



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:32 PM UTC

PDB ID : 3GEQ / pdb_00003geq
Title : Structural basis for the chemical rescue of Src kinase activity
Authors : Muratore, K.E.; Seeliger, M.A.; Wang, Z.; Fomina, D.; Neiswinger, J.; Havranek, J.J.; Baker, D.; Kuriyan, J.; Cole, P.A.
Deposited on : 2009-02-25
Resolution : 2.20 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

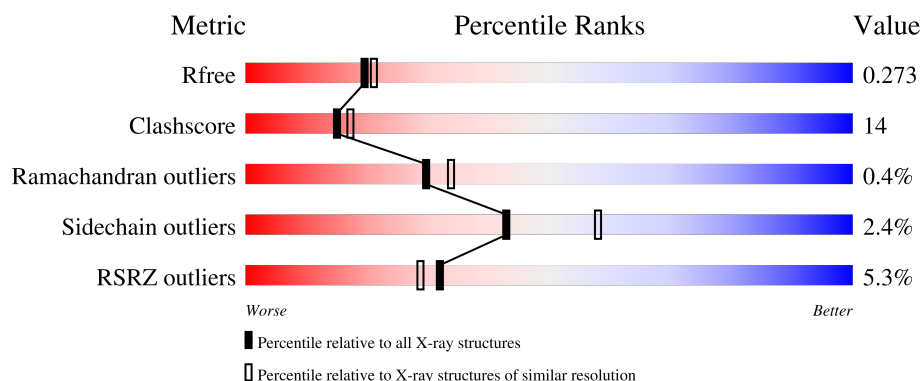
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	286	4% 68% 26% • 5%
1	B	286	6% 64% 30% • 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4619 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.

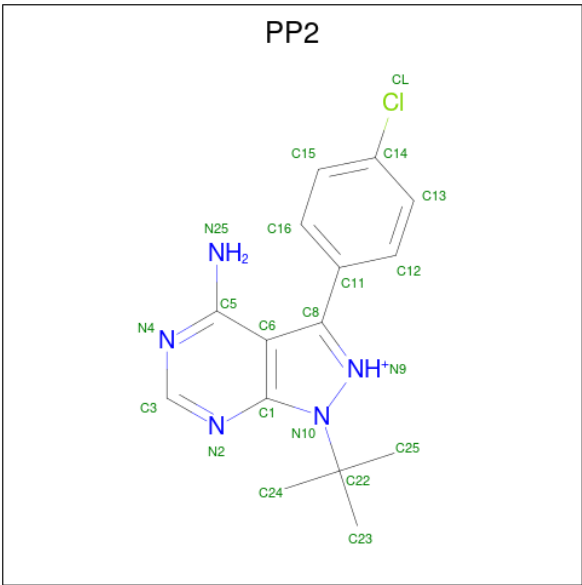
- Molecule 1 is a protein called Proto-oncogene tyrosine-protein kinase Src.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	0	0
			2174	1392	365	399	18			
1	B	273	Total	C	N	O	S	0	0	0
			2182	1398	366	400	18			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	248	GLY	-	expression tag	UNP P00523
A	249	HIS	-	expression tag	UNP P00523
A	250	MET	-	expression tag	UNP P00523
A	388	ALA	ARG	engineered mutation	UNP P00523
B	248	GLY	-	expression tag	UNP P00523
B	249	HIS	-	expression tag	UNP P00523
B	250	MET	-	expression tag	UNP P00523
B	388	ALA	ARG	engineered mutation	UNP P00523

- Molecule 2 is 1-TERT-BUTYL-3-(4-CHLORO-PHENYL)-1H-PYRAZOLO[3,4-D]PYRIMIDIN-4-YLAMINE (CCD ID: PP2) (formula: C₁₅H₁₇ClN₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	0	0
			21	15	1	5		
2	B	1	Total	C	Cl	N	0	0
			21	15	1	5		

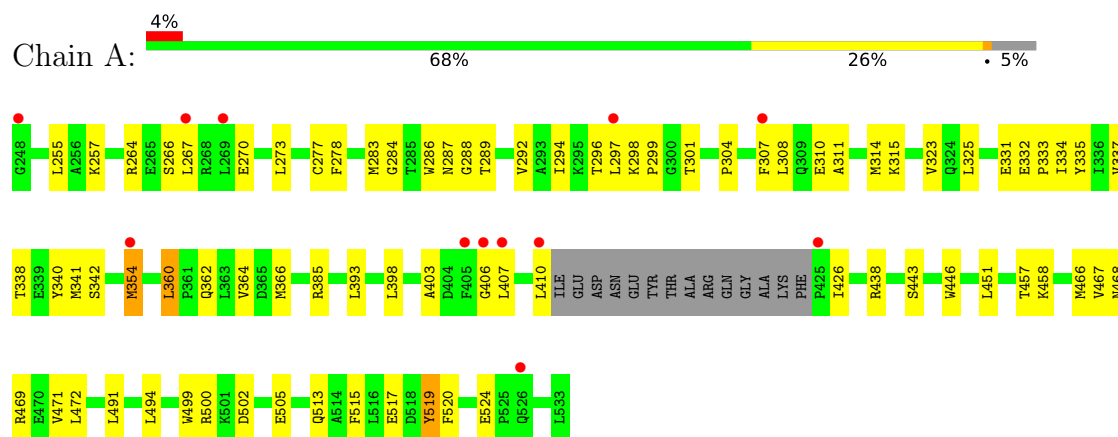
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	105	Total	O	0	0
			105	105		
3	B	116	Total	O	0	0
			116	116		

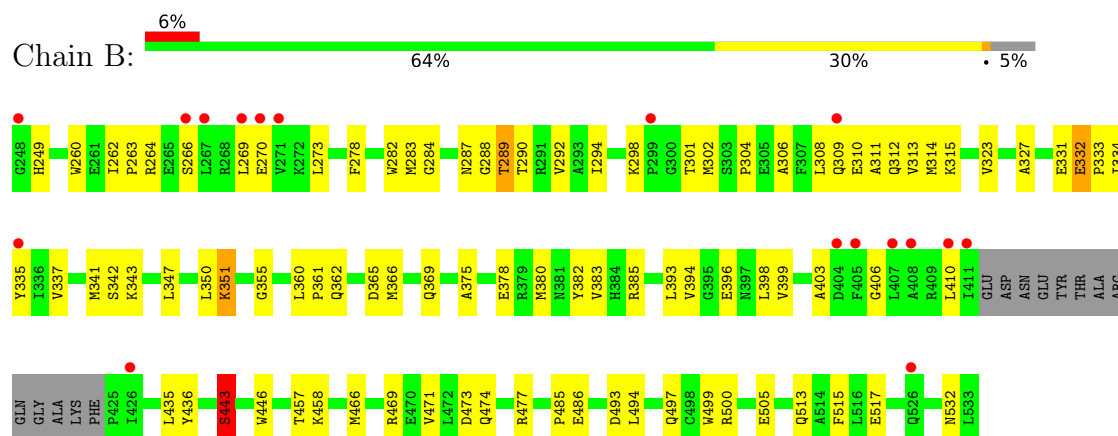
3 Residue-property plots

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: Proto-oncogene tyrosine-protein kinase Src



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.65Å 63.46Å 83.86Å 90.00° 91.45° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	89.0 (50.00-2.20) 89.0 (50.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.226 , 0.273 0.225 , 0.273	Depositor DCC
R_{free} test set	1405 reflections (4.40%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.063 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4619	wwPDB-VP
Average B, all atoms (Å ²)	46.0	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 87.45 % 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 6.8315e-08. 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: PP2

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	0.46	1/2226 (0.0%)	0.89	4/3012 (0.1%)
1	B	0.45	0/2234	0.90	2/3023 (0.1%)
All	All	0.46	1/4460 (0.0%)	0.89	6/6035 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	354	MET	SD-CE	6.23	1.95	1.79

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	342	SER	N-CA-C	8.81	121.84	111.71
1	B	342	SER	N-CA-C	7.16	121.12	112.38
1	A	443	SER	N-CA-C	-5.86	105.39	112.54
1	A	520	PHE	N-CA-C	5.62	119.75	113.01
1	A	519	TYR	N-CA-C	5.37	120.17	111.37
1	B	443	SER	N-CA-C	-5.07	105.94	111.82

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	2174	0	2163	55	0
1	B	2182	0	2174	67	0
2	A	21	0	17	0	0
2	B	21	0	17	0	0
3	A	105	0	0	0	0
3	B	116	0	0	1	0
All	All	4619	0	4371	120	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 14.

All (120) 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:323:VAL:HG21	1:B:393:LEU:HD12	1.39	1.04
1:A:426:ILE:HD11	1:A:472:LEU:HG	1.40	1.01
1:B:500:ARG:HD3	1:B:505:GLU:HB3	1.58	0.85
1:A:385:ARG:HD3	1:A:407:LEU:O	1.82	0.79
1:B:323:VAL:CG2	1:B:393:LEU:HD12	2.13	0.79
1:A:310:GLU:HB2	1:A:406:GLY:HA2	1.66	0.78
1:A:278:PHE:HD1	1:A:301:THR:HG22	1.50	0.75
1:B:260:TRP:CD1	1:B:315:LYS:HE3	2.23	0.73
1:A:502:ASP:HB3	1:A:505:GLU:HG3	1.69	0.72
1:B:443:SER:HB2	3:B:215:HOH:O	1.89	0.70
1:B:294:ILE:HG21	1:B:335:TYR:HD2	1.54	0.70
1:B:310:GLU:HB2	1:B:406:GLY:HA2	1.73	0.69
1:B:327:ALA:H	1:B:337:VAL:HG13	1.58	0.68
1:A:278:PHE:CD1	1:A:301:THR:HG22	2.30	0.66
1:B:304:PRO:O	1:B:308:LEU:HD13	1.96	0.64
1:B:273:LEU:HD11	1:B:283:MET:HB2	1.80	0.64
1:B:311:ALA:O	1:B:315:LYS:HG3	1.99	0.62
1:B:382:TYR:CE1	1:B:410:LEU:HB2	2.34	0.62
1:B:263:PRO:HG2	1:B:266:SER:HB2	1.81	0.62
1:A:323:VAL:HG21	1:A:393:LEU:HD12	1.83	0.61
1:A:426:ILE:HD11	1:A:472:LEU:CG	2.24	0.61
1:B:308:LEU:HD11	1:B:334:ILE:HD13	1.83	0.61
1:A:332:GLU:HA	1:A:333:PRO:C	2.26	0.60
1:A:266:SER:HB2	1:A:287:ASN:OD1	2.02	0.60
1:A:323:VAL:HG21	1:A:403:ALA:HB2	1.83	0.60
1:B:287:ASN:O	1:B:289:THR:HG22	2.01	0.60
1:A:513:GLN:O	1:A:517:GLU:HG3	2.02	0.60
1:B:327:ALA:HB3	1:B:337:VAL:HG11	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:VAL:HG23	1:A:294:ILE:HD11	1.84	0.58
1:B:327:ALA:HB3	1:B:337:VAL:CG1	2.34	0.58
1:B:308:LEU:O	1:B:312:GLN:HG2	2.03	0.58
1:A:273:LEU:HD11	1:A:283:MET:HB2	1.85	0.58
1:B:323:VAL:HG21	1:B:403:ALA:HB2	1.86	0.57
1:A:362:GLN:O	1:A:366:MET:HG3	2.05	0.57
1:B:473:ASP:O	1:B:477:ARG:HG3	2.04	0.57
1:A:500:ARG:HD3	1:A:505:GLU:HB3	1.87	0.56
1:B:278:PHE:HD1	1:B:301:THR:HB	1.70	0.56
1:B:292:VAL:HG23	1:B:294:ILE:HD11	1.89	0.55
1:B:332:GLU:HA	1:B:333:PRO:C	2.32	0.55
1:A:467:VAL:HG22	1:A:468:ASN:N	2.22	0.54
1:B:270:GLU:HG3	1:B:284:GLY:HA2	1.90	0.54
1:B:396:GLU:O	1:B:399:VAL:HG23	2.08	0.53
1:B:466:MET:HE2	1:B:471:VAL:HA	1.90	0.53
1:B:513:GLN:O	1:B:517:GLU:HG3	2.08	0.53
1:A:286:TRP:O	1:A:287:ASN:HB2	2.10	0.52
1:A:298:LYS:HD2	1:A:299:PRO:HD2	1.90	0.52
1:B:289:THR:HG23	1:B:290:THR:H	1.74	0.52
1:B:362:GLN:O	1:B:366:MET:HG3	2.09	0.52
1:A:494:LEU:HD22	1:A:515:PHE:CE1	2.45	0.51
1:B:341:MET:HG3	1:B:393:LEU:HB3	1.93	0.51
1:A:466:MET:CE	1:A:471:VAL:HA	2.40	0.51
1:A:332:GLU:HA	1:A:334:ILE:N	2.25	0.51
1:A:270:GLU:HG3	1:A:284:GLY:HA2	1.92	0.51
1:B:264:ARG:NH2	1:B:331:GLU:O	2.44	0.51
1:B:343:LYS:HB2	1:B:394:VAL:HB	1.93	0.51
1:B:493:ASP:O	1:B:497:GLN:HG3	2.10	0.51
1:A:294:ILE:N	1:A:294:ILE:HD12	2.26	0.51
1:B:485:PRO:O	1:B:486:GLU:HB2	2.11	0.51
1:B:382:TYR:CZ	1:B:410:LEU:HB2	2.46	0.51
1:A:310:GLU:O	1:A:314:MET:HG3	2.10	0.50
1:B:466:MET:CE	1:B:471:VAL:HA	2.41	0.50
1:A:267:LEU:HD21	1:A:337:VAL:HG21	1.94	0.50
1:B:308:LEU:HD11	1:B:334:ILE:HG21	1.94	0.50
1:B:383:VAL:HG12	1:B:385:ARG:HG2	1.94	0.49
1:B:308:LEU:CD1	1:B:334:ILE:HD13	2.43	0.49
1:A:264:ARG:NH2	1:A:331:GLU:O	2.45	0.48
1:A:341:MET:HG3	1:A:393:LEU:HB3	1.94	0.48
1:B:343:LYS:HE2	1:B:396:GLU:OE2	2.13	0.48
1:A:467:VAL:O	1:A:471:VAL:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:347:LEU:O	1:B:351:LYS:HG2	2.13	0.48
1:A:426:ILE:CD1	1:A:472:LEU:HG	2.29	0.47
1:A:304:PRO:O	1:A:308:LEU:HG	2.15	0.47
1:A:451:LEU:C	1:A:451:LEU:HD23	2.40	0.47
1:A:360:LEU:HD22	1:A:364:VAL:HG23	1.97	0.46
1:B:260:TRP:O	1:B:262:ILE:HG23	2.15	0.46
1:A:255:LEU:C	1:A:257:LYS:H	2.23	0.46
1:A:354:MET:SD	1:B:435:LEU:O	2.74	0.46
1:A:323:VAL:CG2	1:A:403:ALA:HB2	2.45	0.46
1:A:273:LEU:HD11	1:A:340:TYR:HE1	1.80	0.46
1:A:297:LEU:HD22	1:A:307:PHE:CG	2.50	0.46
1:B:494:LEU:HD22	1:B:515:PHE:CE1	2.51	0.46
1:A:287:ASN:O	1:A:289:THR:N	2.50	0.45
1:A:311:ALA:O	1:A:315:LYS:HG3	2.16	0.45
1:B:327:ALA:N	1:B:337:VAL:HG13	2.30	0.45
1:B:266:SER:OG	1:B:287:ASN:OD1	2.35	0.45
1:B:446:TRP:CE3	1:B:499:TRP:HA	2.52	0.45
1:A:502:ASP:HB3	1:A:505:GLU:CG	2.44	0.44
1:B:375:ALA:O	1:B:378:GLU:HB3	2.17	0.44
1:A:325:LEU:HA	1:A:338:THR:HG22	2.00	0.44
1:A:277:CYS:O	1:A:301:THR:HG21	2.17	0.44
1:B:310:GLU:O	1:B:314:MET:HG3	2.18	0.44
1:B:466:MET:HE3	1:B:474:GLN:HG3	2.00	0.44
1:B:249:HIS:CD2	1:B:249:HIS:N	2.84	0.43
1:A:296:THR:HG22	1:A:335:TYR:CD1	2.54	0.43
1:A:458:LYS:HA	1:A:458:LYS:HD3	1.89	0.43
1:A:323:VAL:CG2	1:A:393:LEU:HD12	2.48	0.43
1:A:287:ASN:C	1:A:289:THR:H	2.27	0.42
1:A:308:LEU:O	1:A:311:ALA:HB3	2.18	0.42
1:A:466:MET:HE2	1:A:471:VAL:HG22	2.01	0.42
1:B:365:ASP:O	1:B:369:GLN:HG3	2.19	0.42
1:A:292:VAL:HG23	1:A:294:ILE:CD1	2.48	0.42
1:B:332:GLU:HA	1:B:334:ILE:N	2.34	0.42
1:B:294:ILE:HD12	1:B:294:ILE:N	2.33	0.42
1:B:350:LEU:O	1:B:355:GLY:HA3	2.18	0.42
1:B:309:GLN:O	1:B:313:VAL:HG23	2.20	0.42
1:B:298:LYS:HB3	1:B:301:THR:CG2	2.49	0.42
1:B:302:MET:HE2	1:B:306:ALA:C	2.45	0.42
1:B:486:GLU:OE1	1:B:532:ASN:HB2	2.20	0.42
1:B:360:LEU:HB3	1:B:361:PRO:HD3	2.02	0.41
1:A:446:TRP:CE3	1:A:499:TRP:HA	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:THR:O	1:A:458:LYS:HB2	2.21	0.41
1:A:519:TYR:OH	1:A:524:GLU:HG3	2.21	0.41
1:B:351:LYS:HG2	1:B:351:LYS:H	1.76	0.41
1:B:287:ASN:C	1:B:289:THR:H	2.29	0.41
1:B:323:VAL:CG2	1:B:403:ALA:HB2	2.48	0.41
1:A:398:LEU:HD13	1:B:436:TYR:HB3	2.02	0.40
1:B:380:MET:HA	1:B:380:MET:HE2	2.04	0.40
1:A:451:LEU:HD23	1:A:451:LEU:O	2.21	0.40
1:B:457:THR:O	1:B:458:LYS:HB2	2.21	0.40
1:B:269:LEU:HD22	1:B:282:TRP:HB2	2.04	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	268/286 (94%)	252 (94%)	15 (6%)	1 (0%)	30	34
1	B	269/286 (94%)	257 (96%)	11 (4%)	1 (0%)	30	34
All	All	537/572 (94%)	509 (95%)	26 (5%)	2 (0%)	30	34

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	288	GLY
1	B	288	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	233/244 (96%)	228 (98%)	5 (2%)	47	63
1	B	234/244 (96%)	228 (97%)	6 (3%)	40	55
All	All	467/488 (96%)	456 (98%)	11 (2%)	43	58

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	360	LEU
1	A	410	LEU
1	A	438	ARG
1	A	469	ARG
1	A	491	LEU
1	B	289	THR
1	B	332	GLU
1	B	351	LYS
1	B	398	LEU
1	B	443	SER
1	B	469	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	249	HIS
1	A	253	GLN
1	A	309	GLN
1	A	513	GLN
1	B	249	HIS
1	B	492	HIS

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 [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	PP2	A	1	-	23,23,23	1.51	6 (26%)	32,35,35	2.52	9 (28%)
2	PP2	B	2	-	23,23,23	1.54	5 (21%)	32,35,35	2.69	12 (37%)

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	PP2	A	1	-	-	4/10/10/10	0/3/3/3
2	PP2	B	2	-	-	4/10/10/10	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	PP2	C12-C11	2.76	1.43	1.39
2	B	2	PP2	C15-C14	2.71	1.43	1.38
2	B	2	PP2	C1-N10	2.62	1.38	1.35
2	A	1	PP2	C15-C14	2.59	1.42	1.38
2	A	1	PP2	C6-C1	2.50	1.44	1.40
2	A	1	PP2	C1-N10	2.47	1.38	1.35
2	B	2	PP2	C16-C15	2.47	1.42	1.38
2	B	2	PP2	C6-C1	2.46	1.44	1.40
2	A	1	PP2	C12-C11	2.38	1.43	1.39
2	A	1	PP2	C16-C15	2.33	1.42	1.38
2	A	1	PP2	C16-C11	2.21	1.42	1.39

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	PP2	C6-C1-N2	-7.43	119.08	126.97
2	A	1	PP2	C6-C1-N2	-7.27	119.25	126.97
2	B	2	PP2	N2-C1-N10	5.40	134.00	125.48
2	A	1	PP2	N2-C1-N10	5.39	133.98	125.48
2	B	2	PP2	C11-C8-N9	4.70	125.02	118.91
2	A	1	PP2	C6-C8-N9	-4.57	107.05	110.19
2	B	2	PP2	N4-C3-N2	-4.41	121.90	128.58
2	B	2	PP2	C6-C8-N9	-4.41	107.16	110.19
2	A	1	PP2	N4-C3-N2	-4.41	121.91	128.58
2	B	2	PP2	C22-N10-N9	3.95	124.95	117.90
2	A	1	PP2	C22-N10-N9	3.84	124.76	117.90
2	A	1	PP2	C3-N2-C1	3.79	121.09	111.83
2	B	2	PP2	C3-N2-C1	3.79	121.09	111.83
2	A	1	PP2	C11-C8-N9	3.76	123.80	118.91
2	B	2	PP2	C6-C8-C11	-2.92	127.81	131.28
2	B	2	PP2	C5-C6-C8	-2.90	136.13	139.07
2	B	2	PP2	C5-C6-C1	2.54	117.64	115.74
2	A	1	PP2	C5-C6-C8	-2.42	136.62	139.07
2	B	2	PP2	C12-C11-C8	-2.36	117.72	120.64
2	A	1	PP2	C5-C6-C1	2.23	117.41	115.74
2	B	2	PP2	C1-C6-C8	2.13	106.22	104.74

There are no chirality outliers.

All (8) torsion outliers are listed below:

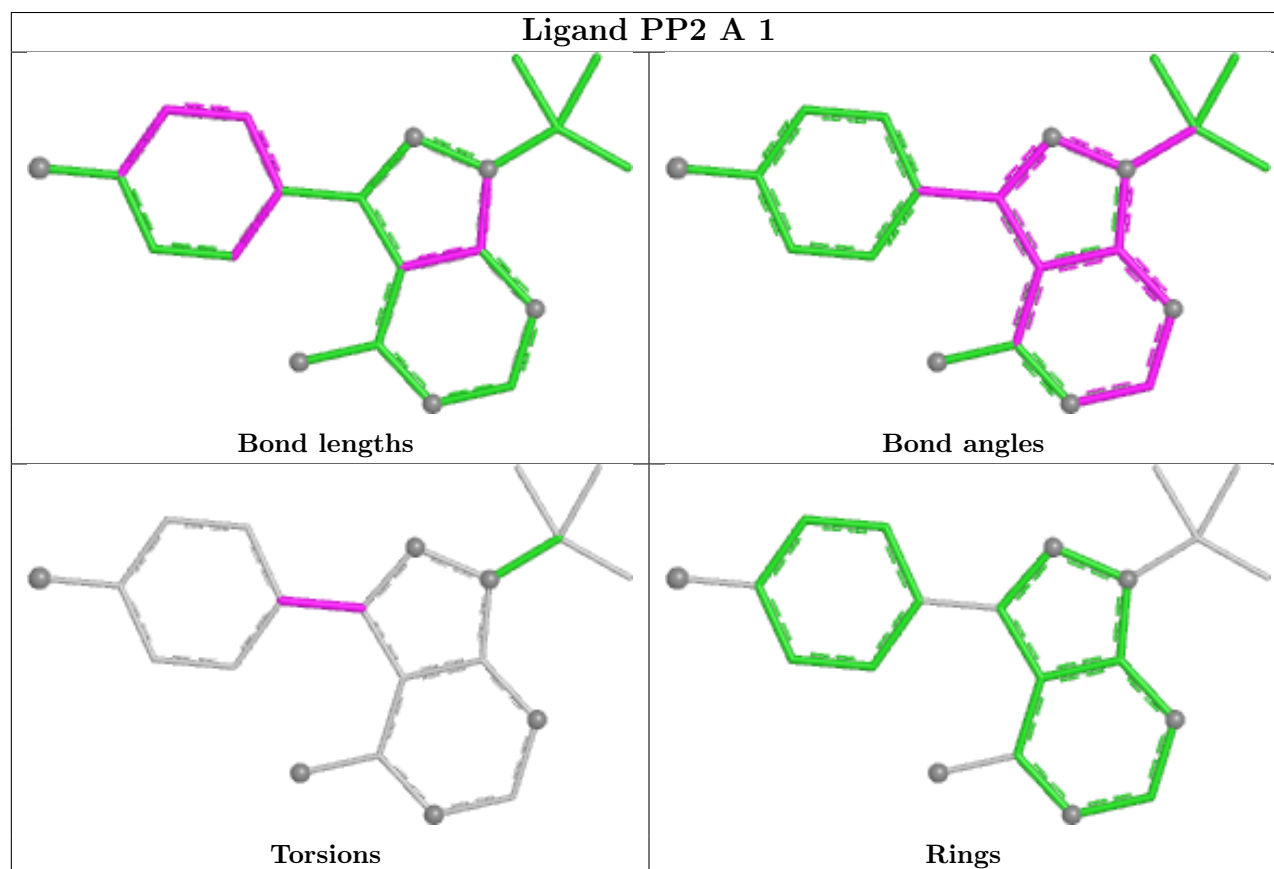
Mol	Chain	Res	Type	Atoms
2	B	2	PP2	C16-C11-C8-C6
2	A	1	PP2	C16-C11-C8-N9
2	B	2	PP2	C12-C11-C8-N9
2	B	2	PP2	C16-C11-C8-N9
2	B	2	PP2	C12-C11-C8-C6
2	A	1	PP2	C12-C11-C8-N9
2	A	1	PP2	C16-C11-C8-C6
2	A	1	PP2	C12-C11-C8-C6

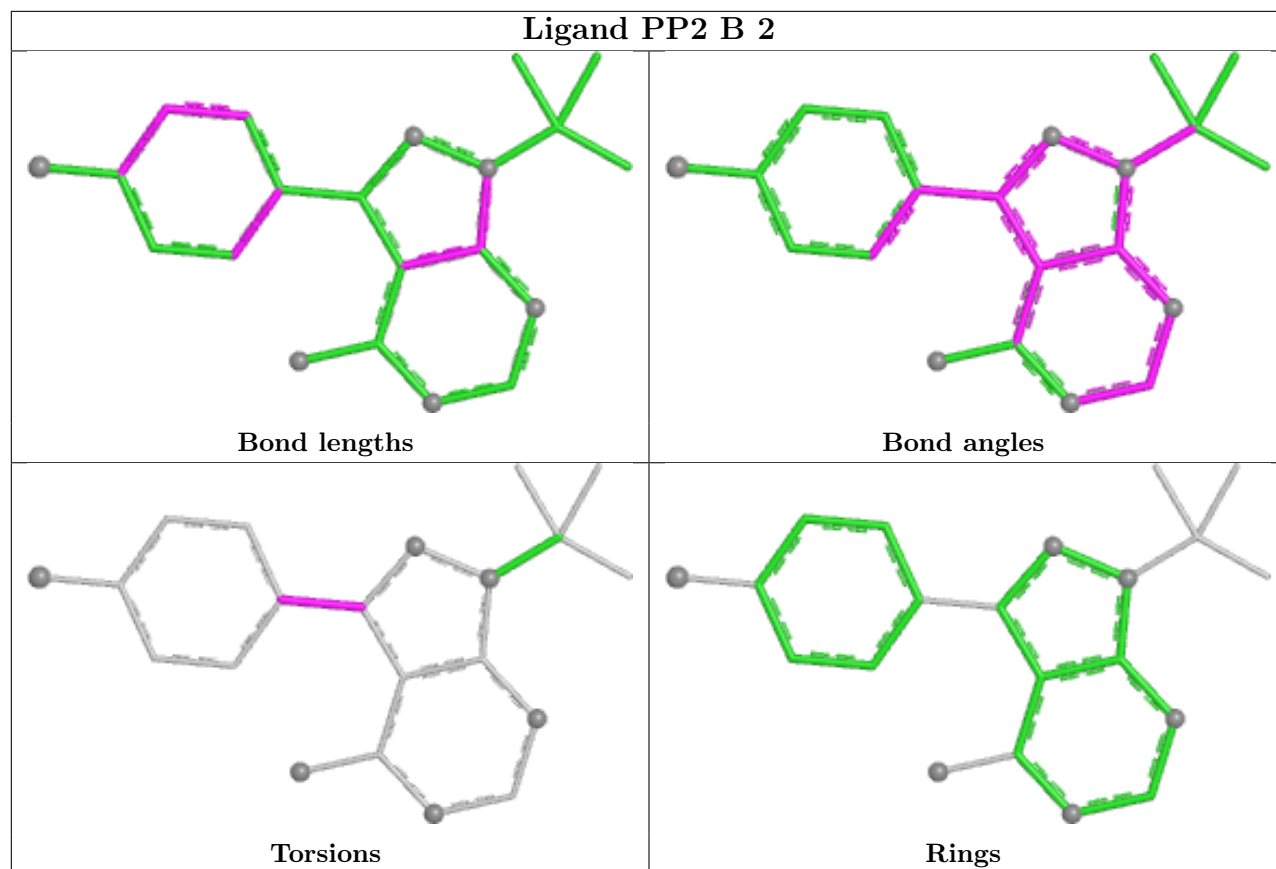
There are no ring outliers.

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 than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	272/286 (95%)	0.48	12 (4%) 39 36	24, 44, 85, 97	0
1	B	273/286 (95%)	0.38	17 (6%) 26 23	23, 41, 83, 90	0
All	All	545/572 (95%)	0.43	29 (5%) 32 29	23, 42, 84, 97	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	407	LEU	4.0
1	A	354	MET	3.9
1	B	248	GLY	3.8
1	B	407	LEU	3.7
1	B	411	ILE	3.5
1	A	248	GLY	3.5
1	B	267	LEU	3.2
1	B	309	GLN	3.2
1	A	269	LEU	2.9
1	A	410	LEU	2.9
1	A	526	GLN	2.8
1	B	426	ILE	2.7
1	A	425	PRO	2.6
1	B	405	PHE	2.5
1	B	270	GLU	2.4
1	B	526	GLN	2.4
1	B	269	LEU	2.4
1	B	404	ASP	2.4
1	A	406	GLY	2.3
1	A	405	PHE	2.2
1	B	410	LEU	2.2
1	B	335	TYR	2.2
1	B	266	SER	2.2
1	A	267	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	297	LEU	2.1
1	B	299	PRO	2.1
1	B	408	ALA	2.1
1	B	271	VAL	2.0
1	A	307	PHE	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](#)

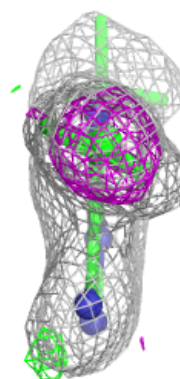
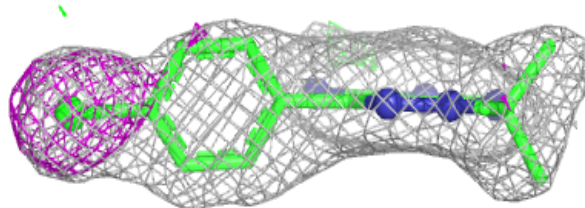
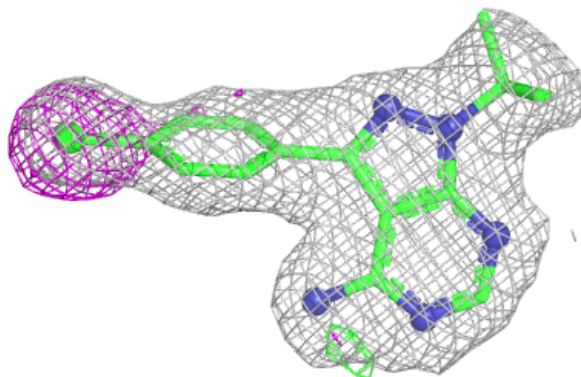
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

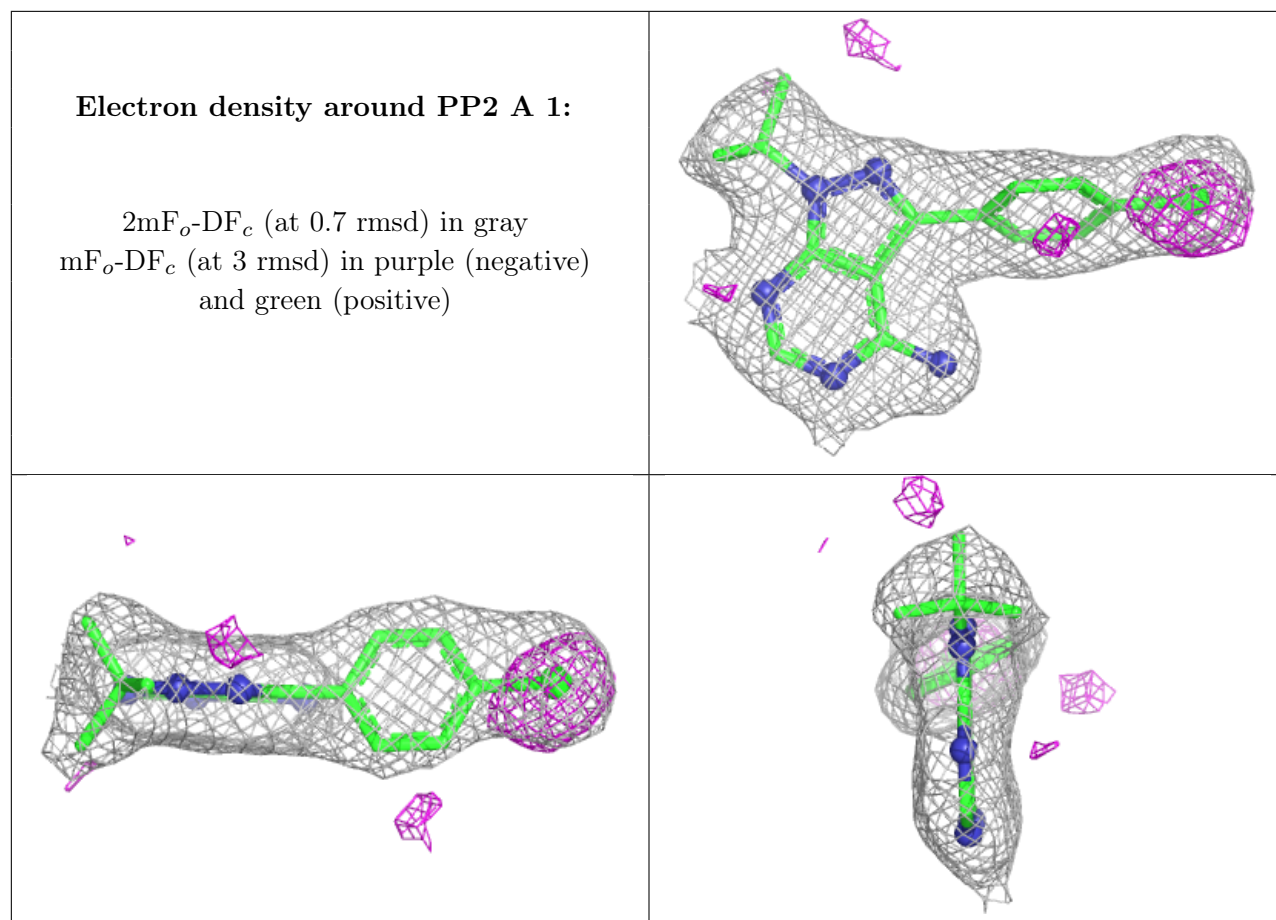
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PP2	B	2	21/21	0.87	0.12	25,34,38,39	0
2	PP2	A	1	21/21	0.89	0.10	34,38,41,42	0

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 PP2 B 2:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.