



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:28 AM UTC

PDB ID : 5U1G / pdb_00005u1g
Title : Structure of TP228 ParA-AMPPNP-ParB complex
Authors : Schumacher, M.A.
Deposited on : 2016-11-28
Resolution : 3.64 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

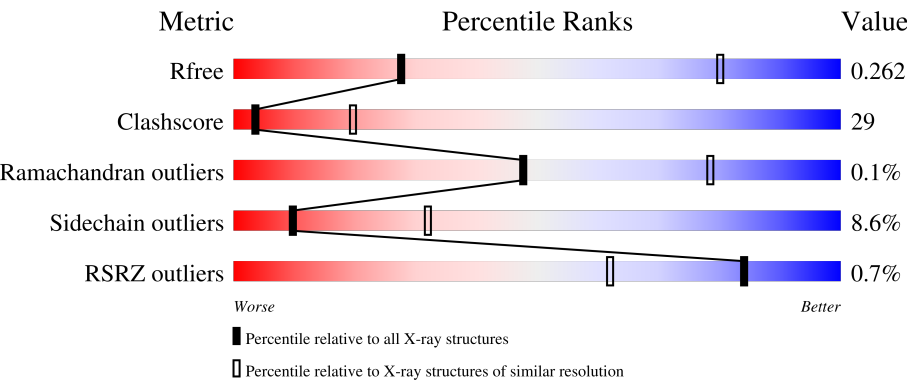
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.64 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	1132 (3.76-3.52)
Clashscore	190562	1014 (3.74-3.54)
Ramachandran outliers	187476	1123 (3.76-3.52)
Sidechain outliers	187428	1121 (3.76-3.52)
RSRZ outliers	180081	1130 (3.76-3.52)

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	214	
1	B	214	
1	C	214	
1	D	214	
2	K	19	

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Mol	Chain	Length	Quality of chain
2	R	19	11%5%16%26%5%47%
2	Y	19	21%16%11%53%
2	Z	19	16%11%11%63%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6734 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 ParA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	211	Total	C	N	O	S	0	0	0
			1578	1003	262	308	5			
1	A	211	Total	C	N	O	S	0	0	0
			1574	1001	262	306	5			
1	B	210	Total	C	N	O	S	0	0	0
			1578	1003	263	307	5			
1	C	209	Total	C	N	O	S	0	0	0
			1565	995	258	307	5			

- Molecule 2 is a protein called TP228 ParB fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Y	9	Total	C	N	O	S	0	0	0
			72	44	14	13	1			
2	Z	7	Total	C	N	O	S	0	0	0
			58	35	11	11	1			
2	K	15	Total	C	N	O	S	0	0	0
			114	69	21	23	1			
2	R	10	Total	C	N	O	S	0	0	0
			71	44	12	14	1			

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).

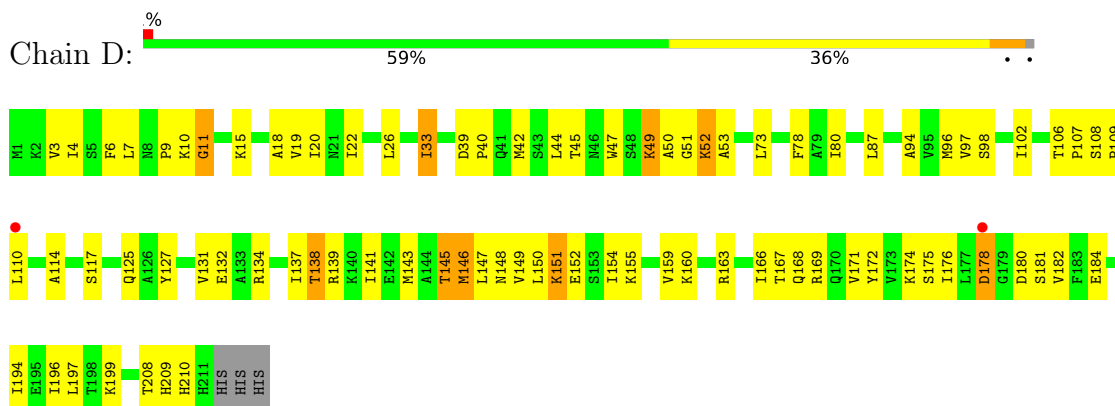


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	A	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	B	1	Total 31	C 10	N 6	O 12	P 3	0	0
3	C	1	Total 31	C 10	N 6	O 12	P 3	0	0

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.

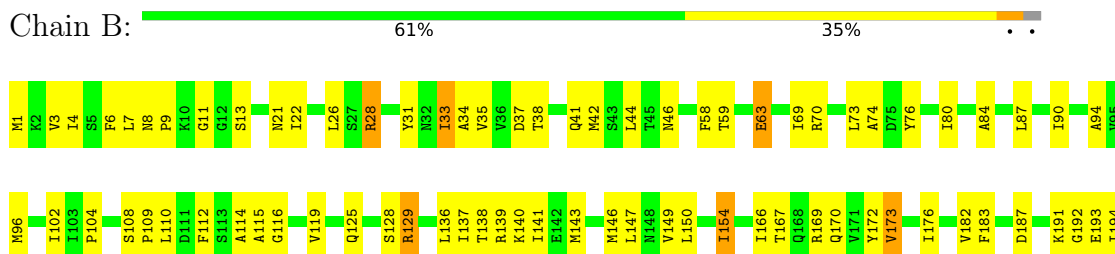
• Molecule 1: ParA

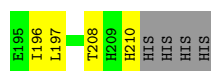


• Molecule 1: ParA



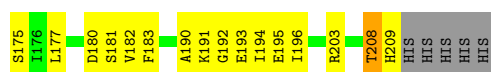
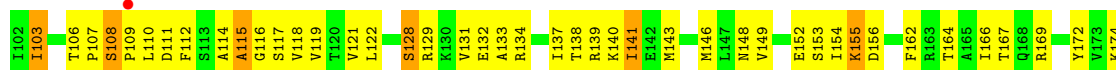
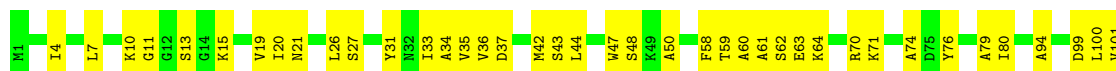
• Molecule 1: ParA





• Molecule 1: ParA

Chain C: 51% 43%



• Molecule 2: TP228 ParB fragment

Chain Y: 21% 16% 11% 53%



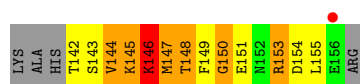
• Molecule 2: TP228 ParB fragment

Chain Z: 16% 11% 11% 63%



• Molecule 2: TP228 ParB fragment

Chain K: 5% 11% 32% 32% 5% 21%



• Molecule 2: TP228 ParB fragment

Chain R: 11% 5% 16% 26% 5% 47%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.36Å 87.51Å 133.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	73.27 – 3.64 73.27 – 3.64	Depositor EDS
% Data completeness (in resolution range)	77.2 (73.27-3.64) 77.3 (73.27-3.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	14.50	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.252 , 0.263 0.245 , 0.262	Depositor DCC
R_{free} test set	1423 reflections (9.94%)	wwPDB-VP
Wilson B-factor (Å ²)	62.0	Xtriage
Anisotropy	0.843	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.124 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6734	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.06 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.6112e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.54	0/1595	0.97	5/2160 (0.2%)
1	B	0.48	0/1600	0.91	5/2166 (0.2%)
1	C	0.47	0/1585	0.91	5/2146 (0.2%)
1	D	0.49	0/1599	0.93	4/2165 (0.2%)
2	K	0.99	0/114	1.81	5/149 (3.4%)
2	R	1.46	1/71 (1.4%)	1.85	3/93 (3.2%)
2	Y	0.90	0/72	1.19	1/93 (1.1%)
2	Z	1.21	0/58	1.37	2/75 (2.7%)
All	All	0.54	1/6694 (0.0%)	0.97	30/9047 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	R	154	GLU	CA-C	7.29	1.62	1.52

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	145	LYS	N-CA-C	9.60	124.31	112.59
2	K	146	LYS	N-CA-C	8.50	123.39	112.34
2	K	144	VAL	N-CA-C	8.15	126.28	109.34
2	R	149	LYS	N-CA-C	-7.31	103.32	111.28
2	R	148	LYS	N-CA-C	-7.22	103.44	111.82
1	A	49	LYS	N-CA-C	7.10	121.20	112.54
1	B	108	SER	CA-C-N	7.07	126.32	118.97
1	B	108	SER	C-N-CA	7.07	126.32	118.97
1	A	150	LEU	N-CA-C	-6.89	103.79	111.71
1	A	39	ASP	CA-C-N	6.52	126.21	119.82
1	A	39	ASP	C-N-CA	6.52	126.21	119.82
2	K	153	ARG	N-CA-C	-6.34	104.42	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	103	ILE	N-CA-C	6.32	114.41	108.15
2	Y	155	ASN	N-CA-C	6.24	118.16	111.36
2	Z	154	GLU	N-CA-C	-6.20	104.44	111.07
1	A	13	SER	N-CA-C	6.09	119.81	112.38
1	C	115	ALA	N-CA-C	-5.83	104.29	111.75
1	B	208	THR	N-CA-C	5.78	118.45	111.40
1	C	108	SER	CA-C-N	5.77	125.25	119.24
1	C	108	SER	C-N-CA	5.77	125.25	119.24
2	K	150	GLY	N-CA-C	-5.69	105.90	112.50
1	D	208	THR	N-CA-C	-5.50	106.74	113.50
1	D	11	GLY	N-CA-C	5.41	126.00	113.18
1	D	209	HIS	N-CA-C	-5.36	106.58	113.23
1	C	99	ASP	N-CA-C	-5.21	106.87	114.12
1	D	51	GLY	N-CA-C	-5.12	103.38	112.54
2	Z	155	ASN	N-CA-C	5.08	117.21	111.11
2	R	152	PHE	N-CA-C	5.08	116.90	111.36
1	B	11	GLY	N-CA-C	5.06	120.28	114.16
1	B	69	ILE	N-CA-C	5.04	115.65	110.36

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	1574	0	1612	105	0
1	B	1578	0	1619	80	0
1	C	1565	0	1612	85	0
1	D	1578	0	1616	94	0
2	K	114	0	108	33	0
2	R	71	0	58	42	0
2	Y	72	0	67	4	0
2	Z	58	0	52	5	0
3	A	31	0	13	0	0
3	B	31	0	13	0	0
3	C	31	0	13	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	31	0	13	3	0
All	All	6734	0	6796	398	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 (398) 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:D:50:ALA:HB1	2:K:145:LYS:HG2	1.19	1.14
2:K:145:LYS:HG3	2:K:148:THR:HG21	1.30	1.11
2:R:149:LYS:HD2	2:R:149:LYS:C	1.77	1.09
1:B:4:ILE:HB	1:B:80:ILE:HD13	1.35	1.07
2:R:151:THR:O	2:R:155:ASN:HB2	1.54	1.06
2:R:149:LYS:HA	2:R:152:PHE:CD1	1.92	1.03
1:C:153:SER:HA	2:K:154:ASP:OD1	1.58	1.03
1:B:150:LEU:O	1:B:154:ILE:HD13	1.60	1.01
1:B:137:ILE:HD12	1:B:150:LEU:HD21	1.44	0.99
2:Z:156:ARG:HG3	2:Z:156:ARG:HH11	1.24	0.98
1:A:22:ILE:HD13	1:A:194:ILE:HG23	1.43	0.98
2:R:146:SER:O	2:R:149:LYS:HG3	1.60	0.98
1:A:112:PHE:O	1:A:112:PHE:CD2	2.17	0.97
1:B:150:LEU:CD1	1:B:154:ILE:HD11	1.99	0.91
2:K:145:LYS:HG3	2:K:148:THR:CG2	2.00	0.91
1:A:149:VAL:HG21	2:R:149:LYS:HB2	1.54	0.88
1:C:149:VAL:HG13	2:K:150:GLY:HA3	1.54	0.87
1:C:20:ILE:HD11	1:C:44:LEU:HG	1.57	0.86
1:A:15:LYS:O	1:A:19:VAL:HG23	1.75	0.86
1:A:137:ILE:H	1:A:137:ILE:HD12	1.40	0.86
1:B:150:LEU:HD12	1:B:154:ILE:HD11	1.57	0.85
1:A:149:VAL:HG11	2:R:149:LYS:HB2	1.59	0.85
2:R:149:LYS:HA	2:R:152:PHE:CE1	2.10	0.84
1:D:137:ILE:HD12	1:D:137:ILE:H	1.42	0.83
1:B:42:MET:HE3	1:B:59:THR:HG21	1.63	0.81
1:D:172:TYR:O	1:D:176:ILE:HD13	1.80	0.79
1:A:42:MET:HE3	1:A:59:THR:HG21	1.65	0.79
1:A:112:PHE:O	1:A:112:PHE:HD2	1.65	0.79
2:R:149:LYS:HA	2:R:152:PHE:HD1	1.44	0.79
1:D:50:ALA:CB	2:K:145:LYS:HG2	2.10	0.79
1:D:178:ASP:O	2:K:145:LYS:CB	2.30	0.79
1:A:22:ILE:HD11	1:A:197:LEU:HD23	1.65	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:150:LEU:HG	1:D:154:ILE:HD11	1.63	0.79
1:C:115:ALA:O	1:C:119:VAL:HG23	1.84	0.78
1:A:149:VAL:HG21	2:R:149:LYS:CB	2.13	0.78
2:R:149:LYS:CA	2:R:152:PHE:HD1	1.96	0.78
1:C:153:SER:CA	2:K:154:ASP:OD1	2.31	0.78
1:A:107:PRO:O	1:A:146:MET:HE3	1.83	0.78
1:C:149:VAL:HG13	2:K:150:GLY:CA	2.13	0.78
1:A:3:VAL:HB	1:A:98:SER:HA	1.66	0.77
1:A:112:PHE:CE2	1:A:115:ALA:HB3	2.18	0.77
1:D:20:ILE:HD11	1:D:44:LEU:HG	1.66	0.77
1:A:4:ILE:HB	1:A:80:ILE:HD13	1.66	0.77
1:A:110:LEU:HD11	1:B:176:ILE:HG21	1.67	0.76
2:R:149:LYS:O	2:R:152:PHE:HB2	1.84	0.76
2:R:149:LYS:HZ1	2:R:150:MET:HG3	1.50	0.76
1:D:174:LYS:O	1:D:178:ASP:OD1	2.03	0.76
2:Z:156:ARG:HG3	2:Z:156:ARG:NH1	2.00	0.75
1:A:149:VAL:CG2	2:R:149:LYS:HB2	2.15	0.74
1:A:49:LYS:HD3	1:A:50:ALA:N	2.02	0.74
1:D:178:ASP:OD2	2:K:142:THR:HG22	1.87	0.74
2:K:146:LYS:C	2:K:147:MET:HE3	2.13	0.74
1:A:20:ILE:HD11	1:A:44:LEU:HG	1.68	0.73
1:B:22:ILE:HD13	1:B:194:ILE:HG23	1.69	0.73
1:B:84:ALA:HB2	1:B:90:ILE:HD12	1.70	0.73
2:R:149:LYS:CA	2:R:152:PHE:CD1	2.70	0.73
1:A:149:VAL:CG1	2:R:149:LYS:HB2	2.18	0.72
2:R:149:LYS:HD2	2:R:149:LYS:O	1.89	0.72
1:D:110:LEU:HD22	1:C:43:SER:HB3	1.71	0.72
1:B:137:ILE:O	1:B:166:ILE:HD12	1.89	0.71
1:D:6:PHE:CZ	1:D:102:ILE:HD12	2.26	0.70
1:D:47:TRP:HA	2:K:149:PHE:HE2	1.56	0.70
1:B:154:ILE:N	1:B:154:ILE:CD1	2.55	0.70
1:C:42:MET:HE3	1:C:59:THR:HG21	1.74	0.70
1:A:20:ILE:HD12	1:A:47:TRP:CZ3	2.26	0.70
1:C:20:ILE:HD12	1:C:47:TRP:CZ3	2.27	0.70
2:R:149:LYS:C	2:R:149:LYS:CD	2.55	0.69
1:B:137:ILE:HD12	1:B:150:LEU:CD2	2.21	0.69
2:R:149:LYS:NZ	2:R:150:MET:HA	2.07	0.69
1:D:26:LEU:HD13	1:D:33:ILE:HD11	1.73	0.69
1:A:143:MET:HG3	1:A:146:MET:HE2	1.74	0.69
1:B:4:ILE:HB	1:B:80:ILE:CD1	2.20	0.69
1:A:7:LEU:HD11	1:A:94:ALA:HB3	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:37:ASP:OD2	1:C:42:MET:HA	1.94	0.68
1:C:21:ASN:HB3	1:C:194:ILE:HD13	1.75	0.67
1:D:15:LYS:HZ2	1:D:15:LYS:HB2	1.59	0.67
2:R:151:THR:HG22	2:R:155:ASN:ND2	2.10	0.67
1:C:63:GLU:CD	1:C:63:GLU:H	2.03	0.67
2:R:149:LYS:C	2:R:152:PHE:HD1	2.03	0.66
1:A:43:SER:HB3	1:B:110:LEU:HD22	1.76	0.66
1:D:3:VAL:C	1:D:4:ILE:HD12	2.20	0.65
1:C:35:VAL:HG22	1:C:80:ILE:HB	1.79	0.65
1:A:207:ALA:HA	1:A:210:HIS:HB2	1.78	0.65
1:D:108:SER:OG	1:D:109:PRO:HD2	1.97	0.65
1:A:31:TYR:HB2	1:A:33:ILE:HD11	1.79	0.64
1:B:38:THR:HB	1:B:90:ILE:HD13	1.79	0.64
1:D:180:ASP:HB3	1:D:184:GLU:HB2	1.79	0.64
1:D:9:PRO:HB3	1:D:87:LEU:HD23	1.78	0.64
1:B:138:THR:OG1	1:B:166:ILE:HD13	1.98	0.63
1:D:6:PHE:CE1	1:D:102:ILE:HD12	2.32	0.63
1:D:178:ASP:HB3	2:K:143:SER:O	1.99	0.63
1:C:172:TYR:HE2	1:C:194:ILE:HD11	1.64	0.62
1:B:84:ALA:CB	1:B:90:ILE:HD12	2.29	0.62
1:D:178:ASP:O	2:K:145:LYS:HB2	1.99	0.62
1:B:112:PHE:CE1	1:B:154:ILE:HD12	2.35	0.62
1:D:49:LYS:C	1:D:49:LYS:HD3	2.26	0.61
1:A:138:THR:HG23	1:A:166:ILE:HB	1.81	0.61
1:B:3:VAL:C	1:B:4:ILE:HD12	2.24	0.61
1:D:11:GLY:HA3	1:C:11:GLY:HA3	1.82	0.61
1:D:174:LYS:NZ	2:K:142:THR:HG21	2.15	0.61
2:R:147:VAL:O	2:R:150:MET:HB2	2.01	0.61
1:A:199:LYS:O	1:A:203:ARG:HG2	2.00	0.61
1:D:3:VAL:HB	1:D:98:SER:HA	1.83	0.61
1:B:22:ILE:HD11	1:B:197:LEU:HD23	1.83	0.61
1:D:20:ILE:HD12	1:D:47:TRP:CZ3	2.35	0.60
1:D:149:VAL:CG1	2:Z:153:GLY:HA3	2.31	0.60
1:C:79:ALA:C	1:C:80:ILE:HD12	2.26	0.60
1:C:44:LEU:O	1:C:44:LEU:HD23	2.01	0.60
1:D:22:ILE:HD13	1:D:194:ILE:HG23	1.82	0.60
1:D:149:VAL:HG13	2:Z:153:GLY:HA3	1.84	0.59
1:A:108:SER:HB3	1:A:111:ASP:HB2	1.84	0.59
2:K:145:LYS:O	2:K:148:THR:HG23	2.03	0.59
1:B:140:LYS:HE3	1:B:147:LEU:HD21	1.84	0.59
1:D:151:LYS:HD3	1:D:152:GLU:HG2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:PHE:CD2	1:A:112:PHE:C	2.78	0.59
2:K:146:LYS:O	2:K:147:MET:HE3	2.01	0.59
1:D:132:GLU:HG3	1:D:134:ARG:HH12	1.68	0.59
1:D:178:ASP:CB	2:K:143:SER:O	2.51	0.59
2:R:149:LYS:NZ	2:R:150:MET:HG3	2.17	0.59
1:C:138:THR:HG23	1:C:166:ILE:HB	1.84	0.59
1:C:172:TYR:CE2	1:C:194:ILE:HD11	2.38	0.59
1:D:47:TRP:HA	2:K:149:PHE:CE2	2.37	0.58
1:D:107:PRO:O	1:D:146:MET:HE3	2.02	0.58
1:C:117:SER:O	1:C:121:VAL:HG23	2.04	0.58
1:A:141:ILE:HD13	1:A:141:ILE:H	1.69	0.58
1:C:31:TYR:HB2	1:C:33:ILE:HD11	1.85	0.58
1:A:49:LYS:HD3	1:A:49:LYS:C	2.29	0.57
2:R:149:LYS:O	2:R:152:PHE:HD1	1.86	0.57
1:C:103:ILE:HD12	1:C:118:VAL:HG11	1.86	0.57
1:C:175:SER:HB2	1:C:181:SER:O	2.04	0.57
1:D:166:ILE:HD11	1:D:197:LEU:HD22	1.85	0.57
1:D:167:THR:HB	1:D:169:ARG:NH1	2.20	0.57
1:C:34:ALA:HB2	1:C:76:TYR:CD1	2.40	0.57
1:C:110:LEU:HD23	1:C:110:LEU:H	1.70	0.57
2:R:149:LYS:HD2	2:R:150:MET:N	2.20	0.57
1:A:140:LYS:HE3	1:A:147:LEU:HD21	1.87	0.57
1:B:44:LEU:HD23	1:B:44:LEU:O	2.05	0.57
1:A:49:LYS:CE	1:A:50:ALA:HB2	2.35	0.56
1:C:4:ILE:HG12	1:C:100:LEU:HD23	1.87	0.56
1:A:107:PRO:HD3	1:A:138:THR:O	2.06	0.56
1:B:150:LEU:HD11	1:B:154:ILE:HD11	1.84	0.56
1:D:174:LYS:HZ3	2:K:142:THR:HG21	1.71	0.56
1:D:178:ASP:OD2	2:K:142:THR:CG2	2.54	0.56
1:D:196:ILE:HA	1:D:199:LYS:HD2	1.88	0.56
1:D:50:ALA:HB1	2:K:145:LYS:CG	2.13	0.56
1:D:143:MET:CE	1:C:174:LYS:HD3	2.36	0.56
1:D:49:LYS:O	1:D:49:LYS:HG2	2.06	0.56
1:B:87:LEU:HD12	1:B:87:LEU:N	2.21	0.56
2:K:145:LYS:HB3	2:K:149:PHE:HE1	1.70	0.56
2:R:149:LYS:HE3	2:R:150:MET:HG3	1.87	0.55
1:C:21:ASN:OD1	1:C:182:VAL:HG13	2.06	0.55
2:R:146:SER:O	2:R:149:LYS:CG	2.45	0.55
1:B:183:PHE:HA	1:B:191:LYS:HD3	1.89	0.55
1:B:146:MET:O	1:B:149:VAL:HB	2.06	0.55
1:C:112:PHE:CE2	1:C:153:SER:HB3	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:ASP:CG	2:K:154:ASP:HB3	2.32	0.55
1:D:44:LEU:HD23	1:D:44:LEU:O	2.06	0.55
1:D:172:TYR:CD2	1:D:182:VAL:HG11	2.42	0.55
1:D:106:THR:HG22	1:D:138:THR:HG23	1.90	0.54
1:D:175:SER:HB2	1:D:181:SER:O	2.06	0.54
2:R:149:LYS:HZ2	2:R:150:MET:HA	1.73	0.54
1:B:7:LEU:HD11	1:B:94:ALA:HB3	1.89	0.54
1:A:107:PRO:HG2	1:A:140:LYS:HA	1.89	0.54
1:B:26:LEU:HD13	1:B:33:ILE:HD11	1.89	0.54
1:C:155:LYS:HB2	1:C:155:LYS:NZ	2.23	0.53
1:B:138:THR:HA	1:B:166:ILE:HB	1.90	0.53
2:R:149:LYS:CE	2:R:150:MET:HG3	2.38	0.53
1:B:34:ALA:HB2	1:B:76:TYR:CE1	2.43	0.53
1:B:150:LEU:HD12	1:B:154:ILE:CD1	2.32	0.53
1:A:149:VAL:HG11	2:R:149:LYS:CB	2.36	0.53
1:B:63:GLU:CD	1:B:63:GLU:H	2.16	0.53
1:B:154:ILE:HD13	1:B:154:ILE:N	2.21	0.53
2:K:147:MET:O	2:K:151:GLU:HG3	2.09	0.53
2:R:151:THR:O	2:R:155:ASN:CB	2.42	0.53
1:B:138:THR:HG23	1:B:166:ILE:HB	1.90	0.53
1:D:145:THR:HG21	1:C:174:LYS:NZ	2.24	0.53
1:A:166:ILE:HD12	1:A:194:ILE:CD1	2.39	0.53
2:K:147:MET:HE2	2:K:147:MET:HA	1.92	0.52
1:A:33:ILE:HD12	1:A:33:ILE:N	2.25	0.52
1:C:48:SER:C	1:C:50:ALA:H	2.16	0.52
1:D:166:ILE:HD11	1:D:197:LEU:CD2	2.40	0.51
1:C:70:ARG:O	1:C:74:ALA:HB2	2.09	0.51
1:C:109:PRO:HD3	1:C:146:MET:HE1	1.91	0.51
1:A:169:ARG:CZ	1:A:189:ALA:HB1	2.39	0.51
1:C:10:LYS:HG2	1:C:111:ASP:OD1	2.11	0.51
1:C:137:ILE:HD11	1:C:154:ILE:HD11	1.91	0.51
1:A:20:ILE:HG23	1:A:47:TRP:CH2	2.46	0.51
2:Y:149:LYS:O	2:Y:152:PHE:HB2	2.11	0.51
1:D:148:ASN:HA	1:D:151:LYS:HD2	1.93	0.51
1:A:100:LEU:HG	1:A:102:ILE:CD1	2.40	0.51
2:R:149:LYS:O	2:R:152:PHE:CD1	2.63	0.51
1:A:44:LEU:HD23	1:A:44:LEU:O	2.11	0.51
1:A:103:ILE:N	1:A:103:ILE:HD12	2.25	0.51
1:D:39:ASP:OD2	1:D:40:PRO:HD2	2.11	0.50
1:D:22:ILE:CD1	1:D:197:LEU:HD23	2.41	0.50
1:C:208:THR:O	1:C:209:HIS:C	2.54	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:175:SER:OG	1:D:176:ILE:HD12	2.11	0.50
1:A:125:GLN:HE21	1:A:129:ARG:HB3	1.77	0.50
1:C:148:ASN:O	1:C:152:GLU:HG3	2.12	0.50
1:C:7:LEU:HD11	1:C:94:ALA:HB3	1.93	0.50
1:C:143:MET:HB3	1:C:146:MET:HB3	1.92	0.50
1:C:34:ALA:HB2	1:C:76:TYR:CE1	2.46	0.49
1:B:13:SER:HB2	1:B:104:PRO:O	2.12	0.49
1:B:128:SER:HB3	1:C:129:ARG:HB2	1.94	0.49
2:K:146:LYS:HB2	2:K:147:MET:HE3	1.93	0.49
1:D:52:LYS:HG3	1:D:53:ALA:O	2.12	0.49
1:A:100:LEU:HG	1:A:102:ILE:HD11	1.94	0.49
1:A:42:MET:HB3	1:A:45:THR:HB	1.95	0.49
1:A:171:VAL:HG22	1:A:182:VAL:HG12	1.94	0.49
2:Z:156:ARG:NH1	2:Z:156:ARG:CG	2.66	0.49
1:B:187:ASP:O	1:B:191:LYS:HG3	2.13	0.49
1:C:137:ILE:HD11	1:C:154:ILE:CD1	2.43	0.49
1:B:26:LEU:O	1:B:31:TYR:HB2	2.13	0.48
1:B:70:ARG:O	1:B:74:ALA:HB2	2.12	0.48
1:D:10:LYS:NZ	3:C:301:ANP:HNB1	2.10	0.48
1:A:191:LYS:O	1:A:195:GLU:HB2	2.13	0.48
1:C:172:TYR:CD2	1:C:182:VAL:HG11	2.48	0.48
1:B:87:LEU:HD22	1:B:90:ILE:HD11	1.95	0.48
1:D:178:ASP:O	2:K:145:LYS:HB3	2.09	0.48
1:A:36:VAL:HG21	1:A:69:ILE:HD12	1.96	0.48
1:C:26:LEU:O	1:C:31:TYR:HB2	2.13	0.48
1:D:22:ILE:HD11	1:D:197:LEU:HD23	1.95	0.48
1:C:33:ILE:HD12	1:C:33:ILE:N	2.28	0.48
1:A:102:ILE:HD12	1:A:102:ILE:N	2.29	0.48
1:C:20:ILE:HG23	1:C:47:TRP:CH2	2.48	0.48
1:D:11:GLY:O	3:D:301:ANP:N3B	2.47	0.47
1:A:174:LYS:HD3	1:B:143:MET:HE3	1.96	0.47
1:B:169:ARG:HB2	1:B:172:TYR:CD1	2.49	0.47
1:C:192:GLY:O	1:C:196:ILE:HG12	2.13	0.47
1:A:1:MET:HE1	1:A:70:ARG:HA	1.97	0.47
1:C:190:ALA:O	1:C:193:GLU:HG2	2.15	0.47
2:R:149:LYS:CD	2:R:150:MET:N	2.76	0.47
2:R:151:THR:O	2:R:155:ASN:ND2	2.44	0.47
1:A:149:VAL:CG2	2:R:149:LYS:CB	2.85	0.47
1:A:150:LEU:O	1:A:154:ILE:HG13	2.14	0.47
1:C:31:TYR:CB	1:C:33:ILE:HD11	2.45	0.47
2:R:149:LYS:O	2:R:152:PHE:CB	2.59	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:VAL:HG13	2:R:149:LYS:HD3	1.95	0.47
1:D:150:LEU:O	1:D:154:ILE:HG13	2.15	0.47
2:K:146:LYS:O	2:K:147:MET:CE	2.62	0.47
1:D:33:ILE:HD13	1:D:78:PHE:HB2	1.97	0.46
1:C:59:THR:HG22	1:C:60:ALA:N	2.30	0.46
2:R:150:MET:C	2:R:152:PHE:N	2.72	0.46
1:A:114:ALA:O	1:A:115:ALA:C	2.58	0.46
3:D:301:ANP:O3'	1:C:108:SER:HB2	2.16	0.46
1:D:178:ASP:O	2:K:145:LYS:CA	2.63	0.46
1:B:136:LEU:CG	1:B:166:ILE:HD11	2.46	0.46
1:A:87:LEU:N	1:A:87:LEU:CD1	2.79	0.46
1:D:137:ILE:H	1:D:137:ILE:CD1	2.18	0.46
1:A:98:SER:N	1:A:129:ARG:HH22	2.14	0.46
1:A:20:ILE:HD11	1:A:44:LEU:CG	2.39	0.46
1:C:80:ILE:HD12	1:C:80:ILE:N	2.31	0.46
1:D:139:ARG:HH12	1:C:139:ARG:HH12	1.64	0.45
1:A:148:ASN:O	1:A:152:GLU:HG3	2.15	0.45
1:A:117:SER:O	1:A:121:VAL:HG23	2.15	0.45
2:R:147:VAL:O	2:R:150:MET:CB	2.65	0.45
1:A:100:LEU:HD12	1:A:101:VAL:N	2.30	0.45
1:A:102:ILE:O	1:A:104:PRO:HD3	2.15	0.45
1:C:101:VAL:HG12	1:C:103:ILE:HD11	1.99	0.45
1:C:110:LEU:H	1:C:110:LEU:CD2	2.29	0.45
1:D:33:ILE:CD1	1:D:78:PHE:HB2	2.46	0.45
1:D:44:LEU:O	1:D:47:TRP:HB3	2.17	0.45
1:D:49:LYS:O	1:D:49:LYS:CG	2.64	0.45
1:D:44:LEU:HD23	1:D:44:LEU:C	2.41	0.45
1:D:171:VAL:HG13	1:D:172:TYR:N	2.31	0.45
1:D:42:MET:CE	1:D:45:THR:HG21	2.46	0.45
1:D:147:LEU:O	1:D:151:LYS:HB3	2.17	0.45
1:A:177:LEU:HG	2:Y:149:LYS:HB3	1.97	0.45
1:B:34:ALA:HB2	1:B:76:TYR:CD1	2.52	0.45
1:B:166:ILE:HD12	1:B:166:ILE:H	1.82	0.45
1:B:84:ALA:HB2	1:B:90:ILE:CD1	2.43	0.44
1:D:145:THR:O	1:D:149:VAL:HG23	2.17	0.44
1:A:143:MET:O	1:A:147:LEU:HB2	2.17	0.44
1:C:4:ILE:O	1:C:80:ILE:HA	2.16	0.44
1:A:116:GLY:O	1:A:119:VAL:HB	2.16	0.44
1:B:4:ILE:O	1:B:80:ILE:HA	2.18	0.44
1:C:112:PHE:CZ	1:C:153:SER:HB3	2.53	0.44
1:B:4:ILE:HD12	1:B:4:ILE:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:MET:HE3	1:C:174:LYS:HD3	2.00	0.44
1:A:141:ILE:HG12	1:A:143:MET:HG2	1.99	0.44
1:D:174:LYS:HE2	2:K:146:LYS:NZ	2.33	0.44
1:B:42:MET:O	1:B:46:ASN:HB2	2.18	0.44
1:C:36:VAL:HA	1:C:58:PHE:O	2.17	0.44
1:C:64:LYS:HB3	1:C:64:LYS:HE2	1.83	0.44
1:D:146:MET:HG3	1:C:177:LEU:HD22	2.00	0.44
1:A:4:ILE:HG23	1:A:102:ILE:CD1	2.47	0.44
1:A:7:LEU:HD22	1:A:91:THR:HA	1.99	0.44
1:A:19:VAL:HG13	1:A:80:ILE:HG21	1.99	0.44
1:A:159:VAL:HG12	1:A:160:LYS:N	2.33	0.44
1:B:35:VAL:HA	1:B:80:ILE:O	2.18	0.44
1:B:154:ILE:N	1:B:154:ILE:HD12	2.32	0.44
1:A:134:ARG:HD3	1:A:160:LYS:HB2	1.99	0.43
1:A:149:VAL:CB	2:R:149:LYS:HB2	2.47	0.43
1:A:167:THR:HG22	1:A:168:GLN:N	2.33	0.43
1:D:7:LEU:HD11	1:D:94:ALA:CB	2.48	0.43
1:C:183:PHE:HA	1:C:191:LYS:HD3	2.00	0.43
1:A:7:LEU:HD11	1:A:94:ALA:CB	2.46	0.43
1:A:49:LYS:HE2	1:A:50:ALA:HB2	1.99	0.43
1:B:1:MET:HG2	1:B:76:TYR:O	2.17	0.43
1:B:138:THR:HG22	1:B:139:ARG:N	2.34	0.43
1:C:20:ILE:CD1	1:C:44:LEU:HG	2.36	0.43
1:A:4:ILE:HB	1:A:80:ILE:CD1	2.41	0.43
1:C:15:LYS:O	1:C:19:VAL:HG23	2.18	0.43
1:A:32:ASN:C	1:A:33:ILE:HD12	2.44	0.43
1:B:102:ILE:O	1:B:104:PRO:HD3	2.19	0.43
1:C:71:LYS:O	1:C:71:LYS:HD3	2.19	0.43
1:D:49:LYS:C	1:D:49:LYS:CD	2.88	0.43
1:D:96:MET:HA	1:D:125:GLN:OE1	2.18	0.43
1:B:167:THR:HG22	1:B:169:ARG:HG2	2.00	0.43
2:R:150:MET:C	2:R:152:PHE:H	2.27	0.43
1:D:150:LEU:HG	1:D:154:ILE:CD1	2.41	0.43
1:D:172:TYR:CE2	1:D:182:VAL:HG11	2.54	0.43
1:B:8:ASN:ND2	1:B:104:PRO:O	2.52	0.43
1:B:109:PRO:HD3	1:B:146:MET:SD	2.59	0.43
1:A:140:LYS:O	1:A:167:THR:HG23	2.18	0.43
1:A:147:LEU:O	1:A:151:LYS:HG2	2.19	0.43
1:D:178:ASP:HB2	2:K:143:SER:O	2.18	0.43
1:A:18:ALA:O	1:A:22:ILE:HG12	2.19	0.43
1:A:112:PHE:O	1:A:113:SER:C	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:109:PRO:HD3	1:C:146:MET:CE	2.49	0.43
1:A:195:GLU:O	1:A:199:LYS:HG3	2.19	0.42
1:B:116:GLY:O	1:B:119:VAL:HB	2.18	0.42
1:C:122:LEU:CD2	1:C:131:VAL:HB	2.50	0.42
1:A:52:LYS:HD3	1:A:53:ALA:O	2.19	0.42
1:C:114:ALA:C	1:C:116:GLY:N	2.76	0.42
1:D:97:VAL:HG23	1:D:97:VAL:O	2.18	0.42
1:D:155:LYS:HA	1:D:155:LYS:HD2	1.93	0.42
1:A:146:MET:HE1	1:B:173:VAL:HG11	2.01	0.42
1:C:140:LYS:HG2	1:C:141:ILE:N	2.33	0.42
1:D:19:VAL:HG13	1:D:80:ILE:HG21	2.00	0.42
1:A:3:VAL:HG21	1:A:97:VAL:O	2.19	0.42
1:A:64:LYS:O	1:A:64:LYS:HG2	2.18	0.42
1:B:28:ARG:NH1	1:B:183:PHE:HB3	2.34	0.42
1:A:136:LEU:HD13	1:A:162:PHE:HD2	1.84	0.42
1:A:143:MET:HE3	1:B:170:GLN:NE2	2.34	0.42
1:B:96:MET:HA	1:B:125:GLN:OE1	2.20	0.42
1:C:122:LEU:HD13	1:C:133:ALA:HB2	2.02	0.42
1:A:7:LEU:O	1:A:8:ASN:HB2	2.20	0.42
1:A:100:LEU:CD2	1:A:102:ILE:HD11	2.50	0.42
1:B:21:ASN:OD1	1:B:182:VAL:HG13	2.20	0.42
1:C:61:ALA:O	1:C:62:SER:HB3	2.19	0.42
1:D:168:GLN:HA	3:D:301:ANP:N1	2.35	0.42
1:A:49:LYS:C	1:A:49:LYS:CD	2.92	0.42
2:K:146:LYS:C	2:K:147:MET:CE	2.90	0.42
1:D:137:ILE:HD12	1:D:137:ILE:N	2.23	0.42
1:D:159:VAL:HG12	1:D:160:LYS:N	2.35	0.42
1:D:167:THR:HG22	1:D:168:GLN:N	2.35	0.42
1:B:128:SER:HB3	1:C:128:SER:HB2	2.01	0.42
1:D:141:ILE:HD13	1:D:143:MET:SD	2.60	0.41
1:A:163:ARG:HB2	1:A:163:ARG:NH1	2.35	0.41
1:A:180:ASP:OD2	1:A:180:ASP:N	2.51	0.41
1:A:207:ALA:CA	1:A:210:HIS:HB2	2.46	0.41
1:B:6:PHE:CZ	1:B:102:ILE:HD12	2.55	0.41
1:B:136:LEU:HG	1:B:166:ILE:HD11	2.01	0.41
2:Y:149:LYS:H	2:Y:149:LYS:HD2	1.85	0.41
1:D:154:ILE:O	1:D:154:ILE:HG22	2.21	0.41
1:A:143:MET:HE2	1:A:143:MET:HA	2.02	0.41
1:B:115:ALA:O	1:B:119:VAL:HG23	2.19	0.41
1:B:125:GLN:HG3	1:B:129:ARG:O	2.20	0.41
1:A:49:LYS:HD3	1:A:50:ALA:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:LEU:HD13	1:B:33:ILE:CD1	2.50	0.41
1:B:192:GLY:O	1:B:196:ILE:HG13	2.19	0.41
2:R:149:LYS:HE3	2:R:150:MET:CG	2.50	0.41
1:A:1:MET:HB2	1:A:74:ALA:O	2.20	0.41
1:B:169:ARG:HG3	1:B:169:ARG:HH11	1.86	0.41
1:C:167:THR:HB	1:C:169:ARG:NH1	2.34	0.41
1:A:87:LEU:N	1:A:87:LEU:HD12	2.35	0.41
1:D:172:TYR:OH	1:D:194:ILE:HD11	2.21	0.41
1:A:8:ASN:O	1:A:15:LYS:HD3	2.20	0.41
1:B:37:ASP:OD2	1:B:42:MET:HA	2.20	0.41
1:B:154:ILE:HD13	1:B:154:ILE:H	1.85	0.41
2:Y:149:LYS:HD2	2:Y:149:LYS:N	2.36	0.41
1:D:18:ALA:O	1:D:22:ILE:HG12	2.21	0.41
1:C:138:THR:HG22	1:C:139:ARG:N	2.36	0.41
1:C:155:LYS:HB2	1:C:155:LYS:HZ2	1.84	0.41
1:D:114:ALA:O	1:D:117:SER:HB3	2.21	0.41
1:B:4:ILE:CB	1:B:80:ILE:HD13	2.26	0.41
1:C:4:ILE:HB	1:C:80:ILE:HG13	2.02	0.41
1:C:107:PRO:O	1:C:108:SER:C	2.64	0.41
1:A:66:VAL:O	1:A:69:ILE:HG12	2.20	0.41
1:A:91:THR:CG2	1:A:118:VAL:HG22	2.51	0.41
1:A:149:VAL:HG22	2:R:149:LYS:HG2	2.03	0.41
1:B:9:PRO:HD2	1:B:114:ALA:HB1	2.03	0.40
1:B:87:LEU:HD12	1:B:87:LEU:H	1.85	0.40
1:C:162:PHE:C	1:C:164:THR:H	2.29	0.40
1:D:125:GLN:HG2	1:D:131:VAL:HG21	2.04	0.40
1:B:9:PRO:HB2	1:B:114:ALA:HB1	2.03	0.40
1:C:13:SER:HA	1:C:106:THR:CG2	2.51	0.40
1:D:160:LYS:HE3	1:D:160:LYS:H	1.87	0.40
1:B:38:THR:CB	1:B:90:ILE:HD13	2.50	0.40
1:B:41:GLN:O	1:B:42:MET:HB2	2.21	0.40
1:B:58:PHE:CE1	1:B:73:LEU:HD22	2.56	0.40
1:C:109:PRO:HD3	1:C:146:MET:SD	2.61	0.40
1:C:132:GLU:HG3	1:C:134:ARG:HH12	1.86	0.40
1:D:127:TYR:HE2	1:A:129:ARG:NH1	2.20	0.40
1:A:49:LYS:HD3	1:A:50:ALA:CB	2.52	0.40
1:C:33:ILE:HG23	1:C:80:ILE:HD13	2.04	0.40
1:C:48:SER:C	1:C:50:ALA:N	2.79	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	209/214 (98%)	190 (91%)	19 (9%)	0	100	100
1	B	208/214 (97%)	180 (86%)	28 (14%)	0	100	100
1	C	207/214 (97%)	184 (89%)	23 (11%)	0	100	100
1	D	209/214 (98%)	181 (87%)	28 (13%)	0	100	100
2	K	13/19 (68%)	10 (77%)	2 (15%)	1 (8%)	1	7
2	R	8/19 (42%)	6 (75%)	2 (25%)	0	100	100
2	Y	7/19 (37%)	7 (100%)	0	0	100	100
2	Z	5/19 (26%)	5 (100%)	0	0	100	100
All	All	866/932 (93%)	763 (88%)	102 (12%)	1 (0%)	48	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	K	144	VAL

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	169/176 (96%)	152 (90%)	17 (10%)	7	27
1	B	171/176 (97%)	162 (95%)	9 (5%)	20	44
1	C	170/176 (97%)	162 (95%)	8 (5%)	23	47
1	D	170/176 (97%)	159 (94%)	11 (6%)	15	40

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	K	12/17 (71%)	7 (58%)	5 (42%)	0	0
2	R	6/17 (35%)	1 (17%)	5 (83%)	0	0
2	Y	7/17 (41%)	3 (43%)	4 (57%)	0	0
2	Z	6/17 (35%)	4 (67%)	2 (33%)	0	1
All	All	711/772 (92%)	650 (91%)	61 (9%)	10	32

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	33	ILE
1	D	49	LYS
1	D	52	LYS
1	D	73	LEU
1	D	138	THR
1	D	145	THR
1	D	146	MET
1	D	151	LYS
1	D	163	ARG
1	D	178	ASP
1	D	210	HIS
1	A	10	LYS
1	A	49	LYS
1	A	70	ARG
1	A	71	LYS
1	A	73	LEU
1	A	86	SER
1	A	87	LEU
1	A	88	SER
1	A	106	THR
1	A	122	LEU
1	A	137	ILE
1	A	141	ILE
1	A	145	THR
1	A	146	MET
1	A	151	LYS
1	A	176	ILE
1	A	195	GLU
1	B	28	ARG
1	B	33	ILE
1	B	63	GLU
1	B	129	ARG

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Mol	Chain	Res	Type
1	B	141	ILE
1	B	154	ILE
1	B	173	VAL
1	B	193	GLU
1	B	210	HIS
1	C	27	SER
1	C	128	SER
1	C	141	ILE
1	C	155	LYS
1	C	180	ASP
1	C	195	GLU
1	C	203	ARG
1	C	208	THR
2	Y	149	LYS
2	Y	150	MET
2	Y	151	THR
2	Y	155	ASN
2	Z	154	GLU
2	Z	156	ARG
2	K	146	LYS
2	K	147	MET
2	K	148	THR
2	K	153	ARG
2	K	155	LEU
2	R	149	LYS
2	R	150	MET
2	R	151	THR
2	R	154	GLU
2	R	155	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	41	GLN
1	D	46	ASN
1	D	148	ASN
1	D	168	GLN
1	A	8	ASN
1	A	125	GLN
1	B	46	ASN
1	B	168	GLN
1	B	170	GLN

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Mol	Chain	Res	Type
1	C	168	GLN
1	C	170	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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
3	ANP	C	301	-	33,33,33	4.12	10 (30%)	45,52,52	2.22	9 (20%)
3	ANP	A	500	-	33,33,33	3.80	8 (24%)	45,52,52	2.36	8 (17%)
3	ANP	B	501	-	33,33,33	4.19	9 (27%)	45,52,52	2.25	8 (17%)
3	ANP	D	301	-	33,33,33	3.85	10 (30%)	45,52,52	2.13	8 (17%)

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
3	ANP	C	301	-	-	5/18/38/38	0/3/3/3
3	ANP	A	500	-	-	3/18/38/38	0/3/3/3
3	ANP	B	501	-	-	5/18/38/38	0/3/3/3
3	ANP	D	301	-	-	4/18/38/38	0/3/3/3

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	ANP	PB-O3A	21.01	1.85	1.59
3	C	301	ANP	PB-O3A	20.38	1.84	1.59
3	A	500	ANP	PB-O3A	20.13	1.84	1.59
3	D	301	ANP	PB-O3A	19.59	1.83	1.59
3	C	301	ANP	PG-O1G	6.44	1.55	1.46
3	B	501	ANP	PG-O1G	5.86	1.55	1.46
3	C	301	ANP	C4-N3	4.49	1.42	1.34
3	A	500	ANP	C4-N3	4.01	1.41	1.34
3	B	501	ANP	C4-N3	3.88	1.41	1.34
3	D	301	ANP	C4-N3	3.85	1.41	1.34
3	D	301	ANP	PB-O2B	-3.69	1.47	1.56
3	D	301	ANP	PB-O1B	3.56	1.51	1.46
3	B	501	ANP	PA-O3A	3.55	1.63	1.59
3	D	301	ANP	PA-O3A	3.46	1.63	1.59
3	B	501	ANP	PG-O2G	-3.45	1.47	1.56
3	B	501	ANP	PB-N3B	-3.42	1.54	1.63
3	C	301	ANP	PG-O2G	-3.28	1.48	1.56
3	C	301	ANP	PA-O3A	3.00	1.62	1.59
3	C	301	ANP	C5-C6	2.99	1.49	1.41
3	D	301	ANP	C2'-C3'	-2.98	1.45	1.53
3	C	301	ANP	PB-N3B	-2.92	1.55	1.63
3	B	501	ANP	C5-C6	2.91	1.49	1.41
3	D	301	ANP	C5'-C4'	-2.83	1.43	1.51
3	C	301	ANP	C5'-C4'	-2.75	1.43	1.51
3	D	301	ANP	C5-C6	2.68	1.48	1.41
3	A	500	ANP	PB-O2B	-2.66	1.49	1.56
3	A	500	ANP	C5-C6	2.62	1.48	1.41
3	D	301	ANP	PB-N3B	-2.58	1.56	1.63
3	B	501	ANP	C2-N3	2.50	1.38	1.33
3	C	301	ANP	C2-N3	2.47	1.38	1.33
3	A	500	ANP	C4-N9	2.40	1.42	1.37
3	A	500	ANP	C2'-C3'	-2.40	1.46	1.53
3	B	501	ANP	C5'-C4'	-2.39	1.44	1.51
3	D	301	ANP	C4-N9	2.36	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	301	ANP	C4-N9	2.33	1.42	1.37
3	A	500	ANP	PB-N3B	-2.10	1.57	1.63
3	A	500	ANP	C5'-C4'	-2.03	1.45	1.51

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	ANP	O2B-PB-O1B	8.33	127.74	109.87
3	C	301	ANP	O2B-PB-O1B	7.95	126.93	109.87
3	D	301	ANP	O2B-PB-O1B	7.84	126.69	109.87
3	B	501	ANP	O2B-PB-O1B	7.83	126.67	109.87
3	D	301	ANP	O3A-PB-N3B	-6.88	87.50	106.59
3	A	500	ANP	O3A-PB-N3B	-6.80	87.72	106.59
3	C	301	ANP	O3A-PB-N3B	-6.64	88.17	106.59
3	A	500	ANP	O1G-PG-N3B	-6.64	102.00	111.77
3	B	501	ANP	O3A-PB-N3B	-6.25	89.25	106.59
3	B	501	ANP	O2B-PB-O3A	-5.16	87.41	104.64
3	D	301	ANP	O1G-PG-N3B	-5.10	104.26	111.77
3	C	301	ANP	O2B-PB-O3A	-4.84	88.50	104.64
3	A	500	ANP	O2B-PB-O3A	-4.67	89.06	104.64
3	B	501	ANP	O1B-PB-N3B	4.66	118.63	111.77
3	C	301	ANP	O1B-PB-N3B	4.27	118.06	111.77
3	D	301	ANP	C3'-C2'-C1'	3.93	108.90	101.46
3	A	500	ANP	O1B-PB-N3B	3.90	117.51	111.77
3	B	501	ANP	O1G-PG-N3B	-3.54	106.55	111.77
3	B	501	ANP	O2G-PG-O3G	3.53	117.07	107.59
3	C	301	ANP	O1G-PG-N3B	-3.45	106.69	111.77
3	A	500	ANP	C4-N9-C8	-3.39	102.18	105.74
3	B	501	ANP	C4-N9-C8	-3.38	102.19	105.74
3	C	301	ANP	O2G-PG-O3G	3.28	116.40	107.59
3	C	301	ANP	C4-N9-C8	-3.26	102.31	105.74
3	D	301	ANP	C4-N9-C8	-3.18	102.40	105.74
3	A	500	ANP	C3'-C2'-C1'	2.71	106.59	101.46
3	D	301	ANP	O2B-PB-O3A	-2.68	95.69	104.64
3	B	501	ANP	C3'-C2'-C1'	2.54	106.26	101.46
3	C	301	ANP	C3'-C2'-C1'	2.39	105.98	101.46
3	D	301	ANP	C2'-C3'-C4'	2.33	107.11	102.61
3	A	500	ANP	C2'-C3'-C4'	2.14	106.74	102.61
3	D	301	ANP	N9-C8-N7	2.03	116.82	113.94
3	C	301	ANP	N9-C8-N7	2.00	116.78	113.94

There are no chirality outliers.

All (17) torsion outliers are listed below:

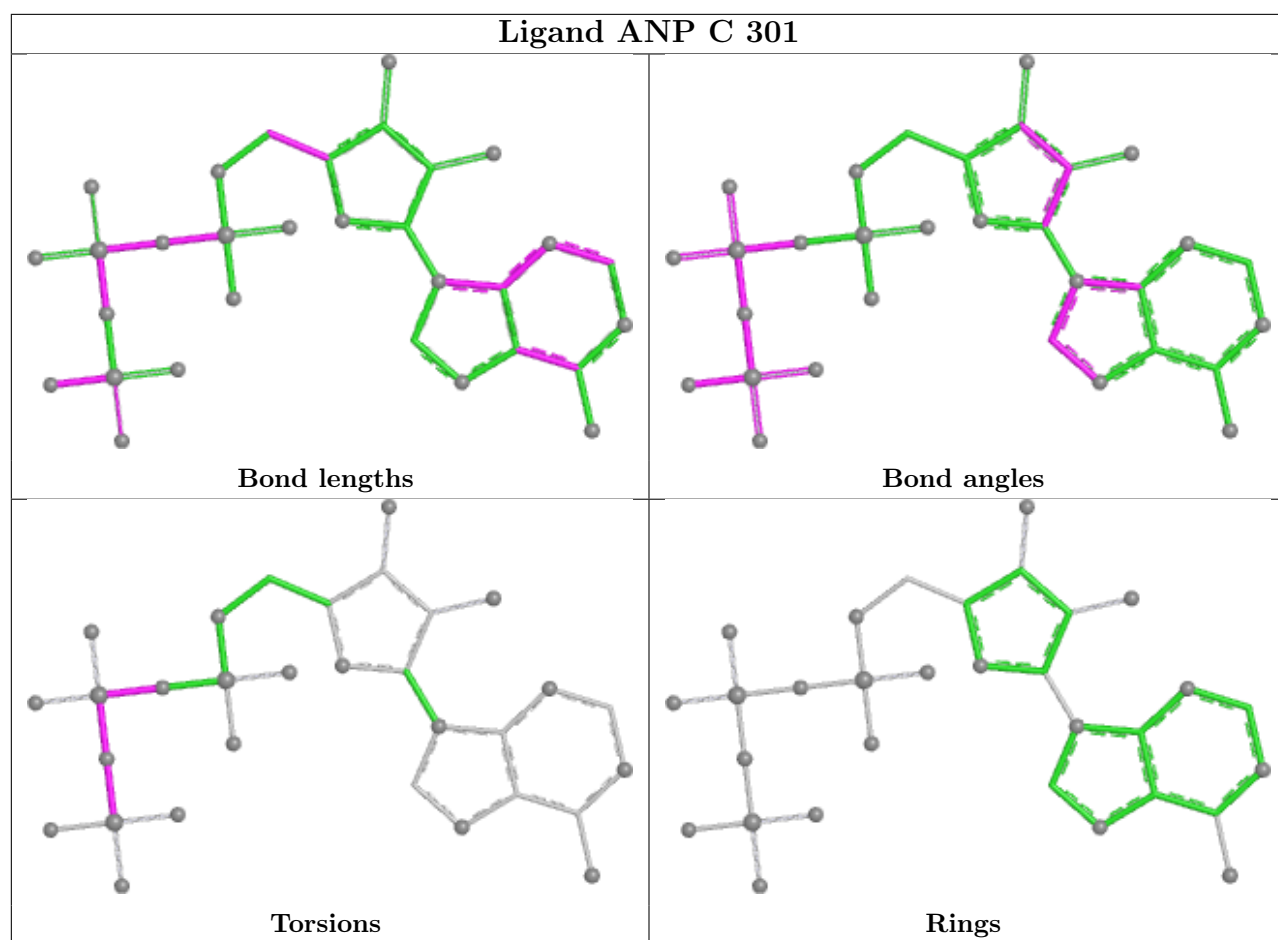
Mol	Chain	Res	Type	Atoms
3	D	301	ANP	PG-N3B-PB-O1B
3	D	301	ANP	PG-N3B-PB-O3A
3	A	500	ANP	PB-N3B-PG-O1G
3	A	500	ANP	PG-N3B-PB-O1B
3	A	500	ANP	PG-N3B-PB-O3A
3	B	501	ANP	PB-N3B-PG-O1G
3	B	501	ANP	PG-N3B-PB-O1B
3	B	501	ANP	PG-N3B-PB-O3A
3	C	301	ANP	PB-N3B-PG-O1G
3	C	301	ANP	PG-N3B-PB-O1B
3	C	301	ANP	PG-N3B-PB-O3A
3	C	301	ANP	PA-O3A-PB-O2B
3	D	301	ANP	O4'-C4'-C5'-O5'
3	B	501	ANP	PA-O3A-PB-O2B
3	D	301	ANP	C3'-C4'-C5'-O5'
3	B	501	ANP	PA-O3A-PB-O1B
3	C	301	ANP	PA-O3A-PB-O1B

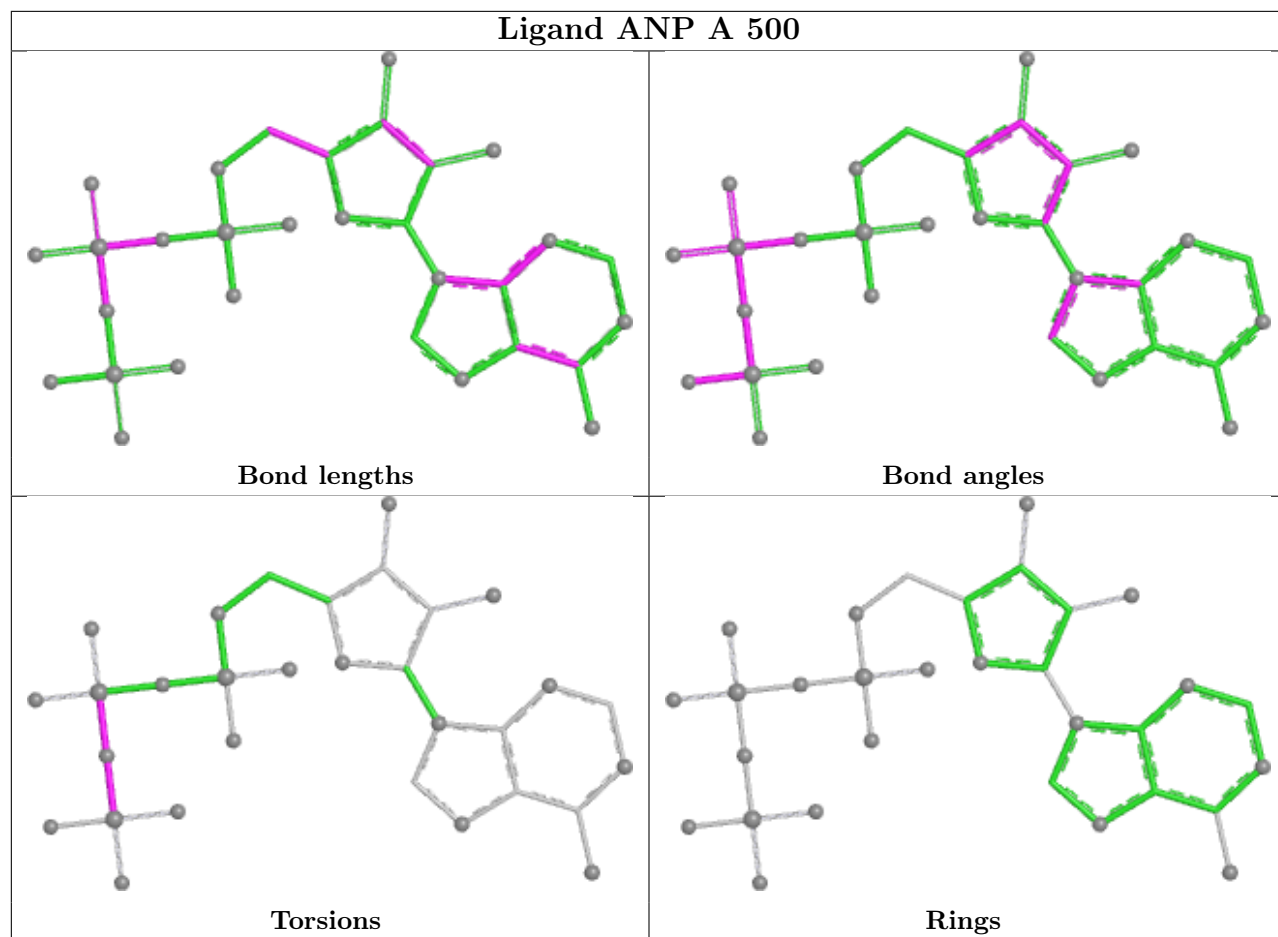
There are no ring outliers.

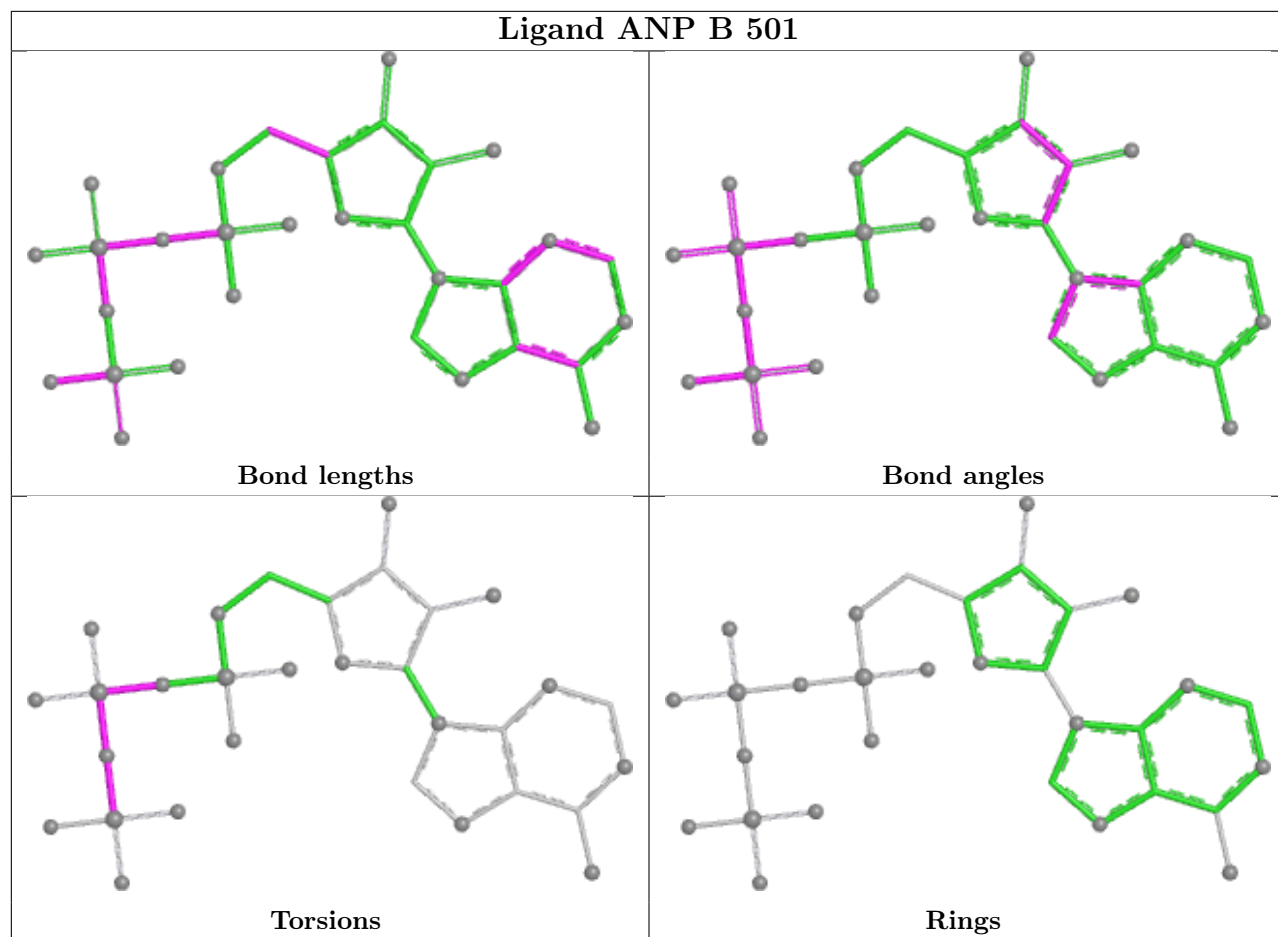
2 monomers are involved in 4 short contacts:

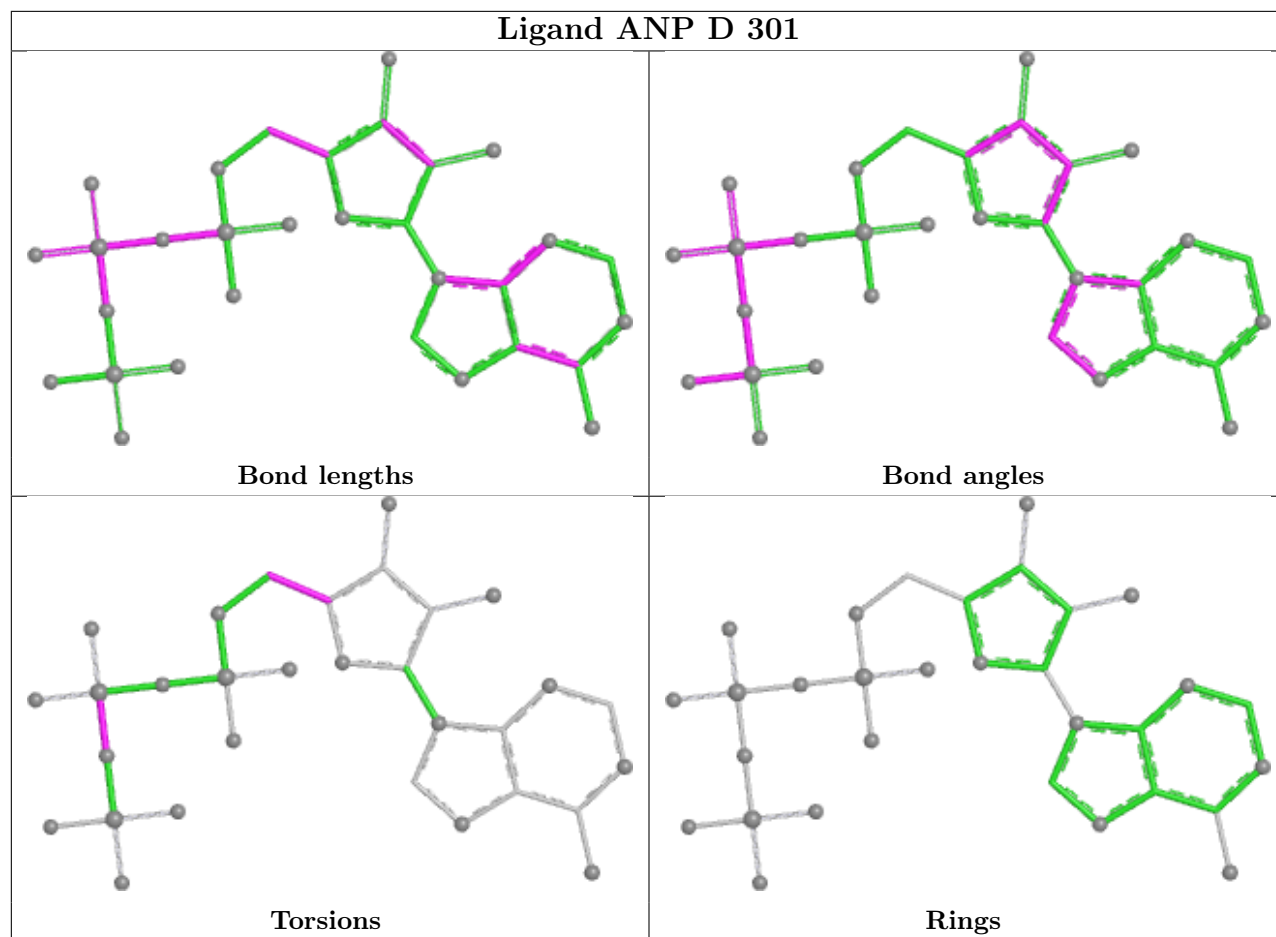
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	301	ANP	1	0
3	D	301	ANP	3	0

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	211/214 (98%)	-0.07	0 100 100	29, 55, 93, 107	0
1	B	210/214 (98%)	-0.02	0 100 100	23, 56, 87, 97	0
1	C	209/214 (97%)	0.08	1 (0%) 87 67	35, 58, 91, 109	0
1	D	211/214 (98%)	-0.11	2 (0%) 81 57	28, 52, 75, 91	0
2	K	15/19 (78%)	0.86	1 (6%) 24 16	11, 55, 64, 64	0
2	R	10/19 (52%)	0.86	2 (20%) 3 3	61, 66, 68, 69	0
2	Y	9/19 (47%)	0.62	0 100 100	59, 61, 67, 67	0
2	Z	7/19 (36%)	0.78	0 100 100	50, 55, 57, 58	0
All	All	882/932 (94%)	0.01	6 (0%) 84 62	11, 56, 87, 109	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	178	ASP	3.1
2	K	156	GLU	2.5
1	C	109	PRO	2.5
2	R	151	THR	2.3
1	D	110	LEU	2.1
2	R	150	MET	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](#)

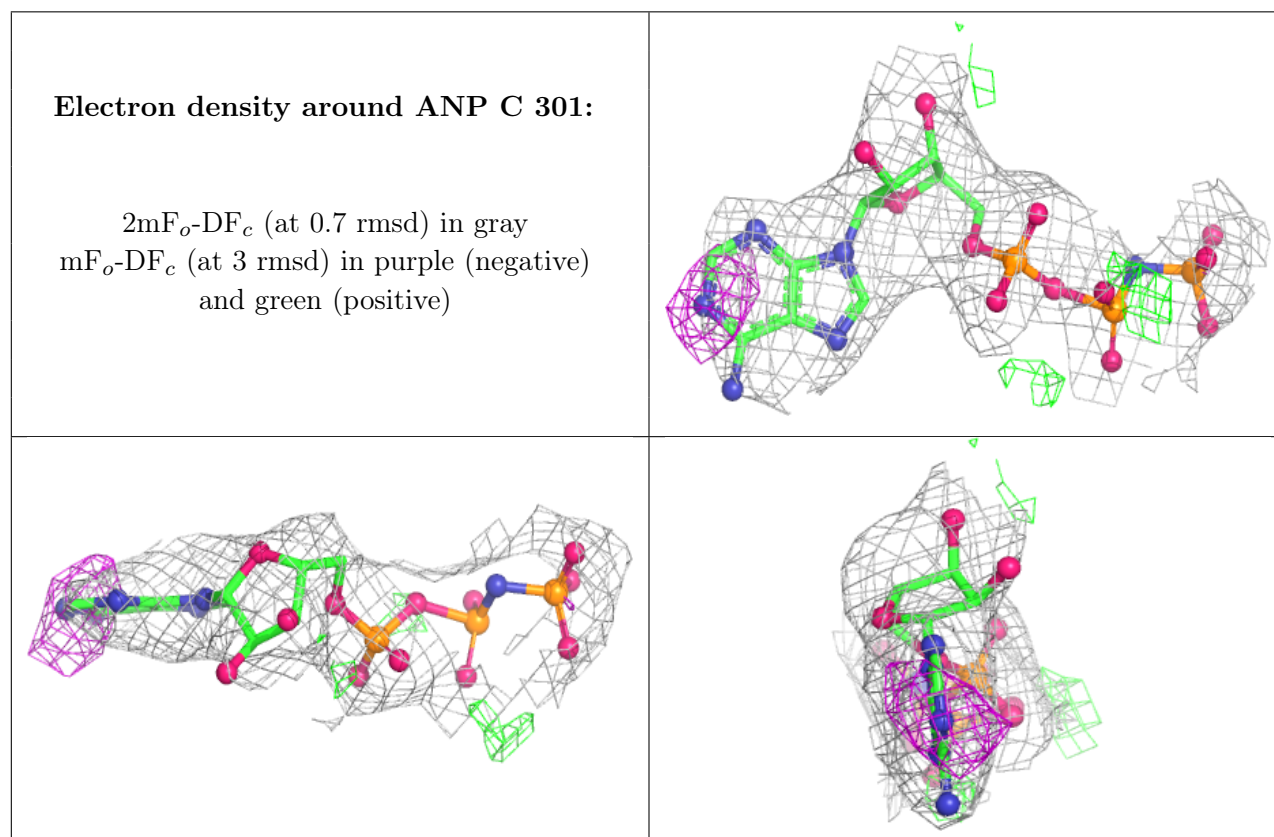
There are no oligosaccharides in this entry.

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.

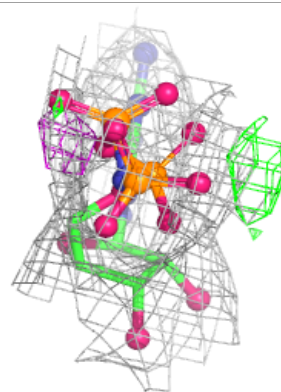
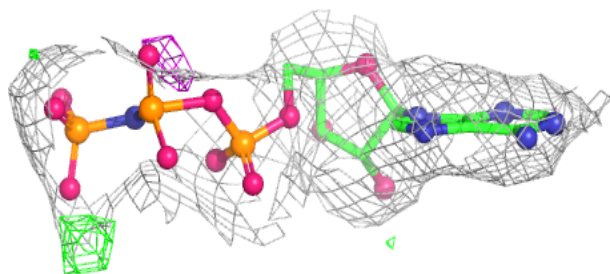
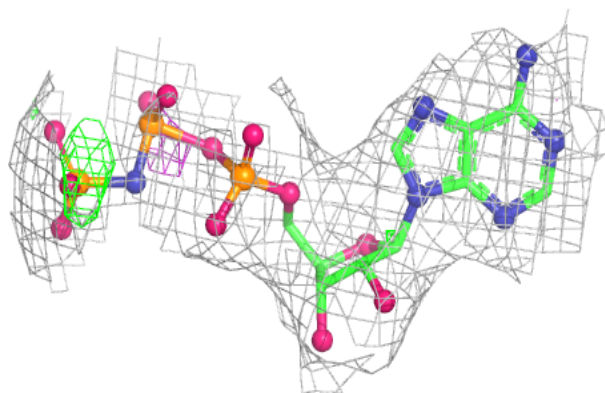
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ANP	C	301	31/31	0.91	0.10	20,27,54,56	0
3	ANP	A	500	31/31	0.92	0.09	26,34,47,48	0
3	ANP	B	501	31/31	0.95	0.07	23,36,54,56	0
3	ANP	D	301	31/31	0.95	0.07	19,29,49,49	0

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.

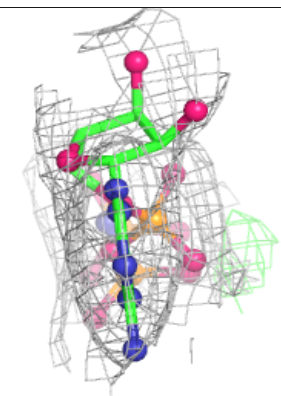
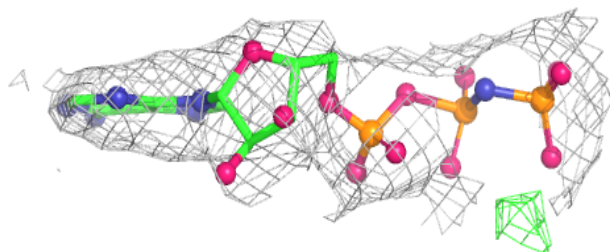
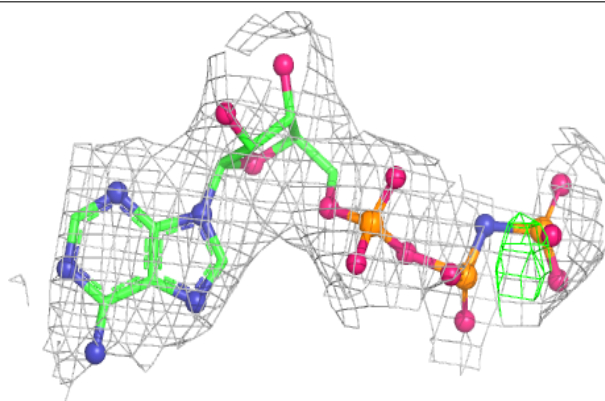


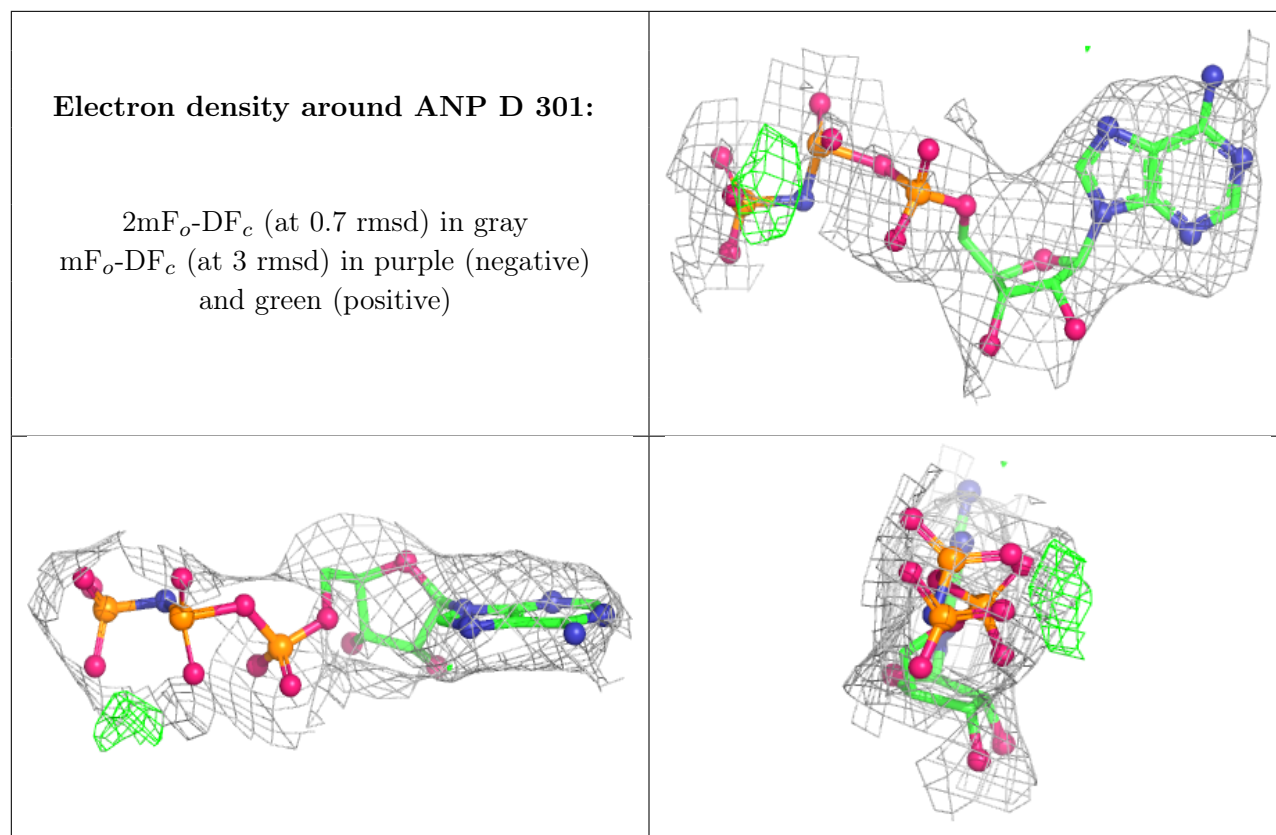
Electron density around ANP A 500:

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

**Electron density around ANP B 501:**

$2mF_o-DF_c$ (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.