



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 10:05 AM EDT

PDB ID : 2QOH / pdb_00002qoh
Title : Crystal Structure of Abl kinase bound with PPY-A
Authors : Zhou, T.; Dalgarno, D.; Zhu, X.
Deposited on : 2007-07-20
Resolution : 1.95 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

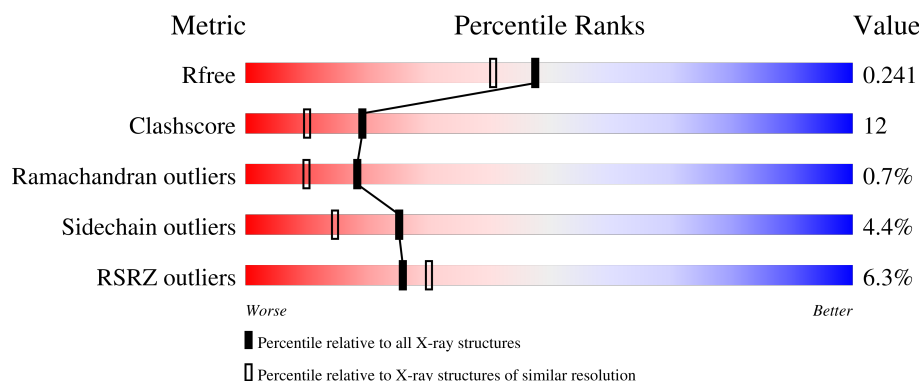
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	288	
1	B	288	

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.

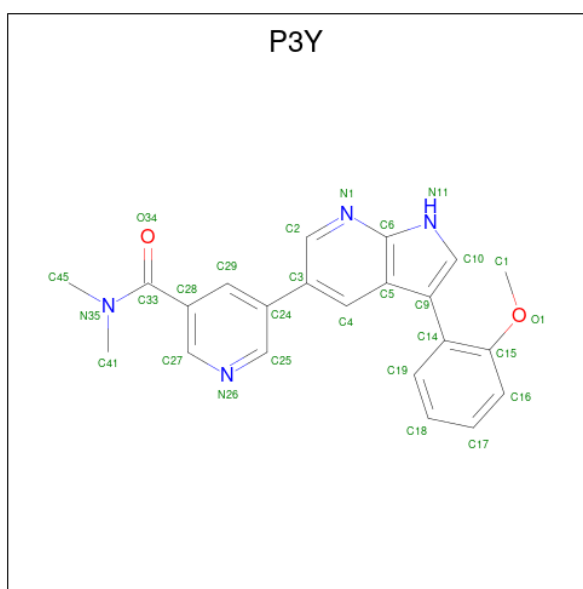
- Molecule 1 is a protein called Proto-oncogene tyrosine-protein kinase ABL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total 2202	C 1418	N 359	O 409	S 16	0	0	0
1	B	281	Total 2278	C 1465	N 370	O 426	S 17	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	GLY	-	expression tag	UNP P00520
B	228	GLY	-	expression tag	UNP P00520

- Molecule 2 is 5-[3-(2-METHOXYPHENYL)-1H-PYRROLO[2,3-B]PYRIDIN-5-YL]-N,N-DIMETHYLPYRIDINE-3-CARBOXAMIDE (CCD ID: P3Y) (formula: $C_{22}H_{20}N_4O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	22	4	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			28	22	4	2		

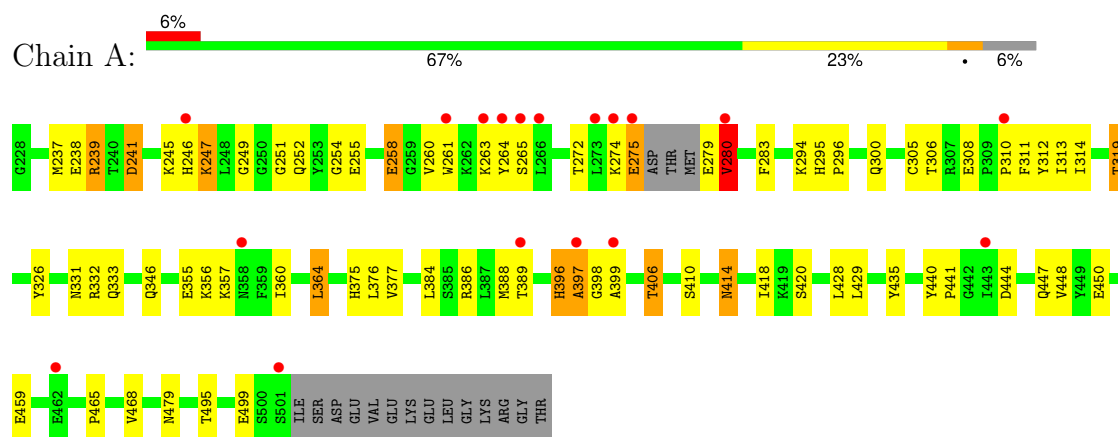
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	125	Total	O	0	0
			125	125		
3	B	132	Total	O	0	0
			132	132		

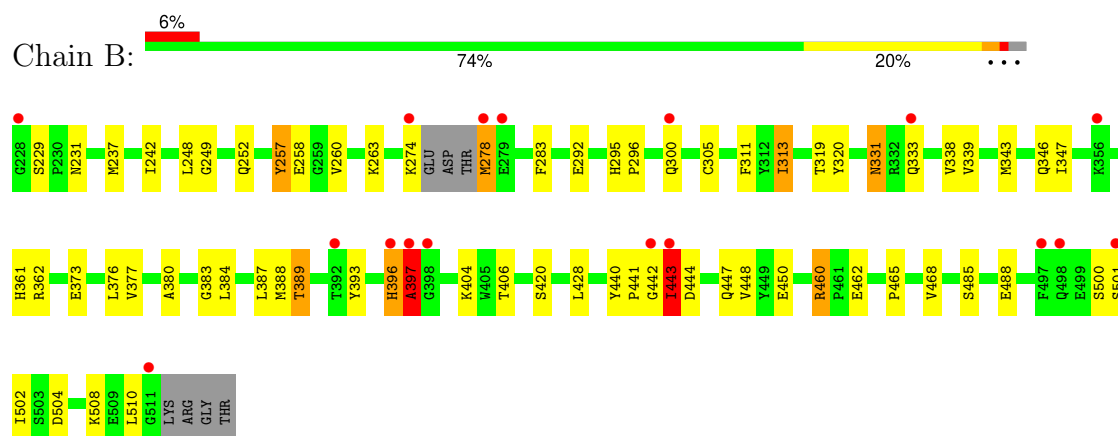
3 Residue-property plots [i](#)

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: Proto-oncogene tyrosine-protein kinase ABL1



- Molecule 1: Proto-oncogene tyrosine-protein kinase ABL1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.01Å 59.13Å 75.94Å 90.00° 118.39° 90.00°	Depositor
Resolution (Å)	50.00 – 1.95 50.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	93.8 (50.00-1.95) 90.7 (50.00-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.94Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.214 , 0.255 0.204 , 0.241	Depositor DCC
R_{free} test set	1893 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.737	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.001 for -h-l,k,h 0.001 for l,k,-h-l 0.033 for h,-k,-h-l 0.027 for -h-l,-k,l 0.025 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4793	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.26 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.0345e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: P3Y

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	0/2259	1.10	15/3056 (0.5%)
1	B	0.53	0/2335	1.03	14/3157 (0.4%)
All	All	0.54	0/4594	1.06	29/6213 (0.5%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	398	GLY	N-CA-C	-13.60	98.89	115.08
1	B	501	SER	N-CA-C	-8.76	101.66	111.82
1	A	280	VAL	N-CA-C	8.05	121.15	111.09
1	B	389	THR	N-CA-C	-7.55	103.06	112.72
1	A	396	HIS	N-CA-C	7.53	120.44	111.33
1	B	319	THR	N-CA-C	7.49	120.33	111.71
1	B	258	GLU	N-CA-C	-6.81	98.16	109.46
1	A	448	VAL	N-CA-C	6.74	117.24	110.23
1	A	333	GLN	N-CA-C	-6.60	104.25	112.90
1	A	275	GLU	N-CA-C	6.42	128.96	111.00
1	B	313	ILE	N-CA-C	-6.33	98.41	107.77
1	A	258	GLU	N-CA-C	-6.30	99.38	109.96
1	A	319	THR	N-CA-C	5.99	119.42	111.75
1	A	300	GLN	N-CA-C	5.82	119.01	109.76
1	A	331	ASN	N-CA-C	-5.63	100.81	109.76
1	A	406	THR	N-CA-C	5.58	118.33	109.24
1	A	389	THR	N-CA-C	-5.54	104.75	112.45
1	B	448	VAL	N-CA-C	5.47	120.72	109.34
1	B	420	SER	N-CA-C	-5.42	105.54	111.82
1	B	397	ALA	N-CA-C	-5.39	99.32	110.80
1	B	300	GLN	N-CA-C	5.35	117.45	109.59
1	B	331	ASN	N-CA-C	-5.28	99.18	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	257	TYR	N-CA-C	5.23	118.27	109.95
1	B	393	TYR	N-CA-C	-5.21	101.20	109.59
1	A	308	GLU	CA-C-N	5.12	123.40	119.66
1	A	308	GLU	C-N-CA	5.12	123.40	119.66
1	A	420	SER	N-CA-C	-5.09	105.81	112.23
1	B	229	SER	N-CA-C	-5.05	98.66	109.81
1	B	406	THR	N-CA-C	5.03	117.81	109.46

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	2202	0	2151	63	0
1	B	2278	0	2228	47	0
2	A	28	0	20	0	0
2	B	28	0	20	1	0
3	A	125	0	0	4	0
3	B	132	0	0	3	0
All	All	4793	0	4419	105	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 12.

All (105) 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:B:444:ASP:HB2	1:B:447:GLN:HE21	1.31	0.95
1:A:237:MET:SD	1:A:314:ILE:HD13	2.14	0.87
1:A:294:LYS:HG2	1:B:274:LYS:NZ	1.95	0.80
1:B:444:ASP:H	1:B:447:GLN:NE2	1.80	0.79
1:B:331:ASN:OD1	1:B:333:GLN:HG2	1.84	0.78
1:B:339:VAL:O	1:B:343:MET:HG3	1.84	0.77
1:B:346:GLN:NE2	1:B:377:VAL:H	1.83	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:444:ASP:CB	1:B:447:GLN:HE21	1.99	0.76
1:A:444:ASP:H	1:A:447:GLN:NE2	1.86	0.73
1:A:238:GLU:O	1:A:241:ASP:HB2	1.89	0.72
1:A:239:ARG:NH2	1:A:310:PRO:O	2.23	0.71
1:A:294:LYS:HG2	1:B:274:LYS:HZ1	1.51	0.71
1:A:249:GLY:O	1:A:252:GLN:HB3	1.93	0.69
1:A:346:GLN:NE2	1:A:377:VAL:H	1.91	0.69
1:A:279:GLU:HG2	1:A:280:VAL:H	1.58	0.68
1:B:460:ARG:HB3	1:B:460:ARG:HH11	1.59	0.67
1:A:396:HIS:O	1:A:397:ALA:CB	2.44	0.65
1:B:346:GLN:HE22	1:B:377:VAL:H	1.44	0.65
1:A:346:GLN:HE22	1:A:376:LEU:HA	1.61	0.65
1:B:249:GLY:O	1:B:252:GLN:HG2	1.97	0.64
1:A:397:ALA:C	1:A:399:ALA:H	2.05	0.62
1:A:396:HIS:O	1:A:397:ALA:HB3	2.00	0.61
1:A:239:ARG:NH1	1:A:305:CYS:HB3	2.18	0.59
1:B:396:HIS:O	1:B:397:ALA:CB	2.51	0.59
1:A:444:ASP:H	1:A:447:GLN:HE21	1.48	0.58
1:A:245:LYS:HE3	1:A:260:VAL:HG23	1.84	0.58
1:A:396:HIS:CE1	1:A:399:ALA:HB2	2.39	0.58
1:A:364:LEU:HD22	1:A:428:LEU:HD22	1.86	0.57
1:B:460:ARG:HH11	1:B:460:ARG:CG	2.16	0.57
1:B:460:ARG:HH11	1:B:460:ARG:CB	2.17	0.57
1:A:295:HIS:CG	1:A:296:PRO:HD2	2.40	0.57
1:A:450:GLU:HG3	3:A:545:HOH:O	2.05	0.57
1:A:237:MET:HE2	1:A:261:TRP:CZ2	2.42	0.55
1:B:346:GLN:HE22	1:B:376:LEU:HA	1.71	0.54
1:A:465:PRO:HG2	1:A:468:VAL:CG2	2.37	0.54
1:B:485:SER:OG	1:B:488:GLU:HG3	2.08	0.54
1:A:239:ARG:HH12	1:A:305:CYS:HB3	1.73	0.53
1:A:396:HIS:CD2	1:A:397:ALA:O	2.61	0.53
1:B:465:PRO:HG2	1:B:468:VAL:CG2	2.38	0.53
1:A:495:THR:O	1:A:499:GLU:HG3	2.08	0.53
1:B:444:ASP:HB2	1:B:447:GLN:NE2	2.13	0.53
1:B:502:ILE:HG23	3:B:624:HOH:O	2.07	0.53
1:A:294:LYS:HD3	1:B:274:LYS:HE2	1.91	0.52
1:A:247:LYS:HD3	1:A:254:GLY:O	2.10	0.52
1:B:450:GLU:HG3	3:B:616:HOH:O	2.11	0.50
1:B:278:MET:SD	1:B:278:MET:N	2.81	0.50
1:A:346:GLN:HE22	1:A:377:VAL:H	1.57	0.50
1:A:319:THR:HG23	3:A:571:HOH:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:LEU:C	1:B:248:LEU:HD12	2.36	0.49
1:B:347:ILE:HD13	1:B:428:LEU:HD21	1.95	0.49
1:A:326:TYR:OH	1:A:375:HIS:CE1	2.66	0.48
1:B:441:PRO:HG3	3:B:549:HOH:O	2.13	0.48
1:B:383:GLY:C	1:B:384:LEU:HD12	2.38	0.48
1:A:255:GLU:HB3	1:A:272:THR:HG22	1.95	0.48
1:A:384:LEU:O	1:A:386:ARG:NH1	2.47	0.48
1:A:283:PHE:CZ	1:A:313:ILE:HG13	2.47	0.48
1:A:294:LYS:HG2	1:B:274:LYS:HZ3	1.78	0.48
1:A:305:CYS:HB2	1:A:312:TYR:HB2	1.95	0.48
1:A:479:ASN:HB3	3:A:527:HOH:O	2.13	0.48
1:A:263:LYS:HG3	1:A:264:TYR:CD1	2.50	0.47
1:A:247:LYS:HD2	1:A:251:GLY:HA2	1.96	0.47
1:B:387:LEU:HD12	1:B:388:MET:N	2.29	0.47
1:A:499:GLU:OE2	1:B:404:LYS:NZ	2.43	0.46
1:A:245:LYS:HE3	1:A:260:VAL:CG2	2.46	0.46
1:A:249:GLY:O	1:A:252:GLN:CB	2.62	0.46
1:A:360:ILE:HG23	1:A:388:MET:HE2	1.98	0.46
1:A:440:TYR:N	1:A:441:PRO:HD3	2.31	0.45
1:A:332:ARG:HG2	1:A:435:TYR:CE2	2.51	0.45
1:B:380:ALA:HB3	2:B:1:P3Y:H13	1.98	0.45
1:B:295:HIS:ND1	1:B:296:PRO:HD2	2.31	0.45
1:A:305:CYS:O	1:A:311:PHE:HA	2.17	0.45
1:B:331:ASN:CG	1:B:333:GLN:HG2	2.42	0.45
1:A:279:GLU:HG2	1:A:280:VAL:HG12	1.99	0.45
1:A:283:PHE:CE2	1:A:311:PHE:HB3	2.52	0.45
1:A:264:TYR:O	1:A:265:SER:C	2.60	0.44
1:A:465:PRO:HG2	1:A:468:VAL:HG23	1.98	0.44
1:B:442:GLY:O	1:B:443:ILE:C	2.60	0.44
1:A:360:ILE:CG2	1:A:388:MET:HE2	2.47	0.44
1:A:306:THR:HA	1:A:311:PHE:CD2	2.52	0.44
1:A:406:THR:CG2	1:A:410:SER:HB2	2.47	0.44
1:A:397:ALA:C	1:A:399:ALA:N	2.71	0.44
1:B:346:GLN:NE2	1:B:376:LEU:HD12	2.33	0.44
1:A:274:LYS:HG2	1:A:275:GLU:H	1.83	0.44
1:A:356:LYS:HB3	3:A:561:HOH:O	2.19	0.43
1:A:414:ASN:HD22	1:A:414:ASN:N	2.14	0.43
1:B:460:ARG:CG	1:B:460:ARG:NH1	2.80	0.43
1:B:504:ASP:O	1:B:508:LYS:HG3	2.18	0.43
1:B:440:TYR:N	1:B:441:PRO:HD3	2.33	0.43
1:A:237:MET:HE2	1:A:261:TRP:CE2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:MET:HE1	1:B:242:ILE:HD11	2.00	0.43
1:B:460:ARG:HH11	1:B:460:ARG:HG2	1.84	0.43
1:B:320:TYR:CE1	1:B:373:GLU:OE2	2.73	0.42
1:B:361:HIS:O	1:B:362:ARG:HB2	2.19	0.42
1:A:241:ASP:OD2	1:A:263:LYS:NZ	2.53	0.41
1:B:305:CYS:O	1:B:311:PHE:HA	2.20	0.41
1:B:460:ARG:NH1	1:B:460:ARG:HG2	2.35	0.41
1:A:384:LEU:N	1:A:384:LEU:HD12	2.35	0.41
1:B:283:PHE:CE2	1:B:313:ILE:HG13	2.56	0.41
1:A:263:LYS:HE2	1:A:264:TYR:HE1	1.84	0.41
1:B:257:TYR:CD1	1:B:257:TYR:N	2.87	0.41
1:B:396:HIS:O	1:B:397:ALA:HB2	2.20	0.41
1:A:246:HIS:CE1	1:A:258:GLU:OE1	2.74	0.41
1:B:346:GLN:HE22	1:B:377:VAL:N	2.14	0.40
1:A:237:MET:SD	1:A:314:ILE:CD1	2.99	0.40
1:A:355:GLU:HA	1:A:418:ILE:HD12	2.03	0.40

There are no symmetry-related clashes.

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	267/288 (93%)	253 (95%)	13 (5%)	1 (0%)	30	21
1	B	277/288 (96%)	264 (95%)	10 (4%)	3 (1%)	11	4
All	All	544/576 (94%)	517 (95%)	23 (4%)	4 (1%)	18	10

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	397	ALA
1	B	397	ALA

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Mol	Chain	Res	Type
1	B	443	ILE
1	B	396	HIS

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	236/251 (94%)	227 (96%)	9 (4%)	29	19
1	B	245/251 (98%)	233 (95%)	12 (5%)	22	11
All	All	481/502 (96%)	460 (96%)	21 (4%)	25	15

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	239	ARG
1	A	241	ASP
1	A	247	LYS
1	A	280	VAL
1	A	357	LYS
1	A	364	LEU
1	A	414	ASN
1	A	429	LEU
1	A	459	GLU
1	B	231	ASN
1	B	260	VAL
1	B	263	LYS
1	B	278	MET
1	B	292	GLU
1	B	338	VAL
1	B	389	THR
1	B	443	ILE
1	B	460	ARG
1	B	462	GLU
1	B	500	SER
1	B	510	LEU

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	252	GLN
1	A	346	GLN
1	A	368	ASN
1	A	375	HIS
1	A	396	HIS
1	A	414	ASN
1	A	447	GLN
1	A	490	HIS
1	B	346	GLN
1	B	368	ASN
1	B	375	HIS
1	B	414	ASN
1	B	447	GLN
1	B	491	GLN
1	B	498	GLN

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](#)

2 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
2	P3Y	B	1	-	31,31,31	1.85	7 (22%)	43,44,44	3.18	8 (18%)
2	P3Y	A	1	-	31,31,31	1.90	9 (29%)	43,44,44	2.97	9 (20%)

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	P3Y	B	1	-	-	6/18/18/18	0/4/4/4
2	P3Y	A	1	-	-	4/18/18/18	0/4/4/4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	P3Y	C10-C9	4.80	1.42	1.37
2	B	1	P3Y	C10-C9	4.31	1.41	1.37
2	B	1	P3Y	C33-N35	3.35	1.39	1.34
2	B	1	P3Y	C4-C5	3.27	1.44	1.39
2	A	1	P3Y	C33-N35	3.21	1.39	1.34
2	A	1	P3Y	C5-C6	-3.15	1.38	1.41
2	A	1	P3Y	C4-C5	3.11	1.44	1.39
2	B	1	P3Y	O1-C15	2.53	1.41	1.37
2	B	1	P3Y	C14-C15	2.48	1.45	1.40
2	B	1	P3Y	C5-C6	-2.45	1.38	1.41
2	A	1	P3Y	O1-C15	2.40	1.41	1.37
2	A	1	P3Y	C16-C15	2.36	1.44	1.39
2	A	1	P3Y	C25-C24	2.23	1.43	1.39
2	A	1	P3Y	C27-C28	2.13	1.42	1.39
2	A	1	P3Y	C14-C15	2.10	1.44	1.40
2	B	1	P3Y	C14-C9	2.01	1.52	1.49

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	P3Y	C5-C6-N11	11.67	112.69	108.42
2	B	1	P3Y	C6-N11-C10	-11.07	103.27	108.48
2	A	1	P3Y	C5-C6-N11	10.69	112.33	108.42
2	A	1	P3Y	C6-N11-C10	-10.14	103.70	108.48
2	B	1	P3Y	C5-C9-C10	-7.00	101.60	105.98
2	A	1	P3Y	C5-C9-C10	-6.39	101.98	105.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	P3Y	C9-C10-N11	6.29	114.53	110.46
2	A	1	P3Y	C9-C10-N11	5.51	114.02	110.46
2	A	1	P3Y	C6-C5-C9	5.01	107.96	106.22
2	B	1	P3Y	C6-C5-C9	4.87	107.91	106.22
2	A	1	P3Y	C5-C6-N1	-4.48	120.89	126.31
2	B	1	P3Y	C5-C6-N1	-4.26	121.15	126.31
2	A	1	P3Y	C14-C9-C5	3.30	131.47	125.97
2	B	1	P3Y	C14-C9-C5	3.19	131.29	125.97
2	A	1	P3Y	C2-N1-C6	2.81	121.69	116.19
2	B	1	P3Y	C2-N1-C6	2.72	121.52	116.19
2	A	1	P3Y	C1-O1-C15	2.29	120.87	117.51

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	P3Y	C28-C33-N35-C41
2	A	1	P3Y	C28-C33-N35-C45
2	A	1	P3Y	O34-C33-N35-C41
2	A	1	P3Y	O34-C33-N35-C45
2	B	1	P3Y	C28-C33-N35-C41
2	B	1	P3Y	C28-C33-N35-C45
2	B	1	P3Y	O34-C33-N35-C41
2	B	1	P3Y	O34-C33-N35-C45
2	B	1	P3Y	C16-C15-O1-C1
2	B	1	P3Y	C14-C15-O1-C1

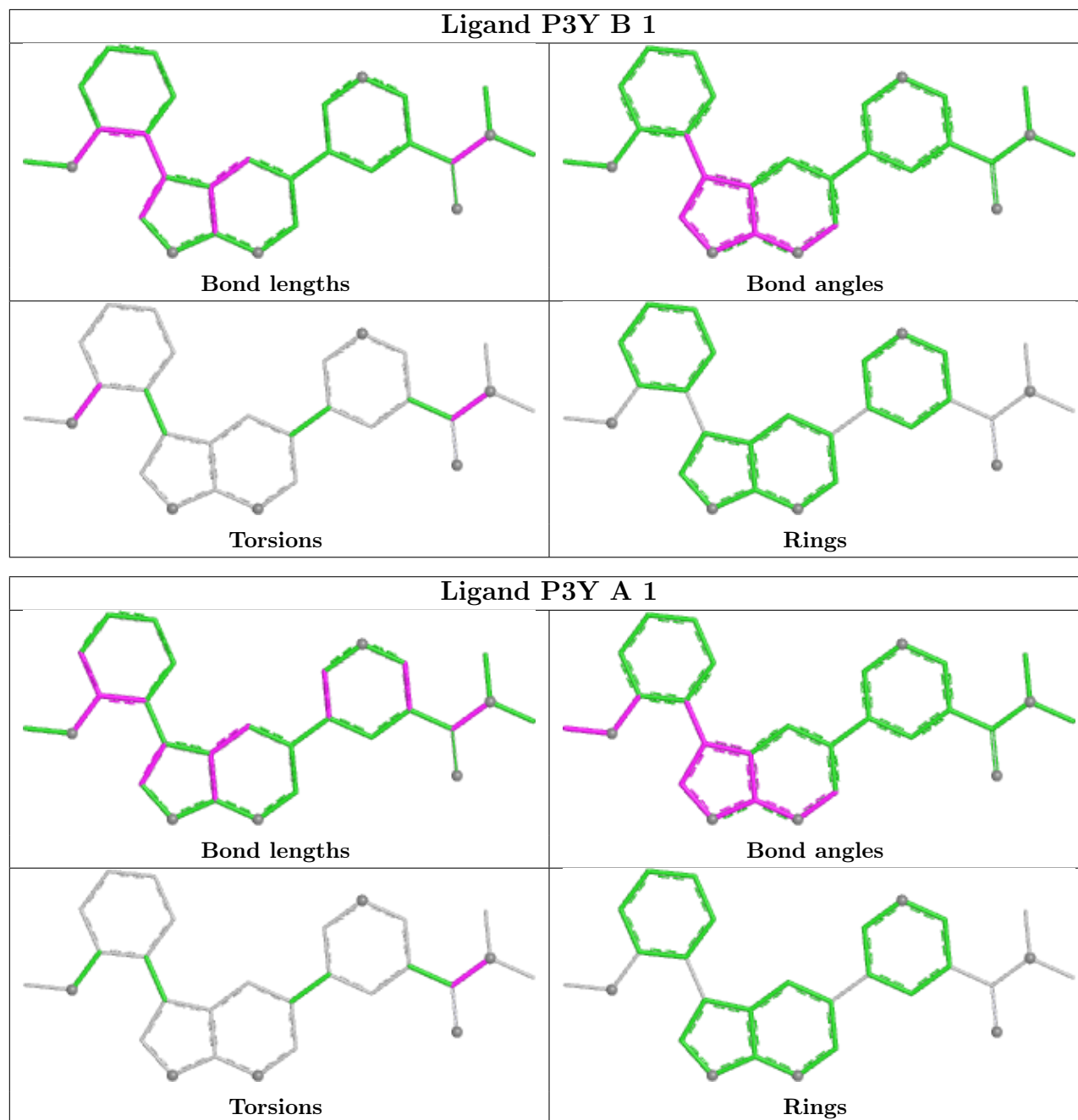
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	P3Y	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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	271/288 (94%)	0.38	18 (6%) 24 28	17, 29, 50, 60	0
1	B	281/288 (97%)	0.43	17 (6%) 27 32	18, 29, 50, 64	0
All	All	552/576 (95%)	0.40	35 (6%) 26 30	17, 29, 50, 64	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	511	GLY	5.8
1	A	397	ALA	3.7
1	B	442	GLY	3.7
1	A	275	GLU	3.4
1	B	396	HIS	3.3
1	A	443	ILE	3.3
1	A	274	LYS	3.3
1	B	443	ILE	3.2
1	B	397	ALA	3.0
1	B	278	MET	2.9
1	B	279	GLU	2.8
1	A	501	SER	2.7
1	A	462	GLU	2.6
1	A	264	TYR	2.6
1	B	398	GLY	2.6
1	A	266	LEU	2.5
1	B	274	LYS	2.5
1	A	273	LEU	2.5
1	A	261	TRP	2.5
1	B	392	THR	2.4
1	A	399	ALA	2.3
1	B	333	GLN	2.3
1	A	358	ASN	2.3
1	A	389	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	263	LYS	2.3
1	A	265	SER	2.3
1	B	497	PHE	2.2
1	A	280	VAL	2.2
1	A	310	PRO	2.2
1	B	356	LYS	2.2
1	B	228	GLY	2.1
1	B	498	GLN	2.1
1	A	246	HIS	2.0
1	B	501	SER	2.0
1	B	300	GLN	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 [i](#)

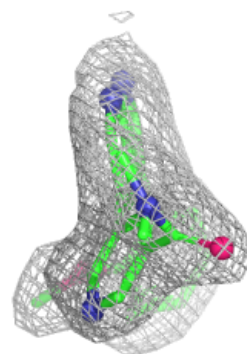
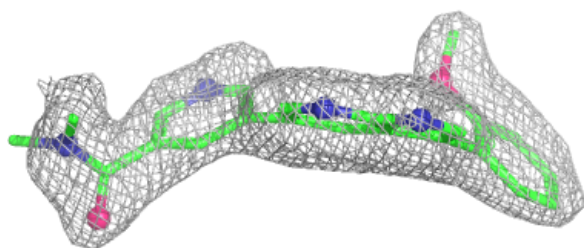
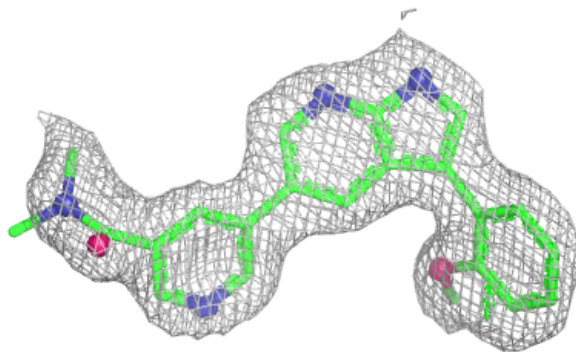
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	P3Y	B	1	28/28	0.89	0.10	30,33,42,45	0
2	P3Y	A	1	28/28	0.92	0.08	24,29,39,42	0

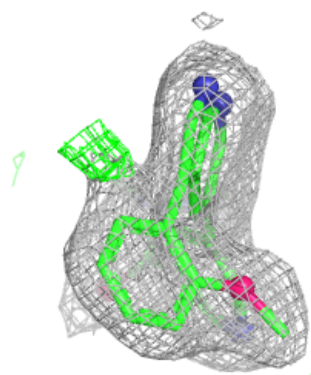
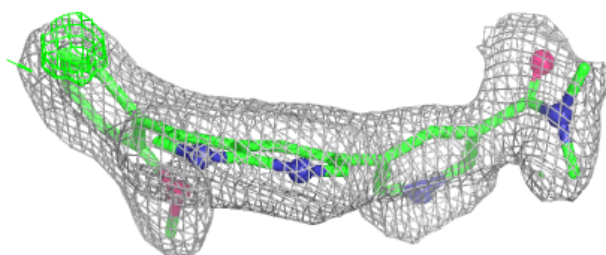
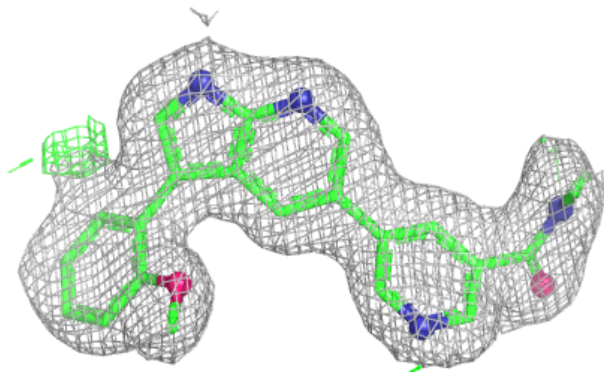
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around P3Y B 1:

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

**Electron density around P3Y A 1:**

$2mF_o-DF_c$ (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.