WORLDWIDE PDB

PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Dec 17, 2023 – 02:30 PM EST

PDB ID	:	4TJW
Title	:	Crystal Structure of human Tankyrase 2 in complex with PJ-34.
Authors	:	Qiu, W.; Lam, R.; Romanov, V.; Gordon, R.; Gebremeskel, S.; Vodsedalek, J.; Thompson, C.; Beletskaya, I.; Battaile, K.P.; Pai, E.F.; Chirgadze, N.Y.
Deposited on	:	2014-05-25
Resolution	:	1.70 Å(reported)

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

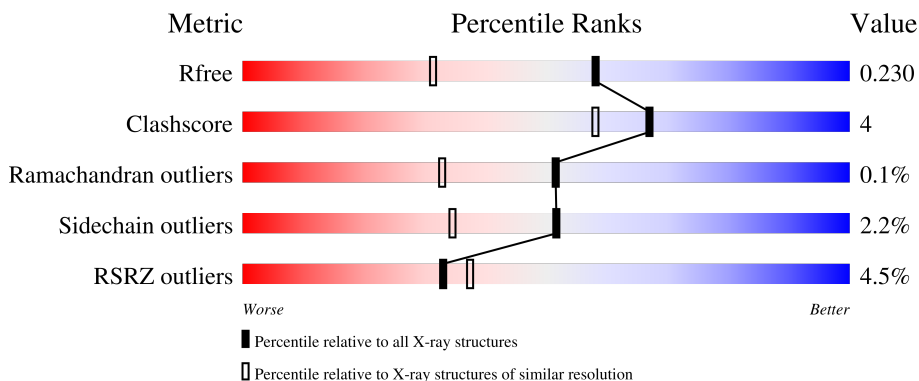|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.70 Å.

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

Metric	Percentile Ranks	Value
Rfree		0.230
Clashscore		4
Ramachandran outliers		0.1%
Sidechain outliers		2.2%
RSRZ outliers		4.5%

Worse .... *Better*

▮ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 <=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	227	3% / 78% 10% 12%
1	B	227	3% / 82% 8% · 9%
1	C	227	3% / 84% 6% · 8%
1	D	227	7% / 73% 11% · 16%

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 7349 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 Tankyrase-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	199	Total	C	N	O	S	0	3	0
			1617	1020	299	286	12			
1	B	207	Total	C	N	O	S	0	3	0
			1684	1064	311	297	12			
1	C	208	Total	C	N	O	S	0	0	0
			1674	1055	307	301	11			
1	D	191	Total	C	N	O	S	0	0	0
			1542	974	287	270	11			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	938	MET	-	initiating methionine	UNP Q9H2K2
A	939	GLY	-	expression tag	UNP Q9H2K2
A	940	SER	-	expression tag	UNP Q9H2K2
A	941	SER	-	expression tag	UNP Q9H2K2
A	942	HIS	-	expression tag	UNP Q9H2K2
A	943	HIS	-	expression tag	UNP Q9H2K2
A	944	HIS	-	expression tag	UNP Q9H2K2
A	945	HIS	-	expression tag	UNP Q9H2K2
A	946	HIS	-	expression tag	UNP Q9H2K2
A	947	HIS	-	expression tag	UNP Q9H2K2
A	948	SER	-	expression tag	UNP Q9H2K2
A	949	SER	-	expression tag	UNP Q9H2K2
A	950	GLY	-	expression tag	UNP Q9H2K2
A	951	ARG	-	expression tag	UNP Q9H2K2
A	952	GLU	-	expression tag	UNP Q9H2K2
A	953	ASN	-	expression tag	UNP Q9H2K2
A	954	LEU	-	expression tag	UNP Q9H2K2
A	955	TYR	-	expression tag	UNP Q9H2K2
A	956	PHE	-	expression tag	UNP Q9H2K2
A	957	GLN	-	expression tag	UNP Q9H2K2
A	958	GLY	-	expression tag	UNP Q9H2K2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	938	MET	-	initiating methionine	UNP Q9H2K2
B	939	GLY	-	expression tag	UNP Q9H2K2
B	940	SER	-	expression tag	UNP Q9H2K2
B	941	SER	-	expression tag	UNP Q9H2K2
B	942	HIS	-	expression tag	UNP Q9H2K2
B	943	HIS	-	expression tag	UNP Q9H2K2
B	944	HIS	-	expression tag	UNP Q9H2K2
B	945	HIS	-	expression tag	UNP Q9H2K2
B	946	HIS	-	expression tag	UNP Q9H2K2
B	947	HIS	-	expression tag	UNP Q9H2K2
B	948	SER	-	expression tag	UNP Q9H2K2
B	949	SER	-	expression tag	UNP Q9H2K2
B	950	GLY	-	expression tag	UNP Q9H2K2
B	951	ARG	-	expression tag	UNP Q9H2K2
B	952	GLU	-	expression tag	UNP Q9H2K2
B	953	ASN	-	expression tag	UNP Q9H2K2
B	954	LEU	-	expression tag	UNP Q9H2K2
B	955	TYR	-	expression tag	UNP Q9H2K2
B	956	PHE	-	expression tag	UNP Q9H2K2
B	957	GLN	-	expression tag	UNP Q9H2K2
B	958	GLY	-	expression tag	UNP Q9H2K2
C	938	MET	-	initiating methionine	UNP Q9H2K2
C	939	GLY	-	expression tag	UNP Q9H2K2
C	940	SER	-	expression tag	UNP Q9H2K2
C	941	SER	-	expression tag	UNP Q9H2K2
C	942	HIS	-	expression tag	UNP Q9H2K2
C	943	HIS	-	expression tag	UNP Q9H2K2
C	944	HIS	-	expression tag	UNP Q9H2K2
C	945	HIS	-	expression tag	UNP Q9H2K2
C	946	HIS	-	expression tag	UNP Q9H2K2
C	947	HIS	-	expression tag	UNP Q9H2K2
C	948	SER	-	expression tag	UNP Q9H2K2
C	949	SER	-	expression tag	UNP Q9H2K2
C	950	GLY	-	expression tag	UNP Q9H2K2
C	951	ARG	-	expression tag	UNP Q9H2K2
C	952	GLU	-	expression tag	UNP Q9H2K2
C	953	ASN	-	expression tag	UNP Q9H2K2
C	954	LEU	-	expression tag	UNP Q9H2K2
C	955	TYR	-	expression tag	UNP Q9H2K2
C	956	PHE	-	expression tag	UNP Q9H2K2
C	957	GLN	-	expression tag	UNP Q9H2K2
C	958	GLY	-	expression tag	UNP Q9H2K2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	938	MET	-	initiating methionine	UNP Q9H2K2
D	939	GLY	-	expression tag	UNP Q9H2K2
D	940	SER	-	expression tag	UNP Q9H2K2
D	941	SER	-	expression tag	UNP Q9H2K2
D	942	HIS	-	expression tag	UNP Q9H2K2
D	943	HIS	-	expression tag	UNP Q9H2K2
D	944	HIS	-	expression tag	UNP Q9H2K2
D	945	HIS	-	expression tag	UNP Q9H2K2
D	946	HIS	-	expression tag	UNP Q9H2K2
D	947	HIS	-	expression tag	UNP Q9H2K2
D	948	SER	-	expression tag	UNP Q9H2K2
D	949	SER	-	expression tag	UNP Q9H2K2
D	950	GLY	-	expression tag	UNP Q9H2K2
D	951	ARG	-	expression tag	UNP Q9H2K2
D	952	GLU	-	expression tag	UNP Q9H2K2
D	953	ASN	-	expression tag	UNP Q9H2K2
D	954	LEU	-	expression tag	UNP Q9H2K2
D	955	TYR	-	expression tag	UNP Q9H2K2
D	956	PHE	-	expression tag	UNP Q9H2K2
D	957	GLN	-	expression tag	UNP Q9H2K2
D	958	GLY	-	expression tag	UNP Q9H2K2

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0
2	D	1	Total 1	Zn 1	0	0

- Molecule 3 is N 2 ,N 2 -DIMETHYL-N 1 -(6-OXO-5,6-DIHYDROPHENANTHRIDIN-2-Y L)GLYCINAMIDE (three-letter code: P34) (formula: $C_{17}H_{17}N_3O_2$).

P34

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			22	17	3	2		
3	B	1	Total	C	N	O	0	0
			22	17	3	2		
3	C	1	Total	C	N	O	0	0
			22	17	3	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	201	Total	O	0	0
			201	201		
4	B	191	Total	O	0	0
			191	191		
4	C	212	Total	O	0	0
			212	212		
4	D	158	Total	O	0	0
			158	158		

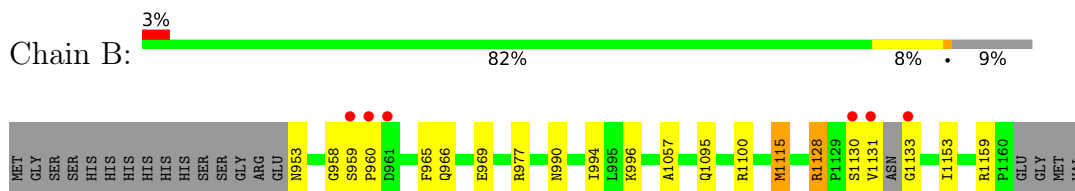# 3 Residue-property plots ⓘ

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.
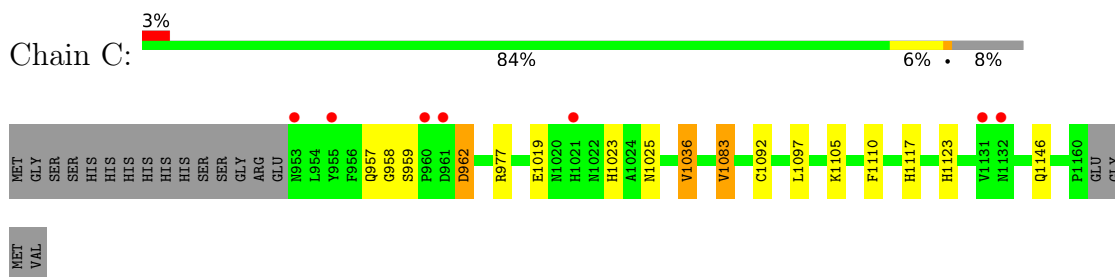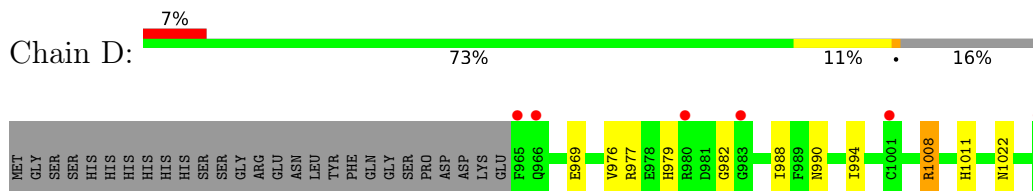
- Molecule 1: Tankyrase-2
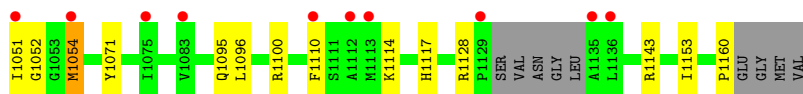
Chain A:

- Molecule 1: Tankyrase-2

Chain B:

- Molecule 1: Tankyrase-2

Chain C:

- Molecule 1: Tankyrase-2

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	73.76Å 79.31Å 153.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.81 – 1.70 19.81 – 1.70	Depositor EDS
% Data completeness (in resolution range)	83.2 (19.81-1.70) 83.2 (19.81-1.70)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.89 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.177 , 0.227 0.184 , 0.230	Depositor DCC
R_{free} test set	1003 reflections (1.21%)	wwPDB-VP
Wilson B-factor (Å^2)	19.5	Xtriage
Anisotropy	0.521	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 49.2	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7349	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.40% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN, P34

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	0.98	0/1669	0.98	3/2244 (0.1%)
1	B	0.93	0/1738	1.00	8/2336 (0.3%)
1	C	0.96	0/1720	0.98	1/2315 (0.0%)
1	D	0.94	0/1584	1.01	7/2128 (0.3%)
All	All	0.95	0/6711	0.99	19/9023 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	1100	ARG	NE-CZ-NH2	-9.46	115.57	120.30
1	B	1128	ARG	NE-CZ-NH2	8.40	124.50	120.30
1	D	1008	ARG	NE-CZ-NH1	6.89	123.75	120.30
1	D	1128	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	B	1128	ARG	NE-CZ-NH1	-6.20	117.20	120.30
1	C	977	ARG	NE-CZ-NH1	6.19	123.40	120.30
1	D	1047	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	B	977	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	D	1008	ARG	NE-CZ-NH2	-5.79	117.41	120.30
1	B	1100	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	A	1100	ARG	CG-CD-NE	-5.67	99.89	111.80
1	D	1143	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	D	1047	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	D	1100	ARG	CB-CA-C	-5.36	99.69	110.40
1	A	1037[A]	ASN	CB-CA-C	5.34	121.09	110.40
1	A	1037[B]	ASN	CB-CA-C	5.34	121.09	110.40
1	B	977	ARG	NE-CZ-NH2	-5.19	117.71	120.30
1	B	1115[A]	MET	CG-SD-CE	5.01	108.21	100.20
1	B	1115[B]	MET	CG-SD-CE	5.01	108.21	100.20

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	1617	0	1557	11	0
1	B	1684	0	1618	12	0
1	C	1674	0	1592	15	0
1	D	1542	0	1479	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	22	0	17	0	0
3	B	22	0	17	0	0
3	C	22	0	17	0	0
4	A	201	0	0	0	1
4	B	191	0	0	3	1
4	C	212	0	0	2	0
4	D	158	0	0	6	0
All	All	7349	0	6297	50	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 4.

All (50) 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:C:1023:HIS:HD2	1:C:1025:ASN:H	1.13	0.88
1:C:1023:HIS:CD2	1:C:1025:ASN:H	2.01	0.77
1:C:1105:LYS:H	1:C:1123:HIS:HD2	1.34	0.75
1:B:1131:VAL:O	1:B:1133:GLY:N	2.28	0.67
1:D:1054:MET:HE2	4:D:1347:HOH:O	1.97	0.65
1:D:1054:MET:N	1:D:1054:MET:SD	2.70	0.61
1:B:1057:ALA:O	1:D:1117:HIS:HE1	1.84	0.60
1:C:1110:PHE:HD2	1:D:1110:PHE:CE2	2.20	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1117:HIS:HD2	4:D:1401:HOH:O	1.85	0.60
1:B:1131:VAL:HG23	4:B:1404:HOH:O	2.02	0.59
1:C:1019:GLU:OE2	1:C:1123:HIS:HE1	1.85	0.59
1:D:1008:ARG:NH2	4:D:1409:HOH:O	2.36	0.59
1:C:1023:HIS:HD2	1:C:1025:ASN:N	1.95	0.58
1:A:1057:ALA:O	1:C:1117:HIS:HE1	1.86	0.58
1:D:1071:TYR:HD2	4:D:1324:HOH:O	1.86	0.57
1:D:979:HIS:O	1:D:982:GLY:N	2.35	0.56
1:C:959:SER:O	1:C:962:ASP:HB2	2.06	0.55
1:C:1083:VAL:HG13	1:C:1092:CYS:SG	2.47	0.55
1:B:1131:VAL:N	4:B:1404:HOH:O	2.40	0.54
1:C:1117:HIS:HD2	4:C:1423:HOH:O	1.89	0.53
1:C:1023:HIS:HE1	1:D:1022:ASN:O	1.93	0.52
1:A:1159[A]:ARG:HB3	1:A:1159[A]:ARG:CZ	2.40	0.51
1:C:1110:PHE:CD2	1:D:1110:PHE:CE2	2.98	0.50
1:A:969:GLU:HG3	1:A:994:ILE:HD12	1.93	0.49
1:B:960:PRO:O	1:B:966:GLN:NE2	2.39	0.49
1:A:1115[A]:MET:HE3	1:A:1115[A]:MET:HB2	1.61	0.47
1:B:1115[B]:MET:HB2	1:B:1115[B]:MET:HE3	1.22	0.47
1:D:976:VAL:O	1:D:977:ARG:NH1	2.48	0.46
1:A:1044:PHE:HB3	1:A:1059:ILE:HD13	1.97	0.46
1:A:975:THR:HB	1:A:1064:ASN:HA	1.98	0.46
1:C:958:GLY:HA2	4:C:1309:HOH:O	2.14	0.46
1:C:1036:VAL:HG22	1:C:1097:LEU:HG	1.98	0.46
1:D:1051:ILE:HG23	1:D:1114:LYS:HB3	1.98	0.45
1:B:1159[A]:ARG:NH2	4:B:1302:HOH:O	2.50	0.45
1:D:1011:HIS:HE1	4:D:1426:HOH:O	2.00	0.45
1:A:1109:GLN:HG3	1:A:1113:MET:CE	2.48	0.44
1:D:1160:PRO:C	4:D:1362:HOH:O	2.55	0.43
1:B:1095:GLN:HA	1:B:1153:ILE:O	2.18	0.43
1:B:1130:SER:O	1:B:1133:GLY:N	2.52	0.43
1:A:1113:MET:CE	1:A:1115[A]:MET:CE	2.97	0.43
1:A:1053:GLY:O	1:A:1114:LYS:HE2	2.19	0.42
1:A:1003:LYS:O	1:A:1007:GLU:HG3	2.19	0.42
1:A:963:LYS:HA	1:A:966:GLN:HB2	2.02	0.42
1:D:1096:LEU:HD23	1:D:1096:LEU:HA	1.83	0.41
1:B:953:ASN:HD22	1:B:953:ASN:HA	1.68	0.41
1:C:1023:HIS:CE1	1:D:1022:ASN:O	2.73	0.41
1:B:969:GLU:HG3	1:B:994:ILE:CD1	2.51	0.41
1:B:958:GLY:HA3	1:B:965:PHE:CD1	2.55	0.41
1:D:1095:GLN:HA	1:D:1153:ILE:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:969:GLU:HG3	1:D:994:ILE:HD12	2.02	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)
4:A:1314:HOH:O	4:B:1324:HOH:O[1_565]	2.19	0.01

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	200/227 (88%)	198 (99%)	2 (1%)	0	100	100
1	B	206/227 (91%)	201 (98%)	5 (2%)	0	100	100
1	C	206/227 (91%)	203 (98%)	3 (2%)	0	100	100
1	D	187/227 (82%)	179 (96%)	7 (4%)	1 (0%)	29	13
All	All	799/908 (88%)	781 (98%)	17 (2%)	1 (0%)	51	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	1052	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	169/192 (88%)	166 (98%)	3 (2%)	59	43
1	B	176/192 (92%)	172 (98%)	4 (2%)	50	33
1	C	175/192 (91%)	170 (97%)	5 (3%)	42	23
1	D	160/192 (83%)	157 (98%)	3 (2%)	57	41
All	All	680/768 (88%)	665 (98%)	15 (2%)	52	34

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

Mol	Chain	Res	Type
1	A	1018	GLU
1	A	1036	VAL
1	A	1161	GLU
1	B	959	SER
1	B	990	ASN
1	B	996	LYS
1	B	1128	ARG
1	C	957	GLN
1	C	962	ASP
1	C	1036	VAL
1	C	1083	VAL
1	C	1146	GLN
1	D	988	ILE
1	D	990	ASN
1	D	1054	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1095	GLN
1	B	953	ASN
1	B	984	HIS
1	B	1037	ASN
1	B	1048	HIS
1	B	1095	GLN
1	C	990	ASN
1	C	1023	HIS
1	C	1117	HIS
1	C	1123	HIS
1	D	990	ASN

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Mol	Chain	Res	Type
1	D	1011	HIS
1	D	1023	HIS
1	D	1117	HIS
1	D	1156	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\#|Z| > 2$
3	P34	B	1202	-	23,24,24	2.17	8 (34%)	30,34,34	2.39	6 (20%)
3	P34	C	1202	-	23,24,24	2.16	8 (34%)	30,34,34	2.42	6 (20%)
3	P34	A	1202	-	23,24,24	2.83	9 (39%)	30,34,34	3.03	11 (36%)

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.

W O R L D W I D E
wwPDB
PROTEIN DATA BANK

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	P34	B	1202	-	-	2/8/8/8	0/3/3/3
3	P34	C	1202	-	-	2/8/8/8	0/3/3/3
3	P34	A	1202	-	-	5/8/8/8	0/3/3/3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1202	P34	CAR-NAN	7.05	1.47	1.35
3	C	1202	P34	CA-C	5.29	1.60	1.52
3	A	1202	P34	CAE-CAF	5.02	1.51	1.38
3	A	1202	P34	CA-N	4.89	1.59	1.47
3	A	1202	P34	CAG-CAP	4.88	1.47	1.39
3	A	1202	P34	CAA-N	4.68	1.60	1.46
3	C	1202	P34	CA-N	4.47	1.58	1.47
3	C	1202	P34	CAE-CAH	4.13	1.46	1.36
3	B	1202	P34	CAQ-NAN	3.93	1.39	1.33
3	B	1202	P34	CAG-CAP	3.79	1.45	1.39
3	B	1202	P34	CAB-N	3.71	1.57	1.46
3	A	1202	P34	CAP-NAM	3.59	1.48	1.41
3	B	1202	P34	CAJ-CAG	3.47	1.43	1.36
3	B	1202	P34	CAR-NAN	3.14	1.40	1.35
3	C	1202	P34	CAB-N	3.10	1.55	1.46
3	B	1202	P34	CA-N	3.00	1.54	1.47
3	B	1202	P34	CAI-CAT	2.68	1.46	1.41
3	B	1202	P34	CAE-CAF	2.63	1.45	1.38
3	C	1202	P34	CAA-N	2.31	1.53	1.46
3	C	1202	P34	CAU-CAR	-2.28	1.37	1.41
3	A	1202	P34	CAJ-CAG	2.07	1.41	1.36
3	A	1202	P34	CA-C	2.04	1.55	1.52
3	A	1202	P34	CAI-CAT	2.04	1.45	1.41
3	C	1202	P34	CAF-CAI	2.04	1.41	1.36
3	C	1202	P34	CAR-NAN	2.02	1.38	1.35

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1202	P34	CAS-CAQ-NAN	-10.81	116.87	124.40
3	C	1202	P34	CAS-CAQ-NAN	-9.18	118.00	124.40
3	B	1202	P34	CAS-CAQ-NAN	-8.24	118.65	124.40
3	B	1202	P34	CAU-CAR-NAN	-6.81	119.90	123.60
3	A	1202	P34	CAR-NAN-CAQ	5.11	123.84	116.83
3	A	1202	P34	CA-C-NAM	5.09	123.69	114.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	1202	P34	CAR-NAN-CAQ	4.65	123.21	116.83
3	A	1202	P34	CAU-CAR-NAN	-4.52	121.14	123.60
3	B	1202	P34	CAR-NAN-CAQ	4.42	122.90	116.83
3	C	1202	P34	CAU-CAR-NAN	-4.19	121.33	123.60
3	A	1202	P34	CAH-CAS-CAQ	-4.03	115.69	121.61
3	A	1202	P34	CAA-N-CA	3.98	116.93	110.38
3	A	1202	P34	CAH-CAS-CAT	3.28	123.11	118.54
3	A	1202	P34	CAB-N-CA	-3.12	105.24	110.38
3	C	1202	P34	CA-C-NAM	3.09	119.92	114.12
3	B	1202	P34	CAH-CAS-CAQ	-2.63	117.76	121.61
3	A	1202	P34	O-C-NAM	-2.62	118.85	123.63
3	C	1202	P34	O-C-NAM	-2.36	119.32	123.63
3	B	1202	P34	CAH-CAS-CAT	2.28	121.71	118.54
3	A	1202	P34	CAP-CAK-CAU	2.19	123.72	120.78
3	B	1202	P34	CAP-CAK-CAU	2.13	123.64	120.78
3	A	1202	P34	O-C-CA	-2.07	117.46	121.08
3	C	1202	P34	CAF-CAI-CAT	2.06	124.33	120.86

There are no chirality outliers.
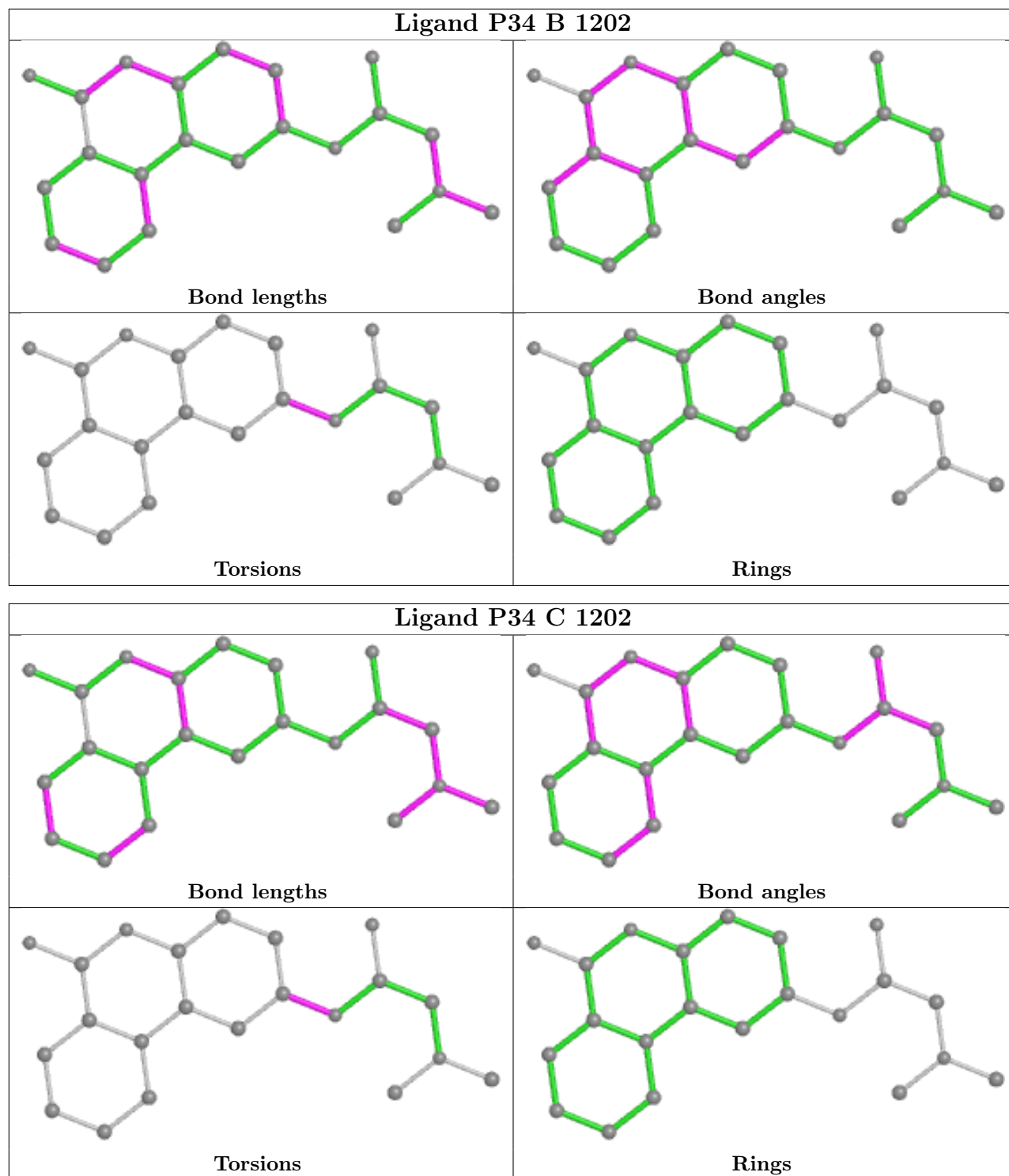
All (9) torsion outliers are listed below:

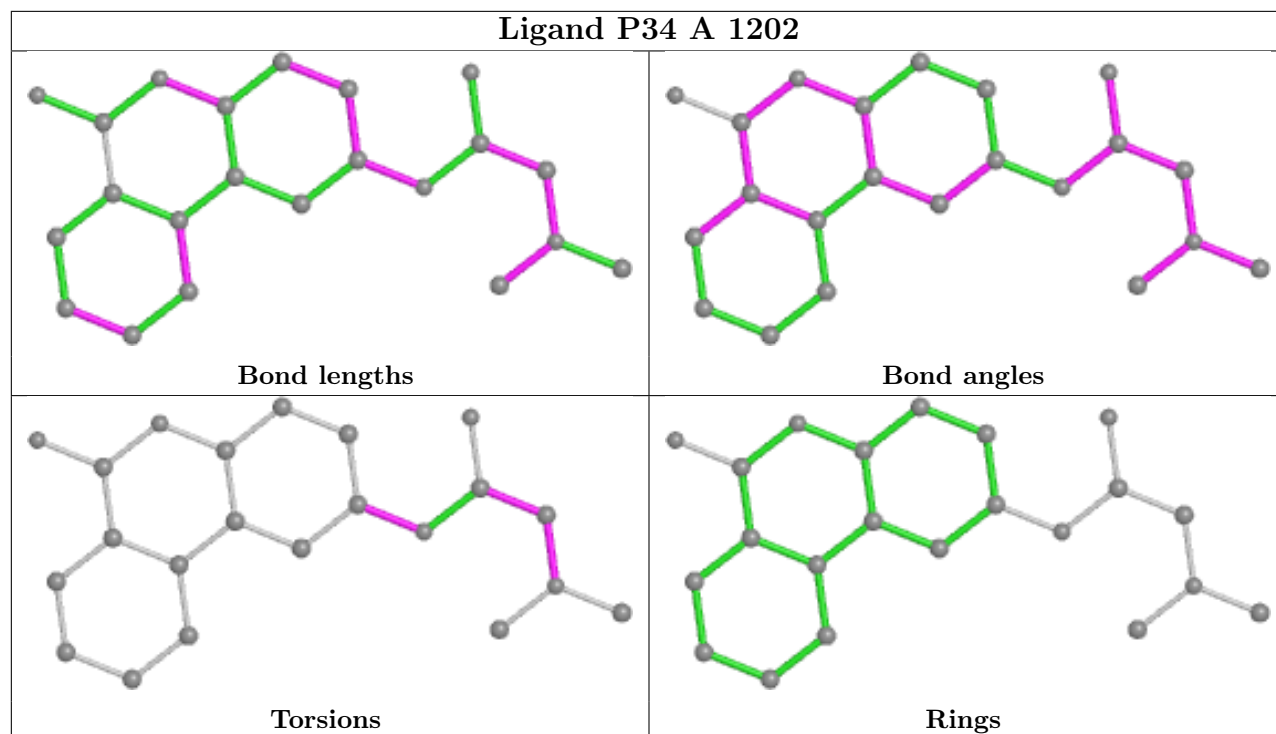| Mol	Chain	Res	Type	Atoms
3	C	1202	P34	CAG-CAP-NAM-C
3	C	1202	P34	CAK-CAP-NAM-C
3	A	1202	P34	NAM-C-CA-N
3	A	1202	P34	O-C-CA-N
3	A	1202	P34	CAK-CAP-NAM-C
3	A	1202	P34	CAG-CAP-NAM-C
3	A	1202	P34	C-CA-N-CAA
3	B	1202	P34	CAK-CAP-NAM-C
3	B	1202	P34	CAG-CAP-NAM-C

There are no ring outliers.

No monomer is involved in short contacts.

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 then 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 P34 B 1202

Bond lengths

Bond angles

Torsions

Rings

Ligand P34 C 1202

Bond lengths

Bond angles

Torsions

Rings

Ligand P34 A 1202

Bond lengths

Bond angles

Torsions

Rings

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å²)	Q<0.9
1	A	199/227 (87%)	-0.07	6 (3%)	50	54	12, 22, 41, 69	0
1	B	207/227 (91%)	-0.04	6 (2%)	51	56	13, 21, 42, 79	0
1	C	208/227 (91%)	-0.07	7 (3%)	45	50	12, 21, 40, 53	0
1	D	191/227 (84%)	0.34	17 (8%)	9	11	14, 28, 53, 67	0
All	All	805/908 (88%)	0.03	36 (4%)	33	37	12, 23, 45, 79	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	960	PRO	5.7
1	D	1135	ALA	5.4
1	D	1136	LEU	5.3
1	B	961	ASP	4.3
1	B	1131	VAL	4.2
1	D	1110	PHE	4.2
1	D	966	GLN	3.9
1	C	955	TYR	3.6
1	D	1051	ILE	3.6
1	A	1131	VAL	3.6
1	D	980	ARG	3.5
1	B	959	SER	3.5
1	D	1129	PRO	3.4
1	A	1161	GLU	3.4
1	D	1050	TYR	3.3
1	D	1083	VAL	3.2
1	D	965	PHE	3.1
1	D	983	GLY	2.9
1	D	1075	ILE	2.9
1	A	1050	TYR	2.9
1	A	1083	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	1113	MET	2.6
1	A	1110	PHE	2.5
1	D	1054	MET	2.5
1	D	1001	CYS	2.5
1	A	1133	GLY	2.5
1	C	961	ASP	2.4
1	C	1021	HIS	2.4
1	C	953	ASN	2.3
1	C	1132	ASN	2.3
1	C	960	PRO	2.3
1	D	1112	ALA	2.1
1	C	1131	VAL	2.1
1	B	1133	GLY	2.1
1	B	1130	SER	2.1
1	D	1037	ASN	2.1

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

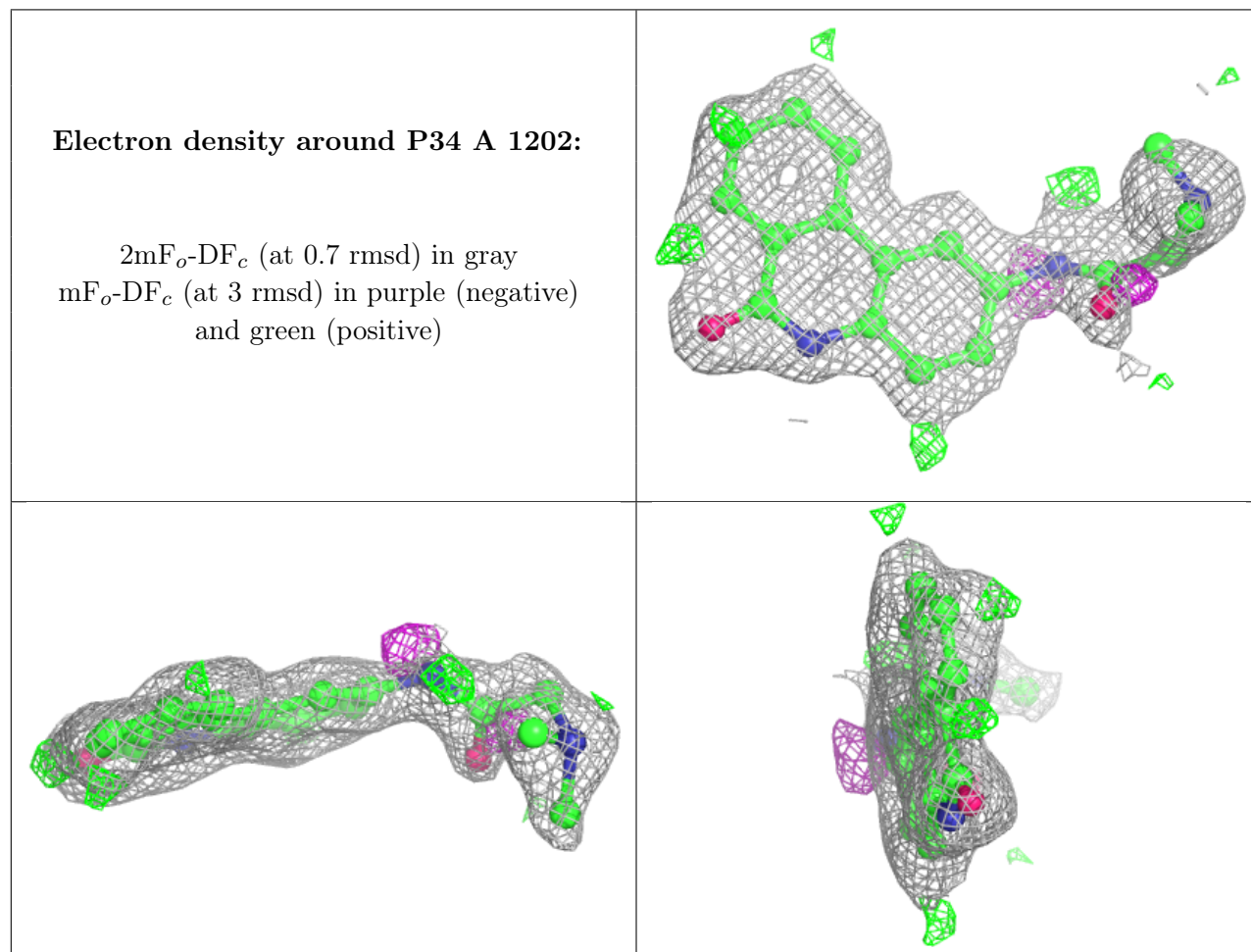There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

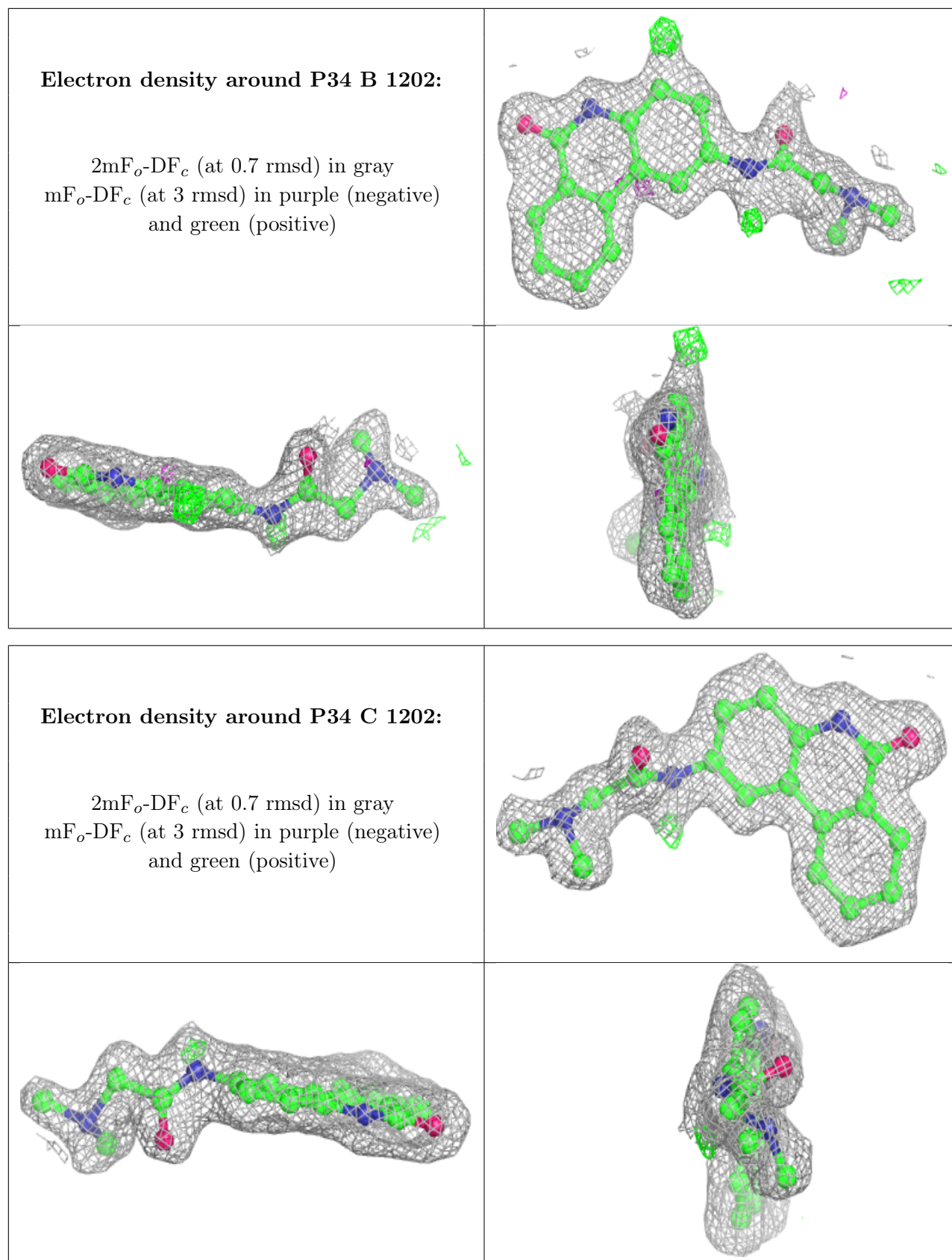## 6.4 Ligands ⓘ

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
3	P34	A	1202	22/22	0.86	0.16	19,24,43,48	0
3	P34	B	1202	22/22	0.90	0.13	21,26,45,49	0
3	P34	C	1202	22/22	0.95	0.09	15,18,42,44	0
2	ZN	C	1201	1/1	0.98	0.04	27,27,27,27	0
2	ZN	B	1201	1/1	0.99	0.03	26,26,26,26	0
2	ZN	A	1201	1/1	0.99	0.03	26,26,26,26	0
2	ZN	D	1200	1/1	0.99	0.02	34,34,34,34	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 P34 A 1202:

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

Electron density around P34 B 1202:

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 P34 C 1202:

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.