



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:25 AM UTC

PDB ID : 3NC9 / pdb_00003nc9
Title : X-ray structure of ketohexokinase complexed with an indazole compound
Authors : Abad, M.C.; Gibbs, A.C.; Spurlino, J.C.
Deposited on : 2010-06-04
Resolution : 2.40 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

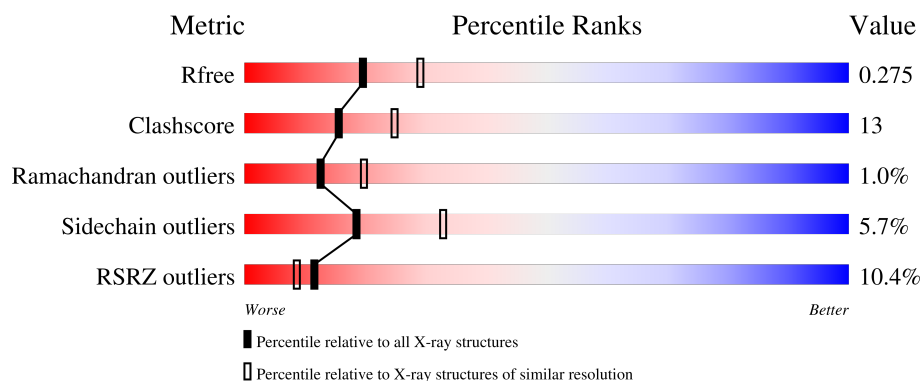
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	313	16% 60% 31% 5%
1	B	313	4% 75% 20% ...

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4656 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 Ketohexokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	9	0	0
			2259	1414	405	428	12			
1	B	301	Total	C	N	O	S	0	0	0
			2296	1438	413	433	12			

There are 38 discrepancies between the modelled and reference sequences:

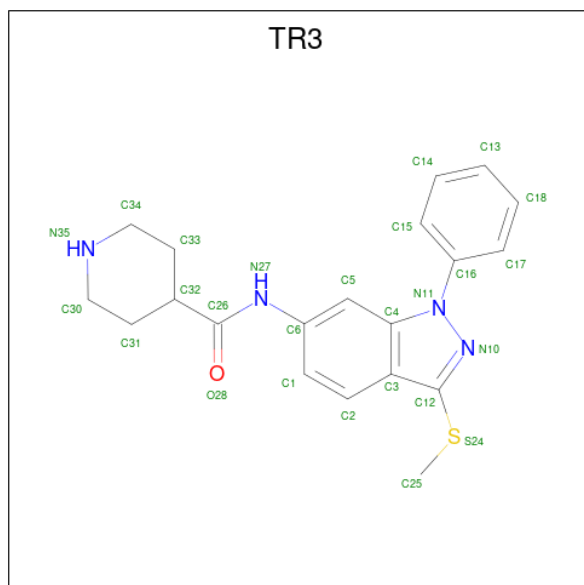
Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	expression tag	UNP P50053-2
A	-13	GLY	-	expression tag	UNP P50053-2
A	-12	SER	-	expression tag	UNP P50053-2
A	-11	SER	-	expression tag	UNP P50053-2
A	-10	HIS	-	expression tag	UNP P50053-2
A	-9	HIS	-	expression tag	UNP P50053-2
A	-8	HIS	-	expression tag	UNP P50053-2
A	-7	HIS	-	expression tag	UNP P50053-2
A	-6	HIS	-	expression tag	UNP P50053-2
A	-5	HIS	-	expression tag	UNP P50053-2
A	-4	SER	-	expression tag	UNP P50053-2
A	-3	SER	-	expression tag	UNP P50053-2
A	-2	GLY	-	expression tag	UNP P50053-2
A	-1	LEU	-	expression tag	UNP P50053-2
A	0	VAL	-	expression tag	UNP P50053-2
A	1	PRO	-	expression tag	UNP P50053-2
A	2	ARG	-	expression tag	UNP P50053-2
A	3	GLY	-	expression tag	UNP P50053-2
A	4	SER	-	expression tag	UNP P50053-2
B	-14	MET	-	expression tag	UNP P50053-2
B	-13	GLY	-	expression tag	UNP P50053-2
B	-12	SER	-	expression tag	UNP P50053-2
B	-11	SER	-	expression tag	UNP P50053-2
B	-10	HIS	-	expression tag	UNP P50053-2
B	-9	HIS	-	expression tag	UNP P50053-2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	HIS	-	expression tag	UNP P50053-2
B	-7	HIS	-	expression tag	UNP P50053-2
B	-6	HIS	-	expression tag	UNP P50053-2
B	-5	HIS	-	expression tag	UNP P50053-2
B	-4	SER	-	expression tag	UNP P50053-2
B	-3	SER	-	expression tag	UNP P50053-2
B	-2B	GLY	-	expression tag	UNP P50053-2
B	-1A	LEU	-	expression tag	UNP P50053-2
B	0	VAL	-	expression tag	UNP P50053-2
B	1	PRO	-	expression tag	UNP P50053-2
B	2	ARG	-	expression tag	UNP P50053-2
B	3	GLY	-	expression tag	UNP P50053-2
B	4	SER	-	expression tag	UNP P50053-2

- Molecule 2 is N-[3-(methylsulfanyl)-1-phenyl-1H-indazol-6-yl]piperidine-4-carboxamide (CCD ID: TR3) (formula: C₂₀H₂₂N₄OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			26	20	4	1	1		
2	B	1	Total	C	N	O	S	0	0
			26	20	4	1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

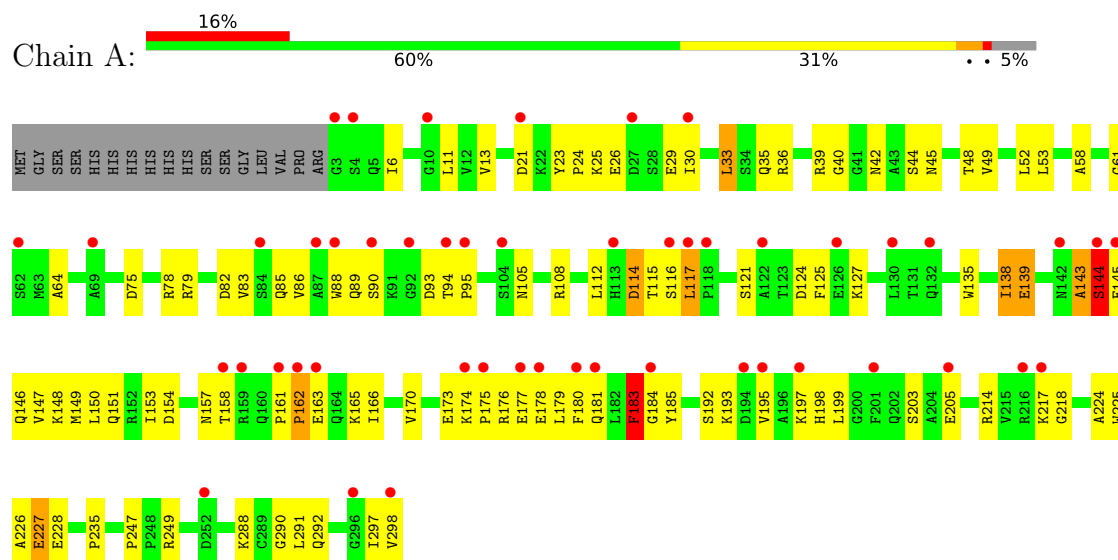
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	13	Total	O	0	0
			13	13		
4	B	26	Total	O	0	0
			26	26		

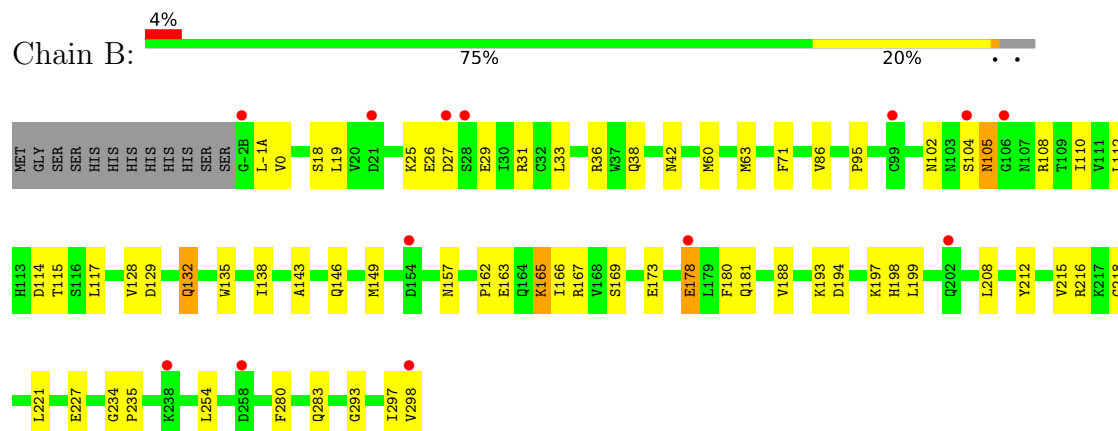
3 Residue-property plots [i](#)

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: Ketohexokinase



• Molecule 1: Ketohexokinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.70Å 85.69Å 136.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.31 – 2.40 36.31 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.3 (36.31-2.40) 95.3 (36.31-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.39Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.237 , 0.285 0.226 , 0.275	Depositor DCC
R_{free} test set	1992 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	60.7	Xtriage
Anisotropy	0.383	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4656	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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: TR3, SO4

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	0.50	0/2301	0.89	5/3111 (0.2%)
1	B	0.65	0/2339	0.91	5/3163 (0.2%)
All	All	0.58	0/4640	0.90	10/6274 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	183	PHE	N-CA-C	6.12	120.81	113.23
1	B	234	GLY	CA-C-N	5.61	125.45	119.28
1	B	234	GLY	C-N-CA	5.61	125.45	119.28
1	B	143	ALA	N-CA-C	5.35	116.80	111.07
1	A	226	ALA	CB-CA-C	-5.17	110.59	116.54
1	A	247	PRO	CA-C-N	5.12	124.91	119.28
1	A	247	PRO	C-N-CA	5.12	124.91	119.28
1	B	138	ILE	N-CA-C	5.12	115.41	107.28
1	A	138	ILE	N-CA-C	5.11	115.40	107.28
1	B	42	ASN	N-CA-C	5.06	116.49	110.97

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	2259	0	2237	78	0
1	B	2296	0	2283	48	0
2	B	52	0	44	8	0
3	B	10	0	0	1	0
4	A	13	0	0	0	0
4	B	26	0	0	0	0
All	All	4656	0	4564	123	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 13.

All (123) 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:146:GLN:HA	1:A:149:MET:HE3	1.46	0.97
1:A:157:ASN:HD21	1:A:166:ILE:H	1.25	0.85
1:A:64:ALA:H	1:A:94:THR:HG22	1.43	0.83
1:B:157:ASN:HD21	1:B:166:ILE:H	1.25	0.83
1:A:183:PHE:N	1:A:184:GLY:HA3	1.93	0.82
1:A:143:ALA:HB3	1:A:178:GLU:HG2	1.60	0.82
1:A:147:VAL:O	1:A:151:GLN:HG2	1.84	0.78
1:A:143:ALA:HB1	1:A:181:GLN:HE21	1.50	0.77
1:B:104:SER:O	1:B:105:ASN:HB3	1.85	0.73
1:B:27:ASP:OD2	2:B:299:TR3:H30A	1.89	0.72
1:A:42:ASN:HB2	1:A:139:GLU:OE1	1.91	0.70
1:A:150:LEU:HD13	1:A:185:TYR:HB2	1.76	0.68
1:A:93:ASP:HB2	1:A:115:THR:HG21	1.74	0.67
1:B:36:ARG:HD3	1:B:38:GLN:HE21	1.58	0.67
1:A:199:LEU:HD23	1:A:199:LEU:O	1.96	0.66
1:A:83:VAL:O	1:A:86:VAL:HG23	1.97	0.65
1:B:162:PRO:HA	1:B:165:LYS:CD	2.28	0.63
1:A:157:ASN:ND2	1:A:166:ILE:H	1.93	0.63
1:A:24:PRO:HG3	1:A:30:ILE:HD13	1.81	0.63
1:A:161:PRO:C	1:A:163:GLU:H	2.06	0.62
1:B:115:THR:HG22	1:B:115:THR:O	1.97	0.62
1:B:102:ASN:HB3	1:B:105:ASN:OD1	1.98	0.62
1:A:197:LYS:C	1:A:199:LEU:H	2.08	0.62
1:A:193:LYS:HD2	1:A:225:TRP:CE2	2.34	0.61
1:A:162:PRO:HA	1:A:165:LYS:HE2	1.82	0.60
1:A:227:GLU:H	1:A:227:GLU:CD	2.08	0.60
1:B:27:ASP:CG	2:B:299:TR3:H30A	2.28	0.59
1:B:178:GLU:O	1:B:181:GLN:HG2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:MET:HE1	1:B:128:VAL:HG21	1.86	0.58
1:B:280:PHE:HA	1:B:283:GLN:HE21	1.68	0.58
1:A:121:SER:HB2	1:A:145:GLU:OE2	2.04	0.58
1:A:13:VAL:HG22	1:A:95:PRO:HB2	1.85	0.58
1:A:116:SER:O	1:A:117:LEU:C	2.47	0.57
1:A:192:SER:OG	1:A:195:VAL:HG23	2.04	0.57
1:B:162:PRO:HA	1:B:165:LYS:HD3	1.86	0.57
1:A:151:GLN:HE21	1:A:151:GLN:HA	1.70	0.56
1:B:293:GLY:HA3	3:B:301:SO4:O2	2.05	0.56
1:A:29:GLU:OE2	1:B:112:LEU:HD22	2.06	0.56
1:B:0:VAL:HG11	1:B:132:GLN:HG2	1.86	0.56
1:A:48:THR:HG23	1:A:79:ARG:HH11	1.71	0.55
1:A:125:PHE:CD2	1:A:149:MET:HG2	2.42	0.55
1:A:53:LEU:HD23	1:A:298:VAL:HG11	1.88	0.55
1:A:205:GLU:H	1:A:205:GLU:CD	2.15	0.54
1:B:115:THR:O	1:B:115:THR:CG2	2.55	0.53
1:B:157:ASN:ND2	1:B:166:ILE:H	1.99	0.53
1:A:146:GLN:CA	1:A:149:MET:HE3	2.30	0.53
1:A:42:ASN:HB2	1:A:139:GLU:CD	2.33	0.53
1:A:127:LYS:O	1:A:127:LYS:HG2	2.09	0.53
1:B:108:ARG:HH21	1:B:110:ILE:HD11	1.75	0.52
1:A:24:PRO:HG3	1:A:30:ILE:CD1	2.39	0.52
1:A:178:GLU:O	1:A:181:GLN:HG2	2.09	0.52
1:A:21:ASP:HB3	1:A:33:LEU:HD21	1.92	0.51
1:A:48:THR:HG23	1:A:79:ARG:NH1	2.25	0.51
1:A:88:TRP:C	1:A:89:GLN:HG2	2.34	0.51
1:B:162:PRO:HA	1:B:165:LYS:HD2	1.93	0.51
1:A:218:GLY:HA2	1:A:235:PRO:HG3	1.93	0.51
1:B:25:LYS:O	1:B:26:GLU:C	2.52	0.51
1:A:173:GLU:O	1:A:195:VAL:HG22	2.11	0.51
1:A:150:LEU:HD13	1:A:185:TYR:CB	2.40	0.50
1:B:-1(A):LEU:C	1:B:-1(A):LEU:HD23	2.35	0.50
1:B:27:ASP:OD1	2:B:299:TR3:N35	2.44	0.50
1:A:58:ALA:HB2	1:A:82:ASP:HB3	1.93	0.50
1:A:176:ARG:HD3	1:A:178:GLU:HB3	1.93	0.50
1:B:193:LYS:O	1:B:197:LYS:HG3	2.13	0.49
1:B:19:LEU:HD23	1:B:33:LEU:HB2	1.95	0.49
1:B:71:PHE:CD2	1:B:71:PHE:C	2.91	0.49
1:A:288:LYS:C	1:A:288:LYS:HD3	2.38	0.48
1:B:169:SER:HA	1:B:188:VAL:O	2.14	0.48
1:A:125:PHE:CG	1:A:149:MET:HG2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:PRO:HD3	1:B:117:LEU:HD22	1.95	0.48
1:A:195:VAL:O	1:A:199:LEU:HB2	2.14	0.48
2:B:299:TR3:O28	2:B:299:TR3:H5	2.14	0.48
1:A:177:GLU:HA	1:A:180:PHE:CD2	2.49	0.47
1:B:178:GLU:H	1:B:178:GLU:CD	2.22	0.47
1:A:48:THR:O	1:A:52:LEU:HG	2.15	0.47
1:B:178:GLU:OE1	1:B:178:GLU:N	2.48	0.47
1:B:27:ASP:OD1	2:B:299:TR3:H30A	2.15	0.46
1:B:194:ASP:O	1:B:198:HIS:HD2	1.98	0.46
1:A:6:ILE:HG23	1:A:135:TRP:HB3	1.97	0.46
1:A:161:PRO:C	1:A:163:GLU:N	2.73	0.46
1:A:174:LYS:O	1:A:179:LEU:HD12	2.15	0.46
1:B:180:PHE:CZ	1:B:199:LEU:HD21	2.50	0.46
1:B:208:LEU:CD2	1:B:221:LEU:HG	2.46	0.46
1:A:39:ARG:NE	1:A:75:ASP:OD2	2.36	0.46
1:B:194:ASP:O	1:B:198:HIS:CD2	2.69	0.46
1:A:180:PHE:HD1	1:A:214:ARG:HH12	1.63	0.46
1:B:63:MET:HE3	1:B:86:VAL:HG21	1.98	0.46
1:B:108:ARG:NH2	1:B:110:ILE:HD11	2.30	0.45
1:A:58:ALA:CB	1:A:82:ASP:HB3	2.46	0.45
1:B:218:GLY:HA2	1:B:235:PRO:HG3	1.99	0.45
1:A:297:ILE:HG13	1:A:298:VAL:N	2.32	0.45
1:A:154:ASP:O	1:A:158:THR:HG23	2.16	0.44
1:A:45:ASN:O	1:A:49:VAL:HG23	2.17	0.44
1:A:143:ALA:O	1:A:144:SER:C	2.61	0.44
1:B:212:TYR:O	1:B:215:VAL:HG22	2.18	0.44
1:A:23:TYR:CD2	1:B:26:GLU:HB2	2.52	0.44
1:B:0:VAL:CG1	1:B:132:GLN:HG2	2.47	0.44
1:A:290:GLY:O	1:A:291:LEU:HD23	2.18	0.43
1:A:25:LYS:O	1:A:26:GLU:C	2.61	0.43
2:B:300:TR3:H17	2:B:300:TR3:H5	2.00	0.43
1:A:174:LYS:C	1:A:176:ARG:H	2.26	0.43
1:A:108:ARG:O	1:B:27:ASP:HA	2.18	0.43
1:A:197:LYS:O	1:A:199:LEU:N	2.51	0.43
1:B:135:TRP:CE3	1:B:167:ARG:HB2	2.54	0.42
1:A:125:PHE:CB	1:A:149:MET:HG2	2.49	0.42
1:A:151:GLN:HA	1:A:151:GLN:NE2	2.33	0.42
1:A:174:LYS:HA	1:A:175:PRO:HD3	1.78	0.42
1:A:40:GLY:O	1:A:44:SER:HB3	2.19	0.42
1:A:23:TYR:HA	1:A:24:PRO:HD3	1.91	0.42
1:A:35:GLN:HG2	1:B:114:ASP:OD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ASP:CB	1:B:31:ARG:NH1	2.83	0.42
1:A:148:LYS:O	1:A:151:GLN:HB2	2.19	0.42
1:A:149:MET:O	1:A:153:ILE:HG13	2.19	0.42
1:B:129:ASP:C	1:B:129:ASP:OD1	2.63	0.41
1:B:27:ASP:OD1	2:B:299:TR3:C30	2.68	0.41
1:A:36:ARG:NH1	1:A:292:GLN:HB3	2.35	0.41
1:A:61:GLY:O	1:A:86:VAL:HG13	2.21	0.41
1:A:150:LEU:HD13	1:A:185:TYR:CG	2.56	0.41
1:B:146:GLN:HA	1:B:149:MET:HG3	2.02	0.41
1:B:297:ILE:O	1:B:298:VAL:OXT	2.40	0.40
1:A:203:SER:HA	1:A:225:TRP:CZ2	2.57	0.40
1:A:224:ALA:HB1	2:B:299:TR3:H25A	2.04	0.40
1:A:125:PHE:C	1:A:127:LYS:H	2.30	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	294/313 (94%)	266 (90%)	23 (8%)	5 (2%)	7	10
1	B	299/313 (96%)	286 (96%)	12 (4%)	1 (0%)	36	50
All	All	593/626 (95%)	552 (93%)	35 (6%)	6 (1%)	12	20

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	105	ASN
1	A	143	ALA
1	A	144	SER
1	A	198	HIS
1	A	117	LEU

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Mol	Chain	Res	Type
1	A	162	PRO

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	245/260 (94%)	227 (93%)	18 (7%)	13	22
1	B	249/260 (96%)	239 (96%)	10 (4%)	28	47
All	All	494/520 (95%)	466 (94%)	28 (6%)	18	33

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

Mol	Chain	Res	Type
1	A	11	LEU
1	A	33	LEU
1	A	78	ARG
1	A	85	GLN
1	A	90	SER
1	A	105	ASN
1	A	112	LEU
1	A	114	ASP
1	A	124	ASP
1	A	138	ILE
1	A	139	GLU
1	A	144	SER
1	A	170	VAL
1	A	183	PHE
1	A	217	LYS
1	A	227	GLU
1	A	228	GLU
1	A	249	ARG
1	B	18	SER
1	B	29	GLU
1	B	132	GLN
1	B	163	GLU

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Mol	Chain	Res	Type
1	B	165	LYS
1	B	173	GLU
1	B	178	GLU
1	B	216	ARG
1	B	227	GLU
1	B	254	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	38	GLN
1	A	151	GLN
1	A	156	HIS
1	A	157	ASN
1	A	181	GLN
1	A	202	GLN
1	B	5	GLN
1	B	38	GLN
1	B	103	ASN
1	B	156	HIS
1	B	157	ASN
1	B	160	GLN
1	B	198	HIS
1	B	275	GLN
1	B	283	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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
3	SO4	B	302	-	4,4,4	0.26	0	6,6,6	0.39	0
3	SO4	B	301	-	4,4,4	0.26	0	6,6,6	0.34	0
2	TR3	B	299	-	28,29,29	1.27	2 (7%)	36,40,40	1.43	6 (16%)
2	TR3	B	300	-	28,29,29	1.20	1 (3%)	36,40,40	1.48	6 (16%)

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	TR3	B	299	-	-	2/14/22/22	0/4/4/4
2	TR3	B	300	-	-	2/14/22/22	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	299	TR3	C3-C12	-4.27	1.41	1.48
2	B	300	TR3	C3-C12	-3.06	1.43	1.48
2	B	299	TR3	N11-N10	-2.51	1.35	1.38

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	300	TR3	C3-C4-N11	-4.24	104.50	106.42
2	B	299	TR3	C3-C4-N11	-3.96	104.62	106.42
2	B	300	TR3	C4-N11-N10	3.95	112.74	111.02
2	B	300	TR3	C2-C3-C12	3.16	137.68	126.81
2	B	299	TR3	C12-N10-N11	-2.99	105.77	109.55
2	B	299	TR3	C4-N11-N10	2.97	112.31	111.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	300	TR3	C3-C12-N10	-2.80	112.11	115.49
2	B	300	TR3	C12-N10-N11	-2.80	106.01	109.55
2	B	299	TR3	C2-C3-C12	2.78	136.36	126.81
2	B	299	TR3	C4-C3-C12	-2.46	104.29	108.20
2	B	300	TR3	C4-C3-C12	-2.23	104.64	108.20
2	B	299	TR3	C3-C12-N10	-2.06	113.00	115.49

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	299	TR3	C3-C12-S24-C25
2	B	299	TR3	N10-C12-S24-C25
2	B	300	TR3	C3-C12-S24-C25
2	B	300	TR3	N10-C12-S24-C25

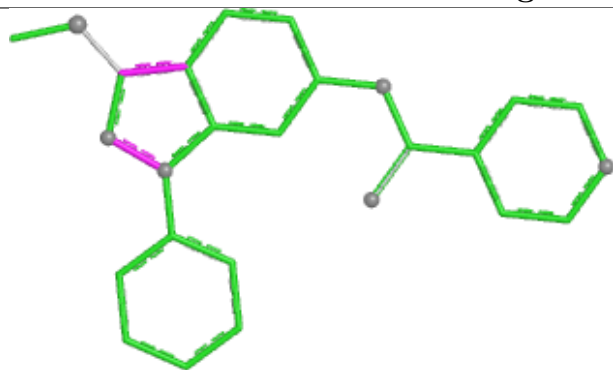
There are no ring outliers.

3 monomers are involved in 9 short contacts:

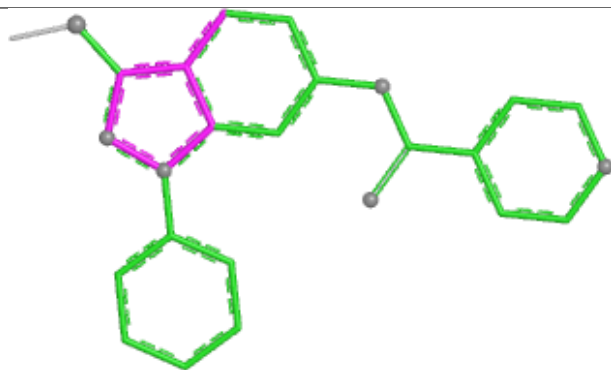
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	SO4	1	0
2	B	299	TR3	7	0
2	B	300	TR3	1	0

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.

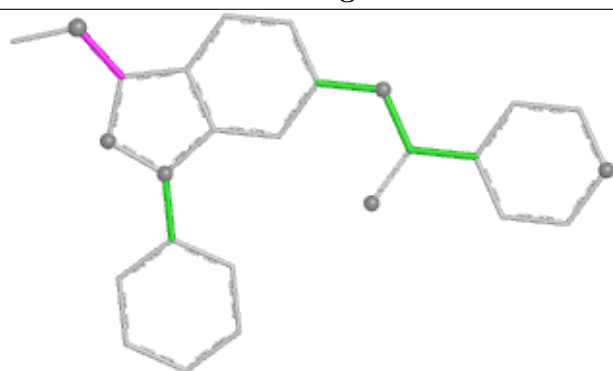
Ligand TR3 B 299



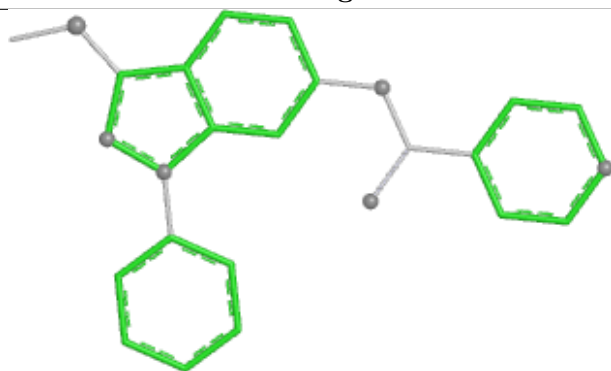
Bond lengths



Bond angles

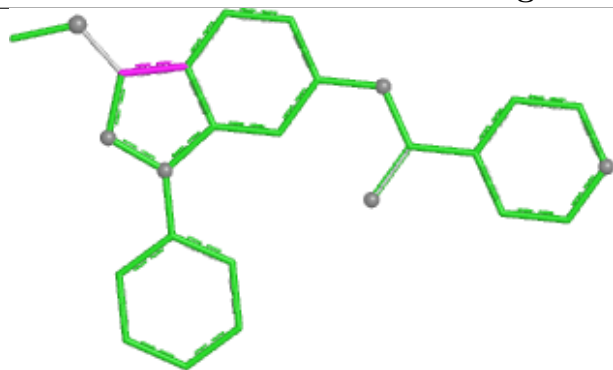


Torsions

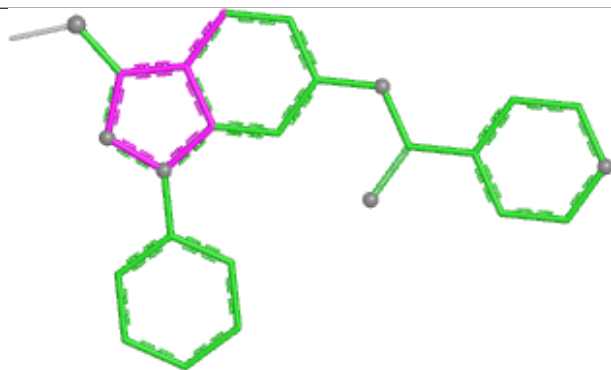


Rings

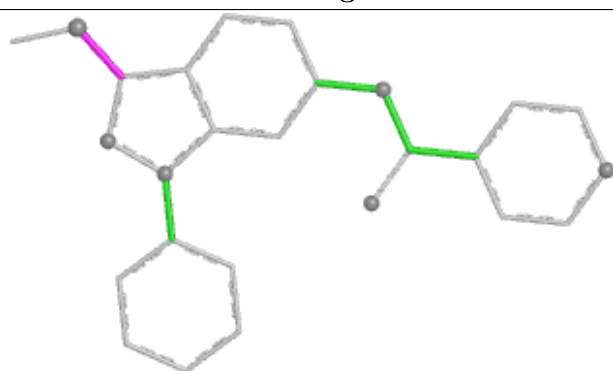
Ligand TR3 B 300



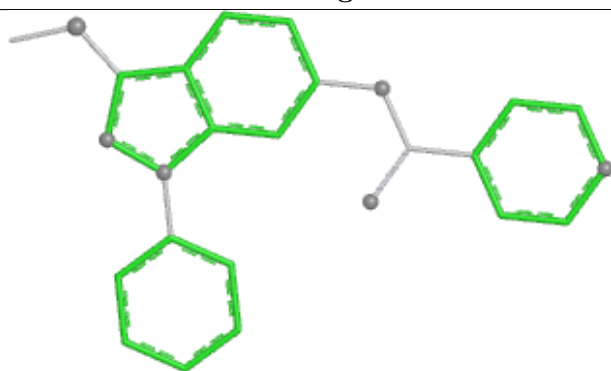
Bond lengths



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, 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	295/313 (94%)	1.12	49 (16%) 4 3	46, 74, 104, 125	0
1	B	301/313 (96%)	0.45	13 (4%) 40 36	41, 55, 86, 101	0
All	All	596/626 (95%)	0.78	62 (10%) 11 9	41, 64, 100, 125	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	298	VAL	4.2
1	A	92	GLY	4.2
1	B	-2(B)	GLY	4.0
1	A	180	PHE	3.8
1	A	298	VAL	3.8
1	B	21	ASP	3.8
1	B	99	CYS	3.8
1	A	161	PRO	3.8
1	A	126	GLU	3.7
1	A	90	SER	3.7
1	A	184	GLY	3.3
1	A	163	GLU	3.3
1	A	4	SER	3.2
1	A	205	GLU	3.0
1	A	84	SER	3.0
1	A	3	GLY	3.0
1	A	217	LYS	2.9
1	A	118	PRO	2.9
1	A	132	GLN	2.9
1	A	252	ASP	2.8
1	B	27	ASP	2.8
1	A	117	LEU	2.8
1	A	130	LEU	2.8
1	A	113	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	178	GLU	2.7
1	A	88	TRP	2.7
1	A	122	ALA	2.7
1	A	197	LYS	2.6
1	A	10	GLY	2.6
1	B	258	ASP	2.6
1	A	95	PRO	2.5
1	B	202	GLN	2.5
1	A	162	PRO	2.5
1	A	177	GLU	2.5
1	B	154	ASP	2.5
1	A	69	ALA	2.4
1	A	62	SER	2.4
1	A	116	SER	2.4
1	A	201	PHE	2.4
1	A	142	ASN	2.4
1	A	174	LYS	2.4
1	A	158	THR	2.3
1	B	28	SER	2.3
1	A	27	ASP	2.3
1	A	175	PRO	2.3
1	A	159	ARG	2.3
1	A	194	ASP	2.3
1	B	238	LYS	2.3
1	A	87	ALA	2.2
1	A	144	SER	2.2
1	B	104	SER	2.2
1	A	21	ASP	2.2
1	A	104	SER	2.2
1	A	94	THR	2.2
1	A	181	GLN	2.1
1	A	296	GLY	2.1
1	B	106	GLY	2.1
1	A	145	GLU	2.1
1	A	30	ILE	2.1
1	A	216	ARG	2.0
1	A	195	VAL	2.0
1	B	178	GLU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

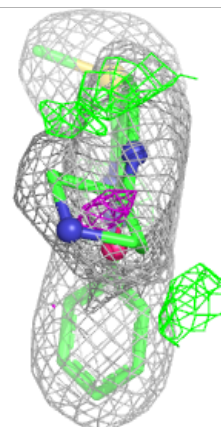
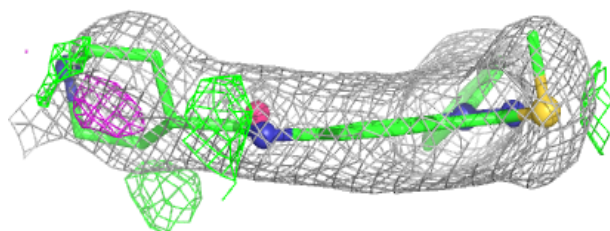
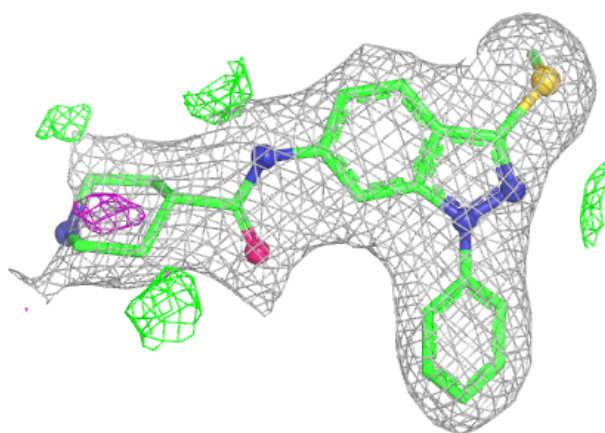
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	SO4	B	302	5/5	0.85	0.11	82,84,90,94	0
3	SO4	B	301	5/5	0.93	0.13	74,75,77,85	0
2	TR3	B	299	26/26	0.93	0.12	53,57,80,80	0
2	TR3	B	300	26/26	0.95	0.09	51,60,78,83	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.

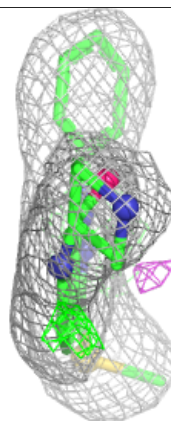
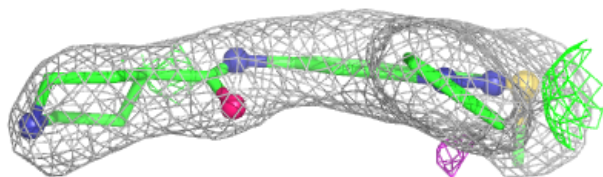
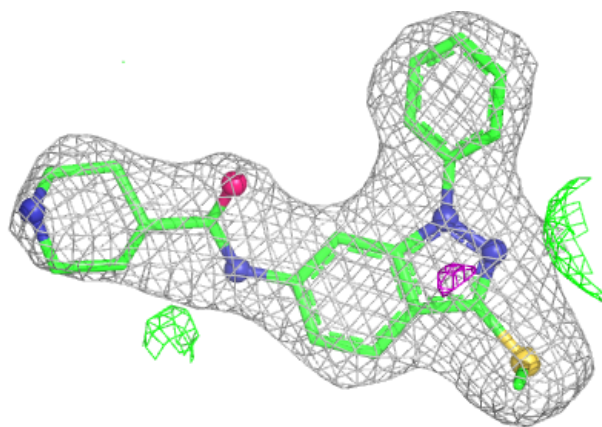
Electron density around TR3 B 299:

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 TR3 B 300:

$2mF_o-DF_c$ (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.