



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 04:08 pm BST

PDB ID : 4XVD
Title : 17beta-HSD5 in complex with 4-nitro-2-({4-[3-(trifluoromethyl)phenyl]piperazin-1-yl)methyl}phenol
Authors : Amano, Y.; Yamaguchi, T.; Niimi, T.; Sakashita, H.
Deposited on : 2015-01-27
Resolution : 2.81 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 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 : 2.11
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) : 2.11

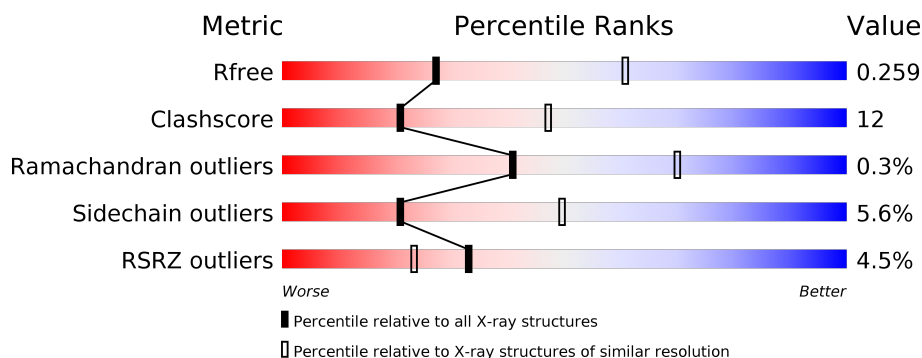
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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 on 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	331	<div> <div> <div></div> <div>71%</div> <div>21%</div> <div>• 5%</div> </div> <div>8%</div> </div>
1	B	331	<div> <div> <div></div> <div>66%</div> <div>26%</div> <div>• 6%</div> </div> <div>8%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5173 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 Aldo-keto reductase family 1 member C3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2519	1606	439	462	12			
1	B	311	Total	C	N	O	S	0	0	0
			2491	1589	435	455	12			

There are 16 discrepancies between the modelled and reference sequences:

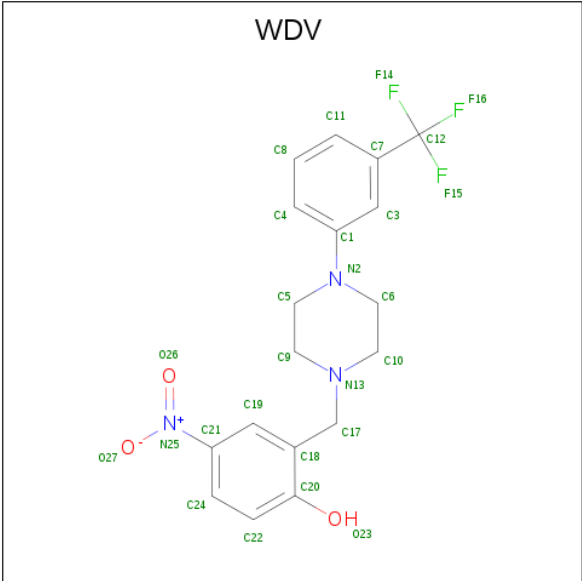
Chain	Residue	Modelled	Actual	Comment	Reference
A	324	LEU	-	expression tag	UNP P42330
A	325	GLU	-	expression tag	UNP P42330
A	326	HIS	-	expression tag	UNP P42330
A	327	HIS	-	expression tag	UNP P42330
A	328	HIS	-	expression tag	UNP P42330
A	329	HIS	-	expression tag	UNP P42330
A	330	HIS	-	expression tag	UNP P42330
A	331	HIS	-	expression tag	UNP P42330
B	324	LEU	-	expression tag	UNP P42330
B	325	GLU	-	expression tag	UNP P42330
B	326	HIS	-	expression tag	UNP P42330
B	327	HIS	-	expression tag	UNP P42330
B	328	HIS	-	expression tag	UNP P42330
B	329	HIS	-	expression tag	UNP P42330
B	330	HIS	-	expression tag	UNP P42330
B	331	HIS	-	expression tag	UNP P42330

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	B	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 3 is 4-nitro-2-({4-[3-(trifluoromethyl)phenyl]piperazin-1-yl}methyl)phenol (three-letter code: WDV) (formula: C₁₈H₁₈F₃N₃O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			27	18	3	3	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	F	N	O	0	0
			27	18	3	3	3		

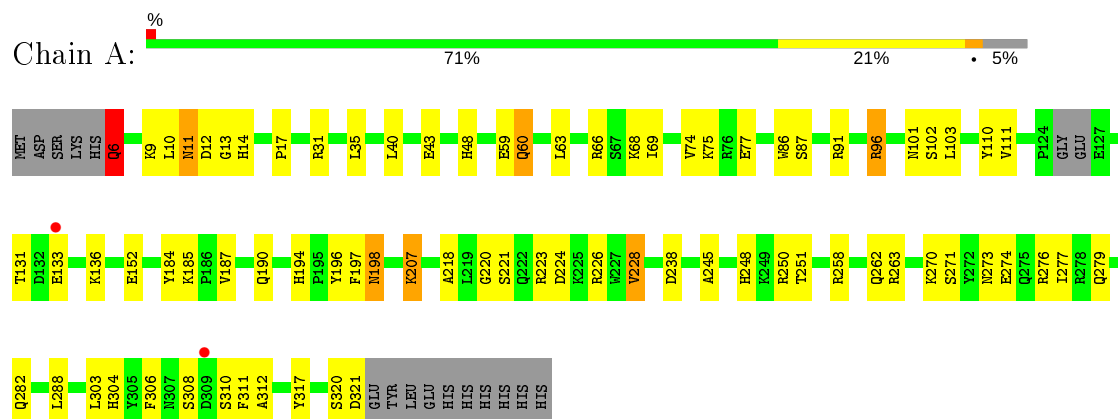
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	10	Total	O	0	0
			10	10		
4	B	3	Total	O	0	0
			3	3		

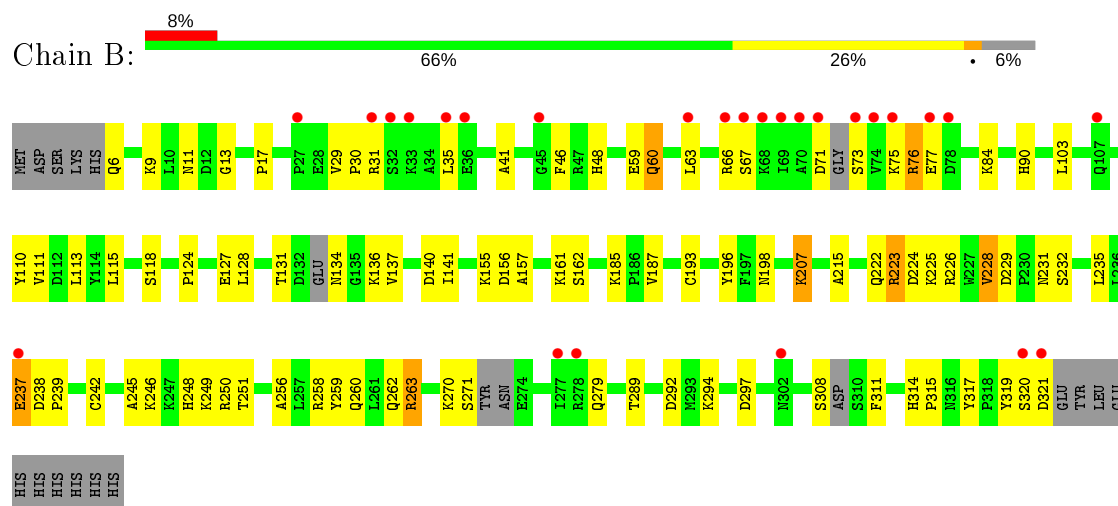
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Aldo-keto reductase family 1 member C3



- Molecule 1: Aldo-keto reductase family 1 member C3



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	145.79 Å 145.79 Å 74.29 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.67 – 2.81 40.67 – 2.81	Depositor EDS
% Data completeness (in resolution range)	90.4 (40.67-2.81) 90.4 (40.67-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.81 Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.209 , 0.259 0.209 , 0.259	Depositor DCC
R_{free} test set	894 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.037 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5173	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% 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: NAP, WDV

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	0.61	0/2576	0.77	1/3485 (0.0%)
1	B	0.60	2/2544 (0.1%)	0.73	3/3436 (0.1%)
All	All	0.60	2/5120 (0.0%)	0.75	4/6921 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	6	GLN	CA-CB	-6.14	1.40	1.53
1	B	6	GLN	CB-CG	-5.10	1.38	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	6	GLN	CA-CB-CG	-6.97	98.06	113.40
1	B	6	GLN	N-CA-CB	-6.76	98.44	110.60
1	B	223	ARG	NE-CZ-NH1	6.74	123.67	120.30
1	A	6	GLN	CA-CB-CG	-5.22	101.91	113.40

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	2519	0	2525	53	0
1	B	2491	0	2503	71	0
2	A	48	0	25	5	0
2	B	48	0	25	5	0
3	A	27	0	18	2	0
3	B	27	0	17	1	0
4	A	10	0	0	0	0
4	B	3	0	0	0	0
All	All	5173	0	5113	125	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (125) 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:60:GLN:N	1:A:60:GLN:HE21	1.44	1.16
1:B:76:ARG:NH2	1:B:110:TYR:CE1	2.14	1.16
1:B:76:ARG:NH2	1:B:110:TYR:CZ	2.18	1.09
1:B:271:SER:HA	2:B:401:NAP:O1X	1.55	1.06
1:A:60:GLN:H	1:A:60:GLN:NE2	1.62	0.98
1:B:75:LYS:HD3	1:B:77:GLU:CD	1.84	0.97
1:B:11:ASN:OD1	1:B:187:VAL:HG12	1.65	0.96
1:B:223:ARG:NH1	1:B:232:SER:O	1.99	0.93
1:A:60:GLN:H	1:A:60:GLN:HE21	0.81	0.81
1:A:40:LEU:HD13	1:A:274:GLU:HG3	1.64	0.79
1:B:76:ARG:NH2	1:B:110:TYR:OH	2.17	0.77
1:A:60:GLN:N	1:A:60:GLN:NE2	2.28	0.74
1:B:75:LYS:HD3	1:B:77:GLU:OE2	1.85	0.74
1:B:196:TYR:OH	1:B:238:ASP:HB2	1.86	0.74
1:A:320:SER:O	1:A:321:ASP:HB2	1.87	0.73
1:B:271:SER:CA	2:B:401:NAP:O1X	2.34	0.73
1:A:220:GLY:O	1:A:221:SER:OG	2.05	0.72
1:B:60:GLN:NE2	1:B:60:GLN:H	1.91	0.69
1:A:6:GLN:NE2	1:A:6:GLN:HA	2.02	0.68
1:B:251:THR:OG1	1:B:279:GLN:NE2	2.26	0.68
1:B:224:ASP:OD2	1:B:226:ARG:NH2	2.27	0.67
1:A:43:GLU:OE2	1:A:68:LYS:NZ	2.27	0.67
1:B:141:ILE:HD11	1:B:315:PRO:HG2	1.77	0.66
1:A:251:THR:OG1	1:A:279:GLN:NE2	2.30	0.64
1:A:306:PHE:CD2	3:A:402:WDV:H9	2.32	0.64
1:A:245:ALA:HB1	1:A:250:ARG:O	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:ARG:HD2	1:A:262:GLN:OE1	2.01	0.61
1:B:207:LYS:NZ	1:B:263:ARG:O	2.34	0.61
1:B:76:ARG:HH21	1:B:76:ARG:CG	2.15	0.60
1:A:10:LEU:HB2	1:A:12:ASP:OD1	2.02	0.60
1:B:17:PRO:HB2	1:B:48:HIS:HB2	1.82	0.60
1:B:59:GLU:HB2	1:B:60:GLN:NE2	2.17	0.60
1:A:9:LYS:HE3	1:A:13:GLY:HA2	1.84	0.58
1:B:314:HIS:ND1	1:B:315:PRO:HD2	2.19	0.58
1:A:224:ASP:OD1	1:A:226:ARG:NH2	2.35	0.58
1:B:110:TYR:CD2	1:B:161:LYS:HG2	2.38	0.58
1:B:193:CYS:HB3	1:B:215:ALA:CB	2.34	0.57
1:A:198:ASN:HB2	1:A:263:ARG:NH1	2.19	0.57
1:B:60:GLN:HE21	1:B:60:GLN:N	2.02	0.57
1:B:271:SER:OG	2:B:401:NAP:O3X	2.22	0.57
1:B:75:LYS:HD3	1:B:77:GLU:OE1	2.05	0.57
1:A:223:ARG:HA	1:A:228:VAL:HG21	1.87	0.57
1:A:11:ASN:OD1	1:A:187:VAL:HG12	2.05	0.57
1:A:207:LYS:NZ	1:A:263:ARG:O	2.37	0.57
1:A:308:SER:HB3	1:A:311:PHE:HB2	1.86	0.57
1:B:222:GLN:O	1:B:223:ARG:HB2	2.05	0.57
1:B:76:ARG:NH2	1:B:110:TYR:HE1	1.97	0.55
1:A:31:ARG:HB3	1:A:60:GLN:OE1	2.06	0.55
1:B:60:GLN:NE2	1:B:60:GLN:N	2.54	0.55
1:B:11:ASN:HD21	1:B:187:VAL:HA	1.73	0.54
1:A:271:SER:OG	2:A:401:NAP:O3X	2.14	0.53
1:B:258:ARG:HD2	1:B:262:GLN:OE1	2.08	0.53
1:B:223:ARG:HA	1:B:228:VAL:HG21	1.91	0.53
1:B:9:LYS:HE3	1:B:13:GLY:HA2	1.90	0.53
1:B:103:LEU:HD21	1:B:111:VAL:HG13	1.90	0.52
1:B:229:ASP:OD1	1:B:231:ASN:HB2	2.09	0.52
1:B:320:SER:O	1:B:321:ASP:HB2	2.10	0.52
1:B:63:LEU:HA	1:B:66:ARG:NH2	2.24	0.52
1:A:306:PHE:HD2	3:A:402:WDV:H9	1.75	0.51
1:A:152:GLU:HG2	1:A:184:TYR:HD2	1.75	0.51
1:A:276:ARG:NH2	2:A:401:NAP:O2X	2.44	0.51
1:A:110:TYR:CD1	1:A:110:TYR:C	2.83	0.51
1:B:134:ASN:HD22	1:B:134:ASN:N	2.09	0.51
1:A:258:ARG:HG3	1:A:288:LEU:HD12	1.92	0.51
1:B:294:LYS:HA	1:B:297:ASP:HB2	1.92	0.51
1:B:113:LEU:HA	1:B:162:SER:O	2.10	0.51
1:B:11:ASN:ND2	1:B:187:VAL:HA	2.25	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:TYR:OH	1:A:238:ASP:HB2	2.11	0.49
1:A:258:ARG:HG3	1:A:288:LEU:CD1	2.42	0.49
1:A:17:PRO:HB2	1:A:48:HIS:HB2	1.93	0.49
1:B:239:PRO:HA	1:B:242:CYS:HB2	1.94	0.49
1:A:270:LYS:O	2:A:401:NAP:H8A	2.13	0.49
1:B:90:HIS:CE1	1:B:118:SER:O	2.67	0.48
1:A:250:ARG:HH21	1:A:282:GLN:HE21	1.61	0.48
1:A:12:ASP:OD2	1:A:14:HIS:HD2	1.96	0.48
1:A:197:PHE:CE2	1:A:304:HIS:HB3	2.49	0.48
1:B:84:LYS:HG2	1:B:115:LEU:HB2	1.95	0.47
1:B:76:ARG:NH2	1:B:76:ARG:CG	2.77	0.47
1:B:75:LYS:HG2	1:B:76:ARG:N	2.28	0.47
1:B:193:CYS:HB3	1:B:215:ALA:HB1	1.97	0.47
1:B:110:TYR:CE2	1:B:161:LYS:HG2	2.50	0.47
1:B:71:ASP:C	1:B:73:SER:N	2.69	0.47
1:A:103:LEU:HD21	1:A:111:VAL:HG13	1.96	0.46
1:A:194:HIS:HB3	1:A:218:ALA:HA	1.97	0.46
1:B:260:GLN:HA	1:B:260:GLN:OE1	2.16	0.46
1:A:10:LEU:C	1:A:12:ASP:H	2.19	0.46
1:B:235:LEU:C	1:B:237:GLU:H	2.19	0.46
1:B:110:TYR:HD2	1:B:161:LYS:HG2	1.80	0.46
1:A:59:GLU:HB2	1:A:60:GLN:NE2	2.31	0.46
1:B:35:LEU:HB2	1:B:60:GLN:HB3	1.97	0.46
1:B:76:ARG:HG2	1:B:76:ARG:HH21	1.79	0.44
1:A:75:LYS:HE3	1:A:77:GLU:OE2	2.16	0.44
1:B:63:LEU:O	1:B:67:SER:N	2.44	0.44
1:B:317:TYR:HE1	3:B:402:WDV:F16	1.90	0.44
2:B:401:NAP:O2N	2:B:401:NAP:H3D	2.18	0.44
1:B:59:GLU:HB2	1:B:60:GLN:HE21	1.81	0.43
1:B:29:VAL:HA	1:B:30:PRO:HD2	1.86	0.43
1:B:270:LYS:O	2:B:401:NAP:H8A	2.18	0.43
1:A:31:ARG:N	1:A:31:ARG:HD2	2.33	0.43
1:B:224:ASP:CG	1:B:226:ARG:HH21	2.21	0.43
1:B:41:ALA:O	1:B:46:PHE:HB2	2.18	0.43
1:B:289:THR:H	1:B:292:ASP:HB2	1.84	0.42
1:A:224:ASP:CG	1:A:226:ARG:HH21	2.21	0.42
1:B:103:LEU:CD2	1:B:111:VAL:HG13	2.49	0.42
1:B:124:PRO:HD3	1:B:140:ASP:OD2	2.20	0.41
1:A:63:LEU:HA	1:A:66:ARG:NH2	2.34	0.41
1:A:69:ILE:HA	1:A:74:VAL:O	2.21	0.41
1:A:152:GLU:HG2	1:A:184:TYR:CD2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:ILE:CD1	1:B:315:PRO:HG2	2.49	0.41
1:B:256:ALA:O	1:B:259:TYR:HB3	2.21	0.41
1:A:35:LEU:HB2	1:A:60:GLN:HB3	2.02	0.41
1:B:245:ALA:HB1	1:B:250:ARG:O	2.20	0.41
1:A:312:ALA:HA	1:A:317:TYR:CD1	2.56	0.41
1:A:190:GLN:OE1	2:A:401:NAP:H2N	2.20	0.41
1:B:308:SER:HB3	1:B:311:PHE:HB2	2.01	0.41
1:B:245:ALA:O	1:B:250:ARG:N	2.53	0.41
1:A:96:ARG:HD2	1:A:96:ARG:HH21	1.77	0.41
1:A:270:LYS:HB2	2:A:401:NAP:H52N	2.03	0.41
1:B:317:TYR:CE2	1:B:319:TYR:O	2.74	0.41
1:A:86:TRP:CG	1:A:87:SER:N	2.89	0.40
1:B:75:LYS:CD	1:B:77:GLU:CD	2.72	0.40
1:A:248:HIS:O	1:A:250:ARG:HG2	2.21	0.40
1:B:248:HIS:O	1:B:249:LYS:HB2	2.20	0.40
1:A:273:ASN:O	1:A:277:ILE:HG13	2.21	0.40
1:B:155:LYS:C	1:B:157:ALA:H	2.25	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	310/331 (94%)	299 (96%)	10 (3%)	1 (0%)	41	70
1	B	301/331 (91%)	285 (95%)	15 (5%)	1 (0%)	41	70
All	All	611/662 (92%)	584 (96%)	25 (4%)	2 (0%)	41	70

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	156	ASP

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Mol	Chain	Res	Type
1	A	11	ASN

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	279/295 (95%)	264 (95%)	15 (5%)	22	51
1	B	276/295 (94%)	260 (94%)	16 (6%)	20	48
All	All	555/590 (94%)	524 (94%)	31 (6%)	21	49

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	60	GLN
1	A	91	ARG
1	A	96	ARG
1	A	101	ASN
1	A	102	SER
1	A	131	THR
1	A	133	GLU
1	A	136	LYS
1	A	185	LYS
1	A	198	ASN
1	A	207	LYS
1	A	228	VAL
1	A	303	LEU
1	A	310	SER
1	B	31	ARG
1	B	60	GLN
1	B	76	ARG
1	B	127	GLU
1	B	128	LEU
1	B	131	THR
1	B	136	LYS

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Mol	Chain	Res	Type
1	B	137	VAL
1	B	185	LYS
1	B	198	ASN
1	B	207	LYS
1	B	225	LYS
1	B	228	VAL
1	B	237	GLU
1	B	246	LYS
1	B	263	ARG

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

Mol	Chain	Res	Type
1	A	6	GLN
1	A	14	HIS
1	A	60	GLN
1	A	134	ASN
1	A	198	ASN
1	A	279	GLN
1	A	282	GLN
1	B	60	GLN
1	B	198	ASN
1	B	279	GLN
1	B	282	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 carbohydrates 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	WDV	B	402	-	28,29,29	0.77	0	40,42,42	1.60	6 (15%)
3	WDV	A	402	-	28,29,29	0.81	0	40,42,42	1.48	7 (17%)
2	NAP	A	401	-	45,52,52	1.41	5 (11%)	56,80,80	1.49	9 (16%)
2	NAP	B	401	-	45,52,52	1.21	4 (8%)	56,80,80	1.46	10 (17%)

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
3	WDV	B	402	-	-	2/16/28/28	0/3/3/3
3	WDV	A	402	-	-	2/16/28/28	0/3/3/3
2	NAP	A	401	-	-	3/31/67/67	0/5/5/5
2	NAP	B	401	-	-	4/31/67/67	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAP	C3N-C7N	-3.90	1.44	1.50
2	B	401	NAP	C3N-C7N	-2.51	1.46	1.50
2	B	401	NAP	P2B-O2B	2.34	1.63	1.59
2	A	401	NAP	C2D-C1D	-2.31	1.50	1.53
2	A	401	NAP	O4D-C4D	-2.15	1.40	1.45
2	B	401	NAP	C2N-N1N	-2.13	1.32	1.35
2	A	401	NAP	C2N-N1N	-2.12	1.32	1.35
2	A	401	NAP	C8A-N7A	-2.07	1.31	1.34
2	B	401	NAP	C2D-C1D	-2.05	1.50	1.53

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	WDV	C24-C21-N25	-5.17	115.49	119.38
2	B	401	NAP	C3N-C7N-N7N	-4.60	112.22	117.75
2	A	401	NAP	C3N-C7N-N7N	-4.57	112.27	117.75
3	B	402	WDV	C19-C21-N25	4.03	122.30	118.75
2	A	401	NAP	C2N-C3N-C4N	3.61	122.35	118.26
3	A	402	WDV	C24-C21-N25	-3.52	116.72	119.38
3	A	402	WDV	C3-C1-N2	-3.44	117.64	121.33
3	A	402	WDV	C6-N2-C5	3.33	118.86	111.52
2	B	401	NAP	C3D-C2D-C1D	3.15	105.72	100.98
2	B	401	NAP	O2N-PN-O1N	2.76	125.86	112.24
2	B	401	NAP	O2A-PA-O1A	2.63	125.25	112.24
3	B	402	WDV	C6-N2-C5	2.53	117.11	111.52
2	A	401	NAP	C3N-C2N-N1N	-2.52	117.96	120.43
2	B	401	NAP	O2D-C2D-C3D	-2.51	103.72	111.82
2	A	401	NAP	O7N-C7N-C3N	2.47	122.59	119.63
2	B	401	NAP	C2N-C3N-C4N	2.46	121.05	118.26
3	B	402	WDV	C5-C9-N13	-2.41	105.70	110.64
2	B	401	NAP	O7N-C7N-C3N	2.29	122.37	119.63
3	A	402	WDV	C6-C10-N13	-2.25	106.02	110.64
2	A	401	NAP	O4D-C4D-C3D	-2.19	100.79	105.11
3	A	402	WDV	C24-C21-C19	2.15	122.83	120.09
2	A	401	NAP	C2D-C3D-C4D	-2.15	98.46	102.64
2	B	401	NAP	PN-O3-PA	2.15	140.21	132.83
2	A	401	NAP	C5N-C4N-C3N	-2.14	117.81	120.34
3	A	402	WDV	C10-C6-N2	-2.14	106.55	110.70
2	A	401	NAP	C3D-C2D-C1D	2.12	104.17	100.98
2	A	401	NAP	O2A-PA-O1A	2.11	122.66	112.24
3	A	402	WDV	C5-N2-C1	-2.11	112.40	118.09
2	B	401	NAP	C3N-C2N-N1N	-2.10	118.38	120.43
2	B	401	NAP	C2D-C3D-C4D	-2.06	98.64	102.64
3	B	402	WDV	C18-C17-N13	2.02	116.28	112.75
3	B	402	WDV	C5-N2-C1	-2.01	112.66	118.09

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	402	WDV	C18-C17-N13-C9
3	A	402	WDV	C18-C17-N13-C9
2	A	401	NAP	O4D-C4D-C5D-O5D
2	B	401	NAP	C2B-O2B-P2B-O3X

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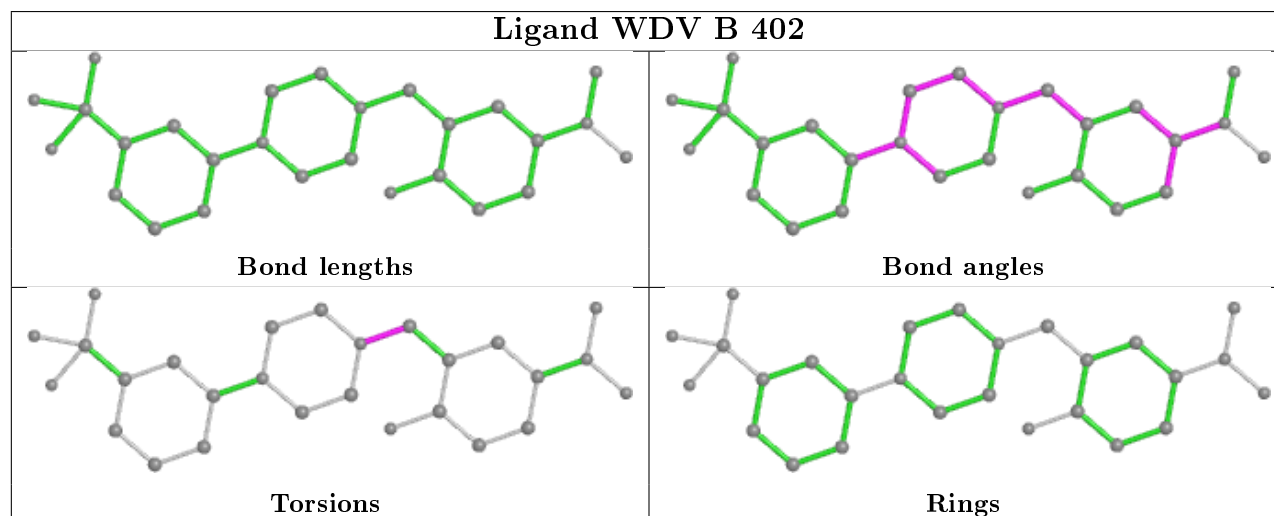
Mol	Chain	Res	Type	Atoms
3	A	402	WDV	C18-C17-N13-C10
3	B	402	WDV	C18-C17-N13-C10
2	A	401	NAP	C3D-C4D-C5D-O5D
2	B	401	NAP	O4D-C4D-C5D-O5D
2	A	401	NAP	PA-O3-PN-O5D
2	B	401	NAP	PA-O3-PN-O5D
2	B	401	NAP	C3D-C4D-C5D-O5D

There are no ring outliers.

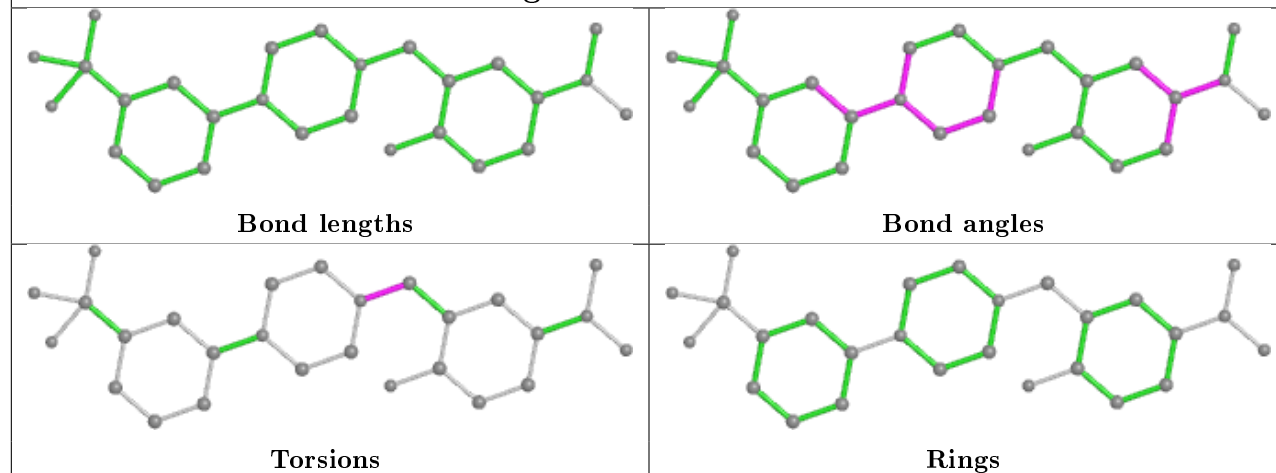
4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	WDV	1	0
3	A	402	WDV	2	0
2	A	401	NAP	5	0
2	B	401	NAP	5	0

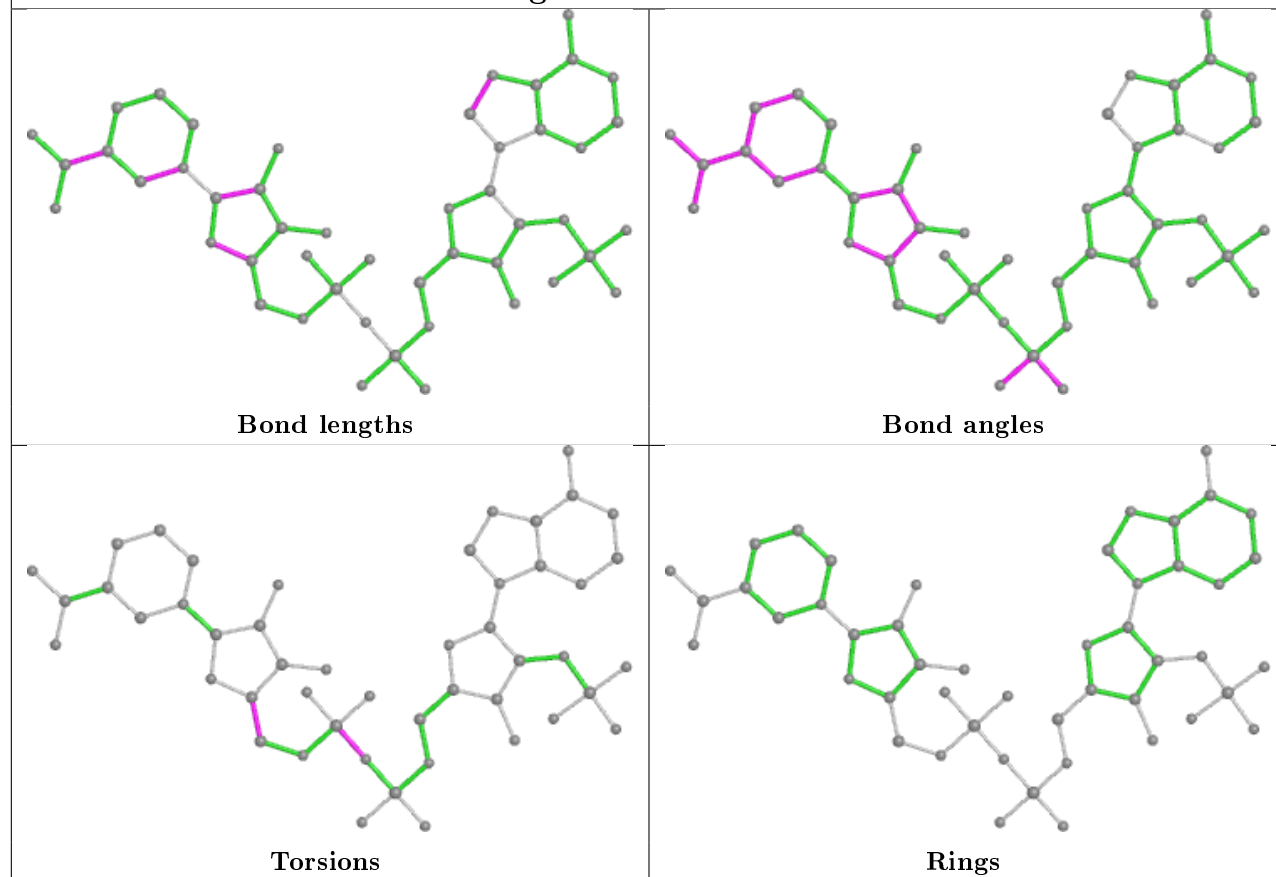
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

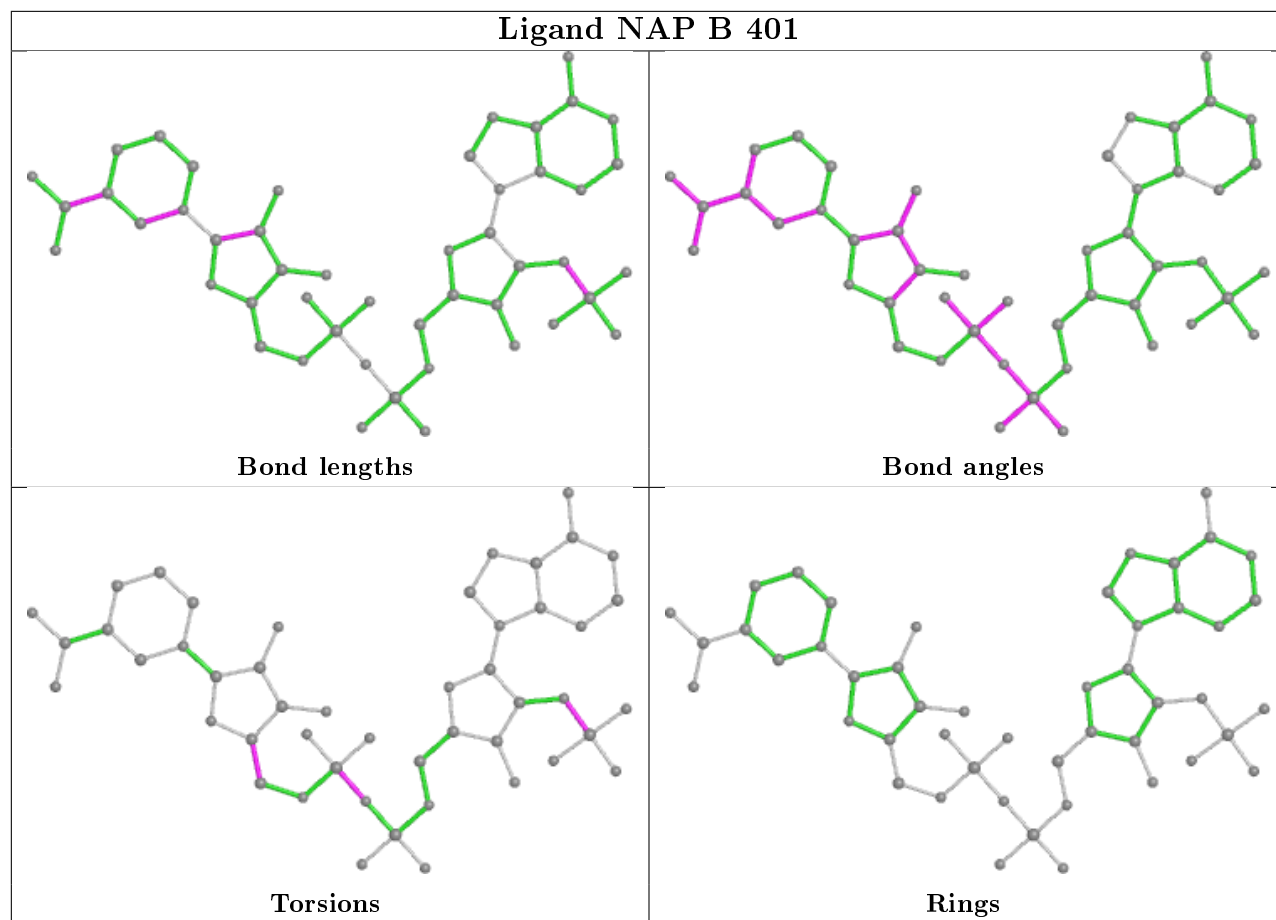


Ligand WDV A 402



Ligand NAP A 401





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	314/331 (94%)	-0.37	2 (0%) 89 86	11, 22, 44, 71	12 (3%)
1	B	311/331 (93%)	0.27	26 (8%) 11 5	15, 40, 70, 105	12 (3%)
All	All	625/662 (94%)	-0.05	28 (4%) 33 23	11, 28, 65, 105	24 (3%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	32	SER	7.2
1	B	67	SER	6.0
1	B	320	SER	4.7
1	B	70	ALA	4.5
1	B	36	GLU	3.8
1	B	33	LYS	3.7
1	B	63	LEU	3.7
1	B	321	ASP	3.6
1	B	27	PRO	3.3
1	B	74	VAL	3.2
1	B	31	ARG	2.7
1	A	133	GLU	2.6
1	B	66	ARG	2.6
1	B	77	GLU	2.5
1	B	237	GLU	2.5
1	B	35	LEU	2.5
1	B	107	GLN	2.5
1	B	278	ARG	2.3
1	B	302	ASN	2.2
1	B	69	ILE	2.1
1	B	73	SER	2.1
1	B	45	GLY	2.1
1	B	75	LYS	2.1
1	B	71	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	78	ASP	2.1
1	B	68	LYS	2.1
1	A	309	ASP	2.1
1	B	277	ILE	2.1

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 carbohydrates in this entry.

6.4 Ligands [i](#)

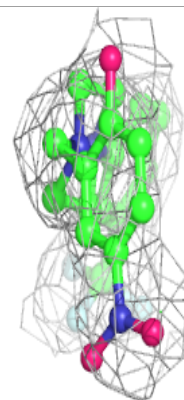
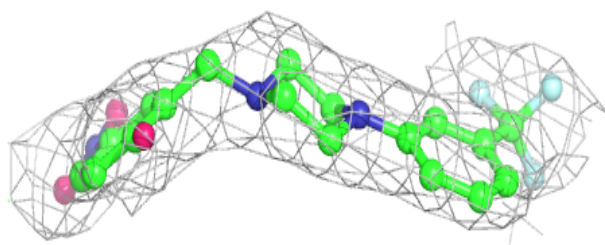
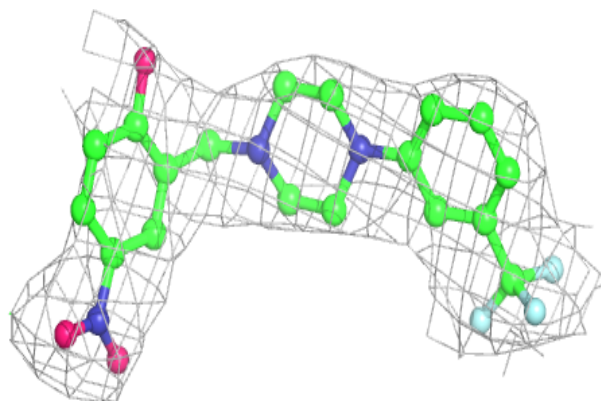
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	WDV	B	402	27/27	0.92	0.20	34,37,53,61	0
2	NAP	B	401	48/48	0.92	0.17	30,42,71,75	0
3	WDV	A	402	27/27	0.97	0.14	16,20,25,28	0
2	NAP	A	401	48/48	0.98	0.13	17,21,24,25	0

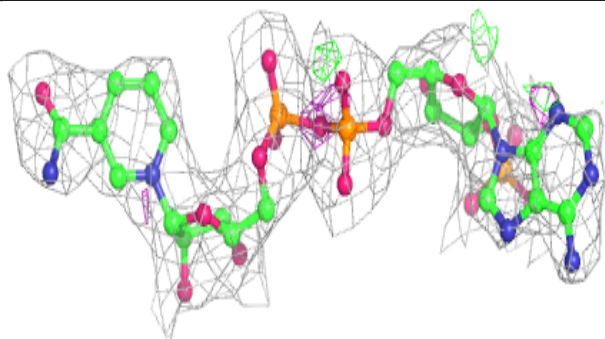
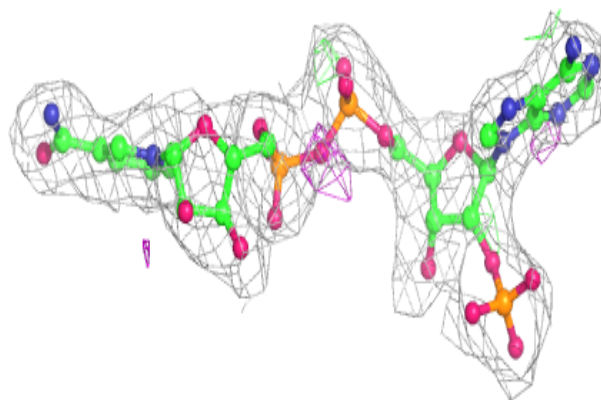
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around WDV B 402:

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

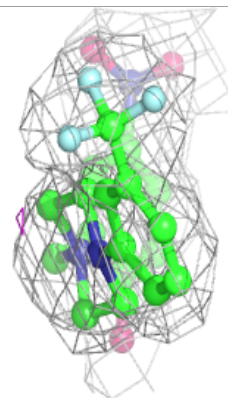
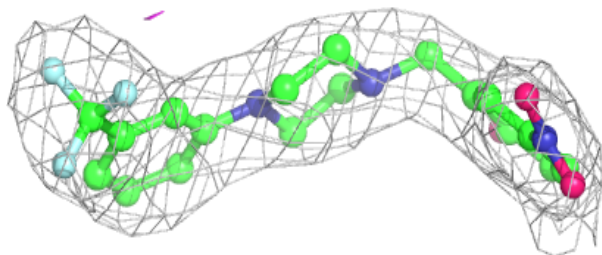
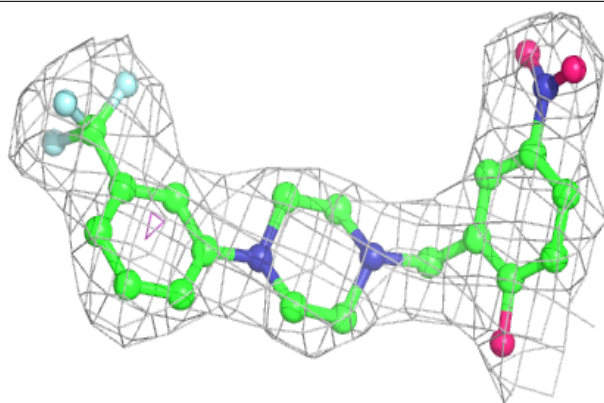
**Electron density around NAP B 401:**

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

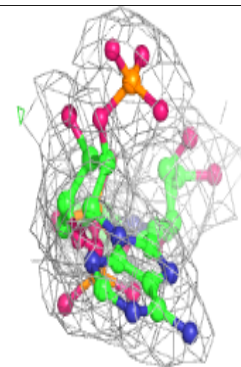
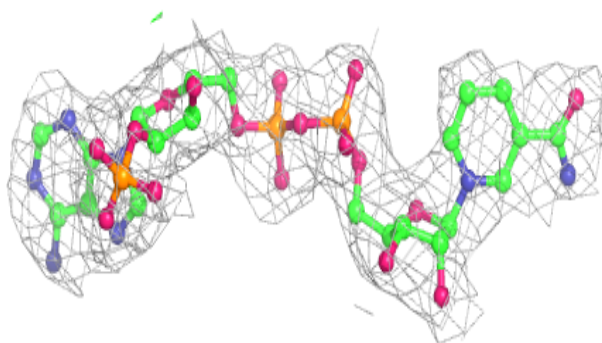
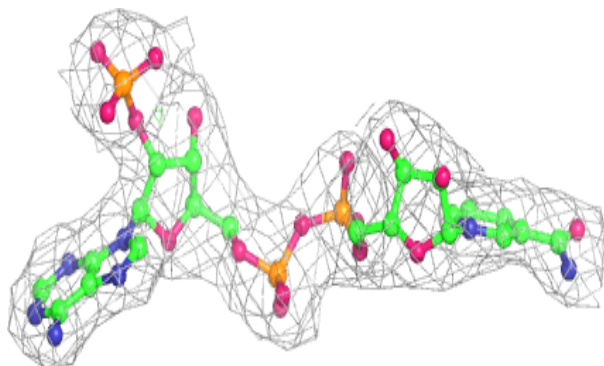


Electron density around WDV A 402:

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

**Electron density around NAP A 401:**

$2mF_o-DF_c$ (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.