



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 07:07 am BST

PDB ID : 2P9U
Title : Crystal structure of bovine Arp2/3 complex co-crystallized with AMP-PNP and calcium
Authors : Nolen, B.J.; Pollard, T.D.
Deposited on : 2007-03-26
Resolution : 2.75 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

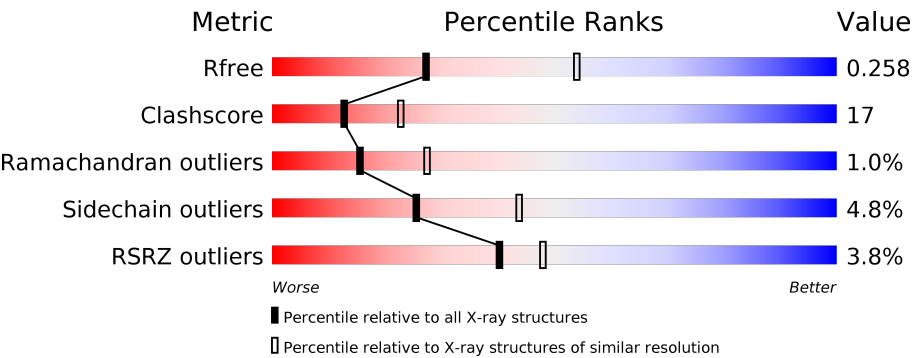
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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 on 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	418	<div><div>5%</div><div><div></div><div>64%</div><div>28%</div><div>• 5%</div></div></div>
2	B	394	<div><div>3%</div><div><div></div><div>31%</div><div>22%</div><div>• 44%</div></div></div>
3	C	372	<div><div>2%</div><div><div></div><div>64%</div><div>25%</div><div>• 8%</div></div></div>
4	D	300	<div><div>%</div><div><div></div><div>70%</div><div>23%</div><div>• 7%</div></div></div>
5	E	178	<div><div>8%</div><div><div></div><div>60%</div><div>34%</div><div>• • •</div></div></div>
6	F	168	<div><div>%</div><div><div></div><div>74%</div><div>24%</div><div>• •</div></div></div>

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Mol	Chain	Length	Quality of chain
7	G	151	<div><div></div><div>5%</div><div>48%</div><div>26%</div><div>•</div><div>23%</div></div>

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 13719 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 Actin-like protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	398	Total	C	N	O	S	0	0	0
			3176	2040	527	594	15			

- Molecule 2 is a protein called Actin-like protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	222	Total	C	N	O	S	0	0	0
			1679	1075	286	313	5			

- Molecule 3 is a protein called Actin-related protein 2/3 complex subunit 1B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	341	Total	C	N	O	S	0	0	0
			2638	1674	460	485	19			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	58	VAL	ILE	CONFLICT	UNP Q58CQ2

- Molecule 4 is a protein called Actin-related protein 2/3 complex subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	280	Total	C	N	O	S	0	0	0
			2256	1436	389	423	8			

- Molecule 5 is a protein called Actin-related protein 2/3 complex subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	174	Total	C	N	O	S	0	0	0
			1415	908	236	262	9			

- Molecule 6 is a protein called Actin-related protein 2/3 complex subunit 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	167	Total	C	N	O	S	0	0	0
			1371	875	239	248	9			

- Molecule 7 is a protein called Actin-related protein 2/3 complex subunit 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	116	Total	C	N	O	S	0	0	0
			868	552	146	168	2			

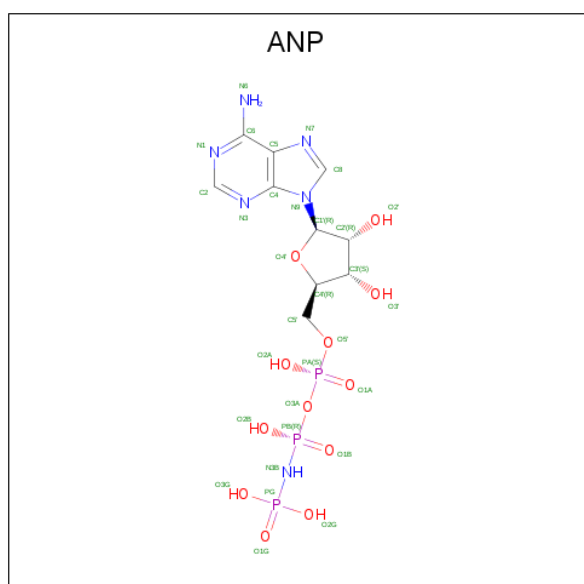
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	17	ASP	GLY	CONFLICT	UNP Q3SYX9
G	28	ASP	GLU	CONFLICT	UNP Q3SYX9

- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Ca	0	0
			1	1		
8	A	1	Total	Ca	0	0
			1	1		

- Molecule 9 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

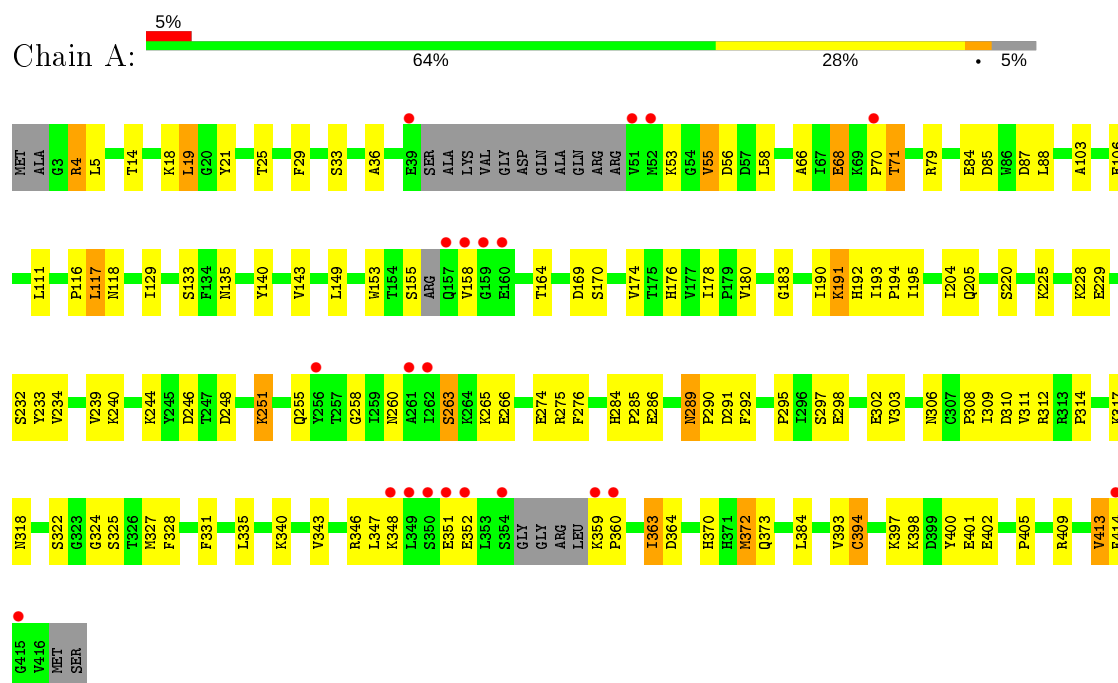
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	42	Total	O	0	0
			42	42		
10	B	24	Total	O	0	0
			24	24		
10	C	73	Total	O	0	0
			73	73		
10	D	58	Total	O	0	0
			58	58		
10	E	4	Total	O	0	0
			4	4		
10	F	45	Total	O	0	0
			45	45		
10	G	6	Total	O	0	0
			6	6		

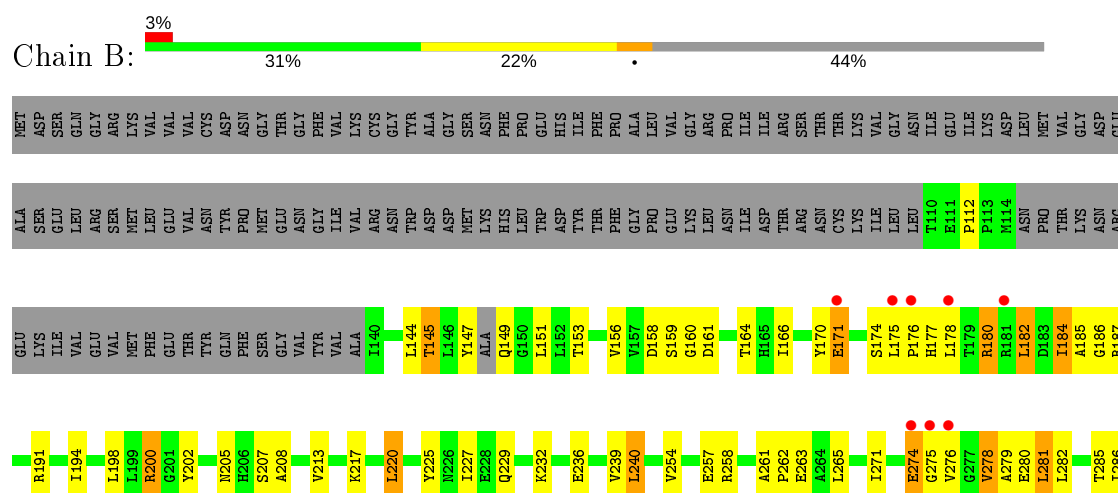
3 Residue-property plots [i](#)

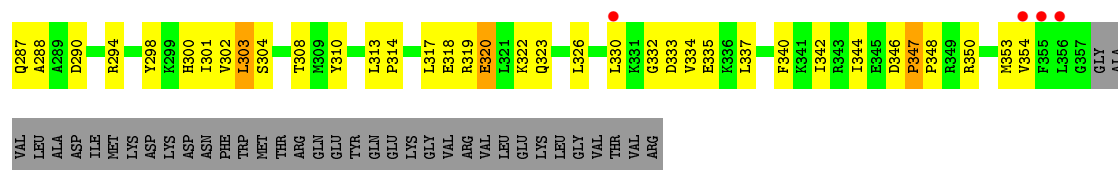
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Actin-like protein 3

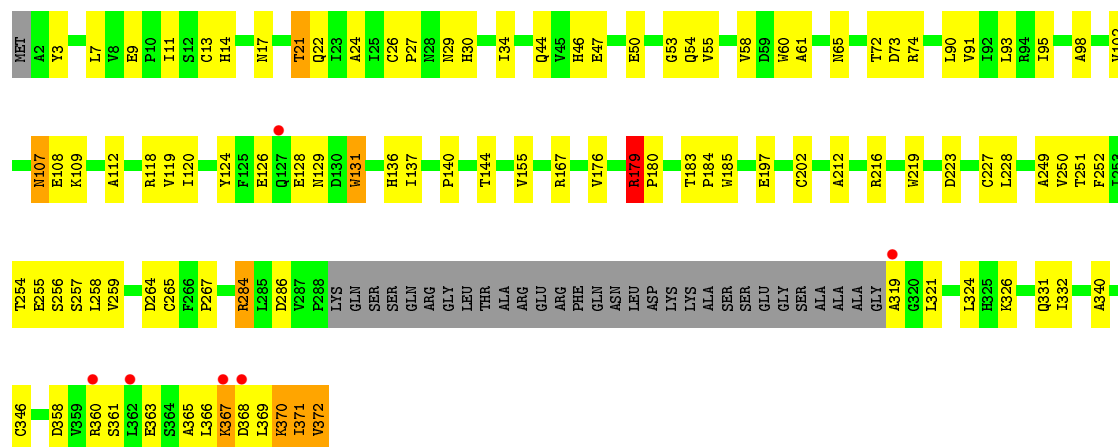


• Molecule 2: Actin-like protein 2

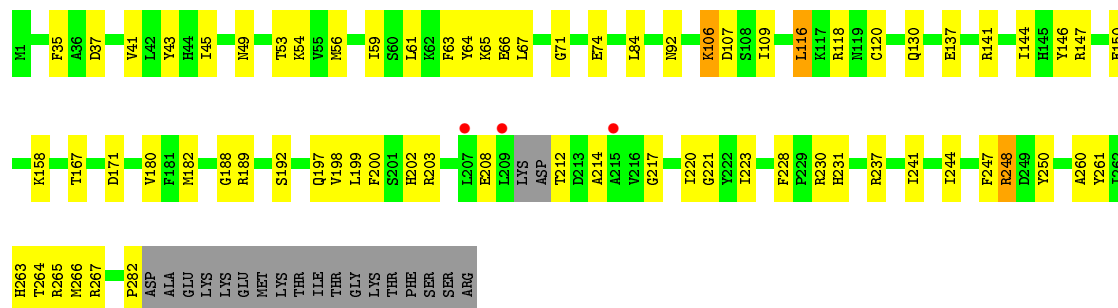




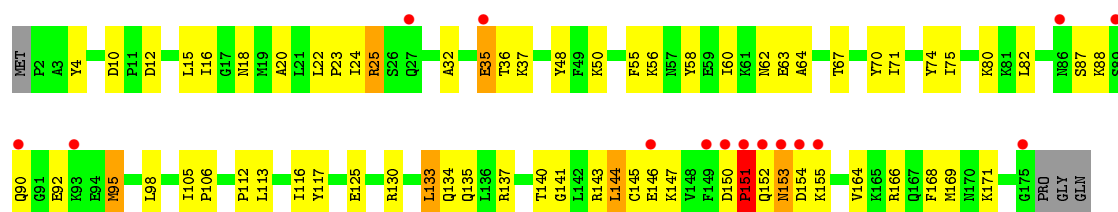
• Molecule 3: Actin-related protein 2/3 complex subunit 1B

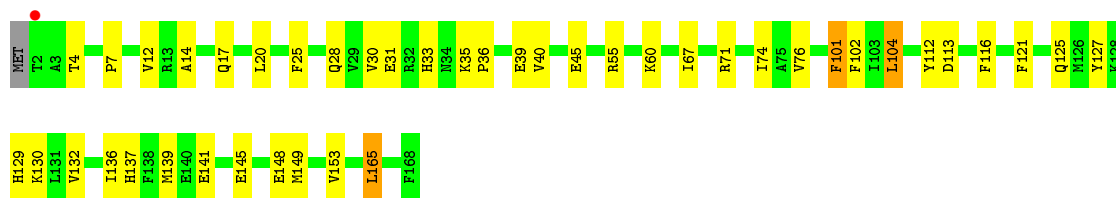
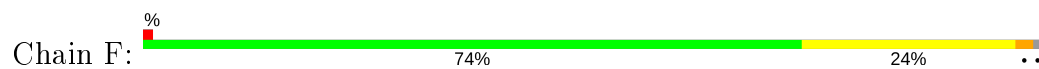


• Molecule 4: Actin-related protein 2/3 complex subunit 2

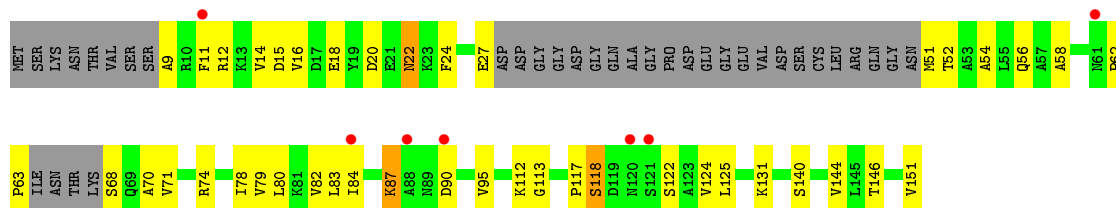


• Molecule 5: Actin-related protein 2/3 complex subunit 3





● Molecule 7: Actin-related protein 2/3 complex subunit 5



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.12Å 129.35Å 203.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.75 47.41 – 2.75	Depositor EDS
% Data completeness (in resolution range)	94.5 (30.00-2.75) 94.5 (47.41-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 2.77Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.219 , 0.265 0.212 , 0.258	Depositor DCC
R_{free} test set	3866 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.4	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	0.92	EDS
Total number of atoms	13719	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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: CA, 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	0.38	0/3256	0.61	1/4418 (0.0%)
2	B	0.39	0/1708	0.65	1/2318 (0.0%)
3	C	0.40	0/2707	0.69	1/3677 (0.0%)
4	D	0.38	0/2304	0.61	0/3111
5	E	0.34	0/1449	0.58	0/1954
6	F	0.41	0/1393	0.62	0/1868
7	G	0.34	0/878	0.55	0/1184
All	All	0.38	0/13695	0.62	3/18530 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
4	D	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	11	ILE	N-CA-C	-6.13	94.44	111.00
2	B	112	PRO	N-CA-CB	5.79	110.25	103.30
1	A	118	ASN	N-CA-C	-5.49	96.17	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	261	TYR	Sidechain

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	3176	0	3106	103	0
2	B	1679	0	1624	93	0
3	C	2638	0	2580	87	0
4	D	2256	0	2219	47	0
5	E	1415	0	1416	70	0
6	F	1371	0	1410	36	0
7	G	868	0	866	37	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
9	A	31	0	13	3	0
9	B	31	0	13	6	0
10	A	42	0	0	1	0
10	B	24	0	0	4	0
10	C	73	0	0	0	0
10	D	58	0	0	1	0
10	E	4	0	0	0	0
10	F	45	0	0	1	0
10	G	6	0	0	0	0
All	All	13719	0	13247	444	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 17.

All (444) 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)
2:B:151:LEU:HD13	2:B:300:HIS:HD2	1.26	1.01
3:C:223:ASP:HB3	7:G:146:THR:HG21	1.42	1.01
2:B:205:ASN:HD22	2:B:208:ALA:H	1.06	1.00
3:C:183:THR:HG22	3:C:185:TRP:H	1.24	0.99
3:C:256:SER:HB2	3:C:372:VAL:HG12	1.48	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:ILE:HA	1:A:312:ARG:HE	1.30	0.95
2:B:337:LEU:HD21	7:G:16:VAL:HG13	1.47	0.95
1:A:363:ILE:H	1:A:363:ILE:HD13	1.31	0.94
3:C:14:HIS:H	3:C:331:GLN:HE22	1.16	0.93
1:A:191:LYS:HE2	1:A:303:VAL:HG22	1.48	0.92
6:F:4:THR:HG23	6:F:55:ARG:HE	1.30	0.92
5:E:75:ILE:HG23	5:E:144:LEU:HD11	1.52	0.90
2:B:205:ASN:ND2	2:B:208:ALA:H	1.72	0.87
2:B:279:ALA:HB1	2:B:320:GLU:HG2	1.57	0.85
6:F:130:LYS:HE2	6:F:130:LYS:HA	1.58	0.85
5:E:15:LEU:HD22	5:E:63:GLU:HG3	1.60	0.84
3:C:167:ARG:HG2	3:C:197:GLU:HG3	1.60	0.83
3:C:367:LYS:HD2	3:C:368:ASP:N	1.96	0.80
5:E:88:LYS:H	5:E:153:ASN:ND2	1.80	0.80
2:B:175:LEU:HD12	2:B:175:LEU:H	1.45	0.80
5:E:152:GLN:HB2	5:E:155:LYS:HD2	1.63	0.79
4:D:65:LYS:HA	4:D:65:LYS:HE2	1.64	0.79
5:E:36:THR:HG23	5:E:37:LYS:HD3	1.65	0.78
2:B:205:ASN:HD22	2:B:208:ALA:N	1.82	0.78
1:A:343:VAL:HG23	1:A:346:ARG:HH21	1.48	0.77
2:B:151:LEU:CD1	2:B:300:HIS:HD2	1.98	0.77
4:D:228:PHE:H	4:D:231:HIS:HD2	1.33	0.77
1:A:343:VAL:HG21	1:A:363:ILE:HG13	1.65	0.76
2:B:279:ALA:CB	2:B:320:GLU:HG2	2.16	0.76
3:C:126:GLU:HB2	3:C:131:TRP:HZ3	1.51	0.75
3:C:179:ARG:HH11	3:C:179:ARG:HG3	1.50	0.75
3:C:202:CYS:HB2	7:G:146:THR:HG22	1.69	0.74
1:A:363:ILE:CD1	1:A:363:ILE:H	2.01	0.73
5:E:15:LEU:CD2	5:E:63:GLU:HG3	2.17	0.73
1:A:309:ILE:HA	1:A:312:ARG:NE	2.02	0.73
1:A:343:VAL:HG23	1:A:346:ARG:NH2	2.04	0.72
2:B:151:LEU:HD13	2:B:300:HIS:CD2	2.18	0.72
7:G:112:LYS:HE2	7:G:151:VAL:O	1.90	0.71
6:F:101:PHE:HB3	6:F:104:LEU:HB2	1.70	0.71
5:E:95:MET:HG2	5:E:141:GLY:O	1.91	0.70
2:B:282:LEU:HD21	2:B:301:ILE:HD13	1.74	0.70
1:A:155:SER:O	1:A:158:VAL:HG22	1.92	0.69
4:D:228:PHE:H	4:D:231:HIS:CD2	2.09	0.69
3:C:370:LYS:HG2	3:C:370:LYS:O	1.92	0.69
5:E:143:ARG:O	5:E:146:GLU:HG2	1.92	0.68
1:A:55:VAL:HG22	1:A:58:LEU:HD12	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:TYR:HB2	2:B:175:LEU:HD11	1.75	0.68
4:D:188:GLY:HA3	6:F:165:LEU:HD23	1.76	0.68
3:C:14:HIS:N	3:C:331:GLN:HE22	1.92	0.67
3:C:34:ILE:HB	3:C:46:HIS:HB2	1.77	0.67
4:D:188:GLY:HA3	6:F:165:LEU:CD2	2.24	0.67
2:B:175:LEU:N	2:B:175:LEU:HD12	2.10	0.67
2:B:322:LYS:HB3	7:G:16:VAL:HG11	1.75	0.67
5:E:25:ARG:NH1	5:E:35:GLU:HB3	2.10	0.67
3:C:107:ASN:ND2	3:C:109:LYS:H	1.94	0.66
3:C:155:VAL:HG21	3:C:180:PRO:HG3	1.78	0.66
1:A:14:THR:CG2	1:A:79:ARG:HH21	2.09	0.66
1:A:183:GLY:HA3	1:A:413:VAL:HG21	1.76	0.66
3:C:119:VAL:HG22	3:C:120:ILE:N	2.10	0.66
3:C:126:GLU:HB2	3:C:131:TRP:CZ3	2.31	0.65
2:B:313:LEU:HB3	2:B:314:PRO:HD3	1.79	0.65
3:C:72:THR:HA	3:C:98:ALA:HB1	1.78	0.65
1:A:363:ILE:N	1:A:363:ILE:HD13	2.09	0.64
7:G:80:LEU:O	7:G:84:ILE:HD13	1.97	0.64
2:B:350:ARG:HA	2:B:353:MET:HG3	1.80	0.64
4:D:180:VAL:HG11	6:F:153:VAL:HG12	1.80	0.63
2:B:346:ASP:OD1	2:B:350:ARG:NH1	2.31	0.63
2:B:229:GLN:NE2	6:F:40:VAL:HB	2.12	0.63
2:B:194:ILE:HG12	2:B:213:VAL:HG21	1.80	0.63
1:A:204:ILE:HD12	1:A:228:LYS:HB2	1.80	0.62
2:B:180:ARG:HD2	2:B:285:THR:OG1	1.99	0.62
5:E:25:ARG:HD2	5:E:35:GLU:HG3	1.81	0.62
5:E:88:LYS:O	5:E:92:GLU:HG3	1.98	0.62
2:B:354:VAL:HA	10:B:903:HOH:O	2.00	0.62
5:E:23:PRO:O	5:E:24:ILE:HG23	1.99	0.62
1:A:409:ARG:HD3	2:B:200:ARG:O	1.99	0.61
2:B:261:ALA:HB3	2:B:262:PRO:HD3	1.82	0.61
2:B:147:TYR:O	2:B:149:GLN:N	2.34	0.61
1:A:116:PRO:O	1:A:117:LEU:HB2	2.01	0.60
4:D:150:GLU:HG2	4:D:167:THR:HA	1.84	0.60
1:A:79:ARG:HB2	1:A:84:GLU:HG3	1.82	0.60
3:C:257:SER:OG	3:C:371:ILE:O	2.16	0.60
3:C:183:THR:HG23	3:C:184:PRO:HD2	1.83	0.60
1:A:55:VAL:O	1:A:55:VAL:HG13	2.01	0.60
2:B:175:LEU:O	2:B:177:HIS:O	2.20	0.60
1:A:359:LYS:N	1:A:360:PRO:HD3	2.17	0.60
2:B:161:ASP:O	2:B:187:ARG:HG2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:248:ARG:C	4:D:248:ARG:HD3	2.23	0.59
1:A:239:VAL:HG11	5:E:48:TYR:HD1	1.67	0.59
4:D:59:ILE:HB	4:D:116:LEU:HD13	1.84	0.59
5:E:95:MET:HA	5:E:95:MET:HE3	1.83	0.59
3:C:14:HIS:H	3:C:331:GLN:NE2	1.95	0.59
5:E:25:ARG:CD	5:E:35:GLU:HG3	2.33	0.59
3:C:107:ASN:HD22	3:C:108:GLU:N	2.00	0.59
6:F:25:PHE:CD1	6:F:67:ILE:HD13	2.38	0.59
2:B:174:SER:O	2:B:176:PRO:HD3	2.03	0.58
7:G:20:ASP:OD1	7:G:22:ASN:HB2	2.02	0.58
5:E:67:THR:O	5:E:71:ILE:HG13	2.03	0.58
3:C:252:PHE:HA	3:C:258:LEU:HD23	1.86	0.58
7:G:117:PRO:O	7:G:118:SER:HB3	2.02	0.58
5:E:18:ASN:O	5:E:63:GLU:HB3	2.04	0.57
4:D:203:ARG:HG3	4:D:217:GLY:O	2.04	0.57
2:B:232:LYS:HE3	2:B:236:GLU:OE2	2.04	0.57
3:C:3:TYR:HB2	3:C:324:LEU:HG	1.85	0.57
5:E:88:LYS:H	5:E:153:ASN:HD21	1.52	0.57
2:B:323:GLN:HG3	7:G:16:VAL:CG2	2.34	0.57
5:E:80:LYS:HG3	5:E:164:VAL:HG23	1.87	0.57
7:G:87:LYS:N	7:G:87:LYS:HD3	2.19	0.57
4:D:53:THR:C	4:D:54:LYS:HD2	2.24	0.57
1:A:274:GLU:OE1	1:A:274:GLU:N	2.35	0.57
1:A:384:LEU:HB3	1:A:414:PHE:CZ	2.40	0.57
3:C:30:HIS:HA	3:C:53:GLY:O	2.04	0.57
5:E:62:ASN:OD1	5:E:64:ALA:HB3	2.05	0.57
2:B:175:LEU:CD1	2:B:175:LEU:H	2.14	0.56
1:A:194:PRO:O	1:A:195:ILE:HD12	2.04	0.56
5:E:87:SER:HA	5:E:153:ASN:OD1	2.06	0.56
2:B:239:VAL:HG23	2:B:240:LEU:HD13	1.88	0.56
3:C:107:ASN:HD22	3:C:107:ASN:C	2.07	0.56
1:A:343:VAL:CG2	1:A:363:ILE:HG13	2.33	0.56
2:B:323:GLN:HG3	7:G:16:VAL:HG21	1.88	0.56
5:E:24:ILE:HD11	5:E:135:GLN:HG2	1.88	0.56
1:A:85:ASP:OD2	1:A:88:LEU:HD22	2.05	0.56
1:A:4:ARG:HB2	1:A:4:ARG:HH11	1.70	0.55
5:E:36:THR:HG23	5:E:37:LYS:H	1.70	0.55
1:A:71:THR:OG1	1:A:71:THR:O	2.25	0.55
1:A:311:VAL:C	1:A:314:PRO:HD2	2.26	0.55
1:A:4:ARG:HD2	1:A:5:LEU:HG	1.88	0.55
7:G:74:ARG:O	7:G:78:ILE:HG13	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:VAL:HG11	5:E:48:TYR:CD1	2.42	0.55
1:A:309:ILE:HG23	1:A:310:ASP:N	2.21	0.55
2:B:254:VAL:HG13	2:B:257:GLU:HG2	1.88	0.55
1:A:308:PRO:O	1:A:311:VAL:HG12	2.07	0.54
4:D:220:ILE:HG22	4:D:221:GLY:N	2.20	0.54
1:A:55:VAL:CG2	1:A:58:LEU:HD12	2.37	0.54
5:E:150:ASP:O	5:E:152:GLN:N	2.40	0.54
5:E:58:TYR:CD1	5:E:168:PHE:HZ	2.24	0.54
3:C:216:ARG:NH1	3:C:255:GLU:O	2.40	0.54
7:G:9:ALA:HA	7:G:11:PHE:CE2	2.42	0.54
3:C:21:THR:HG22	3:C:22:GLN:HG3	1.90	0.54
7:G:11:PHE:O	7:G:14:VAL:HG12	2.08	0.54
1:A:239:VAL:HG13	5:E:4:TYR:CE2	2.44	0.53
1:A:328:PHE:CE1	9:A:901:ANP:H2	2.44	0.53
3:C:358:ASP:OD1	3:C:360:ARG:HG2	2.08	0.53
1:A:87:ASP:OD2	4:D:267:ARG:HD2	2.08	0.53
3:C:119:VAL:HG22	3:C:120:ILE:H	1.72	0.53
2:B:290:ASP:O	2:B:294:ARG:HG3	2.09	0.53
4:D:64:TYR:HB3	4:D:92:ASN:ND2	2.23	0.53
6:F:127:TYR:HB3	6:F:129:HIS:CE1	2.44	0.53
9:A:901:ANP:O1B	9:A:901:ANP:O2G	2.27	0.53
2:B:276:VAL:HB	2:B:280:GLU:HB3	1.89	0.53
7:G:78:ILE:O	7:G:82:VAL:HG23	2.09	0.53
5:E:36:THR:CG2	5:E:37:LYS:HD3	2.37	0.53
7:G:15:ASP:OD2	7:G:18:GLU:HG2	2.09	0.53
6:F:30:VAL:HG11	6:F:33:HIS:ND1	2.24	0.53
6:F:33:HIS:ND1	6:F:35:LYS:NZ	2.56	0.53
3:C:363:GLU:HG2	3:C:369:LEU:HD23	1.91	0.53
3:C:223:ASP:CB	7:G:146:THR:HG21	2.30	0.52
2:B:177:HIS:O	2:B:178:LEU:HB2	2.10	0.52
5:E:25:ARG:HH11	5:E:35:GLU:HB3	1.75	0.52
7:G:79:VAL:O	7:G:83:LEU:HG	2.09	0.52
2:B:177:HIS:CD2	2:B:177:HIS:H	2.26	0.52
4:D:130:GLN:OE1	4:D:130:GLN:HA	2.10	0.52
3:C:126:GLU:CB	3:C:131:TRP:HZ3	2.23	0.52
3:C:109:LYS:HD3	3:C:176:VAL:O	2.10	0.52
3:C:118:ARG:HG3	3:C:118:ARG:HH11	1.75	0.51
3:C:363:GLU:CD	3:C:371:ILE:HD12	2.30	0.51
6:F:74:ILE:HD13	6:F:139:MET:HG2	1.91	0.51
1:A:79:ARG:HB2	1:A:84:GLU:CG	2.40	0.51
4:D:65:LYS:CA	4:D:65:LYS:HE2	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:121:PHE:O	6:F:125:GLN:HG3	2.10	0.51
1:A:397:LYS:HG2	1:A:401:GLU:OE2	2.10	0.51
1:A:149:LEU:HD11	1:A:180:VAL:HB	1.92	0.51
2:B:326:LEU:HD23	7:G:14:VAL:HG11	1.93	0.51
3:C:61:ALA:HB1	3:C:108:GLU:OE1	2.11	0.51
2:B:333:ASP:C	2:B:335:GLU:H	2.15	0.51
3:C:367:LYS:HD2	3:C:367:LYS:C	2.32	0.51
5:E:36:THR:HG23	5:E:37:LYS:N	2.26	0.51
7:G:51:MET:O	7:G:54:ALA:HB3	2.10	0.51
7:G:58:ALA:HB1	7:G:79:VAL:HG22	1.92	0.51
1:A:129:ILE:O	1:A:133:SER:HB2	2.11	0.50
5:E:112:PRO:O	5:E:113:LEU:HB2	2.11	0.50
5:E:150:ASP:OD1	5:E:151:PRO:HD2	2.11	0.50
7:G:87:LYS:H	7:G:87:LYS:HE3	1.77	0.50
2:B:198:LEU:HA	2:B:202:TYR:O	2.11	0.50
4:D:45:ILE:HA	4:D:56:MET:O	2.11	0.50
7:G:124:VAL:HG13	7:G:125:LEU:N	2.26	0.50
1:A:398:LYS:HG3	1:A:402:GLU:OE2	2.12	0.50
7:G:95:VAL:HG21	7:G:131:LYS:HB3	1.93	0.50
4:D:202:HIS:O	4:D:220:ILE:O	2.30	0.50
6:F:30:VAL:HG11	6:F:33:HIS:CE1	2.46	0.50
9:B:902:ANP:O3A	9:B:902:ANP:O3G	2.30	0.50
1:A:289:ASN:HD22	1:A:290:PRO:CD	2.24	0.50
1:A:348:LYS:O	1:A:352:GLU:HG3	2.12	0.50
4:D:197:GLN:HG2	4:D:199:LEU:HD11	1.93	0.50
1:A:263:SER:C	1:A:265:LYS:H	2.15	0.50
2:B:182:LEU:HD13	2:B:184:ILE:HG23	1.93	0.50
6:F:36:PRO:HG2	6:F:39:GLU:HB2	1.93	0.50
2:B:166:ILE:HD13	2:B:282:LEU:HA	1.94	0.49
1:A:205:GLN:HE21	1:A:220:SER:CB	2.26	0.49
4:D:182:MET:HG3	4:D:200:PHE:CD1	2.47	0.49
4:D:189:ARG:HA	4:D:192:SER:O	2.12	0.49
4:D:54:LYS:HD2	4:D:54:LYS:N	2.25	0.49
3:C:370:LYS:CG	3:C:370:LYS:O	2.58	0.49
2:B:217:LYS:HD2	2:B:310:TYR:OH	2.13	0.49
5:E:168:PHE:CE2	5:E:169:MET:HE2	2.48	0.49
5:E:16:ILE:O	5:E:16:ILE:HG23	2.12	0.49
3:C:60:TRP:HE1	3:C:65:ASN:ND2	2.10	0.49
4:D:282:PRO:HG3	6:F:127:TYR:CZ	2.48	0.49
3:C:119:VAL:CG2	3:C:120:ILE:N	2.76	0.49
1:A:153:TRP:HA	1:A:158:VAL:HG21	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:ASN:HD22	1:A:290:PRO:N	2.11	0.49
2:B:254:VAL:CG1	2:B:257:GLU:HG2	2.43	0.49
5:E:60:ILE:CD1	5:E:116:ILE:HG21	2.43	0.49
5:E:88:LYS:N	5:E:153:ASN:HD21	2.11	0.49
3:C:264:ASP:O	3:C:265:CYS:HB2	2.12	0.49
1:A:370:HIS:HE1	10:A:907:HOH:O	1.94	0.49
6:F:20:LEU:HD12	6:F:132:VAL:CG2	2.43	0.49
2:B:170:TYR:CB	2:B:175:LEU:HD11	2.41	0.48
3:C:284:ARG:HD3	3:C:286:ASP:O	2.13	0.48
2:B:182:LEU:HD11	2:B:278:VAL:H	1.78	0.48
4:D:71:GLY:HA2	4:D:74:GLU:OE1	2.14	0.48
5:E:16:ILE:HD12	5:E:125:GLU:HG3	1.95	0.48
4:D:147:ARG:HB2	4:D:150:GLU:HB2	1.94	0.48
3:C:129:ASN:HB2	3:C:131:TRP:CZ3	2.48	0.48
6:F:60:LYS:HE3	6:F:112:TYR:CE2	2.48	0.48
3:C:367:LYS:NZ	3:C:368:ASP:HB3	2.28	0.48
5:E:169:MET:O	5:E:171:LYS:HG2	2.13	0.48
1:A:343:VAL:O	1:A:347:LEU:HD13	2.12	0.48
2:B:159:SER:HB3	2:B:308:THR:HG23	1.94	0.48
6:F:12:VAL:HG12	6:F:136:ILE:HD11	1.96	0.48
1:A:239:VAL:HG23	1:A:240:LYS:N	2.29	0.48
2:B:186:GLY:HA3	9:B:902:ANP:O3'	2.13	0.48
3:C:13:CYS:SG	3:C:58:VAL:HG23	2.53	0.48
1:A:164:THR:HA	1:A:180:VAL:O	2.14	0.48
3:C:212:ALA:HB3	3:C:255:GLU:OE2	2.14	0.48
1:A:14:THR:HG22	1:A:79:ARG:HH21	1.77	0.47
7:G:68:SER:O	7:G:71:VAL:HG12	2.14	0.47
1:A:79:ARG:NE	1:A:79:ARG:HA	2.29	0.47
3:C:179:ARG:HG3	3:C:179:ARG:NH1	2.18	0.47
4:D:106:LYS:HD2	4:D:106:LYS:N	2.29	0.47
5:E:150:ASP:C	5:E:152:GLN:N	2.67	0.47
7:G:87:LYS:N	7:G:87:LYS:CD	2.77	0.47
1:A:239:VAL:HG23	1:A:240:LYS:H	1.79	0.47
4:D:241:ILE:HA	4:D:244:ILE:HG22	1.95	0.47
1:A:170:SER:OG	1:A:325:SER:HB2	2.13	0.47
1:A:183:GLY:CA	1:A:413:VAL:HG21	2.42	0.47
2:B:170:TYR:CD2	2:B:171:GLU:HG2	2.49	0.47
3:C:365:ALA:C	3:C:366:LEU:HD12	2.35	0.47
1:A:372:MET:SD	1:A:372:MET:N	2.88	0.47
1:A:400:TYR:CE1	1:A:405:PRO:HB3	2.49	0.47
1:A:140:TYR:HB2	1:A:394:CYS:SG	2.55	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:70:TYR:CZ	5:E:133:LEU:HG	2.50	0.47
5:E:23:PRO:HB3	5:E:35:GLU:HA	1.97	0.47
1:A:225:LYS:HE3	1:A:229:GLU:OE1	2.14	0.47
5:E:74:TYR:OH	5:E:98:LEU:HD12	2.15	0.47
1:A:21:TYR:OH	1:A:103:ALA:HB2	2.15	0.47
3:C:254:THR:HA	3:C:340:ALA:O	2.14	0.47
4:D:61:LEU:HD23	4:D:63:PHE:CZ	2.49	0.47
5:E:116:ILE:HG22	5:E:117:TYR:CD1	2.50	0.47
5:E:87:SER:OG	5:E:90:GLN:HB2	2.13	0.47
2:B:184:ILE:HD12	2:B:271:ILE:HD11	1.95	0.46
5:E:60:ILE:HD12	5:E:116:ILE:HD13	1.96	0.46
3:C:14:HIS:HA	3:C:24:ALA:O	2.15	0.46
3:C:72:THR:HA	3:C:98:ALA:CB	2.44	0.46
5:E:130:ARG:HG3	5:E:130:ARG:HH11	1.79	0.46
5:E:150:ASP:C	5:E:152:GLN:H	2.19	0.46
2:B:254:VAL:HG12	2:B:258:ARG:HG3	1.97	0.46
3:C:144:THR:H	6:F:28:GLN:NE2	2.14	0.46
3:C:363:GLU:OE2	3:C:371:ILE:HD12	2.16	0.46
1:A:289:ASN:ND2	1:A:291:ASP:H	2.13	0.46
2:B:182:LEU:HD11	2:B:278:VAL:N	2.30	0.46
5:E:56:LYS:HD2	5:E:56:LYS:N	2.31	0.46
2:B:229:GLN:NE2	6:F:40:VAL:CB	2.78	0.46
2:B:160:GLY:O	2:B:185:ALA:HB1	2.16	0.46
2:B:303:LEU:CD2	2:B:314:PRO:HG3	2.46	0.46
3:C:90:LEU:HD12	3:C:91:VAL:H	1.79	0.46
3:C:7:LEU:HD12	3:C:9:GLU:HB2	1.98	0.46
4:D:263:HIS:HD2	4:D:266:MET:CE	2.29	0.46
4:D:66:GLU:OE1	4:D:144:ILE:HA	2.16	0.46
2:B:161:ASP:H	9:B:902:ANP:PG	2.39	0.46
4:D:265:ARG:NH2	6:F:148:GLU:HB3	2.30	0.46
1:A:169:ASP:HA	1:A:322:SER:O	2.15	0.46
1:A:328:PHE:CZ	9:A:901:ANP:H2	2.51	0.46
2:B:161:ASP:OD2	2:B:187:ARG:HB3	2.16	0.46
2:B:151:LEU:CD1	2:B:300:HIS:CD2	2.89	0.45
1:A:246:ASP:OD1	5:E:50:LYS:HE3	2.16	0.45
2:B:160:GLY:HA3	9:B:902:ANP:O3G	2.16	0.45
3:C:250:VAL:HG22	3:C:251:THR:N	2.31	0.45
5:E:95:MET:HA	5:E:95:MET:CE	2.46	0.45
2:B:227:ILE:HD11	2:B:263:GLU:OE2	2.16	0.45
2:B:182:LEU:HG	2:B:281:LEU:HD12	1.98	0.45
1:A:258:GLY:O	1:A:266:GLU:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:182:LEU:HD23	2:B:182:LEU:HA	1.85	0.45
5:E:63:GLU:H	5:E:63:GLU:CD	2.20	0.45
1:A:18:LYS:HD3	1:A:18:LYS:N	2.32	0.45
1:A:232:SER:HA	1:A:275:ARG:O	2.16	0.45
2:B:184:ILE:HD13	2:B:265:LEU:HA	1.97	0.45
2:B:340:PHE:CE2	2:B:342:ILE:HD11	2.52	0.45
2:B:317:LEU:HB3	2:B:344:ILE:HD13	1.98	0.45
3:C:119:VAL:HG23	3:C:137:ILE:O	2.17	0.45
3:C:319:ALA:HB2	6:F:127:TYR:CZ	2.51	0.45
1:A:191:LYS:HB2	1:A:303:VAL:HG22	1.97	0.45
1:A:324:GLY:O	1:A:327:MET:HG2	2.17	0.45
1:A:68:GLU:O	1:A:70:PRO:HD3	2.17	0.45
1:A:55:VAL:HG22	1:A:58:LEU:CD1	2.45	0.45
3:C:219:TRP:CE2	3:C:227:CYS:HB2	2.52	0.44
4:D:137:GLU:OE2	4:D:158:LYS:HE2	2.17	0.44
6:F:4:THR:CG2	6:F:55:ARG:HE	2.16	0.44
1:A:193:ILE:HG23	1:A:292:PHE:CE2	2.52	0.44
1:A:240:LYS:O	1:A:244:LYS:HG3	2.17	0.44
1:A:311:VAL:O	1:A:314:PRO:HD2	2.18	0.44
1:A:317:LYS:HE3	1:A:364:ASP:HB3	1.99	0.44
2:B:318:GLU:HG3	2:B:344:ILE:HD12	1.99	0.44
3:C:107:ASN:C	3:C:107:ASN:ND2	2.71	0.44
7:G:140:SER:O	7:G:144:VAL:HG23	2.16	0.44
1:A:260:ASN:HB3	1:A:263:SER:HB3	1.99	0.44
1:A:302:GLU:O	1:A:306:ASN:ND2	2.50	0.44
1:A:393:VAL:HG21	1:A:414:PHE:CD2	2.53	0.44
2:B:240:LEU:HB2	10:B:908:HOH:O	2.16	0.44
6:F:145:GLU:HG3	6:F:149:MET:CE	2.47	0.44
6:F:71:ARG:HA	6:F:116:PHE:O	2.17	0.44
7:G:68:SER:OG	7:G:71:VAL:HG12	2.17	0.44
2:B:353:MET:O	2:B:354:VAL:C	2.55	0.44
4:D:223:ILE:HD12	4:D:223:ILE:N	2.32	0.44
1:A:190:ILE:O	1:A:191:LYS:HD3	2.17	0.44
2:B:225:TYR:CZ	2:B:319:ARG:HD2	2.52	0.44
7:G:52:THR:O	7:G:56:GLN:HG3	2.18	0.44
1:A:106:GLU:HG2	1:A:135:ASN:HB2	1.99	0.44
2:B:286:ILE:C	2:B:288:ALA:H	2.21	0.44
2:B:151:LEU:HD22	2:B:300:HIS:CD2	2.53	0.44
3:C:284:ARG:CD	3:C:286:ASP:O	2.65	0.44
1:A:87:ASP:CG	4:D:264:THR:HG22	2.38	0.44
1:A:191:LYS:HB2	1:A:303:VAL:CG2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:346:ASP:HB3	3:C:74:ARG:HH12	1.83	0.44
2:B:350:ARG:HG2	2:B:350:ARG:O	2.16	0.44
1:A:111:LEU:C	1:A:111:LEU:HD23	2.38	0.44
5:E:140:THR:O	5:E:144:LEU:HB2	2.18	0.44
1:A:158:VAL:O	1:A:158:VAL:HG23	2.17	0.43
1:A:359:LYS:NZ	1:A:359:LYS:HB3	2.33	0.43
4:D:230:ARG:NH1	4:D:230:ARG:HB2	2.34	0.43
5:E:153:ASN:OD1	5:E:154:ASP:N	2.52	0.43
5:E:55:PHE:C	5:E:56:LYS:HD2	2.38	0.43
6:F:45:GLU:HB3	7:G:24:PHE:CG	2.53	0.43
3:C:259:VAL:HG12	3:C:332:ILE:CD1	2.49	0.43
3:C:257:SER:OG	3:C:372:VAL:HA	2.19	0.43
5:E:166:ARG:HH11	5:E:166:ARG:HG2	1.83	0.43
6:F:45:GLU:HB3	7:G:24:PHE:CD2	2.53	0.43
1:A:193:ILE:HA	1:A:194:PRO:HD3	1.86	0.43
7:G:62:PRO:HA	7:G:63:PRO:HD3	1.93	0.43
3:C:119:VAL:CG2	3:C:120:ILE:H	2.31	0.43
4:D:220:ILE:CG2	4:D:221:GLY:N	2.81	0.43
1:A:19:LEU:HG	1:A:29:PHE:HB2	2.01	0.43
2:B:161:ASP:N	9:B:902:ANP:O3G	2.51	0.43
3:C:326:LYS:HA	3:C:326:LYS:HD3	1.90	0.43
3:C:267:PRO:HD2	3:C:286:ASP:HB2	2.00	0.42
5:E:95:MET:HG2	5:E:141:GLY:C	2.40	0.42
3:C:254:THR:OG1	3:C:372:VAL:HG13	2.19	0.42
5:E:10:ASP:HB3	5:E:12:ASP:OD1	2.19	0.42
2:B:178:LEU:O	2:B:285:THR:HG23	2.19	0.42
3:C:155:VAL:HG21	3:C:180:PRO:CG	2.47	0.42
4:D:67:LEU:HD13	4:D:120:CYS:O	2.20	0.42
5:E:20:ALA:HB3	5:E:22:LEU:CD2	2.49	0.42
5:E:23:PRO:O	5:E:24:ILE:CG2	2.68	0.42
3:C:183:THR:HG23	3:C:184:PRO:CD	2.49	0.42
7:G:113:GLY:HA3	7:G:125:LEU:HD11	2.01	0.42
1:A:251:LYS:HD2	1:A:251:LYS:N	2.34	0.42
1:A:284:HIS:N	1:A:285:PRO:HD3	2.35	0.42
2:B:175:LEU:CD1	2:B:175:LEU:N	2.80	0.42
4:D:109:ILE:HG23	10:D:312:HOH:O	2.20	0.42
5:E:134:GLN:HA	5:E:137:ARG:CZ	2.49	0.42
1:A:397:LYS:HE2	1:A:401:GLU:OE2	2.20	0.42
3:C:17:ASN:HB3	3:C:60:TRP:CZ2	2.55	0.42
5:E:20:ALA:HB3	5:E:22:LEU:HD21	2.01	0.42
6:F:137:HIS:CE1	6:F:141:GLU:HG3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:164:THR:HB	2:B:182:LEU:HB3	2.01	0.42
4:D:247:PHE:O	4:D:250:TYR:HB3	2.19	0.42
9:B:902:ANP:H1'	10:B:906:HOH:O	2.19	0.42
3:C:93:LEU:HB2	3:C:95:ILE:HG12	2.02	0.42
3:C:360:ARG:HG3	3:C:361:SER:N	2.35	0.42
5:E:24:ILE:HG12	5:E:32:ALA:HB1	2.01	0.42
6:F:14:ALA:O	6:F:17:GLN:HB2	2.19	0.42
2:B:322:LYS:CB	7:G:16:VAL:HG11	2.48	0.42
1:A:295:PRO:HD2	1:A:298:GLU:CD	2.41	0.42
2:B:254:VAL:HG13	2:B:257:GLU:CG	2.49	0.42
3:C:228:LEU:HD23	3:C:228:LEU:C	2.40	0.42
5:E:60:ILE:HD11	5:E:116:ILE:HG21	2.02	0.42
10:F:186:HOH:O	7:G:117:PRO:HG2	2.19	0.42
2:B:323:GLN:NE2	7:G:12:ARG:HA	2.34	0.41
4:D:199:LEU:HD12	4:D:199:LEU:N	2.34	0.41
1:A:116:PRO:HG2	1:A:178:ILE:HD13	2.01	0.41
3:C:183:THR:HG21	3:C:185:TRP:HD1	1.86	0.41
5:E:74:TYR:CE1	5:E:137:ARG:HD2	2.56	0.41
1:A:286:GLU:O	5:E:56:LYS:HD3	2.21	0.41
1:A:53:LYS:O	1:A:56:ASP:OD2	2.38	0.41
6:F:76:VAL:HG23	6:F:113:ASP:HB2	2.03	0.41
3:C:102:VAL:HA	3:C:112:ALA:O	2.21	0.41
3:C:44:GLN:NE2	3:C:47:GLU:OE1	2.52	0.41
2:B:144:LEU:O	2:B:145:THR:C	2.59	0.41
1:A:233:TYR:O	1:A:276:PHE:HA	2.21	0.41
2:B:191:ARG:O	2:B:194:ILE:HB	2.20	0.41
3:C:128:GLU:HG3	3:C:128:GLU:H	1.68	0.41
3:C:27:PRO:HG2	3:C:29:ASN:OD1	2.21	0.41
4:D:63:PHE:CD2	4:D:146:TYR:HA	2.56	0.41
1:A:174:VAL:HG13	1:A:193:ILE:O	2.21	0.41
1:A:311:VAL:HG22	1:A:311:VAL:O	2.20	0.41
2:B:274:GLU:OE1	2:B:275:GLY:N	2.52	0.41
3:C:124:TYR:HB2	3:C:176:VAL:HG11	2.02	0.41
4:D:37:ASP:HB3	4:D:41:VAL:HB	2.00	0.41
2:B:220:LEU:HD11	10:B:916:HOH:O	2.20	0.41
4:D:198:VAL:C	4:D:199:LEU:HD12	2.41	0.41
5:E:166:ARG:NH1	5:E:166:ARG:HG2	2.35	0.41
6:F:20:LEU:HD12	6:F:132:VAL:HG22	2.03	0.41
1:A:176:HIS:CD2	1:A:192:HIS:CD2	3.08	0.41
2:B:156:VAL:HG22	2:B:302:VAL:HG13	2.03	0.41
2:B:322:LYS:HG3	2:B:342:ILE:HD12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:119:VAL:HG21	3:C:136:HIS:HB3	2.03	0.41
3:C:73:ASP:O	3:C:74:ARG:HB2	2.21	0.41
5:E:152:GLN:O	5:E:155:LYS:HB2	2.21	0.41
2:B:229:GLN:HE22	6:F:40:VAL:HB	1.82	0.41
1:A:234:VAL:HG21	1:A:331:PHE:HA	2.02	0.40
1:A:248:ASP:OD2	1:A:251:LYS:HD3	2.21	0.40
1:A:36:ALA:HB2	1:A:66:ALA:HB1	2.03	0.40
2:B:347:PRO:HA	2:B:348:PRO:HD2	1.99	0.40
3:C:29:ASN:O	3:C:54:GLN:HA	2.21	0.40
2:B:330:LEU:C	2:B:332:GLY:H	2.23	0.40
4:D:35:PHE:N	4:D:35:PHE:CD2	2.89	0.40
5:E:60:ILE:CD1	5:E:116:ILE:CG2	2.98	0.40
1:A:263:SER:C	1:A:265:LYS:N	2.74	0.40
2:B:175:LEU:HD22	2:B:178:LEU:HD12	2.03	0.40
2:B:205:ASN:ND2	2:B:207:SER:H	2.19	0.40
5:E:146:GLU:CG	5:E:147:LYS:N	2.84	0.40
6:F:145:GLU:HG3	6:F:149:MET:HE2	2.03	0.40
2:B:158:ASP:HB2	2:B:304:SER:HB3	2.04	0.40
3:C:131:TRP:O	3:C:131:TRP:HE3	2.03	0.40
3:C:249:ALA:HB1	3:C:332:ILE:HG22	2.03	0.40
4:D:212:THR:C	4:D:214:ALA:H	2.25	0.40
4:D:37:ASP:HB2	4:D:43:TYR:CE1	2.57	0.40
7:G:68:SER:C	7:G:70:ALA:H	2.25	0.40
2:B:287:GLN:HG2	2:B:298:TYR:OH	2.22	0.40
3:C:26:CYS:SG	3:C:55:VAL:HB	2.61	0.40
4:D:260:ALA:O	4:D:263:HIS:HB2	2.20	0.40
5:E:105:ILE:HB	5:E:106:PRO:HD2	2.03	0.40
3:C:74:ARG:HH11	6:F:31:GLU:CD	2.25	0.40
6:F:4:THR:C	6:F:7:PRO:HD2	2.42	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	390/418 (93%)	357 (92%)	31 (8%)	2 (0%)	29	47
2	B	216/394 (55%)	189 (88%)	21 (10%)	6 (3%)	5	7
3	C	337/372 (91%)	316 (94%)	19 (6%)	2 (1%)	25	42
4	D	276/300 (92%)	264 (96%)	10 (4%)	2 (1%)	22	39
5	E	172/178 (97%)	154 (90%)	16 (9%)	2 (1%)	13	23
6	F	165/168 (98%)	157 (95%)	8 (5%)	0	100	100
7	G	110/151 (73%)	103 (94%)	5 (4%)	2 (2%)	8	15
All	All	1666/1981 (84%)	1540 (92%)	110 (7%)	16 (1%)	15	27

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	263	SER
5	E	153	ASN
7	G	118	SER
2	B	184	ILE
2	B	278	VAL
4	D	237	ARG
2	B	334	VAL
3	C	50	GLU
4	D	208	GLU
1	A	33	SER
2	B	145	THR
2	B	171	GLU
3	C	179	ARG
5	E	151	PRO
7	G	22	ASN
2	B	347	PRO

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	346/363 (95%)	324 (94%)	22 (6%)	17	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	169/345 (49%)	159 (94%)	10 (6%)	19	34
3	C	288/313 (92%)	276 (96%)	12 (4%)	30	49
4	D	245/264 (93%)	236 (96%)	9 (4%)	34	54
5	E	156/159 (98%)	148 (95%)	8 (5%)	24	41
6	F	154/155 (99%)	150 (97%)	4 (3%)	46	66
7	G	90/124 (73%)	86 (96%)	4 (4%)	28	47
All	All	1448/1723 (84%)	1379 (95%)	69 (5%)	25	44

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	19	LEU
1	A	25	THR
1	A	55	VAL
1	A	68	GLU
1	A	71	THR
1	A	117	LEU
1	A	143	VAL
1	A	191	LYS
1	A	251	LYS
1	A	255	GLN
1	A	289	ASN
1	A	297	SER
1	A	318	ASN
1	A	335	LEU
1	A	340	LYS
1	A	351	GLU
1	A	363	ILE
1	A	372	MET
1	A	373	GLN
1	A	394	CYS
1	A	413	VAL
2	B	153	THR
2	B	180	ARG
2	B	182	LEU
2	B	200	ARG
2	B	220	LEU
2	B	240	LEU
2	B	274	GLU

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Mol	Chain	Res	Type
2	B	281	LEU
2	B	303	LEU
2	B	320	GLU
3	C	21	THR
3	C	107	ASN
3	C	131	TRP
3	C	140	PRO
3	C	179	ARG
3	C	284	ARG
3	C	321	LEU
3	C	346	CYS
3	C	367	LYS
3	C	370	LYS
3	C	371	ILE
3	C	372	VAL
4	D	49	ASN
4	D	84	LEU
4	D	106	LYS
4	D	107	ASP
4	D	116	LEU
4	D	118	ARG
4	D	141	ARG
4	D	171	ASP
4	D	248	ARG
5	E	25	ARG
5	E	35	GLU
5	E	82	LEU
5	E	95	MET
5	E	133	LEU
5	E	144	LEU
5	E	145	CYS
5	E	151	PRO
6	F	101	PHE
6	F	102	PHE
6	F	104	LEU
6	F	165	LEU
7	G	27	GLU
7	G	87	LYS
7	G	90	ASP
7	G	122	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	ASN
1	A	157	GLN
1	A	192	HIS
1	A	205	GLN
1	A	255	GLN
1	A	289	ASN
1	A	306	ASN
1	A	318	ASN
2	B	205	ASN
2	B	229	GLN
2	B	231	GLN
2	B	284	ASN
2	B	300	HIS
2	B	323	GLN
3	C	33	HIS
3	C	65	ASN
3	C	107	ASN
3	C	331	GLN
4	D	21	ASN
4	D	26	ASN
4	D	140	ASN
4	D	231	HIS
4	D	263	HIS
5	E	90	GLN
5	E	134	GLN
5	E	138	GLN
5	E	167	GLN
6	F	129	HIS
6	F	137	HIS
6	F	167	ASN
7	G	96	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 [i](#)

There are no carbohydrates 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$
9	ANP	A	901	8	29,33,33	1.15	1 (3%)	31,52,52	2.17	6 (19%)
9	ANP	B	902	8	29,33,33	1.18	3 (10%)	31,52,52	1.54	3 (9%)

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
9	ANP	A	901	8	-	0/14/38/38	0/3/3/3
9	ANP	B	902	8	-	6/14/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	901	ANP	PB-O2B	-3.16	1.48	1.56
9	B	902	ANP	PB-O2B	-3.01	1.48	1.56
9	B	902	ANP	PB-O3A	2.32	1.62	1.59
9	B	902	ANP	O4'-C1'	2.28	1.44	1.41

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	901	ANP	O2B-PB-O1B	7.33	125.30	109.92
9	A	901	ANP	O3A-PB-N3B	-4.81	93.23	106.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	902	ANP	N3-C2-N1	-4.48	121.68	128.68
9	A	901	ANP	N3-C2-N1	-4.46	121.71	128.68
9	B	902	ANP	O2B-PB-O1B	4.11	118.54	109.92
9	A	901	ANP	O4'-C1'-C2'	-3.94	101.17	106.93
9	A	901	ANP	PA-O3A-PB	-3.16	121.50	132.62
9	B	902	ANP	PA-O3A-PB	-2.88	122.48	132.62
9	A	901	ANP	O2B-PB-O3A	-2.31	96.94	104.64

There are no chirality outliers.

All (6) torsion outliers are listed below:

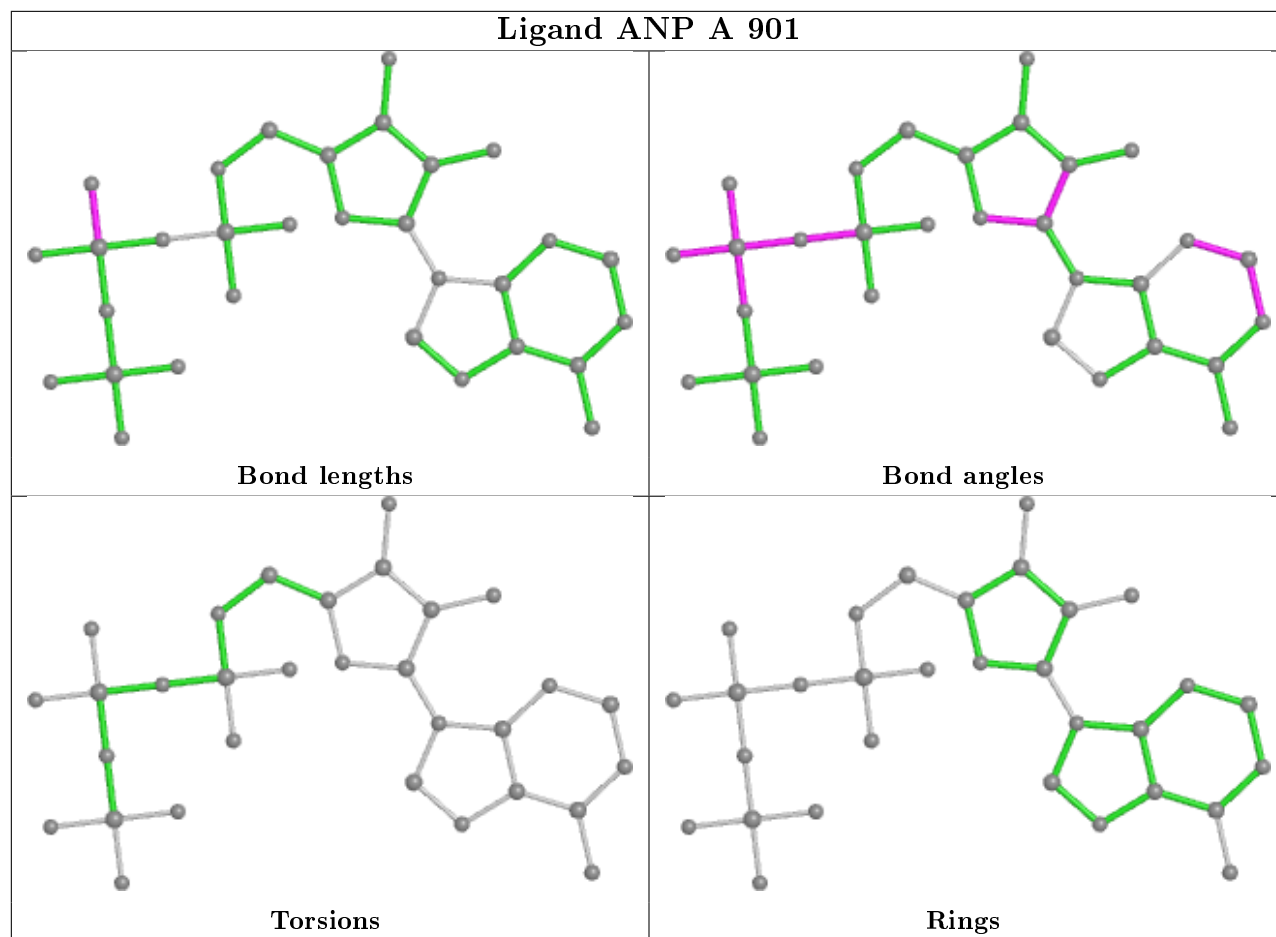
Mol	Chain	Res	Type	Atoms
9	B	902	ANP	PG-N3B-PB-O1B
9	B	902	ANP	C5'-O5'-PA-O1A
9	B	902	ANP	C5'-O5'-PA-O2A
9	B	902	ANP	PB-O3A-PA-O1A
9	B	902	ANP	PB-O3A-PA-O2A
9	B	902	ANP	C5'-O5'-PA-O3A

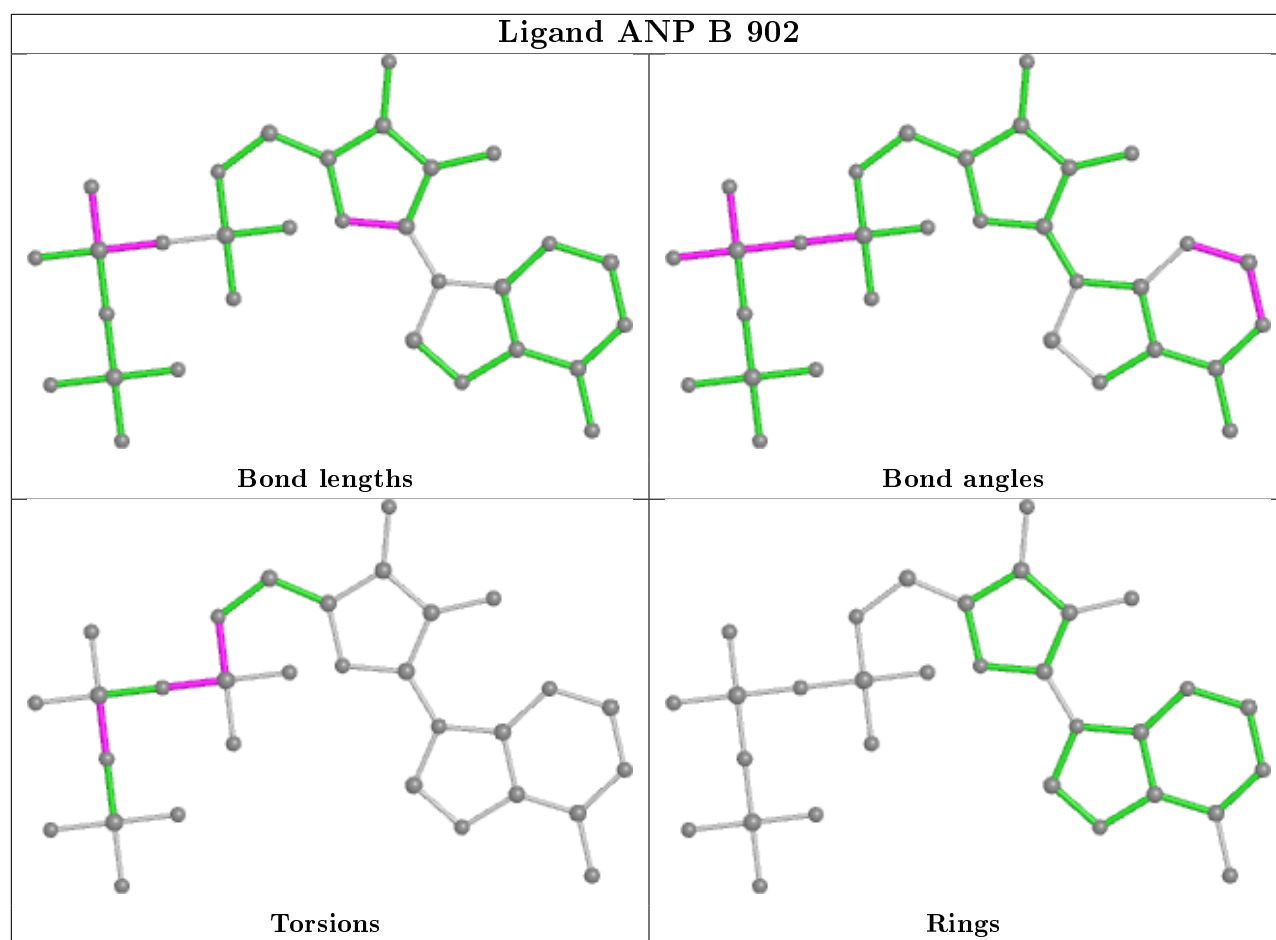
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	901	ANP	3	0
9	B	902	ANP	6	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	398/418 (95%)	0.06	21 (5%) 26 31	19, 45, 89, 127	0
2	B	222/394 (56%)	0.16	12 (5%) 25 31	27, 52, 82, 97	0
3	C	341/372 (91%)	-0.16	6 (1%) 68 76	24, 36, 65, 100	0
4	D	280/300 (93%)	-0.20	3 (1%) 80 86	23, 38, 70, 86	0
5	E	174/178 (97%)	0.41	15 (8%) 10 13	38, 59, 85, 108	0
6	F	167/168 (99%)	-0.33	1 (0%) 89 92	23, 35, 50, 87	0
7	G	116/151 (76%)	0.42	7 (6%) 21 26	30, 67, 86, 94	0
All	All	1698/1981 (85%)	0.01	65 (3%) 40 48	19, 43, 82, 127	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	359	LYS	7.4
5	E	154	ASP	6.5
1	A	351	GLU	5.3
1	A	360	PRO	4.8
2	B	356	LEU	4.6
7	G	120	ASN	4.3
2	B	355	PHE	4.2
1	A	350	SER	4.2
1	A	51	VAL	4.0
1	A	159	GLY	4.0
2	B	178	LEU	4.0
1	A	354	SER	3.9
6	F	2	THR	3.6
3	C	367	LYS	3.6
1	A	39	GLU	3.5
2	B	181	ARG	3.4
1	A	157	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
5	E	153	ASN	3.2
5	E	175	GLY	3.2
1	A	352	GLU	3.2
1	A	70	PRO	3.2
1	A	348	LYS	3.1
2	B	171	GLU	3.1
1	A	262	ILE	3.1
1	A	414	PHE	3.1
7	G	88	ALA	3.1
3	C	319	ALA	3.0
2	B	176	PRO	3.0
3	C	362	LEU	3.0
3	C	368	ASP	2.9
5	E	86	ASN	2.9
1	A	160	GLU	2.8
7	G	121	SER	2.8
4	D	209	LEU	2.8
7	G	84	ILE	2.7
3	C	360	ARG	2.7
2	B	276	VAL	2.7
4	D	207	LEU	2.7
2	B	274	GLU	2.7
5	E	155	LYS	2.6
7	G	11	PHE	2.6
5	E	149	PHE	2.6
5	E	35	GLU	2.5
5	E	89	SER	2.5
5	E	152	GLN	2.5
2	B	175	LEU	2.4
5	E	150	ASP	2.4
2	B	330	LEU	2.3
5	E	146	GLU	2.3
1	A	158	VAL	2.3
5	E	151	PRO	2.3
7	G	61	ASN	2.3
5	E	93	LYS	2.3
1	A	349	LEU	2.3
3	C	127	GLN	2.3
1	A	52	MET	2.2
1	A	256	TYR	2.2
7	G	90	ASP	2.2
2	B	275	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	261	ALA	2.2
5	E	27	GLN	2.1
5	E	90	GLN	2.1
4	D	215	ALA	2.1
1	A	415	GLY	2.1
2	B	354	VAL	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 carbohydrates in this entry.

6.4 Ligands [i](#)

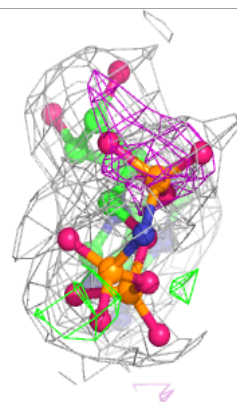
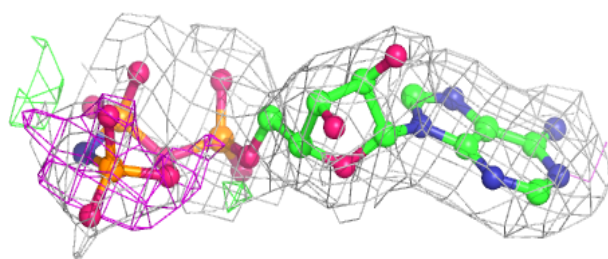
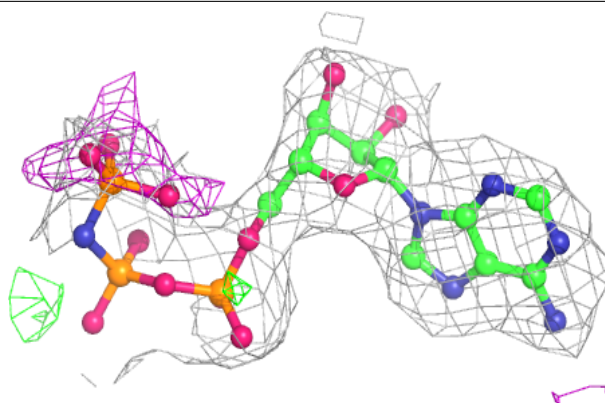
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	CA	B	802	1/1	0.91	0.13	70,70,70,70	0
9	ANP	B	902	31/31	0.92	0.14	43,54,86,87	0
8	CA	A	801	1/1	0.93	0.26	41,41,41,41	0
9	ANP	A	901	31/31	0.95	0.18	46,50,54,57	0

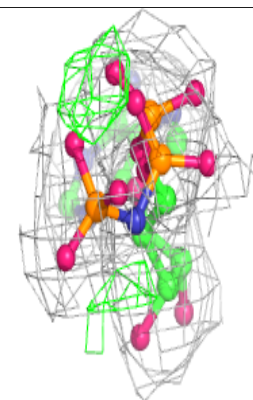
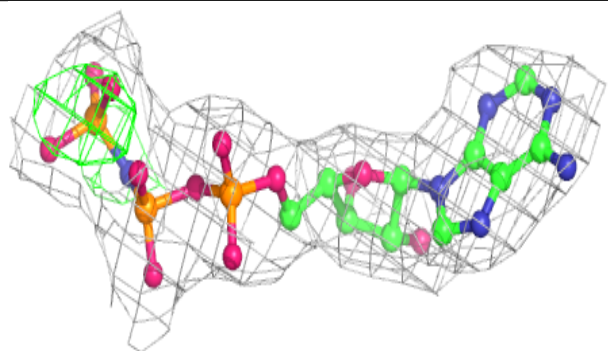
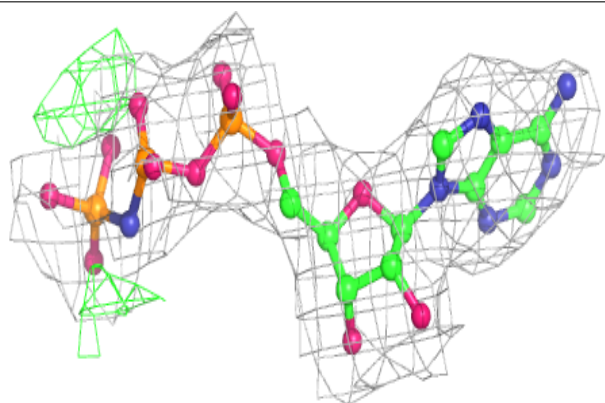
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ANP B 902:

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

**Electron density around ANP A 901:**

$2mF_o-DF_c$ (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.