



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:40 PM UTC

PDB ID : 3NJO / pdb_00003njo
Title : X-ray crystal structure of the Pyr1-pyrabactin A complex
Authors : Burgie, E.S.; Bingman, C.A.; Phillips Jr., G.N.; Peterson, F.C.; Volkman, B.F.; Cutler, S.R.; Jensen, D.R.; Center for Eukaryotic Structural Genomics (CESG)
Deposited on : 2010-06-17
Resolution : 2.47 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

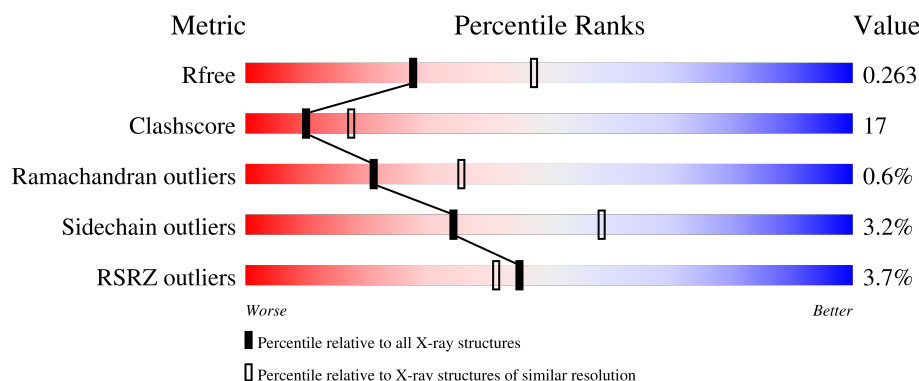
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	194	3% 54% 29% • 13%
1	B	194	3% 64% 23% • 9%
1	C	194	5% 59% 22% 6% • 12%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4368 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.

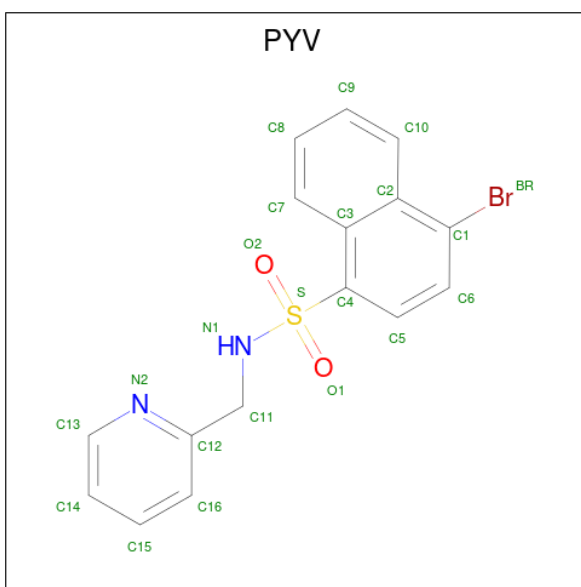
- Molecule 1 is a protein called Absciscic acid receptor PYR1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	169	Total	C	N	O	S	0	2	0
			1370	856	245	261	8			
1	B	177	Total	C	N	O	S	0	0	0
			1418	884	253	274	7			
1	C	171	Total	C	N	O	S	0	3	0
			1404	878	253	266	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP O49686
A	-1	SER	-	expression tag	UNP O49686
A	0	HIS	-	expression tag	UNP O49686
A	88	SER	PRO	engineered mutation	UNP O49686
B	-2	GLY	-	expression tag	UNP O49686
B	-1	SER	-	expression tag	UNP O49686
B	0	HIS	-	expression tag	UNP O49686
B	88	SER	PRO	engineered mutation	UNP O49686
C	-2	GLY	-	expression tag	UNP O49686
C	-1	SER	-	expression tag	UNP O49686
C	0	HIS	-	expression tag	UNP O49686
C	88	SER	PRO	engineered mutation	UNP O49686

- Molecule 2 is 4-bromo-N-(pyridin-2-ylmethyl)naphthalene-1-sulfonamide (CCD ID: PYV) (formula: C₁₆H₁₃BrN₂O₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	Br	C	N	O	S	0	1
			31	1	22	2	4	2		
2	C	1	Total	Br	C	N	O	S	1	1
			44	2	32	4	4	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

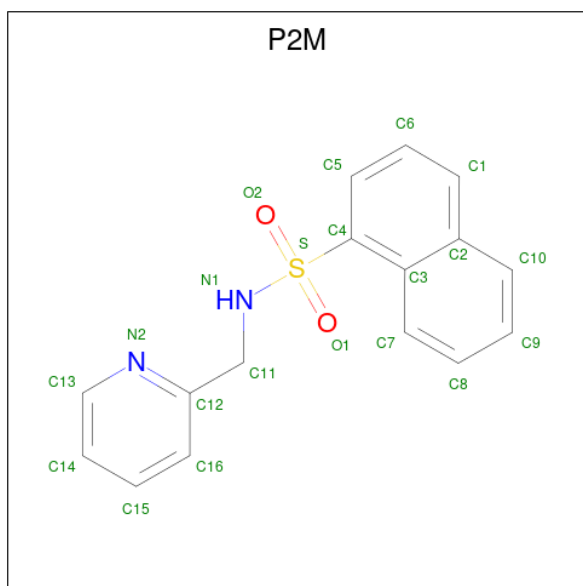


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 5 is N-(pyridin-2-ylmethyl)naphthalene-1-sulfonamide (CCD ID: P2M) (formula: C₁₆H₁₄N₂O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			21	16	2	2	1		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Cl	0	0
			1	1		

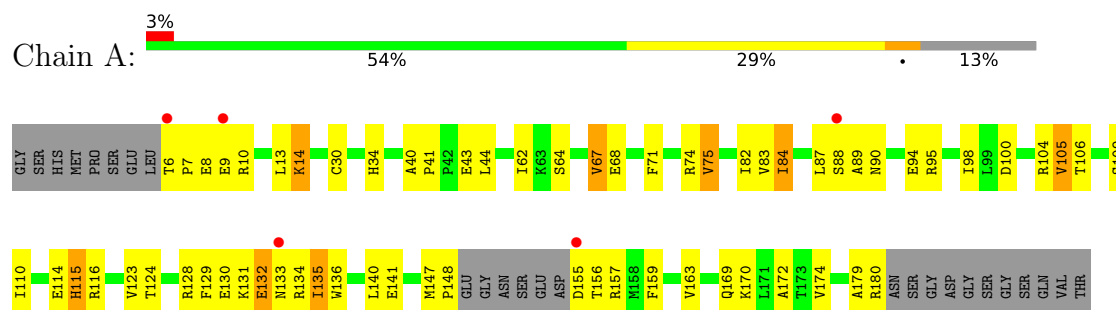
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	25	Total	O	0	3
			25	25		
7	B	12	Total	O	0	1
			12	12		
7	C	31	Total	O	0	0
			31	31		

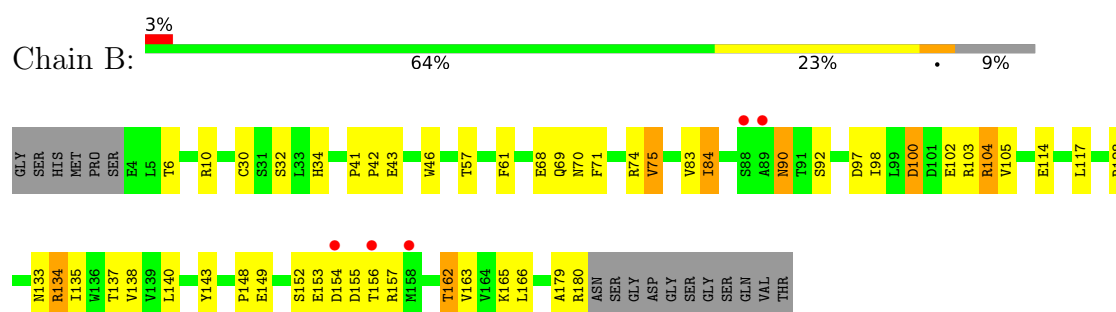
3 Residue-property plots [i](#)

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.

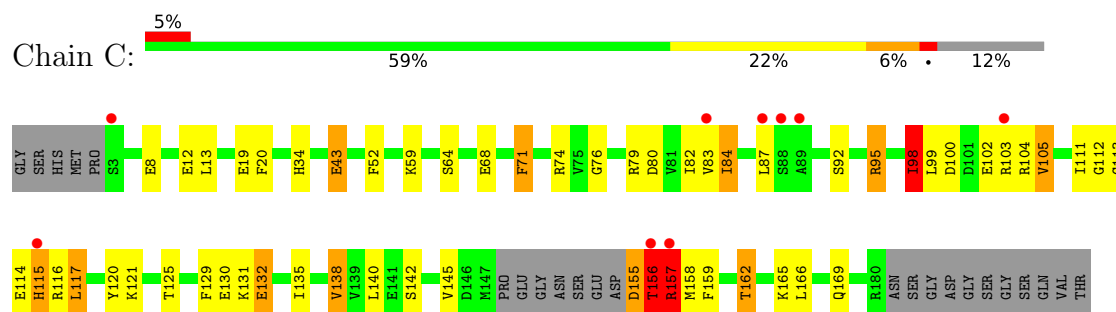
• Molecule 1: Absciscic acid receptor PYR1



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• Molecule 1: Absciscic acid receptor PYR1



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	60.31Å 60.31Å 527.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.05 – 2.47 32.05 – 2.47	Depositor EDS
% Data completeness (in resolution range)	89.8 (32.05-2.47) 89.7 (32.05-2.47)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.222 , 0.265 0.220 , 0.263	Depositor DCC
R_{free} test set	2000 reflections (9.54%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.560	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 64.1	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	0.93	EDS
Total number of atoms	4368	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.79% 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: CL, SO4, P2M, PYV, NA

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	1.75	10/1400 (0.7%)	1.15	10/1893 (0.5%)
1	B	1.59	15/1444 (1.0%)	1.06	9/1954 (0.5%)
1	C	1.46	14/1428 (1.0%)	1.02	11/1928 (0.6%)
All	All	1.60	39/4272 (0.9%)	1.08	30/5775 (0.5%)

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	41	PRO	C-O	-7.10	1.19	1.25
1	B	100	ASP	CA-C	-7.00	1.45	1.53
1	B	75	VAL	CA-CB	-6.93	1.47	1.54
1	C	105	VAL	C-O	-6.83	1.17	1.24
1	B	41	PRO	CA-C	-6.64	1.48	1.51
1	A	140	LEU	C-O	-6.49	1.16	1.24
1	C	117	LEU	CA-C	-6.40	1.45	1.53
1	B	149	GLU	N-CA	6.25	1.54	1.46
1	B	162	THR	C-O	-6.21	1.15	1.24
1	C	138	VAL	C-O	-6.17	1.17	1.24
1	C	19	GLU	N-CA	-6.13	1.38	1.46
1	A	123	VAL	C-O	-6.11	1.17	1.24
1	C	98	ILE	C-O	-6.10	1.16	1.23
1	B	92	SER	C-O	-6.04	1.17	1.24
1	C	162	THR	C-O	-6.03	1.17	1.24
1	B	30	CYS	C-O	-5.91	1.16	1.23
1	C	84	ILE	CA-CB	-5.79	1.48	1.54
1	C	95[A]	ARG	C-O	-5.68	1.17	1.24
1	C	95[B]	ARG	C-O	-5.68	1.17	1.24
1	A	105	VAL	C-O	-5.62	1.18	1.24
1	C	142	SER	C-O	-5.58	1.16	1.23
1	C	140	LEU	C-O	-5.52	1.17	1.24
1	B	105	VAL	C-O	-5.50	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	140	LEU	C-O	-5.38	1.17	1.24
1	A	75	VAL	C-O	-5.38	1.17	1.24
1	A	95	ARG	C-O	-5.21	1.17	1.23
1	B	70	ASN	CA-C	-5.21	1.47	1.53
1	C	34	HIS	C-O	-5.19	1.17	1.23
1	B	165	LYS	C-O	-5.19	1.18	1.24
1	A	141	GLU	C-O	-5.18	1.18	1.24
1	B	148	PRO	N-CD	5.15	1.54	1.47
1	A	34	HIS	C-O	-5.12	1.17	1.23
1	B	163	VAL	C-O	-5.11	1.18	1.24
1	A	124	THR	C-O	-5.10	1.17	1.23
1	C	121[A]	LYS	C-O	-5.05	1.17	1.23
1	C	121[B]	LYS	C-O	-5.05	1.17	1.23
1	A	8	GLU	C-N	5.04	1.40	1.33
1	A	172	ALA	CA-CB	-5.04	1.45	1.53
1	B	34	HIS	C-O	-5.02	1.17	1.23

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	PRO	CA-C-N	11.27	135.76	120.44
1	A	7	PRO	C-N-CA	11.27	135.76	120.44
1	B	148	PRO	CA-N-CD	-9.53	98.66	112.00
1	B	41	PRO	O-C-N	8.30	125.13	121.31
1	C	155	ASP	CB-CA-C	7.47	124.30	110.10
1	B	134	ARG	N-CA-C	7.36	120.60	109.41
1	C	19	GLU	N-CA-CB	-7.31	99.80	110.47
1	A	132	GLU	N-CA-C	-7.23	95.40	110.80
1	C	132	GLU	CB-CA-C	-6.90	106.68	116.34
1	C	157	ARG	N-CA-C	-6.69	104.84	112.87
1	C	156	THR	N-CA-C	-6.45	105.06	113.12
1	B	148	PRO	CB-CA-C	6.42	119.37	110.95
1	C	115	HIS	N-CA-C	-6.08	102.51	110.53
1	C	84	ILE	N-CA-C	-5.84	101.33	110.09
1	C	130	GLU	N-CA-C	5.73	118.74	109.40
1	C	114	GLU	N-CA-C	-5.65	106.61	112.93
1	B	153	GLU	N-CA-C	-5.61	104.38	111.11
1	A	115	HIS	N-CA-C	5.59	115.41	108.19
1	A	8	GLU	CA-C-O	-5.59	114.94	120.70
1	A	135	ILE	CB-CA-C	-5.54	102.20	111.29
1	B	41	PRO	CA-C-N	5.51	125.03	119.19
1	B	41	PRO	C-N-CA	5.51	125.03	119.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	109	SER	CA-C-N	-5.43	115.36	122.37
1	A	109	SER	C-N-CA	-5.43	115.36	122.37
1	C	71	PHE	N-CA-C	5.42	118.33	110.42
1	A	156	THR	N-CA-C	5.41	119.59	112.89
1	A	41	PRO	O-C-N	5.27	123.63	121.15
1	B	152	SER	N-CA-C	-5.22	106.89	113.20
1	B	90	ASN	CB-CA-C	-5.17	104.21	111.65
1	C	116	ARG	N-CA-C	-5.05	106.43	112.59

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	1370	0	1355	54	0
1	B	1418	0	1395	33	0
1	C	1404	0	1398	52	0
2	A	31	0	24	5	0
2	C	44	0	26	9	0
3	A	5	0	0	1	0
3	C	5	0	0	0	0
4	A	1	0	0	0	0
5	B	21	0	14	2	0
6	B	1	0	0	0	0
7	A	25	0	0	1	1
7	B	12	0	0	1	1
7	C	31	0	0	2	0
All	All	4368	0	4212	148	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 17.

All (148) 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:67:VAL:HG23	1:A:68:GLU:H	1.13	1.14
1:C:157:ARG:HH21	1:C:157:ARG:CG	1.74	1.00
1:C:117:LEU:CD1	1:C:145:VAL:HG21	1.98	0.92
2:A:300[A]:PYV:H7	2:A:300[A]:PYV:O2	1.69	0.90
2:A:300[B]:PYV:O2	2:A:300[B]:PYV:H7	1.69	0.90
1:A:130:GLU:HG2	1:A:135:ILE:HG12	1.55	0.88
1:A:67:VAL:HG23	1:A:68:GLU:N	1.89	0.88
1:C:157:ARG:HG3	1:C:157:ARG:NH2	1.91	0.85
1:C:87:LEU:HD13	2:C:300[A]:PYV:BR	2.36	0.81
1:C:117:LEU:HD11	1:C:145:VAL:HG21	1.61	0.81
1:C:157:ARG:HH21	1:C:157:ARG:HG2	1.48	0.79
1:A:104:ARG:HH12	1:A:128:ARG:HH21	1.29	0.78
1:C:157:ARG:HH21	1:C:157:ARG:HG3	1.44	0.78
1:A:135:ILE:HG22	1:A:135:ILE:O	1.81	0.77
1:A:104:ARG:HH12	1:A:128:ARG:NH2	1.81	0.77
1:A:128:ARG:O	1:A:129:PHE:CD1	2.38	0.76
1:C:131:LYS:O	1:C:132:GLU:HB2	1.86	0.76
1:C:155:ASP:HA	1:C:158:MET:HE2	1.67	0.76
1:A:104:ARG:NH1	1:A:128:ARG:HH21	1.83	0.75
1:C:157:ARG:CG	1:C:157:ARG:NH2	2.40	0.75
2:A:300[A]:PYV:O2	2:A:300[A]:PYV:C7	2.30	0.74
2:A:300[B]:PYV:O2	2:A:300[B]:PYV:C7	2.30	0.74
1:C:117:LEU:HD12	1:C:145:VAL:HG21	1.70	0.74
1:B:74:ARG:HB2	7:B:195:HOH:O	1.86	0.73
1:C:83:VAL:HB	1:C:87:LEU:HD12	1.71	0.73
1:A:131:LYS:HG2	1:A:132:GLU:H	1.55	0.71
2:C:300[A]:PYV:O2	2:C:300[A]:PYV:H7	1.90	0.71
2:C:300[B]:PYV:H7	2:C:300[B]:PYV:O2	1.90	0.71
1:B:155:ASP:OD1	1:B:155:ASP:C	2.33	0.70
2:C:300[A]:PYV:O2	2:C:300[A]:PYV:C7	2.38	0.68
2:C:300[B]:PYV:O2	2:C:300[B]:PYV:C7	2.38	0.68
1:B:68:GLU:O	1:B:71:PHE:HB2	1.94	0.67
1:C:98:ILE:HG13	1:C:99:LEU:N	2.08	0.67
5:B:300:P2M:O2	5:B:300:P2M:H7	1.93	0.66
1:A:104:ARG:NH1	1:A:128:ARG:NH2	2.42	0.66
1:C:83:VAL:HG23	1:C:84:ILE:O	1.96	0.66
1:C:159:PHE:HZ	2:C:300[A]:PYV:BR	2.34	0.65
1:B:155:ASP:OD1	1:B:156:THR:N	2.30	0.65
1:C:117:LEU:CD1	1:C:145:VAL:CG2	2.73	0.64
1:A:10:ARG:HG3	7:A:202:HOH:O	1.98	0.64
1:B:68:GLU:HG3	1:B:69:GLN:H	1.62	0.63
1:C:117:LEU:HD12	1:C:145:VAL:CG2	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:VAL:C	1:A:84:ILE:O	2.40	0.62
1:B:68:GLU:HG3	1:B:69:GLN:N	2.15	0.62
1:C:98:ILE:CG1	1:C:99:LEU:N	2.61	0.62
1:C:20:PHE:CZ	1:C:103:ARG:HG2	2.35	0.62
1:A:94:GLU:HG2	1:A:110:ILE:HG12	1.84	0.60
1:C:8:GLU:O	1:C:12:GLU:HG2	2.02	0.60
1:C:59:LYS:NZ	7:C:219:HOH:O	2.34	0.60
1:A:157:ARG:HD2	1:C:157:ARG:HD3	1.85	0.59
1:B:32:SER:HB2	1:B:143:TYR:CE2	2.38	0.59
1:B:42:PRO:CG	1:B:128:ARG:HG3	2.33	0.58
1:B:117:LEU:HD12	1:B:117:LEU:O	2.03	0.58
1:A:128:ARG:C	1:A:129:PHE:CD1	2.82	0.58
1:A:131:LYS:O	1:A:132:GLU:C	2.47	0.58
1:C:117:LEU:HG	1:C:120:TYR:HB2	1.85	0.58
1:B:100:ASP:OD1	1:B:102:GLU:HB3	2.04	0.58
1:A:83:VAL:HG12	1:A:84:ILE:N	2.18	0.57
1:A:89:ALA:HB1	1:A:115:HIS:HB3	1.87	0.57
1:C:117:LEU:HD11	1:C:145:VAL:CG2	2.33	0.56
1:A:68:GLU:O	1:A:71:PHE:CB	2.53	0.56
1:B:162:THR:O	1:B:166:LEU:HG	2.06	0.56
1:C:82:ILE:HD12	1:C:82:ILE:N	2.20	0.56
1:C:103:ARG:HG3	1:C:105:VAL:CG2	2.36	0.55
1:C:105:VAL:HG13	1:C:125:THR:HG22	1.86	0.55
1:C:117:LEU:HD21	1:C:120:TYR:CD1	2.41	0.55
1:C:98:ILE:HG13	1:C:99:LEU:H	1.72	0.55
5:B:300:P2M:O2	5:B:300:P2M:C7	2.53	0.55
1:A:40:ALA:HB1	1:A:44:LEU:HD23	1.90	0.54
1:A:132:GLU:O	1:A:134:ARG:N	2.34	0.54
1:A:131:LYS:O	1:A:133:ASN:N	2.40	0.54
1:A:159:PHE:HB2	1:B:84:ILE:CD1	2.37	0.54
1:B:42:PRO:HG2	1:B:128:ARG:HG3	1.89	0.54
1:A:74:ARG:O	1:A:75:VAL:C	2.49	0.53
1:B:97:ASP:C	1:B:98:ILE:HG13	2.33	0.53
1:C:156:THR:C	1:C:158:MET:N	2.65	0.52
1:A:159:PHE:HB2	1:B:84:ILE:HD11	1.92	0.52
1:C:162:THR:O	1:C:166:LEU:HG	2.11	0.51
1:C:12:GLU:HG3	1:C:129:PHE:HE2	1.75	0.51
1:A:43:GLU:OE2	1:A:128:ARG:NH2	2.37	0.51
1:A:67:VAL:CG2	1:A:68:GLU:N	2.61	0.51
1:B:103:ARG:O	1:B:104:ARG:HB2	2.11	0.50
1:B:133:ASN:OD1	1:B:134:ARG:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ASP:HB3	1:A:105:VAL:HG12	1.93	0.50
1:B:154:ASP:OD1	1:B:154:ASP:O	2.30	0.50
1:A:83:VAL:O	1:A:84:ILE:C	2.54	0.50
1:A:83:VAL:O	1:A:84:ILE:O	2.29	0.50
1:A:170:LYS:O	1:A:174:VAL:HG23	2.12	0.50
1:C:68:GLU:O	1:C:71:PHE:HB2	2.12	0.50
1:B:61:PHE:O	1:B:83:VAL:HG23	2.12	0.50
1:A:131:LYS:HB3	1:A:136:TRP:HZ3	1.76	0.50
1:A:179:ALA:O	1:A:180:ARG:O	2.30	0.49
1:B:83:VAL:O	1:B:84:ILE:C	2.55	0.49
1:A:131:LYS:HB3	1:A:136:TRP:CZ3	2.48	0.49
1:A:89:ALA:CB	1:A:115:HIS:HB3	2.42	0.48
1:A:90:ASN:HB2	1:A:114:GLU:HB2	1.95	0.48
1:A:88:SER:O	1:A:89:ALA:C	2.55	0.48
1:C:100:ASP:OD1	1:C:102:GLU:HB3	2.14	0.48
1:A:83:VAL:CG1	1:A:84:ILE:N	2.78	0.47
1:A:68:GLU:O	1:A:71:PHE:HB2	2.12	0.47
1:B:155:ASP:OD1	1:B:157:ARG:N	2.30	0.47
1:C:165:LYS:O	1:C:169:GLN:HG3	2.15	0.47
1:C:43:GLU:CD	1:C:43:GLU:H	2.23	0.46
1:B:6:THR:O	1:B:10:ARG:HG3	2.16	0.46
1:C:156:THR:C	1:C:158:MET:H	2.23	0.46
1:B:46:TRP:CG	1:B:104:ARG:HD3	2.51	0.46
1:B:154:ASP:O	1:B:154:ASP:CG	2.58	0.46
1:C:103:ARG:HG3	1:C:105:VAL:HG23	1.98	0.46
1:C:155:ASP:O	1:C:158:MET:HB3	2.16	0.45
1:B:83:VAL:O	1:B:84:ILE:O	2.33	0.45
1:B:83:VAL:C	1:B:84:ILE:O	2.56	0.45
1:B:90:ASN:OD1	1:B:114:GLU:OE1	2.34	0.45
1:C:20:PHE:HZ	1:C:103:ARG:HG2	1.80	0.45
1:C:103:ARG:O	1:C:104:ARG:HB2	2.15	0.45
1:B:42:PRO:HG3	1:B:128:ARG:HG3	1.99	0.44
1:A:87:LEU:HD13	2:A:300[A]:PYV:BR	2.72	0.44
1:A:147:MET:HA	1:A:148:PRO:HD3	1.69	0.44
1:A:14:LYS:HZ3	1:A:14:LYS:HG3	1.45	0.44
1:B:179:ALA:C	1:B:180:ARG:O	2.60	0.43
1:B:43:GLU:H	1:B:43:GLU:CD	2.27	0.43
1:A:6:THR:OG1	1:A:9:GLU:HG3	2.19	0.43
1:C:112:GLY:O	1:C:115:HIS:HD2	2.01	0.43
1:A:115:HIS:HB2	1:A:116:ARG:H	1.73	0.43
1:A:131:LYS:HB2	1:A:136:TRP:CH2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:LYS:HB2	1:A:136:TRP:HH2	1.83	0.43
1:C:64:SER:HB2	7:C:208:HOH:O	2.19	0.42
1:A:30[A]:CYS:SG	1:A:147:MET:CG	3.07	0.42
1:C:76:GLY:HA2	1:C:95[A]:ARG:NH1	2.34	0.42
1:A:159:PHE:O	1:A:163:VAL:HG23	2.18	0.42
1:A:98:ILE:O	1:A:106:THR:HA	2.19	0.42
1:A:170:LYS:NZ	3:A:192:SO4:O1	2.52	0.42
1:C:113:GLY:C	1:C:115:HIS:H	2.27	0.41
1:A:67:VAL:CG2	1:A:68:GLU:H	1.95	0.41
1:B:137:THR:HG22	1:B:138:VAL:N	2.34	0.41
1:C:82:ILE:HD12	1:C:82:ILE:H	1.85	0.41
1:C:131:LYS:O	1:C:132:GLU:CB	2.60	0.41
1:B:134:ARG:C	1:B:135:ILE:HG13	2.46	0.41
2:C:300[A]:PYV:H7	2:C:300[A]:PYV:N1	2.34	0.41
2:C:300[B]:PYV:H7	2:C:300[B]:PYV:N1	2.34	0.41
1:A:9:GLU:HA	1:A:129:PHE:HE2	1.86	0.41
1:C:52:PHE:O	1:C:79:ARG:NH1	2.50	0.41
1:C:87:LEU:HD23	1:C:87:LEU:HA	1.82	0.41
1:C:111:ILE:H	1:C:111:ILE:HG13	1.66	0.40
1:A:64:SER:OG	1:A:82:ILE:HD12	2.20	0.40
1:B:74:ARG:HG3	1:B:75:VAL:O	2.21	0.40
1:A:43:GLU:HG3	1:A:104:ARG:HH22	1.86	0.40
1:C:13:LEU:HD11	1:C:138:VAL:HG11	2.03	0.40
1:C:87:LEU:CD1	2:C:300[A]:PYV:BR	3.17	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)
7:A:206:HOH:O	7:B:203:HOH:O[8_555]	2.18	0.02

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	167/194 (86%)	147 (88%)	18 (11%)	2 (1%)	10	18
1	B	175/194 (90%)	165 (94%)	9 (5%)	1 (1%)	21	35
1	C	170/194 (88%)	158 (93%)	12 (7%)	0	100	100
All	All	512/582 (88%)	470 (92%)	39 (8%)	3 (1%)	21	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	ILE
1	B	84	ILE
1	A	67	VAL

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	156/174 (90%)	151 (97%)	5 (3%)	34	58
1	B	160/174 (92%)	158 (99%)	2 (1%)	61	80
1	C	159/174 (91%)	151 (95%)	8 (5%)	22	41
All	All	475/522 (91%)	460 (97%)	15 (3%)	34	58

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	14	LYS
1	A	62	ILE
1	A	155	ASP
1	A	169	GLN
1	B	57	THR
1	B	104	ARG
1	C	43	GLU

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Mol	Chain	Res	Type
1	C	74	ARG
1	C	80	ASP
1	C	92	SER
1	C	98	ILE
1	C	135	ILE
1	C	156	THR
1	C	157	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	90	ASN
1	C	60	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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
2	PYV	A	300[A]	-	24,24,24	5.38	10 (41%)	33,34,34	2.49	10 (30%)
5	P2M	B	300	-	23,23,23	4.24	5 (21%)	31,32,32	1.74	5 (16%)
2	PYV	C	300[B]	-	24,24,24	5.98	11 (45%)	33,34,34	3.24	16 (48%)
2	PYV	A	300[B]	-	24,23,24	5.38	10 (41%)	33,32,34	2.49	10 (30%)
3	SO4	C	192	-	4,4,4	0.41	0	6,6,6	0.51	0
2	PYV	C	300[A]	-	24,24,24	5.97	11 (45%)	33,34,34	3.24	16 (48%)
3	SO4	A	192	-	4,4,4	0.55	0	6,6,6	0.82	0

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
2	PYV	A	300[A]	-	-	0/12/12/12	0/3/3/3
5	P2M	B	300	-	-	0/12/12/12	0/3/3/3
2	PYV	C	300[B]	-	-	0/12/12/12	0/3/3/3
2	PYV	A	300[B]	-	-	0/12/12/12	0/3/3/3
2	PYV	C	300[A]	-	-	0/12/12/12	0/3/3/3

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	300[B]	PYV	C4-S	-26.47	1.50	1.77
2	C	300[A]	PYV	C4-S	-26.41	1.50	1.77
2	A	300[A]	PYV	C4-S	-24.24	1.52	1.77
2	A	300[B]	PYV	C4-S	-24.21	1.52	1.77
5	B	300	P2M	C4-S	-17.43	1.59	1.77
5	B	300	P2M	O1-S	-6.64	1.35	1.43
5	B	300	P2M	O2-S	-6.42	1.36	1.43
2	C	300[A]	PYV	C3-C2	-5.33	1.33	1.43
2	C	300[B]	PYV	C3-C2	-5.33	1.33	1.43
2	C	300[B]	PYV	O1-S	-5.27	1.37	1.43
2	C	300[A]	PYV	O1-S	-5.23	1.37	1.43
2	C	300[B]	PYV	O2-S	-5.04	1.37	1.43
2	C	300[A]	PYV	O2-S	-5.00	1.37	1.43
2	A	300[A]	PYV	C3-C2	-4.61	1.34	1.43
2	A	300[B]	PYV	C3-C2	-4.61	1.34	1.43
2	C	300[A]	PYV	C4-C3	-4.52	1.35	1.43
2	C	300[B]	PYV	C4-C3	-4.47	1.35	1.43
2	A	300[B]	PYV	O2-S	-3.83	1.39	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300[A]	PYV	O2-S	-3.76	1.39	1.43
2	A	300[B]	PYV	C4-C3	-3.68	1.36	1.43
2	C	300[A]	PYV	C1-C2	-3.65	1.34	1.42
2	C	300[B]	PYV	C1-C2	-3.65	1.34	1.42
2	A	300[A]	PYV	C4-C3	-3.65	1.36	1.43
2	A	300[A]	PYV	O1-S	-3.49	1.39	1.43
2	A	300[B]	PYV	O1-S	-3.49	1.39	1.43
2	A	300[A]	PYV	S-N1	-3.33	1.56	1.61
2	A	300[B]	PYV	S-N1	-3.33	1.56	1.61
2	C	300[A]	PYV	C7-C3	-2.89	1.36	1.42
2	C	300[B]	PYV	C7-C3	-2.89	1.36	1.42
2	C	300[A]	PYV	BR-C1	-2.83	1.82	1.90
2	C	300[B]	PYV	BR-C1	-2.83	1.82	1.90
2	C	300[A]	PYV	C10-C2	-2.62	1.37	1.42
2	C	300[B]	PYV	C10-C2	-2.62	1.37	1.42
2	A	300[A]	PYV	C1-C2	-2.55	1.36	1.42
2	A	300[B]	PYV	C1-C2	-2.55	1.36	1.42
2	C	300[B]	PYV	C11-N1	-2.55	1.43	1.47
2	C	300[A]	PYV	C11-N1	-2.52	1.43	1.47
2	C	300[A]	PYV	S-N1	-2.33	1.58	1.61
2	C	300[B]	PYV	S-N1	-2.33	1.58	1.61
2	A	300[B]	PYV	C10-C2	-2.31	1.37	1.42
2	A	300[A]	PYV	C10-C2	-2.31	1.37	1.42
2	A	300[A]	PYV	C11-N1	-2.29	1.44	1.47
2	A	300[B]	PYV	C11-N1	-2.29	1.44	1.47
2	A	300[A]	PYV	C7-C3	-2.20	1.37	1.42
2	A	300[B]	PYV	C7-C3	-2.16	1.38	1.42
5	B	300	P2M	C5-C4	-2.15	1.34	1.37
5	B	300	P2M	C3-C2	-2.09	1.39	1.43

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300[B]	PYV	O2-S-O1	-9.08	108.50	119.52
2	C	300[A]	PYV	O2-S-O1	-9.05	108.52	119.52
2	A	300[A]	PYV	C3-C4-S	-7.54	112.78	121.66
2	A	300[B]	PYV	C3-C4-S	-7.53	112.80	121.66
2	C	300[B]	PYV	C3-C4-S	-6.66	113.82	121.66
2	C	300[A]	PYV	C3-C4-S	-6.65	113.84	121.66
2	C	300[A]	PYV	O1-S-N1	5.80	116.06	107.03
2	C	300[B]	PYV	O1-S-N1	5.79	116.04	107.03
2	C	300[A]	PYV	O2-S-C4	-5.52	98.46	108.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	300[B]	PYV	O2-S-C4	-5.52	98.47	108.09
2	C	300[A]	PYV	BR-C1-C2	-4.98	110.23	119.75
2	C	300[B]	PYV	BR-C1-C2	-4.98	110.23	119.75
5	B	300	P2M	C5-C4-S	4.76	122.46	116.92
2	A	300[B]	PYV	C6-C5-C4	-4.56	116.10	121.66
2	A	300[A]	PYV	C6-C5-C4	-4.55	116.12	121.66
5	B	300	P2M	C3-C4-S	-4.49	116.37	121.66
2	C	300[A]	PYV	C4-S-N1	4.48	115.06	106.78
2	C	300[B]	PYV	C4-S-N1	4.48	115.06	106.78
2	C	300[A]	PYV	C5-C4-C3	4.41	125.88	120.95
2	C	300[A]	PYV	C6-C5-C4	-4.37	116.33	121.66
2	C	300[B]	PYV	C5-C4-C3	4.37	125.83	120.95
2	C	300[B]	PYV	C6-C5-C4	-4.32	116.39	121.66
2	A	300[A]	PYV	C5-C4-S	4.23	121.85	116.92
2	A	300[B]	PYV	C5-C4-S	4.20	121.81	116.92
2	A	300[B]	PYV	O2-S-O1	-4.14	114.49	119.52
2	A	300[A]	PYV	O2-S-O1	-4.13	114.50	119.52
5	B	300	P2M	C11-N1-S	4.01	128.42	120.26
2	A	300[B]	PYV	C5-C4-C3	3.96	125.37	120.95
2	A	300[A]	PYV	C5-C4-C3	3.94	125.35	120.95
2	A	300[A]	PYV	O2-S-N1	3.80	112.95	107.03
2	A	300[B]	PYV	O2-S-N1	3.80	112.95	107.03
2	A	300[A]	PYV	BR-C1-C2	-3.60	112.88	119.75
2	A	300[B]	PYV	BR-C1-C2	-3.60	112.88	119.75
2	C	300[A]	PYV	C10-C2-C3	3.30	122.52	118.42
2	C	300[B]	PYV	C10-C2-C3	3.30	122.52	118.42
2	C	300[B]	PYV	C5-C4-S	2.81	120.19	116.92
2	C	300[A]	PYV	C5-C4-S	2.76	120.14	116.92
2	C	300[A]	PYV	C6-C1-C2	2.75	126.04	121.87
2	C	300[B]	PYV	C6-C1-C2	2.74	126.03	121.87
2	C	300[B]	PYV	C9-C10-C2	-2.45	117.61	120.91
2	C	300[A]	PYV	C9-C10-C2	-2.42	117.64	120.91
2	C	300[B]	PYV	C5-C6-C1	-2.41	116.41	120.69
2	C	300[A]	PYV	C5-C6-C1	-2.40	116.42	120.69
2	C	300[B]	PYV	BR-C1-C6	2.36	123.72	116.88
2	C	300[A]	PYV	BR-C1-C6	2.36	123.70	116.88
2	C	300[B]	PYV	O2-S-N1	2.31	110.62	107.03
2	C	300[A]	PYV	C10-C2-C1	-2.29	119.74	124.93
2	C	300[B]	PYV	C10-C2-C1	-2.29	119.74	124.93
2	C	300[A]	PYV	O2-S-N1	2.29	110.59	107.03
2	A	300[B]	PYV	C14-C13-N2	-2.20	119.94	123.42
2	A	300[A]	PYV	C14-C13-N2	-2.19	119.95	123.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	300[B]	PYV	C13-N2-C12	2.15	120.43	117.41
2	A	300[A]	PYV	C13-N2-C12	2.14	120.41	117.41
2	A	300[A]	PYV	C10-C2-C3	2.10	121.03	118.42
2	A	300[B]	PYV	C10-C2-C3	2.07	120.99	118.42
5	B	300	P2M	C6-C1-C2	2.04	123.32	120.48
5	B	300	P2M	C7-C3-C4	-2.02	120.54	123.58

There are no chirality outliers.

There are no torsion outliers.

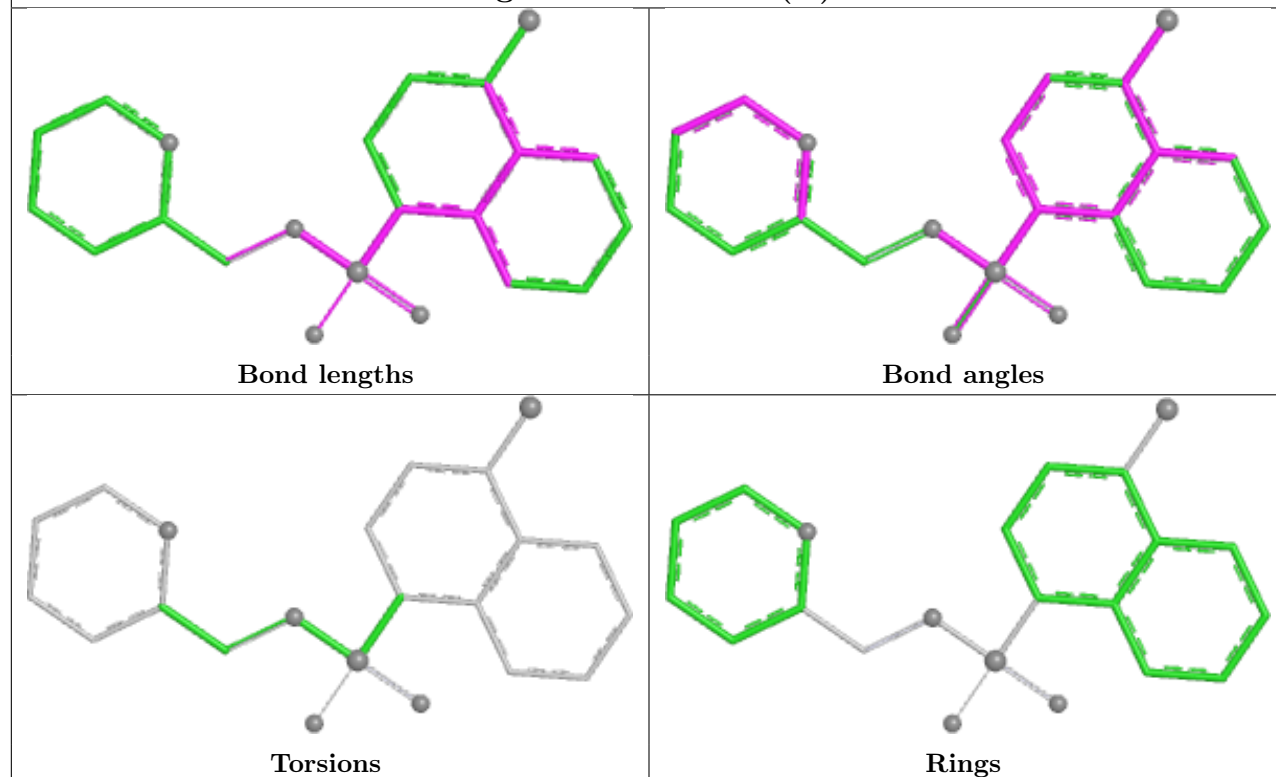
There are no ring outliers.

6 monomers are involved in 17 short contacts:

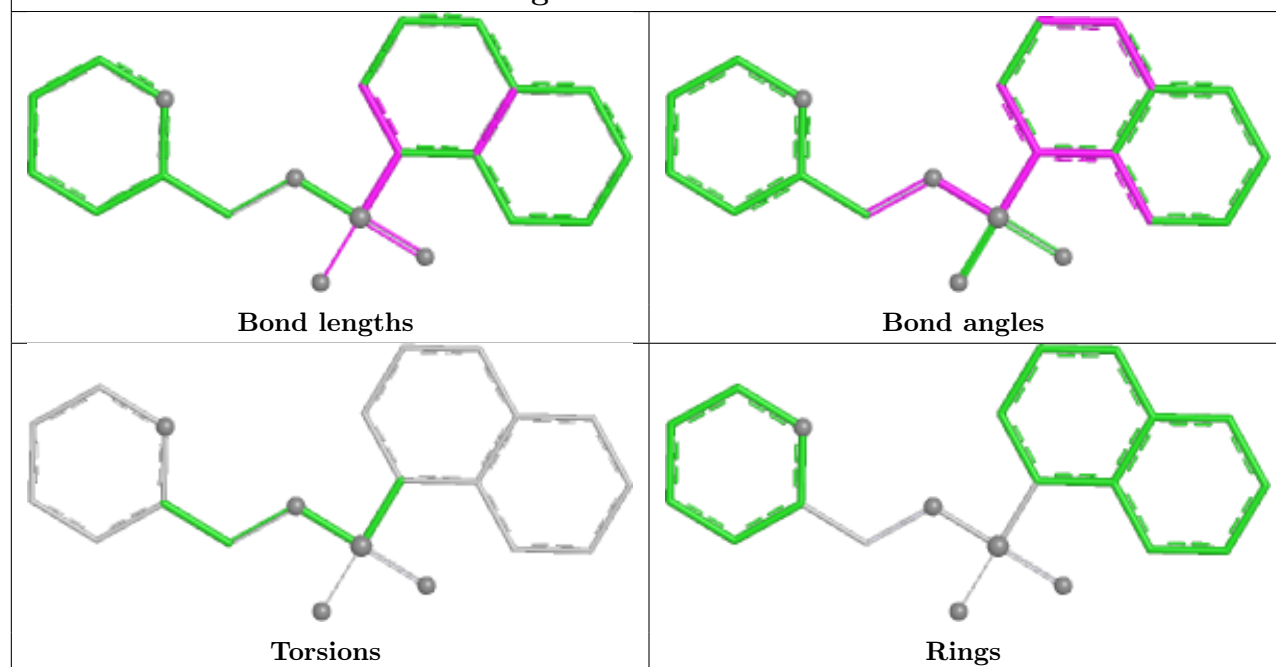
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300[A]	PYV	3	0
5	B	300	P2M	2	0
2	C	300[B]	PYV	3	0
2	A	300[B]	PYV	2	0
2	C	300[A]	PYV	6	0
3	A	192	SO4	1	0

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.

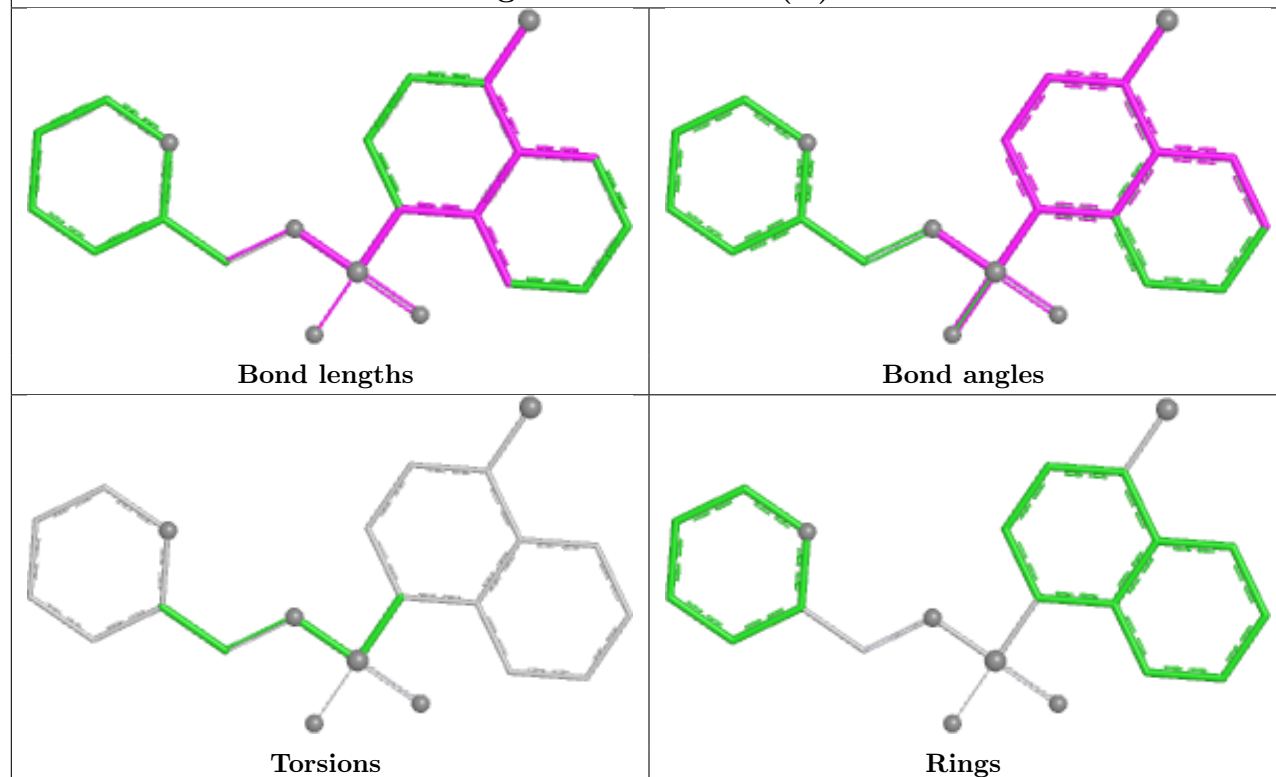
Ligand PYV A 300 (A)



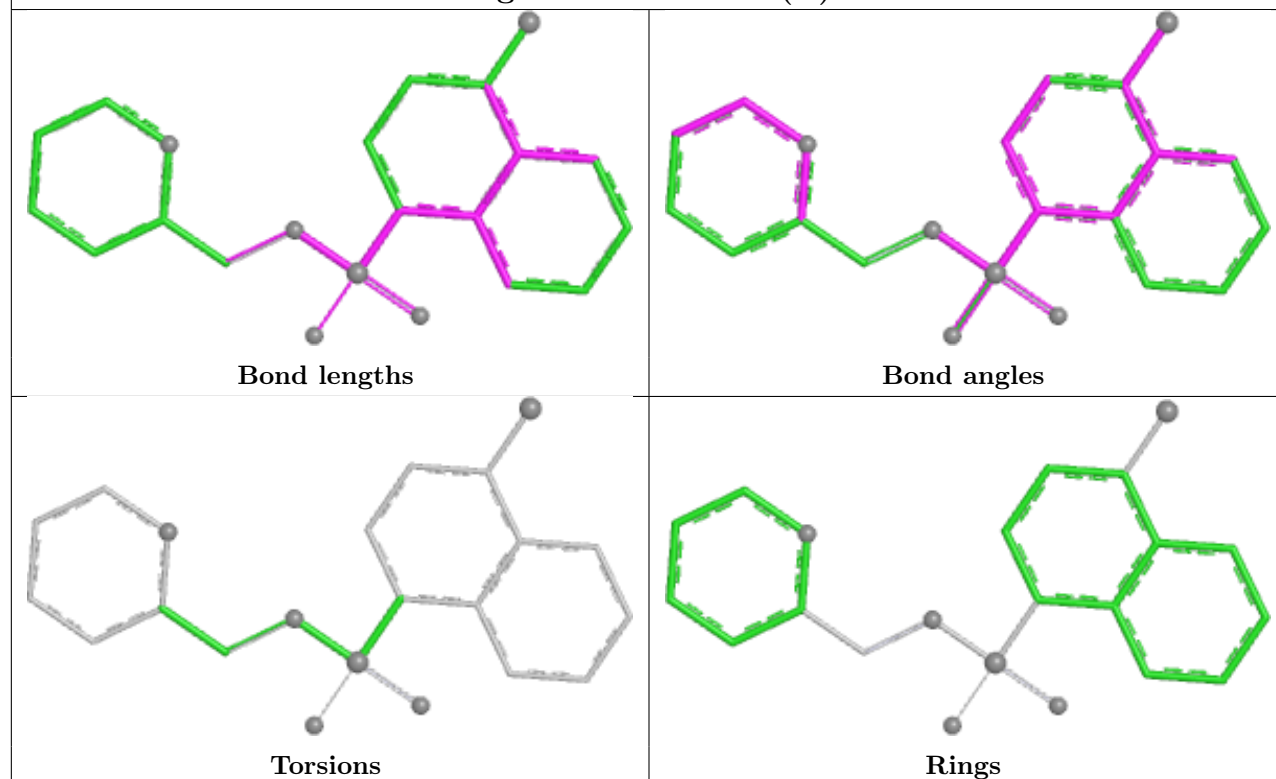
Ligand P2M B 300

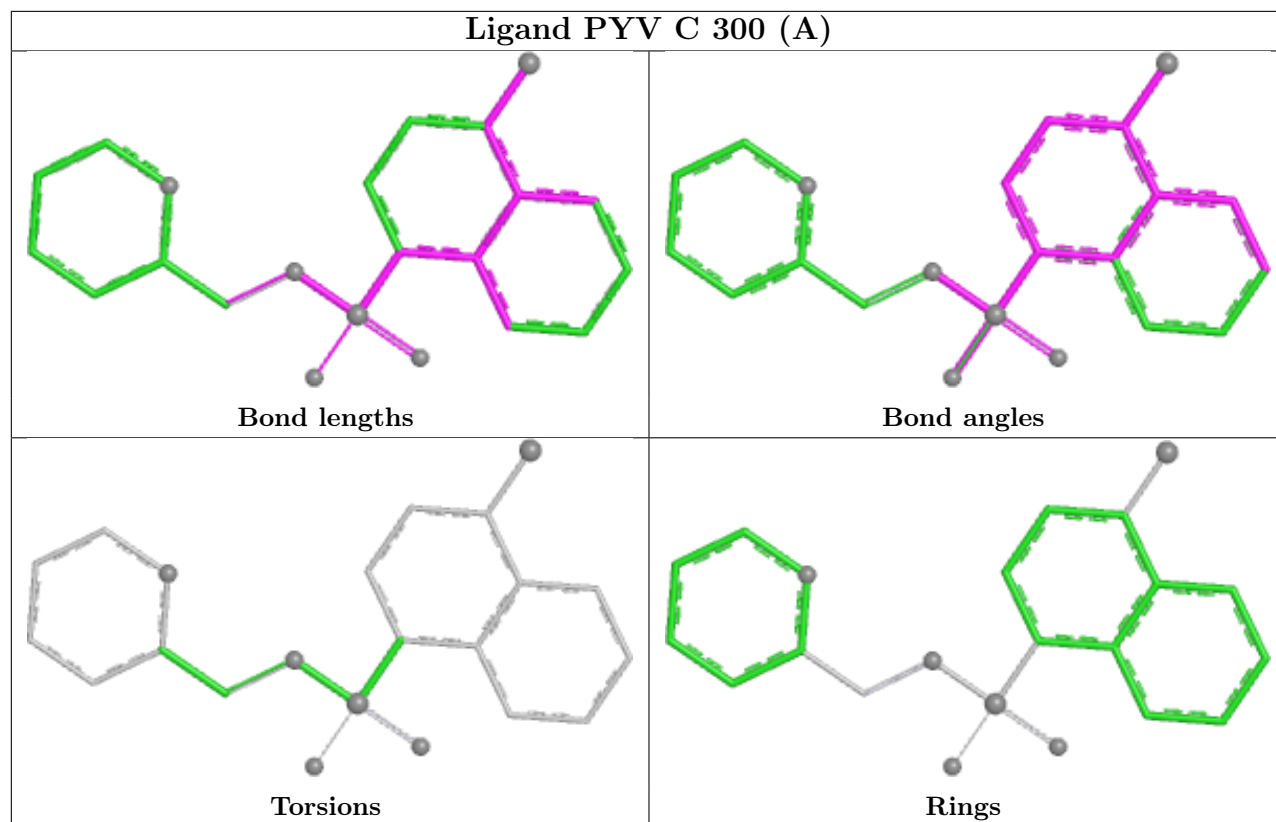


Ligand PYV C 300 (B)



Ligand PYV A 300 (B)





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	169/194 (87%)	0.16	5 (2%)	52	49	21, 46, 101, 148	2 (1%)
1	B	177/194 (91%)	0.23	5 (2%)	55	51	28, 50, 96, 153	0
1	C	171/194 (88%)	0.14	9 (5%)	32	28	14, 43, 104, 137	3 (1%)
All	All	517/582 (88%)	0.18	19 (3%)	45	41	14, 47, 101, 153	5 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	3	SER	3.6
1	A	6	THR	3.4
1	C	103	ARG	3.1
1	A	88	SER	3.0
1	C	89	ALA	2.8
1	C	83	VAL	2.6
1	A	133	ASN	2.6
1	B	154	ASP	2.5
1	B	158	MET	2.5
1	A	155	ASP	2.5
1	B	89	ALA	2.5
1	C	88	SER	2.3
1	C	115	HIS	2.3
1	C	87	LEU	2.3
1	C	156	THR	2.2
1	B	156	THR	2.1
1	B	88	SER	2.1
1	A	9	GLU	2.1
1	C	157	ARG	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](#)

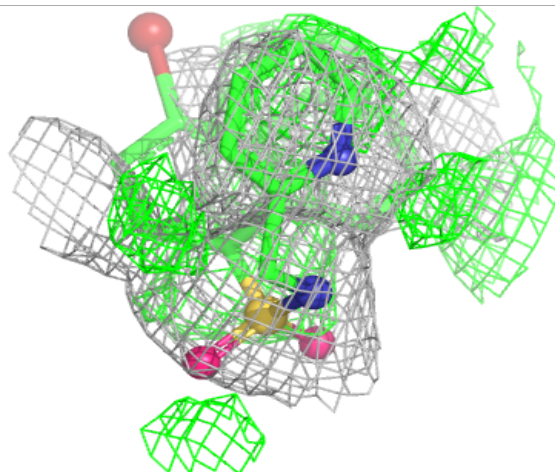
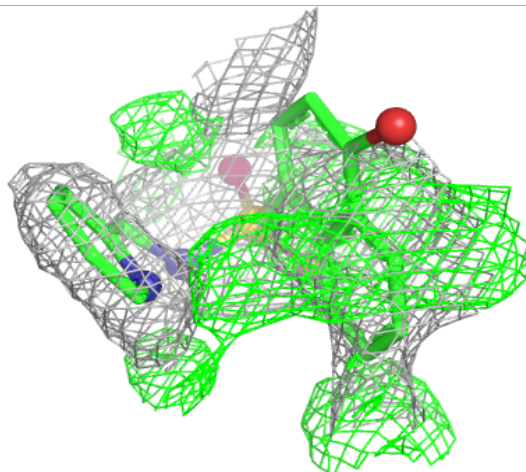
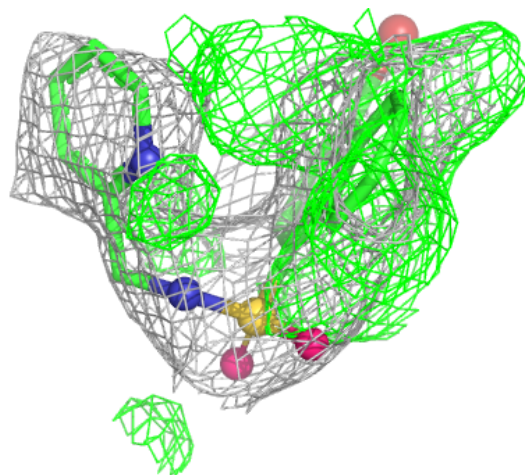
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	C	192	5/5	0.66	0.20	69,83,90,92	0
6	CL	B	192	1/1	0.76	0.18	95,95,95,95	0
2	PYV	A	300[A]	22/22	0.84	0.27	28,55,56,73	22
2	PYV	A	300[B]	21/22	0.84	0.27	28,54,56,57	21
5	P2M	B	300	21/21	0.86	0.18	45,53,64,65	21
4	NA	A	193	1/1	0.89	0.13	60,60,60,60	0
2	PYV	C	300[B]	22/22	0.89	0.14	42,65,69,112	22
2	PYV	C	300[A]	22/22	0.89	0.14	42,65,69,73	22
3	SO4	A	192	5/5	0.91	0.09	52,57,66,73	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.

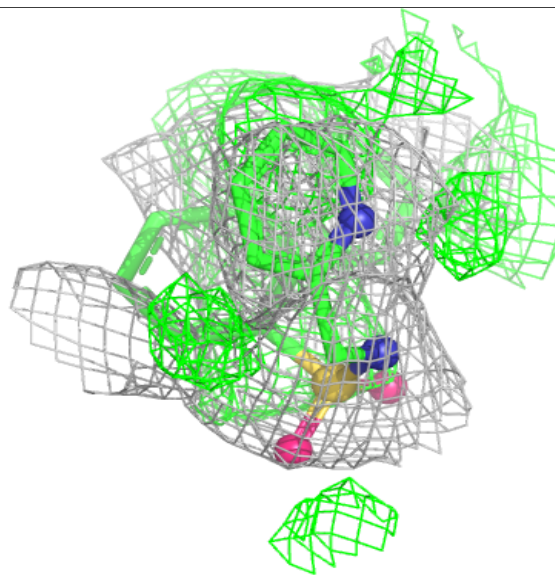
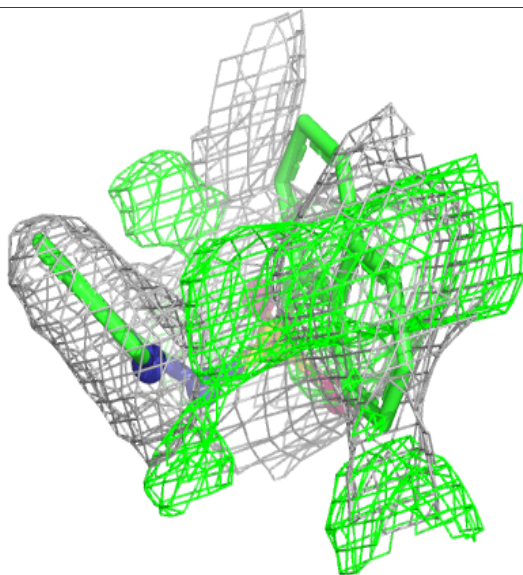
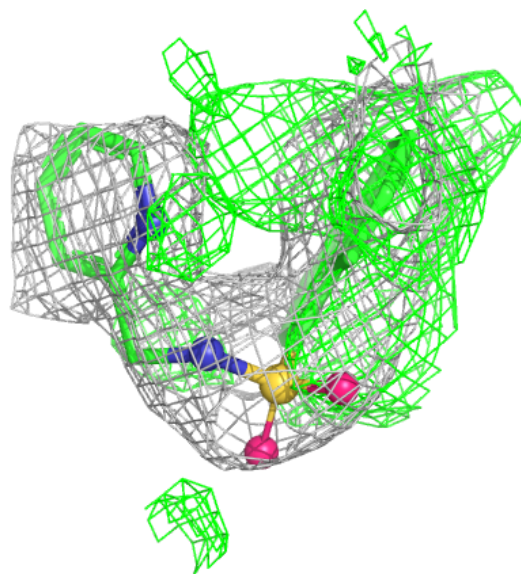
Electron density around PYV A 300 (A):

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



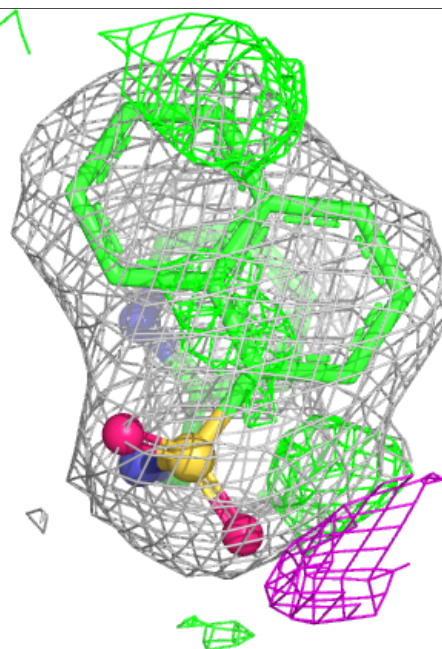
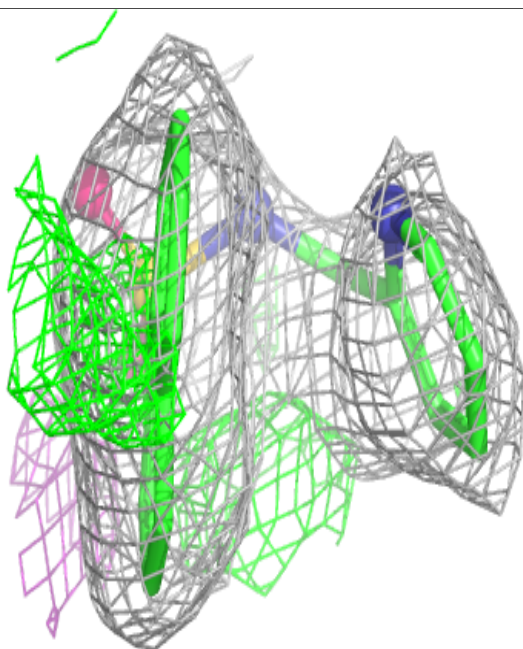
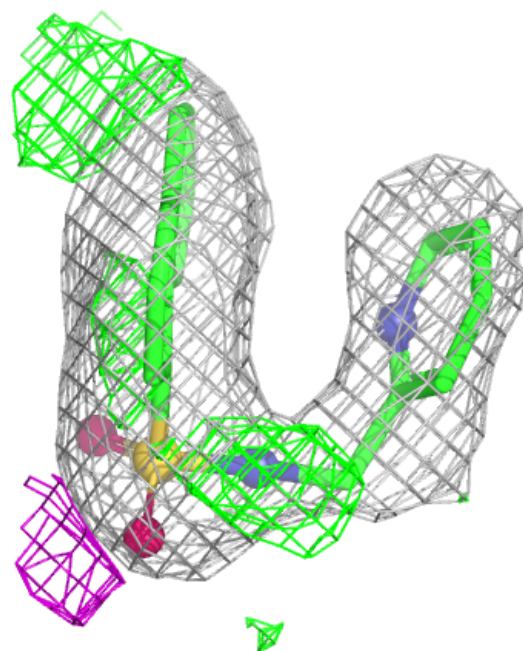
Electron density around PYV A 300 (B):

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



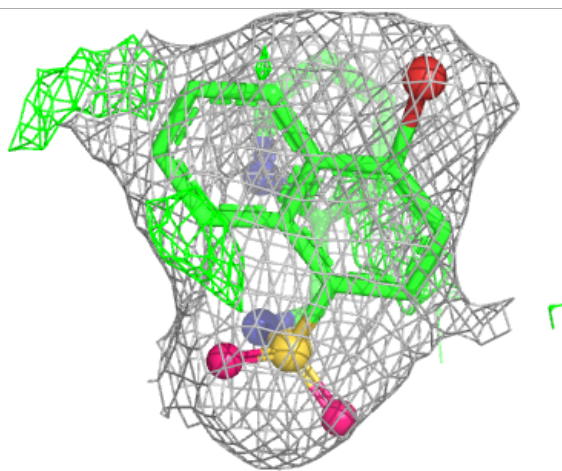
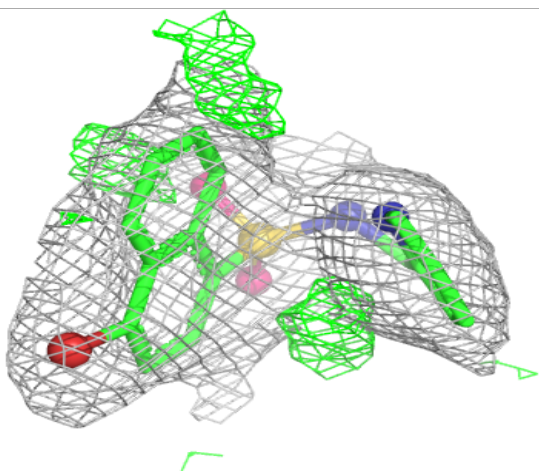
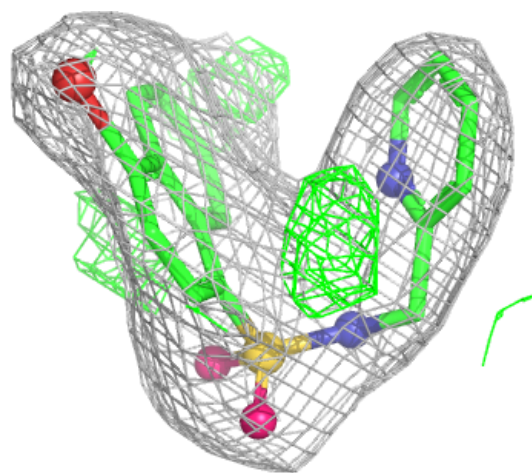
Electron density around P2M B 300:

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



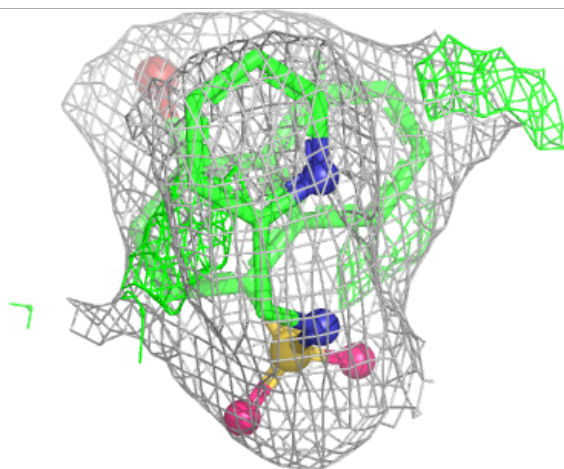
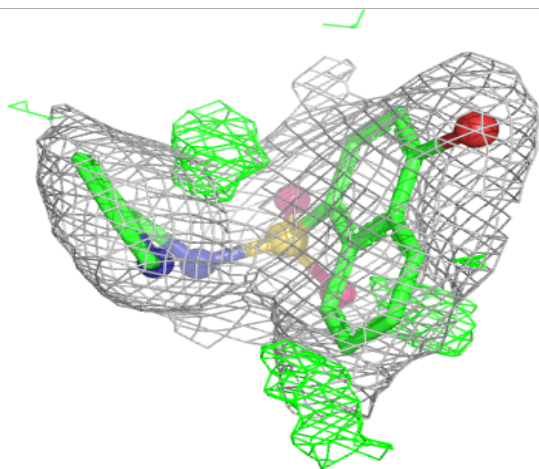
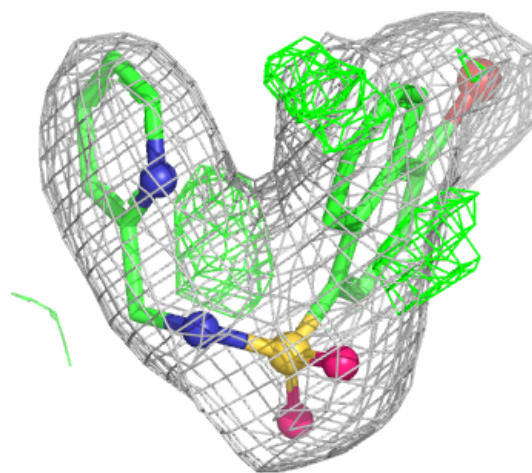
Electron density around PYV C 300 (B):

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



Electron density around PYV C 300 (A):

$2mF_o-DF_c$ (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.