



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:31 PM UTC

PDB ID : 3JVV / pdb_00003jvv
Title : Crystal Structure of P. aeruginosa PilT with bound AMP-PCP
Authors : Misic, A.M.; Satyshur, K.A.; Forest, K.T.
Deposited on : 2009-09-17
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

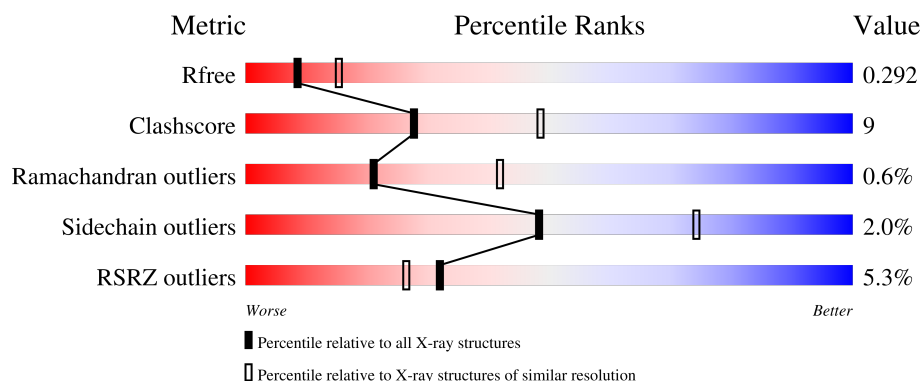
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

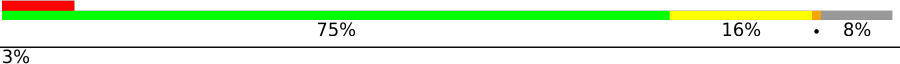
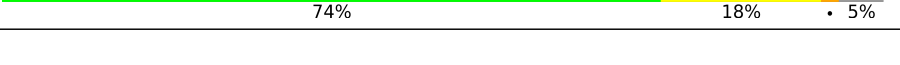
The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	356	
1	B	356	
1	C	356	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7939 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 Twitching mobility protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	0	0
			2577	1617	458	487	15			
1	B	328	Total	C	N	O	S	0	0	0
			2551	1599	453	484	15			
1	C	337	Total	C	N	O	S	0	0	0
			2613	1641	464	493	15			

There are 36 discrepancies between the modelled and reference sequences:

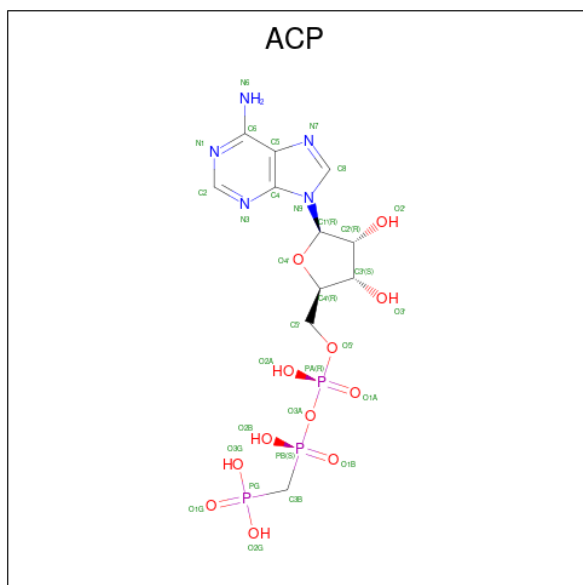
Chain	Residue	Modelled	Actual	Comment	Reference
A	345	GLY	-	expression tag	UNP P24559
A	346	ALA	-	expression tag	UNP P24559
A	347	ALA	-	expression tag	UNP P24559
A	348	ALA	-	expression tag	UNP P24559
A	349	LEU	-	expression tag	UNP P24559
A	350	GLU	-	expression tag	UNP P24559
A	351	HIS	-	expression tag	UNP P24559
A	352	HIS	-	expression tag	UNP P24559
A	353	HIS	-	expression tag	UNP P24559
A	354	HIS	-	expression tag	UNP P24559
A	355	HIS	-	expression tag	UNP P24559
A	356	HIS	-	expression tag	UNP P24559
B	345	GLY	-	expression tag	UNP P24559
B	346	ALA	-	expression tag	UNP P24559
B	347	ALA	-	expression tag	UNP P24559
B	348	ALA	-	expression tag	UNP P24559
B	349	LEU	-	expression tag	UNP P24559
B	350	GLU	-	expression tag	UNP P24559
B	351	HIS	-	expression tag	UNP P24559
B	352	HIS	-	expression tag	UNP P24559
B	353	HIS	-	expression tag	UNP P24559
B	354	HIS	-	expression tag	UNP P24559
B	355	HIS	-	expression tag	UNP P24559

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Chain	Residue	Modelled	Actual	Comment	Reference
B	356	HIS	-	expression tag	UNP P24559
C	345	GLY	-	expression tag	UNP P24559
C	346	ALA	-	expression tag	UNP P24559
C	347	ALA	-	expression tag	UNP P24559
C	348	ALA	-	expression tag	UNP P24559
C	349	LEU	-	expression tag	UNP P24559
C	350	GLU	-	expression tag	UNP P24559
C	351	HIS	-	expression tag	UNP P24559
C	352	HIS	-	expression tag	UNP P24559
C	353	HIS	-	expression tag	UNP P24559
C	354	HIS	-	expression tag	UNP P24559
C	355	HIS	-	expression tag	UNP P24559
C	356	HIS	-	expression tag	UNP P24559

- Molecule 2 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (CCD ID: ACP) (formula: $C_{11}H_{18}N_5O_{12}P_3$).

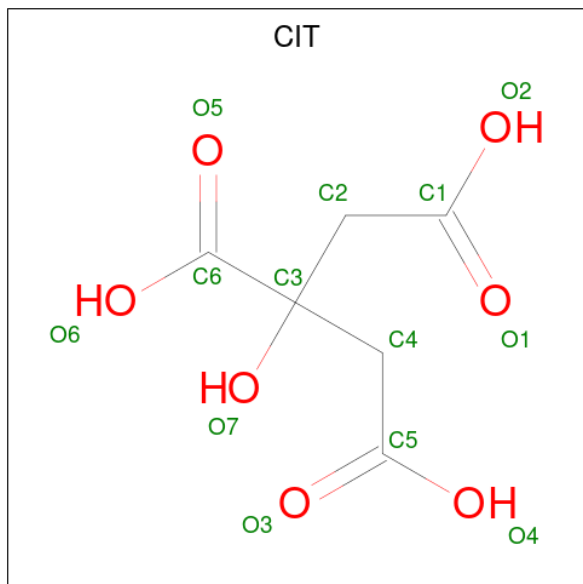


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 11	N 5	O 12	P 3	0	0
2	B	1	Total 31	C 11	N 5	O 12	P 3	0	0
2	C	1	Total 31	C 11	N 5	O 12	P 3	0	0

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

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

- Molecule 4 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 13 6 7	0	0

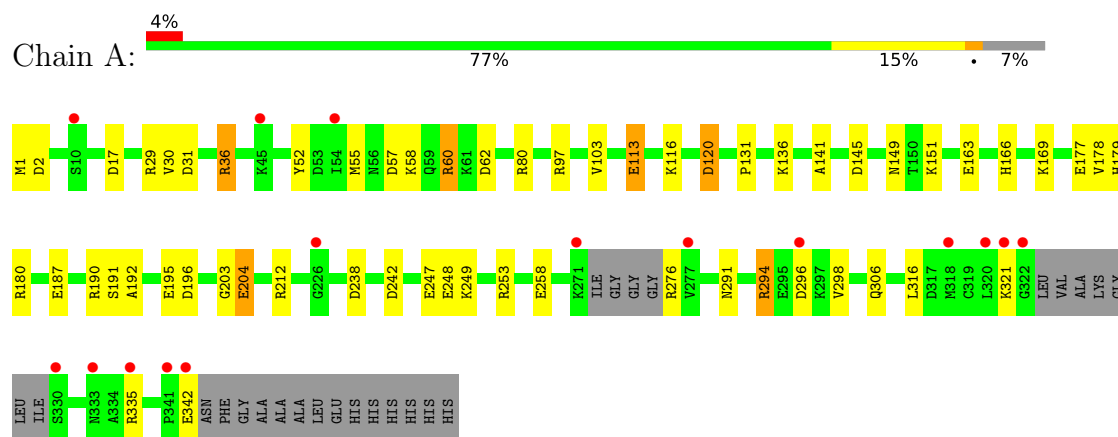
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	32	Total O 32 32	0	0
5	B	29	Total O 29 29	0	0
5	C	28	Total O 28 28	0	0

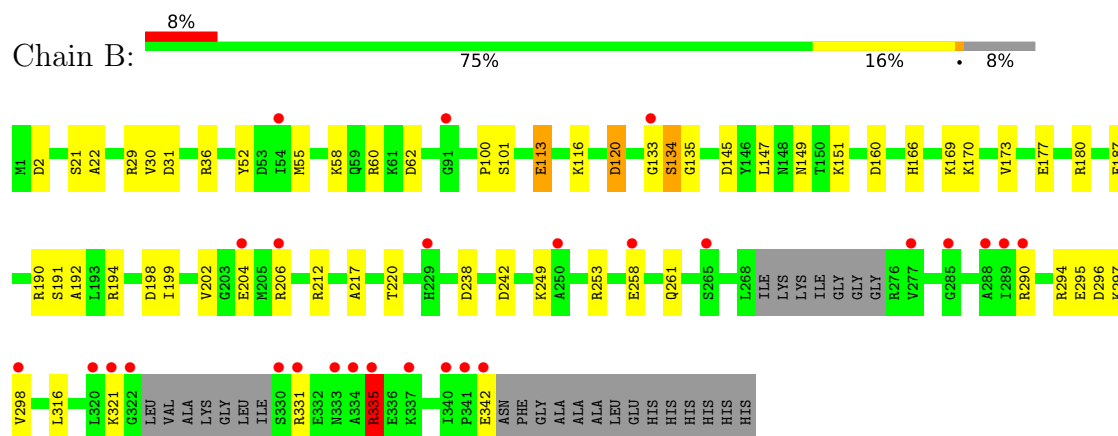
3 Residue-property plots [i](#)

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 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.

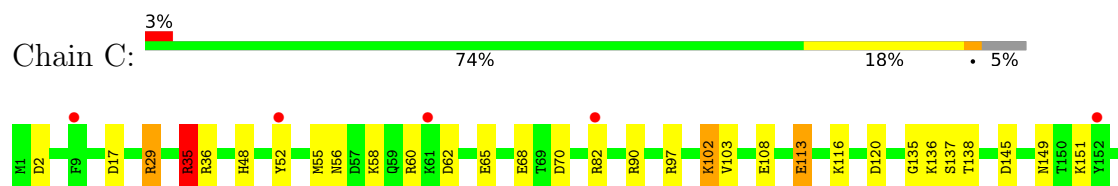
• Molecule 1: Twitching mobility protein

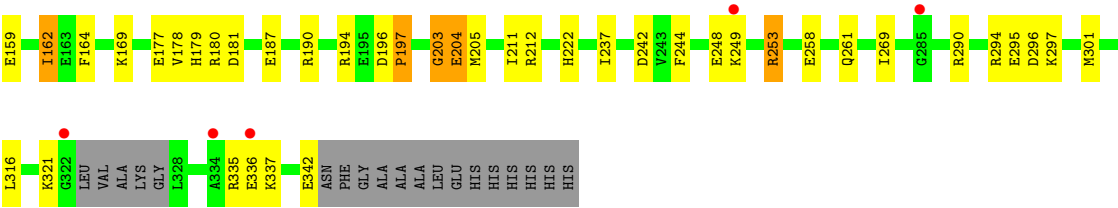


• Molecule 1: Twitching mobility protein



• Molecule 1: Twitching mobility protein





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	108.48Å 119.55Å 185.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.12 – 2.60 29.12 – 2.60	Depositor EDS
% Data completeness (in resolution range)	94.2 (29.12-2.60) 94.4 (29.12-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.244 , 0.291 0.247 , 0.292	Depositor DCC
R_{free} test set	1759 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	60.7	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7939	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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: CIT, ACP, MG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.65	2/2614 (0.1%)	0.93	13/3523 (0.4%)
1	B	0.67	1/2588 (0.0%)	0.96	11/3490 (0.3%)
1	C	0.75	4/2651 (0.2%)	0.92	12/3574 (0.3%)
All	All	0.69	7/7853 (0.1%)	0.94	36/10587 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	48	HIS	CE1-NE2	14.44	1.47	1.32
1	C	48	HIS	CG-CD2	8.54	1.45	1.35
1	A	1	MET	N-CA	6.80	1.59	1.46
1	C	48	HIS	ND1-CE1	6.66	1.39	1.32
1	C	48	HIS	CG-ND1	5.92	1.44	1.38
1	B	166	HIS	ND1-CE1	5.27	1.37	1.32
1	A	166	HIS	ND1-CE1	5.11	1.37	1.32

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ARG	NE-CZ-NH1	-8.89	112.61	121.50
1	B	212	ARG	NE-CZ-NH1	-8.27	113.23	121.50
1	B	331	ARG	NE-CZ-NH2	8.11	126.50	119.20
1	A	60	ARG	NE-CZ-NH2	7.99	126.39	119.20
1	B	331	ARG	NE-CZ-NH1	-7.88	113.62	121.50
1	B	212	ARG	NE-CZ-NH2	7.63	126.07	119.20
1	A	166	HIS	CB-CG-CD2	-6.98	122.12	131.20
1	B	166	HIS	CB-CG-CD2	-6.68	122.51	131.20
1	C	29	ARG	NE-CZ-NH2	-6.49	113.36	119.20
1	C	253	ARG	NE-CZ-NH2	-6.48	113.37	119.20
1	A	294	ARG	NE-CZ-NH2	6.31	124.88	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	ARG	NE-CZ-NH2	6.30	124.87	119.20
1	A	190	ARG	NE-CZ-NH2	-6.26	113.56	119.20
1	C	253	ARG	CD-NE-CZ	6.12	132.96	124.40
1	C	48	HIS	CB-CG-CD2	-6.02	123.37	131.20
1	A	276	ARG	NE-CZ-NH1	-5.98	115.52	121.50
1	A	166	HIS	CB-CG-ND1	5.94	131.60	122.70
1	C	253	ARG	NE-CZ-NH1	5.75	127.25	121.50
1	B	166	HIS	CB-CG-ND1	5.72	131.27	122.70
1	C	162	ILE	CB-CA-C	-5.69	102.31	111.59
1	A	36	ARG	NE-CZ-NH2	-5.68	114.08	119.20
1	C	29	ARG	CD-NE-CZ	5.68	132.35	124.40
1	C	35	ARG	CG-CD-NE	5.64	124.41	112.00
1	A	29	ARG	NE-CZ-NH1	-5.50	116.00	121.50
1	C	36	ARG	NE-CZ-NH1	-5.49	116.01	121.50
1	B	134	SER	N-CA-C	5.48	122.48	110.80
1	A	294	ARG	NE-CZ-NH1	-5.46	116.04	121.50
1	A	60	ARG	CD-NE-CZ	5.42	131.99	124.40
1	C	36	ARG	NE-CZ-NH2	5.40	124.06	119.20
1	B	331	ARG	CD-NE-CZ	5.34	131.88	124.40
1	B	190	ARG	NE-CZ-NH1	-5.26	116.24	121.50
1	B	190	ARG	NE-CZ-NH2	5.20	123.88	119.20
1	C	29	ARG	NE-CZ-NH1	5.12	126.62	121.50
1	A	190	ARG	CD-NE-CZ	5.09	131.52	124.40
1	C	190	ARG	NE-CZ-NH1	-5.04	116.46	121.50
1	B	335	ARG	NE-CZ-NH2	-5.03	114.67	119.20

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	2577	0	2628	40	1
1	B	2551	0	2591	49	0
1	C	2613	0	2671	62	1
2	A	31	0	14	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	31	0	13	0	0
2	C	31	0	14	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	13	0	5	0	0
5	A	32	0	0	4	0
5	B	29	0	0	7	0
5	C	28	0	0	6	0
All	All	7939	0	7936	142	1

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 9.

All (142) 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:36:ARG:NH1	5:A:366:HOH:O	2.14	0.81
1:A:321:LYS:NZ	1:A:342:GLU:OE2	2.18	0.75
1:C:321:LYS:NZ	1:C:342:GLU:OE2	2.20	0.73
1:C:65:GLU:O	5:C:363:HOH:O	2.05	0.73
1:B:194:ARG:HE	1:C:82:ARG:HH21	1.36	0.72
1:B:100:PRO:HA	5:B:373:HOH:O	1.88	0.72
1:B:321:LYS:NZ	1:B:342:GLU:OE2	2.20	0.69
1:B:238:ASP:HB2	5:B:368:HOH:O	1.92	0.68
1:A:298:VAL:HG12	5:A:365:HOH:O	1.93	0.68
1:C:222:HIS:HA	5:C:372:HOH:O	1.96	0.65
1:B:170:LYS:O	5:B:376:HOH:O	2.15	0.64
1:B:194:ARG:HE	1:C:82:ARG:NH2	1.95	0.64
1:B:173:VAL:O	1:C:90:ARG:NH2	2.29	0.64
1:B:100:PRO:CA	5:B:373:HOH:O	2.43	0.64
1:C:35:ARG:HG2	1:C:35:ARG:HH21	1.63	0.62
1:B:194:ARG:NE	1:C:82:ARG:NH2	2.47	0.62
1:A:80:ARG:NH1	1:A:163:GLU:O	2.33	0.61
1:C:102:LYS:NZ	1:C:108:GLU:OE2	2.33	0.61
1:C:120:ASP:OD1	1:C:151:LYS:HE3	2.01	0.60
1:C:180:ARG:HD2	1:C:181:ASP:OD1	2.01	0.60
1:B:258:GLU:OE2	1:B:294:ARG:NH1	2.31	0.60
1:C:97:ARG:HG3	5:C:371:HOH:O	2.02	0.60
1:B:335:ARG:HD2	1:B:342:GLU:O	2.02	0.59
1:B:58:LYS:NZ	1:B:62:ASP:OD2	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LYS:NZ	1:A:62:ASP:OD2	2.36	0.59
1:C:70:ASP:OD2	5:C:376:HOH:O	2.17	0.58
1:C:103:VAL:HG12	1:C:164:PHE:CD1	2.38	0.58
1:B:187:GLU:CD	1:B:187:GLU:N	2.62	0.58
1:A:120:ASP:OD1	1:A:151:LYS:HE3	2.04	0.57
1:A:187:GLU:N	1:A:187:GLU:CD	2.63	0.57
1:C:55:MET:O	1:C:60:ARG:NH1	2.37	0.57
1:C:145:ASP:OD1	1:C:169:LYS:NZ	2.34	0.57
1:C:187:GLU:N	1:C:187:GLU:CD	2.64	0.56
1:A:212:ARG:HD3	1:A:248:GLU:OE2	2.06	0.56
1:C:82:ARG:NE	5:C:376:HOH:O	2.26	0.56
1:A:253:ARG:CZ	1:A:296:ASP:CG	2.79	0.55
1:A:247:GLU:HG3	5:A:376:HOH:O	2.05	0.55
1:B:191:SER:OG	1:C:68:GLU:OE1	2.20	0.55
1:C:258:GLU:OE2	1:C:294:ARG:NH1	2.31	0.55
1:C:58:LYS:NZ	1:C:62:ASP:OD2	2.39	0.54
1:A:149:ASN:OD1	1:A:169:LYS:HD2	2.08	0.54
1:C:253:ARG:CZ	1:C:296:ASP:CG	2.81	0.54
1:C:103:VAL:HG12	1:C:164:PHE:HD1	1.72	0.54
1:C:295:GLU:OE1	1:C:297:LYS:NZ	2.25	0.54
1:B:198:ASP:OD2	1:C:29:ARG:NH2	2.38	0.53
1:B:194:ARG:HD2	1:C:82:ARG:CZ	2.39	0.53
1:A:177:GLU:OE2	1:A:180:ARG:NH2	2.42	0.53
1:B:120:ASP:OD1	1:B:151:LYS:HE3	2.09	0.52
1:B:253:ARG:CZ	1:B:296:ASP:CG	2.81	0.52
1:C:244:PHE:O	1:C:249:LYS:NZ	2.32	0.52
1:C:203:GLY:O	1:C:204:GLU:HB3	2.10	0.52
1:B:160:ASP:OD2	1:B:206:ARG:NH2	2.40	0.51
1:C:177:GLU:OE2	1:C:180:ARG:NH2	2.44	0.51
1:B:55:MET:O	1:B:60:ARG:NH1	2.44	0.51
1:B:149:ASN:OD1	1:B:169:LYS:HD2	2.11	0.50
1:B:295:GLU:OE1	1:B:297:LYS:NZ	2.29	0.50
1:A:57:ASP:OD1	1:A:60:ARG:NH2	2.44	0.50
1:A:103:VAL:HG13	1:A:141:ALA:HB1	1.94	0.50
1:A:169:LYS:O	1:B:36:ARG:NH1	2.45	0.49
1:A:335:ARG:NE	1:A:342:GLU:O	2.40	0.49
1:C:336:GLU:CD	1:C:337:LYS:NZ	2.70	0.49
1:A:17:ASP:OD1	1:A:97:ARG:NE	2.35	0.49
1:B:177:GLU:OE2	1:B:180:ARG:NH2	2.45	0.49
1:A:253:ARG:NH1	1:A:296:ASP:CG	2.70	0.49
1:B:113:GLU:OE2	1:B:116:LYS:NZ	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:GLU:CD	1:B:187:GLU:H	2.19	0.49
1:C:253:ARG:NH1	1:C:296:ASP:CG	2.70	0.48
1:A:291:ASN:OD1	1:A:294:ARG:NH2	2.46	0.48
1:A:242:ASP:O	1:A:249:LYS:NZ	2.47	0.48
1:A:187:GLU:CD	1:A:187:GLU:H	2.22	0.47
1:C:212:ARG:HD3	1:C:248:GLU:OE2	2.13	0.47
1:B:101:SER:N	5:B:373:HOH:O	2.48	0.47
1:A:131:PRO:O	1:A:136:LYS:NZ	2.48	0.47
1:A:52:TYR:C	1:A:52:TYR:CD1	2.92	0.47
1:C:187:GLU:CD	1:C:187:GLU:H	2.23	0.47
1:A:204:GLU:O	5:A:383:HOH:O	2.20	0.46
1:B:238:ASP:OD1	1:B:298:VAL:HG11	2.15	0.46
1:C:120:ASP:OD1	1:C:151:LYS:CE	2.62	0.46
1:C:17:ASP:OD1	1:C:97:ARG:HG2	2.16	0.46
1:A:55:MET:O	1:A:60:ARG:NH1	2.49	0.46
1:A:145:ASP:CG	1:A:169:LYS:HZ1	2.16	0.46
1:C:335:ARG:NE	1:C:342:GLU:O	2.43	0.46
1:C:135:GLY:O	1:C:136:LYS:C	2.58	0.46
1:C:269:ILE:HG12	1:C:337:LYS:HD2	1.98	0.46
1:A:258:GLU:CD	1:A:294:ARG:HH11	2.23	0.45
1:B:120:ASP:OD1	1:B:151:LYS:CE	2.64	0.45
1:C:149:ASN:OD1	1:C:169:LYS:HD2	2.15	0.45
1:C:253:ARG:NH1	1:C:296:ASP:OD2	2.50	0.45
1:C:159:GLU:HB2	1:C:162:ILE:HA	1.98	0.45
1:A:120:ASP:OD1	1:A:151:LYS:CE	2.64	0.45
1:C:336:GLU:OE2	1:C:337:LYS:NZ	2.50	0.45
1:C:336:GLU:CD	1:C:337:LYS:HZ2	2.25	0.44
1:A:113:GLU:OE2	1:A:116:LYS:NZ	2.50	0.44
1:B:253:ARG:NH1	1:B:296:ASP:CG	2.75	0.44
1:B:261:GLN:OE1	1:B:290:ARG:NH1	2.51	0.44
1:C:138:THR:OG1	2:C:400:ACP:O2A	2.36	0.44
1:C:113:GLU:OE2	1:C:116:LYS:NZ	2.51	0.44
1:B:217:ALA:O	1:B:220:THR:HG22	2.18	0.44
1:C:97:ARG:HD2	5:C:378:HOH:O	2.18	0.43
1:C:159:GLU:CB	1:C:162:ILE:HA	2.48	0.43
1:B:52:TYR:O	1:B:60:ARG:NH1	2.49	0.43
1:C:145:ASP:CG	1:C:169:LYS:HZ1	2.24	0.43
1:C:269:ILE:HG12	1:C:337:LYS:CD	2.48	0.43
1:B:113:GLU:OE1	1:B:116:LYS:NZ	2.49	0.43
1:C:35:ARG:HH21	1:C:35:ARG:CG	2.32	0.43
1:C:178:VAL:O	1:C:179:HIS:HB2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:238:ASP:OD1	1:A:298:VAL:HG11	2.19	0.43
1:C:70:ASP:CG	1:C:82:ARG:HH21	2.27	0.43
1:A:253:ARG:NH1	1:A:296:ASP:OD2	2.52	0.42
1:B:133:GLY:O	1:B:135:GLY:N	2.52	0.42
1:A:178:VAL:O	1:A:179:HIS:HB2	2.20	0.42
1:B:58:LYS:NZ	1:B:62:ASP:OD1	2.52	0.42
1:B:242:ASP:O	1:B:249:LYS:NZ	2.52	0.42
1:A:113:GLU:OE1	1:A:116:LYS:NZ	2.52	0.42
1:B:147:LEU:HD13	1:B:199:ILE:HD13	2.02	0.42
1:A:58:LYS:NZ	1:A:62:ASP:OD1	2.53	0.42
1:A:30:VAL:O	1:A:31:ASP:C	2.62	0.42
1:A:191:SER:O	1:A:192:ALA:C	2.62	0.42
1:B:100:PRO:C	5:B:373:HOH:O	2.63	0.42
1:B:21:SER:O	1:B:22:ALA:C	2.63	0.42
1:B:204:GLU:OE1	1:B:206:ARG:NH2	2.52	0.41
1:C:196:ASP:N	1:C:197:PRO:HD3	2.35	0.41
1:B:191:SER:O	1:B:192:ALA:C	2.63	0.41
1:B:145:ASP:OD1	1:B:169:LYS:NZ	2.26	0.41
1:C:336:GLU:OE1	1:C:337:LYS:NZ	2.47	0.41
1:A:58:LYS:NZ	1:A:62:ASP:CG	2.78	0.41
1:B:194:ARG:HD2	1:C:82:ARG:NH2	2.36	0.41
1:C:242:ASP:O	1:C:249:LYS:NZ	2.54	0.41
1:C:261:GLN:OE1	1:C:290:ARG:NH1	2.53	0.41
2:A:400:ACP:O1A	2:A:400:ACP:H3B2	2.21	0.41
1:B:30:VAL:O	1:B:31:ASP:C	2.62	0.41
1:A:335:ARG:NH1	1:A:342:GLU:O	2.52	0.41
1:C:56:ASN:OD1	1:C:56:ASN:C	2.64	0.41
1:B:58:LYS:NZ	1:B:62:ASP:CG	2.78	0.41
1:B:194:ARG:CD	1:C:82:ARG:NH2	2.84	0.41
1:A:196:ASP:CG	1:B:29:ARG:HH11	2.28	0.40
2:C:400:ACP:H3B1	2:C:400:ACP:O1A	2.22	0.40
1:B:202:VAL:O	5:B:374:HOH:O	2.21	0.40
1:A:195:GLU:O	1:A:196:ASP:C	2.64	0.40
1:C:205:MET:O	1:C:211:ILE:HD11	2.21	0.40
1:C:237:ILE:HD12	1:C:301:MET:HE1	2.03	0.40
1:C:58:LYS:NZ	1:C:62:ASP:OD1	2.53	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:GLN:O	1:C:60:ARG:NH2[5_545]	2.03	0.17

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	325/356 (91%)	309 (95%)	14 (4%)	2 (1%)	21	42
1	B	322/356 (90%)	307 (95%)	14 (4%)	1 (0%)	36	58
1	C	333/356 (94%)	314 (94%)	16 (5%)	3 (1%)	14	30
All	All	980/1068 (92%)	930 (95%)	44 (4%)	6 (1%)	21	42

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	134	SER
1	A	204	GLU
1	C	204	GLU
1	C	197	PRO
1	C	203	GLY
1	A	203	GLY

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	283/299 (95%)	279 (99%)	4 (1%)	59	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	280/299 (94%)	275 (98%)	5 (2%)	51	76
1	C	286/299 (96%)	278 (97%)	8 (3%)	38	66
All	All	849/897 (95%)	832 (98%)	17 (2%)	48	74

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	113	GLU
1	A	120	ASP
1	A	316	LEU
1	B	2	ASP
1	B	113	GLU
1	B	120	ASP
1	B	316	LEU
1	B	335	ARG
1	C	2	ASP
1	C	35	ARG
1	C	52	TYR
1	C	102	LYS
1	C	113	GLU
1	C	137	SER
1	C	194	ARG
1	C	316	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	84	ASN
1	B	13	GLN
1	B	84	ASN
1	B	300	GLN
1	C	13	GLN
1	C	87	ASN
1	C	166	HIS
1	C	300	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 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$
4	CIT	A	357	-	12,12,12	1.18	0	17,17,17	1.63	2 (11%)
2	ACP	A	400	3	31,33,33	1.54	8 (25%)	47,52,52	1.99	12 (25%)
2	ACP	B	400	3	31,33,33	1.55	7 (22%)	47,52,52	1.90	13 (27%)
2	ACP	C	400	3	31,33,33	1.65	6 (19%)	47,52,52	1.77	9 (19%)

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
4	CIT	A	357	-	-	3/16/16/16	-
2	ACP	A	400	3	-	0/19/38/38	0/3/3/3
2	ACP	B	400	3	-	2/19/38/38	0/3/3/3
2	ACP	C	400	3	-	3/19/38/38	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	400	ACP	C5-C4	4.91	1.47	1.39
2	A	400	ACP	C5-C4	4.86	1.47	1.39
2	B	400	ACP	C5-C4	4.65	1.47	1.39
2	C	400	ACP	PB-O3A	3.47	1.62	1.58
2	C	400	ACP	PA-O3A	3.15	1.62	1.59
2	A	400	ACP	PB-O3A	2.91	1.61	1.58
2	A	400	ACP	C5-C6	2.69	1.48	1.41
2	B	400	ACP	PA-O3A	2.68	1.62	1.59
2	C	400	ACP	C5-C6	2.62	1.48	1.41
2	B	400	ACP	C5-C6	2.50	1.47	1.41
2	C	400	ACP	C5-N7	-2.50	1.34	1.39
2	A	400	ACP	PA-O3A	2.40	1.62	1.59
2	B	400	ACP	C5-N7	-2.35	1.34	1.39
2	B	400	ACP	C8-N7	2.30	1.36	1.31
2	B	400	ACP	C4-N9	-2.25	1.33	1.37
2	B	400	ACP	PB-O2B	-2.24	1.50	1.56
2	A	400	ACP	C8-N7	2.20	1.35	1.31
2	A	400	ACP	C5-N7	-2.20	1.35	1.39
2	C	400	ACP	C8-N7	2.17	1.35	1.31
2	A	400	ACP	C4-N9	-2.11	1.33	1.37
2	A	400	ACP	PB-O2B	-2.09	1.51	1.56

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	400	ACP	C5-C4-N3	-5.71	118.85	126.72
2	A	400	ACP	C5-C4-N3	-5.50	119.14	126.72
2	B	400	ACP	C5-C4-N3	-5.34	119.37	126.72
2	A	400	ACP	PB-O3A-PA	-5.32	115.00	132.37
2	B	400	ACP	PB-O3A-PA	-4.89	116.41	132.37
2	C	400	ACP	N3-C4-N9	4.62	135.02	127.17
2	A	400	ACP	N3-C4-N9	4.49	134.81	127.17
2	B	400	ACP	N3-C4-N9	4.39	134.64	127.17
2	B	400	ACP	C2-N3-C4	3.53	120.44	111.83
2	A	400	ACP	C2-N3-C4	3.44	120.24	111.83
2	C	400	ACP	C2-N3-C4	3.37	120.06	111.83
4	A	357	CIT	O6-C6-C3	3.35	119.57	113.14
2	B	400	ACP	N3-C2-N1	-3.33	123.53	128.58
2	C	400	ACP	O1G-PG-C3B	-3.17	104.46	111.37
2	A	400	ACP	O1G-PG-C3B	-3.17	104.47	111.37
4	A	357	CIT	O7-C3-C6	3.13	113.40	108.96
2	A	400	ACP	O2B-PB-O1B	3.01	119.75	109.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	ACP	C4-C5-N7	-2.99	107.16	110.58
2	B	400	ACP	C4-C5-N7	-2.99	107.16	110.58
2	A	400	ACP	N3-C2-N1	-2.98	124.07	128.58
2	C	400	ACP	N3-C2-N1	-2.90	124.20	128.58
2	C	400	ACP	C4-C5-N7	-2.88	107.29	110.58
2	C	400	ACP	PB-O3A-PA	-2.84	123.09	132.37
2	C	400	ACP	O2B-PB-O1B	2.84	119.21	109.95
2	B	400	ACP	C4-N9-C8	2.83	108.71	105.74
2	C	400	ACP	O3G-PG-O2G	2.79	115.91	107.96
2	A	400	ACP	C4-N9-C8	2.53	108.39	105.74
2	A	400	ACP	O4'-C4'-C3'	2.47	110.05	105.15
2	B	400	ACP	O3G-PG-O2G	2.39	114.78	107.96
2	B	400	ACP	O2B-PB-O1B	2.25	117.26	109.95
2	A	400	ACP	C2'-C1'-N9	-2.21	107.81	113.30
2	B	400	ACP	O1G-PG-C3B	-2.18	106.63	111.37
2	B	400	ACP	O2G-PG-C3B	-2.12	101.26	106.40
2	B	400	ACP	C2-N1-C6	2.05	122.09	118.73
2	B	400	ACP	C5-N7-C8	2.04	106.66	103.45
2	A	400	ACP	C5-N7-C8	2.00	106.60	103.45

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	400	ACP	PG-C3B-PB-O1B
2	B	400	ACP	PG-C3B-PB-O3A
2	C	400	ACP	PG-C3B-PB-O1B
2	C	400	ACP	PG-C3B-PB-O3A
4	A	357	CIT	C2-C3-C4-C5
4	A	357	CIT	C6-C3-C4-C5
4	A	357	CIT	O7-C3-C4-C5
2	C	400	ACP	PG-C3B-PB-O2B

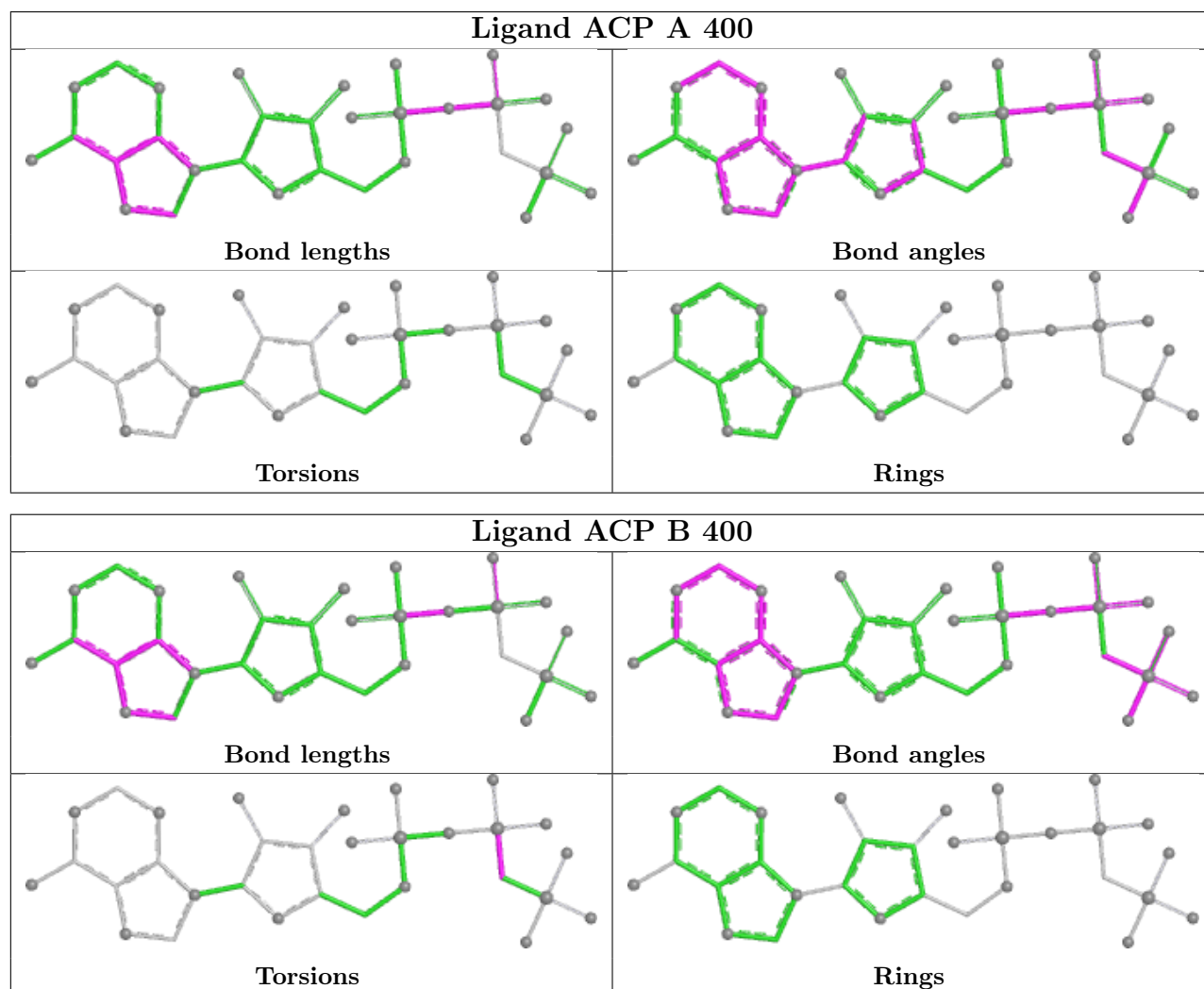
There are no ring outliers.

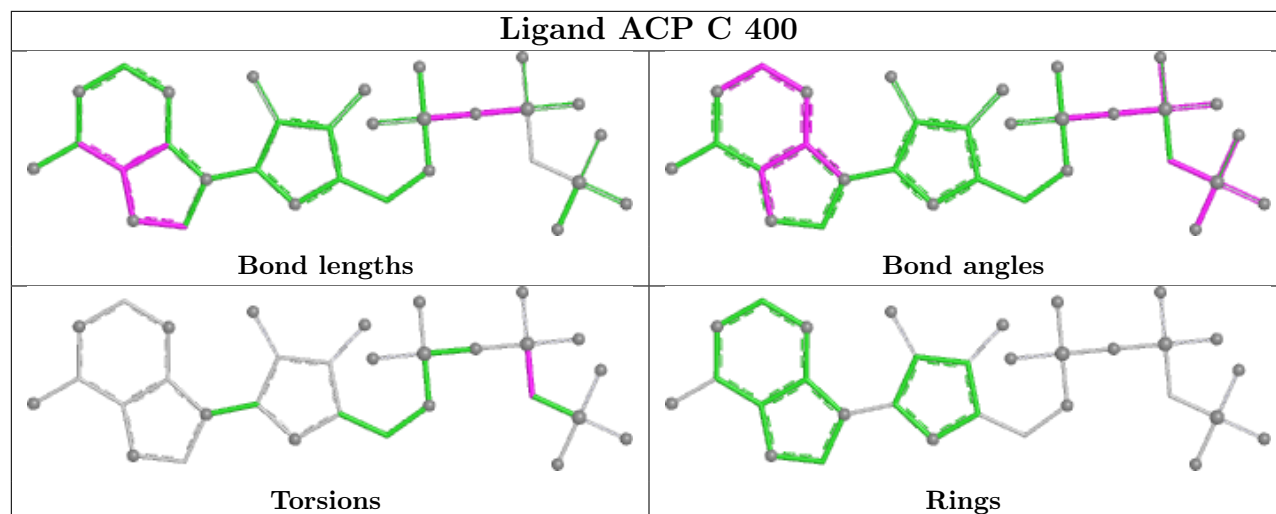
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	ACP	1	0
2	C	400	ACP	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 [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	331/356 (92%)	0.40	16 (4%)	35 30	38, 58, 89, 118	0
1	B	328/356 (92%)	0.59	27 (8%)	17 14	34, 57, 94, 121	0
1	C	337/356 (94%)	0.37	10 (2%)	52 47	35, 58, 86, 101	0
All	All	996/1068 (93%)	0.45	53 (5%)	32 26	34, 58, 89, 121	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	322	GLY	6.1
1	B	322	GLY	4.9
1	B	285	GLY	4.7
1	B	340	ILE	4.3
1	B	330	SER	3.9
1	B	321	LYS	3.9
1	C	285	GLY	3.8
1	A	271	LYS	3.5
1	B	335	ARG	3.5
1	B	206	ARG	3.5
1	A	296	ASP	3.3
1	A	330	SER	3.3
1	A	320	LEU	3.2
1	B	341	PRO	3.2
1	B	334	ALA	3.0
1	B	298	VAL	2.9
1	B	320	LEU	2.8
1	B	204	GLU	2.8
1	B	91	GLY	2.7
1	B	290	ARG	2.7
1	B	229	HIS	2.7
1	B	288	ALA	2.7
1	A	341	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	342	GLU	2.6
1	A	318	MET	2.5
1	B	277	VAL	2.5
1	A	321	LYS	2.5
1	B	337	LYS	2.5
1	A	45	LYS	2.5
1	C	61	LYS	2.5
1	B	265	SER	2.4
1	C	152	TYR	2.4
1	C	249	LYS	2.4
1	B	258	GLU	2.3
1	A	10	SER	2.3
1	A	277	VAL	2.3
1	C	82	ARG	2.3
1	A	335	ARG	2.3
1	C	336	GLU	2.2
1	B	133	GLY	2.2
1	B	54	ILE	2.2
1	B	331	ARG	2.2
1	C	322	GLY	2.2
1	B	250	ALA	2.2
1	A	226	GLY	2.2
1	A	54	ILE	2.2
1	A	342	GLU	2.1
1	B	333	ASN	2.1
1	C	9	PHE	2.1
1	C	52	TYR	2.1
1	C	334	ALA	2.1
1	A	333	ASN	2.1
1	B	289	ILE	2.1

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

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

6.3 Carbohydrates [i](#)

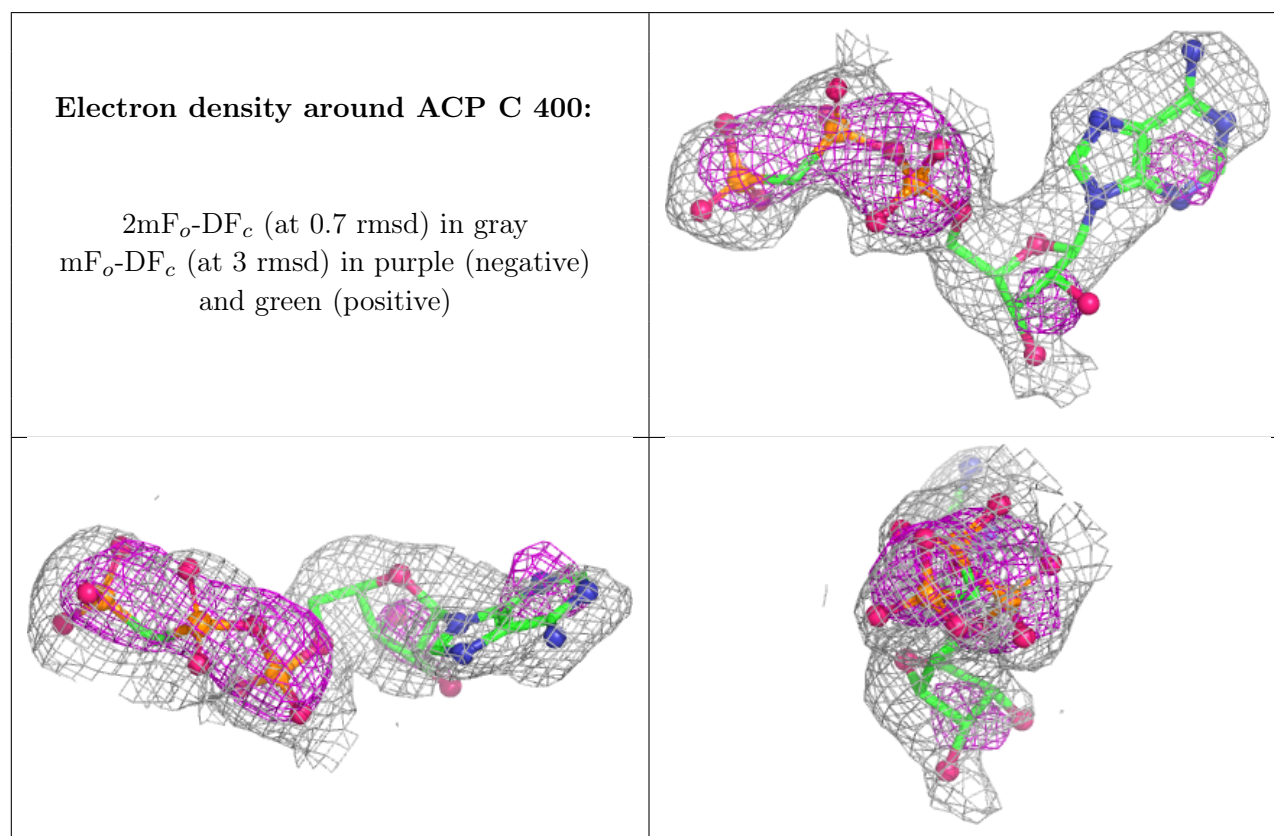
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

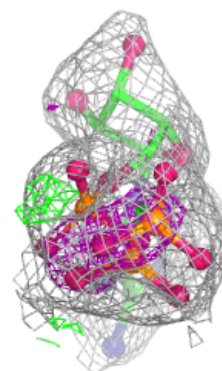
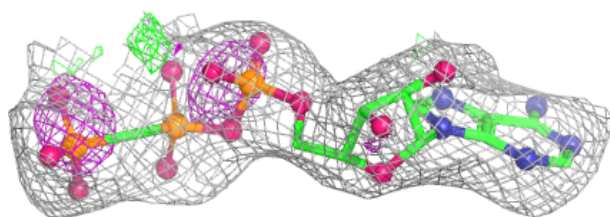
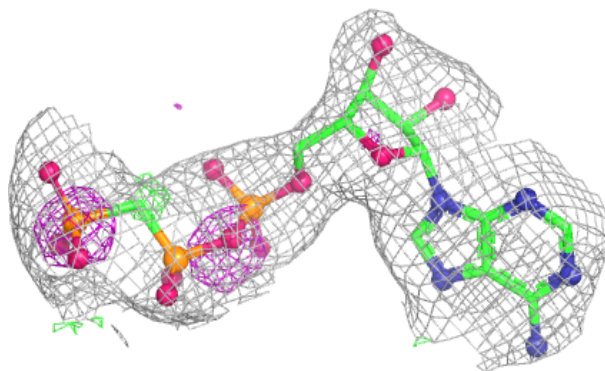
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	MG	C	401	1/1	0.75	0.13	64,64,64,64	0
2	ACP	C	400	31/31	0.80	0.14	68,71,71,72	0
3	MG	A	401	1/1	0.84	0.11	55,55,55,55	0
2	ACP	A	400	31/31	0.87	0.12	67,72,73,73	0
4	CIT	A	357	13/13	0.87	0.11	53,54,58,58	0
2	ACP	B	400	31/31	0.90	0.11	58,62,63,63	0
3	MG	B	401	1/1	0.97	0.10	44,44,44,44	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.

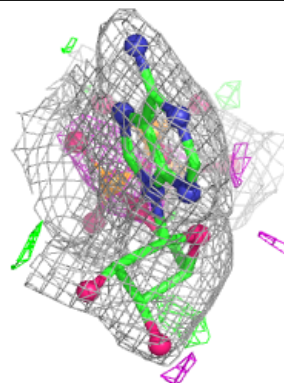
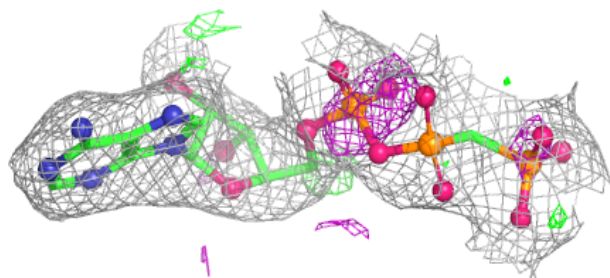
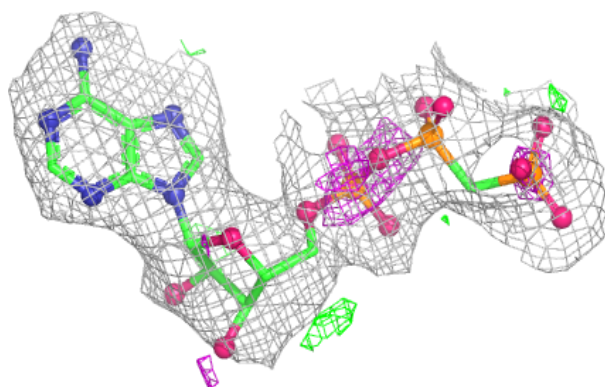


Electron density around ACP A 400:

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

**Electron density around ACP B 400:**

$2mF_o-DF_c$ (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.