



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:20 AM UTC

PDB ID : 2JBT / pdb_00002jbt
Title : Structure of the monooxygenase component of p-hydroxyphenylacetate hydroxylase from *Acinetobacter baumannii*
Authors : Alfieri, A.; Mattevi, A.
Deposited on : 2006-12-11
Resolution : 2.80 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

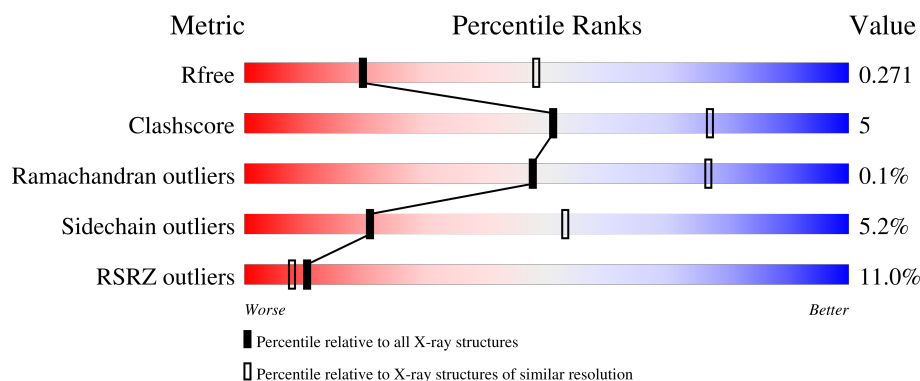
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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.

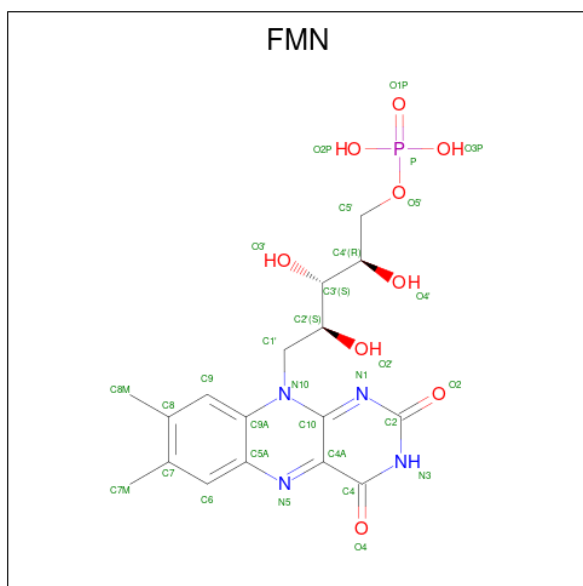
Mol	Chain	Length	Quality of chain
1	A	422	5% 83% 10% • 5%
1	B	422	5% 83% 10% • 5%
1	C	422	5% 82% 12% • 5%
1	D	422	27% 82% 11% • 5%

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 P-HYDROXYPHENYLACETATE HYDROXYLASE C2\OXYGENASE COMPONENT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	400	Total 3129	C 1986	N 539	O 583	S 21	0	0	0
1	B	399	Total 3121	C 1980	N 538	O 582	S 21	0	0	0
1	C	399	Total 3121	C 1980	N 538	O 582	S 21	0	0	0
1	D	399	Total 3121	C 1980	N 538	O 582	S 21	0	0	0

- Molecule 2 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



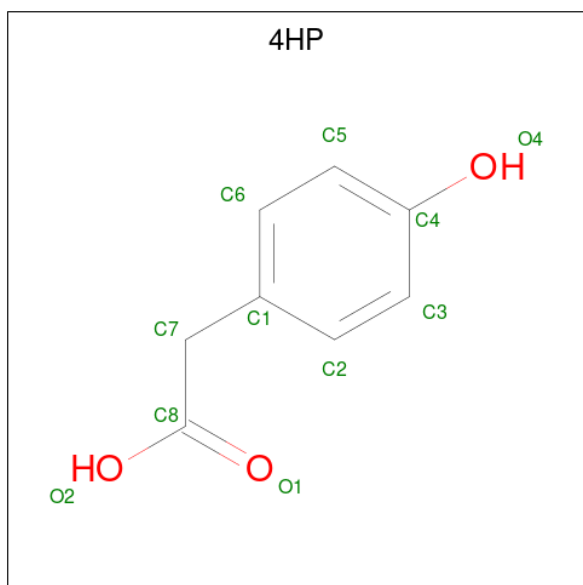
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 17	N 4	O 9	P 1	0	0
2	B	1	Total 31	C 17	N 4	O 9	P 1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	D	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is 4-HYDROXYPHENYLACETATE (CCD ID: 4HP) (formula: $C_8H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	8	3		
3	B	1	Total	C	O	0	0
			11	8	3		
3	C	1	Total	C	O	0	0
			11	8	3		
3	D	1	Total	C	O	0	0
			11	8	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	O	0	0
			7	7		
4	B	10	Total	O	0	0
			10	10		
4	C	6	Total	O	0	0
			6	6		

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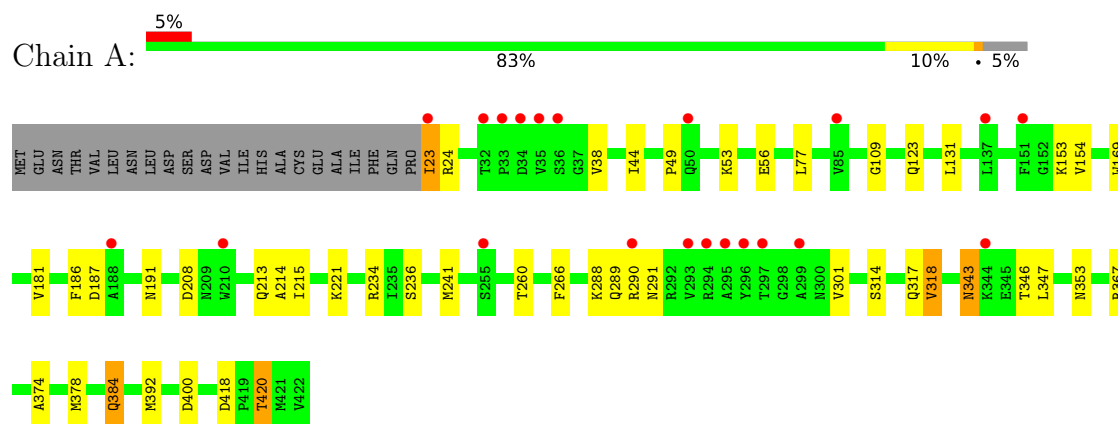
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	2	Total	O	0	0
			2	2		

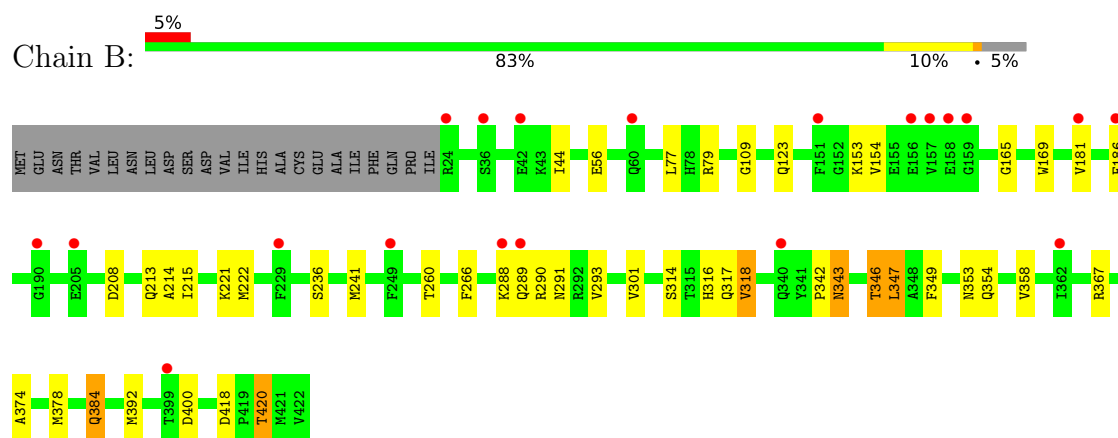
3 Residue-property plots [i](#)

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.

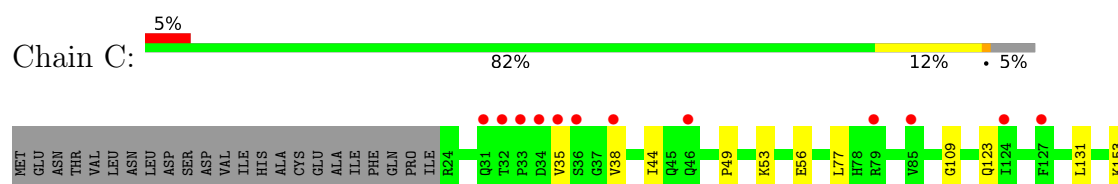
• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\ :OXYGENASE COMPONENT

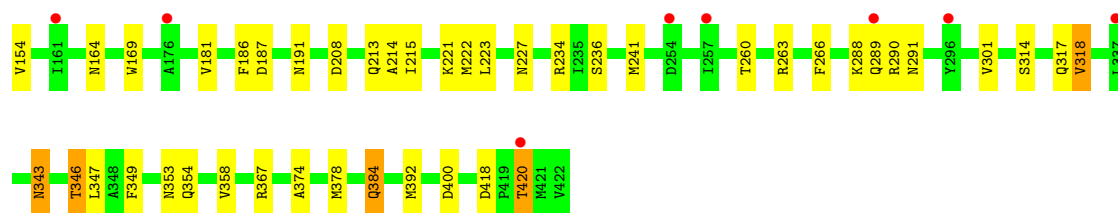


• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\ :OXYGENASE COMPONENT

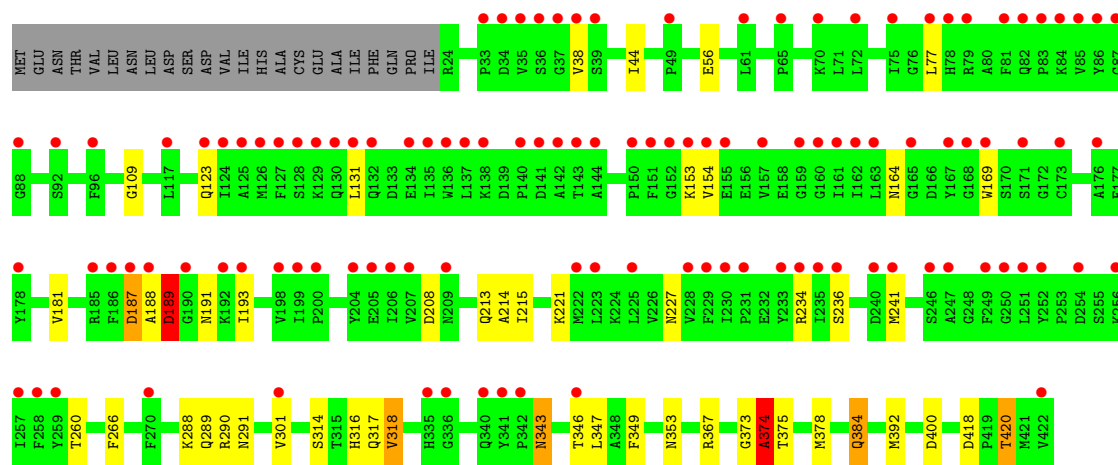
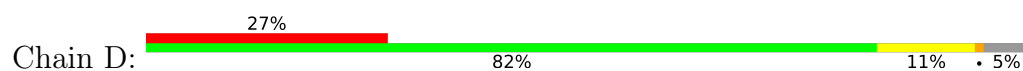


• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\ :OXYGENASE COMPONENT





● Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\ :OXYGENASE COMPONENT



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	91.99Å 181.26Å 286.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.51 – 2.80 84.51 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (84.51-2.80) 100.0 (84.51-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.82Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.217 , 0.235 0.253 , 0.271	Depositor DCC
R_{free} test set	616 reflections (1.04%)	wwPDB-VP
Wilson B-factor (Å ²)	58.0	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	12685	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 4HP, FMN

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	0.55	0/3201	0.82	1/4329 (0.0%)
1	B	0.56	0/3193	0.81	1/4318 (0.0%)
1	C	0.50	0/3193	0.79	0/4318
1	D	0.92	8/3193 (0.3%)	1.00	14/4318 (0.3%)
All	All	0.65	8/12780 (0.1%)	0.86	16/17283 (0.1%)

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	B	0	1
1	C	0	1
1	D	0	1
All	All	0	3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	187	ASP	CG-OD1	-22.67	0.82	1.25
1	D	374	ALA	C-N	-20.60	1.07	1.33
1	D	374	ALA	N-CA	16.71	1.67	1.46
1	D	189	ASP	CG-OD2	16.59	1.56	1.25
1	D	187	ASP	CG-OD2	14.25	1.52	1.25
1	D	373	GLY	C-N	11.65	1.49	1.34
1	D	189	ASP	CG-OD1	8.01	1.40	1.25
1	D	187	ASP	CA-CB	6.39	1.64	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	187	ASP	CB-CA-C	-12.40	86.67	109.37
1	D	374	ALA	CA-C-N	12.31	137.25	120.38
1	D	374	ALA	C-N-CA	12.31	137.25	120.38
1	D	187	ASP	N-CA-C	-12.26	90.46	109.95
1	D	373	GLY	O-C-N	11.94	135.13	122.60
1	D	373	GLY	CA-C-N	-11.73	103.39	120.28
1	D	373	GLY	C-N-CA	-11.73	103.39	120.28
1	D	375	THR	O-C-N	10.20	132.93	122.12
1	D	187	ASP	N-CA-CB	-9.12	96.25	110.85
1	D	188	ALA	N-CA-CB	-7.98	97.00	110.49
1	D	187	ASP	OD1-CG-OD2	-6.62	107.01	122.90
1	D	374	ALA	N-CA-C	6.44	119.12	111.71
1	D	374	ALA	N-CA-CB	-6.23	100.62	110.22
1	D	374	ALA	CA-C-O	-5.73	113.63	120.10
1	A	186	PHE	N-CA-C	5.31	117.48	109.41
1	B	165	GLY	N-CA-C	5.13	117.06	110.96

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	186	PHE	Peptide
1	C	186	PHE	Peptide
1	D	374	ALA	Mainchain

5.2 Too-close contacts [i](#)

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	3129	0	3071	36	0
1	B	3121	0	3060	36	0
1	C	3121	0	3060	39	0
1	D	3121	0	3059	38	0
2	A	31	0	19	2	0
2	B	31	0	19	2	0
2	C	31	0	19	1	0
2	D	31	0	19	1	0
3	A	11	0	6	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	11	0	6	2	0
3	C	11	0	6	3	0
3	D	11	0	7	2	0
4	A	7	0	0	0	0
4	B	10	0	0	3	0
4	C	6	0	0	0	0
4	D	2	0	0	1	0
All	All	12685	0	12351	130	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 5.

All (130) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:187:ASP:OD1	1:D:187:ASP:CG	0.82	1.06
1:D:187:ASP:OD1	1:D:187:ASP:CB	2.05	1.04
1:D:187:ASP:OD1	1:D:187:ASP:OD2	1.93	0.87
1:A:23:ILE:HG13	1:A:23:ILE:O	1.81	0.81
1:B:384:GLN:HG3	1:C:213:GLN:HG3	1.66	0.77
1:A:384:GLN:HG3	1:D:213:GLN:HG3	1.69	0.75
1:B:384:GLN:CG	1:C:213:GLN:HG3	2.17	0.74
1:B:213:GLN:HG3	1:C:384:GLN:HG3	1.69	0.74
1:A:213:GLN:HG3	1:D:384:GLN:HG3	1.70	0.73
1:A:123:GLN:HG2	1:A:241:MET:CE	2.19	0.72
1:B:213:GLN:HG3	1:C:384:GLN:CG	2.21	0.71
1:D:123:GLN:HG2	1:D:241:MET:CE	2.21	0.70
1:A:384:GLN:CG	1:D:213:GLN:HG3	2.21	0.70
1:A:213:GLN:HG3	1:D:384:GLN:CG	2.22	0.69
1:A:317:GLN:OE1	1:A:367:ARG:NH2	2.25	0.67
1:B:123:GLN:HG2	1:B:241:MET:CE	2.25	0.67
1:A:123:GLN:HG2	1:A:241:MET:HE1	1.74	0.67
1:D:123:GLN:HG2	1:D:241:MET:HE1	1.79	0.64
1:B:79:ARG:HD2	4:B:2001:HOH:O	1.96	0.64
1:C:317:GLN:OE1	1:C:367:ARG:NH2	2.26	0.62
1:C:123:GLN:HG2	1:C:241:MET:CE	2.30	0.61
1:B:123:GLN:HG2	1:B:241:MET:HE1	1.83	0.59
1:B:346:THR:HB	4:B:2007:HOH:O	2.03	0.59
1:D:189:ASP:OD2	1:D:189:ASP:N	2.37	0.57
1:C:208:ASP:HA	1:C:221:LYS:HG2	1.87	0.57
1:B:418:ASP:OD1	1:B:420:THR:HB	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:SER:O	1:A:318:VAL:HG13	2.05	0.56
1:D:314:SER:O	1:D:318:VAL:HG13	2.05	0.56
1:D:123:GLN:HG2	1:D:241:MET:HE2	1.88	0.56
1:B:317:GLN:OE1	1:B:367:ARG:NH2	2.29	0.56
1:A:418:ASP:OD1	1:A:420:THR:HB	2.05	0.56
1:A:214:ALA:HB3	1:A:392:MET:HB2	1.88	0.55
1:B:343:ASN:C	1:B:343:ASN:HD22	2.15	0.55
1:C:214:ALA:HB3	1:C:392:MET:HB2	1.88	0.55
1:C:343:ASN:HD22	1:C:343:ASN:C	2.14	0.55
1:A:343:ASN:C	1:A:343:ASN:HD22	2.15	0.55
1:D:109:GLY:HA2	1:D:215:ILE:HG23	1.88	0.55
1:C:123:GLN:HG2	1:C:241:MET:HE1	1.88	0.54
1:C:314:SER:O	1:C:318:VAL:HG13	2.07	0.54
1:C:418:ASP:OD1	1:C:420:THR:HB	2.08	0.54
1:A:109:GLY:HA2	1:A:215:ILE:HG23	1.89	0.54
1:D:187:ASP:HB2	1:D:191:ASN:O	2.08	0.54
1:A:208:ASP:HA	1:A:221:LYS:HG2	1.89	0.54
1:D:208:ASP:HA	1:D:221:LYS:HG2	1.88	0.54
1:D:418:ASP:OD1	1:D:420:THR:HB	2.08	0.53
1:C:109:GLY:HA2	1:C:215:ILE:HG23	1.90	0.53
1:C:215:ILE:HG22	1:C:215:ILE:O	2.09	0.53
1:A:266:PHE:CD2	3:A:1424:4HP:C3	2.92	0.53
1:B:109:GLY:HA2	1:B:215:ILE:HG23	1.90	0.53
1:D:317:GLN:OE1	1:D:367:ARG:NH2	2.30	0.52
1:D:266:PHE:CD2	3:D:1424:4HP:C3	2.93	0.52
1:A:123:GLN:HG2	1:A:241:MET:HE2	1.90	0.52
1:B:266:PHE:CD2	3:B:1424:4HP:C3	2.93	0.52
1:B:208:ASP:HA	1:B:221:LYS:HG2	1.91	0.52
1:B:314:SER:O	1:B:318:VAL:HG13	2.10	0.51
1:D:44:ILE:HD11	1:D:77:LEU:HD22	1.92	0.51
1:A:215:ILE:HG22	1:A:215:ILE:O	2.11	0.51
1:B:346:THR:HG21	4:B:2009:HOH:O	2.10	0.51
1:B:123:GLN:HG2	1:B:241:MET:HE2	1.93	0.51
1:B:215:ILE:HG22	1:B:215:ILE:O	2.11	0.51
1:A:44:ILE:HD11	1:A:77:LEU:HD22	1.93	0.50
1:B:44:ILE:HD11	1:B:77:LEU:HD22	1.93	0.50
1:C:266:PHE:CD2	3:C:1424:4HP:C3	2.93	0.50
1:A:23:ILE:O	1:A:23:ILE:CG1	2.57	0.50
1:C:44:ILE:HD11	1:C:77:LEU:HD22	1.94	0.50
1:A:214:ALA:CB	1:A:392:MET:HB2	2.41	0.49
1:A:378:MET:HE3	1:D:169:TRP:CH2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:214:ALA:CB	1:C:392:MET:HB2	2.42	0.49
1:D:215:ILE:O	1:D:215:ILE:HG22	2.12	0.49
1:D:214:ALA:HB3	1:D:392:MET:HB2	1.94	0.48
1:D:343:ASN:HD22	1:D:343:ASN:C	2.20	0.48
1:B:374:ALA:N	2:C:1423:FMN:O1P	2.47	0.48
1:A:374:ALA:N	2:D:1423:FMN:O1P	2.47	0.47
1:A:169:TRP:CH2	1:D:378:MET:HE3	2.49	0.47
1:B:169:TRP:CH2	1:C:378:MET:HE3	2.48	0.47
1:B:354:GLN:O	1:B:358:VAL:HG23	2.15	0.47
1:B:213:GLN:HG3	1:C:384:GLN:CD	2.40	0.47
1:C:123:GLN:HG2	1:C:241:MET:HE2	1.96	0.46
1:C:400:ASP:OD1	1:C:400:ASP:C	2.57	0.46
1:B:349:PHE:HD2	1:B:353:ASN:HD21	1.62	0.46
1:B:384:GLN:CD	1:C:213:GLN:HG3	2.39	0.46
1:B:214:ALA:HB3	1:B:392:MET:HB2	1.97	0.45
1:C:354:GLN:O	1:C:358:VAL:HG23	2.16	0.45
2:B:1423:FMN:O1P	1:C:374:ALA:N	2.49	0.45
1:A:187:ASP:HB2	1:A:191:ASN:H	1.80	0.45
1:D:214:ALA:CB	1:D:392:MET:HB2	2.46	0.45
1:A:187:ASP:HB2	1:A:191:ASN:N	2.31	0.45
1:B:214:ALA:CB	1:B:392:MET:HB2	2.47	0.44
1:C:346:THR:HB	4:D:2001:HOH:O	2.16	0.44
1:B:208:ASP:CG	1:B:221:LYS:HE3	2.42	0.44
1:A:384:GLN:CD	1:D:213:GLN:HG3	2.43	0.44
1:B:289:GLN:O	1:B:290:ARG:C	2.61	0.44
1:A:353:ASN:HD22	1:B:316:HIS:CD2	2.35	0.44
1:B:400:ASP:C	1:B:400:ASP:OD1	2.60	0.44
1:C:164:ASN:ND2	1:C:227:ASN:H	2.16	0.43
1:A:400:ASP:OD1	1:A:400:ASP:C	2.61	0.43
1:C:289:GLN:O	1:C:290:ARG:C	2.61	0.43
1:D:131:LEU:HD13	1:D:234:ARG:HD2	2.01	0.43
1:A:266:PHE:CD2	3:A:1424:4HP:C2	3.02	0.43
1:D:400:ASP:OD1	1:D:400:ASP:C	2.62	0.43
1:B:342:PRO:HG2	1:B:347:LEU:HD22	2.01	0.43
1:A:208:ASP:CG	1:A:221:LYS:HE3	2.45	0.42
1:C:131:LEU:HD13	1:C:234:ARG:HD2	2.02	0.42
1:C:187:ASP:OD2	1:C:191:ASN:HB2	2.18	0.42
1:D:289:GLN:O	1:D:290:ARG:C	2.62	0.42
1:D:349:PHE:HD2	1:D:353:ASN:HD21	1.68	0.42
1:A:187:ASP:CB	1:A:191:ASN:H	2.32	0.42
1:D:187:ASP:OD2	1:D:193:ILE:HD11	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:ASP:CG	1:D:221:LYS:HE3	2.45	0.42
1:B:378:MET:HE3	1:C:169:TRP:CH2	2.55	0.42
1:C:208:ASP:CG	1:C:221:LYS:HE3	2.45	0.42
1:B:266:PHE:CD2	3:B:1424:4HP:C2	3.03	0.41
1:D:314:SER:HA	1:D:367:ARG:HH21	1.85	0.41
1:B:169:TRP:HH2	1:C:378:MET:HE3	1.85	0.41
1:C:349:PHE:HD2	1:C:353:ASN:HD21	1.69	0.41
1:A:289:GLN:O	1:A:290:ARG:C	2.64	0.41
2:B:1423:FMN:H1'2	2:B:1423:FMN:H4'	1.89	0.41
1:A:131:LEU:HD13	1:A:234:ARG:HD2	2.03	0.41
1:A:169:TRP:HH2	1:D:378:MET:HE3	1.84	0.41
2:A:1423:FMN:O1P	1:D:374:ALA:N	2.53	0.41
1:C:222:MET:HG2	1:C:223:LEU:N	2.36	0.41
1:C:263:ARG:HE	3:C:1424:4HP:C8	2.34	0.41
1:C:266:PHE:CD2	3:C:1424:4HP:C2	3.04	0.41
1:D:164:ASN:ND2	1:D:227:ASN:H	2.19	0.41
1:A:169:TRP:O	2:A:1423:FMN:C4	2.70	0.40
1:C:49:PRO:O	1:C:53:LYS:HG2	2.22	0.40
1:C:353:ASN:HD22	1:D:316:HIS:CD2	2.39	0.40
1:B:384:GLN:HG3	1:C:213:GLN:CG	2.46	0.40
1:D:266:PHE:CD2	3:D:1424:4HP:C2	3.04	0.40
1:A:49:PRO:O	1:A:53:LYS:HG2	2.20	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	398/422 (94%)	389 (98%)	8 (2%)	1 (0%)	36	66
1	B	397/422 (94%)	386 (97%)	11 (3%)	0	100	100
1	C	397/422 (94%)	386 (97%)	11 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	397/422 (94%)	384 (97%)	13 (3%)	0	100	100
All	All	1589/1688 (94%)	1545 (97%)	43 (3%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	ARG

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	325/345 (94%)	308 (95%)	17 (5%)	21	53
1	B	324/345 (94%)	307 (95%)	17 (5%)	21	53
1	C	324/345 (94%)	307 (95%)	17 (5%)	21	53
1	D	324/345 (94%)	307 (95%)	17 (5%)	21	53
All	All	1297/1380 (94%)	1229 (95%)	68 (5%)	21	53

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ILE
1	A	38	VAL
1	A	56	GLU
1	A	153	LYS
1	A	154	VAL
1	A	181	VAL
1	A	236	SER
1	A	260	THR
1	A	288	LYS
1	A	291	ASN
1	A	301	VAL
1	A	318	VAL
1	A	343	ASN

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Mol	Chain	Res	Type
1	A	346	THR
1	A	347	LEU
1	A	384	GLN
1	A	420	THR
1	B	56	GLU
1	B	153	LYS
1	B	154	VAL
1	B	181	VAL
1	B	222	MET
1	B	236	SER
1	B	260	THR
1	B	288	LYS
1	B	291	ASN
1	B	293	VAL
1	B	301	VAL
1	B	318	VAL
1	B	343	ASN
1	B	346	THR
1	B	347	LEU
1	B	384	GLN
1	B	420	THR
1	C	35	VAL
1	C	38	VAL
1	C	56	GLU
1	C	153	LYS
1	C	154	VAL
1	C	181	VAL
1	C	236	SER
1	C	260	THR
1	C	288	LYS
1	C	291	ASN
1	C	301	VAL
1	C	318	VAL
1	C	343	ASN
1	C	346	THR
1	C	347	LEU
1	C	384	GLN
1	C	420	THR
1	D	38	VAL
1	D	56	GLU
1	D	153	LYS
1	D	154	VAL

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Mol	Chain	Res	Type
1	D	181	VAL
1	D	189	ASP
1	D	236	SER
1	D	260	THR
1	D	288	LYS
1	D	291	ASN
1	D	301	VAL
1	D	318	VAL
1	D	343	ASN
1	D	346	THR
1	D	347	LEU
1	D	384	GLN
1	D	420	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	164	ASN
1	A	184	ASN
1	A	227	ASN
1	A	340	GLN
1	A	343	ASN
1	A	353	ASN
1	B	46	GLN
1	B	164	ASN
1	B	184	ASN
1	B	191	ASN
1	B	227	ASN
1	B	340	GLN
1	B	343	ASN
1	B	353	ASN
1	C	46	GLN
1	C	130	GLN
1	C	164	ASN
1	C	184	ASN
1	C	213	GLN
1	C	227	ASN
1	C	340	GLN
1	C	343	ASN
1	C	353	ASN
1	D	46	GLN

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Mol	Chain	Res	Type
1	D	164	ASN
1	D	184	ASN
1	D	213	GLN
1	D	227	ASN
1	D	340	GLN
1	D	343	ASN
1	D	353	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 ⓘ

8 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	4HP	C	1424	-	11,11,11	0.86	0	14,14,14	0.89	0
3	4HP	A	1424	-	11,11,11	0.80	0	14,14,14	0.81	0
3	4HP	D	1424	-	11,11,11	0.80	0	14,14,14	0.77	0
2	FMN	C	1423	-	33,33,33	1.04	2 (6%)	48,50,50	1.37	10 (20%)
2	FMN	A	1423	-	33,33,33	2.50	5 (15%)	48,50,50	1.52	11 (22%)
2	FMN	B	1423	-	33,33,33	1.02	2 (6%)	48,50,50	1.38	9 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	4HP	B	1424	-	11,11,11	0.71	0	14,14,14	0.78	0
2	FMN	D	1423	-	33,33,33	1.02	2 (6%)	48,50,50	1.33	7 (14%)

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	4HP	C	1424	-	-	2/4/4/4	0/1/1/1
3	4HP	A	1424	-	-	2/4/4/4	0/1/1/1
3	4HP	D	1424	-	-	2/4/4/4	0/1/1/1
2	FMN	C	1423	-	-	5/18/18/18	0/3/3/3
2	FMN	A	1423	-	-	4/18/18/18	0/3/3/3
2	FMN	B	1423	-	-	5/18/18/18	0/3/3/3
3	4HP	B	1424	-	-	2/4/4/4	0/1/1/1
2	FMN	D	1423	-	-	5/18/18/18	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1423	FMN	P-O2P	-8.91	1.21	1.54
2	A	1423	FMN	P-O3P	-8.74	1.22	1.54
2	A	1423	FMN	P-O1P	-4.18	1.37	1.50
2	C	1423	FMN	C4A-N5	3.95	1.39	1.30
2	A	1423	FMN	C4A-N5	3.94	1.39	1.30
2	B	1423	FMN	C4A-N5	3.84	1.39	1.30
2	D	1423	FMN	C4A-N5	3.78	1.38	1.30
2	D	1423	FMN	C10-N1	2.68	1.38	1.33
2	A	1423	FMN	C10-N1	2.62	1.38	1.33
2	B	1423	FMN	C10-N1	2.57	1.38	1.33
2	C	1423	FMN	C10-N1	2.43	1.38	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1423	FMN	O2P-P-O5'	3.85	116.70	106.67
2	A	1423	FMN	C4A-C10-N10	3.46	121.44	116.48
2	B	1423	FMN	C4A-C10-N10	3.36	121.29	116.48
2	D	1423	FMN	C4A-C10-N10	3.27	121.17	116.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1423	FMN	C4A-C10-N10	3.21	121.08	116.48
2	B	1423	FMN	C4-N3-C2	-3.19	119.98	125.64
2	C	1423	FMN	C4-N3-C2	-3.17	120.01	125.64
2	A	1423	FMN	C4-N3-C2	-3.06	120.21	125.64
2	D	1423	FMN	C4-N3-C2	-3.04	120.25	125.64
2	C	1423	FMN	C4'-C3'-C2'	-3.02	108.55	113.57
2	A	1423	FMN	C10-C4A-N5	-2.74	119.22	124.81
2	D	1423	FMN	C10-C4A-N5	-2.74	119.22	124.81
2	C	1423	FMN	C4A-C4-N3	2.70	120.12	113.25
2	B	1423	FMN	C10-C4A-N5	-2.64	119.42	124.81
2	B	1423	FMN	C4'-C3'-C2'	-2.63	109.19	113.57
2	D	1423	FMN	C4A-C4-N3	2.61	119.89	113.25
2	A	1423	FMN	C4A-C4-N3	2.59	119.85	113.25
2	C	1423	FMN	C10-C4A-N5	-2.55	119.61	124.81
2	B	1423	FMN	C4A-C4-N3	2.52	119.67	113.25
2	C	1423	FMN	O4-C4-C4A	-2.47	120.00	126.53
2	A	1423	FMN	C4-C4A-N5	2.39	121.51	118.21
2	B	1423	FMN	O2-C2-N1	-2.38	117.84	121.80
2	D	1423	FMN	C4-C4A-N5	2.35	121.46	118.21
2	A	1423	FMN	C4'-C3'-C2'	-2.34	109.67	113.57
2	A	1423	FMN	O4-C4-C4A	-2.28	120.51	126.53
2	B	1423	FMN	O4-C4-C4A	-2.20	120.73	126.53
2	B	1423	FMN	C5A-C9A-N10	2.19	119.95	117.97
2	D	1423	FMN	C4'-C3'-C2'	-2.19	109.92	113.57
2	C	1423	FMN	C5A-C9A-N10	2.18	119.94	117.97
2	B	1423	FMN	C4-C4A-N5	2.18	121.22	118.21
2	C	1423	FMN	C4-C4A-N5	2.17	121.21	118.21
2	A	1423	FMN	O2-C2-N1	-2.17	118.19	121.80
2	D	1423	FMN	O4-C4-C4A	-2.17	120.82	126.53
2	A	1423	FMN	O2P-P-O1P	2.14	119.18	110.83
2	A	1423	FMN	C4A-C10-N1	-2.04	119.58	124.59
2	C	1423	FMN	O2-C2-N1	-2.01	118.47	121.80
2	C	1423	FMN	C4A-C10-N1	-2.00	119.68	124.59

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1423	FMN	C3'-C4'-C5'-O5'
2	A	1423	FMN	O4'-C4'-C5'-O5'
2	B	1423	FMN	C3'-C4'-C5'-O5'
2	B	1423	FMN	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
2	B	1423	FMN	C5'-O5'-P-O1P
2	B	1423	FMN	C5'-O5'-P-O2P
2	B	1423	FMN	C5'-O5'-P-O3P
2	C	1423	FMN	C3'-C4'-C5'-O5'
2	C	1423	FMN	O4'-C4'-C5'-O5'
2	C	1423	FMN	C5'-O5'-P-O1P
2	C	1423	FMN	C5'-O5'-P-O2P
2	C	1423	FMN	C5'-O5'-P-O3P
2	D	1423	FMN	C3'-C4'-C5'-O5'
2	D	1423	FMN	O4'-C4'-C5'-O5'
2	D	1423	FMN	C5'-O5'-P-O1P
2	D	1423	FMN	C5'-O5'-P-O2P
2	D	1423	FMN	C5'-O5'-P-O3P
2	A	1423	FMN	C5'-O5'-P-O1P
3	A	1424	4HP	C1-C7-C8-O2
3	B	1424	4HP	C1-C7-C8-O2
3	C	1424	4HP	C1-C7-C8-O2
3	D	1424	4HP	C1-C7-C8-O2
3	A	1424	4HP	C1-C7-C8-O1
3	B	1424	4HP	C1-C7-C8-O1
3	D	1424	4HP	C1-C7-C8-O1
2	A	1423	FMN	C5'-O5'-P-O3P
3	C	1424	4HP	C1-C7-C8-O1

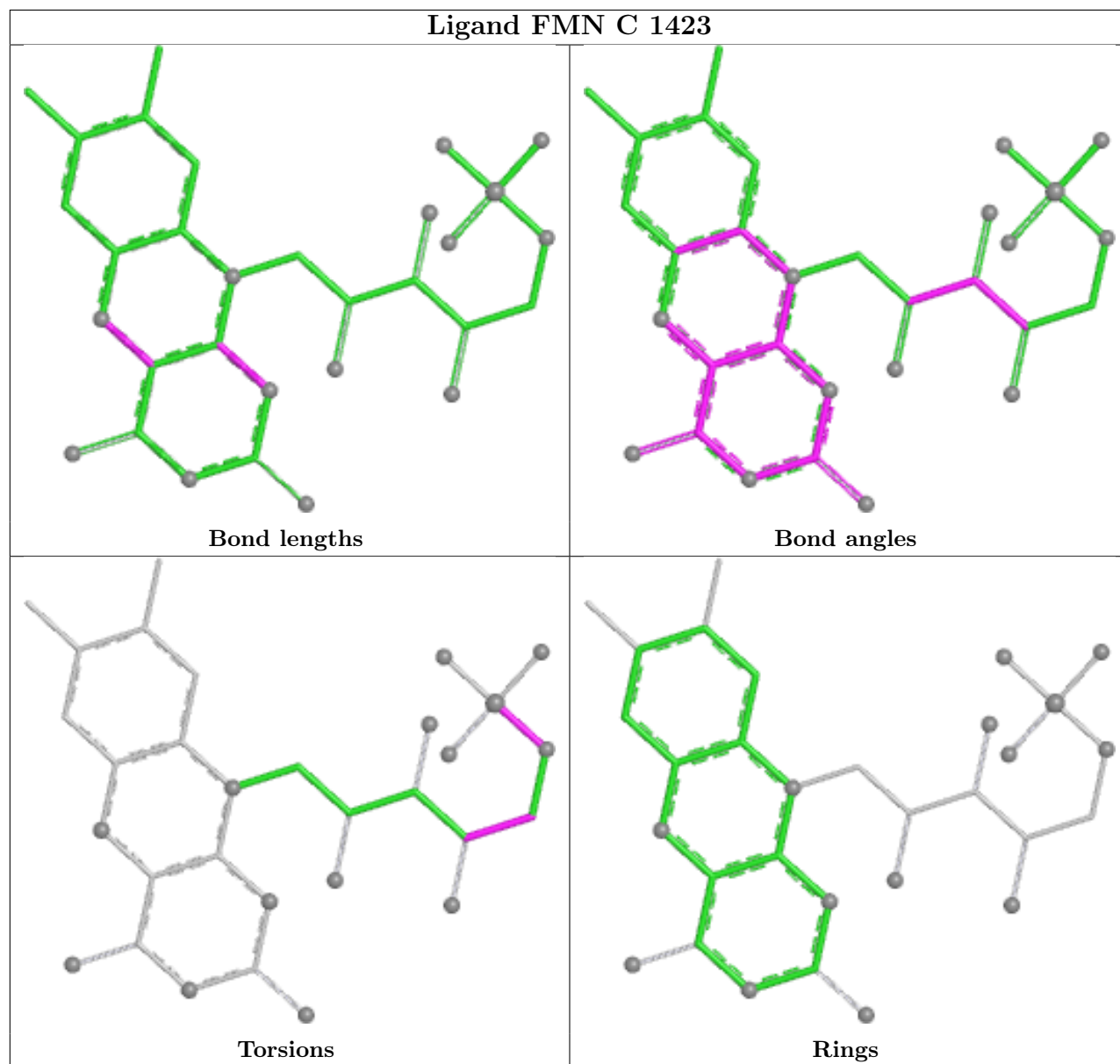
There are no ring outliers.

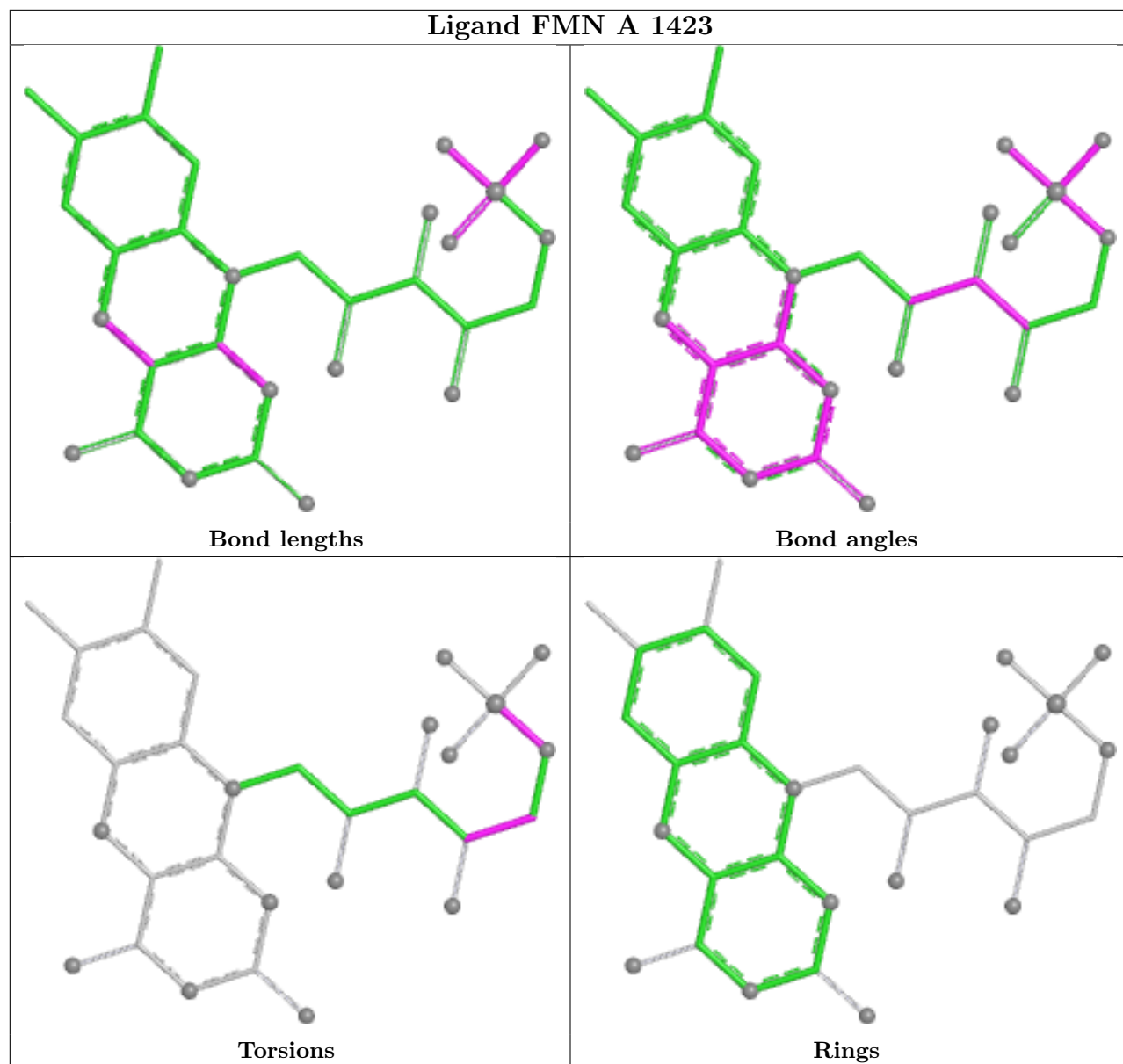
8 monomers are involved in 15 short contacts:

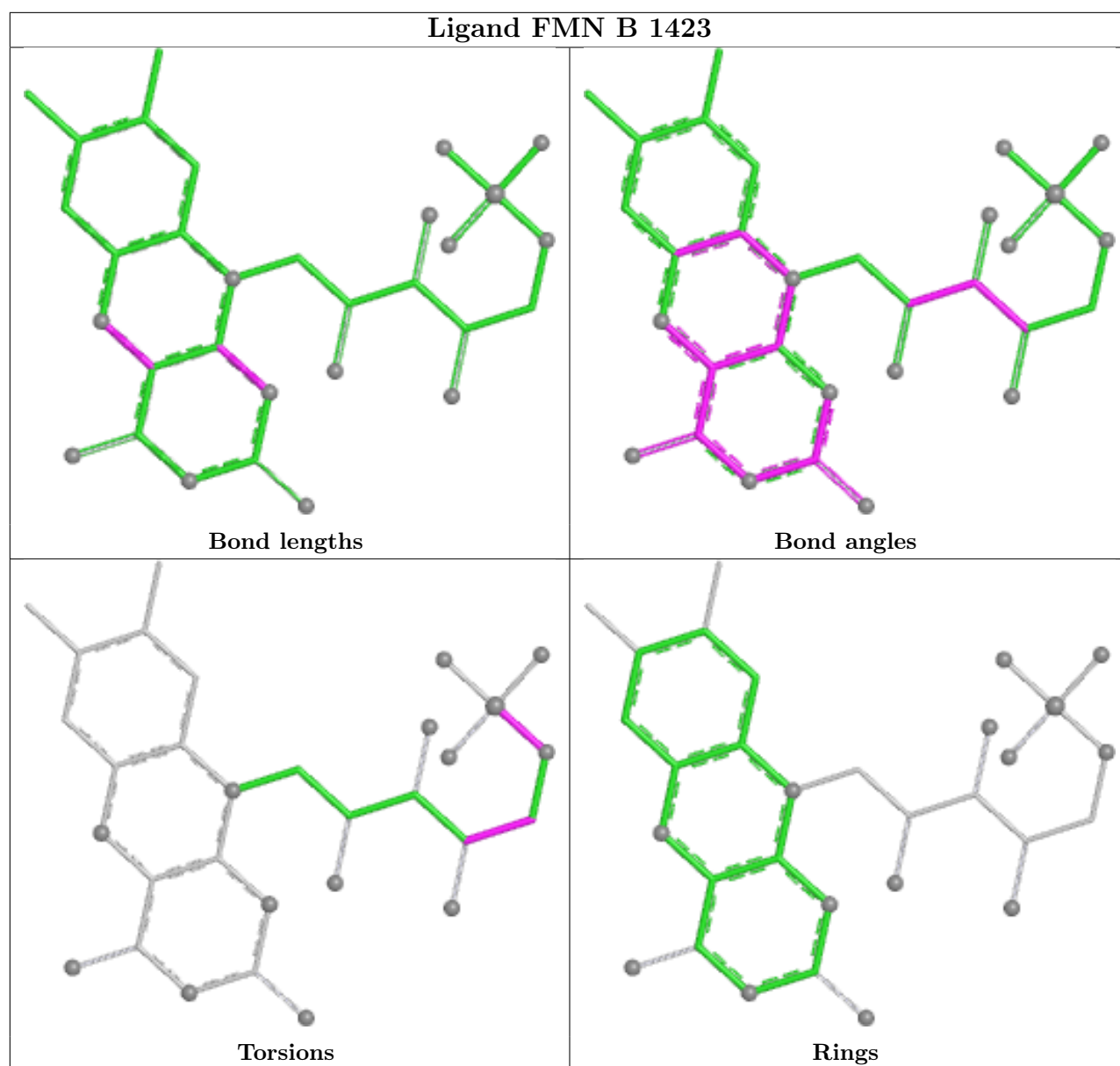
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1424	4HP	3	0
3	A	1424	4HP	2	0
3	D	1424	4HP	2	0
2	C	1423	FMN	1	0
2	A	1423	FMN	2	0
2	B	1423	FMN	2	0
3	B	1424	4HP	2	0
2	D	1423	FMN	1	0

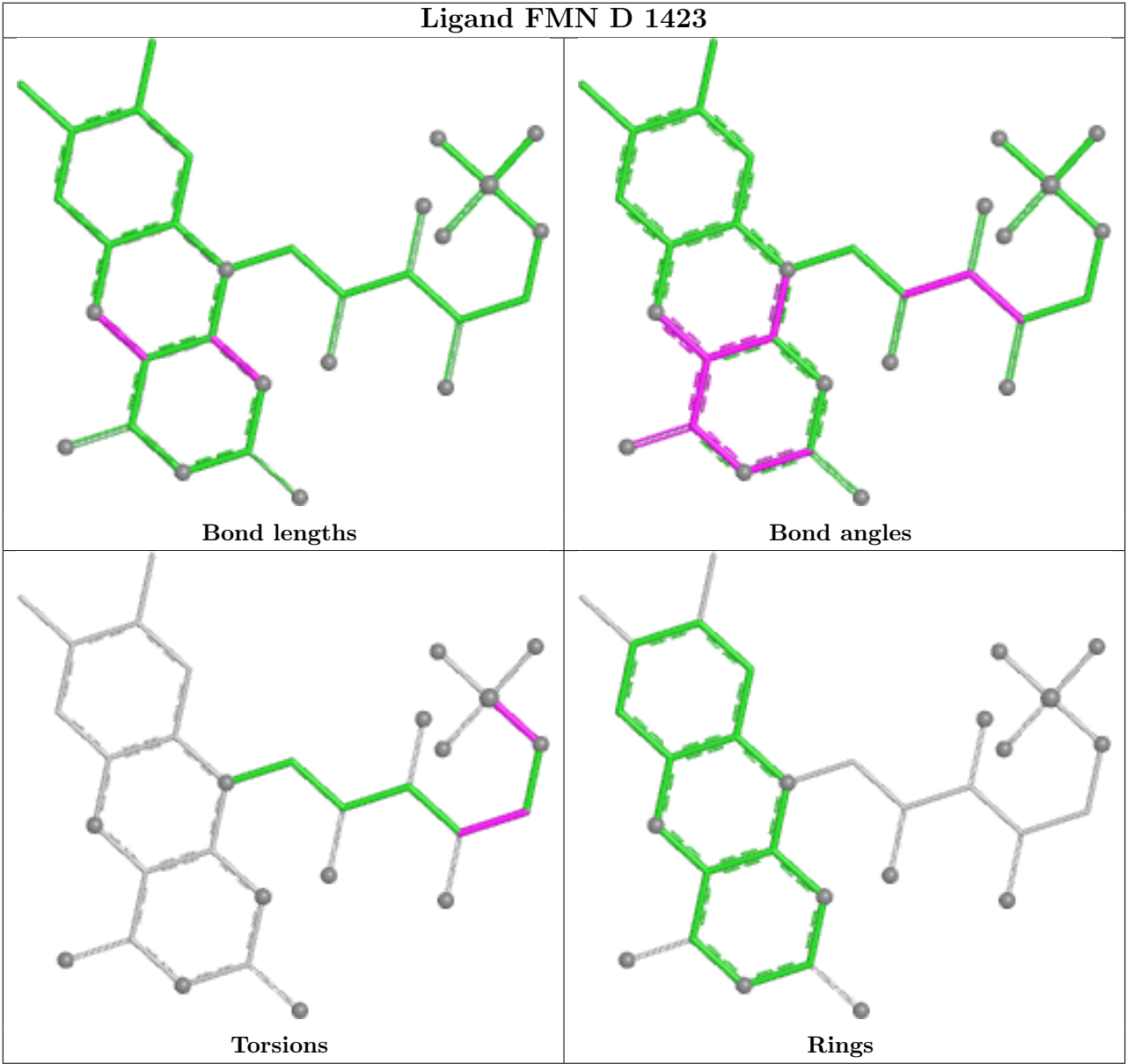
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	374:ALA	C	375:THR	N	1.07

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	400/422 (94%)	0.63	21 (5%)	32	24	43, 56, 75, 85	0
1	B	399/422 (94%)	0.82	20 (5%)	34	26	43, 56, 72, 83	0
1	C	399/422 (94%)	0.67	20 (5%)	34	26	43, 56, 74, 84	0
1	D	399/422 (94%)	1.37	115 (28%)	1	1	43, 56, 72, 83	0
All	All	1597/1688 (94%)	0.87	176 (11%)	10	8	43, 56, 74, 85	0

All (176) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	23	ILE	5.5
1	A	35	VAL	5.1
1	D	130	GLN	5.0
1	D	151	PHE	4.7
1	C	33	PRO	4.4
1	D	35	VAL	4.3
1	D	81	PHE	4.2
1	C	34	ASP	4.2
1	D	136	TRP	4.0
1	B	158	GLU	3.9
1	D	128	SER	3.9
1	D	154	VAL	3.9
1	B	190	GLY	3.8
1	B	60	GLN	3.8
1	D	157	VAL	3.8
1	D	85	VAL	3.7
1	D	259	TYR	3.7
1	D	258	PHE	3.7
1	C	35	VAL	3.6
1	D	38	VAL	3.6
1	D	142	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	209	ASN	3.5
1	A	34	ASP	3.4
1	D	162	ILE	3.4
1	D	143	THR	3.3
1	D	83	PRO	3.3
1	D	168	GLY	3.3
1	D	252	TYR	3.2
1	C	254	ASP	3.2
1	D	254	ASP	3.2
1	D	137	LEU	3.2
1	D	176	ALA	3.2
1	C	46	GLN	3.2
1	D	86	TYR	3.2
1	D	167	TYR	3.2
1	C	79	ARG	3.1
1	D	193	ILE	3.1
1	D	159	GLY	3.1
1	D	36	SER	3.1
1	A	33	PRO	3.0
1	D	198	VAL	3.0
1	D	126	MET	3.0
1	D	222	MET	3.0
1	D	178	TYR	3.0
1	D	233	TYR	3.0
1	D	123	GLN	3.0
1	D	135	ILE	3.0
1	A	294	ARG	3.0
1	D	230	ILE	3.0
1	B	289	GLN	3.0
1	D	301	VAL	3.0
1	D	152	GLY	2.9
1	D	33	PRO	2.9
1	D	144	ALA	2.9
1	A	36	SER	2.9
1	D	160	GLY	2.8
1	C	36	SER	2.8
1	A	293	VAL	2.8
1	D	246	SER	2.8
1	C	31	GLN	2.8
1	C	32	THR	2.7
1	D	79	ARG	2.7
1	B	159	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	336	GLY	2.7
1	C	124	ILE	2.7
1	A	290	ARG	2.7
1	D	205	GLU	2.7
1	D	199	ILE	2.7
1	D	247	ALA	2.7
1	D	70	LYS	2.7
1	A	297	THR	2.7
1	B	36	SER	2.7
1	D	131	LEU	2.7
1	D	188	ALA	2.6
1	D	341	TYR	2.6
1	D	342	PRO	2.6
1	A	295	ALA	2.6
1	D	161	ILE	2.6
1	D	340	GLN	2.6
1	D	127	PHE	2.6
1	D	251	LEU	2.6
1	A	85	VAL	2.6
1	B	340	GLN	2.6
1	D	187	ASP	2.6
1	B	24	ARG	2.5
1	D	234	ARG	2.5
1	D	231	PRO	2.5
1	C	38	VAL	2.5
1	D	257	ILE	2.5
1	D	249	PHE	2.5
1	D	82	GLN	2.5
1	D	229	PHE	2.5
1	D	78	HIS	2.5
1	D	140	PRO	2.5
1	D	250	GLY	2.5
1	B	205	GLU	2.5
1	D	207	VAL	2.5
1	D	171	SER	2.4
1	A	50	GLN	2.4
1	D	84	LYS	2.4
1	C	127	PHE	2.4
1	A	296	TYR	2.4
1	D	422	VAL	2.4
1	B	229	PHE	2.4
1	D	169	TRP	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	124	ILE	2.4
1	D	129	LYS	2.4
1	D	141	ASP	2.4
1	D	39	SER	2.4
1	A	32	THR	2.4
1	D	163	LEU	2.4
1	B	399	THR	2.3
1	D	92	SER	2.3
1	D	49	PRO	2.3
1	C	85	VAL	2.3
1	D	37	GLY	2.3
1	A	151	PHE	2.3
1	D	186	PHE	2.3
1	D	155	GLU	2.3
1	D	77	LEU	2.3
1	D	72	LEU	2.3
1	D	88	GLY	2.3
1	B	42	GLU	2.2
1	D	134	GLU	2.2
1	D	61	LEU	2.2
1	D	335	HIS	2.2
1	C	420	THR	2.2
1	D	150	PRO	2.2
1	D	346	THR	2.2
1	D	165	GLY	2.2
1	D	153	LYS	2.2
1	D	200	PRO	2.2
1	B	151	PHE	2.2
1	A	299	ALA	2.2
1	B	157	VAL	2.2
1	D	173	CYS	2.2
1	D	223	LEU	2.2
1	B	156	GLU	2.2
1	D	206	ILE	2.2
1	D	235	ILE	2.2
1	D	132	GLN	2.2
1	D	34	ASP	2.2
1	C	176	ALA	2.2
1	B	288	LYS	2.1
1	D	96	PHE	2.1
1	D	75	ILE	2.1
1	C	289	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	344	LYS	2.1
1	D	138	LYS	2.1
1	B	249	PHE	2.1
1	D	225	LEU	2.1
1	D	185	ARG	2.1
1	D	204	TYR	2.1
1	D	228	VAL	2.1
1	C	257	ILE	2.1
1	D	125	ALA	2.1
1	C	296	TYR	2.1
1	D	256	LYS	2.1
1	D	65	PRO	2.1
1	A	137	LEU	2.1
1	D	87	GLY	2.1
1	D	270	PHE	2.1
1	A	188	ALA	2.1
1	C	337	LEU	2.0
1	D	117	LEU	2.0
1	B	186	PHE	2.0
1	D	190	GLY	2.0
1	A	210	TRP	2.0
1	D	192	LYS	2.0
1	D	241	MET	2.0
1	A	255	SER	2.0
1	B	181	VAL	2.0
1	D	236	SER	2.0
1	B	362	ILE	2.0
1	C	161	ILE	2.0
1	D	240	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

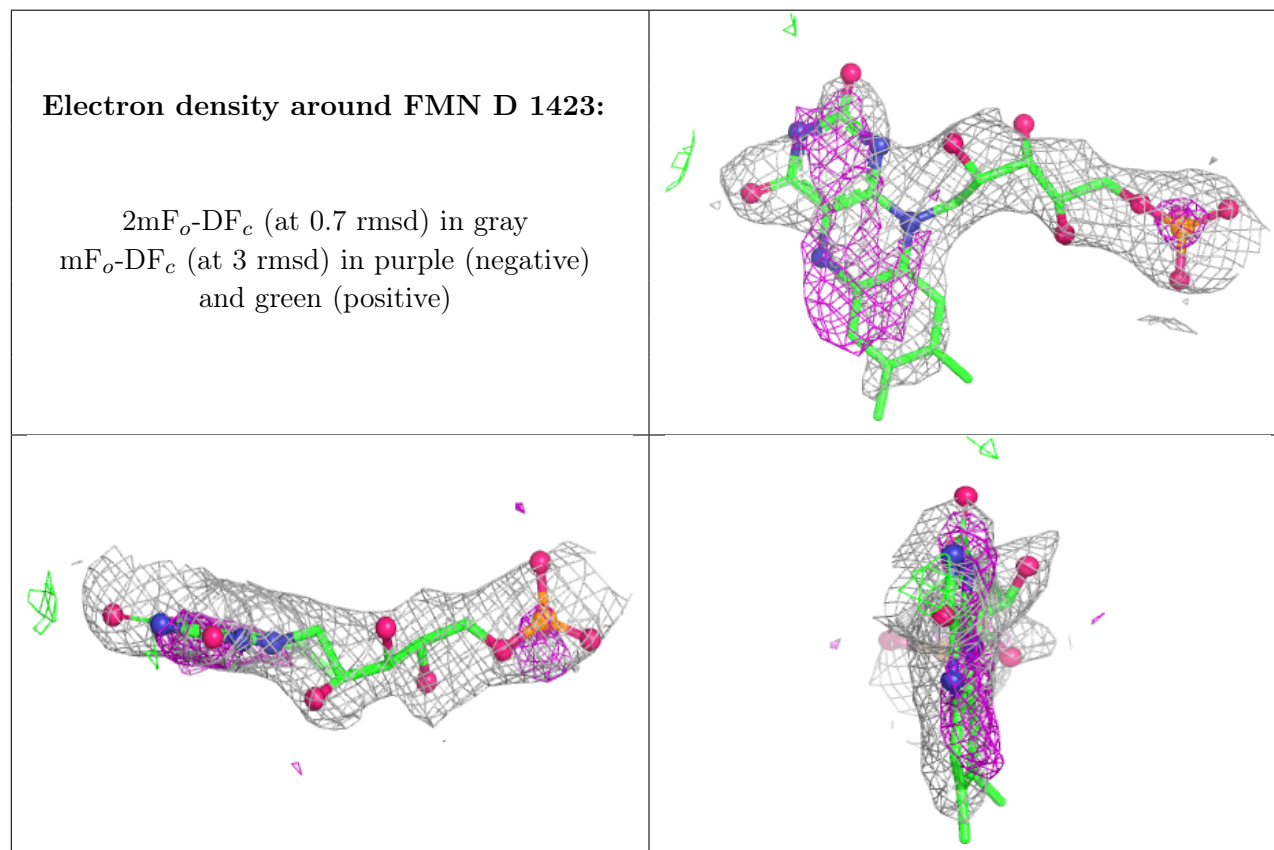
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

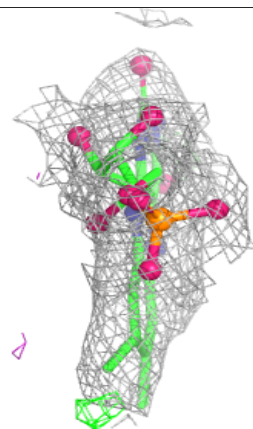
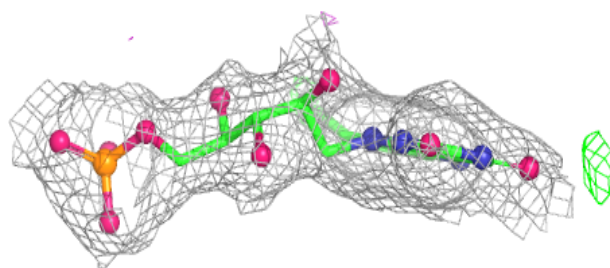
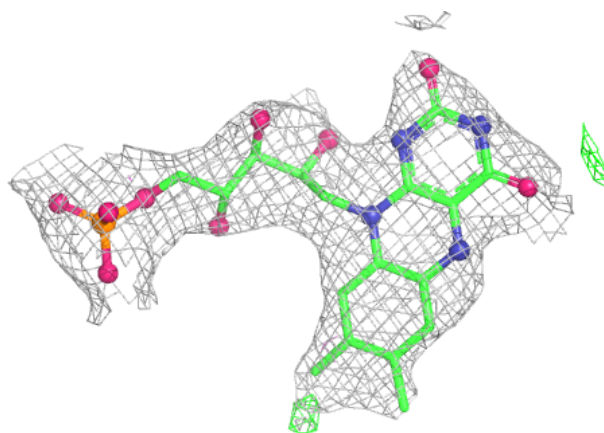
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FMN	D	1423	31/31	0.77	0.17	70,71,81,82	0
3	4HP	D	1424	11/11	0.79	0.20	78,79,81,82	0
3	4HP	A	1424	11/11	0.80	0.23	78,79,81,82	0
3	4HP	C	1424	11/11	0.85	0.17	78,79,81,82	0
2	FMN	C	1423	31/31	0.89	0.14	70,71,80,83	0
2	FMN	B	1423	31/31	0.90	0.13	70,71,80,83	0
2	FMN	A	1423	31/31	0.91	0.12	70,71,80,82	0
3	4HP	B	1424	11/11	0.92	0.17	78,79,81,82	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.



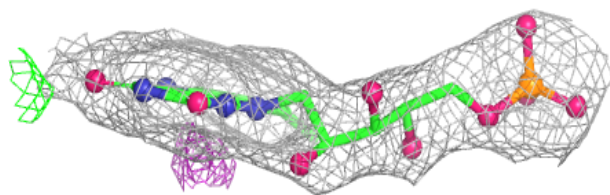
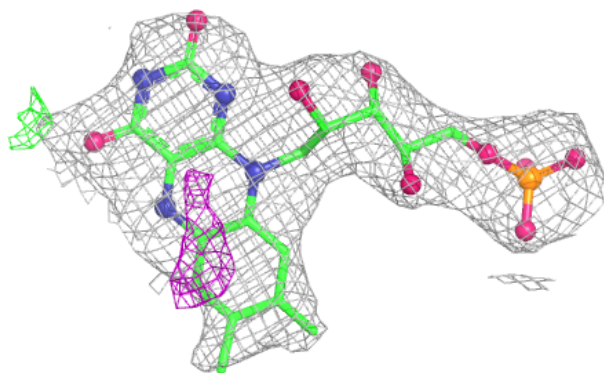
Electron density around FMN C 1423:

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

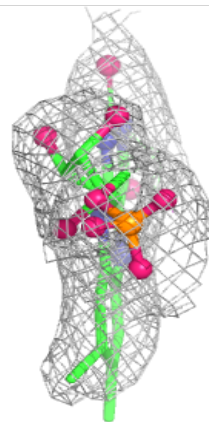
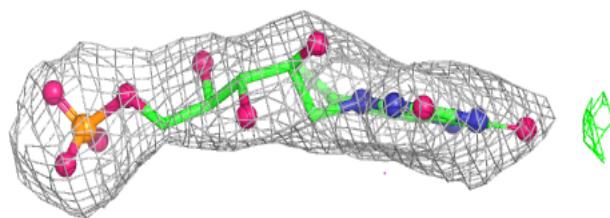
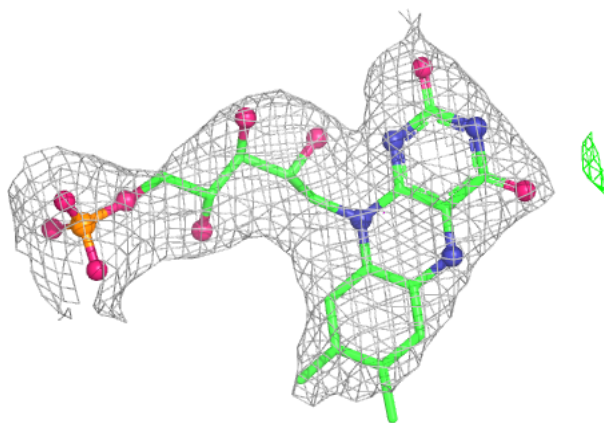


Electron density around FMN B 1423:

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

**Electron density around FMN A 1423:**

$2mF_o-DF_c$ (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.