



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:28 PM UTC

PDB ID : 6BN5 / pdb_00006bn5
Title : Non-receptor Protein Tyrosine Phosphatase SHP2 F285S in Complex with Allosteric Inhibitor JLR-2
Authors : Blacklow, S.C.; Stams, T.; Fodor, M.; LaRochelle, J.R.
Deposited on : 2017-11-16
Resolution : 2.22 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

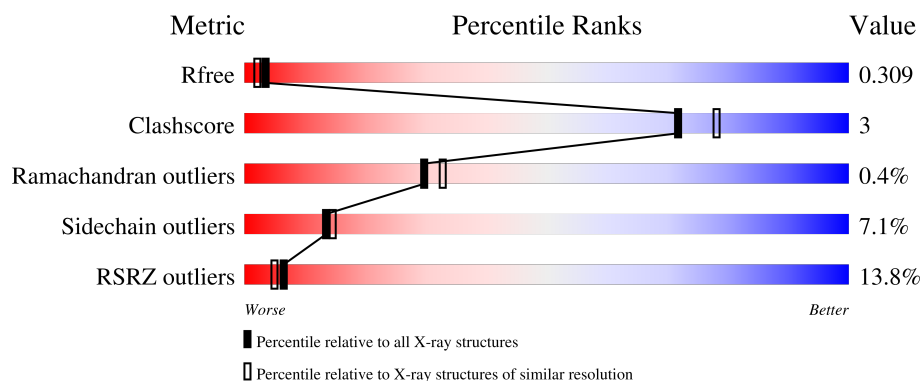
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	526	11% 72% 13% 15%
1	B	526	12% 68% 17% 15%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7517 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.

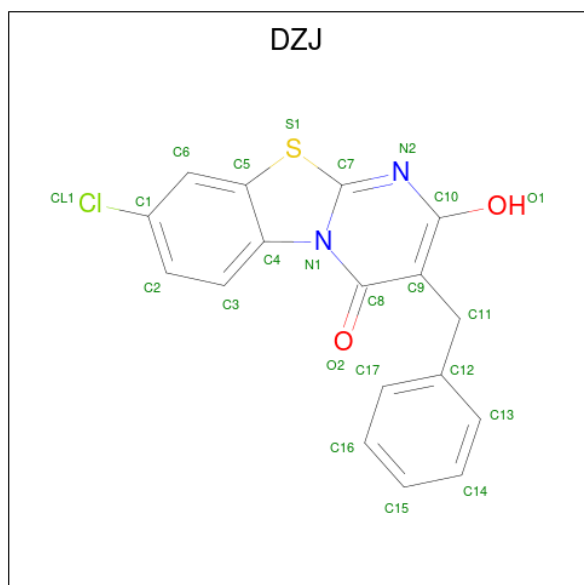
- Molecule 1 is a protein called Tyrosine-protein phosphatase non-receptor type 11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3631	2295	640	678	18			
1	B	448	Total	C	N	O	S	0	0	0
			3649	2307	650	674	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q06124
A	285	SER	PHE	engineered mutation	UNP Q06124
B	0	SER	-	expression tag	UNP Q06124
B	285	SER	PHE	engineered mutation	UNP Q06124

- Molecule 2 is 3-benzyl-8-chloro-2-hydroxy-4H-pyrimido[2,1-b][1,3]benzothiazol-4-one (CCD ID: DZJ) (formula: C₁₇H₁₁ClN₂O₂S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 23	C 17	Cl 1	N 2	O 2	S 1	0	0
2	B	1	Total 23	C 17	Cl 1	N 2	O 2	S 1	0	0

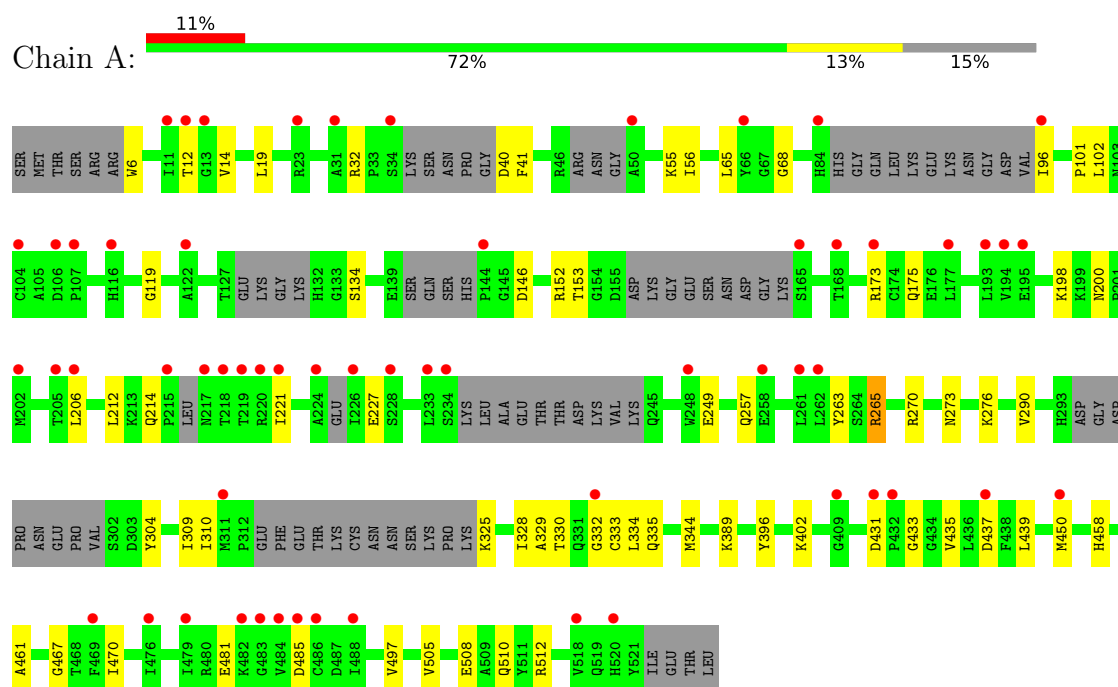
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	104	Total 104	O 104	0	0
3	B	87	Total 87	O 87	0	0

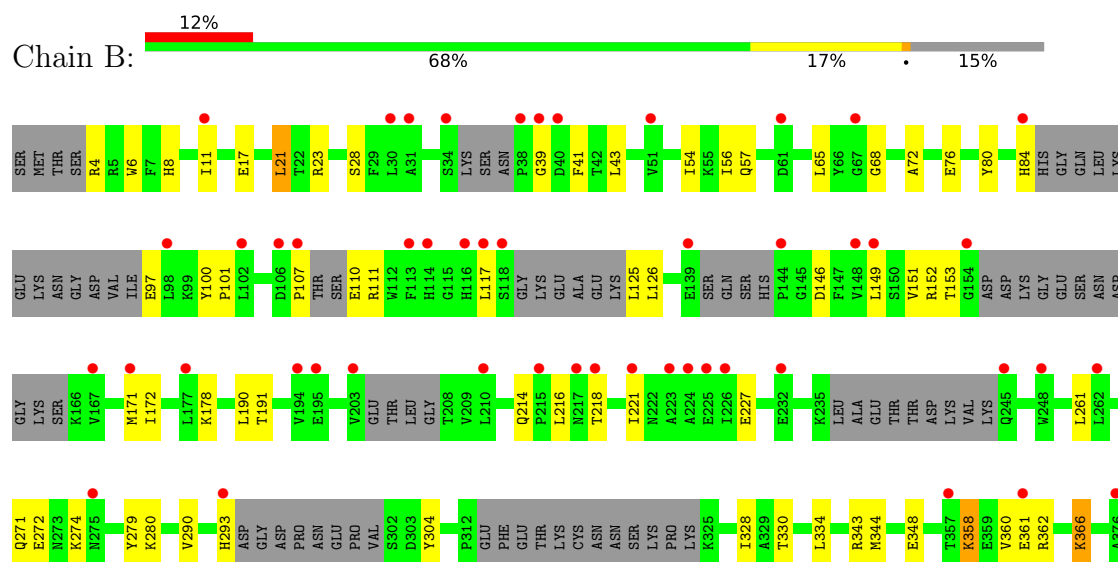
3 Residue-property plots [i](#)

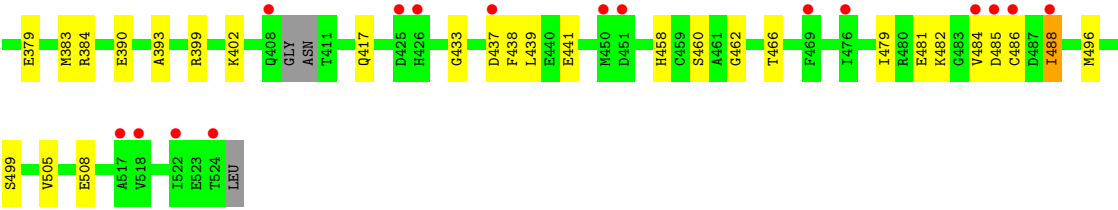
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Tyrosine-protein phosphatase non-receptor type 11



- Molecule 1: Tyrosine-protein phosphatase non-receptor type 11





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.19Å 212.23Å 55.47Å 90.00° 96.23° 90.00°	Depositor
Resolution (Å)	22.38 – 2.22 22.38 – 2.22	Depositor EDS
% Data completeness (in resolution range)	92.0 (22.38-2.22) 91.9 (22.38-2.22)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.21Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, R_{free}	0.236 , 0.292 0.250 , 0.309	Depositor DCC
R_{free} test set	2413 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.575	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.8	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	0.91	EDS
Total number of atoms	7517	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.42% 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: DZJ

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	0.88	0/3701	1.35	18/4987 (0.4%)
1	B	0.88	0/3720	1.34	22/5008 (0.4%)
All	All	0.88	0/7421	1.35	40/9995 (0.4%)

There are no bond length outliers.

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	437	ASP	CA-CB-CG	7.29	119.89	112.60
1	B	437	ASP	CA-CB-CG	6.28	118.88	112.60
1	A	334	LEU	N-CA-C	-6.17	101.07	110.14
1	B	39	GLY	N-CA-C	-6.08	107.02	114.92
1	B	304	TYR	N-CA-C	5.97	118.97	109.24
1	B	393	ALA	N-CA-C	-5.96	101.10	109.69
1	B	358	LYS	N-CA-C	-5.95	102.52	110.55
1	B	334	LEU	N-CA-C	-5.83	101.11	110.14
1	B	433	GLY	CA-C-N	5.80	126.54	119.99
1	B	433	GLY	C-N-CA	5.80	126.54	119.99
1	A	433	GLY	CA-C-N	5.71	126.31	119.98
1	A	433	GLY	C-N-CA	5.71	126.31	119.98
1	A	257	GLN	CA-C-N	5.67	127.81	120.44
1	A	257	GLN	C-N-CA	5.67	127.81	120.44
1	B	438	PHE	CA-CB-CG	5.54	119.34	113.80
1	B	279	TYR	CA-C-N	5.43	127.82	120.38
1	B	279	TYR	C-N-CA	5.43	127.82	120.38
1	B	437	ASP	CA-C-N	5.38	128.26	120.79
1	B	437	ASP	C-N-CA	5.38	128.26	120.79
1	A	332	GLY	CA-C-N	5.27	128.34	120.90
1	A	332	GLY	C-N-CA	5.27	128.34	120.90
1	A	119	GLY	CA-C-N	5.26	127.27	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	GLY	C-N-CA	5.26	127.27	120.44
1	A	265	ARG	CA-C-N	5.25	128.09	120.79
1	A	265	ARG	C-N-CA	5.25	128.09	120.79
1	B	481	GLU	CA-C-N	5.25	129.08	120.63
1	B	481	GLU	C-N-CA	5.25	129.08	120.63
1	B	146	ASP	CA-CB-CG	5.20	117.80	112.60
1	A	396	TYR	N-CA-C	5.17	116.67	108.96
1	A	304	TYR	N-CA-C	5.17	118.24	109.76
1	B	466	THR	CA-C-N	5.16	125.67	119.94
1	B	466	THR	C-N-CA	5.16	125.67	119.94
1	B	110	GLU	CA-C-N	5.15	131.38	121.54
1	B	110	GLU	C-N-CA	5.15	131.38	121.54
1	A	510	GLN	CA-C-N	5.15	127.49	120.54
1	A	510	GLN	C-N-CA	5.15	127.49	120.54
1	A	270	ARG	CA-C-N	5.09	127.41	120.54
1	A	270	ARG	C-N-CA	5.09	127.41	120.54
1	B	462	GLY	CA-C-N	-5.06	117.46	122.77
1	B	462	GLY	C-N-CA	-5.06	117.46	122.77

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	3631	0	3558	19	0
1	B	3649	0	3595	28	0
2	A	23	0	0	0	0
2	B	23	0	0	0	0
3	A	104	0	0	3	0
3	B	87	0	0	0	0
All	All	7517	0	7153	47	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 3.

All (47) 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:310:ILE:HD11	1:A:329:ALA:HB2	1.60	0.81
1:B:290:VAL:HG11	1:B:344:MET:HG3	1.70	0.73
1:B:390:GLU:HG2	1:B:399:ARG:HG2	1.71	0.72
1:B:65:LEU:HD23	1:B:68:GLY:HA3	1.78	0.65
1:A:310:ILE:HD13	1:A:470:ILE:HD13	1.77	0.65
1:A:330:THR:HG23	1:A:458:HIS:HB3	1.79	0.64
1:B:379:GLU:HG2	1:B:384:ARG:HG3	1.82	0.59
1:B:330:THR:HG23	1:B:458:HIS:HB3	1.85	0.59
1:A:497:VAL:HG11	3:A:757:HOH:O	2.03	0.58
1:B:271:GLN:HA	1:B:274:LYS:HD2	1.87	0.57
1:B:261:LEU:HD22	1:B:499:SER:HB3	1.87	0.56
1:B:17:GLU:O	1:B:21:LEU:HB2	2.07	0.55
1:A:41:PHE:HB2	1:A:56:ILE:HB	1.90	0.54
1:A:330:THR:CG2	1:A:458:HIS:HB3	2.38	0.53
1:A:198:LYS:HA	1:A:212:LEU:HB2	1.91	0.52
1:B:6:TRP:HB3	1:B:101:PRO:HB3	1.90	0.52
1:A:310:ILE:HD11	1:A:329:ALA:CB	2.35	0.52
1:B:107:PRO:HG3	1:B:190:LEU:HD12	1.93	0.51
1:B:328:ILE:HD11	1:B:348:GLU:HG3	1.91	0.51
1:A:309:ILE:HD13	1:A:328:ILE:HG12	1.92	0.50
1:A:467:GLY:HA2	3:A:757:HOH:O	2.10	0.50
1:A:461:ALA:HB3	3:A:735:HOH:O	2.12	0.49
1:B:482:LYS:HB3	1:B:486:CYS:HB3	1.95	0.48
1:A:134:SER:HA	1:A:214:GLN:O	2.13	0.48
1:A:65:LEU:HD23	1:A:68:GLY:HA3	1.94	0.47
1:B:149:LEU:HB2	1:B:172:ILE:HD11	1.95	0.47
1:A:6:TRP:HB3	1:A:101:PRO:HB3	1.96	0.47
1:A:263:TYR:HE2	1:A:265:ARG:HD2	1.81	0.46
1:B:361:GLU:HB2	1:B:366:LYS:HD3	1.97	0.46
1:A:290:VAL:HG11	1:A:344:MET:HG3	1.98	0.46
1:B:482:LYS:HD3	1:B:486:CYS:HB2	1.98	0.45
1:B:125:LEU:HD22	1:B:216:LEU:HD21	1.98	0.45
1:B:80:TYR:CE2	1:B:84:HIS:HE1	2.35	0.45
1:B:41:PHE:HB2	1:B:56:ILE:HB	1.99	0.45
1:B:126:LEU:HD23	1:B:216:LEU:HD13	1.98	0.44
1:B:28:SER:HA	1:B:100:TYR:O	2.18	0.44
1:B:8:HIS:HB3	1:B:11:ILE:HG12	2.00	0.43
1:B:43:LEU:HB3	1:B:54:ILE:HB	2.01	0.42
1:B:496:MET:HE3	1:B:496:MET:HB3	1.99	0.42
1:B:399:ARG:HD2	1:B:417:GLN:OE1	2.20	0.42
1:A:508:GLU:CD	1:A:512:ARG:HH12	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LEU:HD22	1:A:102:LEU:HD21	2.02	0.41
1:B:366:LYS:HG3	1:B:460:SER:HB3	2.02	0.41
1:B:343:ARG:HA	1:B:383:MET:HE3	2.02	0.41
1:B:479:ILE:HD11	1:B:488:ILE:HD13	2.02	0.41
1:B:72:ALA:HB3	1:B:76:GLU:OE1	2.21	0.41
1:A:273:ASN:HA	1:A:276:LYS:HD2	2.02	0.41

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	424/526 (81%)	402 (95%)	21 (5%)	1 (0%)	43	50
1	B	424/526 (81%)	404 (95%)	18 (4%)	2 (0%)	24	26
All	All	848/1052 (81%)	806 (95%)	39 (5%)	3 (0%)	30	33

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	111	ARG
1	A	505	VAL
1	B	505	VAL

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	398/468 (85%)	371 (93%)	27 (7%)	14	16
1	B	400/468 (86%)	370 (92%)	30 (8%)	12	13
All	All	798/936 (85%)	741 (93%)	57 (7%)	13	14

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	THR
1	A	14	VAL
1	A	32	ARG
1	A	40	ASP
1	A	55	LYS
1	A	96	ILE
1	A	146	ASP
1	A	152	ARG
1	A	153	THR
1	A	173	ARG
1	A	175	GLN
1	A	200	ASN
1	A	206	LEU
1	A	221	ILE
1	A	227	GLU
1	A	249	GLU
1	A	325	LYS
1	A	333	CYS
1	A	335	GLN
1	A	389	LYS
1	A	402	LYS
1	A	431	ASP
1	A	435	VAL
1	A	439	LEU
1	A	450	MET
1	A	481	GLU
1	A	485	ASP
1	B	4	ARG
1	B	21	LEU
1	B	23	ARG
1	B	57	GLN
1	B	97	GLU
1	B	117	LEU
1	B	151	VAL
1	B	152	ARG

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Mol	Chain	Res	Type
1	B	153	THR
1	B	171	MET
1	B	178	LYS
1	B	191	THR
1	B	214	GLN
1	B	218	THR
1	B	221	ILE
1	B	227	GLU
1	B	272	GLU
1	B	280	LYS
1	B	293	HIS
1	B	358	LYS
1	B	360	VAL
1	B	362	ARG
1	B	366	LYS
1	B	402	LYS
1	B	439	LEU
1	B	441	GLU
1	B	484	VAL
1	B	485	ASP
1	B	488	ILE
1	B	508	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	HIS
1	A	175	GLN
1	A	200	ASN
1	A	256	GLN
1	A	520	HIS
1	B	48	ASN
1	B	84	HIS
1	B	200	ASN
1	B	217	ASN
1	B	281	ASN
1	B	408	GLN

5.3.3 RNA ⓘ

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	DZJ	A	601	-	25,26,26	1.86	13 (52%)	36,38,38	1.77	9 (25%)
2	DZJ	B	601	-	25,26,26	1.84	12 (48%)	36,38,38	1.77	10 (27%)

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	DZJ	A	601	-	-	0/4/4/4	0/4/4/4
2	DZJ	B	601	-	-	0/4/4/4	0/4/4/4

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	DZJ	C4-N1	2.86	1.47	1.41
2	B	601	DZJ	C3-C2	2.82	1.43	1.38
2	A	601	DZJ	C3-C4	2.69	1.43	1.39
2	A	601	DZJ	C4-N1	2.68	1.46	1.41
2	A	601	DZJ	C3-C2	2.67	1.43	1.38
2	B	601	DZJ	C17-C12	2.60	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	DZJ	C17-C12	2.42	1.43	1.38
2	A	601	DZJ	C2-C1	2.39	1.42	1.38
2	B	601	DZJ	C3-C4	2.38	1.43	1.39
2	A	601	DZJ	C8-C9	2.33	1.50	1.44
2	B	601	DZJ	C8-C9	2.30	1.49	1.44
2	A	601	DZJ	C16-C17	2.22	1.42	1.38
2	B	601	DZJ	C2-C1	2.19	1.42	1.38
2	A	601	DZJ	C7-N1	2.17	1.43	1.38
2	B	601	DZJ	C7-S1	2.16	1.80	1.75
2	B	601	DZJ	C13-C12	2.15	1.43	1.38
2	B	601	DZJ	C7-N1	2.11	1.43	1.38
2	B	601	DZJ	C6-C1	2.10	1.41	1.38
2	A	601	DZJ	C11-C9	2.10	1.54	1.51
2	A	601	DZJ	C13-C12	2.10	1.43	1.38
2	B	601	DZJ	C11-C9	2.09	1.54	1.51
2	A	601	DZJ	C16-C15	2.09	1.42	1.38
2	B	601	DZJ	C14-C13	2.09	1.42	1.38
2	A	601	DZJ	C6-C1	2.07	1.41	1.38
2	A	601	DZJ	C14-C13	2.07	1.42	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	DZJ	O1-C10-N2	-4.98	108.89	114.83
2	A	601	DZJ	O1-C10-C9	3.73	127.87	120.69
2	B	601	DZJ	O1-C10-N2	-3.57	110.58	114.83
2	B	601	DZJ	S1-C7-N2	3.54	127.04	121.08
2	A	601	DZJ	S1-C7-N2	3.25	126.55	121.08
2	B	601	DZJ	C3-C4-C5	-3.20	115.90	120.06
2	B	601	DZJ	C6-C5-C4	3.20	124.54	120.83
2	B	601	DZJ	C5-S1-C7	2.90	93.36	89.37
2	A	601	DZJ	C6-C5-C4	2.90	124.19	120.83
2	B	601	DZJ	S1-C7-N1	-2.82	108.68	111.99
2	B	601	DZJ	O1-C10-C9	2.49	125.48	120.69
2	B	601	DZJ	C11-C9-C8	2.45	121.41	116.95
2	A	601	DZJ	C5-S1-C7	2.43	92.71	89.37
2	A	601	DZJ	C3-C4-C5	-2.41	116.92	120.06
2	A	601	DZJ	S1-C7-N1	-2.35	109.23	111.99
2	A	601	DZJ	C7-N2-C10	2.26	118.69	114.14
2	A	601	DZJ	N1-C7-N2	-2.17	124.21	126.62
2	B	601	DZJ	C2-C1-C6	-2.14	118.74	121.53
2	B	601	DZJ	N1-C7-N2	-2.10	124.28	126.62

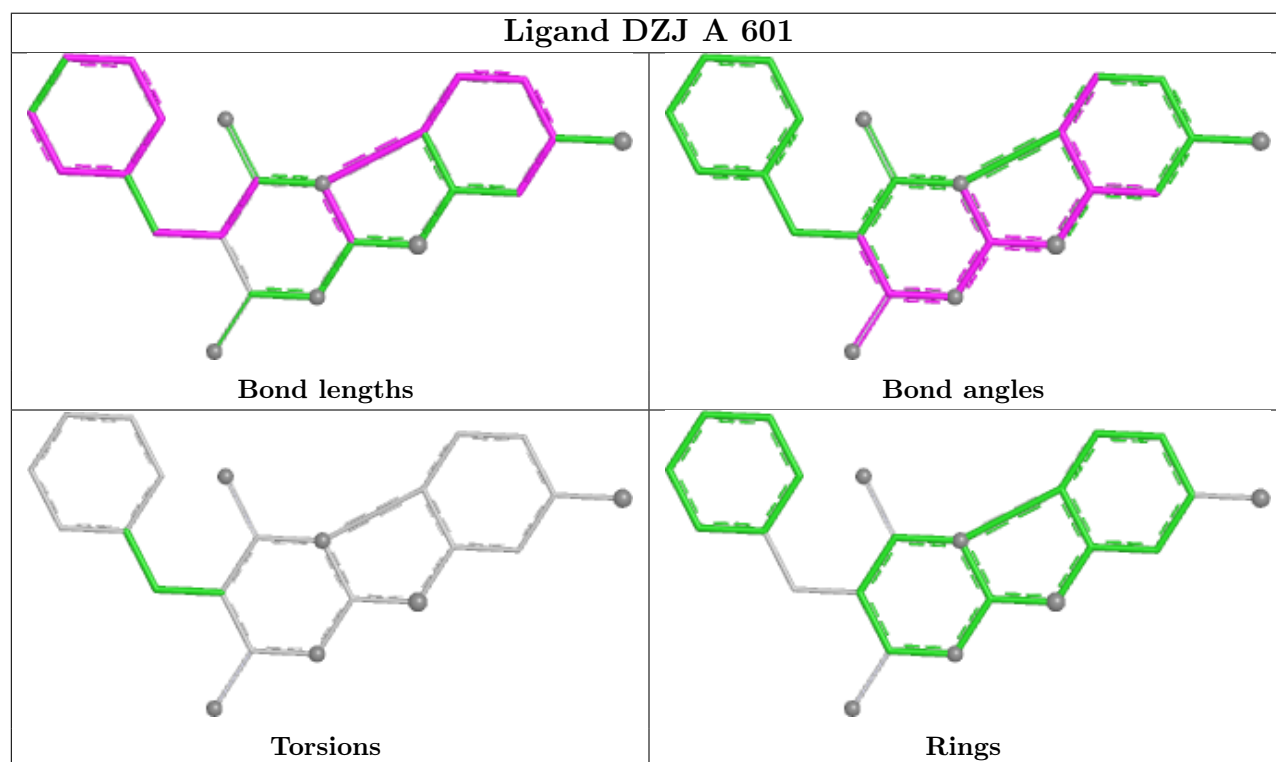
There are no chirality outliers.

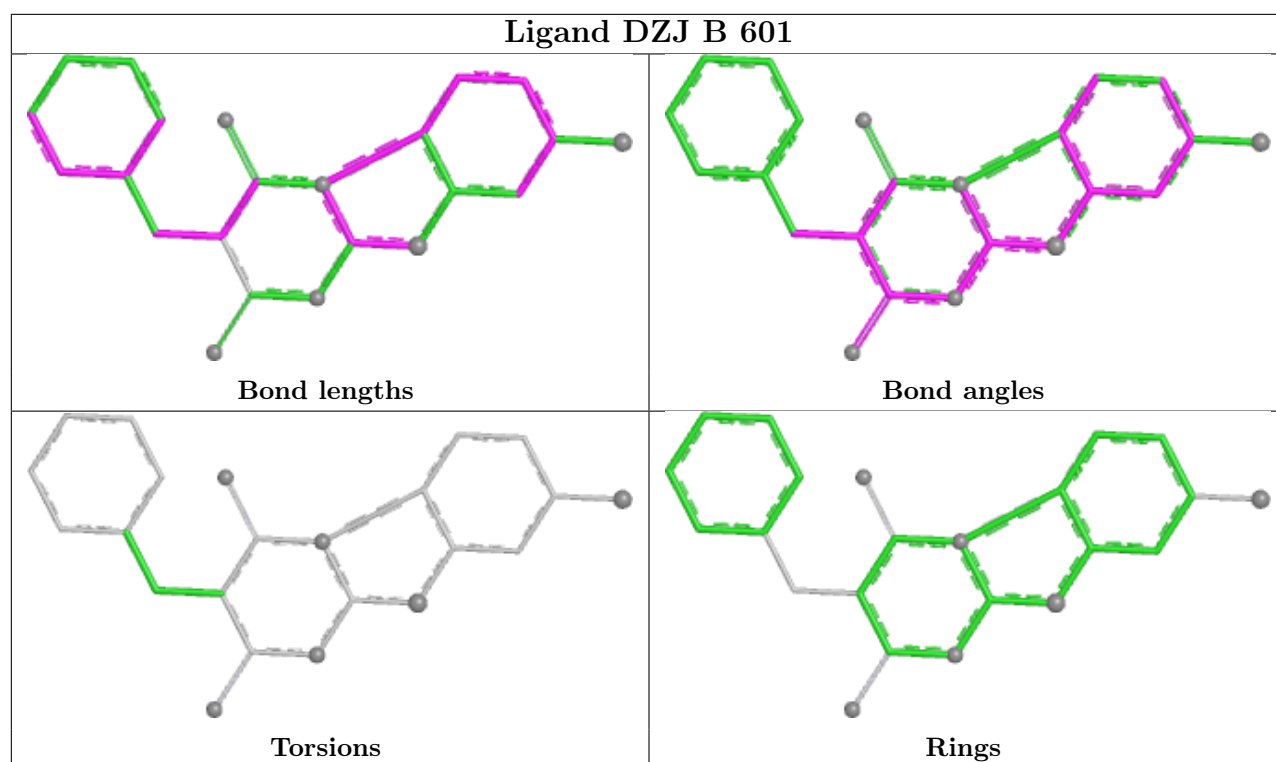
There are no torsion outliers.

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	448/526 (85%)	1.03	59 (13%) 7 5	20, 44, 68, 87	0
1	B	448/526 (85%)	1.07	65 (14%) 6 4	20, 45, 68, 88	0
All	All	896/1052 (85%)	1.05	124 (13%) 6 5	20, 44, 68, 88	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	476	ILE	4.5
1	A	215	PRO	4.1
1	B	177	LEU	4.0
1	A	484	VAL	3.8
1	A	50	ALA	3.8
1	B	522	ILE	3.6
1	B	485	ASP	3.6
1	B	426	HIS	3.4
1	A	116	HIS	3.4
1	B	225	GLU	3.4
1	A	409	GLY	3.4
1	A	165	SER	3.3
1	A	106	ASP	3.3
1	B	144	PRO	3.3
1	B	357	THR	3.3
1	A	224	ALA	3.2
1	B	116	HIS	3.2
1	A	13	GLY	3.2
1	B	484	VAL	3.2
1	A	206	LEU	3.2
1	A	233	LEU	3.1
1	B	524	THR	3.1
1	A	437	ASP	3.1
1	A	144	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	203	VAL	3.1
1	B	262	LEU	3.1
1	B	11	ILE	3.0
1	B	275	ASN	3.0
1	B	437	ASP	3.0
1	A	31	ALA	2.9
1	A	177	LEU	2.9
1	B	194	VAL	2.9
1	A	11	ILE	2.9
1	B	450	MET	2.9
1	A	226	ILE	2.9
1	A	228	SER	2.9
1	B	34	SER	2.9
1	B	118	SER	2.8
1	A	469	PHE	2.8
1	A	219	THR	2.8
1	B	218	THR	2.8
1	B	226	ILE	2.8
1	B	107	PRO	2.8
1	B	31	ALA	2.8
1	A	84	HIS	2.8
1	A	431	ASP	2.8
1	A	66	TYR	2.8
1	B	486	CYS	2.7
1	A	218	THR	2.7
1	B	408	GLN	2.7
1	A	332	GLY	2.7
1	A	107	PRO	2.7
1	A	485	ASP	2.6
1	B	39	GLY	2.6
1	B	117	LEU	2.6
1	A	12	THR	2.6
1	A	488	ILE	2.6
1	B	517	ALA	2.6
1	A	220	ARG	2.6
1	A	482	LYS	2.6
1	A	261	LEU	2.6
1	B	67	GLY	2.6
1	B	221	ILE	2.5
1	A	258	GLU	2.5
1	A	518	VAL	2.5
1	B	488	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	34	SER	2.4
1	A	122	ALA	2.4
1	B	195	GLU	2.4
1	A	205	THR	2.4
1	B	223	ALA	2.4
1	A	96	ILE	2.4
1	B	148	VAL	2.4
1	B	476	ILE	2.4
1	A	23	ARG	2.4
1	B	114	HIS	2.4
1	B	232	GLU	2.4
1	A	168	THR	2.3
1	A	311	MET	2.3
1	A	450	MET	2.3
1	A	520	HIS	2.3
1	B	38	PRO	2.3
1	B	376	ALA	2.3
1	B	139	GLU	2.3
1	B	245	GLN	2.3
1	A	221	ILE	2.3
1	A	479	ILE	2.3
1	A	248	TRP	2.2
1	B	248	TRP	2.2
1	B	171	MET	2.2
1	A	173	ARG	2.2
1	A	486	CYS	2.2
1	B	106	ASP	2.2
1	B	425	ASP	2.2
1	B	293	HIS	2.2
1	B	98	LEU	2.1
1	B	51	VAL	2.1
1	B	518	VAL	2.1
1	B	224	ALA	2.1
1	A	202	MET	2.1
1	B	167	VAL	2.1
1	B	113	PHE	2.1
1	B	149	LEU	2.1
1	B	84	HIS	2.1
1	A	194	VAL	2.1
1	B	215	PRO	2.1
1	B	469	PHE	2.1
1	A	104	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	234	SER	2.1
1	A	262	LEU	2.1
1	B	30	LEU	2.1
1	B	61	ASP	2.1
1	B	217	ASN	2.1
1	A	432	PRO	2.1
1	A	483	GLY	2.1
1	B	154	GLY	2.1
1	A	195	GLU	2.0
1	B	361	GLU	2.0
1	B	102	LEU	2.0
1	B	210	LEU	2.0
1	A	193	LEU	2.0
1	A	217	ASN	2.0
1	B	40	ASP	2.0
1	B	451	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

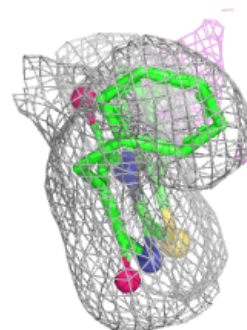
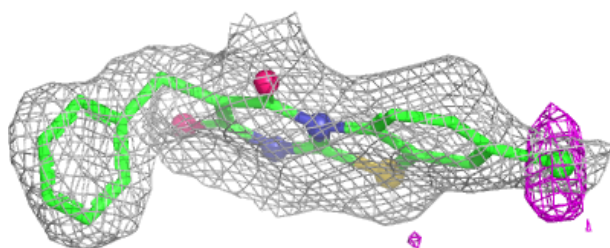
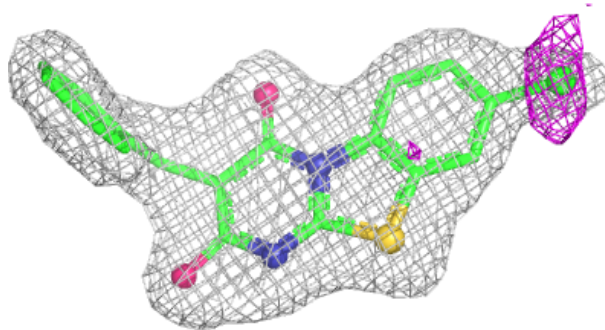
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	DZJ	B	601	23/23	0.83	0.12	41,43,46,51	0
2	DZJ	A	601	23/23	0.84	0.13	47,50,54,60	0

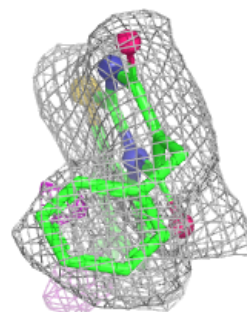
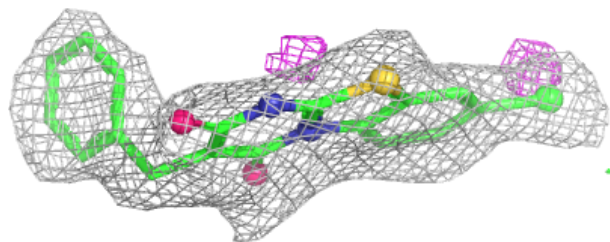
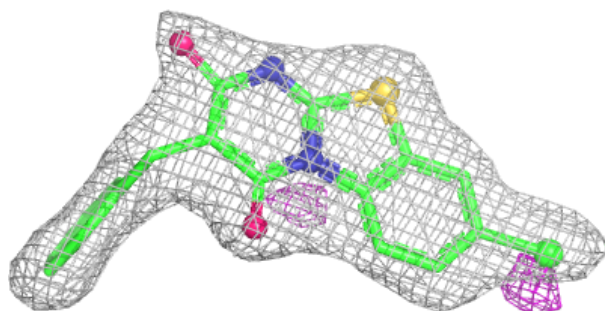
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DZJ B 601:

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

**Electron density around DZJ A 601:**

$2mF_o-DF_c$ (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.