



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 09:40 PM UTC

PDB ID : 5ZMF / pdb_00005zmf
Title : AMPPNP complex of C. reinhardtii ArsA1
Authors : Lin, T.W.; Hsiao, C.D.; Chang, H.Y.
Deposited on : 2018-04-03
Resolution : 3.56 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

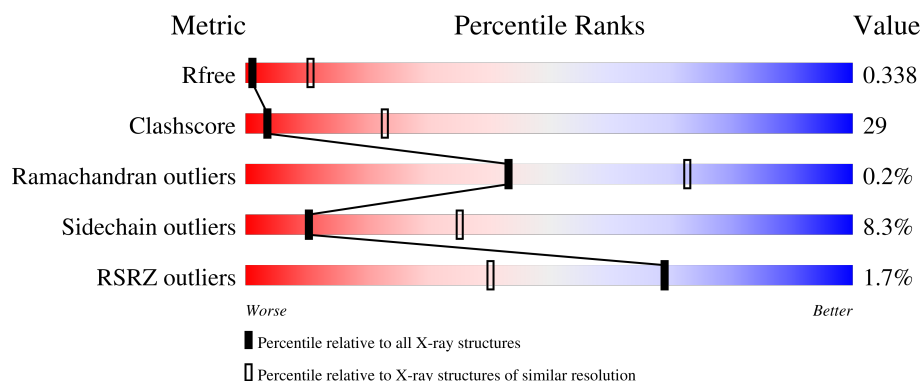
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.56 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	1410 (3.62-3.50)
Clashscore	190562	1480 (3.62-3.50)
Ramachandran outliers	187476	1440 (3.62-3.50)
Sidechain outliers	187428	1441 (3.62-3.50)
RSRZ outliers	180081	1409 (3.62-3.50)

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	687	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4911 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 ATPase ARSA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	637	Total	C	N	O	S	0	0	0
			4847	3080	838	912	17			

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

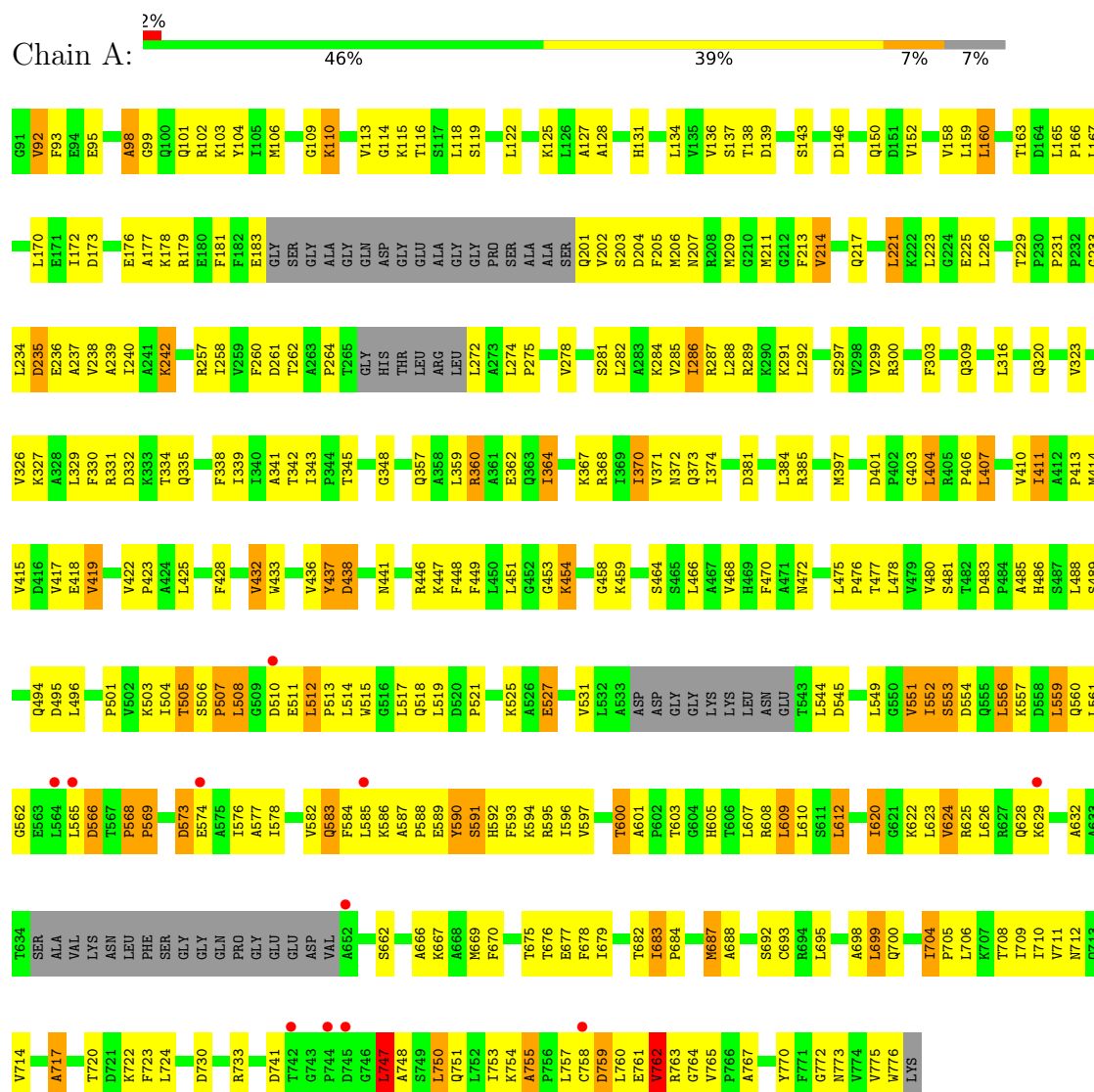
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		

3 Residue-property plots

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: ATPase ARSA1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	90.46Å 90.46Å 306.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.72 – 3.56 29.72 – 3.56	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.72-3.56) 99.3 (29.72-3.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 3.55Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.275 , 0.327 0.279 , 0.338	Depositor DCC
R_{free} test set	1609 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	137.1	Xtriage
Anisotropy	0.408	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 69.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4911	wwPDB-VP
Average B, all atoms (Å ²)	133.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: ANP, MG

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	0.78	3/4925 (0.1%)	1.24	35/6672 (0.5%)

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	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	603	THR	CA-CB	7.90	1.61	1.54
1	A	683	ILE	CA-C	7.02	1.58	1.52
1	A	569	PRO	CA-C	6.10	1.57	1.52

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	683	ILE	N-CA-C	9.28	118.14	107.84
1	A	417	VAL	CB-CA-C	-8.76	103.51	111.74
1	A	569	PRO	CA-C-N	8.08	129.94	119.84
1	A	569	PRO	C-N-CA	8.08	129.94	119.84
1	A	98	ALA	N-CA-C	8.01	120.72	111.11
1	A	712	ASN	CA-C-N	-7.95	111.80	122.85
1	A	712	ASN	C-N-CA	-7.95	111.80	122.85
1	A	755	ALA	CA-C-N	7.92	127.94	119.78
1	A	755	ALA	C-N-CA	7.92	127.94	119.78
1	A	568	PRO	N-CA-C	7.89	120.33	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	747	LEU	N-CA-C	7.80	119.54	108.00
1	A	554	ASP	N-CA-C	7.63	123.64	113.72
1	A	483	ASP	CA-C-N	7.24	127.40	119.87
1	A	483	ASP	C-N-CA	7.24	127.40	119.87
1	A	437	TYR	N-CA-C	6.85	119.73	109.25
1	A	566	ASP	N-CA-C	6.83	120.08	111.69
1	A	765	VAL	CB-CA-C	-6.52	105.46	114.00
1	A	761	GLU	N-CA-C	-6.45	104.33	111.36
1	A	553	SER	N-CA-C	6.43	118.63	108.67
1	A	419	VAL	CB-CA-C	-6.37	103.08	111.81
1	A	527	GLU	N-CA-C	-6.35	103.88	111.69
1	A	92	VAL	CB-CA-C	-6.02	104.15	112.04
1	A	717	ALA	N-CA-C	5.94	118.24	111.11
1	A	612	LEU	CA-C-N	-5.88	113.62	119.56
1	A	612	LEU	C-N-CA	-5.88	113.62	119.56
1	A	603	THR	N-CA-CB	5.79	115.41	109.51
1	A	360	ARG	N-CA-C	-5.62	105.92	112.89
1	A	364	ILE	CB-CA-C	-5.48	104.72	110.53
1	A	404	LEU	N-CA-C	-5.42	105.40	112.23
1	A	214	VAL	N-CA-C	5.31	116.09	110.72
1	A	508	LEU	N-CA-C	-5.31	103.39	110.55
1	A	438	ASP	N-CA-C	-5.12	106.62	112.92
1	A	747	LEU	CB-CA-C	-5.12	109.17	116.34
1	A	137	SER	CA-C-N	-5.08	115.26	122.77
1	A	137	SER	C-N-CA	-5.08	115.26	122.77

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	110	LYS	Peptide
1	A	505	THR	Peptide
1	A	507	PRO	Peptide

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	4847	0	4969	291	0
2	A	62	0	26	15	0
3	A	2	0	0	0	0
All	All	4911	0	4995	291	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 29.

All (291) 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:99:GLY:H	1:A:257:ARG:HH22	1.10	0.93
1:A:214:VAL:HG22	1:A:632:ALA:HB1	1.52	0.92
1:A:675:THR:HG23	1:A:676:THR:HG23	1.56	0.88
1:A:521:PRO:HB3	1:A:574:GLU:HG3	1.57	0.87
1:A:418:GLU:HG3	2:A:803:ANP:O2'	1.75	0.86
1:A:448:PHE:HB2	1:A:676:THR:HG22	1.58	0.86
1:A:384:LEU:HD21	1:A:724:LEU:HB2	1.59	0.84
1:A:504:ILE:HG13	1:A:514:LEU:HD22	1.60	0.83
1:A:556:LEU:HD13	1:A:557:LYS:HG2	1.63	0.80
1:A:620:ILE:HA	1:A:623:LEU:HG	1.62	0.80
1:A:404:LEU:HD21	1:A:693:CYS:HA	1.64	0.79
1:A:458:GLY:HA2	2:A:803:ANP:C5'	2.12	0.79
1:A:586:LYS:H	1:A:590:TYR:HB3	1.48	0.79
1:A:612:LEU:HD13	1:A:667:LYS:HG3	1.65	0.78
1:A:201:GLN:HG3	1:A:202:VAL:H	1.51	0.76
1:A:357:GLN:HG2	1:A:747:LEU:HG	1.68	0.76
1:A:760:LEU:HB2	1:A:762:VAL:HG22	1.67	0.75
1:A:99:GLY:H	1:A:257:ARG:NH2	1.84	0.74
1:A:470:PHE:HB3	1:A:477:THR:HG21	1.70	0.74
1:A:240:ILE:HG12	1:A:330:PHE:HZ	1.53	0.73
1:A:272:LEU:HD21	1:A:327:LYS:HG3	1.70	0.73
1:A:95:GLU:CD	1:A:368:ARG:HH22	1.98	0.72
1:A:165:LEU:HD12	1:A:166:PRO:HD2	1.71	0.71
1:A:116:THR:N	2:A:801:ANP:O2B	2.20	0.70
1:A:118:LEU:HD22	1:A:370:ILE:HD12	1.75	0.69
1:A:287:ARG:HD2	1:A:557:LYS:HZ1	1.55	0.69
1:A:110:LYS:HA	1:A:264:PRO:HB3	1.75	0.68
1:A:480:VAL:HG22	1:A:517:LEU:HD11	1.76	0.67
1:A:291:LYS:HD3	1:A:552:ILE:HG12	1.76	0.67
1:A:367:LYS:HB2	1:A:368:ARG:HH11	1.60	0.67
1:A:106:MET:HG3	1:A:338:PHE:CE1	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:SER:HA	1:A:300:ARG:HH11	1.60	0.66
1:A:458:GLY:HA2	2:A:803:ANP:H5'1	1.77	0.65
1:A:233:GLY:HA3	1:A:274:LEU:HD11	1.79	0.64
1:A:95:GLU:OE1	1:A:368:ARG:NH2	2.21	0.64
1:A:131:HIS:CE1	1:A:257:ARG:HE	2.15	0.64
1:A:214:VAL:HG13	1:A:629:LYS:HE2	1.78	0.64
1:A:600:THR:HB	1:A:605:HIS:ND1	2.14	0.63
1:A:458:GLY:HA2	2:A:803:ANP:H5'2	1.80	0.63
1:A:339:ILE:HG12	1:A:368:ARG:HB2	1.81	0.62
1:A:507:PRO:HB2	1:A:508:LEU:O	1.98	0.62
1:A:119:SER:OG	1:A:261:ASP:HB2	2.00	0.61
1:A:609:LEU:HD22	1:A:610:LEU:H	1.65	0.61
1:A:699:LEU:HG	1:A:704:ILE:HD12	1.82	0.61
1:A:454:LYS:CD	1:A:454:LYS:H	2.13	0.61
1:A:414:MET:HG2	1:A:415:VAL:H	1.64	0.61
1:A:475:LEU:HG	1:A:594:LYS:HD2	1.83	0.61
1:A:565:LEU:HD13	1:A:622:LYS:NZ	2.16	0.61
1:A:177:ALA:HB3	1:A:235:ASP:HB3	1.82	0.60
1:A:730:ASP:HA	1:A:733:ARG:HD2	1.84	0.60
1:A:103:LYS:HB2	1:A:257:ARG:HH11	1.66	0.59
1:A:159:LEU:HD12	1:A:160:LEU:H	1.66	0.59
1:A:370:ILE:HG22	1:A:753:ILE:HB	1.83	0.59
1:A:447:LYS:HB2	1:A:595:ARG:HG2	1.84	0.59
1:A:669:MET:HG3	1:A:670:PHE:N	2.17	0.59
1:A:134:LEU:HD11	1:A:170:LEU:HB2	1.85	0.58
1:A:118:LEU:HD13	1:A:370:ILE:HD11	1.85	0.58
1:A:419:VAL:HG21	1:A:428:PHE:HD2	1.69	0.58
1:A:741:ASP:O	1:A:748:ALA:HB2	2.04	0.58
1:A:478:LEU:HB2	1:A:593:PHE:CE2	2.38	0.58
1:A:566:ASP:O	1:A:569:PRO:HD2	2.04	0.58
1:A:601:ALA:H	1:A:605:HIS:CE1	2.20	0.58
1:A:289:ARG:O	1:A:309:GLN:HG3	2.04	0.58
1:A:278:VAL:HG13	1:A:323:VAL:HG11	1.86	0.57
1:A:373:GLN:HB3	1:A:757:LEU:HD23	1.87	0.57
1:A:678:PHE:HB2	1:A:706:LEU:HA	1.85	0.57
1:A:586:LYS:HB3	1:A:588:PRO:HD2	1.86	0.57
1:A:211:MET:HE2	1:A:211:MET:HA	1.87	0.57
1:A:299:VAL:HG21	1:A:551:VAL:HG11	1.86	0.57
1:A:505:THR:HG22	1:A:512:LEU:HD22	1.87	0.57
1:A:372:ASN:HA	1:A:755:ALA:HB3	1.87	0.57
1:A:556:LEU:CD1	1:A:557:LYS:HG2	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:342:THR:O	1:A:371:VAL:HA	2.05	0.56
1:A:128:ALA:HA	1:A:165:LEU:HD11	1.88	0.56
1:A:131:HIS:CE1	1:A:257:ARG:HH21	2.23	0.56
1:A:525:LYS:HA	1:A:576:ILE:HD12	1.88	0.56
1:A:451:LEU:O	1:A:600:THR:HG23	2.06	0.56
1:A:669:MET:HE2	1:A:670:PHE:CD1	2.40	0.55
1:A:418:GLU:CG	2:A:803:ANP:O2'	2.51	0.55
1:A:574:GLU:HG2	1:A:605:HIS:CE1	2.41	0.55
1:A:274:LEU:O	1:A:278:VAL:HG12	2.07	0.55
1:A:329:LEU:O	1:A:332:ASP:HB3	2.07	0.55
1:A:341:ALA:HA	1:A:370:ILE:HG12	1.89	0.54
1:A:113:VAL:HG21	1:A:342:THR:HA	1.87	0.54
1:A:327:LYS:HE2	1:A:331:ARG:NH1	2.22	0.54
1:A:343:ILE:HG13	1:A:345:THR:HG22	1.90	0.54
1:A:446:ARG:HA	1:A:594:LYS:O	2.06	0.54
1:A:299:VAL:CG2	1:A:551:VAL:HG11	2.37	0.54
1:A:231:PRO:HB3	1:A:281:SER:HB3	1.89	0.54
1:A:370:ILE:HA	1:A:753:ILE:HB	1.90	0.54
1:A:419:VAL:HG21	1:A:428:PHE:CD2	2.43	0.54
1:A:291:LYS:NZ	1:A:553:SER:O	2.41	0.54
1:A:374:ILE:HG22	1:A:755:ALA:O	2.07	0.53
1:A:260:PHE:HE1	1:A:262:THR:HA	1.73	0.53
1:A:237:ALA:HB2	1:A:274:LEU:HD23	1.90	0.53
1:A:92:VAL:HG13	1:A:751:GLN:HE22	1.73	0.53
1:A:287:ARG:HD2	1:A:557:LYS:NZ	2.24	0.53
1:A:92:VAL:HG13	1:A:751:GLN:NE2	2.24	0.53
1:A:449:PHE:HD1	1:A:677:GLU:HG2	1.72	0.53
1:A:240:ILE:HG12	1:A:330:PHE:CZ	2.41	0.52
1:A:506:SER:N	1:A:507:PRO:HD3	2.24	0.52
1:A:521:PRO:HB2	1:A:573:ASP:HB2	1.91	0.52
1:A:489:SER:HB3	1:A:495:ASP:HA	1.92	0.52
1:A:545:ASP:O	1:A:549:LEU:HG	2.08	0.52
1:A:173:ASP:OD1	1:A:173:ASP:N	2.37	0.52
1:A:288:LEU:HD11	1:A:557:LYS:HG3	1.91	0.52
1:A:221:LEU:HB3	1:A:223:LEU:CD1	2.41	0.51
1:A:503:LYS:HG3	1:A:515:TRP:NE1	2.26	0.51
1:A:584:PHE:HB3	1:A:589:GLU:OE1	2.10	0.51
1:A:272:LEU:O	1:A:275:PRO:HD2	2.09	0.51
1:A:589:GLU:O	1:A:592:HIS:N	2.41	0.51
1:A:103:LYS:HB2	1:A:257:ARG:NH1	2.25	0.51
1:A:204:ASP:O	1:A:207:ASN:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:THR:HG23	1:A:348:GLY:H	1.76	0.51
1:A:368:ARG:N	1:A:750:LEU:HD12	2.26	0.51
1:A:272:LEU:HD11	1:A:327:LYS:HE3	1.92	0.51
1:A:285:VAL:CG2	1:A:561:LEU:HD21	2.40	0.51
1:A:605:HIS:CD2	1:A:605:HIS:H	2.29	0.51
1:A:139:ASP:OD2	1:A:143:SER:HB2	2.11	0.50
1:A:404:LEU:HD13	1:A:407:LEU:HD11	1.92	0.50
1:A:213:PHE:O	1:A:217:GLN:HG2	2.12	0.50
1:A:181:PHE:HE2	1:A:242:LYS:HZ3	1.59	0.50
1:A:477:THR:HA	1:A:595:ARG:O	2.11	0.50
1:A:286:ILE:HD11	1:A:316:LEU:C	2.37	0.50
1:A:297:SER:HA	1:A:300:ARG:HE	1.76	0.50
1:A:272:LEU:HB3	1:A:362:GLU:OE1	2.11	0.49
1:A:238:VAL:O	1:A:242:LYS:HG3	2.13	0.49
1:A:104:TYR:CE2	1:A:258:ILE:HD13	2.47	0.49
1:A:370:ILE:HG22	1:A:753:ILE:CB	2.42	0.49
1:A:260:PHE:CE1	1:A:262:THR:HA	2.47	0.49
1:A:286:ILE:HD11	1:A:316:LEU:HB3	1.94	0.49
1:A:415:VAL:N	2:A:803:ANP:N1	2.61	0.49
1:A:441:ASN:HB3	1:A:595:ARG:NH2	2.28	0.49
1:A:760:LEU:HB2	1:A:762:VAL:CG2	2.42	0.48
1:A:767:ALA:O	1:A:770:TYR:HB3	2.12	0.48
1:A:587:ALA:O	1:A:591:SER:OG	2.31	0.48
1:A:428:PHE:O	1:A:432:VAL:HG22	2.13	0.48
1:A:758:CYS:HB2	1:A:759:ASP:HB3	1.96	0.48
1:A:684:PRO:HD3	1:A:711:VAL:HG12	1.96	0.48
1:A:103:LYS:CB	1:A:257:ARG:HH11	2.25	0.48
1:A:127:ALA:HB2	1:A:167:LEU:HD12	1.96	0.47
1:A:226:LEU:HD21	1:A:561:LEU:HB3	1.96	0.47
1:A:519:LEU:HD12	1:A:577:ALA:HB1	1.97	0.47
1:A:679:ILE:HG12	1:A:708:THR:HB	1.96	0.47
1:A:285:VAL:HG22	1:A:561:LEU:HD21	1.97	0.47
1:A:494:GLN:HE22	1:A:504:ILE:HG23	1.77	0.47
1:A:414:MET:HE1	1:A:723:PHE:HE2	1.79	0.47
1:A:578:ILE:O	1:A:582:VAL:HG23	2.14	0.47
1:A:109:GLY:O	1:A:264:PRO:HA	2.15	0.47
1:A:173:ASP:OD2	1:A:176:GLU:HB2	2.14	0.47
1:A:425:LEU:HD22	1:A:464:SER:HB2	1.95	0.47
1:A:122:LEU:HB2	1:A:776:TRP:CZ2	2.50	0.47
1:A:330:PHE:O	1:A:364:ILE:HG23	2.14	0.47
1:A:436:VAL:HG12	1:A:437:TYR:H	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:LEU:HD22	1:A:622:LYS:HE3	1.97	0.47
1:A:762:VAL:HG21	2:A:801:ANP:N3	2.30	0.47
1:A:116:THR:HG21	1:A:143:SER:OG	2.14	0.47
1:A:414:MET:N	1:A:714:VAL:O	2.48	0.47
1:A:449:PHE:CD1	1:A:677:GLU:HG2	2.49	0.47
1:A:454:LYS:H	1:A:454:LYS:HD3	1.80	0.47
1:A:478:LEU:HD12	1:A:515:TRP:O	2.14	0.47
1:A:544:LEU:O	1:A:544:LEU:HD23	2.15	0.47
1:A:586:LYS:N	1:A:590:TYR:HB3	2.23	0.47
1:A:496:LEU:HD21	1:A:504:ILE:HD11	1.95	0.46
1:A:556:LEU:HD12	1:A:557:LYS:H	1.81	0.46
1:A:717:ALA:HA	1:A:720:THR:OG1	2.15	0.46
1:A:373:GLN:CB	1:A:757:LEU:HD23	2.46	0.46
1:A:403:GLY:C	1:A:406:PRO:HD2	2.40	0.46
1:A:747:LEU:N	1:A:747:LEU:HD13	2.30	0.46
1:A:152:VAL:HG12	1:A:158:VAL:HB	1.96	0.46
1:A:476:PRO:HA	1:A:513:PRO:O	2.14	0.46
1:A:687:MET:HG3	1:A:688:ALA:N	2.31	0.46
1:A:438:ASP:OD1	1:A:441:ASN:ND2	2.49	0.46
1:A:544:LEU:HD11	1:A:629:LYS:HG3	1.98	0.46
1:A:757:LEU:HD13	1:A:758:CYS:HB3	1.97	0.46
1:A:411:ILE:HG23	1:A:714:VAL:HG21	1.98	0.46
1:A:709:ILE:HG12	1:A:710:ILE:N	2.31	0.46
1:A:559:LEU:O	1:A:562:GLY:N	2.49	0.46
1:A:624:VAL:O	1:A:628:GLN:HG2	2.16	0.46
1:A:565:LEU:HD13	1:A:622:LYS:HZ2	1.81	0.46
1:A:759:ASP:OD1	1:A:759:ASP:N	2.49	0.46
1:A:441:ASN:O	1:A:595:ARG:NH2	2.49	0.45
1:A:125:LYS:HD3	1:A:773:ASN:ND2	2.31	0.45
1:A:573:ASP:OD1	1:A:573:ASP:N	2.48	0.45
1:A:114:GLY:HA2	2:A:801:ANP:O1A	2.16	0.45
1:A:284:LYS:HD3	1:A:560:GLN:OE1	2.16	0.45
1:A:234:LEU:O	1:A:238:VAL:HG23	2.17	0.45
1:A:415:VAL:H	2:A:803:ANP:H2	1.81	0.45
1:A:371:VAL:HG23	1:A:754:LYS:HA	1.99	0.45
1:A:466:LEU:HD23	1:A:597:VAL:HG21	1.99	0.45
1:A:115:LYS:N	2:A:801:ANP:O2B	2.50	0.45
1:A:374:ILE:HG22	1:A:755:ALA:C	2.42	0.45
1:A:625:ARG:O	1:A:628:GLN:HB2	2.17	0.45
1:A:284:LYS:NZ	1:A:560:GLN:HB3	2.32	0.45
1:A:678:PHE:C	1:A:679:ILE:HG13	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:764:GLY:O	1:A:767:ALA:HB3	2.17	0.45
1:A:178:LYS:HB2	1:A:234:LEU:HD23	1.98	0.44
1:A:203:SER:HB2	1:A:205:PHE:CD2	2.52	0.44
1:A:578:ILE:HG21	1:A:670:PHE:HZ	1.83	0.44
1:A:150:GLN:NE2	1:A:160:LEU:HB3	2.32	0.44
1:A:225:GLU:O	1:A:229:THR:HG23	2.17	0.44
1:A:607:LEU:HD23	1:A:607:LEU:HA	1.71	0.44
1:A:670:PHE:O	1:A:705:PRO:HD2	2.18	0.44
1:A:759:ASP:C	1:A:760:LEU:HD12	2.42	0.44
1:A:125:LYS:HD2	1:A:125:LYS:HA	1.82	0.44
1:A:695:LEU:HD23	1:A:699:LEU:HD22	1.99	0.44
1:A:441:ASN:HB3	1:A:595:ARG:HH22	1.83	0.44
1:A:511:GLU:C	1:A:513:PRO:HD3	2.43	0.44
1:A:583:GLN:OE1	1:A:662:SER:HA	2.18	0.44
1:A:695:LEU:O	1:A:699:LEU:HB2	2.18	0.44
1:A:303:PHE:CD1	1:A:303:PHE:N	2.86	0.43
1:A:432:VAL:HG23	1:A:433:TRP:H	1.82	0.43
1:A:517:LEU:HD13	1:A:519:LEU:HD21	1.99	0.43
1:A:136:VAL:HG22	1:A:170:LEU:HD23	2.00	0.43
1:A:217:GLN:NE2	1:A:628:GLN:OE1	2.46	0.43
1:A:239:ALA:HA	1:A:242:LYS:HD2	2.00	0.43
1:A:682:THR:OG1	1:A:683:ILE:N	2.51	0.43
1:A:102:ARG:HE	1:A:102:ARG:HB2	1.65	0.43
1:A:453:GLY:O	1:A:459:LYS:NZ	2.31	0.43
1:A:494:GLN:HB2	1:A:496:LEU:HD12	2.01	0.43
1:A:146:ASP:HB2	1:A:763:ARG:NH1	2.34	0.43
1:A:209:MET:HE2	1:A:209:MET:HB3	1.63	0.43
1:A:292:LEU:HD23	1:A:292:LEU:HA	1.82	0.43
1:A:496:LEU:HD21	1:A:504:ILE:CD1	2.49	0.43
1:A:585:LEU:O	1:A:585:LEU:HD12	2.18	0.43
1:A:131:HIS:HE1	1:A:257:ARG:HH21	1.64	0.43
1:A:411:ILE:CG2	1:A:714:VAL:HG21	2.49	0.43
1:A:582:VAL:HG11	1:A:666:ALA:HA	2.00	0.42
1:A:203:SER:HB2	1:A:205:PHE:CE2	2.53	0.42
1:A:282:LEU:HD23	1:A:282:LEU:HA	1.80	0.42
1:A:385:ARG:HA	1:A:385:ARG:HD3	1.68	0.42
1:A:750:LEU:N	1:A:750:LEU:HD23	2.34	0.42
1:A:403:GLY:O	1:A:700:GLN:NE2	2.52	0.42
1:A:422:VAL:N	1:A:423:PRO:HD2	2.34	0.42
1:A:549:LEU:HA	1:A:552:ILE:O	2.18	0.42
1:A:177:ALA:O	1:A:181:PHE:HD2	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ILE:HG21	1:A:330:PHE:CZ	2.55	0.42
1:A:332:ASP:OD1	1:A:332:ASP:C	2.63	0.42
1:A:367:LYS:C	1:A:750:LEU:HD12	2.45	0.42
1:A:204:ASP:C	1:A:206:MET:N	2.77	0.42
1:A:334:THR:HG23	1:A:335:GLN:HG3	2.00	0.42
1:A:163:THR:OG1	1:A:165:LEU:N	2.36	0.42
1:A:178:LYS:HG2	1:A:179:ARG:N	2.35	0.42
1:A:451:LEU:HG	1:A:679:ILE:HB	2.02	0.42
1:A:397:MET:HE3	1:A:397:MET:HB2	1.86	0.42
1:A:372:ASN:OD1	2:A:801:ANP:N6	2.52	0.42
1:A:339:ILE:HA	1:A:368:ARG:O	2.20	0.41
1:A:413:PRO:HA	1:A:714:VAL:HB	2.02	0.41
1:A:297:SER:HA	1:A:300:ARG:NH1	2.30	0.41
1:A:568:PRO:HB2	1:A:569:PRO:HD3	2.02	0.41
1:A:609:LEU:HD22	1:A:610:LEU:N	2.33	0.41
1:A:730:ASP:O	1:A:733:ARG:HB2	2.20	0.41
1:A:401:ASP:HB3	1:A:404:LEU:HB2	2.02	0.41
1:A:501:PRO:HG3	1:A:517:LEU:HD23	2.02	0.41
1:A:233:GLY:O	1:A:274:LEU:HD21	2.21	0.41
1:A:286:ILE:CG1	1:A:316:LEU:HB3	2.51	0.41
1:A:464:SER:O	1:A:468:VAL:HG23	2.20	0.41
1:A:608:ARG:HB2	1:A:698:ALA:HB1	2.00	0.41
1:A:226:LEU:HD11	1:A:562:GLY:HA2	2.02	0.41
1:A:236:GLU:O	1:A:240:ILE:HD12	2.21	0.41
1:A:360:ARG:HH12	1:A:367:LYS:NZ	2.19	0.41
1:A:481:SER:HB2	1:A:488:LEU:HD13	2.03	0.41
1:A:136:VAL:HG22	1:A:170:LEU:HB3	2.01	0.41
1:A:165:LEU:CD1	1:A:166:PRO:HD2	2.46	0.41
1:A:99:GLY:O	1:A:257:ARG:NH2	2.54	0.41
1:A:201:GLN:HG3	1:A:202:VAL:N	2.29	0.41
1:A:221:LEU:HB3	1:A:223:LEU:HD12	2.02	0.41
1:A:260:PHE:CD1	1:A:262:THR:HG22	2.55	0.41
1:A:762:VAL:HB	2:A:801:ANP:O2'	2.21	0.41
1:A:510:ASP:O	1:A:512:LEU:HD23	2.21	0.41
1:A:605:HIS:CD2	1:A:605:HIS:N	2.89	0.41
1:A:93:PHE:HZ	1:A:125:LYS:O	2.03	0.41
1:A:95:GLU:O	1:A:98:ALA:HB2	2.21	0.41
1:A:204:ASP:HB2	1:A:207:ASN:OD1	2.21	0.41
1:A:217:GLN:NE2	1:A:628:GLN:HB3	2.36	0.41
1:A:503:LYS:HA	1:A:515:TRP:CD1	2.56	0.41
1:A:574:GLU:CG	1:A:605:HIS:CE1	3.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:688:ALA:O	1:A:692:SER:OG	2.39	0.41
1:A:762:VAL:HG21	2:A:801:ANP:C2	2.51	0.41
1:A:485:ALA:O	1:A:486:HIS:HB2	2.20	0.41
1:A:384:LEU:HD13	1:A:720:THR:HG22	2.02	0.40
1:A:410:VAL:HG22	1:A:436:VAL:HG11	2.03	0.40
1:A:415:VAL:H	2:A:803:ANP:C2	2.34	0.40
1:A:448:PHE:HA	1:A:596:ILE:O	2.20	0.40
1:A:494:GLN:HB2	1:A:496:LEU:CD1	2.51	0.40
1:A:282:LEU:HD13	1:A:320:GLN:HB2	2.03	0.40
1:A:559:LEU:HD12	1:A:559:LEU:HA	1.87	0.40
1:A:605:HIS:H	1:A:605:HIS:HD2	1.68	0.40
1:A:414:MET:HG2	1:A:415:VAL:N	2.34	0.40
1:A:772:GLY:O	1:A:776:TRP:HD1	2.05	0.40
1:A:172:ILE:HG21	1:A:242:LYS:HD3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	627/687 (91%)	582 (93%)	44 (7%)	1 (0%)	43 73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	762	VAL

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	521/554 (94%)	478 (92%)	43 (8%)	10	35

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	GLN
1	A	138	THR
1	A	160	LEU
1	A	183	GLU
1	A	221	LEU
1	A	235	ASP
1	A	242	LYS
1	A	286	ILE
1	A	326	VAL
1	A	359	LEU
1	A	370	ILE
1	A	381	ASP
1	A	407	LEU
1	A	411	ILE
1	A	432	VAL
1	A	454	LYS
1	A	472	ASN
1	A	512	LEU
1	A	518	GLN
1	A	527	GLU
1	A	531	VAL
1	A	551	VAL
1	A	552	ILE
1	A	556	LEU
1	A	559	LEU
1	A	573	ASP
1	A	583	GLN
1	A	590	TYR
1	A	591	SER
1	A	600	THR
1	A	609	LEU
1	A	620	ILE
1	A	624	VAL
1	A	626	LEU

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Mol	Chain	Res	Type
1	A	687	MET
1	A	699	LEU
1	A	704	ILE
1	A	722	LYS
1	A	747	LEU
1	A	750	LEU
1	A	759	ASP
1	A	762	VAL
1	A	775	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	HIS
1	A	335	GLN
1	A	494	GLN
1	A	555	GLN
1	A	605	HIS
1	A	713	GLN

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ANP	A	803	-	33,33,33	1.87	6 (18%)	45,52,52	1.85	9 (20%)
2	ANP	A	801	-	33,33,33	1.71	6 (18%)	45,52,52	0.79	1 (2%)

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	A	803	-	-	5/18/38/38	0/3/3/3
2	ANP	A	801	-	-	6/18/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	ANP	PB-O1B	7.13	1.57	1.46
2	A	803	ANP	C5-C4	4.74	1.47	1.39
2	A	803	ANP	PG-N3B	4.63	1.75	1.63
2	A	803	ANP	PB-N3B	4.62	1.75	1.63
2	A	801	ANP	PG-N3B	3.46	1.72	1.63
2	A	801	ANP	PG-O1G	2.95	1.50	1.46
2	A	801	ANP	PB-N3B	2.92	1.71	1.63
2	A	803	ANP	C5-C6	2.78	1.48	1.41
2	A	801	ANP	PB-O2B	-2.64	1.49	1.56
2	A	801	ANP	PA-O3A	2.42	1.62	1.59
2	A	803	ANP	C8-N7	2.36	1.36	1.31
2	A	803	ANP	C5-N7	-2.26	1.35	1.39

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	803	ANP	C5-C4-N3	-5.80	118.73	126.72
2	A	803	ANP	N3-C4-N9	4.60	134.98	127.17
2	A	803	ANP	C2-N3-C4	3.68	120.83	111.83
2	A	803	ANP	C4-C5-N7	-3.47	106.61	110.58
2	A	803	ANP	N3-C2-N1	-3.21	123.72	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	803	ANP	C4-N9-C8	2.60	108.47	105.74
2	A	803	ANP	C5-N7-C8	2.57	107.49	103.45
2	A	803	ANP	C3'-C2'-C1'	2.54	106.27	101.46
2	A	801	ANP	O3G-PG-O1G	-2.10	108.19	113.45
2	A	803	ANP	C6-C5-N7	2.09	136.13	132.09

There are no chirality outliers.

All (11) torsion outliers are listed below:

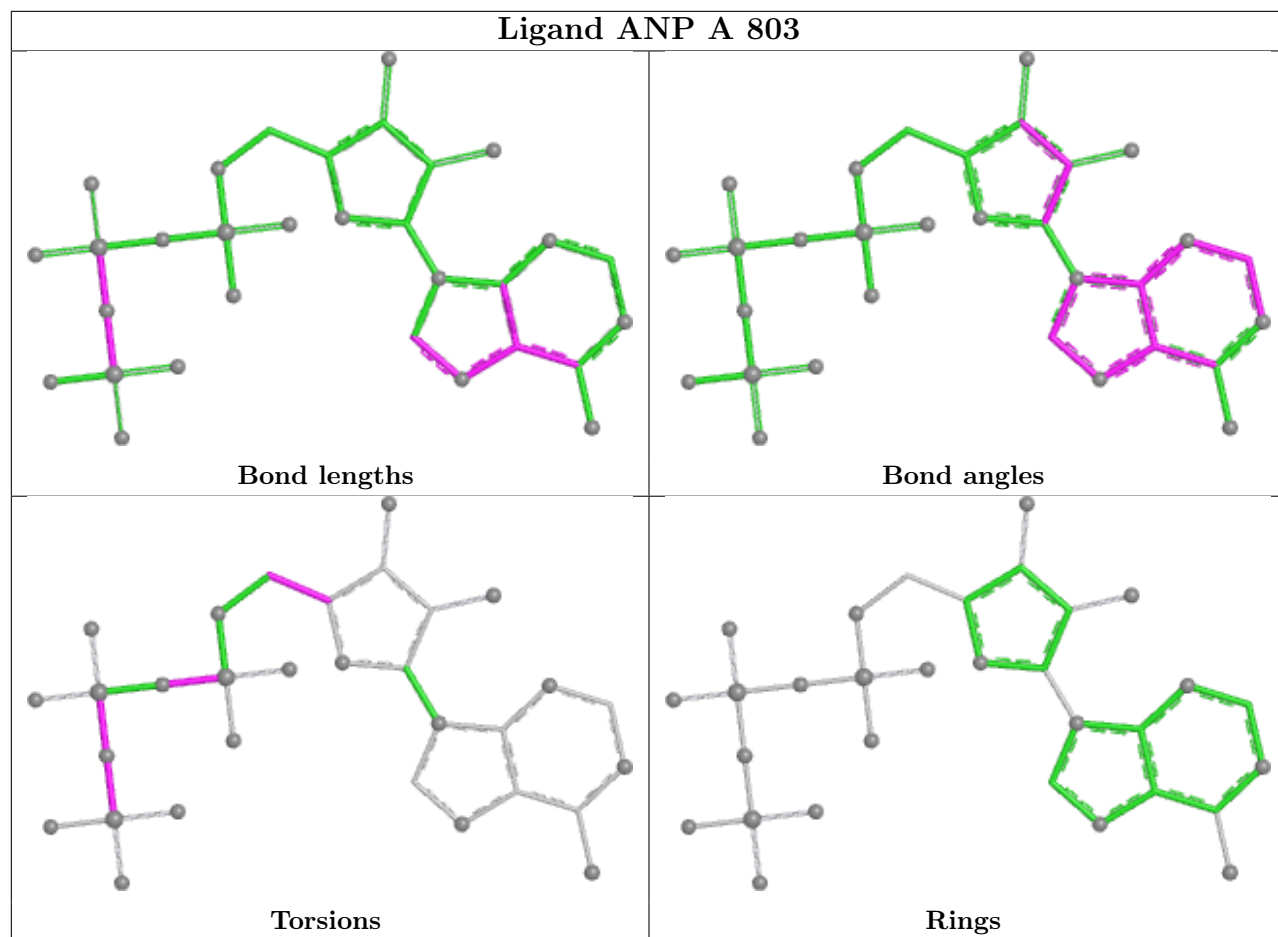
Mol	Chain	Res	Type	Atoms
2	A	801	ANP	PB-N3B-PG-O1G
2	A	801	ANP	C5'-O5'-PA-O1A
2	A	801	ANP	C5'-O5'-PA-O2A
2	A	801	ANP	C5'-O5'-PA-O3A
2	A	801	ANP	O4'-C4'-C5'-O5'
2	A	803	ANP	PB-N3B-PG-O1G
2	A	803	ANP	PG-N3B-PB-O1B
2	A	803	ANP	PB-O3A-PA-O5'
2	A	801	ANP	C3'-C4'-C5'-O5'
2	A	803	ANP	O4'-C4'-C5'-O5'
2	A	803	ANP	C3'-C4'-C5'-O5'

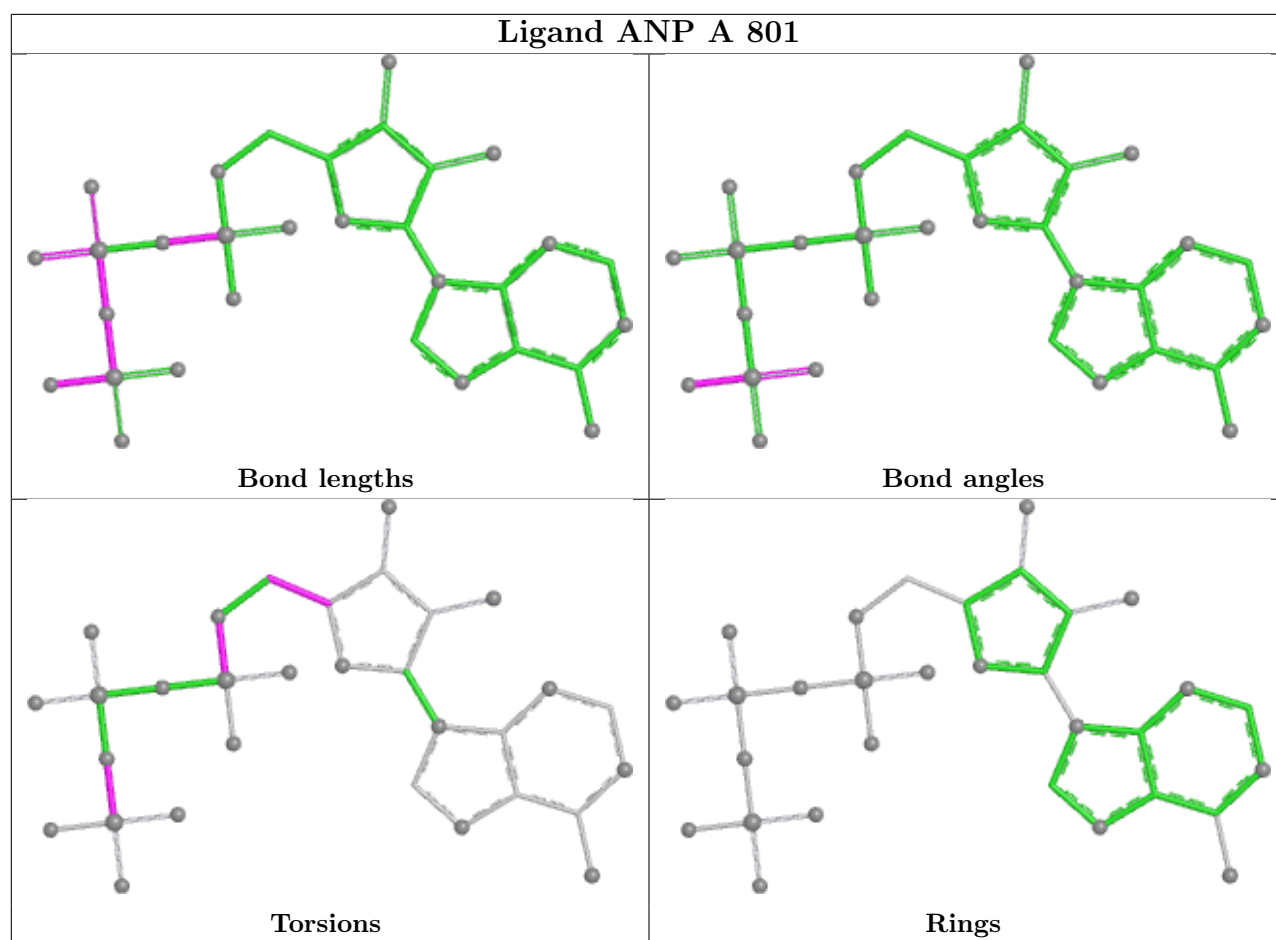
There are no ring outliers.

2 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	803	ANP	8	0
2	A	801	ANP	7	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 [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	637/687 (92%)	-0.10	11 (1%) 69 40	96, 132, 164, 184	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	758	CYS	4.8
1	A	742	THR	3.2
1	A	744	PRO	2.7
1	A	652	ALA	2.6
1	A	745	ASP	2.5
1	A	585	LEU	2.4
1	A	565	LEU	2.3
1	A	564	LEU	2.2
1	A	510	ASP	2.2
1	A	629	LYS	2.1
1	A	574	GLU	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.

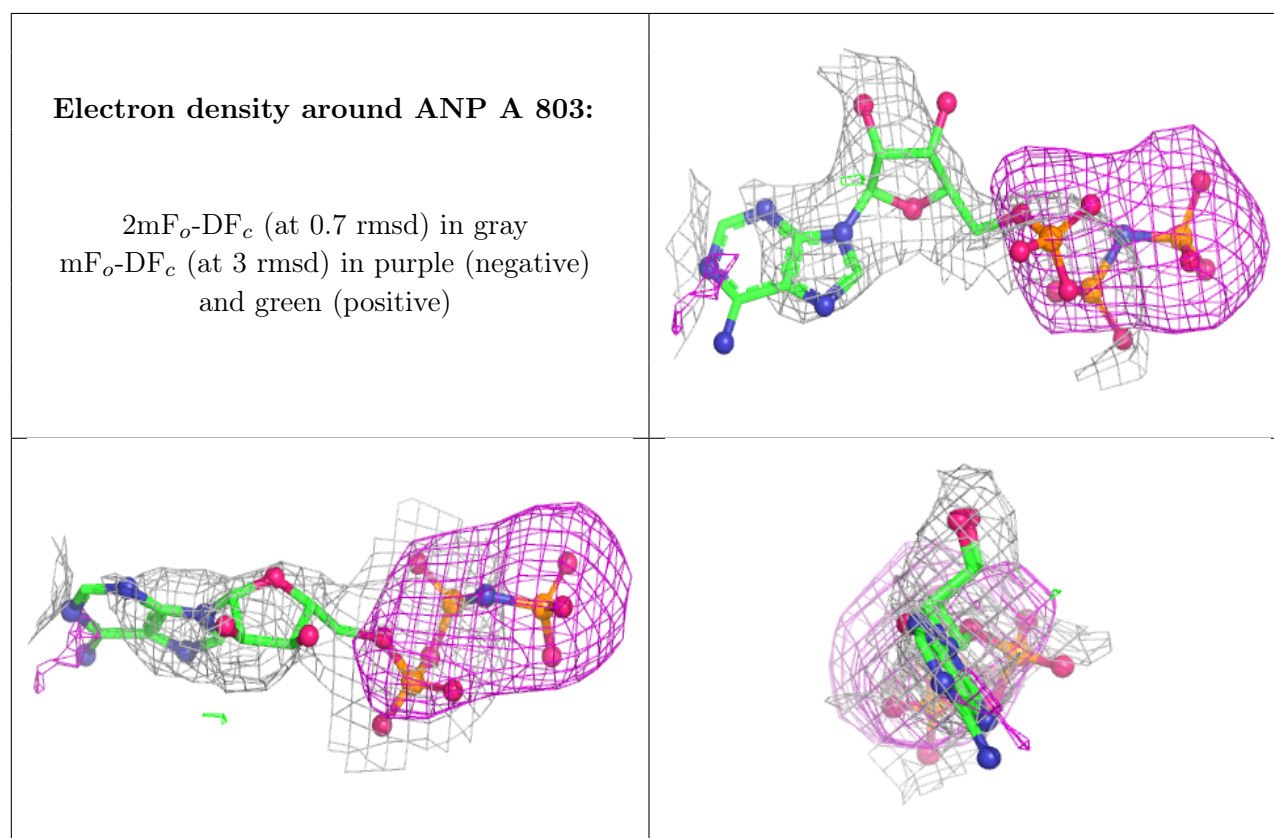
6.4 Ligands [i](#)

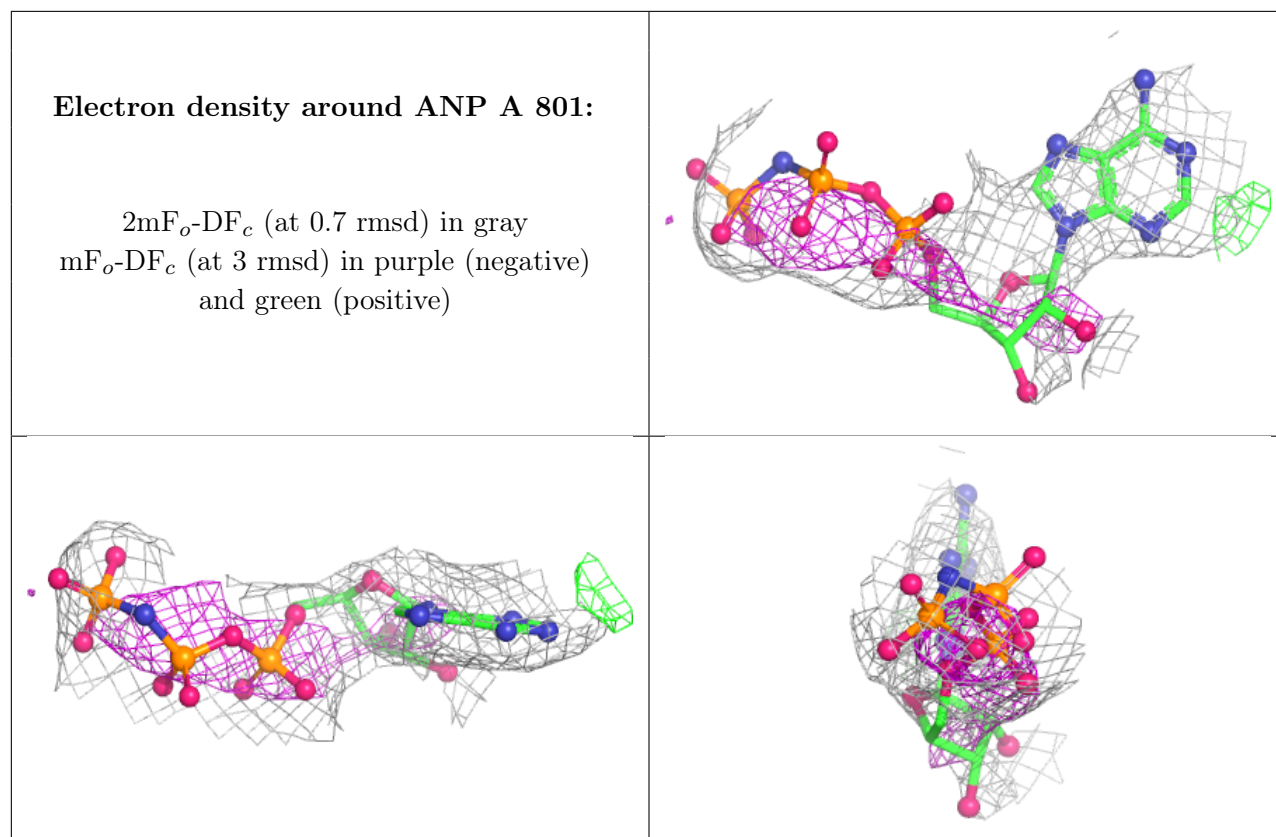
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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(Å ²)	Q<0.9
2	ANP	A	803	31/31	0.51	0.17	140,157,167,169	0
2	ANP	A	801	31/31	0.78	0.11	140,157,167,169	0
3	MG	A	802	1/1	0.83	0.11	126,126,126,126	0
3	MG	A	804	1/1	0.94	0.10	126,126,126,126	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.