



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 07:31 pm BST

PDB ID : 2IEH
Title : Crystal structure of human kinesin Eg5 in complex with (R)-mon97, a new monastrol-based inhibitor that binds as (R)-enantiomer
Authors : Garcia-Saez, I.; Kozielski, F.
Deposited on : 2006-09-19
Resolution : 2.70 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

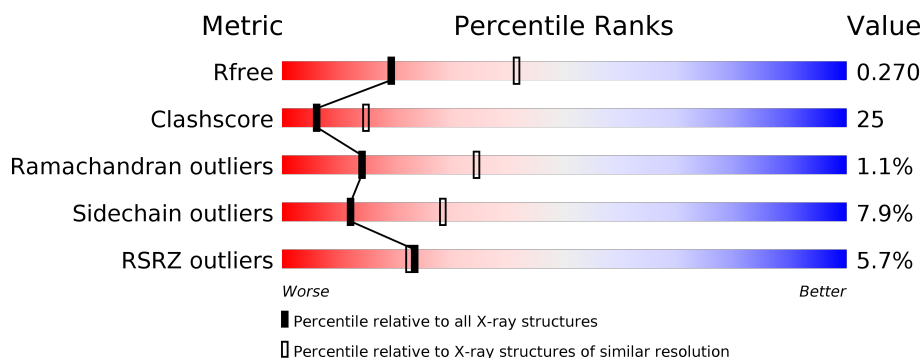
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	367	<div> <div>6%</div> <div> <div></div> <div>51%</div> <div>35%</div> <div>• • 9%</div> </div> </div>
1	B	367	<div> <div>5%</div> <div> <div></div> <div>59%</div> <div>28%</div> <div>• 9%</div> </div> </div>

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
7	PG4	A	606	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5607 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 Kinesin-like protein KIF11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	2	0
			2644	1655	462	517	10			
1	B	335	Total	C	N	O	S	0	0	0
			2630	1647	458	515	10			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

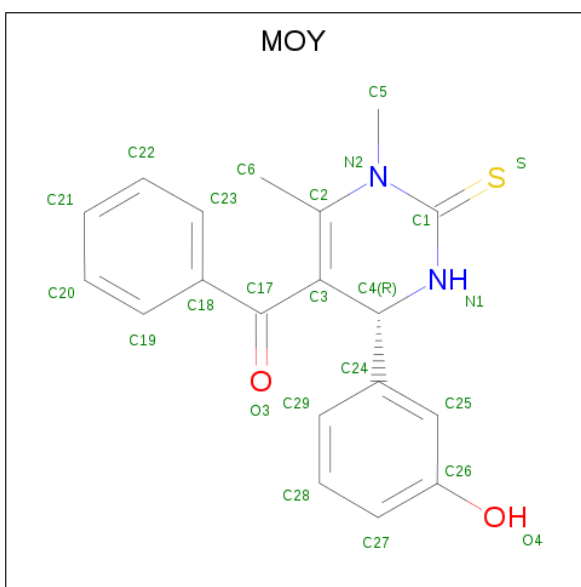
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

- Molecule 6 is [(4R)-4-(3-HYDROXYPHENYL)-1,6-DIMETHYL-2-THIOXO-1,2,3,4-TETRAHYDROPYRIMIDIN-5-YL](PHENYL)METHANONE (three-letter code: MOY) (formula: C₁₉H₁₈N₂O₂S).



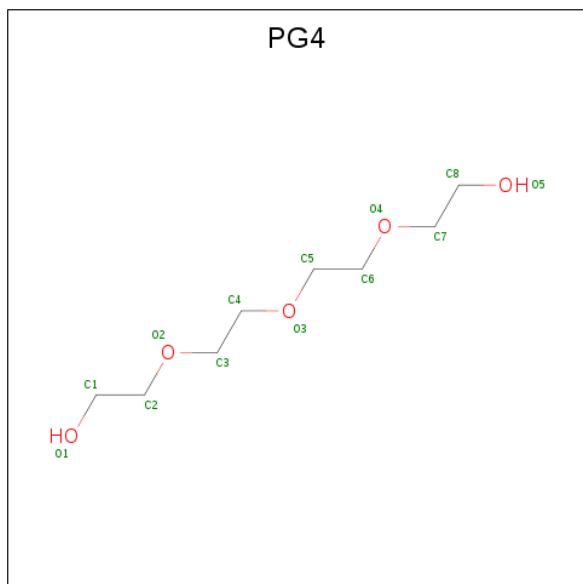
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	S	0	0
			24	19	2	2	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	S	0	0
			24	19	2	2	1		

- Molecule 7 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			13	8	5		

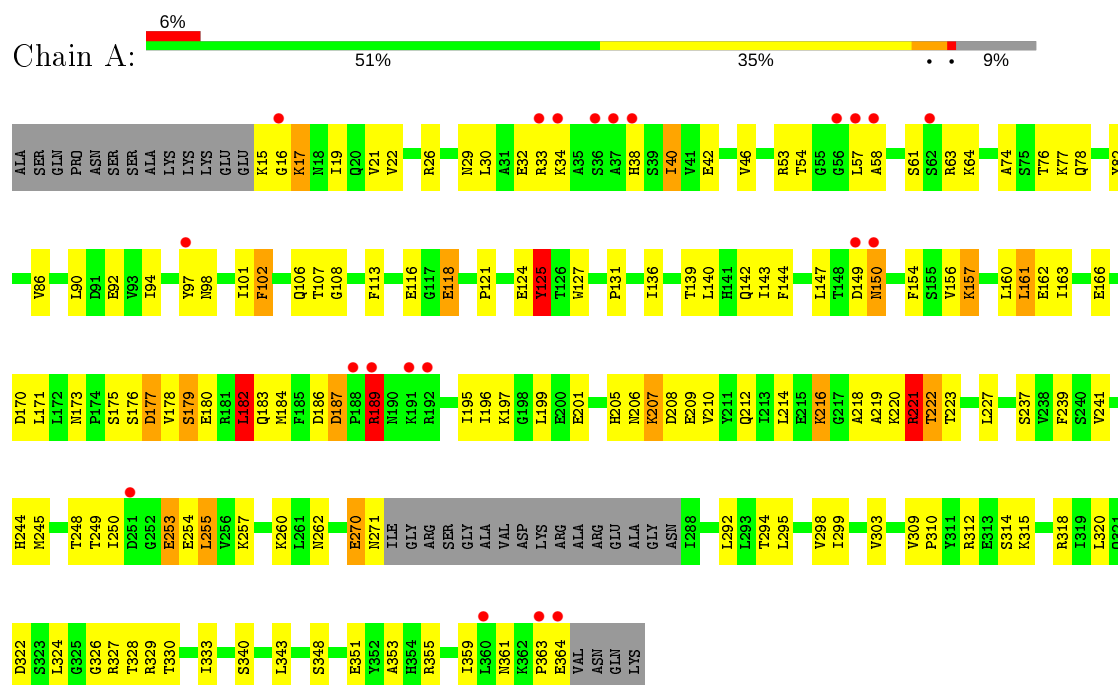
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	114	Total	O	0	0
			114	114		
8	B	100	Total	O	0	0
			100	100		

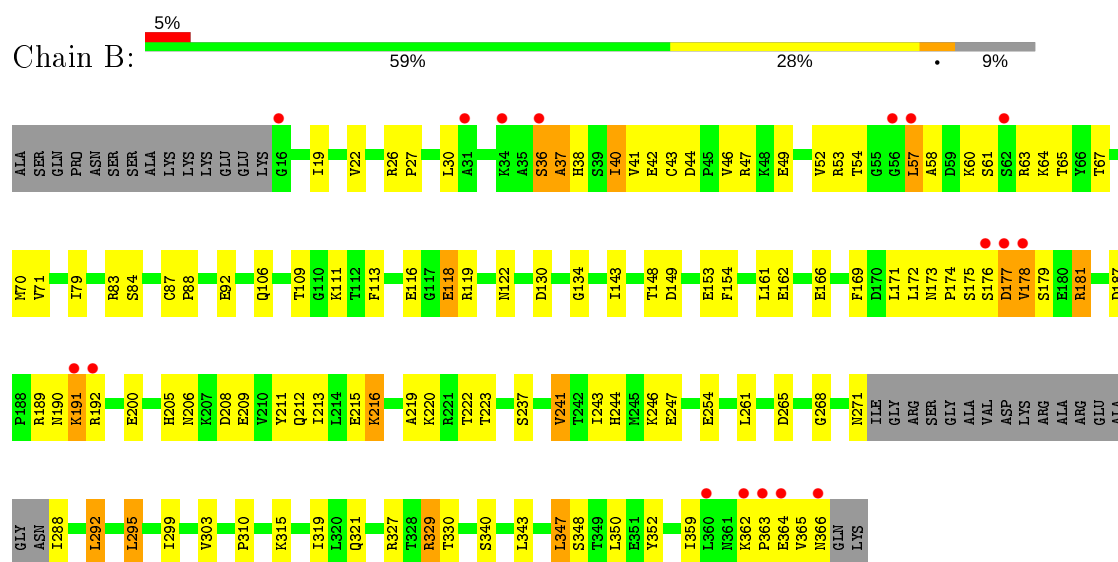
3 Residue-property plots

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: Kinesin-like protein KIF11



• Molecule 1: Kinesin-like protein KIF11



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.37Å 79.90Å 159.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.70 44.33 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.4 (25.00-2.70) 99.1 (44.33-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.61Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.234 , 0.281 0.230 , 0.270	Depositor DCC
R_{free} test set	2613 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5607	wwPDB-VP
Average B, all atoms (Å ²)	27.0	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 43.32 % 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 1.7940e-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: MG, ADP, CL, K, MOY, PG4

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	0.63	7/2683 (0.3%)	1.00	14/3625 (0.4%)
1	B	0.51	2/2669 (0.1%)	0.77	2/3609 (0.1%)
All	All	0.57	9/5352 (0.2%)	0.89	16/7234 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	102	PHE	CB-CG	-8.16	1.37	1.51
1	B	36	SER	C-O	7.78	1.38	1.23
1	B	37	ALA	CA-CB	-7.51	1.36	1.52
1	A	189	ARG	CG-CD	7.21	1.70	1.51
1	A	16	GLY	CA-C	6.29	1.61	1.51
1	A	102	PHE	CD1-CE1	-6.05	1.27	1.39
1	A	189	ARG	CB-CG	6.04	1.68	1.52
1	A	157	LYS	CB-CG	-5.90	1.36	1.52
1	A	16	GLY	N-CA	5.53	1.54	1.46

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	ARG	NE-CZ-NH2	-13.70	113.45	120.30
1	A	221	ARG	NE-CZ-NH1	13.50	127.05	120.30
1	A	189	ARG	NE-CZ-NH1	12.31	126.46	120.30
1	A	189	ARG	NE-CZ-NH2	-12.18	114.21	120.30
1	A	221	ARG	CG-CD-NE	11.05	135.00	111.80
1	A	221	ARG	CD-NE-CZ	9.11	136.36	123.60
1	B	37	ALA	C-N-CA	8.86	143.84	121.70
1	A	15	LYS	CA-C-N	-7.53	101.13	116.20
1	A	182	LEU	CA-CB-CG	7.21	131.90	115.30
1	A	227	LEU	CB-CG-CD1	-6.42	100.08	111.00
1	A	16	GLY	N-CA-C	-6.06	97.96	113.10
1	A	189	ARG	CD-NE-CZ	5.70	131.57	123.60
1	A	102	PHE	CB-CA-C	-5.63	99.13	110.40
1	B	36	SER	CA-C-N	-5.52	105.05	117.20
1	A	15	LYS	N-CA-C	5.43	125.65	111.00
1	A	125	TYR	CB-CG-CD2	-5.18	117.89	121.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	125	TYR	Sidechain
1	B	36	SER	Mainchain

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	2644	0	2671	147	0
1	B	2630	0	2655	117	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	27	0	12	2	0
5	B	27	0	12	2	0
6	A	24	0	18	1	0
6	B	24	0	18	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	13	0	18	19	0
8	A	114	0	0	20	0
8	B	100	0	0	12	0
All	All	5607	0	5404	266	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 25.

All (266) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:ILE:HG21	8:A:720:HOH:O	1.57	1.03
1:A:171:LEU:HD22	8:A:714:HOH:O	1.59	1.00
1:B:83:ARG:HD2	8:B:705:HOH:O	1.62	0.98
1:B:49:GLU:HG2	1:B:67:THR:HG22	1.44	0.95
1:A:183:GLN:O	8:A:717:HOH:O	1.84	0.95
1:A:77:LYS:HD2	7:A:606:PG4:H41	1.47	0.95
1:B:191:LYS:HG3	1:B:192:ARG:HG3	1.46	0.94
1:A:197:LYS:N	8:A:717:HOH:O	2.01	0.93
1:A:221:ARG:CD	8:A:714:HOH:O	2.16	0.93
1:A:57:LEU:O	1:A:61:SER:HB3	1.68	0.92
1:A:328:THR:O	1:A:361:ASN:ND2	2.02	0.92
1:A:77:LYS:HB3	7:A:606:PG4:H81	1.54	0.90
1:B:30:LEU:HD12	8:B:679:HOH:O	1.73	0.89
1:A:189:ARG:HD3	8:A:630:HOH:O	1.77	0.83
1:A:30:LEU:HD13	1:A:33[A]:ARG:NH2	1.94	0.82
1:B:153:GLU:HG2	8:B:680:HOH:O	1.78	0.82
1:A:78:GLN:H	7:A:606:PG4:H52	1.45	0.81
1:B:57:LEU:O	1:B:61:SER:HB3	1.81	0.81
1:A:30:LEU:CD1	1:A:33[A]:ARG:NH2	2.44	0.80
1:A:221:ARG:NE	8:A:714:HOH:O	2.12	0.80
1:B:52:VAL:HG11	1:B:347:LEU:CD1	2.11	0.80
1:A:221:ARG:HD3	8:A:714:HOH:O	1.81	0.77
1:B:87:CYS:HB3	1:B:88:PRO:HD3	1.64	0.77
1:A:97:TYR:HE2	1:A:364:GLU:HA	1.50	0.77
1:B:46:VAL:HG23	1:B:47:ARG:HG2	1.66	0.76
1:A:162:GLU:OE1	8:A:714:HOH:O	2.04	0.75
1:A:218:ALA:O	1:A:221:ARG:HB3	1.87	0.75
1:A:54:THR:HG21	1:A:64:LYS:HG3	1.68	0.75
1:A:166:GLU:O	1:A:315:LYS:HE3	1.86	0.74
1:A:177:ASP:OD1	1:A:179:SER:OG	2.05	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:ASN:HD22	1:B:200:GLU:HG2	1.53	0.73
1:B:220:LYS:NZ	8:B:698:HOH:O	2.20	0.73
1:B:177:ASP:HB3	1:B:179:SER:OG	1.89	0.73
1:A:40:ILE:HG13	1:A:40:ILE:O	1.88	0.73
1:A:78:GLN:H	7:A:606:PG4:C5	2.02	0.72
1:B:57:LEU:CD2	1:B:58:ALA:H	2.02	0.72
1:A:139:THR:O	1:A:143:ILE:HG13	1.90	0.72
1:A:351:GLU:HG2	8:A:718:HOH:O	1.90	0.72
1:A:329:ARG:HG3	8:A:703:HOH:O	1.89	0.71
1:B:177:ASP:C	1:B:179:SER:H	1.93	0.71
1:A:131:PRO:HB2	7:A:606:PG4:C7	2.22	0.70
1:A:131:PRO:HB2	7:A:606:PG4:H71	1.73	0.69
1:A:147:LEU:HD23	1:A:154:PHE:CG	2.28	0.69
1:B:237:SER:HB3	1:B:265:ASP:HB3	1.75	0.69
1:A:299:ILE:O	1:A:303:VAL:HG22	1.93	0.69
1:A:249:THR:HB	1:A:253[A]:GLU:HG3	1.73	0.68
1:A:299:ILE:HG23	1:A:359:ILE:HD11	1.75	0.68
1:A:180:GLU:HG3	1:A:180:GLU:O	1.92	0.68
1:A:38:HIS:CD2	8:A:642:HOH:O	2.46	0.68
1:B:189:ARG:HG2	1:B:189:ARG:HH11	1.58	0.67
1:B:241:VAL:HG13	1:B:261:LEU:HB3	1.76	0.67
1:B:53:ARG:HB2	1:B:63:ARG:NH1	2.09	0.67
1:B:271:ASN:O	1:B:292:LEU:HD12	1.95	0.66
1:B:189:ARG:NH1	1:B:189:ARG:HG2	2.11	0.66
1:A:107:THR:OG1	1:A:270:GLU:OE2	2.14	0.66
1:B:52:VAL:HG11	1:B:347:LEU:HD11	1.77	0.66
1:B:119:ARG:NH2	1:B:130:ASP:OD2	2.29	0.66
1:A:40:ILE:HG21	1:A:340:SER:HA	1.76	0.66
5:A:600:ADP:N1	7:A:606:PG4:H31	2.11	0.65
1:B:327:ARG:HD3	1:B:364:GLU:HB3	1.77	0.65
1:B:61:SER:OG	8:B:696:HOH:O	2.07	0.65
1:A:33[A]:ARG:NH2	8:A:664:HOH:O	2.27	0.64
1:A:320:LEU:O	1:A:324:LEU:HD12	1.98	0.64
1:B:54:THR:HG21	1:B:64:LYS:HG2	1.79	0.64
1:B:22:VAL:CG1	1:B:70:MET:HB2	2.28	0.63
1:B:327:ARG:HG3	1:B:327:ARG:HH11	1.63	0.63
1:B:22:VAL:HG12	1:B:70:MET:HB2	1.80	0.63
1:A:249:THR:HG22	1:A:250:ILE:N	2.13	0.63
1:A:40:ILE:HD12	1:A:343:LEU:HD13	1.81	0.62
1:A:74:ALA:O	7:A:606:PG4:H11	2.00	0.62
1:B:54:THR:CG2	1:B:64:LYS:HG2	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:HD11	1:A:33[A]:ARG:NH2	2.15	0.61
1:A:249:THR:HG22	1:A:250:ILE:H	1.65	0.61
1:B:177:ASP:CB	1:B:179:SER:OG	2.49	0.61
1:A:351:GLU:OE1	8:A:713:HOH:O	2.16	0.61
1:A:38:HIS:H	1:A:38:HIS:CD2	2.18	0.61
1:A:351:GLU:CG	8:A:718:HOH:O	2.48	0.60
1:A:210:VAL:O	1:A:214:LEU:HG	2.01	0.60
1:B:327:ARG:O	1:B:363:PRO:HA	2.00	0.60
1:A:254:GLU:C	1:A:255:LEU:HD23	2.22	0.60
1:A:30:LEU:HD13	1:A:33[A]:ARG:HH21	1.66	0.60
1:B:172:LEU:O	1:B:174:PRO:HD3	2.02	0.59
1:A:249:THR:CB	1:A:253[A]:GLU:HG3	2.32	0.59
1:A:17:LYS:HD3	8:A:652:HOH:O	2.02	0.58
1:B:181:ARG:HG3	1:B:181:ARG:NH1	2.17	0.58
1:A:113:PHE:CE1	1:A:118:GLU:HG3	2.39	0.58
1:A:249:THR:HB	1:A:253[A]:GLU:CG	2.33	0.58
1:B:57:LEU:HD23	1:B:58:ALA:H	1.68	0.58
1:B:70:MET:CE	1:B:84:SER:HB3	2.34	0.58
1:A:107:THR:CB	1:A:270:GLU:OE2	2.51	0.57
1:A:74:ALA:O	7:A:606:PG4:C1	2.53	0.57
1:A:184:MET:HG3	1:A:196:ILE:HD13	1.85	0.57
1:B:143:ILE:HD13	1:B:243:ILE:HD11	1.87	0.57
1:B:244:HIS:CD2	8:B:665:HOH:O	2.58	0.57
1:A:162:GLU:HG2	1:A:237:SER:HA	1.87	0.56
1:A:40:ILE:CD1	1:A:343:LEU:HD13	2.36	0.56
1:B:113:PHE:HD1	1:B:118:GLU:OE2	1.87	0.56
1:A:78:GLN:N	7:A:606:PG4:H52	2.18	0.56
1:A:147:LEU:HB3	1:A:154:PHE:CE2	2.41	0.55
1:A:106:GLN:HG3	1:A:270:GLU:HG2	1.88	0.55
1:B:192:ARG:HD3	1:B:321:GLN:HE21	1.72	0.55
1:A:113:PHE:HE1	1:A:118:GLU:HG3	1.71	0.55
1:A:77:LYS:HB3	7:A:606:PG4:H52	1.88	0.55
1:B:92:GLU:HG3	1:B:329:ARG:NH1	2.22	0.55
1:A:77:LYS:CB	7:A:606:PG4:H81	2.34	0.55
1:A:178:VAL:HG12	1:A:220:LYS:HE2	1.88	0.55
1:A:216:LYS:NZ	8:A:685:HOH:O	2.31	0.55
1:B:173:ASN:ND2	1:B:200:GLU:HG2	2.20	0.55
1:B:271:ASN:O	1:B:292:LEU:CD1	2.55	0.54
1:B:40:ILE:HD12	1:B:340:SER:CB	2.37	0.54
1:B:205:HIS:HB2	1:B:209:GLU:OE2	2.07	0.54
1:B:327:ARG:HG3	1:B:327:ARG:NH1	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:GLU:CG	1:B:67:THR:HG22	2.30	0.54
1:A:219:ALA:O	1:A:222:THR:HG23	2.06	0.54
1:A:157:LYS:HE3	1:A:201:GLU:OE1	2.07	0.54
5:B:603:ADP:H3'	8:B:702:HOH:O	2.07	0.53
1:A:170:ASP:HB2	1:A:182:LEU:HD21	1.89	0.53
1:B:244:HIS:HD2	8:B:665:HOH:O	1.92	0.53
1:B:57:LEU:HD22	1:B:58:ALA:H	1.70	0.53
1:A:150:ASN:HB3	8:A:682:HOH:O	2.07	0.53
1:A:54:THR:HG21	1:A:64:LYS:CG	2.38	0.52
1:A:53:ARG:NH1	1:A:58:ALA:HB2	2.24	0.52
1:A:303:VAL:HG11	1:A:355:ARG:O	2.09	0.52
1:A:131:PRO:HB2	7:A:606:PG4:H72	1.91	0.52
1:A:260:LYS:HE2	1:A:262:ASN:HD21	1.73	0.52
1:A:42:GLU:HG3	1:A:63:ARG:NH2	2.24	0.52
1:B:154:PHE:HA	1:B:244:HIS:O	2.09	0.52
1:B:187:ASP:OD1	1:B:189:ARG:N	2.41	0.52
1:B:46:VAL:HG23	1:B:47:ARG:N	2.25	0.52
1:A:162:GLU:CD	8:A:714:HOH:O	2.47	0.52
1:B:52:VAL:HG11	1:B:347:LEU:HD13	1.88	0.52
1:B:111:LYS:HZ3	1:B:268:GLY:HA2	1.75	0.51
1:B:70:MET:HE2	1:B:84:SER:HB3	1.93	0.51
1:B:211:TYR:O	1:B:215:GLU:HG3	2.10	0.51
1:B:92:GLU:CG	1:B:329:ARG:NH1	2.74	0.51
1:B:365:VAL:CG1	1:B:366:ASN:N	2.74	0.51
1:A:19:ILE:HD12	1:A:330:THR:HB	1.93	0.50
1:B:166:GLU:OE1	1:B:315:LYS:HG2	2.12	0.50
1:B:83:ARG:HG2	1:B:83:ARG:HH11	1.76	0.50
1:B:111:LYS:HE3	5:B:603:ADP:O2B	2.11	0.50
1:B:220:LYS:O	1:B:223:THR:N	2.41	0.49
1:A:205:HIS:HB2	1:A:209:GLU:OE2	2.11	0.49
1:B:212:GLN:HG2	1:B:216:LYS:NZ	2.28	0.49
1:A:178:VAL:CG1	1:A:220:LYS:HE2	2.42	0.49
1:A:124:GLU:HG2	1:A:125:TYR:CD1	2.47	0.49
1:A:322:ASP:OD1	1:A:327:ARG:N	2.34	0.49
1:B:44:ASP:OD2	1:B:47:ARG:HG3	2.12	0.49
1:A:160:LEU:HD13	1:A:239:PHE:HD2	1.76	0.49
1:A:255:LEU:HD23	1:A:255:LEU:N	2.28	0.49
1:A:363:PRO:O	1:A:364:GLU:HG3	2.13	0.49
1:B:92:GLU:HA	1:B:92:GLU:OE1	2.11	0.49
1:B:173:ASN:ND2	1:B:200:GLU:CG	2.75	0.49
1:A:22:VAL:CG2	1:A:333:ILE:HG23	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:ILE:HG23	1:B:359:ILE:HD11	1.95	0.48
1:A:127:TRP:HB2	6:A:602:MOY:H27	1.95	0.48
1:B:143:ILE:CD1	1:B:243:ILE:HD11	2.43	0.48
1:A:176:SER:HB3	1:A:180:GLU:OE1	2.12	0.48
1:B:206:ASN:OD1	1:B:208:ASP:HB2	2.14	0.48
1:A:177:ASP:O	1:A:180:GLU:HG2	2.14	0.48
1:B:111:LYS:NZ	1:B:268:GLY:HA2	2.28	0.48
1:A:207:LYS:HG2	1:A:208:ASP:N	2.29	0.48
1:B:41:VAL:HG22	1:B:52:VAL:HG23	1.96	0.48
1:B:43:CYS:HB3	1:B:71:VAL:HG12	1.96	0.48
1:A:124:GLU:HB3	1:A:125:TYR:HD1	1.79	0.48
1:A:77:LYS:HD2	7:A:606:PG4:H61	1.96	0.48
1:A:94:ILE:O	1:A:257:LYS:HE2	2.14	0.48
1:A:19:ILE:CG1	1:A:359:ILE:HB	2.44	0.47
1:B:181:ARG:HH11	1:B:181:ARG:HG3	1.78	0.47
1:A:322:ASP:O	1:A:326:GLY:HA3	2.14	0.47
1:A:312:ARG:HA	1:A:318:ARG:CG	2.43	0.47
1:A:223:THR:O	1:A:223:THR:HG22	2.14	0.47
1:B:111:LYS:HZ3	1:B:268:GLY:CA	2.26	0.47
1:B:38:HIS:ND1	1:B:38:HIS:N	2.62	0.47
1:A:29:ASN:OD1	1:A:32:GLU:HG3	2.14	0.47
1:B:19:ILE:HD12	1:B:359:ILE:HB	1.97	0.47
1:B:176:SER:CB	8:B:704:HOH:O	2.62	0.47
1:B:365:VAL:HG12	1:B:366:ASN:N	2.29	0.47
1:A:78:GLN:HG2	7:A:606:PG4:H51	1.96	0.47
1:A:97:TYR:CE2	1:A:364:GLU:HA	2.40	0.47
1:B:219:ALA:O	1:B:222:THR:HB	2.14	0.46
1:A:248:THR:HA	1:A:253[B]:GLU:O	2.14	0.46
1:B:49:GLU:OE1	1:B:65:THR:HG21	2.15	0.46
1:A:312:ARG:HG2	1:A:312:ARG:NH1	2.31	0.46
1:A:186:ASP:OD2	1:A:312:ARG:NH2	2.49	0.46
1:B:176:SER:HB3	8:B:704:HOH:O	2.14	0.46
1:A:108:GLY:H	5:A:600:ADP:PB	2.38	0.46
1:A:312:ARG:HG2	1:A:312:ARG:HH11	1.81	0.45
1:B:191:LYS:HB2	1:B:191:LYS:HE2	1.81	0.45
1:A:90:LEU:HD13	1:A:139:THR:HG23	1.97	0.45
1:A:161:LEU:HD23	1:A:199:LEU:HD13	1.98	0.45
1:B:119:ARG:HD2	6:B:605:MOY:C26	2.46	0.45
1:B:299:ILE:O	1:B:303:VAL:HG23	2.17	0.45
1:A:248:THR:HA	1:A:253[A]:GLU:O	2.16	0.45
1:A:248:THR:HG23	1:A:248:THR:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:79:ILE:HG22	8:B:628:HOH:O	2.16	0.45
1:A:22:VAL:HG22	1:A:333:ILE:HA	1.98	0.45
1:B:205:HIS:ND1	1:B:209:GLU:OE2	2.45	0.45
1:A:187:ASP:HA	1:A:195:ILE:CD1	2.47	0.45
1:A:294:THR:OG1	1:A:314:SER:HB3	2.17	0.45
1:B:187:ASP:OD1	1:B:187:ASP:C	2.54	0.45
1:B:192:ARG:HB3	1:B:321:GLN:NE2	2.31	0.45
1:A:82:TYR:OH	1:A:142:GLN:HG3	2.18	0.44
1:A:171:LEU:HD22	1:A:221:ARG:HD3	1.99	0.44
1:B:181:ARG:HH11	1:B:181:ARG:CG	2.31	0.44
1:A:245:MET:HE3	1:A:257:LYS:HB2	1.99	0.44
1:B:190:ASN:HB3	1:B:191:LYS:H	1.68	0.44
1:B:177:ASP:C	1:B:179:SER:N	2.65	0.44
1:A:245:MET:CE	1:A:257:LYS:HB2	2.48	0.44
1:A:184:MET:HG3	1:A:196:ILE:CD1	2.48	0.44
1:A:82:TYR:CD2	1:A:86:VAL:HB	2.53	0.44
1:A:40:ILE:HG21	1:A:340:SER:CA	2.46	0.43
1:A:249:THR:CG2	1:A:250:ILE:N	2.80	0.43
1:A:154:PHE:HA	1:A:244:HIS:O	2.19	0.43
1:A:22:VAL:HG22	1:A:333:ILE:HG12	2.00	0.43
1:A:144:PHE:CZ	1:A:156:VAL:HG21	2.54	0.43
1:B:148:THR:HG23	1:B:149:ASP:N	2.33	0.43
1:A:98:ASN:O	1:A:328:THR:HG23	2.19	0.43
1:A:19:ILE:HG13	1:A:359:ILE:HB	2.00	0.43
1:A:249:THR:CG2	1:A:250:ILE:H	2.30	0.43
1:B:191:LYS:CG	1:B:192:ARG:N	2.82	0.43
1:B:40:ILE:HD13	8:B:620:HOH:O	2.19	0.43
1:A:21:VAL:HG11	1:A:353:ALA:HB1	2.01	0.42
1:B:46:VAL:CG2	1:B:47:ARG:N	2.81	0.42
1:B:209:GLU:O	1:B:213:ILE:HG13	2.20	0.42
1:B:37:ALA:HA	1:B:38:HIS:ND1	2.34	0.42
1:A:19:ILE:HG12	1:A:359:ILE:O	2.19	0.42
1:A:171:LEU:CD2	1:A:221:ARG:HD3	2.50	0.42
1:A:26:ARG:HH11	1:A:26:ARG:HG3	1.84	0.42
1:A:77:LYS:CD	7:A:606:PG4:H41	2.34	0.42
1:B:148:THR:HG23	1:B:149:ASP:OD1	2.20	0.42
1:A:101:ILE:HG22	1:A:102:PHE:N	2.33	0.42
1:B:116:GLU:O	1:B:134:GLY:N	2.52	0.41
1:A:144:PHE:CE1	1:A:206:ASN:HA	2.55	0.41
1:A:162:GLU:CD	1:A:171:LEU:HD13	2.40	0.41
1:A:76:THR:O	7:A:606:PG4:H32	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:ARG:CG	1:B:189:ARG:HH11	2.28	0.41
1:A:116:GLU:HG2	1:A:136:ILE:HD12	2.01	0.41
1:B:247:GLU:O	1:B:254:GLU:HA	2.20	0.41
1:A:30:LEU:HD11	1:A:33[A]:ARG:HH22	1.86	0.41
1:A:315:LYS:HB2	8:A:720:HOH:O	2.20	0.41
1:B:161:LEU:HD21	1:B:319:ILE:HD13	2.03	0.41
1:A:298:VAL:HG13	1:A:309:VAL:CG1	2.51	0.41
1:B:106:GLN:HG3	1:B:109:THR:CG2	2.50	0.41
1:B:162:GLU:CD	1:B:171:LEU:CD1	2.89	0.41
1:A:270:GLU:H	1:A:270:GLU:HG3	1.18	0.41
1:B:19:ILE:HG12	1:B:330:THR:HB	2.03	0.41
1:B:92:GLU:CG	1:B:329:ARG:HH12	2.33	0.41
1:A:38:HIS:CD2	1:A:38:HIS:N	2.87	0.41
1:B:111:LYS:NZ	1:B:268:GLY:CA	2.84	0.41
1:B:40:ILE:HD12	1:B:340:SER:HA	2.03	0.41
1:B:347:LEU:HA	1:B:347:LEU:HD12	1.90	0.40
1:B:169:PHE:CD2	1:B:178:VAL:O	2.74	0.40
1:B:26:ARG:NH1	1:B:27:PRO:O	2.54	0.40
1:B:83:ARG:HG2	1:B:83:ARG:NH1	2.37	0.40
1:A:241:VAL:O	1:A:241:VAL:HG13	2.22	0.40
1:B:119:ARG:HG2	6:B:605:MOY:S	2.61	0.40
1:B:162:GLU:CD	1:B:171:LEU:HD11	2.41	0.40
1:B:220:LYS:O	1:B:223:THR:HB	2.21	0.40
1:B:288:ILE:O	1:B:288:ILE:HG22	2.20	0.40
1:B:295:LEU:HD13	1:B:352:TYR:OH	2.21	0.40
1:B:40:ILE:HD12	1:B:340:SER:HB3	2.02	0.40
1:A:160:LEU:HD13	1:A:239:PHE:CD2	2.56	0.40
1:A:17:LYS:N	1:A:361:ASN:O	2.54	0.40
1:A:78:GLN:HG2	7:A:606:PG4:C5	2.52	0.40
1:A:124:GLU:HG2	1:A:125:TYR:CE1	2.56	0.40
1:A:74:ALA:O	7:A:606:PG4:H12	2.22	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	332/367 (90%)	311 (94%)	16 (5%)	5 (2%)	10	26
1	B	331/367 (90%)	313 (95%)	16 (5%)	2 (1%)	25	50
All	All	663/734 (90%)	624 (94%)	32 (5%)	7 (1%)	14	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	ASP
1	A	150	ASN
1	A	17	LYS
1	A	121	PRO
1	B	178	VAL
1	A	310	PRO
1	B	310	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	298/321 (93%)	271 (91%)	27 (9%)	9	21
1	B	297/321 (92%)	276 (93%)	21 (7%)	14	34
All	All	595/642 (93%)	547 (92%)	48 (8%)	12	27

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LYS
1	A	40	ILE
1	A	46	VAL
1	A	92	GLU
1	A	118	GLU
1	A	140	LEU

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Mol	Chain	Res	Type
1	A	161	LEU
1	A	173	ASN
1	A	175	SER
1	A	177	ASP
1	A	179	SER
1	A	182	LEU
1	A	187	ASP
1	A	189	ARG
1	A	207	LYS
1	A	212	GLN
1	A	216	LYS
1	A	221	ARG
1	A	222	THR
1	A	253[A]	GLU
1	A	253[B]	GLU
1	A	255	LEU
1	A	270	GLU
1	A	271	ASN
1	A	292	LEU
1	A	295	LEU
1	A	348	SER
1	B	40	ILE
1	B	42	GLU
1	B	57	LEU
1	B	60	LYS
1	B	118	GLU
1	B	122	ASN
1	B	175	SER
1	B	177	ASP
1	B	181	ARG
1	B	191	LYS
1	B	216	LYS
1	B	241	VAL
1	B	246	LYS
1	B	292	LEU
1	B	295	LEU
1	B	329	ARG
1	B	343	LEU
1	B	347	LEU
1	B	348	SER
1	B	350	LEU
1	B	362	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	229	ASN
1	A	262	ASN
1	A	342	ASN
1	B	18	ASN
1	B	98	ASN
1	B	106	GLN
1	B	212	GLN
1	B	244	HIS
1	B	262	ASN
1	B	321	GLN
1	B	342	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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
6	MOY	A	602	-	26,26,26	3.59	12 (46%)	32,37,37	1.76	5 (15%)
7	PG4	A	606	-	12,12,12	1.02	0	11,11,11	0.72	0
6	MOY	B	605	-	26,26,26	3.53	14 (53%)	32,37,37	1.75	7 (21%)
5	ADP	B	603	-	24,29,29	1.98	10 (41%)	29,45,45	1.82	4 (13%)
5	ADP	A	600	-	24,29,29	1.79	8 (33%)	29,45,45	1.78	4 (13%)

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
6	MOY	A	602	-	-	0/12/32/32	0/3/3/3
7	PG4	A	606	-	-	8/10/10/10	-
6	MOY	B	605	-	-	0/12/32/32	0/3/3/3
5	ADP	B	603	-	-	5/12/32/32	0/3/3/3
5	ADP	A	600	-	-	6/12/32/32	0/3/3/3

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	602	MOY	C1-N1	12.46	1.43	1.33
6	B	605	MOY	C1-N1	10.49	1.42	1.33
6	B	605	MOY	C4-N1	5.93	1.52	1.47
6	B	605	MOY	C4-C3	5.84	1.56	1.51
6	A	602	MOY	C4-N1	5.81	1.52	1.47
6	A	602	MOY	C4-C3	5.50	1.56	1.51
6	B	605	MOY	C2-N2	4.89	1.47	1.39
6	A	602	MOY	C2-C3	4.82	1.41	1.35
6	B	605	MOY	C2-C3	4.30	1.41	1.35
6	B	605	MOY	C24-C4	4.14	1.59	1.52
6	A	602	MOY	C24-C4	4.08	1.59	1.52
5	B	603	ADP	C2-N3	3.95	1.38	1.32
6	A	602	MOY	C2-N2	3.68	1.45	1.39
6	B	605	MOY	C25-C26	3.60	1.44	1.39
5	A	600	ADP	C4-N3	3.56	1.40	1.35
6	B	605	MOY	C25-C24	3.39	1.44	1.39
5	B	603	ADP	C4-N3	3.31	1.40	1.35
5	A	600	ADP	C2-N3	3.24	1.37	1.32
6	A	602	MOY	C25-C24	3.14	1.44	1.39
6	B	605	MOY	C29-C24	3.01	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	603	ADP	C2'-C1'	-2.97	1.49	1.53
5	A	600	ADP	PB-O3B	-2.92	1.43	1.54
5	B	603	ADP	O4'-C1'	2.86	1.45	1.41
5	A	600	ADP	O4'-C1'	2.79	1.45	1.41
6	A	602	MOY	C29-C24	2.75	1.43	1.39
5	B	603	ADP	C2-N1	2.73	1.39	1.33
6	B	605	MOY	C23-C18	2.68	1.43	1.39
6	A	602	MOY	C25-C26	2.60	1.43	1.39
5	B	603	ADP	PA-O2A	-2.59	1.43	1.55
5	A	600	ADP	C5-N7	-2.58	1.30	1.39
5	B	603	ADP	C5'-C4'	2.57	1.59	1.51
5	A	600	ADP	C2-N1	2.51	1.38	1.33
5	B	603	ADP	O4'-C4'	-2.51	1.39	1.45
5	B	603	ADP	PB-O3B	-2.50	1.45	1.54
6	B	605	MOY	C1-N2	2.47	1.40	1.37
6	A	602	MOY	C23-C18	2.45	1.43	1.39
6	B	605	MOY	C20-C19	2.43	1.44	1.38
5	B	603	ADP	C5-N7	-2.35	1.31	1.39
6	A	602	MOY	C21-C20	2.34	1.44	1.38
6	B	605	MOY	C21-C20	2.34	1.44	1.38
6	A	602	MOY	C20-C19	2.29	1.43	1.38
5	A	600	ADP	PA-O2A	-2.27	1.44	1.55
6	B	605	MOY	C19-C18	2.24	1.43	1.39
5	A	600	ADP	C5'-C4'	2.02	1.57	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	600	ADP	N3-C2-N1	-7.52	116.92	128.68
5	B	603	ADP	N3-C2-N1	-7.42	117.08	128.68
6	A	602	MOY	C4-N1-C1	-5.90	116.14	125.36
6	B	605	MOY	C4-N1-C1	-5.72	116.42	125.36
6	A	602	MOY	C3-C4-N1	-5.12	105.01	109.09
6	B	605	MOY	S-C1-N2	3.85	126.88	123.53
6	B	605	MOY	C3-C2-N2	-3.34	115.64	119.04
6	B	605	MOY	C3-C4-N1	-3.32	106.45	109.09
6	A	602	MOY	C6-C2-C3	3.20	129.08	125.17
6	A	602	MOY	C3-C2-N2	-3.09	115.89	119.04
5	B	603	ADP	C2'-C3'-C4'	2.83	108.14	102.64
6	A	602	MOY	C29-C24-C4	2.72	125.27	120.80
6	B	605	MOY	S-C1-N1	-2.71	117.71	121.10
5	A	600	ADP	O4'-C4'-C5'	-2.61	100.79	109.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	600	ADP	C2-N1-C6	2.59	123.19	118.75
6	B	605	MOY	C29-C24-C4	2.58	125.04	120.80
5	B	603	ADP	C2-N1-C6	2.40	122.86	118.75
5	B	603	ADP	O4'-C4'-C5'	-2.33	101.70	109.37
6	B	605	MOY	C6-C2-C3	2.30	127.98	125.17
5	A	600	ADP	C2'-C3'-C4'	2.02	106.58	102.64

There are no chirality outliers.

All (19) torsion outliers are listed below:

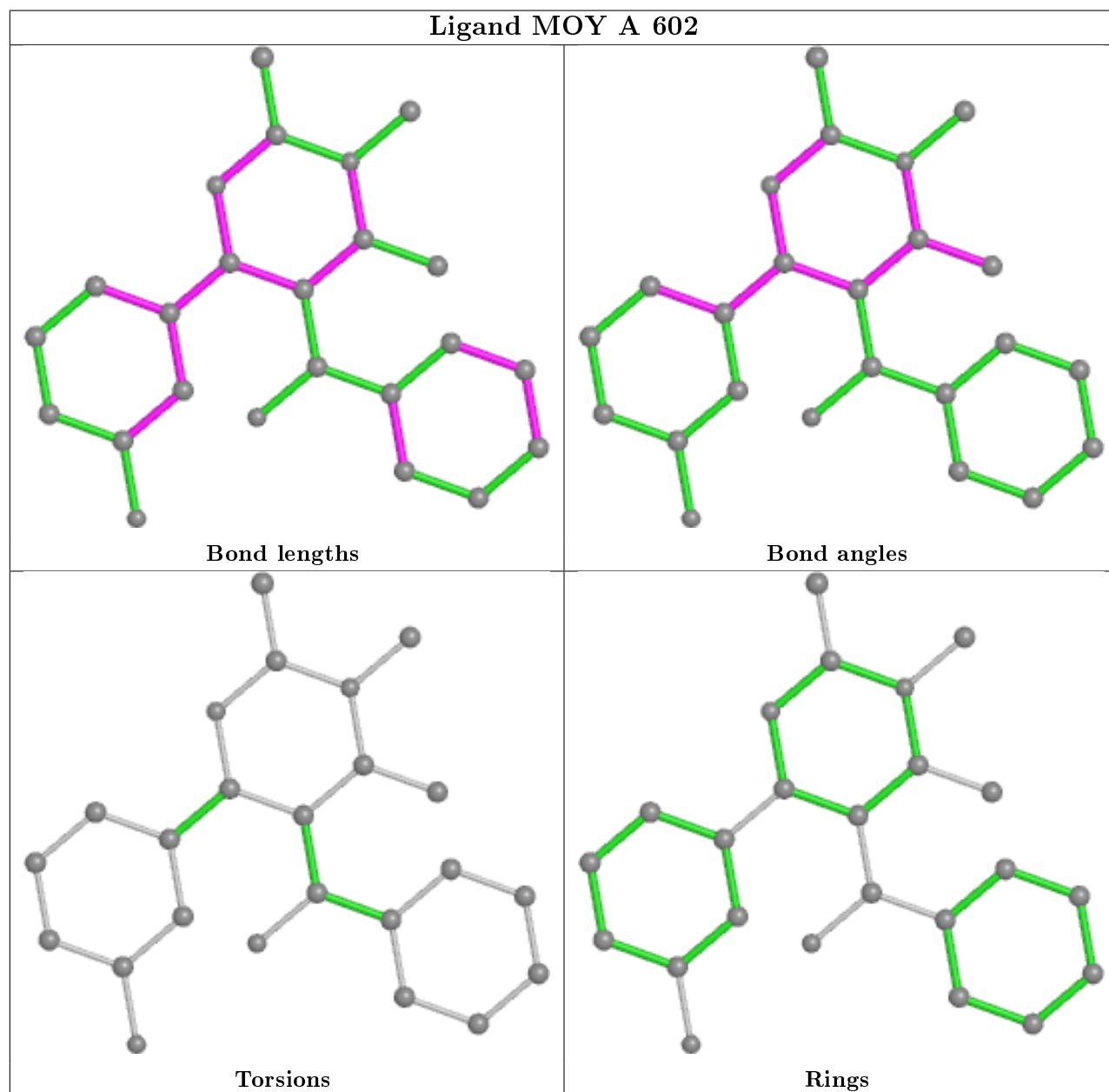
Mol	Chain	Res	Type	Atoms
5	A	600	ADP	C5'-O5'-PA-O1A
5	A	600	ADP	C5'-O5'-PA-O2A
5	A	600	ADP	C5'-O5'-PA-O3A
5	B	603	ADP	C5'-O5'-PA-O1A
5	B	603	ADP	C5'-O5'-PA-O2A
5	B	603	ADP	C5'-O5'-PA-O3A
7	A	606	PG4	C6-C5-O3-C4
7	A	606	PG4	O3-C5-C6-O4
5	A	600	ADP	O4'-C4'-C5'-O5'
5	A	600	ADP	C3'-C4'-C5'-O5'
5	B	603	ADP	O4'-C4'-C5'-O5'
5	B	603	ADP	C3'-C4'-C5'-O5'
7	A	606	PG4	O4-C7-C8-O5
7	A	606	PG4	O2-C3-C4-O3
7	A	606	PG4	O1-C1-C2-O2
7	A	606	PG4	C5-C6-O4-C7
5	A	600	ADP	PA-O3A-PB-O2B
7	A	606	PG4	C8-C7-O4-C6
7	A	606	PG4	C3-C4-O3-C5

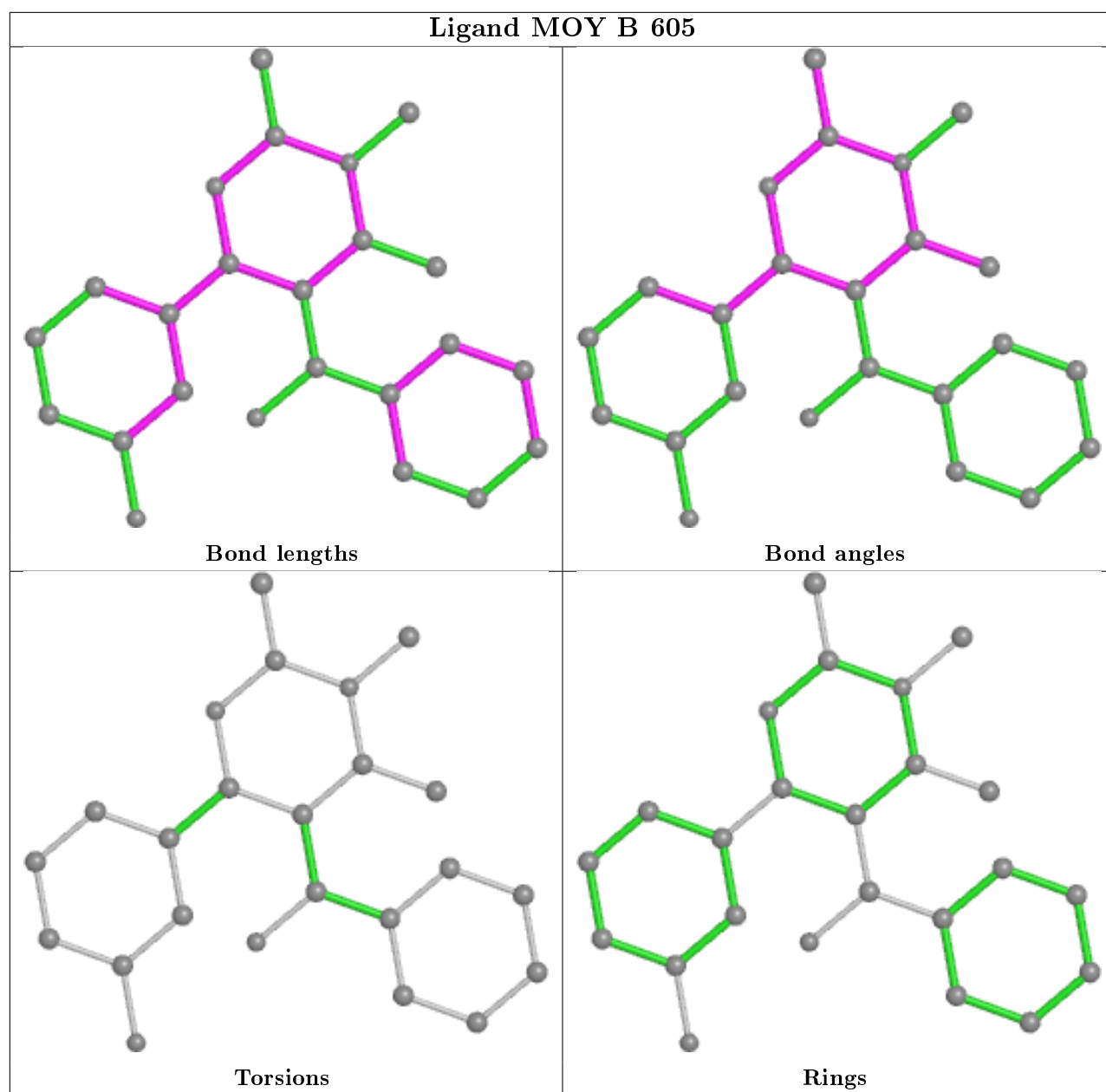
There are no ring outliers.

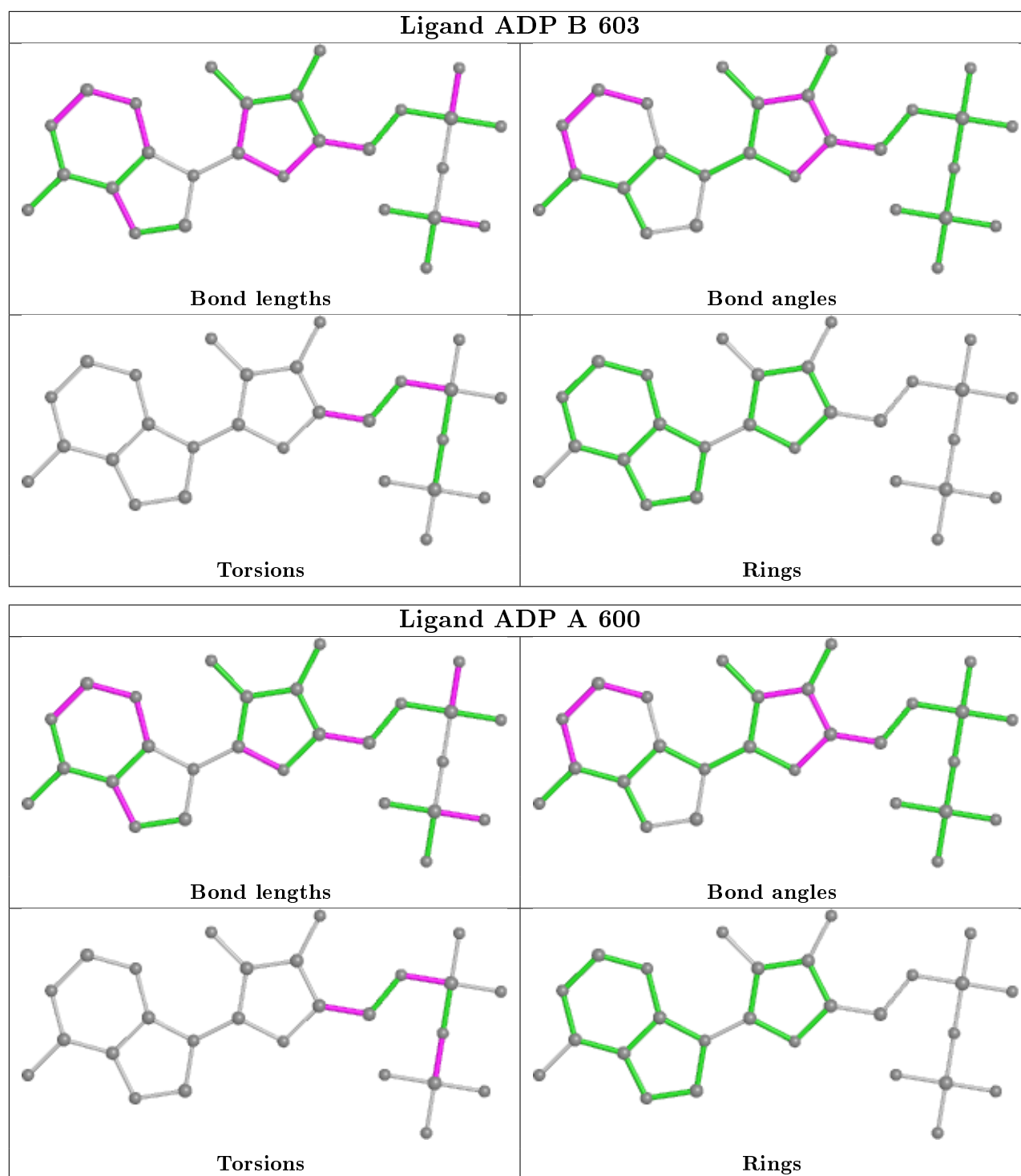
5 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	602	MOY	1	0
7	A	606	PG4	19	0
6	B	605	MOY	2	0
5	B	603	ADP	2	0
5	A	600	ADP	2	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 ⓘ

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	334/367 (91%)	0.21	21 (6%)	20 19	5, 21, 59, 75	3 (0%)
1	B	335/367 (91%)	0.15	17 (5%)	28 26	3, 24, 63, 84	1 (0%)
All	All	669/734 (91%)	0.18	38 (5%)	23 22	3, 23, 61, 84	4 (0%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	16	GLY	6.1
1	A	34	LYS	5.8
1	A	149	ASP	5.2
1	B	56	GLY	4.2
1	A	364	GLU	3.9
1	A	192	ARG	3.8
1	B	364	GLU	3.8
1	A	150	ASN	3.6
1	B	363	PRO	3.6
1	A	363	PRO	3.5
1	B	192	ARG	3.5
1	B	366	ASN	3.5
1	B	34	LYS	3.2
1	A	57	LEU	3.2
1	A	56	GLY	3.1
1	A	36	SER	3.1
1	A	58	ALA	3.1
1	B	31	ALA	3.0
1	B	176	SER	2.8
1	B	57	LEU	2.8
1	A	188	PRO	2.7
1	B	178	VAL	2.7
1	B	177	ASP	2.7
1	A	189	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	251	ASP	2.6
1	B	62	SER	2.5
1	A	33[A]	ARG	2.4
1	A	37	ALA	2.3
1	B	36	SER	2.3
1	A	16	GLY	2.2
1	A	38	HIS	2.2
1	A	191	LYS	2.2
1	A	97	TYR	2.2
1	A	360	LEU	2.2
1	B	191	LYS	2.1
1	A	62	SER	2.1
1	B	362	LYS	2.0
1	B	360	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

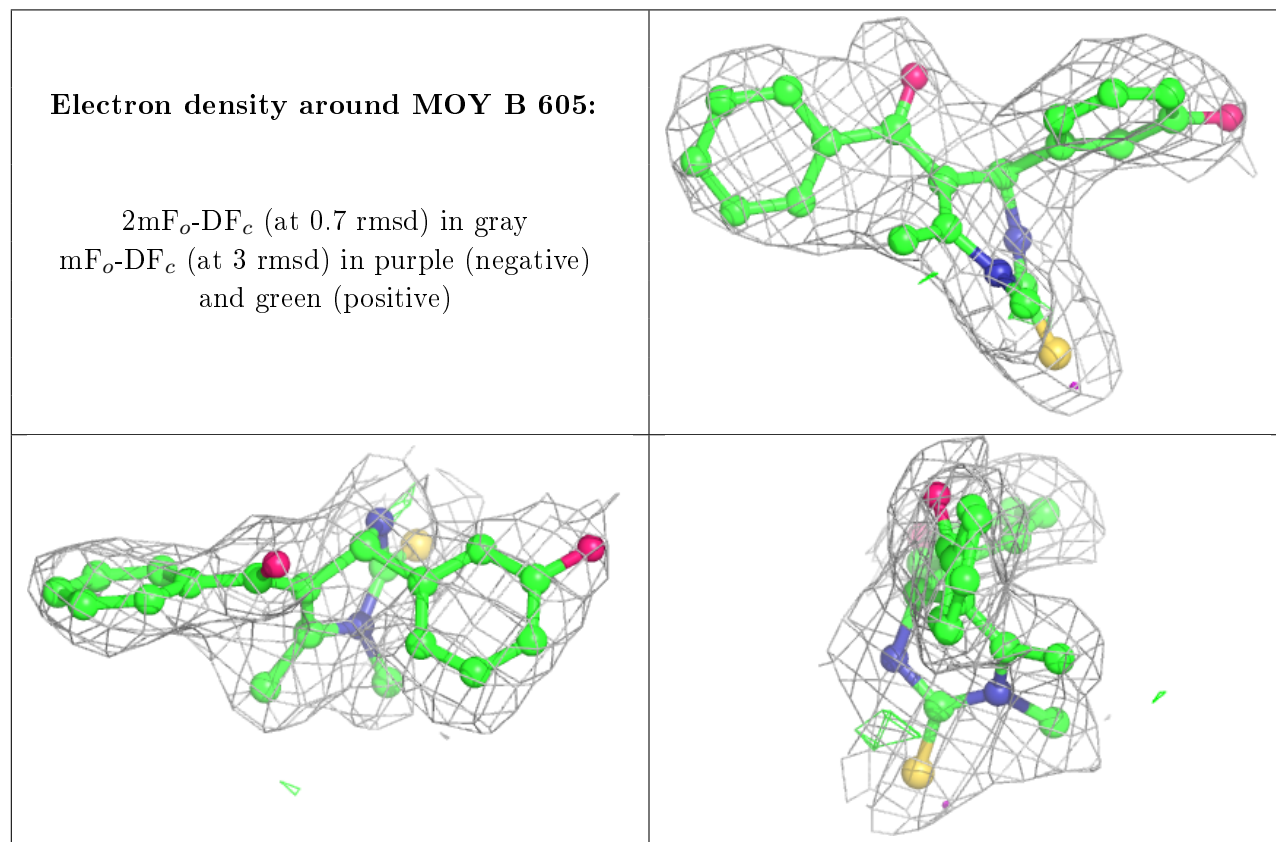
There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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.

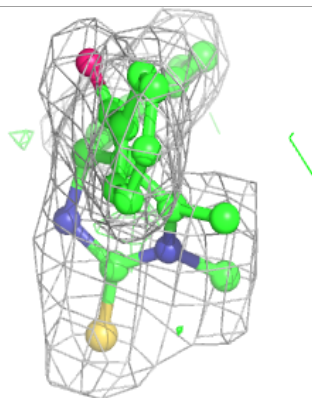
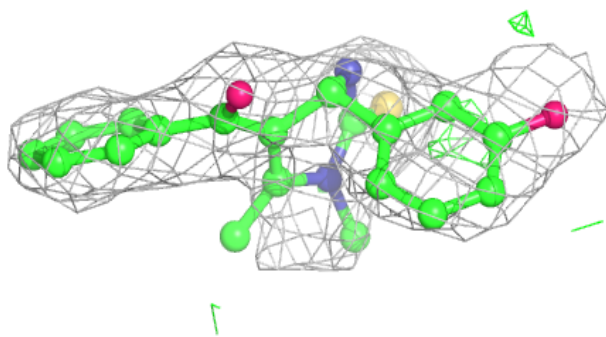
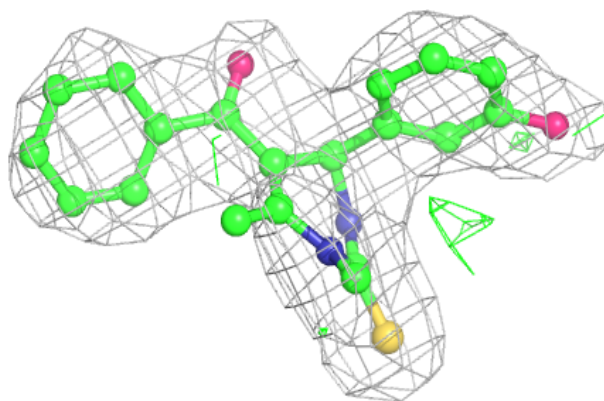
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	PG4	A	606	13/13	0.84	0.35	35,49,54,54	0
6	MOY	B	605	24/24	0.89	0.20	14,27,38,40	0
4	CL	A	608	1/1	0.93	0.10	33,33,33,33	0
2	MG	A	601	1/1	0.93	0.21	11,11,11,11	0
3	K	A	607	1/1	0.93	0.07	35,35,35,35	0
6	MOY	A	602	24/24	0.94	0.18	15,25,30,45	0
2	MG	B	604	1/1	0.96	0.17	11,11,11,11	0
5	ADP	A	600	27/27	0.96	0.13	4,17,27,29	0
5	ADP	B	603	27/27	0.96	0.15	10,21,36,44	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.

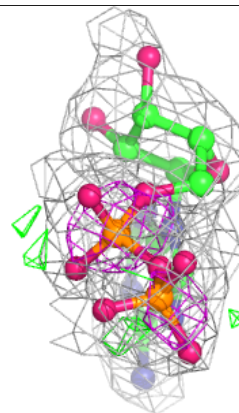
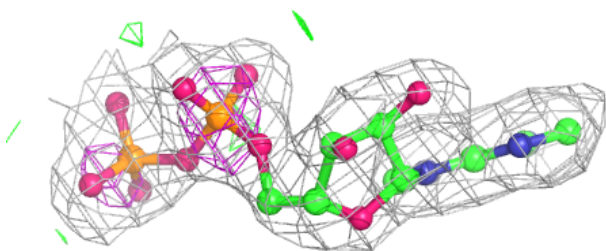
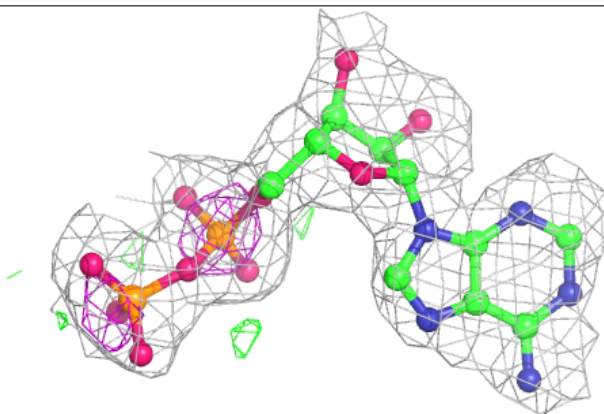


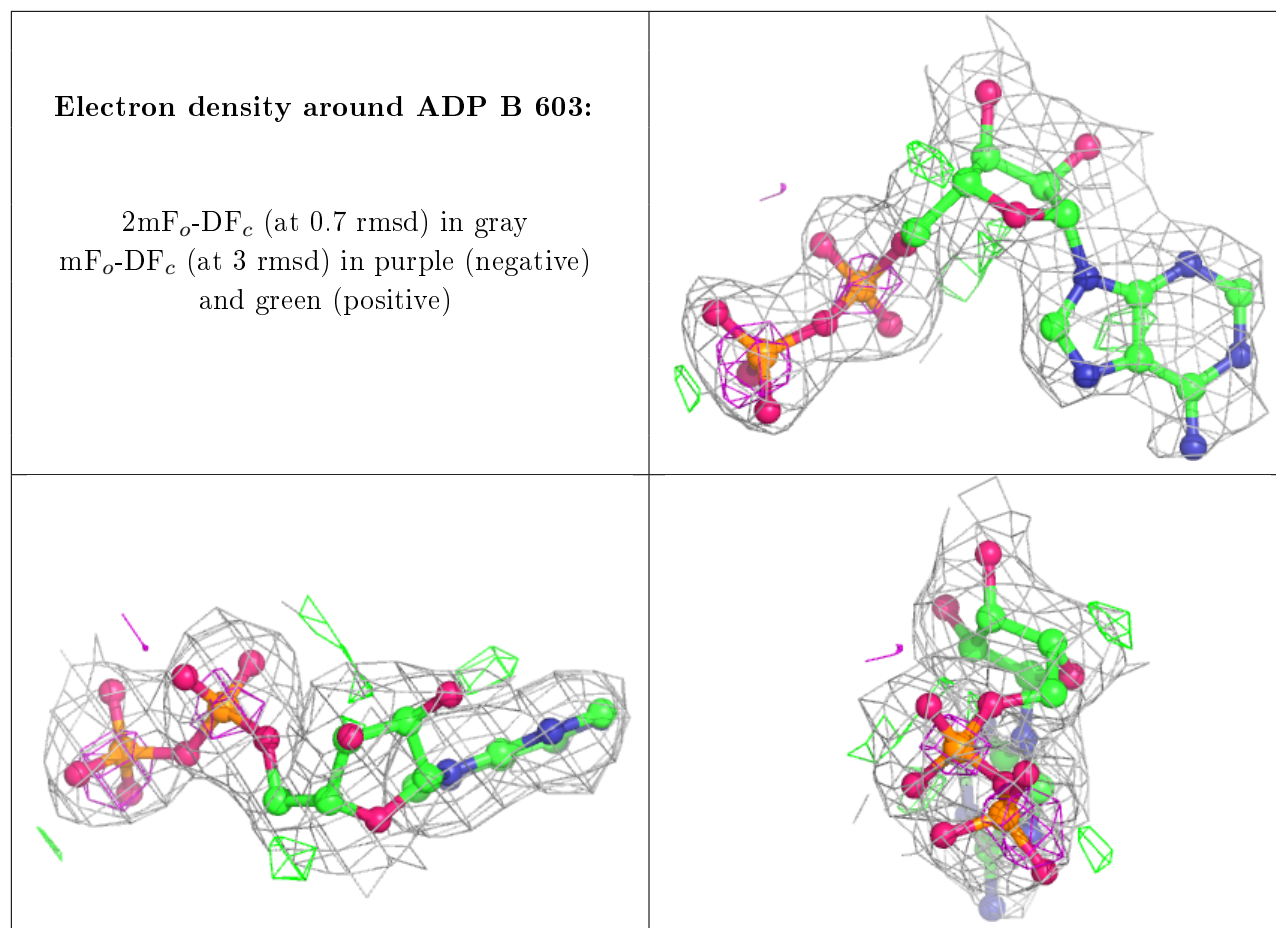
Electron density around MOY A 602:

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

**Electron density around ADP A 600:**

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





6.5 Other polymers ⓘ

There are no such residues in this entry.