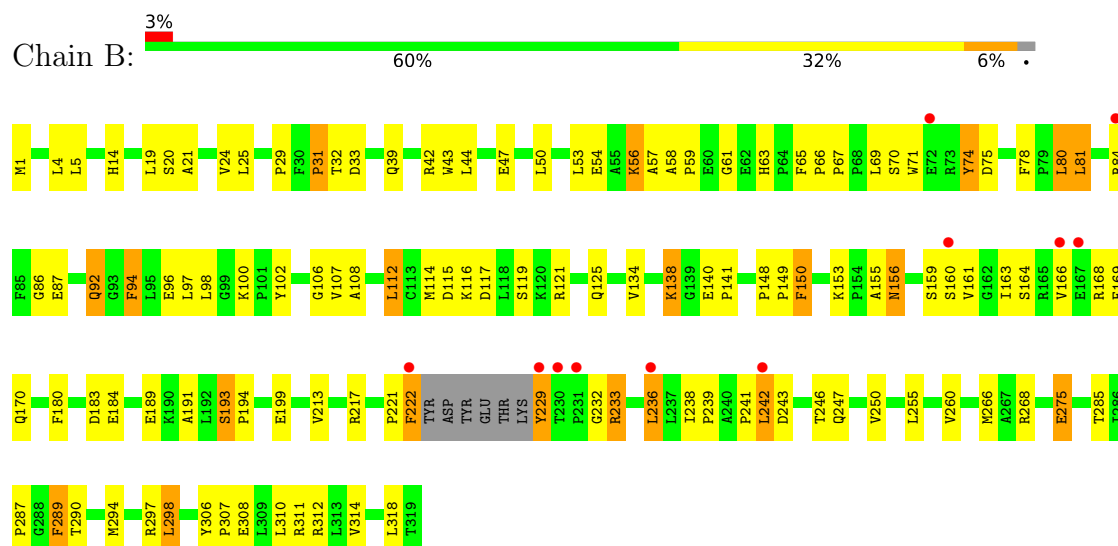
3 Residue-property plots [i](#)

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: D-alanine-D-alanine ligase

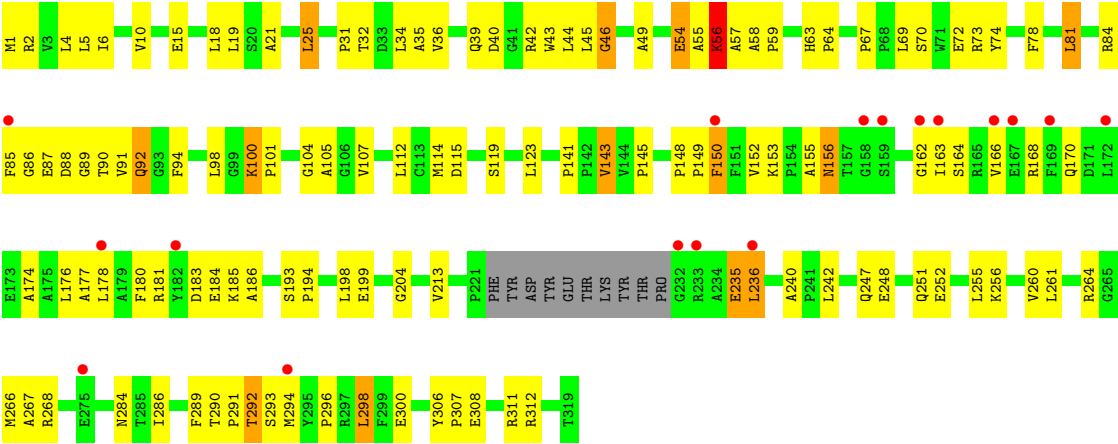


• Molecule 1: D-alanine-D-alanine ligase



• Molecule 1: D-alanine-D-alanine ligase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	167.72Å 56.71Å 141.83Å 90.00° 108.56° 90.00°	Depositor
Resolution (Å)	47.91 – 2.60 47.91 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.4 (47.91-2.60) 96.5 (47.91-2.60)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.88 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.280 0.229 , 0.280	Depositor DCC
R_{free} test set	3856 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	44.6	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7372	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.49	0/2507	1.01	13/3418 (0.4%)
1	B	0.46	0/2458	0.95	9/3351 (0.3%)
1	C	0.51	0/2418	1.03	10/3295 (0.3%)
All	All	0.49	0/7383	1.00	32/10064 (0.3%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	105	ALA	N-CA-C	11.80	125.27	111.11
1	B	66	PRO	CA-C-N	7.99	125.41	119.66
1	B	66	PRO	C-N-CA	7.99	125.41	119.66
1	A	141	PRO	CA-C-N	7.89	127.65	119.76
1	A	141	PRO	C-N-CA	7.89	127.65	119.76
1	A	227	THR	N-CA-C	7.73	119.34	111.07
1	C	300	GLU	N-CA-C	-7.57	102.97	111.07
1	A	300	GLU	N-CA-C	-7.05	103.53	111.14
1	C	268	ARG	N-CA-C	-6.28	98.50	108.73
1	B	290	THR	N-CA-C	-6.17	101.31	110.20
1	C	141	PRO	CA-C-N	6.17	126.18	119.89
1	C	141	PRO	C-N-CA	6.17	126.18	119.89
1	A	232	GLY	N-CA-C	-5.89	105.85	112.33
1	A	67	PRO	N-CA-C	5.83	116.07	110.47
1	A	63	HIS	N-CA-C	5.76	119.68	108.85
1	B	74	TYR	N-CA-C	5.73	118.88	109.72
1	C	32	THR	N-CA-C	5.56	118.16	108.76
1	A	155	ALA	N-CA-C	5.45	117.30	111.36
1	B	31	PRO	N-CA-C	-5.39	102.04	110.50
1	B	289	PHE	N-CA-C	5.34	119.40	111.87
1	A	66	PRO	CA-C-N	5.33	123.55	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	PRO	C-N-CA	5.33	123.55	119.66
1	A	105	ALA	N-CA-C	5.23	117.82	110.23
1	C	58	ALA	CA-C-N	5.21	126.36	119.84
1	C	58	ALA	C-N-CA	5.21	126.36	119.84
1	A	290	THR	N-CA-C	-5.17	102.66	109.84
1	A	74	TYR	N-CA-C	5.16	117.66	109.50
1	C	107	VAL	N-CA-C	5.13	115.56	110.23
1	B	63	HIS	N-CA-C	5.13	117.06	109.62
1	B	94	PHE	N-CA-C	-5.07	105.64	111.07
1	C	104	GLY	N-CA-C	5.07	118.19	111.70
1	B	32	THR	N-CA-C	5.04	117.28	108.76

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	2443	0	2461	101	0
1	B	2396	0	2421	119	0
1	C	2359	0	2389	103	0
2	A	31	0	13	1	0
2	B	31	0	13	0	0
2	C	31	0	13	0	0
3	A	24	0	0	1	0
3	B	22	0	0	0	0
3	C	35	0	0	2	0
All	All	7372	0	7310	317	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 22.

All (317) 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:C:85:PHE:HA	1:C:90:THR:HG21	1.34	1.08
1:C:89:GLY:HA2	1:C:92:GLN:HG3	1.42	1.01
1:C:290:THR:HG22	1:C:292:THR:H	1.29	0.97
1:A:89:GLY:HA2	1:A:92:GLN:HE21	1.34	0.93
1:C:55:ALA:O	1:C:56:LYS:HD3	1.68	0.92
1:A:70:SER:HB2	1:A:73:ARG:HD3	1.55	0.87
1:A:195:VAL:HG12	1:A:274:ALA:HB2	1.58	0.85
1:B:236:LEU:HD22	1:B:297:ARG:HE	1.42	0.85
1:B:92:GLN:H	1:B:92:GLN:NE2	1.80	0.79
1:B:153:LYS:HB3	1:B:163:ILE:HG12	1.64	0.79
1:C:85:PHE:HA	1:C:90:THR:CG2	2.09	0.79
1:A:224:ASP:HB3	1:A:227:THR:HB	1.63	0.78
1:A:19:LEU:HG	1:A:56:LYS:HZ1	1.46	0.78
1:B:5:LEU:HD12	1:B:78:PHE:HB3	1.65	0.77
1:B:236:LEU:HD22	1:B:297:ARG:NE	2.00	0.77
1:A:89:GLY:HA2	1:A:92:GLN:NE2	1.99	0.76
1:C:115:ASP:HA	1:C:156:ASN:HB3	1.67	0.75
1:B:115:ASP:HA	1:B:156:ASN:HB3	1.67	0.75
1:A:3:VAL:HG22	1:A:76:VAL:HG13	1.69	0.75
1:B:87:GLU:HG3	1:B:114:MET:CE	2.18	0.74
1:B:138:LYS:HD3	1:B:138:LYS:N	2.03	0.74
1:C:168:ARG:NH2	1:C:168:ARG:HB2	2.03	0.74
1:C:236:LEU:HG	1:C:294:MET:HE2	1.69	0.72
1:B:114:MET:HE2	1:B:114:MET:HA	1.71	0.72
1:A:290:THR:HG22	1:A:292:THR:H	1.54	0.71
1:B:14:HIS:CE1	1:B:58:ALA:H	2.10	0.70
1:B:199:GLU:HG3	1:B:294:MET:SD	2.31	0.70
1:A:24:VAL:HG21	1:A:80:LEU:HD11	1.74	0.69
1:A:319:THR:HG22	1:A:319:THR:O	1.92	0.69
1:A:224:ASP:CB	1:A:227:THR:HB	2.22	0.69
1:A:306:TYR:HB3	1:A:307:PRO:HD3	1.73	0.69
1:B:100:LYS:HA	1:B:100:LYS:HE2	1.73	0.69
1:B:306:TYR:HB3	1:B:307:PRO:HD3	1.75	0.68
1:A:236:LEU:HD22	1:A:297:ARG:HH21	1.58	0.68
1:B:42:ARG:HG3	1:B:42:ARG:HH11	1.58	0.68
1:C:89:GLY:CA	1:C:92:GLN:HG3	2.23	0.67
1:B:87:GLU:HG3	1:B:114:MET:HE3	1.74	0.67
1:A:19:LEU:HG	1:A:56:LYS:NZ	2.09	0.67
1:A:45:LEU:HD22	1:A:69:LEU:HD21	1.75	0.66
1:B:153:LYS:CB	1:B:163:ILE:HG12	2.26	0.66
1:A:153:LYS:HG2	1:A:163:ILE:HG23	1.78	0.66
1:A:14:HIS:HE1	1:A:58:ALA:H	1.42	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:PHE:CE1	1:B:232:GLY:HA2	2.32	0.65
1:B:308:GLU:O	1:B:312:ARG:HG3	1.96	0.65
1:C:153:LYS:HB3	1:C:163:ILE:HG12	1.77	0.65
1:A:116:LYS:HE2	1:A:153:LYS:HD2	1.77	0.65
1:B:24:VAL:HG13	1:B:310:LEU:HD22	1.79	0.65
1:A:76:VAL:HG11	1:A:314:VAL:HG13	1.78	0.64
1:C:74:TYR:O	1:C:100:LYS:HE3	1.99	0.63
1:A:227:THR:O	1:A:228:LYS:HB2	1.98	0.63
1:C:90:THR:HG23	1:C:91:VAL:N	2.14	0.62
1:B:163:ILE:HG22	1:B:164:SER:N	2.15	0.62
1:A:9:GLY:O	1:A:14:HIS:HD2	1.82	0.62
1:B:94:PHE:CZ	1:B:98:LEU:HD11	2.35	0.61
1:B:238:ILE:HA	1:B:239:PRO:C	2.24	0.61
1:A:153:LYS:HB3	1:A:163:ILE:HG12	1.81	0.61
1:B:161:VAL:O	1:B:161:VAL:HG13	2.01	0.61
1:B:69:LEU:HD23	1:B:70:SER:N	2.15	0.61
1:B:236:LEU:HD22	1:B:297:ARG:NH2	2.15	0.61
1:C:290:THR:HG21	1:C:292:THR:HG22	1.83	0.61
1:C:6:ILE:HG22	1:C:81:LEU:HD21	1.81	0.61
1:A:150:PHE:HB2	1:A:189:GLU:O	2.01	0.60
1:B:1:MET:HA	1:B:75:ASP:OD2	2.02	0.60
1:B:222:PHE:N	1:B:222:PHE:HD1	2.00	0.60
1:B:238:ILE:HD12	1:B:298:LEU:HD12	1.83	0.60
1:A:70:SER:CB	1:A:73:ARG:HD3	2.29	0.60
1:A:149:PRO:HG2	1:A:191:ALA:HB3	1.83	0.60
1:A:89:GLY:CA	1:A:92:GLN:HE21	2.11	0.60
1:A:192:LEU:HB3	1:A:195:VAL:HG11	1.84	0.59
1:B:236:LEU:HD22	1:B:297:ARG:HH21	1.65	0.59
1:B:14:HIS:HE1	1:B:58:ALA:H	1.50	0.59
1:B:236:LEU:HD12	1:B:236:LEU:H	1.67	0.59
1:B:29:PRO:HG3	1:B:311:ARG:HB2	1.84	0.59
1:C:55:ALA:C	1:C:56:LYS:HD3	2.28	0.59
1:B:42:ARG:HG3	1:B:42:ARG:NH1	2.17	0.59
1:A:294:MET:HB2	3:A:404:HOH:O	2.03	0.59
1:B:71:TRP:HB3	1:B:100:LYS:HD2	1.84	0.59
1:A:294:MET:HG3	1:A:298:LEU:HD13	1.85	0.58
1:B:92:GLN:H	1:B:92:GLN:HE21	1.51	0.58
1:C:2:ARG:HD2	1:C:73:ARG:O	2.03	0.58
1:B:138:LYS:HD2	1:B:184:GLU:HB3	1.84	0.58
1:C:236:LEU:HG	1:C:294:MET:CE	2.33	0.58
1:C:56:LYS:HD3	1:C:56:LYS:C	2.27	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:ALA:O	1:A:25:LEU:HB2	2.04	0.57
1:B:222:PHE:N	1:B:222:PHE:CD1	2.71	0.57
1:B:236:LEU:HD13	1:B:297:ARG:NH2	2.18	0.57
1:A:266:MET:HB3	1:A:313:LEU:HD21	1.86	0.57
1:B:108:ALA:O	1:B:112:LEU:HD22	2.04	0.57
1:A:27:HIS:HD2	1:A:306:TYR:HE2	1.53	0.57
1:B:138:LYS:HD3	1:B:138:LYS:H	1.69	0.57
1:B:268:ARG:HB2	1:B:287:PRO:HB3	1.86	0.57
1:B:78:PHE:CE1	1:B:80:LEU:HD13	2.39	0.57
1:A:65:PHE:HD1	1:B:98:LEU:HD23	1.70	0.56
1:B:236:LEU:HD22	1:B:297:ARG:CZ	2.34	0.56
1:A:277:GLU:OE2	1:A:277:GLU:HA	2.05	0.56
1:C:256:LYS:O	1:C:260:VAL:HG23	2.06	0.56
1:A:148:PRO:HG3	1:A:150:PHE:CE2	2.41	0.56
1:B:233:ARG:NH2	1:B:233:ARG:HB2	2.20	0.56
1:C:156:ASN:O	1:C:156:ASN:ND2	2.38	0.56
1:C:168:ARG:HH22	1:C:170:GLN:HB2	1.70	0.55
1:C:181:ARG:HG3	1:C:181:ARG:HH11	1.70	0.55
1:B:233:ARG:HB2	1:B:233:ARG:HH21	1.70	0.55
1:A:268:ARG:HH21	1:A:268:ARG:HG3	1.71	0.55
1:A:121:ARG:O	1:B:125:GLN:HG3	2.07	0.55
1:B:193:SER:HA	1:B:194:PRO:C	2.32	0.55
1:A:56:LYS:HZ3	1:A:56:LYS:HB3	1.71	0.54
1:C:1:MET:O	1:C:31:PRO:HD2	2.07	0.54
1:B:4:LEU:HB2	1:B:74:TYR:CD2	2.42	0.54
1:B:50:LEU:HD12	1:B:50:LEU:H	1.71	0.54
1:A:293:SER:O	1:A:297:ARG:HG3	2.07	0.54
1:B:117:ASP:O	1:B:121:ARG:HG3	2.08	0.54
1:C:248:GLU:HA	1:C:251:GLN:HE21	1.72	0.54
1:B:116:LYS:HG3	1:B:155:ALA:O	2.07	0.54
1:C:5:LEU:HD12	1:C:78:PHE:HB3	1.89	0.54
1:B:20:SER:O	1:B:24:VAL:HG23	2.07	0.54
1:C:55:ALA:O	1:C:56:LYS:CD	2.51	0.54
1:C:87:GLU:HB3	1:C:286:ILE:HD12	1.90	0.54
1:A:56:LYS:O	1:A:57:ALA:HB2	2.08	0.53
1:B:236:LEU:HD13	1:B:297:ARG:HH21	1.73	0.53
1:A:14:HIS:CE1	1:A:58:ALA:H	2.23	0.53
1:B:5:LEU:CD1	1:B:78:PHE:HB3	2.38	0.53
1:A:87:GLU:O	1:A:110:SER:HB3	2.08	0.53
1:C:155:ALA:HB3	1:C:183:ASP:OD1	2.09	0.53
1:C:89:GLY:HA2	1:C:92:GLN:CG	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:PRO:HB3	1:B:229:TYR:HE1	1.72	0.52
1:C:193:SER:HA	1:C:194:PRO:C	2.33	0.52
1:B:50:LEU:HD12	1:B:50:LEU:N	2.23	0.52
1:C:88:ASP:C	1:C:88:ASP:OD2	2.52	0.52
1:A:89:GLY:CA	1:A:92:GLN:NE2	2.70	0.52
1:A:227:THR:O	1:A:228:LYS:CB	2.58	0.52
1:C:6:ILE:CG2	1:C:81:LEU:HD21	2.39	0.52
1:C:162:GLY:HA2	1:C:178:LEU:HG	1.92	0.52
1:A:143:VAL:O	1:A:145:PRO:HD3	2.10	0.52
1:C:266:MET:SD	1:C:289:PHE:HE2	2.33	0.52
1:B:148:PRO:HB3	1:B:149:PRO:HA	1.92	0.51
1:C:10:VAL:HG23	1:C:39:GLN:O	2.10	0.51
1:A:116:LYS:HD2	2:A:401:ANP:O1G	2.11	0.51
1:C:4:LEU:HB2	1:C:74:TYR:CD2	2.46	0.51
1:C:54:GLU:C	1:C:56:LYS:H	2.18	0.51
1:A:114:MET:HG3	1:A:114:MET:O	2.09	0.51
1:A:268:ARG:NE	1:A:270:ASP:OD2	2.43	0.51
1:C:56:LYS:HZ2	1:C:57:ALA:HB2	1.76	0.51
1:B:47:GLU:CD	1:C:181:ARG:HG2	2.36	0.51
1:A:27:HIS:HD2	1:A:306:TYR:CE2	2.29	0.50
1:B:246:THR:O	1:B:250:VAL:HG23	2.12	0.50
1:A:150:PHE:HA	1:A:191:ALA:HB2	1.93	0.50
1:B:112:LEU:O	1:B:119:SER:HB3	2.10	0.50
1:B:233:ARG:HH21	1:B:233:ARG:CB	2.24	0.50
1:A:10:VAL:HG21	1:A:84:ARG:NH1	2.26	0.50
1:C:148:PRO:HG3	1:C:150:PHE:CE1	2.46	0.50
1:B:50:LEU:O	1:B:54:GLU:HG3	2.12	0.50
1:A:251:GLN:O	1:A:255:LEU:HD22	2.12	0.49
1:C:36:VAL:HG23	1:C:49:ALA:HB1	1.93	0.49
1:A:56:LYS:O	1:A:57:ALA:CB	2.60	0.49
1:B:1:MET:HE1	1:B:318:LEU:HD11	1.94	0.49
1:B:236:LEU:CD2	1:B:297:ARG:HE	2.21	0.49
1:C:88:ASP:OD2	1:C:90:THR:HG22	2.12	0.49
1:C:39:GLN:HE21	1:C:59:PRO:HA	1.77	0.49
1:C:290:THR:CG2	1:C:292:THR:HG22	2.42	0.49
1:A:148:PRO:HB3	1:A:149:PRO:HA	1.95	0.49
1:C:44:LEU:O	1:C:49:ALA:HB2	2.13	0.49
1:A:73:ARG:HG2	1:A:73:ARG:HH11	1.78	0.49
1:A:290:THR:O	1:A:296:PRO:HG3	2.12	0.49
1:B:87:GLU:HG3	1:B:114:MET:HE1	1.93	0.49
1:B:92:GLN:O	1:B:96:GLU:HG2	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ALA:HB2	1:A:255:LEU:HD21	1.94	0.48
1:C:56:LYS:NZ	1:C:57:ALA:HB2	2.28	0.48
1:C:149:PRO:HA	1:C:166:VAL:O	2.14	0.48
1:B:21:ALA:O	1:B:25:LEU:HB2	2.14	0.48
1:B:33:ASP:HB2	1:B:74:TYR:CE2	2.49	0.48
1:C:168:ARG:HB2	1:C:168:ARG:HH21	1.78	0.48
1:B:114:MET:HE2	1:B:114:MET:CA	2.39	0.48
1:A:27:HIS:CD2	1:A:306:TYR:CE2	3.01	0.48
1:B:81:LEU:HD13	1:B:81:LEU:N	2.28	0.48
1:B:138:LYS:H	1:B:138:LYS:CD	2.27	0.48
1:C:43:TRP:CE3	1:C:67:PRO:HG3	2.48	0.48
1:B:150:PHE:CE2	1:B:166:VAL:HB	2.49	0.48
1:C:15:GLU:O	1:C:19:LEU:HG	2.14	0.47
1:C:90:THR:CG2	1:C:91:VAL:N	2.78	0.47
1:C:312:ARG:HD2	3:C:419:HOH:O	2.13	0.47
1:A:290:THR:HB	1:A:293:SER:OG	2.14	0.47
1:B:138:LYS:N	1:B:138:LYS:CD	2.75	0.47
1:C:213:VAL:CG2	1:C:247:GLN:HG3	2.45	0.47
1:C:290:THR:O	1:C:296:PRO:HG3	2.14	0.47
1:A:224:ASP:CG	1:A:227:THR:HB	2.39	0.47
1:C:290:THR:HG23	1:C:291:PRO:HD2	1.95	0.47
1:A:115:ASP:OD2	1:A:115:ASP:C	2.57	0.47
1:A:311:ARG:HG2	1:A:311:ARG:HH11	1.80	0.47
1:C:90:THR:HG23	1:C:91:VAL:H	1.77	0.47
1:B:150:PHE:HA	1:B:191:ALA:HB2	1.96	0.47
1:B:150:PHE:HB2	1:B:189:GLU:O	2.15	0.47
1:C:100:LYS:HE2	3:C:424:HOH:O	2.14	0.47
1:C:294:MET:HG3	1:C:298:LEU:HD22	1.96	0.47
1:C:168:ARG:HB2	1:C:168:ARG:CZ	2.45	0.47
1:A:56:LYS:HE2	1:A:56:LYS:HA	1.95	0.46
1:A:269:VAL:HG22	1:A:283:LEU:HG	1.97	0.46
1:B:163:ILE:CG2	1:B:164:SER:N	2.78	0.46
1:A:280:LEU:HD23	1:A:280:LEU:C	2.40	0.46
1:B:236:LEU:HD12	1:B:236:LEU:N	2.28	0.46
1:C:152:VAL:CG1	1:C:176:LEU:HD23	2.45	0.46
1:C:199:GLU:OE1	1:C:294:MET:HG2	2.15	0.46
1:A:1:MET:HG2	1:A:30:PHE:CD2	2.51	0.46
1:A:290:THR:C	1:A:292:THR:H	2.24	0.46
1:C:290:THR:HB	1:C:293:SER:OG	2.15	0.46
1:C:70:SER:HB3	1:C:72:GLU:OE2	2.15	0.46
1:C:84:ARG:NH1	1:C:84:ARG:HB3	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:240:ALA:O	1:C:242:LEU:N	2.46	0.45
1:B:115:ASP:OD2	1:B:115:ASP:C	2.59	0.45
1:A:131:VAL:O	1:A:132:PRO:C	2.60	0.45
1:A:28:ILE:HD12	1:A:29:PRO:HD2	1.97	0.45
1:B:160:SER:HA	1:B:163:ILE:HD12	1.98	0.45
1:C:4:LEU:HD21	1:C:69:LEU:HD21	1.99	0.45
1:C:213:VAL:HG23	1:C:247:GLN:HG3	1.99	0.45
1:C:155:ALA:HB1	1:C:185:LYS:HE3	1.98	0.45
1:C:306:TYR:HB3	1:C:307:PRO:HD3	1.97	0.45
1:A:72:GLU:CD	1:A:72:GLU:H	2.24	0.45
1:A:56:LYS:NZ	1:A:56:LYS:HB3	2.31	0.45
1:A:311:ARG:HG2	1:A:311:ARG:NH1	2.32	0.45
1:B:213:VAL:HB	1:B:247:GLN:HG3	1.99	0.45
1:C:40:ASP:OD2	1:C:40:ASP:C	2.59	0.45
1:C:204:GLY:O	1:C:264:ARG:O	2.35	0.45
1:B:33:ASP:HB2	1:B:74:TYR:HE2	1.81	0.45
1:C:84:ARG:HH11	1:C:84:ARG:CB	2.29	0.45
1:C:91:VAL:O	1:C:94:PHE:HB3	2.17	0.45
1:A:148:PRO:HD3	1:A:169:PHE:CE1	2.51	0.45
1:A:268:ARG:HG3	1:A:268:ARG:NH2	2.32	0.45
1:A:156:ASN:HB2	1:A:183:ASP:OD2	2.17	0.44
1:B:24:VAL:HG21	1:B:80:LEU:HD21	1.98	0.44
1:C:148:PRO:HB3	1:C:149:PRO:HA	1.98	0.44
1:C:5:LEU:CD2	1:C:25:LEU:HD13	2.47	0.44
1:C:42:ARG:HG2	1:C:64:PRO:HA	1.99	0.44
1:C:155:ALA:H	1:C:186:ALA:HA	1.81	0.44
1:A:192:LEU:HB3	1:A:195:VAL:CG1	2.47	0.44
1:B:53:LEU:HD13	1:B:53:LEU:O	2.18	0.44
1:B:156:ASN:ND2	1:B:183:ASP:OD1	2.51	0.44
1:A:290:THR:C	1:A:292:THR:N	2.76	0.44
1:B:148:PRO:HD3	1:B:169:PHE:CD2	2.53	0.44
1:C:156:ASN:C	1:C:156:ASN:HD22	2.26	0.44
1:B:294:MET:O	1:B:298:LEU:HD22	2.17	0.44
1:C:311:ARG:HD3	1:C:311:ARG:C	2.42	0.44
1:A:166:VAL:HG13	1:A:171:ASP:HB2	1.98	0.43
1:B:242:LEU:N	1:B:242:LEU:HD22	2.33	0.43
1:A:156:ASN:N	1:A:183:ASP:OD2	2.51	0.43
1:A:196:ARG:HH21	1:A:215:GLU:CD	2.26	0.43
1:B:4:LEU:HD21	1:B:69:LEU:HD11	1.99	0.43
1:C:84:ARG:HG2	1:C:85:PHE:CD2	2.53	0.43
1:A:52:ALA:HB1	1:A:58:ALA:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:PHE:CD2	1:B:166:VAL:HB	2.53	0.43
1:A:90:THR:HG22	1:B:107:VAL:HG21	2.01	0.43
1:A:286:ILE:O	1:A:286:ILE:HG22	2.18	0.43
1:C:90:THR:CG2	1:C:91:VAL:H	2.32	0.43
1:C:94:PHE:CZ	1:C:98:LEU:HD11	2.54	0.43
1:A:192:LEU:HD21	1:A:279:TYR:CD2	2.53	0.43
1:C:21:ALA:O	1:C:25:LEU:HB2	2.19	0.43
1:C:174:ALA:O	1:C:177:ALA:HB3	2.18	0.43
1:C:180:PHE:HA	1:C:183:ASP:O	2.18	0.43
1:C:198:LEU:HD13	1:C:242:LEU:HD11	2.01	0.43
1:B:87:GLU:HB2	1:B:285:THR:OG1	2.18	0.43
1:B:310:LEU:O	1:B:314:VAL:HG23	2.18	0.43
1:B:1:MET:O	1:B:31:PRO:HD2	2.19	0.42
1:C:143:VAL:O	1:C:145:PRO:HD3	2.18	0.42
1:A:133:TRP:HA	1:A:188:VAL:O	2.18	0.42
1:A:236:LEU:HD12	1:A:236:LEU:N	2.34	0.42
1:B:275:GLU:OE1	1:B:275:GLU:HA	2.19	0.42
1:C:63:HIS:HA	1:C:64:PRO:HD3	1.80	0.42
1:A:210:ALA:HB2	1:A:255:LEU:CD2	2.50	0.42
1:B:297:ARG:HG2	1:B:297:ARG:HH11	1.85	0.42
1:C:5:LEU:HD23	1:C:25:LEU:HD13	2.00	0.42
1:A:92:GLN:O	1:A:96:GLU:HG2	2.19	0.42
1:A:97:LEU:HB3	1:B:65:PHE:CE1	2.55	0.42
1:B:24:VAL:HG13	1:B:310:LEU:CD2	2.48	0.42
1:B:148:PRO:CB	1:B:149:PRO:HA	2.50	0.42
1:B:56:LYS:O	1:B:57:ALA:HB2	2.20	0.42
1:B:78:PHE:C	1:B:78:PHE:CD1	2.97	0.42
1:C:235:GLU:C	1:C:236:LEU:HD12	2.44	0.42
1:B:14:HIS:HE1	1:B:57:ALA:HA	1.85	0.42
1:B:69:LEU:HD22	1:B:71:TRP:CD2	2.55	0.42
1:B:102:TYR:OH	1:B:106:GLY:HA2	2.20	0.42
1:B:266:MET:SD	1:B:289:PHE:HE2	2.43	0.42
1:B:44:LEU:HD21	1:B:61:GLY:HA2	2.01	0.42
1:C:72:GLU:H	1:C:72:GLU:CD	2.28	0.42
1:A:145:PRO:HG2	1:A:146:PHE:HD1	1.85	0.42
1:A:146:PHE:CD1	1:A:146:PHE:N	2.88	0.42
1:B:241:PRO:C	1:B:242:LEU:HD22	2.45	0.42
1:A:5:LEU:HD12	1:A:78:PHE:HB3	2.02	0.41
1:A:202:VAL:HG11	1:A:258:TYR:CG	2.55	0.41
1:C:119:SER:O	1:C:123:LEU:HG	2.20	0.41
1:C:45:LEU:HD22	1:C:69:LEU:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:ILE:HG22	1:C:164:SER:N	2.35	0.41
1:B:148:PRO:HD3	1:B:169:PHE:HD2	1.84	0.41
1:C:248:GLU:O	1:C:252:GLU:HG3	2.20	0.41
1:C:100:LYS:HA	1:C:101:PRO:HD3	1.96	0.41
1:B:168:ARG:HB3	1:B:170:GLN:HG2	2.03	0.41
1:B:78:PHE:C	1:B:78:PHE:HD1	2.29	0.41
1:C:267:ALA:HA	1:C:284:ASN:O	2.21	0.41
1:C:45:LEU:O	1:C:46:GLY:C	2.64	0.41
1:C:290:THR:C	1:C:292:THR:N	2.77	0.41
1:B:138:LYS:HA	1:B:180:PHE:CZ	2.56	0.41
1:B:306:TYR:N	1:B:307:PRO:CD	2.84	0.41
1:A:43:TRP:CD1	1:A:43:TRP:N	2.87	0.41
1:A:45:LEU:HD13	1:A:69:LEU:HD23	2.03	0.41
1:A:295:TYR:CE1	1:A:309:LEU:HD21	2.56	0.41
1:B:39:GLN:NE2	1:B:59:PRO:HA	2.36	0.41
1:B:140:GLU:HA	1:B:141:PRO:HD3	1.91	0.41
1:B:222:PHE:HE1	1:B:232:GLY:HA2	1.84	0.41
1:C:34:LEU:HD12	1:C:35:ALA:H	1.85	0.41
1:C:56:LYS:C	1:C:56:LYS:CD	2.92	0.40
1:A:9:GLY:O	1:A:14:HIS:CD2	2.70	0.40
1:A:150:PHE:CB	1:A:189:GLU:O	2.67	0.40
1:A:201:GLY:N	1:A:298:LEU:HD23	2.36	0.40
1:A:57:ALA:O	1:A:59:PRO:HD3	2.21	0.40
1:A:66:PRO:HD3	1:B:98:LEU:HD21	2.03	0.40
1:B:43:TRP:CE3	1:B:67:PRO:HG3	2.57	0.40
1:C:168:ARG:NH2	1:C:170:GLN:CG	2.85	0.40
1:B:53:LEU:HD13	1:B:53:LEU:C	2.47	0.40
1:B:308:GLU:OE2	1:B:308:GLU:HA	2.22	0.40
1:C:168:ARG:HH21	1:C:168:ARG:CB	2.34	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	313/319 (98%)	289 (92%)	20 (6%)	4 (1%)	9	21
1	B	309/319 (97%)	287 (93%)	19 (6%)	3 (1%)	12	28
1	C	305/319 (96%)	278 (91%)	23 (8%)	4 (1%)	9	21
All	All	927/957 (97%)	854 (92%)	62 (7%)	11 (1%)	10	23

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	57	ALA
1	A	228	LYS
1	C	56	LYS
1	A	86	GLY
1	C	46	GLY
1	B	159	SER
1	C	143	VAL
1	C	86	GLY
1	A	244	PRO
1	B	86	GLY
1	B	193	SER

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	255/256 (100%)	240 (94%)	15 (6%)	18	38
1	B	250/256 (98%)	227 (91%)	23 (9%)	8	19
1	C	246/256 (96%)	227 (92%)	19 (8%)	12	27
All	All	751/768 (98%)	694 (92%)	57 (8%)	12	27

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	15	GLU

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Mol	Chain	Res	Type
1	A	72	GLU
1	A	76	VAL
1	A	92	GLN
1	A	97	LEU
1	A	176	LEU
1	A	199	GLU
1	A	219	GLU
1	A	226	GLU
1	A	236	LEU
1	A	247	GLN
1	A	253	LEU
1	A	255	LEU
1	A	261	LEU
1	B	19	LEU
1	B	56	LYS
1	B	80	LEU
1	B	81	LEU
1	B	84	ARG
1	B	92	GLN
1	B	97	LEU
1	B	112	LEU
1	B	134	VAL
1	B	138	LYS
1	B	150	PHE
1	B	156	ASN
1	B	217	ARG
1	B	222	PHE
1	B	229	TYR
1	B	233	ARG
1	B	236	LEU
1	B	242	LEU
1	B	243	ASP
1	B	255	LEU
1	B	260	VAL
1	B	275	GLU
1	B	298	LEU
1	C	18	LEU
1	C	25	LEU
1	C	54	GLU
1	C	56	LYS
1	C	81	LEU
1	C	92	GLN

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Mol	Chain	Res	Type
1	C	100	LYS
1	C	112	LEU
1	C	114	MET
1	C	150	PHE
1	C	156	ASN
1	C	184	GLU
1	C	235	GLU
1	C	236	LEU
1	C	255	LEU
1	C	261	LEU
1	C	292	THR
1	C	298	LEU
1	C	308	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	HIS
1	A	27	HIS
1	A	92	GLN
1	A	251	GLN
1	B	14	HIS
1	B	39	GLN
1	B	63	HIS
1	B	82	HIS
1	B	92	GLN
1	B	156	ASN
1	B	205	ASN
1	B	251	GLN
1	B	281	ASN
1	C	39	GLN
1	C	82	HIS
1	C	92	GLN
1	C	156	ASN
1	C	205	ASN
1	C	247	GLN
1	C	251	GLN

5.3.3 RNA ⓘ

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](#)

3 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	ANP	C	403	-	33,33,33	1.65	6 (18%)	45,52,52	1.85	8 (17%)
2	ANP	B	402	-	33,33,33	1.59	7 (21%)	45,52,52	1.82	8 (17%)
2	ANP	A	401	-	33,33,33	1.70	7 (21%)	45,52,52	1.85	9 (20%)

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	ANP	C	403	-	-	5/18/38/38	0/3/3/3
2	ANP	B	402	-	-	2/18/38/38	0/3/3/3
2	ANP	A	401	-	-	1/18/38/38	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	ANP	PA-O3A	4.37	1.64	1.59
2	A	401	ANP	PB-O3A	4.14	1.64	1.59
2	C	403	ANP	PA-O3A	3.91	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	ANP	PB-O3A	3.42	1.63	1.59
2	B	402	ANP	PA-O3A	3.36	1.63	1.59
2	C	403	ANP	PB-O3A	3.33	1.63	1.59
2	C	403	ANP	PB-N3B	3.29	1.71	1.63
2	C	403	ANP	PG-N3B	3.16	1.71	1.63
2	A	401	ANP	PG-N3B	3.14	1.71	1.63
2	B	402	ANP	PG-N3B	3.09	1.71	1.63
2	A	401	ANP	PB-N3B	2.96	1.71	1.63
2	B	402	ANP	PB-N3B	2.88	1.70	1.63
2	C	403	ANP	C5-C4	2.48	1.43	1.39
2	B	402	ANP	O4'-C1'	2.45	1.47	1.42
2	B	402	ANP	C5-C4	2.21	1.43	1.39
2	A	401	ANP	O4'-C1'	2.20	1.47	1.42
2	C	403	ANP	O4'-C1'	2.19	1.47	1.42
2	A	401	ANP	C5-C4	2.03	1.42	1.39
2	A	401	ANP	PG-O1G	2.03	1.49	1.46
2	B	402	ANP	PG-O2G	2.01	1.62	1.56

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	ANP	C5-C4-N3	-5.38	119.31	126.72
2	C	403	ANP	C5-C4-N3	-5.38	119.31	126.72
2	A	401	ANP	C5-C4-N3	-5.23	119.52	126.72
2	A	401	ANP	N3-C2-N1	-4.78	121.35	128.58
2	B	402	ANP	N3-C2-N1	-4.69	121.49	128.58
2	C	403	ANP	N3-C2-N1	-4.59	121.63	128.58
2	C	403	ANP	N3-C4-N9	4.58	134.96	127.17
2	B	402	ANP	N3-C4-N9	4.50	134.83	127.17
2	A	401	ANP	N3-C4-N9	4.44	134.72	127.17
2	B	402	ANP	C2-N3-C4	3.83	121.18	111.83
2	C	403	ANP	C5-N7-C8	3.79	109.40	103.45
2	C	403	ANP	C2-N3-C4	3.77	121.03	111.83
2	A	401	ANP	C2-N3-C4	3.74	120.96	111.83
2	B	402	ANP	C5-N7-C8	3.70	109.26	103.45
2	A	401	ANP	C5-N7-C8	3.68	109.24	103.45
2	C	403	ANP	C4-C5-N7	-2.93	107.23	110.58
2	B	402	ANP	C4-C5-N7	-2.89	107.28	110.58
2	A	401	ANP	C4-C5-N7	-2.71	107.49	110.58
2	A	401	ANP	N9-C8-N7	-2.39	110.54	113.94
2	A	401	ANP	PA-O5'-C5'	-2.37	107.78	121.35
2	B	402	ANP	PA-O5'-C5'	-2.30	108.18	121.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	ANP	N9-C8-N7	-2.24	110.76	113.94
2	C	403	ANP	N9-C8-N7	-2.23	110.78	113.94
2	C	403	ANP	PA-O5'-C5'	-2.20	108.73	121.35
2	A	401	ANP	O1G-PG-N3B	-2.01	108.81	111.77

There are no chirality outliers.

All (8) torsion outliers are listed below:

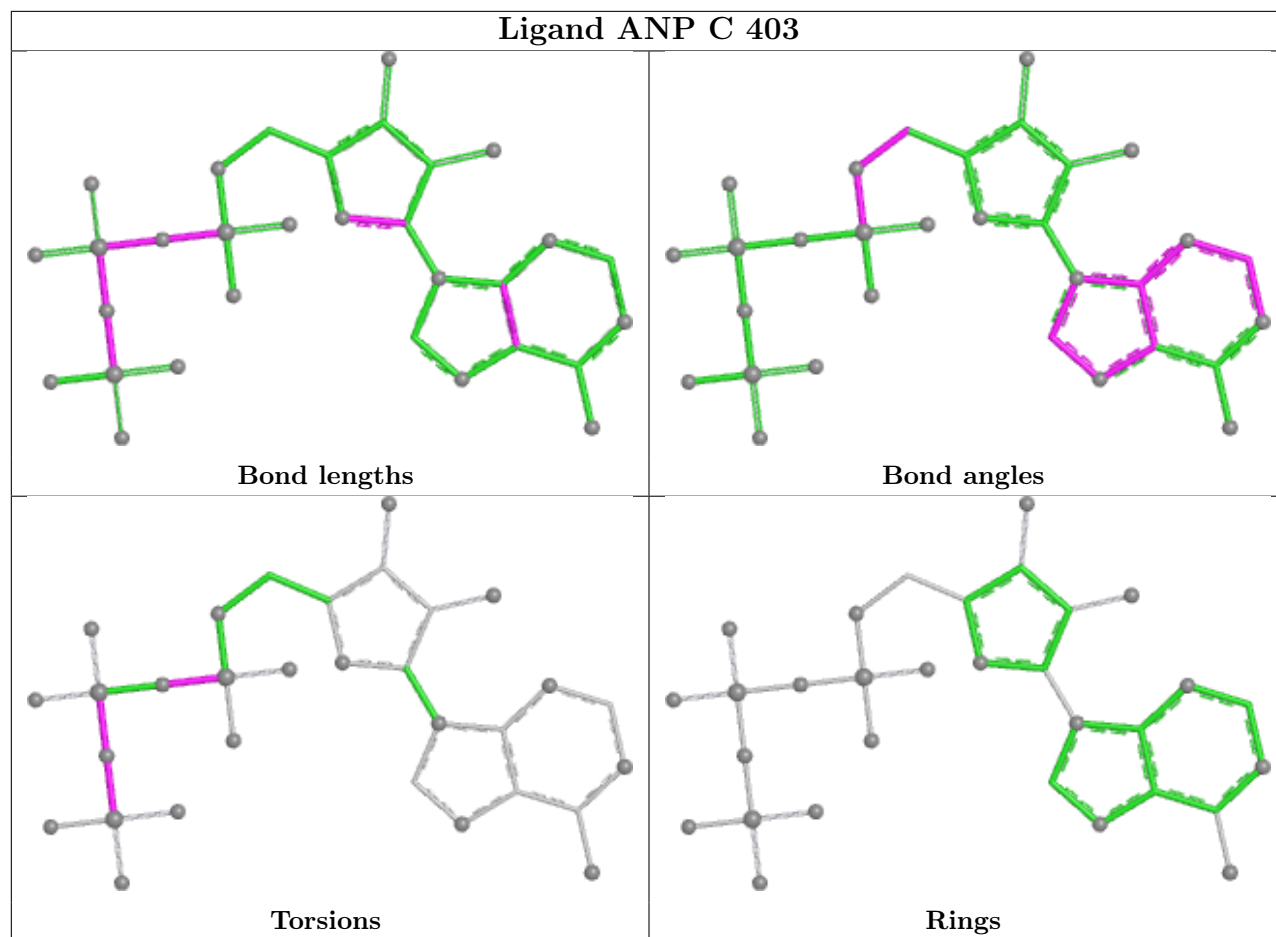
Mol	Chain	Res	Type	Atoms
2	A	401	ANP	PB-N3B-PG-O1G
2	B	402	ANP	PB-N3B-PG-O1G
2	B	402	ANP	PG-N3B-PB-O1B
2	C	403	ANP	PB-N3B-PG-O1G
2	C	403	ANP	PG-N3B-PB-O1B
2	C	403	ANP	PG-N3B-PB-O3A
2	C	403	ANP	PB-O3A-PA-O2A
2	C	403	ANP	PB-O3A-PA-O1A

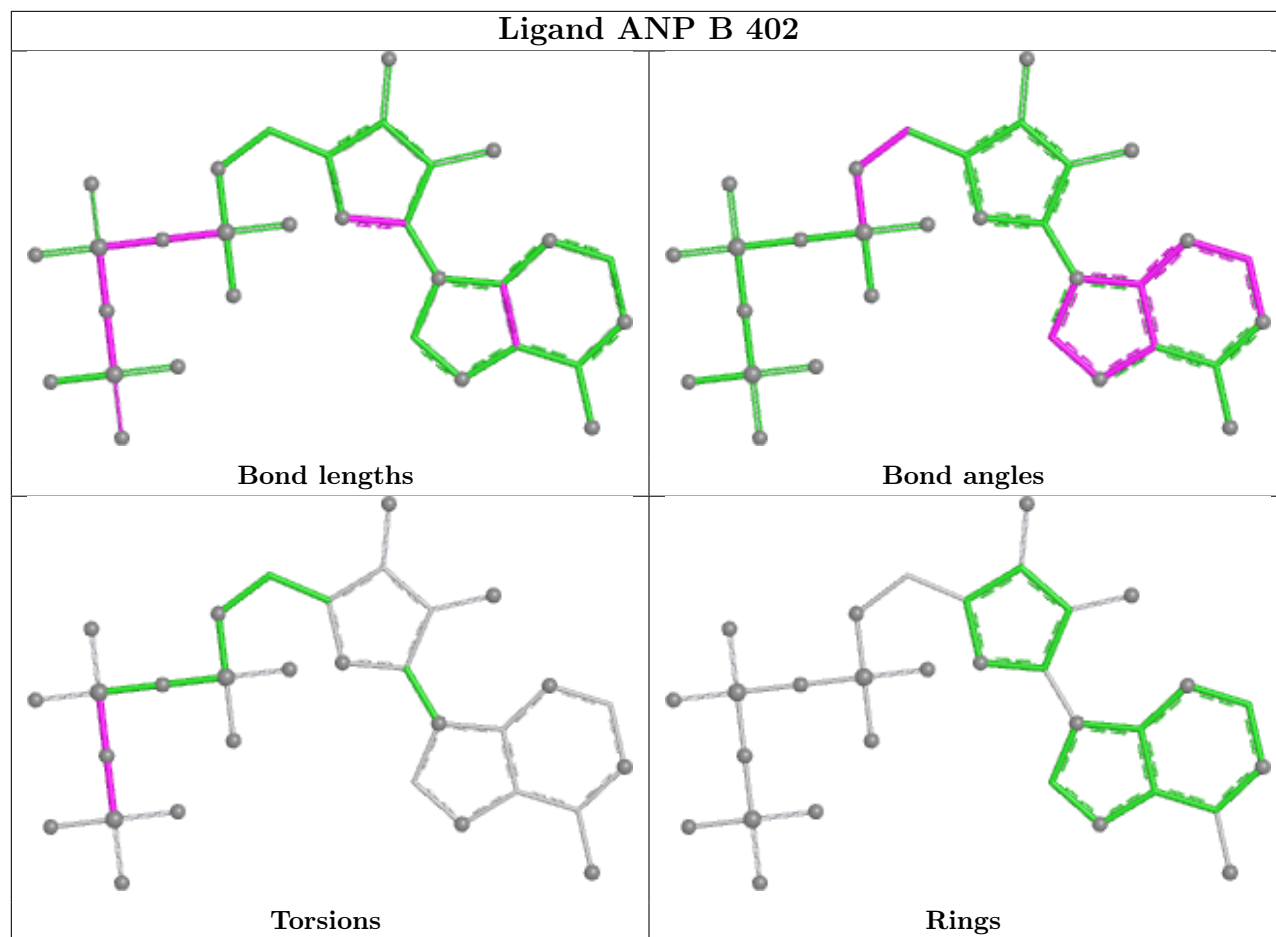
There are no ring outliers.

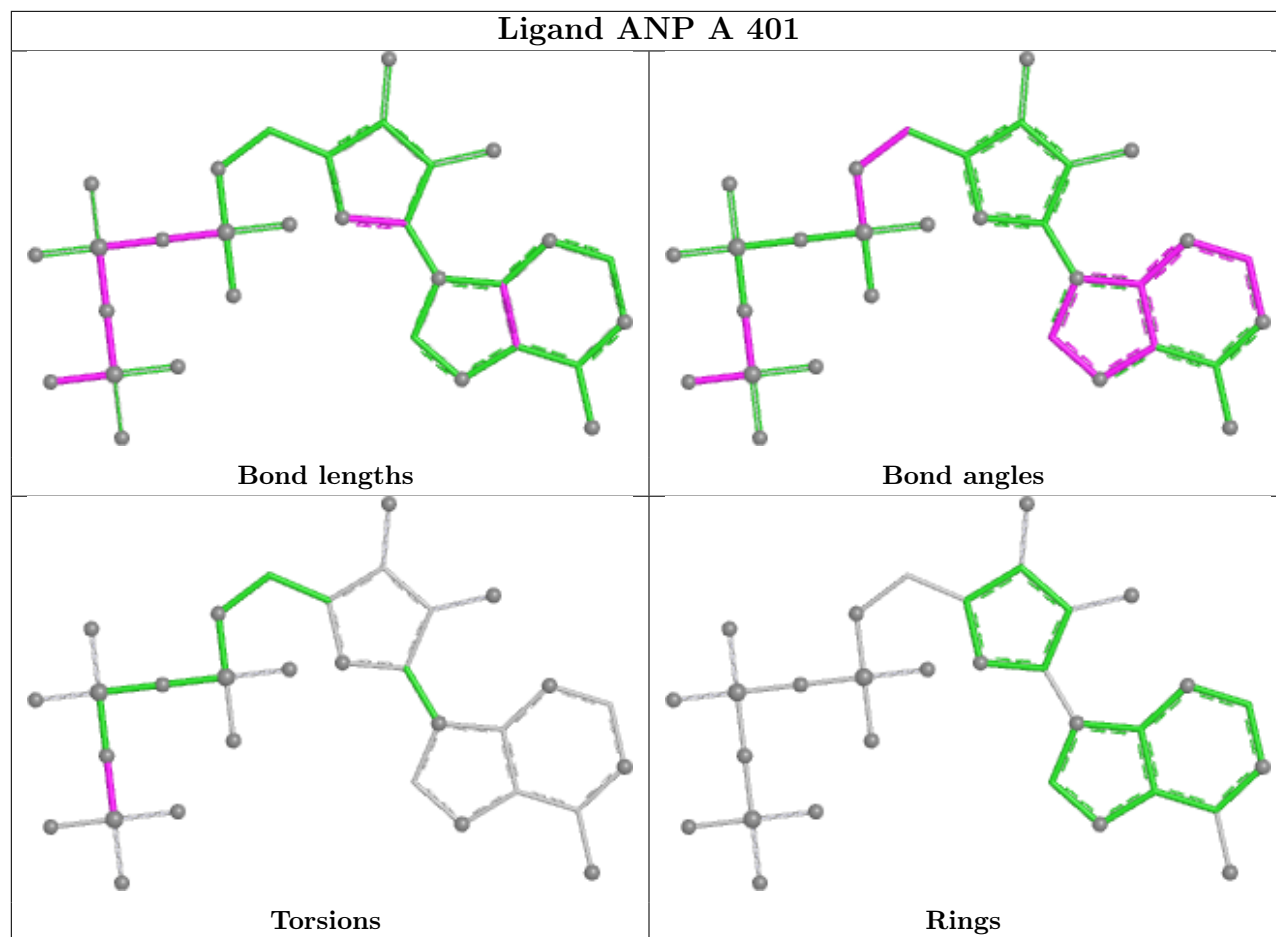
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	ANP	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 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	317/319 (99%)	0.29	10 (3%) 50 44	23, 43, 68, 77	0
1	B	313/319 (98%)	0.45	11 (3%) 47 41	27, 50, 78, 86	0
1	C	309/319 (96%)	0.31	17 (5%) 30 25	22, 42, 74, 89	0
All	All	939/957 (98%)	0.35	38 (4%) 42 37	22, 45, 73, 89	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	182	TYR	3.8
1	C	232	GLY	3.6
1	C	159	SER	3.5
1	B	229	TYR	3.5
1	B	236	LEU	3.5
1	C	158	GLY	3.3
1	C	162	GLY	3.3
1	A	57	ALA	3.1
1	A	160	SER	3.0
1	B	230	THR	2.9
1	C	275	GLU	2.7
1	C	166	VAL	2.7
1	C	169	PHE	2.7
1	C	294	MET	2.6
1	A	162	GLY	2.6
1	B	231	PRO	2.6
1	C	163	ILE	2.5
1	A	182	TYR	2.5
1	A	161	VAL	2.5
1	C	167	GLU	2.5
1	A	163	ILE	2.4
1	B	166	VAL	2.4
1	A	225	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	167	GLU	2.3
1	A	294	MET	2.3
1	B	222	PHE	2.3
1	C	150	PHE	2.3
1	C	236	LEU	2.3
1	B	72	GLU	2.2
1	A	150	PHE	2.2
1	B	160	SER	2.2
1	B	242	LEU	2.2
1	C	172	LEU	2.1
1	C	233	ARG	2.1
1	C	178	LEU	2.1
1	C	85	PHE	2.1
1	A	319	THR	2.0
1	B	84	ARG	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](#)

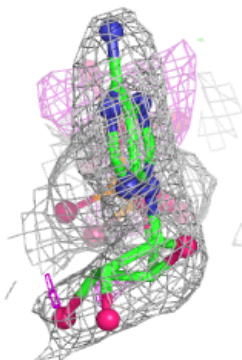
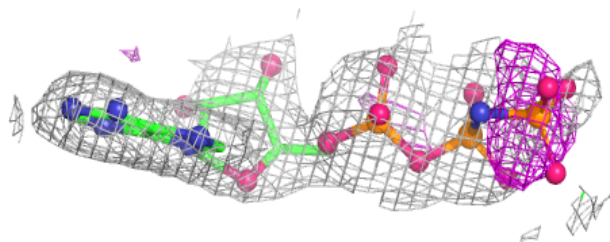
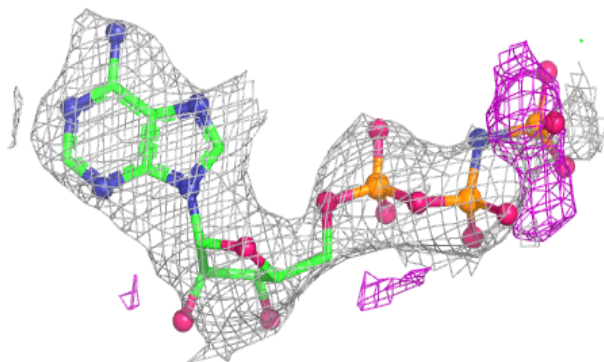
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	ANP	B	402	31/31	0.80	0.18	84,93,118,118	0
2	ANP	A	401	31/31	0.82	0.18	105,109,130,130	0
2	ANP	C	403	31/31	0.83	0.15	76,84,102,104	0

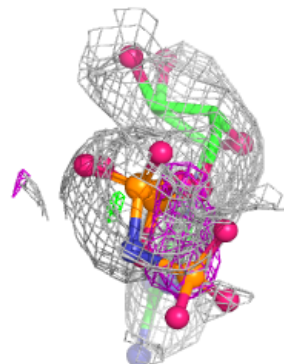
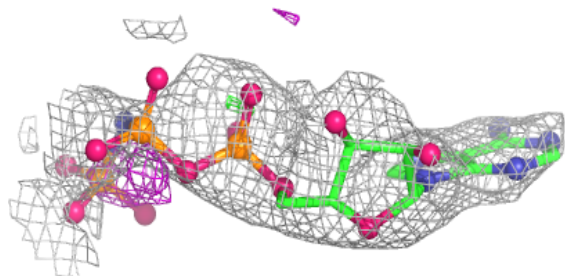
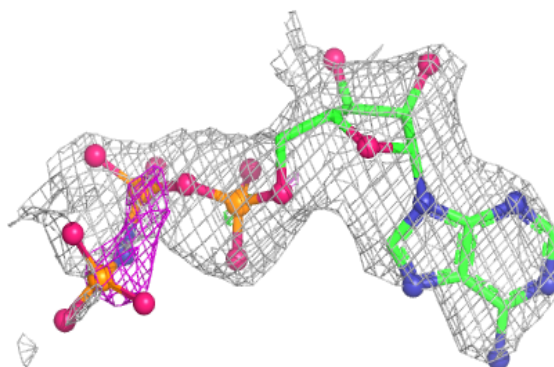
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

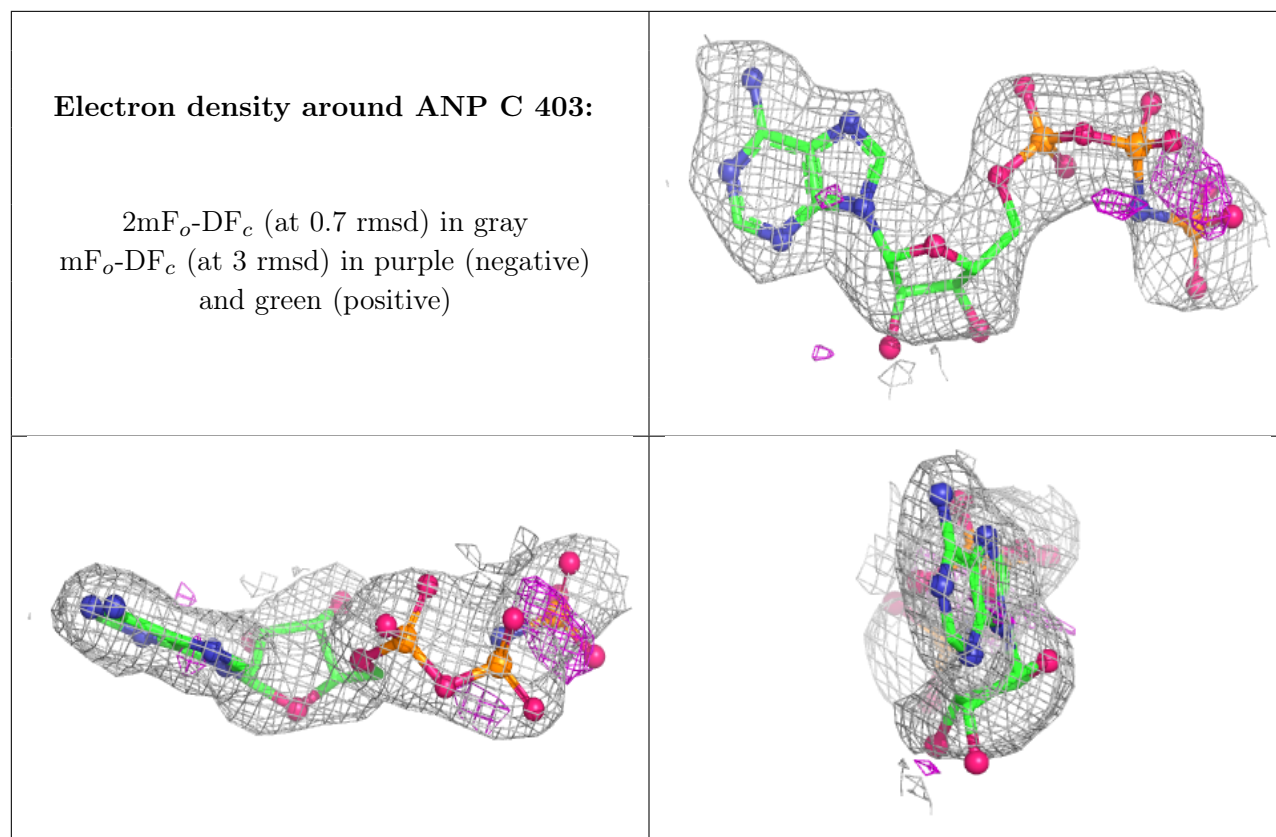
Electron density around ANP B 402:

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

**Electron density around ANP A 401:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.