Full wwPDB X-ray Structure Validation Report ⓘ

Jun 25, 2024 – 11:47 PM EDT

PDB ID	:	6RSE
Title	:	Structure based optimization of JAK1-ATP binding pocket Inhibitors in the aminopyrazole class
Authors	:	Brown, D.G.; Lupardus, P.J.
Deposited on	:	2019-05-21
Resolution	:	1.80 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:

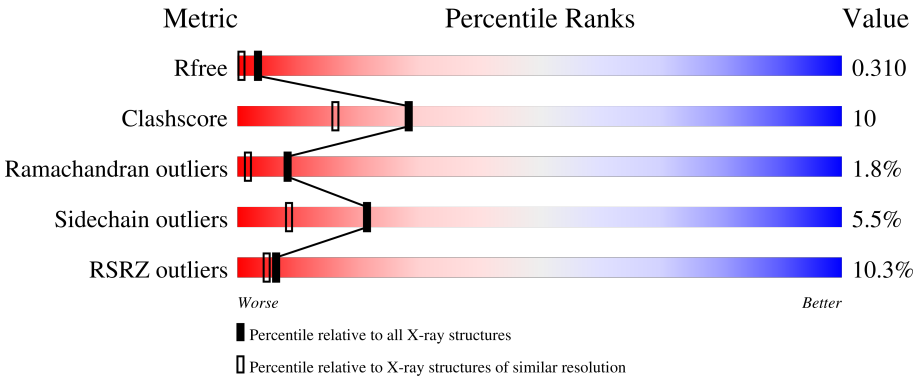
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	302	9% 76% 14% • • 5%
1	B	302	11% 72% 19% • • 7%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5000 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.

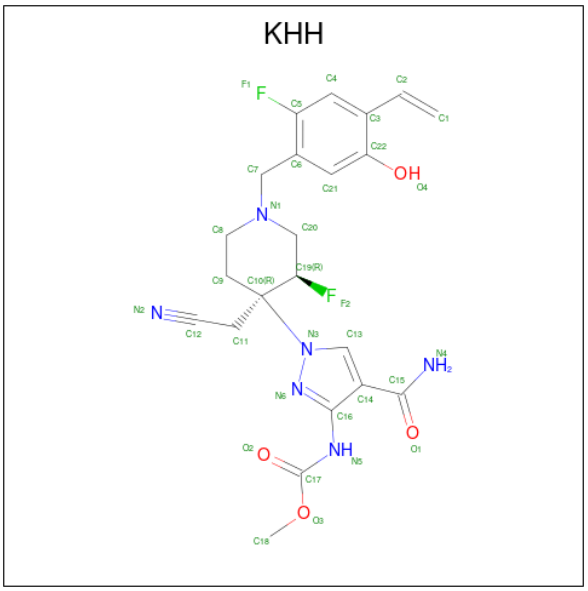
- Molecule 1 is a protein called Tyrosine-protein kinase JAK1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	286	Total	C	N	O	P	S	0	1	0
			2331	1482	395	437	2	15			
1	B	281	Total	C	N	O	P	S	0	0	0
			2291	1461	388	425	2	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	853	GLY	-	expression tag	UNP P23458
B	853	GLY	-	expression tag	UNP P23458

- Molecule 2 is methyl {N}-[4-aminocarbonyl-1-[(3 {R},4 {R})-4-(cyanomethyl)-1-[(4-ethenyl-2-fluoranyl-5-oxidanyl-phenyl)methyl]-3-fluoranyl-piperidin-4-yl]pyrazol-3-yl]carbamate (three-letter code: KHH) (formula: $C_{22}H_{24}F_2N_6O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			34	22	2	6	4		
2	B	1	Total	C	F	N	O	0	0
			34	22	2	6	4		

- Molecule 3 is water.

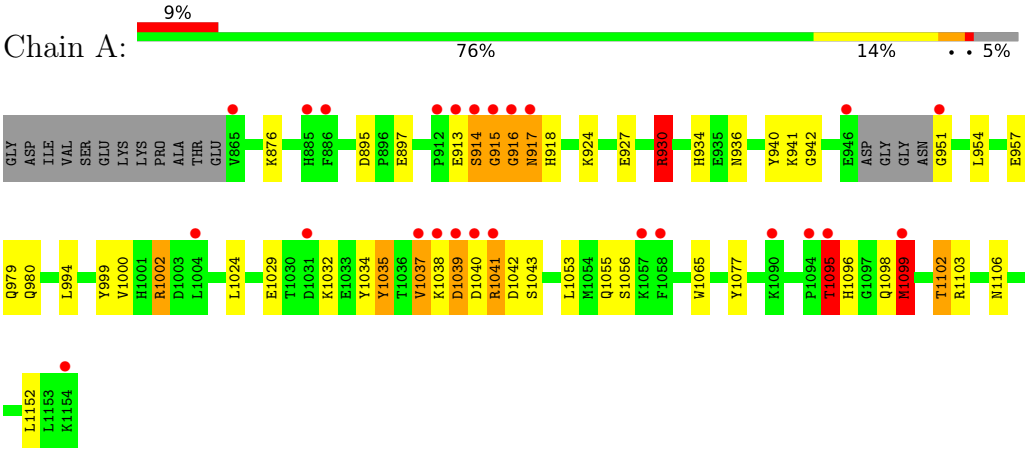
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	149	Total	O	0	0
			149	149		
3	B	161	Total	O	0	0
			161	161		

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.

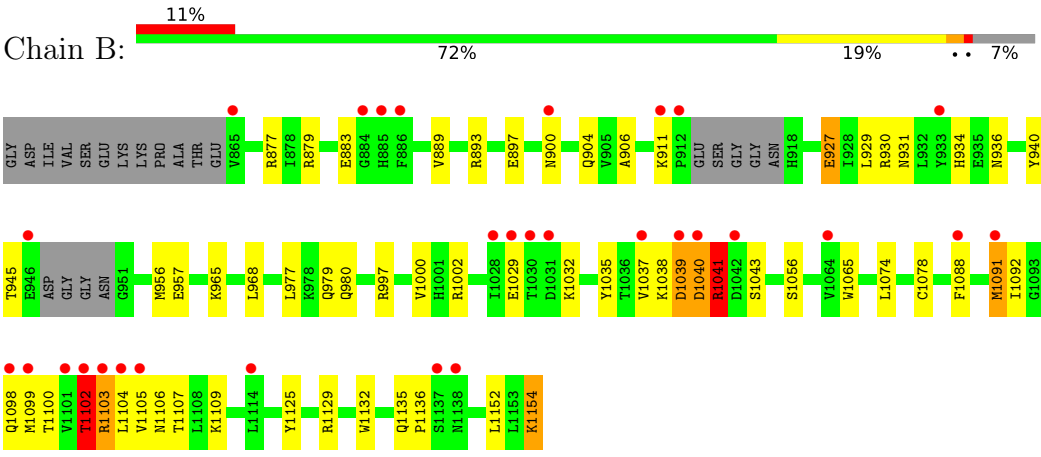
• Molecule 1: Tyrosine-protein kinase JAK1

Chain A:



• Molecule 1: Tyrosine-protein kinase JAK1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	42.89Å 88.96Å 173.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.04 – 1.80 39.04 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.3 (39.04-1.80) 99.3 (39.04-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.78 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.263 , 0.303 0.270 , 0.310	Depositor DCC
R_{free} test set	3063 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	25.1	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 43.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.55$, $< L^2 > = 0.39$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5000	wwPDB-VP
Average B, all atoms (Å^2)	35.0	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 63.42 % 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 9.6906e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: KHH, PTR

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	0.90	2/2347 (0.1%)	1.02	2/3159 (0.1%)
1	B	1.04	8/2306 (0.3%)	1.02	1/3103 (0.0%)
All	All	0.97	10/4653 (0.2%)	1.02	3/6262 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
All	All	0	6

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1102	THR	C-O	16.54	1.54	1.23
1	B	1104	LEU	C-O	8.77	1.40	1.23
1	B	1105	VAL	C-O	8.53	1.39	1.23
1	A	897	GLU	CD-OE2	-7.77	1.17	1.25
1	B	957	GLU	CD-OE2	-7.14	1.17	1.25
1	B	1109	LYS	C-O	6.83	1.36	1.23
1	B	897	GLU	CD-OE2	-6.55	1.18	1.25
1	B	1100	THR	C-O	6.47	1.35	1.23
1	B	956	MET	C-O	5.03	1.32	1.23
1	A	1099	MET	CG-SD	5.02	1.94	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	930	ARG	NE-CZ-NH1	-8.16	116.22	120.30
1	B	879	ARG	NE-CZ-NH1	6.17	123.39	120.30
1	A	1002	ARG	NE-CZ-NH1	5.98	123.29	120.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1041	ARG	Peptide
1	A	1102	THR	Mainchain
1	A	916	GLY	Peptide
1	B	1039	ASP	Peptide
1	B	1041	ARG	Peptide
1	B	1102	THR	Mainchain

5.2 Too-close contacts (i)

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	2331	0	2315	50	0
1	B	2291	0	2286	53	0
2	A	34	0	0	0	0
2	B	34	0	0	0	0
3	A	149	0	0	7	0
3	B	161	0	0	6	0
All	All	5000	0	4601	91	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 10.

All (91) 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:900:ASN:HB2	3:B:1446:HOH:O	1.67	0.93
1:A:1103:ARG:HE	1:B:1041:ARG:HH22	1.08	0.91
1:A:1041:ARG:HH11	1:B:1096:HIS:CD2	1.90	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:906:ALA:HB1	3:B:1434:HOH:O	1.75	0.85
1:A:927:GLU:OE1	1:A:930:ARG:NH1	2.10	0.84
1:B:977:LEU:HA	1:B:980:GLN:HE21	1.45	0.79
1:A:977:LEU:HA	1:A:980:GLN:HE21	1.47	0.79
1:A:1041:ARG:HH11	1:B:1096:HIS:CG	2.04	0.74
1:A:934:HIS:HD2	1:A:936:ASN:H	1.36	0.73
1:B:934:HIS:HD2	1:B:936:ASN:H	1.38	0.72
1:A:927:GLU:OE2	1:A:930:ARG:NH1	2.21	0.71
1:A:1095:THR:HB	1:B:1040:ASP:HB3	1.72	0.71
1:A:927:GLU:CD	1:A:930:ARG:NH1	2.44	0.71
1:A:927:GLU:OE2	3:A:1301:HOH:O	2.08	0.70
1:A:1032:LYS:NZ	1:A:1034:PTR:O2P	2.26	0.69
1:A:914:SER:HB2	1:A:918:HIS:HB2	1.75	0.68
1:A:1041:ARG:HH12	1:B:1103:ARG:HD2	1.57	0.68
1:A:1103:ARG:HE	1:B:1041:ARG:NH2	1.89	0.68
1:A:1041:ARG:NH1	1:B:1096:HIS:CD2	2.62	0.68
1:A:1103:ARG:NE	1:B:1041:ARG:HH22	1.89	0.66
1:A:927:GLU:CD	1:A:930:ARG:HH12	1.99	0.65
1:A:1037:VAL:O	1:A:1056:SER:HB2	1.97	0.65
1:B:934:HIS:CD2	1:B:936:ASN:H	2.15	0.64
1:A:1053:LEU:O	1:B:1102:THR:HG21	1.97	0.64
1:A:934:HIS:CD2	1:A:936:ASN:H	2.15	0.63
1:A:954:LEU:HD13	3:A:1379:HOH:O	1.99	0.62
1:B:965:LYS:HD2	1:B:1078:CYS:HB3	1.80	0.62
1:A:1041:ARG:HH12	1:B:1103:ARG:CD	2.13	0.61
1:A:1035:PTR:HE1	1:A:1037:VAL:CG2	2.31	0.60
1:A:914:SER:HA	1:A:918:HIS:HD2	1.66	0.60
1:B:1039:ASP:O	1:B:1041:ARG:HD3	2.02	0.60
1:A:1035:PTR:HE1	1:A:1037:VAL:HG22	1.85	0.57
1:B:1041:ARG:HG3	1:B:1043:SER:HB3	1.85	0.57
1:A:1098:GLN:C	1:A:1099:MET:HE2	2.24	0.56
1:B:904:GLN:NE2	3:B:1305:HOH:O	2.38	0.56
1:B:1029:GLU:OE2	1:B:1032:LYS:HE3	2.06	0.56
1:A:914:SER:CB	1:A:918:HIS:HB2	2.35	0.56
1:B:1041:ARG:HG3	1:B:1043:SER:CB	2.36	0.56
1:B:1065:TRP:CE3	1:B:1132:TRP:HA	2.42	0.54
1:A:940:TYR:CZ	1:A:942:GLY:HA2	2.45	0.52
1:B:1041:ARG:HG3	1:B:1043:SER:OG	2.10	0.51
1:B:1002:ARG:NH2	1:B:1037:VAL:HG21	2.25	0.51
1:A:915:GLY:O	1:A:917:ASN:N	2.43	0.51
1:B:943:ILE:HD12	1:B:945:THR:CG2	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:893:ARG:NE	1:B:900:ASN:HD22	2.10	0.49
1:B:1154:LYS:HA	1:B:1154:LYS:CE	2.43	0.49
1:B:1029:GLU:OE2	1:B:1032:LYS:CE	2.60	0.49
1:A:954:LEU:CD1	3:A:1379:HOH:O	2.59	0.48
1:B:889:VAL:HG11	3:B:1434:HOH:O	2.12	0.48
1:A:1103:ARG:HH11	1:B:1041:ARG:NH2	2.10	0.48
1:B:1040:ASP:OD1	1:B:1040:ASP:N	2.40	0.48
1:A:972:LYS:HE3	1:A:1077:TYR:CD2	2.49	0.47
1:B:1102:THR:O	1:B:1106:ASN:HB2	2.14	0.47
1:A:1102:THR:HG23	3:A:1311:HOH:O	2.13	0.47
1:B:979:GLN:NE2	3:B:1314:HOH:O	2.48	0.47
1:B:893:ARG:NE	1:B:900:ASN:ND2	2.62	0.46
1:B:965:LYS:HE2	3:B:1441:HOH:O	2.15	0.46
1:A:941:LYS:HD2	1:A:957:GLU:HA	1.98	0.46
1:B:1092:ILE:O	1:B:1096:HIS:HE1	1.99	0.46
1:A:1041:ARG:HG3	1:A:1043:SER:OG	2.15	0.45
1:A:1035:PTR:CE1	1:A:1037:VAL:HG22	2.45	0.45
1:A:951:GLY:C	3:A:1345:HOH:O	2.55	0.45
1:B:931:ASN:O	1:B:997:ARG:NH2	2.46	0.45
1:A:979:GLN:NE2	3:A:1310:HOH:O	2.49	0.45
1:B:1035:PTR:HE1	1:B:1037:VAL:HG22	1.98	0.45
1:A:1099:MET:CA	1:A:1099:MET:CE	2.95	0.45
1:A:1098:GLN:HE21	1:B:1041:ARG:HB3	1.81	0.44
1:B:1125:TYR:CZ	1:B:1129:ARG:HD3	2.53	0.44
1:B:930:ARG:HG3	1:B:940:TYR:CE2	2.53	0.44
1:B:1039:ASP:O	1:B:1041:ARG:CD	2.65	0.44
1:B:1088:PHE:HA	1:B:1091:MET:HE3	1.98	0.44
1:B:1041:ARG:HD3	1:B:1041:ARG:N	2.33	0.44
1:B:980:GLN:O	1:B:1074:LEU:HD21	2.18	0.43
1:B:927:GLU:OE1	1:B:927:GLU:HA	2.18	0.43
1:B:940:TYR:CZ	1:B:942:GLY:HA2	2.53	0.43
1:B:943:ILE:HD12	1:B:945:THR:HG23	2.00	0.43
1:A:994:LEU:HG	1:A:999:TYR:HB2	2.01	0.42
1:B:1065:TRP:CZ3	1:B:1132:TRP:HA	2.54	0.42
1:B:1092:ILE:O	1:B:1096:HIS:CE1	2.72	0.42
1:A:1039:ASP:CG	1:A:1040:ASP:H	2.22	0.42
1:A:1039:ASP:OD2	1:A:1040:ASP:N	2.44	0.42
1:A:1065:TRP:CE3	1:A:1132:TRP:HA	2.54	0.42
1:B:1135:GLN:HA	1:B:1136:PRO:HD2	1.93	0.42
1:A:940:TYR:HE2	3:A:1379:HOH:O	2.03	0.42
1:A:1002:ARG:NH1	1:A:1024:LEU:O	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1055:GLN:C	1:A:1056:SER:OG	2.58	0.41
1:A:1029:GLU:HB3	1:A:1032:LYS:HG3	2.02	0.41
1:A:876:LYS:HE2	1:A:895:ASP:HB3	2.02	0.41
1:A:1099:MET:HE2	1:A:1099:MET:N	2.35	0.41
1:B:965:LYS:CD	1:B:1078:CYS:HB3	2.48	0.41
1:B:968:LEU:HD11	1:B:1074:LEU: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	281/302 (93%)	263 (94%)	9 (3%)	9 (3%)	4 0
1	B	273/302 (90%)	263 (96%)	9 (3%)	1 (0%)	34 21
All	All	554/604 (92%)	526 (95%)	18 (3%)	10 (2%)	8 2

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	913	GLU
1	A	914	SER
1	A	916	GLY
1	A	1039	ASP
1	A	917	ASN
1	B	1095	THR
1	A	1095	THR
1	A	1096	HIS
1	A	1042	ASP
1	A	915	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	256/267 (96%)	246 (96%)	10 (4%)	32 17
1	B	252/267 (94%)	234 (93%)	18 (7%)	14 5
All	All	508/534 (95%)	480 (94%)	28 (6%)	21 8

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	924	LYS
1	A	930	ARG
1	A	1000	VAL
1	A	1037	VAL
1	A	1038	LYS
1	A	1095	THR
1	A	1099	MET
1	A	1106	ASN
1	A	1112	LYS
1	A	1152	LEU
1	B	877	ARG
1	B	883	GLU
1	B	911	LYS
1	B	927	GLU
1	B	929	LEU
1	B	1000	VAL
1	B	1038	LYS
1	B	1040	ASP
1	B	1041	ARG
1	B	1056	SER
1	B	1091	MET
1	B	1095	THR
1	B	1098	GLN
1	B	1099	MET
1	B	1103	ARG
1	B	1107	THR
1	B	1152	LEU

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Mol	Chain	Res	Type
1	B	1154	LYS

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	918	HIS
1	A	934	HIS
1	A	936	ASN
1	A	979	GLN
1	A	980	GLN
1	A	1098	GLN
1	A	1135	GLN
1	B	900	ASN
1	B	934	HIS
1	B	936	ASN
1	B	979	GLN
1	B	980	GLN
1	B	1096	HIS
1	B	1106	ASN
1	B	1138	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

4 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	PTR	A	1035	1	15,16,17	0.53	0	19,22,24	1.23	3 (15%)
1	PTR	A	1034	1	15,16,17	0.60	0	19,22,24	0.76	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
1	PTR	B	1035	1	15,16,17	0.72	0	19,22,24	0.71	0
1	PTR	B	1034	1	15,16,17	0.62	0	19,22,24	0.68	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PTR	A	1035	1	-	2/10/11/13	0/1/1/1
1	PTR	A	1034	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1035	1	-	0/10/11/13	0/1/1/1
1	PTR	B	1034	1	-	0/10/11/13	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	1035	PTR	O2P-P-OH	2.85	114.14	105.24
1	A	1035	PTR	OH-CZ-CE2	2.18	125.71	119.23
1	A	1035	PTR	OH-CZ-CE1	-2.10	112.97	119.23

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1035	PTR	CE1-CZ-OH-P
1	A	1035	PTR	CE2-CZ-OH-P

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1035	PTR	3	0
1	A	1034	PTR	1	0
1	B	1035	PTR	1	0

5.5 Carbohydrates (i)

There are no monosaccharides 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	KHH	B	1201	-	31,36,36	0.72	0	35,52,52	2.56	9 (25%)
2	KHH	A	1201	-	31,36,36	0.67	1 (3%)	35,52,52	2.66	7 (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	KHH	B	1201	-	-	4/16/42/42	0/3/3/3
2	KHH	A	1201	-	-	6/16/42/42	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	KHH	C10-C19	-2.26	1.50	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1201	KHH	C13-N3-C10	10.64	132.27	124.18
2	A	1201	KHH	C13-N3-C10	10.55	132.20	124.18
2	A	1201	KHH	C8-N1-C20	6.62	119.64	109.52
2	B	1201	KHH	C8-N1-C20	4.98	117.13	109.52
2	A	1201	KHH	C7-N1-C8	4.42	120.91	111.06
2	B	1201	KHH	C13-N3-N6	-4.16	109.19	111.69
2	A	1201	KHH	C7-N1-C20	4.15	118.09	111.36
2	A	1201	KHH	C6-C7-N1	3.76	119.33	112.75
2	B	1201	KHH	C7-N1-C8	3.52	118.89	111.06
2	A	1201	KHH	C13-N3-N6	-3.34	109.69	111.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1201	KHH	C9-C10-N3	-3.29	105.74	109.35
2	B	1201	KHH	C7-N1-C20	3.26	116.64	111.36
2	B	1201	KHH	C19-C10-N3	2.93	113.25	109.06
2	A	1201	KHH	C9-C8-N1	2.91	114.10	111.23
2	B	1201	KHH	C6-C7-N1	2.18	116.56	112.75
2	B	1201	KHH	C9-C8-N1	2.03	113.24	111.23

There are no chirality outliers.

All (10) torsion outliers are listed below:

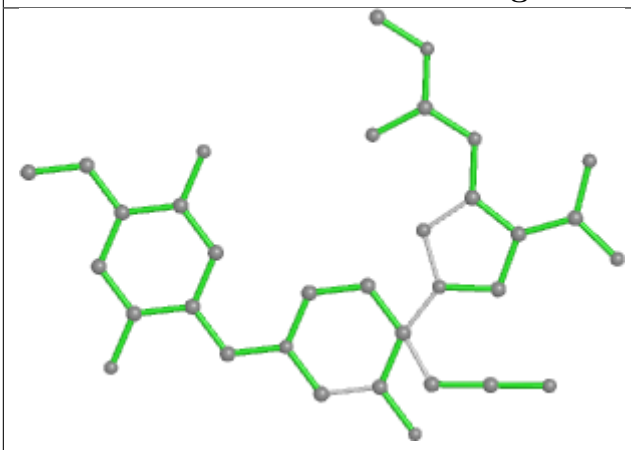
Mol	Chain	Res	Type	Atoms
2	A	1201	KHH	C9-C10-N3-N6
2	A	1201	KHH	C11-C10-N3-N6
2	B	1201	KHH	C9-C10-N3-N6
2	B	1201	KHH	C11-C10-N3-C13
2	B	1201	KHH	C11-C10-N3-N6
2	B	1201	KHH	C6-C7-N1-C20
2	A	1201	KHH	C6-C7-N1-C8
2	A	1201	KHH	C9-C10-N3-C13
2	A	1201	KHH	C6-C7-N1-C20
2	A	1201	KHH	C11-C10-N3-C13

There are no ring outliers.

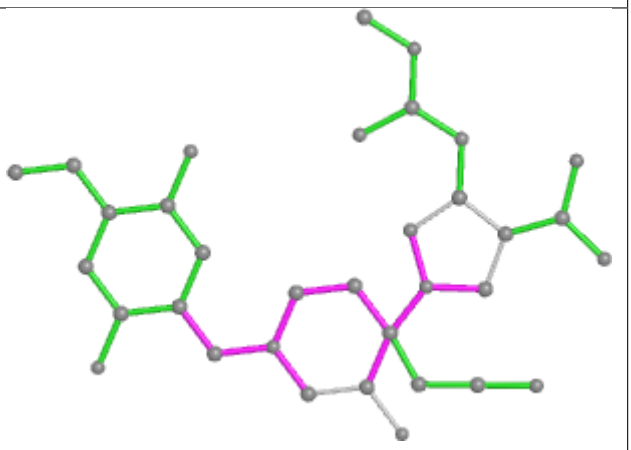
No monomer is involved in short contacts.

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.

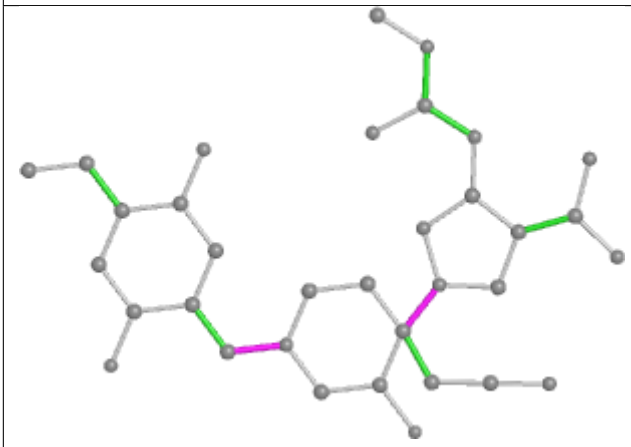
Ligand KHH B 1201



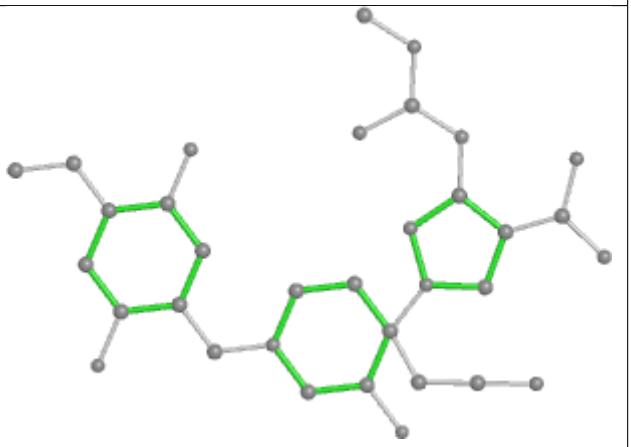
Bond lengths



Bond angles

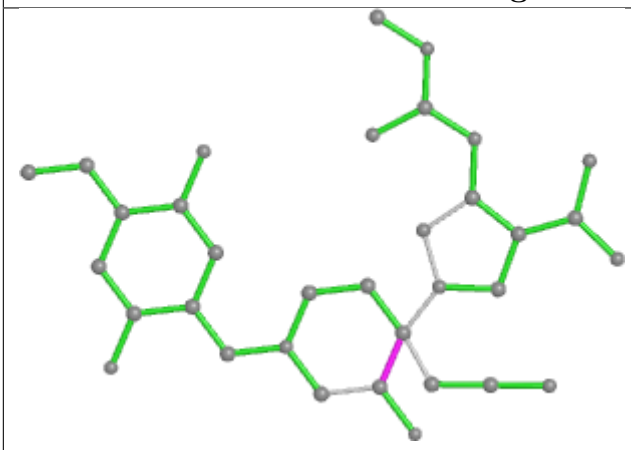


Torsions

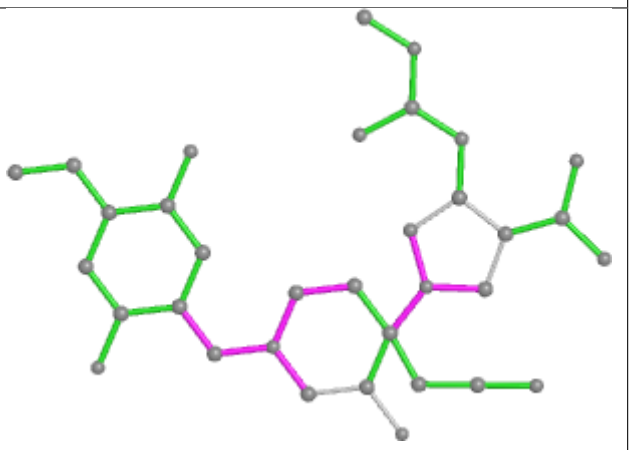


Rings

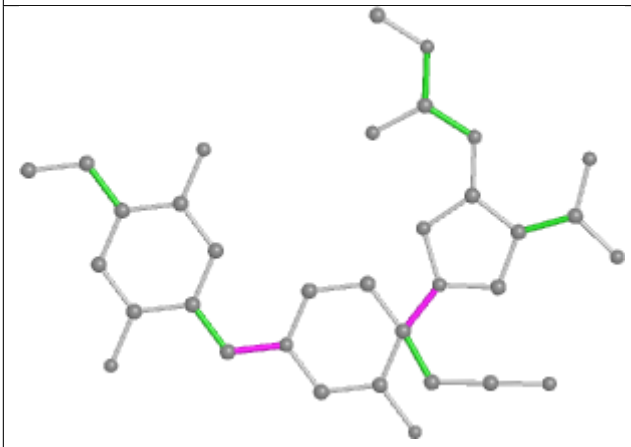
Ligand KHH A 1201



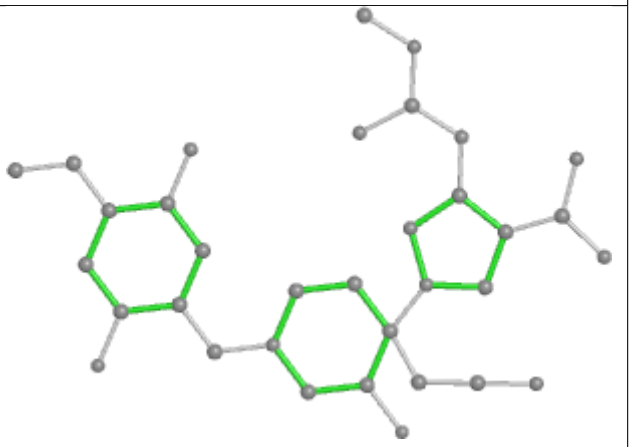
Bond lengths



Bond angles



Torsions



Rings

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, 95^{th} 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	284/302 (94%)	0.83	26 (9%)	9	6	16, 31, 62, 98	0
1	B	279/302 (92%)	0.82	32 (11%)	4	3	16, 32, 65, 83	0
All	All	563/604 (93%)	0.82	58 (10%)	6	5	16, 32, 64, 98	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	916	GLY	9.8
1	A	915	GLY	7.5
1	A	1095	THR	7.1
1	A	914	SER	7.1
1	A	1037	VAL	6.8
1	B	1040	ASP	6.6
1	A	885	HIS	5.8
1	A	917	ASN	5.4
1	B	885	HIS	5.3
1	A	1040	ASP	5.3
1	A	913	GLU	5.3
1	B	1031	ASP	4.9
1	A	1039	ASP	4.9
1	A	865	VAL	4.7
1	B	946	GLU	4.1
1	B	1030	THR	4.1
1	A	946	GLU	4.0
1	A	1038	LYS	3.7
1	B	1039	ASP	3.7
1	A	912	PRO	3.5
1	B	1037	VAL	3.5
1	B	1029	GLU	3.4
1	A	1090	LYS	3.3
1	B	886	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	1099	MET	3.1
1	A	1094	PRO	3.0
1	B	912	PRO	3.0
1	B	1094	PRO	2.9
1	B	1138	ASN	2.8
1	A	951	GLY	2.8
1	A	886	PHE	2.7
1	B	911	LYS	2.7
1	B	865	VAL	2.6
1	A	1154	LYS	2.6
1	B	1103	ARG	2.6
1	B	900	ASN	2.6
1	A	1041	ARG	2.5
1	B	1102	THR	2.5
1	B	1101	VAL	2.4
1	B	933	TYR	2.4
1	B	1088	PHE	2.4
1	A	1031	ASP	2.4
1	B	1098	GLN	2.4
1	B	1105	VAL	2.4
1	A	1004	LEU	2.3
1	A	1057	LYS	2.2
1	A	1099	MET	2.2
1	B	1137	SER	2.2
1	B	884	GLY	2.2
1	A	1137	SER	2.2
1	B	1114	LEU	2.2
1	B	1104	LEU	2.1
1	B	1042	ASP	2.1
1	B	1064	VAL	2.1
1	B	1028	ILE	2.1
1	B	1095	THR	2.1
1	A	1058	PHE	2.0
1	B	1091	MET	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95^{th} 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
1	PTR	B	1034	16/17	0.73	0.24	52,62,73,75	0
1	PTR	A	1035	16/17	0.77	0.18	39,48,63,64	0
1	PTR	B	1035	16/17	0.77	0.18	46,54,67,67	0
1	PTR	A	1034	16/17	0.78	0.23	45,54,67,67	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

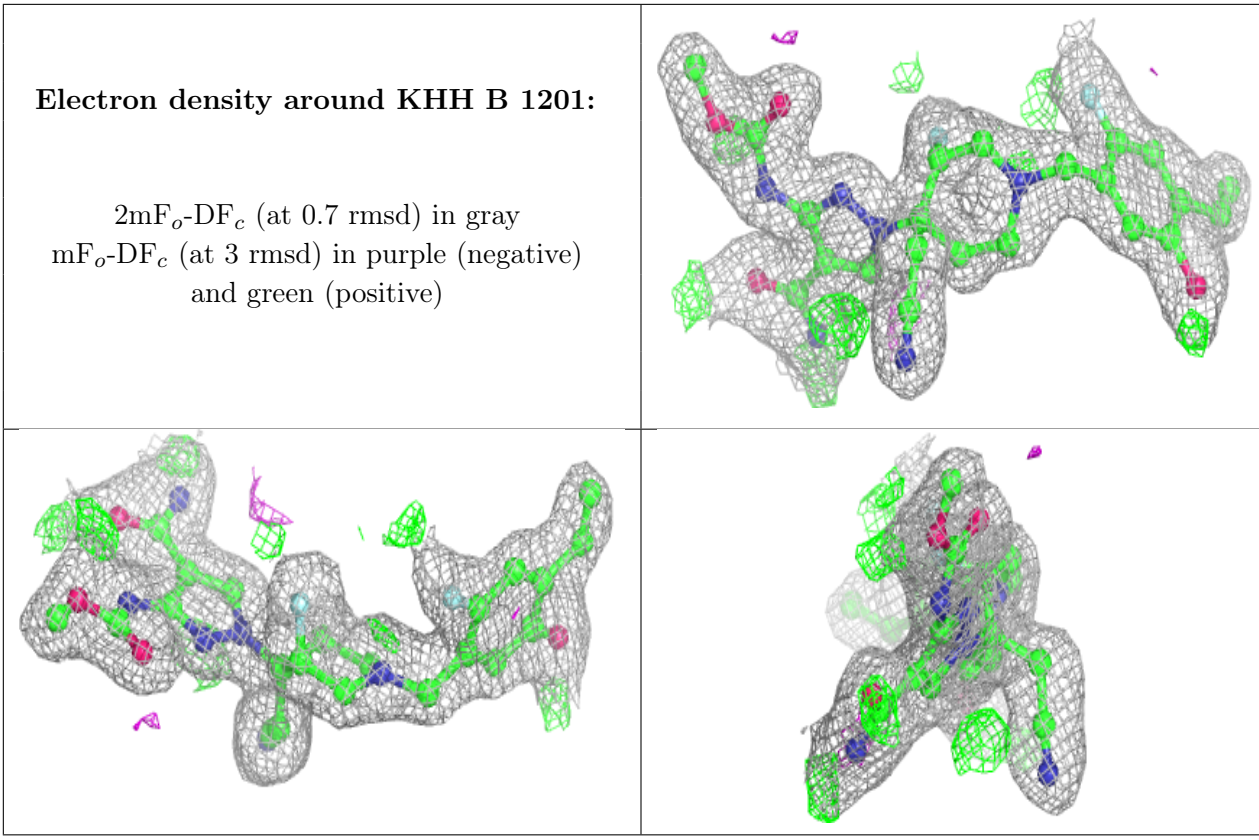
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
2	KHH	B	1201	34/34	0.92	0.12	17,21,31,31	0
2	KHH	A	1201	34/34	0.94	0.10	16,19,29,31	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.

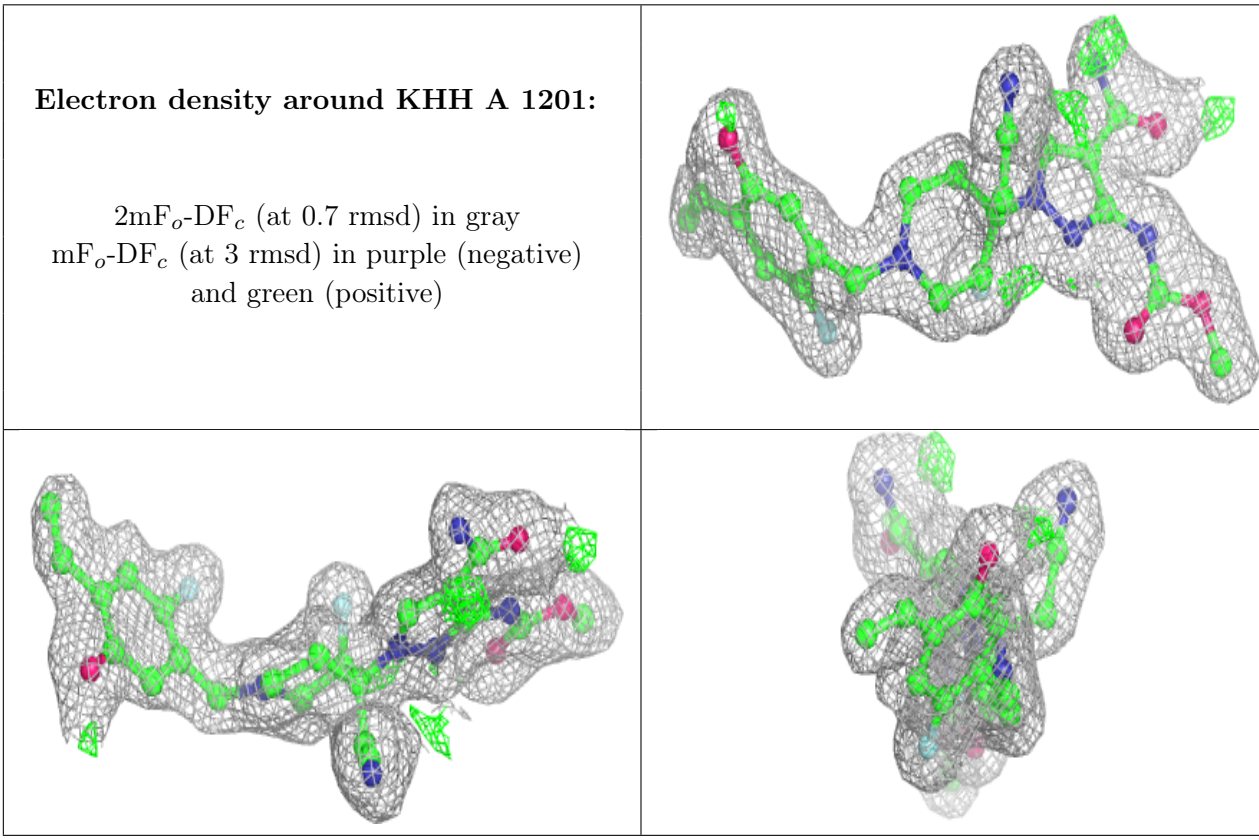
Electron density around KHH B 1201:

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



Electron density around KHH A 1201:

$2mF_o$-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.