



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:35 AM UTC

PDB ID : 1G6O / pdb_00001g6o
Title : CRYSTAL STRUCTURE OF THE HELICOBACTER PYLORI ATPASE, HP0525, IN COMPLEX WITH ADP
Authors : Yeo, H.J.; Savvides, S.N.; Herr, A.B.; Lanka, E.; Waksman, G.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2000-11-07
Resolution : 2.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

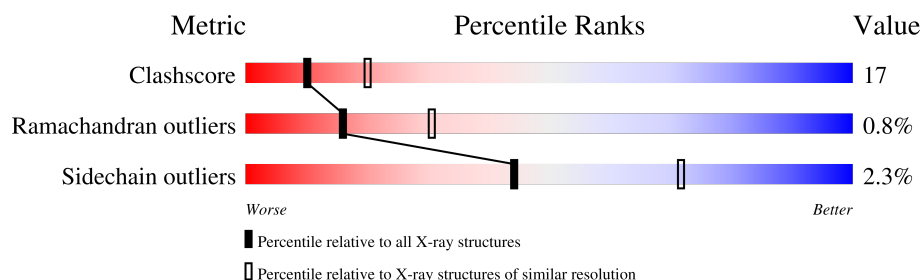
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	330	
1	B	330	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PEG	B	501	-	X	-	-
3	PEG	B	502	-	X	-	-
3	PEG	B	503	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5482 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.

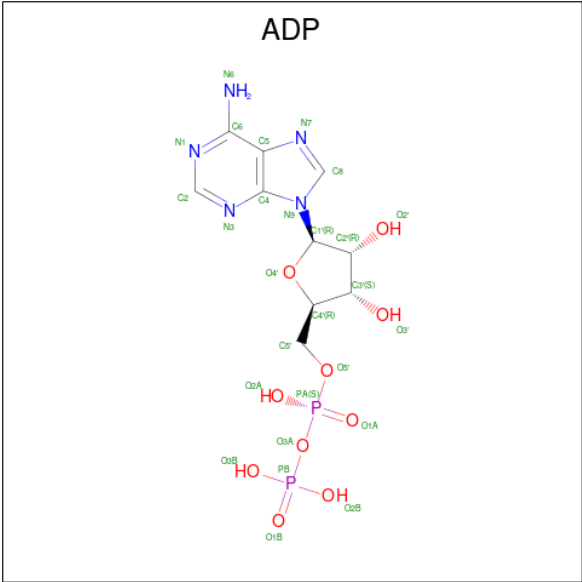
- Molecule 1 is a protein called CAG-ALPHA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	Se	0	0	0
			2554	1620	436	484	8	6			
1	B	323	Total	C	N	O	S	Se	0	0	0
			2551	1617	436	484	8	6			

There are 12 discrepancies between the modelled and reference sequences:

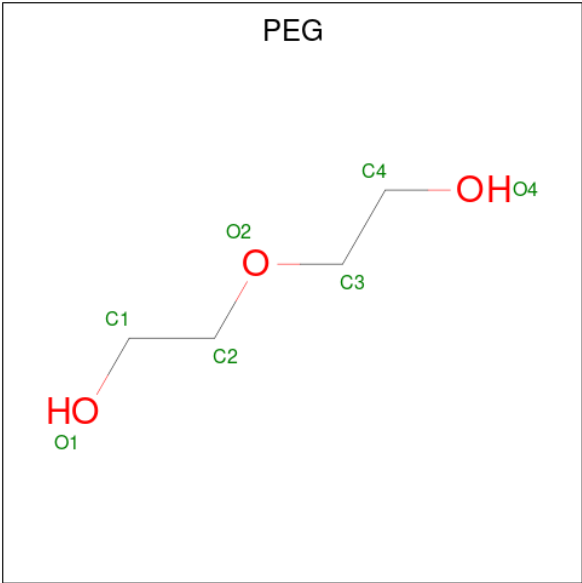
Chain	Residue	Modelled	Actual	Comment	Reference
A	42	MSE	MET	modified residue	GB 8843975
A	82	MSE	MET	modified residue	GB 8843975
A	192	MSE	MET	modified residue	GB 8843975
A	239	MSE	MET	modified residue	GB 8843975
A	287	MSE	MET	modified residue	GB 8843975
A	312	MSE	MET	modified residue	GB 8843975
B	42	MSE	MET	modified residue	GB 8843975
B	82	MSE	MET	modified residue	GB 8843975
B	192	MSE	MET	modified residue	GB 8843975
B	239	MSE	MET	modified residue	GB 8843975
B	287	MSE	MET	modified residue	GB 8843975
B	312	MSE	MET	modified residue	GB 8843975

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



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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			6	4	2		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			3	2	1		

- Molecule 4 is water.

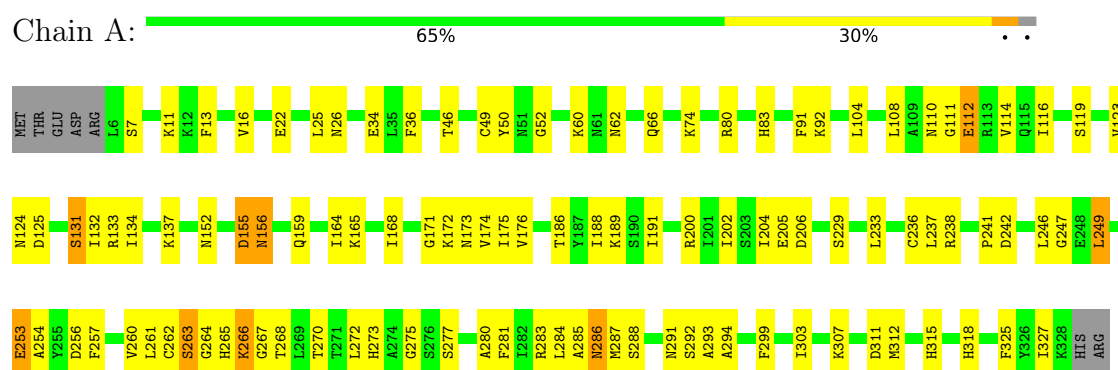
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	155	Total	O	0	0
			155	155		
4	B	140	Total	O	0	0
			140	140		

3 Residue-property plots [i](#)

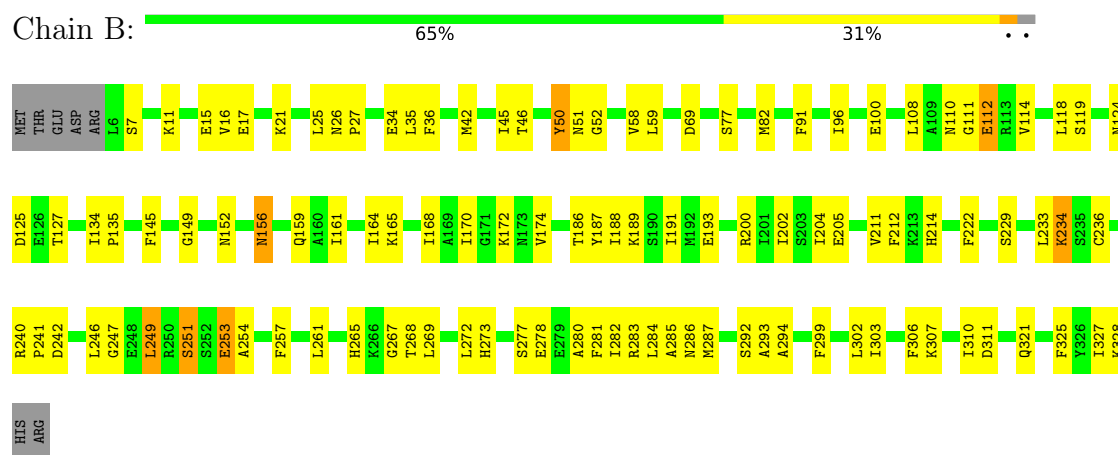
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

• Molecule 1: CAG-ALPHA



• Molecule 1: CAG-ALPHA



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	112.62Å 112.62Å 234.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.50	Depositor
% Data completeness (in resolution range)	88.8 (30.00-2.50)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5482	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, ADP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.71	1/2597 (0.0%)	1.08	15/3491 (0.4%)
1	B	0.68	0/2594	1.07	16/3487 (0.5%)
All	All	0.70	1/5191 (0.0%)	1.08	31/6978 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	46	THR	CA-CB	5.52	1.62	1.54

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	242	ASP	N-CA-C	-8.04	103.42	113.55
1	A	263	SER	N-CA-C	-7.70	104.30	112.93
1	B	124	ASN	CA-C-N	7.53	132.12	120.82
1	B	124	ASN	C-N-CA	7.53	132.12	120.82
1	A	266	LYS	N-CA-C	7.28	120.25	110.35
1	A	242	ASP	N-CA-C	-6.72	104.72	113.12
1	A	265	HIS	N-CA-C	6.57	119.77	110.50
1	A	16	VAL	N-CA-C	6.39	117.18	110.72
1	A	111	GLY	N-CA-C	6.31	121.17	113.79
1	A	83	HIS	N-CA-C	-6.20	104.21	110.97
1	B	119	SER	N-CA-C	-6.07	101.70	110.08
1	A	123	VAL	N-CA-C	5.99	117.45	110.62
1	A	206	ASP	N-CA-C	-5.94	105.69	113.12
1	B	265	HIS	N-CA-C	5.86	118.55	110.35
1	A	119	SER	N-CA-C	-5.80	102.46	109.83
1	B	251	SER	N-CA-C	5.76	123.07	110.80
1	B	36	PHE	N-CA-C	5.74	119.60	112.59
1	A	152	ASN	N-CA-C	5.60	118.15	111.71

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	124	ASN	N-CA-C	5.48	115.77	108.38
1	B	77	SER	N-CA-C	-5.46	102.40	110.48
1	B	27	PRO	N-CA-C	-5.23	106.59	113.65
1	B	127	THR	N-CA-C	5.17	117.14	107.99
1	B	26	ASN	CA-C-N	-5.17	113.60	119.28
1	B	26	ASN	C-N-CA	-5.17	113.60	119.28
1	B	50	TYR	N-CA-C	-5.13	100.53	108.90
1	A	124	ASN	CA-C-N	5.12	131.32	121.54
1	A	124	ASN	C-N-CA	5.12	131.32	121.54
1	B	16	VAL	N-CA-C	5.12	115.89	110.72
1	B	111	GLY	N-CA-C	5.11	119.77	113.79
1	A	36	PHE	N-CA-C	5.02	120.02	112.94
1	B	15	GLU	N-CA-C	5.02	117.13	111.11

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	2554	0	2518	91	0
1	B	2551	0	2509	89	0
2	A	27	0	12	1	0
2	B	27	0	12	0	0
3	B	28	0	18	1	0
4	A	155	0	0	5	0
4	B	140	0	0	6	0
All	All	5482	0	5069	176	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 (176) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:MSE:HE1	1:B:58:VAL:HG21	1.40	1.02
1:B:172:LYS:HE3	1:B:311:ASP:HB3	1.50	0.94
1:A:318:HIS:HB2	4:A:1107:HOH:O	1.66	0.94
1:A:299:PHE:HE2	1:A:303:ILE:HD11	1.42	0.85
1:A:229:SER:HB3	1:A:253:GLU:HG2	1.61	0.83
1:A:299:PHE:CE2	1:A:303:ILE:HD11	2.15	0.82
1:B:168:ILE:HD12	1:B:191:ILE:CG2	2.10	0.80
1:A:257:PHE:HE1	1:A:268:THR:HG21	1.47	0.80
1:B:168:ILE:HD12	1:B:191:ILE:HG21	1.65	0.78
1:A:156:ASN:HD22	1:A:156:ASN:H	1.30	0.76
1:A:22:GLU:HG2	1:A:26:ASN:ND2	2.01	0.75
1:A:25:LEU:HD13	1:A:91:PHE:CE1	2.21	0.75
1:A:202:ILE:HD12	1:A:241:PRO:HB3	1.69	0.74
1:B:229:SER:HB3	1:B:253:GLU:HG2	1.69	0.73
1:B:108:LEU:HB2	1:B:112:GLU:HG2	1.70	0.73
1:B:42:MSE:HE1	1:B:58:VAL:CG2	2.19	0.71
1:B:42:MSE:CE	1:B:58:VAL:HG11	2.20	0.71
1:A:156:ASN:HB2	1:A:159:GLN:HB3	1.72	0.71
1:B:156:ASN:HB2	1:B:159:GLN:HB3	1.72	0.70
1:B:321:GLN:HB3	4:B:1218:HOH:O	1.92	0.69
1:B:306:PHE:O	1:B:310:ILE:HD13	1.92	0.68
1:A:156:ASN:H	1:A:156:ASN:ND2	1.91	0.68
1:A:283:ARG:HG2	1:A:283:ARG:HH11	1.58	0.68
1:B:108:LEU:HD12	1:B:112:GLU:HG3	1.76	0.67
1:B:168:ILE:CD1	1:B:191:ILE:HG21	2.24	0.67
1:A:200:ARG:HH12	1:B:51:ASN:HD21	1.44	0.65
1:A:137:LYS:HG2	4:A:1038:HOH:O	1.98	0.64
1:A:156:ASN:HB2	1:A:159:GLN:HE21	1.61	0.64
1:B:42:MSE:HE1	1:B:58:VAL:HG11	1.80	0.62
1:B:34:GLU:HG2	1:B:35:LEU:HD12	1.82	0.62
1:B:69:ASP:HB2	3:B:500:PEG:H12	1.82	0.62
1:A:164:ILE:HG23	1:A:165:LYS:N	2.13	0.62
1:A:238:ARG:HH11	1:A:238:ARG:HA	1.65	0.61
1:B:284:LEU:HD13	1:B:306:PHE:CD2	2.35	0.61
1:B:278:GLU:O	1:B:282:ILE:HG13	2.01	0.61
1:A:22:GLU:CG	1:A:26:ASN:ND2	2.63	0.61
1:A:238:ARG:HA	1:A:238:ARG:NH1	2.15	0.61
1:B:204:ILE:HB	1:B:246:LEU:HD12	1.82	0.60
1:B:164:ILE:CG2	1:B:165:LYS:N	2.65	0.60
1:B:193:GLU:HG2	1:B:214:HIS:NE2	2.17	0.59
1:A:281:PHE:CE2	1:A:307:LYS:HA	2.38	0.59
1:A:229:SER:CB	1:A:253:GLU:HG2	2.32	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:PHE:CE2	1:B:303:ILE:HD11	2.38	0.59
1:B:174:VAL:HG13	1:B:269:LEU:HD13	1.85	0.59
1:B:42:MSE:HE3	1:B:58:VAL:HG11	1.85	0.58
1:A:74:LYS:O	1:A:80:ARG:HG3	2.03	0.58
1:B:164:ILE:HG23	1:B:165:LYS:N	2.18	0.58
1:A:108:LEU:HB2	1:A:112:GLU:HG2	1.86	0.57
1:B:202:ILE:HD12	1:B:241:PRO:HB3	1.86	0.57
1:B:229:SER:CB	1:B:253:GLU:HG2	2.35	0.57
1:A:172:LYS:HD3	1:A:312:MSE:HE3	1.86	0.56
1:B:188:ILE:HG22	1:B:269:LEU:HD11	1.87	0.56
1:A:283:ARG:HG2	1:A:283:ARG:NH1	2.21	0.56
1:B:254:ALA:O	1:B:257:PHE:HB3	2.04	0.56
1:A:156:ASN:HD22	1:A:156:ASN:N	1.96	0.56
1:B:7:SER:O	1:B:11:LYS:HG3	2.05	0.56
1:A:164:ILE:CG2	1:A:165:LYS:N	2.68	0.56
1:A:60:LYS:HB2	1:A:62:ASN:OD1	2.06	0.56
1:A:325:PHE:CE2	1:A:327:ILE:HD11	2.41	0.56
1:A:108:LEU:CD2	1:A:114:VAL:HG21	2.35	0.56
1:A:273:HIS:O	1:A:283:ARG:NH1	2.39	0.56
1:A:229:SER:HB3	1:A:253:GLU:CG	2.35	0.55
1:B:156:ASN:HD22	1:B:156:ASN:H	1.54	0.55
1:A:7:SER:O	1:A:11:LYS:HG3	2.06	0.55
1:B:281:PHE:CE2	1:B:307:LYS:HA	2.41	0.55
1:A:50:TYR:CE2	1:A:52:GLY:HA2	2.42	0.54
1:A:22:GLU:CG	1:A:26:ASN:HD21	2.21	0.54
1:B:168:ILE:CD1	1:B:191:ILE:CG2	2.81	0.54
1:B:240:ARG:O	1:B:240:ARG:HG3	2.07	0.54
1:A:125:ASP:HB3	4:A:1153:HOH:O	2.06	0.54
1:A:34:GLU:OE2	1:A:80:ARG:HD3	2.07	0.53
1:B:50:TYR:CE2	1:B:52:GLY:HA2	2.44	0.53
1:B:25:LEU:HD13	1:B:91:PHE:CE1	2.44	0.53
1:A:284:LEU:O	1:A:288:SER:HB2	2.08	0.53
1:B:200:ARG:NH1	1:B:202:ILE:HD11	2.24	0.53
1:B:292:SER:C	1:B:294:ALA:H	2.16	0.52
1:A:277:SER:O	1:A:280:ALA:HB3	2.09	0.52
1:B:112:GLU:CD	1:B:112:GLU:H	2.18	0.52
1:A:13:PHE:HZ	1:B:82:MSE:HG2	1.73	0.52
1:A:292:SER:C	1:A:294:ALA:H	2.17	0.52
1:B:233:LEU:O	1:B:236:CYS:HB3	2.11	0.51
1:A:133:ARG:HD3	4:A:1023:HOH:O	2.10	0.51
1:A:263:SER:HB2	1:B:273:HIS:HB3	1.91	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:LYS:HE2	1:A:66:GLN:NE2	2.26	0.50
1:B:257:PHE:HE1	1:B:268:THR:HG21	1.74	0.50
1:A:104:LEU:HD23	1:A:116:ILE:HD12	1.94	0.50
1:B:51:ASN:ND2	4:B:1045:HOH:O	2.42	0.50
1:A:186:THR:OG1	2:A:401:ADP:O2A	2.26	0.49
1:B:156:ASN:HB2	1:B:159:GLN:HE21	1.78	0.49
1:B:257:PHE:CE2	1:B:261:LEU:HD11	2.47	0.49
1:A:175:ILE:HG12	1:A:270:THR:HG22	1.94	0.49
1:B:328:LYS:HA	4:B:1185:HOH:O	2.12	0.49
1:A:254:ALA:O	1:A:257:PHE:HB3	2.12	0.49
1:B:17:GLU:O	1:B:21:LYS:HG2	2.13	0.49
1:B:110:ASN:HB2	1:B:112:GLU:OE2	2.13	0.48
1:A:168:ILE:CD1	1:A:191:ILE:HD12	2.43	0.48
1:A:172:LYS:HE3	1:A:311:ASP:HB3	1.95	0.48
1:B:125:ASP:HB2	4:B:1013:HOH:O	2.11	0.48
1:B:156:ASN:H	1:B:156:ASN:ND2	2.10	0.48
1:A:257:PHE:CE1	1:A:268:THR:HG21	2.37	0.48
1:B:286:ASN:OD1	1:B:299:PHE:HZ	1.97	0.48
1:A:112:GLU:H	1:A:112:GLU:CD	2.20	0.48
1:A:286:ASN:ND2	1:A:299:PHE:HZ	2.12	0.48
1:A:188:ILE:HG13	1:A:189:LYS:N	2.29	0.48
1:B:164:ILE:CG2	1:B:165:LYS:H	2.26	0.48
1:A:205:GLU:HG3	4:A:1190:HOH:O	2.13	0.47
1:A:249:LEU:HD22	1:A:270:THR:OG1	2.15	0.47
1:A:257:PHE:CZ	1:A:261:LEU:HD21	2.49	0.47
1:A:275:GLY:O	1:A:315:HIS:NE2	2.47	0.47
1:A:91:PHE:CE2	1:A:92:LYS:HE2	2.49	0.47
1:A:168:ILE:HD11	1:A:191:ILE:HD12	1.97	0.47
1:A:292:SER:O	1:A:294:ALA:N	2.44	0.47
1:B:108:LEU:CD2	1:B:114:VAL:HG21	2.45	0.47
1:B:302:LEU:O	1:B:306:PHE:HD1	1.98	0.47
1:A:74:LYS:O	1:A:74:LYS:HG2	2.13	0.47
1:B:42:MSE:HE3	1:B:45:ILE:HG21	1.96	0.47
1:B:283:ARG:HG2	1:B:283:ARG:HH11	1.81	0.46
1:B:292:SER:O	1:B:294:ALA:N	2.42	0.46
1:B:325:PHE:CD2	1:B:327:ILE:HD11	2.51	0.46
1:A:108:LEU:HD21	1:A:114:VAL:HG21	1.98	0.46
1:A:325:PHE:CD2	1:A:327:ILE:HD11	2.51	0.46
1:A:155:ASP:OD2	1:A:155:ASP:N	2.43	0.46
1:A:110:ASN:HB2	1:A:112:GLU:OE2	2.16	0.45
1:B:156:ASN:HB2	1:B:159:GLN:NE2	2.31	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:ILE:HD12	1:A:134:ILE:N	2.32	0.45
1:B:170:ILE:O	1:B:170:ILE:HG22	2.16	0.45
1:A:49:CYS:SG	1:A:131:SER:HB3	2.56	0.45
1:B:172:LYS:HB3	1:B:311:ASP:CB	2.47	0.45
1:A:174:VAL:HG22	1:A:312:MSE:HG3	1.99	0.45
1:A:246:LEU:HD23	1:A:249:LEU:HD22	1.99	0.45
1:B:191:ILE:O	1:B:191:ILE:HG22	2.17	0.45
1:A:171:GLY:HA3	1:A:266:LYS:HD2	2.00	0.44
1:A:237:LEU:HD11	1:A:260:VAL:HG22	1.99	0.44
1:B:42:MSE:HE1	1:B:58:VAL:CG1	2.46	0.44
1:A:286:ASN:ND2	1:A:299:PHE:CZ	2.85	0.44
1:B:145:PHE:HZ	1:B:186:THR:HG22	1.83	0.44
1:A:285:ALA:HB2	1:A:303:ILE:HG12	1.99	0.44
1:A:256:ASP:O	1:A:260:VAL:HG23	2.18	0.43
1:B:156:ASN:HD22	1:B:156:ASN:N	2.14	0.43
1:B:188:ILE:HG13	1:B:189:LYS:N	2.33	0.43
1:B:134:ILE:HA	1:B:135:PRO:HD3	1.86	0.43
1:B:187:TYR:O	1:B:191:ILE:HG12	2.18	0.43
1:B:234:LYS:HG3	4:B:1161:HOH:O	2.17	0.43
1:B:249:LEU:HD13	1:B:249:LEU:HA	1.88	0.43
1:A:283:ARG:O	1:A:287:MSE:HG3	2.19	0.43
1:A:108:LEU:HD11	1:A:132:ILE:HD13	2.00	0.42
1:A:156:ASN:CB	1:A:159:GLN:HE21	2.29	0.42
1:A:164:ILE:CG2	1:A:165:LYS:H	2.31	0.42
1:A:262:CYS:C	1:A:264:GLY:H	2.25	0.42
1:B:303:ILE:HG22	1:B:307:LYS:HE3	2.01	0.42
1:B:246:LEU:O	1:B:247:GLY:C	2.62	0.42
1:B:188:ILE:CG2	1:B:269:LEU:HD11	2.49	0.42
1:B:211:VAL:HG12	1:B:212:PHE:N	2.35	0.42
1:A:291:ASN:OD1	1:A:291:ASN:C	2.63	0.42
1:B:205:GLU:O	1:B:222:PHE:HA	2.20	0.42
1:A:246:LEU:O	1:A:247:GLY:C	2.63	0.41
1:B:96:ILE:HG12	1:B:118:LEU:HD11	2.02	0.41
1:B:100:GLU:HB2	4:B:1265:HOH:O	2.20	0.41
1:B:285:ALA:HB2	1:B:303:ILE:HG12	2.01	0.41
1:B:272:LEU:HD13	1:B:287:MSE:HE1	2.02	0.41
1:A:325:PHE:HE2	1:A:327:ILE:HD11	1.84	0.41
1:B:161:ILE:O	1:B:161:ILE:HG22	2.21	0.41
1:B:172:LYS:HE3	1:B:311:ASP:CB	2.35	0.41
1:B:149:GLY:HA3	1:B:152:ASN:ND2	2.35	0.41
1:B:325:PHE:CE2	1:B:327:ILE:HD11	2.55	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:VAL:HG21	1:A:188:ILE:HG22	2.02	0.41
1:A:233:LEU:O	1:A:236:CYS:HB3	2.21	0.41
1:A:286:ASN:HD21	1:A:299:PHE:HZ	1.68	0.41
1:A:202:ILE:CD1	1:A:241:PRO:HB3	2.46	0.41
1:A:204:ILE:HB	1:A:246:LEU:HD12	2.02	0.40
1:B:277:SER:O	1:B:280:ALA:HB3	2.20	0.40
1:B:108:LEU:HD12	1:B:112:GLU:CG	2.49	0.40
1:A:200:ARG:HH12	1:B:51:ASN:ND2	2.15	0.40
1:A:246:LEU:HD23	1:A:249:LEU:CD2	2.52	0.40
1:A:173:ASN:ND2	1:A:261:LEU:HD12	2.37	0.40
1:B:46:THR:CG2	1:B:59:LEU:HD23	2.52	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	321/330 (97%)	301 (94%)	18 (6%)	2 (1%)	21	38
1	B	321/330 (97%)	295 (92%)	23 (7%)	3 (1%)	14	27
All	All	642/660 (97%)	596 (93%)	41 (6%)	5 (1%)	16	31

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	GLY
1	A	293	ALA
1	B	267	GLY
1	B	293	ALA
1	B	251	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	280/287 (98%)	272 (97%)	8 (3%)	37	65
1	B	279/287 (97%)	274 (98%)	5 (2%)	51	77
All	All	559/574 (97%)	546 (98%)	13 (2%)	44	72

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	GLU
1	A	131	SER
1	A	155	ASP
1	A	156	ASN
1	A	249	LEU
1	A	253	GLU
1	A	272	LEU
1	A	286	ASN
1	B	112	GLU
1	B	156	ASN
1	B	234	LYS
1	B	249	LEU
1	B	253	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	51	ASN
1	A	66	GLN
1	A	156	ASN
1	A	159	GLN
1	A	226	ASN
1	A	286	ASN
1	A	321	GLN
1	B	51	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	61	ASN
1	B	156	ASN
1	B	159	GLN
1	B	215	HIS
1	B	265	HIS
1	B	273	HIS
1	B	286	ASN

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 ⓘ

7 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	PEG	B	503	3	5,5,6	15.75	4 (80%)	4,4,5	11.99	4 (100%)
2	ADP	B	402	-	28,29,29	2.07	6 (21%)	43,45,45	1.89	9 (20%)
3	PEG	B	501	3	5,5,6	10.62	2 (40%)	4,4,5	15.26	4 (100%)
3	PEG	B	502	3	5,5,6	16.99	4 (80%)	4,4,5	11.70	4 (100%)
3	PEG	B	504	-	2,2,6	0.60	0	1,1,5	0.66	0
3	PEG	B	500	3	6,6,6	1.24	1 (16%)	5,5,5	0.87	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ADP	A	401	-	28,29,29	2.08	7 (25%)	43,45,45	1.92	7 (16%)

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	PEG	B	503	3	-	3/3/3/4	-
2	ADP	B	402	-	-	3/16/32/32	0/3/3/3
3	PEG	B	501	3	-	2/3/3/4	-
3	PEG	B	502	3	-	3/3/3/4	-
3	PEG	B	500	3	-	0/4/4/4	-
2	ADP	A	401	-	-	3/16/32/32	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	PEG	O4-C4	30.26	2.96	1.42
3	B	503	PEG	O4-C4	27.46	2.82	1.42
3	B	501	PEG	C3-C4	19.11	2.50	1.49
3	B	503	PEG	C3-C4	18.31	2.46	1.49
3	B	502	PEG	C3-C4	18.29	2.46	1.49
3	B	501	PEG	O2-C2	13.96	2.58	1.42
3	B	502	PEG	O2-C2	13.56	2.55	1.42
3	B	503	PEG	O2-C2	11.99	2.42	1.42
2	B	402	ADP	C5-N7	-5.97	1.28	1.39
2	A	401	ADP	C5-N7	-5.60	1.28	1.39
2	A	401	ADP	PA-O3A	-4.31	1.54	1.59
2	A	401	ADP	C2-N3	3.57	1.40	1.33
2	B	402	ADP	C2-N1	3.50	1.40	1.33
2	B	402	ADP	C4-N9	-3.49	1.30	1.37
2	B	402	ADP	C5-C4	3.47	1.45	1.39
2	A	401	ADP	C2-N1	3.42	1.40	1.33
2	B	402	ADP	C1'-N9	3.23	1.55	1.46
2	B	402	ADP	C2-N3	3.11	1.39	1.33
2	A	401	ADP	C5-C4	3.10	1.44	1.39
3	B	502	PEG	O2-C3	2.87	1.54	1.42
2	A	401	ADP	C4-N9	-2.86	1.31	1.37
3	B	503	PEG	O2-C3	2.46	1.52	1.42
2	A	401	ADP	C8-N9	-2.10	1.34	1.37

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	500	PEG	O2-C3	2.02	1.50	1.42

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	PEG	C2-O2-C3	-24.65	26.55	113.06
3	B	503	PEG	O2-C3-C4	-17.33	33.74	110.11
3	B	501	PEG	O2-C3-C4	-17.15	34.52	110.11
3	B	502	PEG	O2-C3-C4	-17.15	34.55	110.11
3	B	502	PEG	O4-C4-C3	-13.99	29.44	111.82
3	B	503	PEG	O4-C4-C3	-13.63	31.60	111.82
3	B	503	PEG	C2-O2-C3	9.05	144.81	113.06
3	B	502	PEG	C2-O2-C3	7.26	138.53	113.06
2	A	401	ADP	N3-C2-N1	-6.32	119.01	128.58
2	B	402	ADP	N3-C2-N1	-6.24	119.14	128.58
2	A	401	ADP	C5-C4-N3	-5.08	119.73	126.72
2	B	402	ADP	C5-C4-N3	-5.02	119.81	126.72
3	B	501	PEG	O4-C4-C3	-4.96	82.63	111.82
2	B	402	ADP	C2-N3-C4	4.23	122.17	111.83
2	A	401	ADP	C2-N3-C4	4.19	122.06	111.83
2	A	401	ADP	N3-C4-N9	3.90	133.80	127.17
2	B	402	ADP	N3-C4-N9	3.56	133.22	127.17
2	B	402	ADP	C5-N7-C8	3.13	108.37	103.45
2	B	402	ADP	C4-C5-N7	-3.07	107.07	110.58
2	A	401	ADP	C5-N7-C8	2.94	108.07	103.45
2	A	401	ADP	C4-C5-N7	-2.85	107.32	110.58
3	B	503	PEG	O2-C2-C1	2.72	138.76	111.79
2	A	401	ADP	O3B-PB-O3A	2.60	113.34	104.64
2	B	402	ADP	O3B-PB-O3A	2.35	112.52	104.64
3	B	501	PEG	O2-C2-C1	2.31	134.71	111.79
3	B	502	PEG	O2-C2-C1	2.26	134.15	111.79
2	B	402	ADP	C2'-C3'-C4'	2.24	106.94	102.61
2	B	402	ADP	N9-C8-N7	-2.09	110.96	113.94

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	402	ADP	PA-O3A-PB-O3B
3	B	503	PEG	C4-C3-O2-C2
3	B	502	PEG	C4-C3-O2-C2
3	B	502	PEG	C1-C2-O2-C3

Continued on next page...

Continued from previous page...

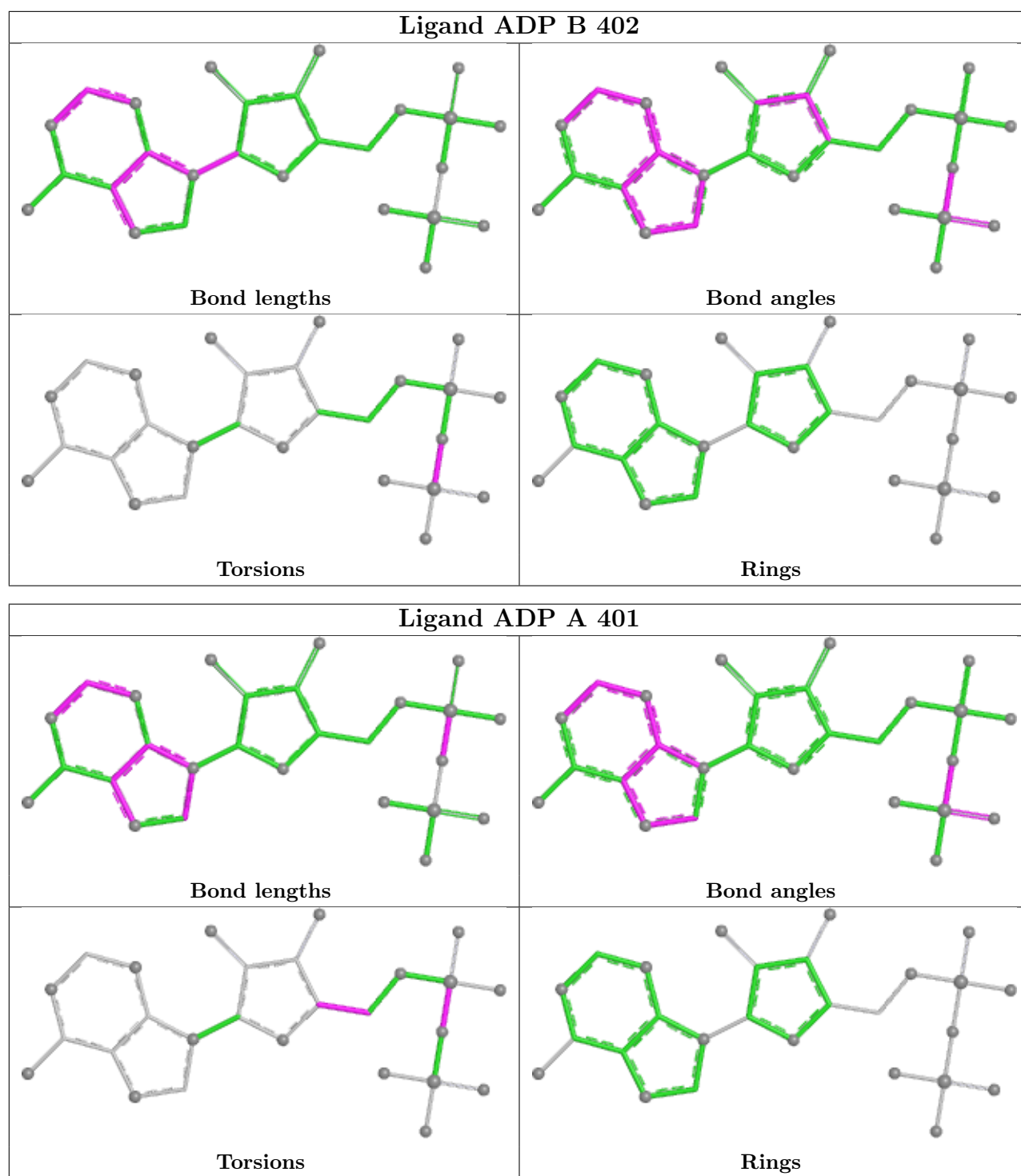
Mol	Chain	Res	Type	Atoms
3	B	503	PEG	C1-C2-O2-C3
3	B	502	PEG	O2-C3-C4-O4
2	A	401	ADP	O4'-C4'-C5'-O5'
3	B	501	PEG	C1-C2-O2-C3
3	B	503	PEG	O2-C3-C4-O4
3	B	501	PEG	O2-C3-C4-O4
2	B	402	ADP	PA-O3A-PB-O1B
2	B	402	ADP	PA-O3A-PB-O2B
2	A	401	ADP	C3'-C4'-C5'-O5'
2	A	401	ADP	PB-O3A-PA-O2A

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	500	PEG	1	0
2	A	401	ADP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.