



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 03:33 PM EDT

PDB ID : 8OOD / pdb_00008ood
Title : Crystal structure of human DCAF1 WD40 repeats (Q1250L) in complex with compound 15
Authors : Schroeder, M.; Renatus, M.
Deposited on : 2023-04-05
Resolution : 1.50 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

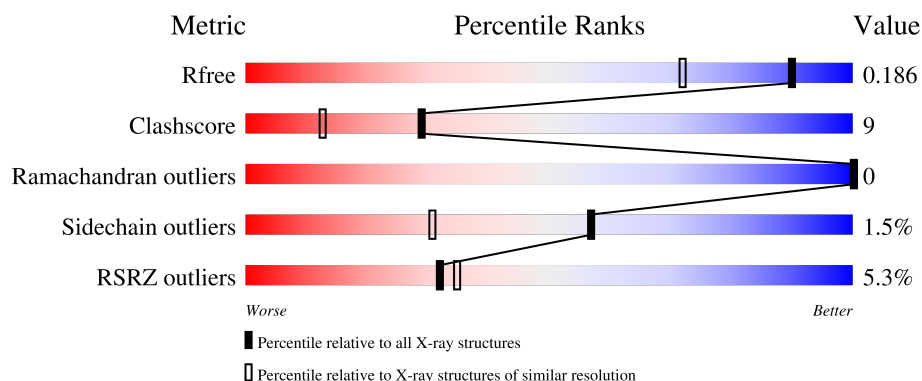
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	367	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	EDO	A	1505	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5579 atoms, of which 2557 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 DDB1- and CUL4-associated factor 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	300	Total	C	H	N	O	S	90	22	0
			5091	1627	2521	441	480	22			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1035	GLY	-	expression tag	UNP Q9Y4B6
A	1036	GLY	-	expression tag	UNP Q9Y4B6
A	1037	GLY	-	expression tag	UNP Q9Y4B6
A	1038	ARG	-	expression tag	UNP Q9Y4B6
A	1250	LEU	GLN	engineered mutation	UNP Q9Y4B6

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			7	2	3	2		
3	A	1	Total	C	H	O	0	0
			7	2	3	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



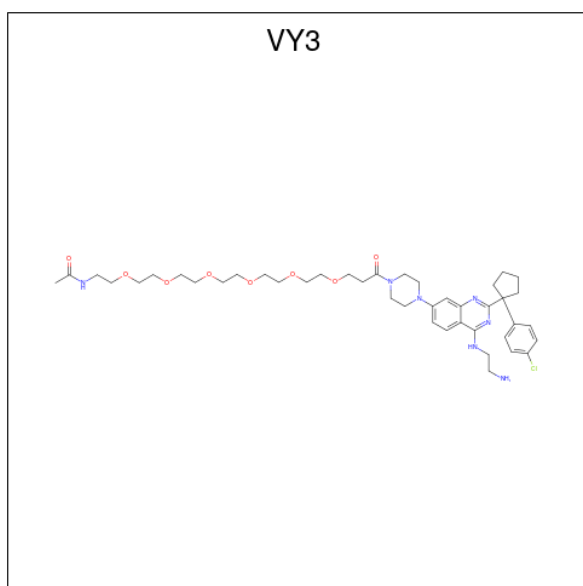
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	1	0
			10	2	6	2		
4	A	1	Total	C	H	O	1	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	1	0
			10	2	6	2		
4	A	1	Total	C	H	O	1	0
			10	2	6	2		
4	A	1	Total	C	H	O	1	0
			10	2	6	2		

- Molecule 5 is {N}-[2-[2-[2-[2-[2-[3-[4-[4-(2-azanylethylamino)-2-[1-(4-chlorophenyl)cyclopentyl]quinazolin-7-yl]piperazin-1-yl]-3-oxidanylidene-propoxy]ethoxy]ethoxy]ethoxy]ethoxy]ethoxy]ethyl]ethanamide (CCD ID: VY3) (formula: C₄₂H₆₂ClN₇O₈) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	Cl	N	O	0	0
			43	32	1	6	4		
5	A	1	Total	C	Cl	N		0	1
			64	50	2	12			

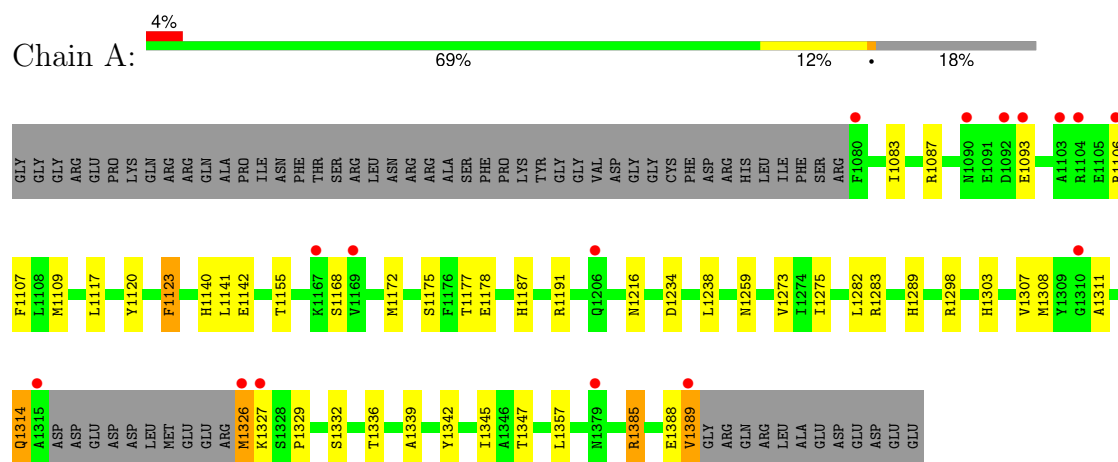
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	316	Total	O	0	0
			316	316		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: DDB1- and CUL4-associated factor 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	81.50Å 81.50Å 232.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	52.26 – 1.50 52.26 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (52.26-1.50) 99.5 (52.26-1.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.170 , 0.186 0.170 , 0.186	Depositor DCC
R_{free} test set	3672 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5579	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: VY3, EDO, MG, ACT

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	1.26	8/2627 (0.3%)	1.33	9/3558 (0.3%)

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	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1303	HIS	CE1-NE2	7.30	1.39	1.32
1	A	1093	GLU	C-O	-6.71	1.16	1.24
1	A	1311	ALA	CA-CB	-6.02	1.44	1.53
1	A	1107	PHE	N-CA	5.78	1.53	1.46
1	A	1216	ASN	C-O	-5.51	1.20	1.25
1	A	1106	ARG	C-O	5.47	1.34	1.24
1	A	1289	HIS	ND1-CE1	5.25	1.37	1.32
1	A	1187	HIS	CE1-NE2	5.04	1.37	1.32

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1329	PRO	N-CA-CB	-7.40	95.68	103.52
1	A	1385	ARG	CG-CD-NE	-6.97	96.66	112.00
1	A	1120	TYR	O-C-N	6.40	131.11	123.24
1	A	1347	THR	CA-CB-OG1	-5.58	101.22	109.60
1	A	1234	ASP	CA-CB-CG	5.57	118.17	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1123	PHE	CA-CB-CG	-5.37	108.43	113.80
1	A	1389	VAL	CA-C-O	-5.25	111.88	120.80
1	A	1106	ARG	O-C-N	5.17	127.30	121.87
1	A	1314	GLN	CB-CA-C	5.11	117.50	109.84

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1178	GLU	Sidechain

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	2570	2521	2487	41	0
2	A	1	0	0	0	0
3	A	8	6	6	0	0
4	A	20	30	29	10	0
5	A	107	0	0	1	0
6	A	316	0	0	17	4
All	All	3022	2557	2522	46	4

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 9.

All (46) 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:1238[B]:LEU:HD11	1:A:1273[B]:VAL:CG2	1.89	1.00
1:A:1238[B]:LEU:HD11	1:A:1273[B]:VAL:HG23	1.51	0.93
4:A:1508:EDO:H11	6:A:1860:HOH:O	1.77	0.85
1:A:1238[B]:LEU:HD11	1:A:1273[B]:VAL:HG21	1.59	0.83
1:A:1140:HIS:HE2	4:A:1505:EDO:H11	1.44	0.81
1:A:1172[A]:MET:HE2	1:A:1175:SER:HB3	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1283[B]:ARG:NE	6:A:1601:HOH:O	2.08	0.78
1:A:1140:HIS:HE2	4:A:1505:EDO:C1	1.98	0.76
1:A:1177[B]:THR:HG23	6:A:1832:HOH:O	1.91	0.71
4:A:1506:EDO:H12	6:A:1612:HOH:O	1.95	0.67
1:A:1140:HIS:NE2	4:A:1505:EDO:H11	2.12	0.64
1:A:1332:SER:H	4:A:1508:EDO:H12	1.64	0.62
1:A:1191[B]:ARG:CZ	6:A:1620:HOH:O	2.49	0.60
1:A:1238[B]:LEU:CD1	1:A:1273[B]:VAL:CG2	2.73	0.58
1:A:1238[B]:LEU:CD1	1:A:1273[B]:VAL:HG23	2.28	0.57
1:A:1336[B]:THR:HG21	6:A:1773:HOH:O	2.06	0.55
1:A:1307:VAL:HG13	1:A:1336[B]:THR:HG23	1.89	0.55
1:A:1298[A]:ARG:NH2	6:A:1608:HOH:O	2.38	0.54
1:A:1238[A]:LEU:HG	6:A:1679:HOH:O	2.07	0.54
1:A:1307:VAL:CG1	1:A:1336[B]:THR:HG23	2.40	0.52
1:A:1155[B]:THR:HG23	6:A:1890:HOH:O	2.09	0.51
1:A:1326[A]:MET:O	6:A:1602:HOH:O	2.14	0.51
1:A:1308:MET:CE	1:A:1339:ALA:HA	2.41	0.50
1:A:1177[B]:THR:HG22	6:A:1725:HOH:O	2.12	0.49
1:A:1336[B]:THR:HG22	1:A:1345:ILE:HB	1.95	0.49
1:A:1273[A]:VAL:HG21	1:A:1282:LEU:HD21	1.95	0.47
1:A:1273[A]:VAL:HG12	1:A:1275:ILE:HG13	1.97	0.46
4:A:1505:EDO:C2	6:A:1604:HOH:O	2.62	0.46
1:A:1273[A]:VAL:CG2	1:A:1282:LEU:HD21	2.46	0.46
1:A:1083[A]:ILE:HG13	1:A:1388:GLU:HB2	1.97	0.46
1:A:1385:ARG:NE	6:A:1612:HOH:O	2.47	0.45
1:A:1109[A]:MET:HE2	1:A:1117[A]:LEU:HD11	1.98	0.45
4:A:1505:EDO:O2	6:A:1604:HOH:O	2.21	0.45
4:A:1505:EDO:H11	6:A:1604:HOH:O	2.15	0.45
1:A:1326[A]:MET:HB3	1:A:1326[A]:MET:HE2	1.66	0.44
1:A:1357:LEU:C	1:A:1357:LEU:HD12	2.43	0.43
1:A:1238[B]:LEU:CD1	1:A:1273[B]:VAL:HG21	2.39	0.43
1:A:1191[B]:ARG:NH1	6:A:1620:HOH:O	2.52	0.43
1:A:1117[B]:LEU:CD1	1:A:1141:LEU:HD22	2.49	0.42
1:A:1087[B]:ARG:HG2	6:A:1621:HOH:O	2.21	0.41
1:A:1109[B]:MET:HE3	1:A:1109[B]:MET:HB3	1.94	0.41
1:A:1142:GLU:CD	4:A:1505:EDO:H12	2.45	0.41
1:A:1083[A]:ILE:CG2	1:A:1123:PHE:HE1	2.34	0.40
1:A:1117[A]:LEU:HD23	5:A:1510[A]:VY3:CL55	2.59	0.40
1:A:1308:MET:HE1	1:A:1342:TYR:CE1	2.56	0.40
1:A:1314:GLN:HG3	1:A:1327:LYS:HE3	2.03	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:1900:HOH:O	6:A:1900:HOH:O[12_545]	0.94	1.26
6:A:1665:HOH:O	6:A:1665:HOH:O[12_545]	0.96	1.24
6:A:1822:HOH:O	6:A:1822:HOH:O[12_555]	1.02	1.18
6:A:1786:HOH:O	6:A:1803:HOH:O[12_555]	2.13	0.07

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	318/367 (87%)	310 (98%)	8 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	288/321 (90%)	283 (98%)	5 (2%)	53	26

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1168	SER
1	A	1259	ASN
1	A	1326[A]	MET

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Mol	Chain	Res	Type
1	A	1326[B]	MET
1	A	1389	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1090	ASN
1	A	1353	ASN

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 ⓘ

Of 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 for Mogul analysis.

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	$\# Z > 2$
5	VY3	A	1509	-	47,47,62	1.71	10 (21%)	53,64,80	2.11	17 (32%)
4	EDO	A	1508	-	3,3,3	0.10	0	2,2,2	0.41	0
3	ACT	A	1502	-	3,3,3	0.92	0	3,3,3	0.76	0
4	EDO	A	1504	-	3,3,3	0.08	0	2,2,2	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	VY3	A	1510[A]	-	36,36,62	1.31	3 (8%)	41,51,80	3.29	18 (43%)
5	VY3	A	1510[B]	-	36,36,62	1.11	2 (5%)	41,51,80	3.09	18 (43%)
4	EDO	A	1505	-	3,3,3	0.96	0	2,2,2	0.54	0
4	EDO	A	1506	-	3,3,3	0.39	0	2,2,2	0.42	0
3	ACT	A	1503	-	3,3,3	1.90	1 (33%)	3,3,3	0.73	0
4	EDO	A	1507	-	3,3,3	0.25	0	2,2,2	0.11	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
5	VY3	A	1509	-	-	7/29/52/67	0/5/5/5
4	EDO	A	1508	-	-	1/1/1/1	-
4	EDO	A	1504	-	-	0/1/1/1	-
5	VY3	A	1510[A]	-	-	4/16/37/67	0/5/5/5
5	VY3	A	1510[B]	-	-	3/16/37/67	0/5/5/5
4	EDO	A	1505	-	-	1/1/1/1	-
4	EDO	A	1506	-	-	0/1/1/1	-
4	EDO	A	1507	-	-	0/1/1/1	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1509	VY3	C12-C13	-6.08	1.45	1.54
5	A	1509	VY3	C11-N10	4.12	1.38	1.33
3	A	1503	ACT	OXT-C	-3.10	1.16	1.30
5	A	1509	VY3	C5-C4	-2.87	1.38	1.42
5	A	1510[A]	VY3	C11-N8	2.72	1.37	1.33
5	A	1509	VY3	C21-N22	-2.61	1.42	1.47
5	A	1509	VY3	C23-N22	2.50	1.51	1.47
5	A	1510[A]	VY3	C11-N10	2.45	1.36	1.33
5	A	1509	VY3	C7-N8	2.43	1.36	1.33
5	A	1509	VY3	C54-C12	2.23	1.59	1.56
5	A	1509	VY3	C7-C5	-2.23	1.42	1.44
5	A	1509	VY3	C6-C1	2.20	1.41	1.36
5	A	1509	VY3	C24-N19	2.13	1.50	1.46
5	A	1510[B]	VY3	C11-N8	2.12	1.36	1.33
5	A	1510[B]	VY3	C11-N10	2.09	1.35	1.33
5	A	1510[A]	VY3	C20-N19	2.02	1.50	1.46

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1510[A]	VY3	C54-C12-C13	-11.18	95.56	113.30
5	A	1510[B]	VY3	C54-C12-C13	-9.12	98.83	113.30
5	A	1510[A]	VY3	C51-C12-C13	8.80	127.27	113.30
5	A	1510[B]	VY3	C54-C12-C51	-7.43	88.26	102.40
5	A	1510[B]	VY3	C51-C12-C13	7.05	124.49	113.30
5	A	1510[B]	VY3	C5-C7-N9	6.29	127.02	120.76
5	A	1510[A]	VY3	C54-C12-C51	-6.28	90.44	102.40
5	A	1510[A]	VY3	C5-C7-N9	5.61	126.35	120.76
5	A	1509	VY3	C14-C15-C16	-4.90	114.31	119.24
5	A	1509	VY3	C7-C5-C4	4.90	118.75	115.86
5	A	1509	VY3	C3-C4-C5	4.70	124.83	119.58
5	A	1510[B]	VY3	C18-C13-C12	-4.56	112.14	121.45
5	A	1510[A]	VY3	C18-C13-C12	-4.43	112.39	121.45
5	A	1509	VY3	C54-C12-C13	-4.16	106.70	113.30
5	A	1510[A]	VY3	C11-N8-C7	4.06	125.30	114.54
5	A	1510[A]	VY3	C11-N10-C4	3.97	119.37	115.47
5	A	1510[A]	VY3	C51-C52-C53	-3.84	95.49	105.78
5	A	1510[B]	VY3	C11-N8-C7	3.80	124.61	114.54
5	A	1510[B]	VY3	C51-C52-C53	-3.75	95.74	105.78
5	A	1510[A]	VY3	C17-C16-CL55	-3.69	113.91	119.36
5	A	1509	VY3	C11-N10-C4	3.68	119.09	115.47
5	A	1509	VY3	C23-C24-N19	-3.64	103.13	110.78
5	A	1510[B]	VY3	C17-C16-C15	3.60	125.70	121.24
5	A	1509	VY3	C14-C13-C12	-3.56	114.18	121.45
5	A	1509	VY3	C18-C13-C12	3.47	128.53	121.45
5	A	1510[A]	VY3	C14-C15-C16	-3.42	115.80	119.24
5	A	1510[A]	VY3	C23-N22-C21	3.40	120.00	110.40
5	A	1509	VY3	C5-C4-N10	-3.34	119.27	122.80
5	A	1509	VY3	O32-C31-C29	-3.24	99.33	109.53
5	A	1510[B]	VY3	C23-N22-C21	3.05	119.01	110.40
5	A	1510[B]	VY3	C11-N10-C4	3.03	118.44	115.47
5	A	1510[B]	VY3	C5-C7-N8	-2.98	117.61	122.07
5	A	1510[A]	VY3	C5-C7-N8	-2.94	117.66	122.07
5	A	1509	VY3	C3-C2-N19	-2.92	118.80	122.11
5	A	1510[B]	VY3	C24-N19-C20	2.77	117.79	111.57
5	A	1510[A]	VY3	C24-N19-C20	2.73	117.70	111.57
5	A	1510[B]	VY3	C14-C15-C16	-2.72	116.50	119.24
5	A	1509	VY3	C2-C3-C4	-2.72	117.08	120.57
5	A	1510[B]	VY3	C1-C2-N19	-2.71	117.62	121.39
5	A	1509	VY3	C17-C16-C15	2.69	124.57	121.24
5	A	1510[A]	VY3	C17-C16-C15	2.68	124.56	121.24
5	A	1510[B]	VY3	C18-C17-C16	-2.62	116.60	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1510[B]	VY3	C7-C5-C4	2.60	117.40	115.86
5	A	1510[B]	VY3	C18-C13-C14	2.60	121.91	118.03
5	A	1509	VY3	C24-C23-N22	-2.55	105.35	110.42
5	A	1509	VY3	C20-N19-C2	-2.37	111.66	118.11
5	A	1510[A]	VY3	C6-C1-C2	-2.31	117.21	120.70
5	A	1509	VY3	C15-C14-C13	2.24	124.25	121.17
5	A	1510[A]	VY3	C14-C13-C12	2.20	125.94	121.45
5	A	1510[A]	VY3	C3-C2-N19	-2.18	119.64	122.11
5	A	1509	VY3	C3-C4-N10	-2.07	115.75	118.78
5	A	1510[A]	VY3	C1-C2-N19	-2.02	118.58	121.39
5	A	1510[B]	VY3	C14-C13-C12	2.02	125.58	121.45

There are no chirality outliers.

All (16) torsion outliers are listed below:

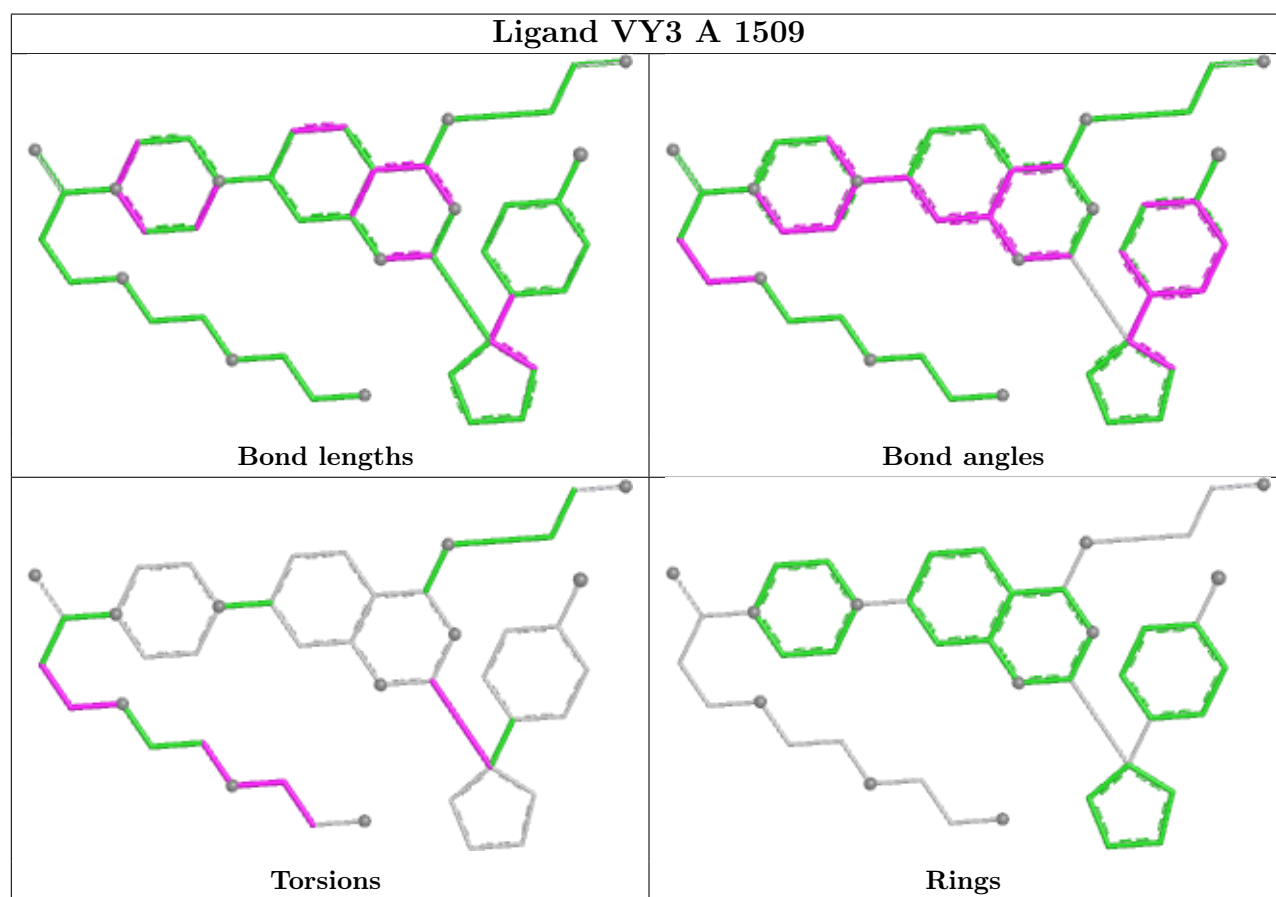
Mol	Chain	Res	Type	Atoms
5	A	1509	VY3	N10-C11-C12-C13
5	A	1509	VY3	C28-C29-C31-O32
5	A	1510[B]	VY3	N9-C25-C26-N27
5	A	1510[B]	VY3	C5-C7-N9-C25
5	A	1510[B]	VY3	N8-C7-N9-C25
5	A	1509	VY3	O35-C36-C37-O1
5	A	1510[A]	VY3	N8-C7-N9-C25
4	A	1508	EDO	O1-C1-C2-O2
5	A	1510[A]	VY3	N9-C25-C26-N27
5	A	1510[A]	VY3	C5-C7-N9-C25
5	A	1509	VY3	C29-C31-O32-C33
5	A	1509	VY3	C37-C36-O35-C34
5	A	1509	VY3	C33-C34-O35-C36
4	A	1505	EDO	O1-C1-C2-O2
5	A	1509	VY3	N8-C11-C12-C13
5	A	1510[A]	VY3	C26-C25-N9-C7

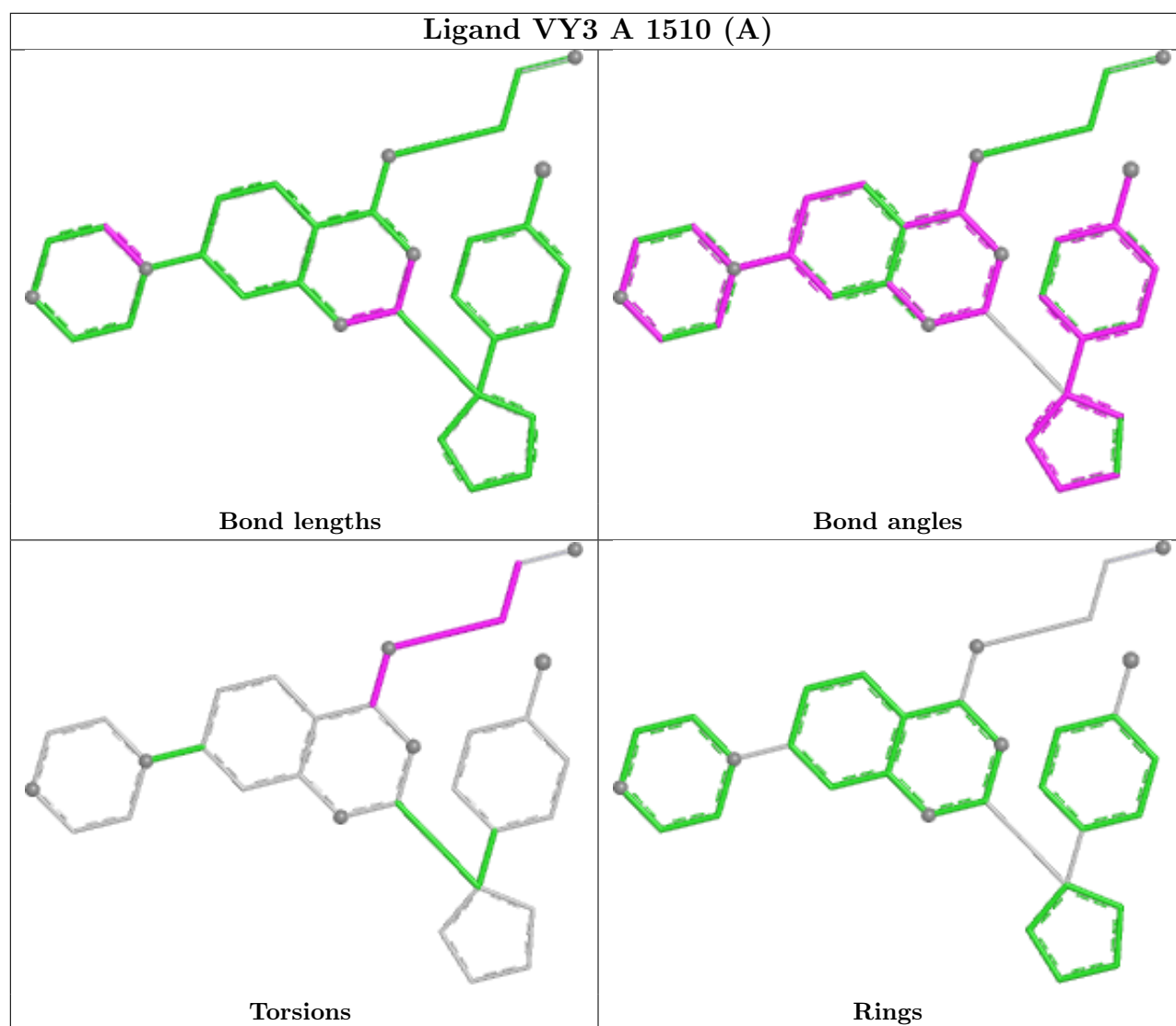
There are no ring outliers.

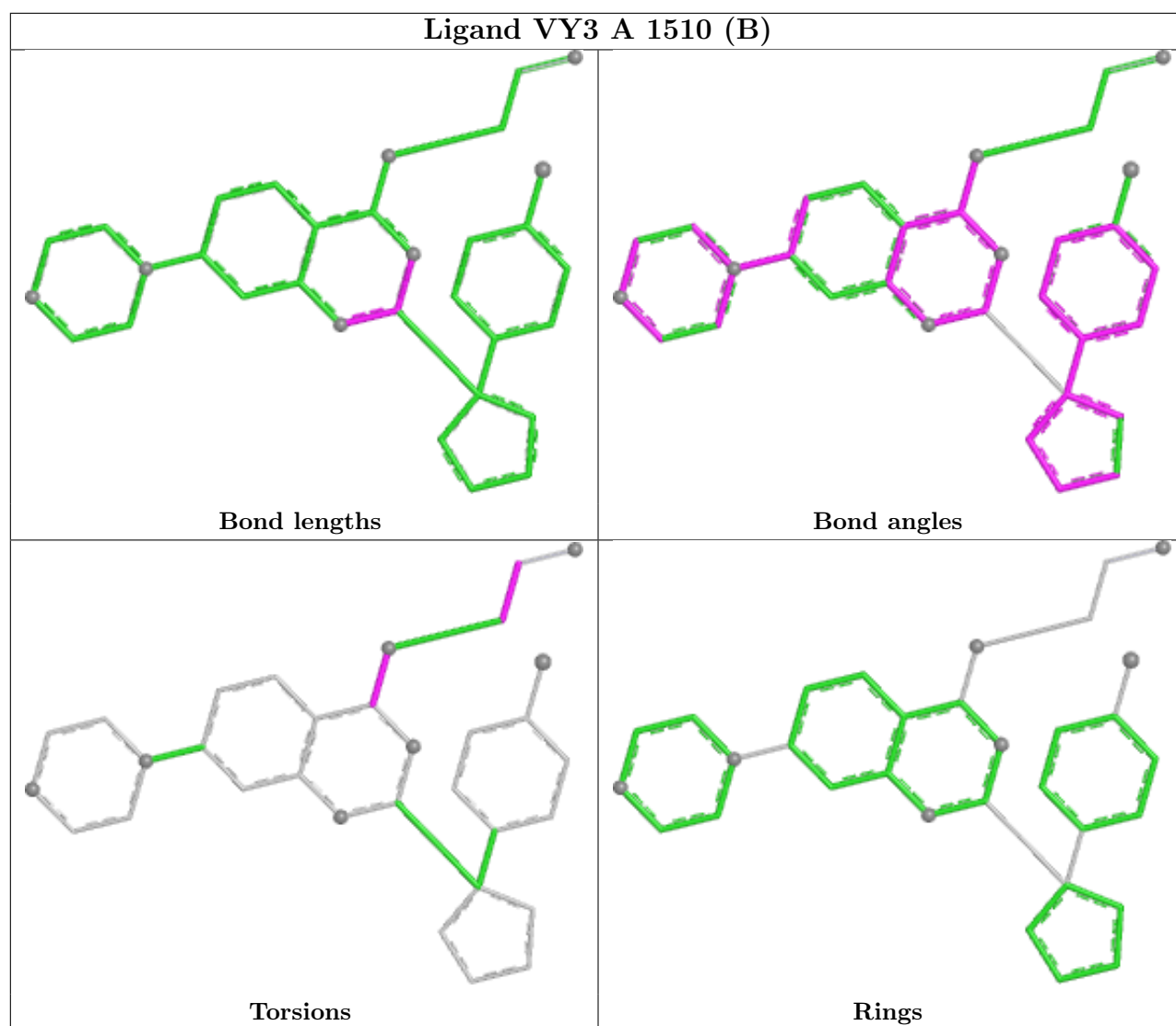
4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1508	EDO	2	0
5	A	1510[A]	VY3	1	0
4	A	1505	EDO	7	0
4	A	1506	EDO	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.







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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	300/367 (81%)	-0.16	16 (5%) 32 35	5, 18, 39, 71	22 (7%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1315	ALA	6.0
1	A	1326[A]	MET	4.3
1	A	1090	ASN	4.1
1	A	1080	PHE	3.9
1	A	1389	VAL	3.8
1	A	1103	ALA	3.8
1	A	1092	ASP	3.1
1	A	1093	GLU	2.9
1	A	1106	ARG	2.8
1	A	1167	LYS	2.7
1	A	1310	GLY	2.5
1	A	1327	LYS	2.4
1	A	1104	ARG	2.3
1	A	1169	VAL	2.2
1	A	1206	GLN	2.1
1	A	1379	ASN	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

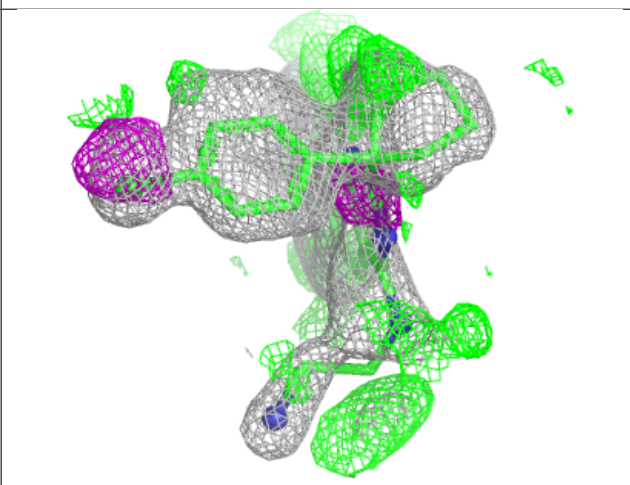
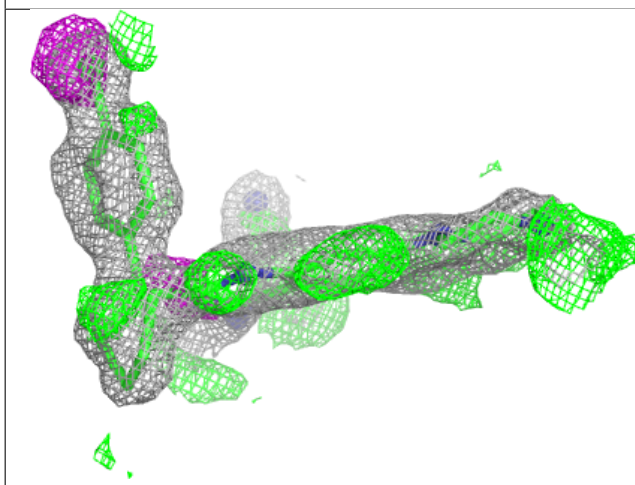
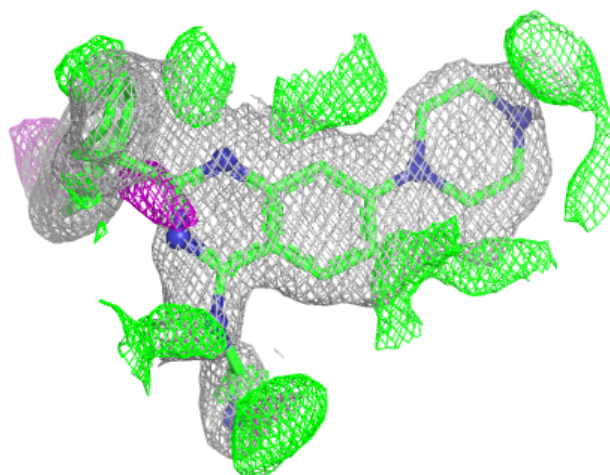
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EDO	A	1508	4/4	0.81	0.17	37,45,47,50	1
4	EDO	A	1506	4/4	0.84	0.17	37,45,46,48	1
5	VY3	A	1510[A]	32/58	0.86	0.20	20,36,44,47	32
5	VY3	A	1510[B]	32/58	0.86	0.20	26,36,51,55	32
4	EDO	A	1507	4/4	0.87	0.14	35,40,42,44	1
2	MG	A	1501	1/1	0.91	0.41	43,43,43,43	0
4	EDO	A	1505	4/4	0.93	0.09	22,26,37,44	1
3	ACT	A	1503	4/4	0.96	0.12	21,23,24,32	0
4	EDO	A	1504	4/4	0.97	0.08	15,19,22,37	1
5	VY3	A	1509	43/58	0.98	0.09	11,15,64,68	0
3	ACT	A	1502	4/4	0.99	0.04	12,12,13,13	7

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.

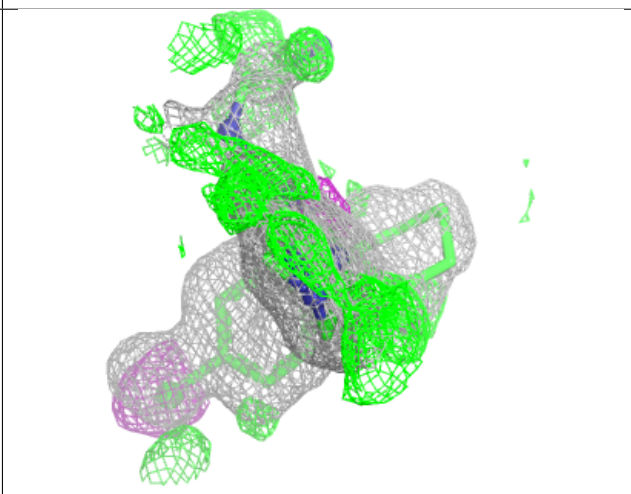
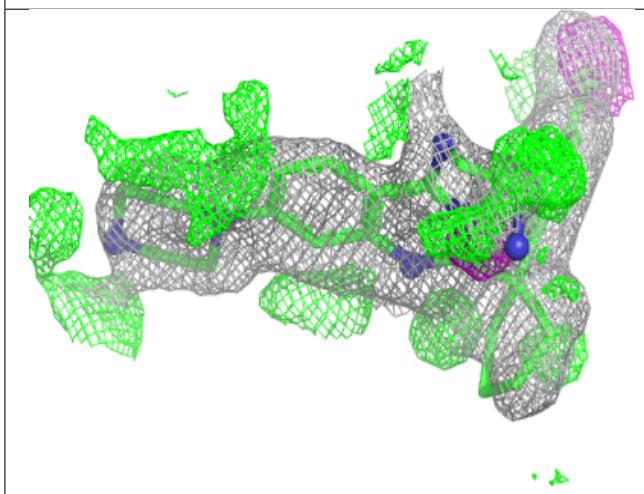
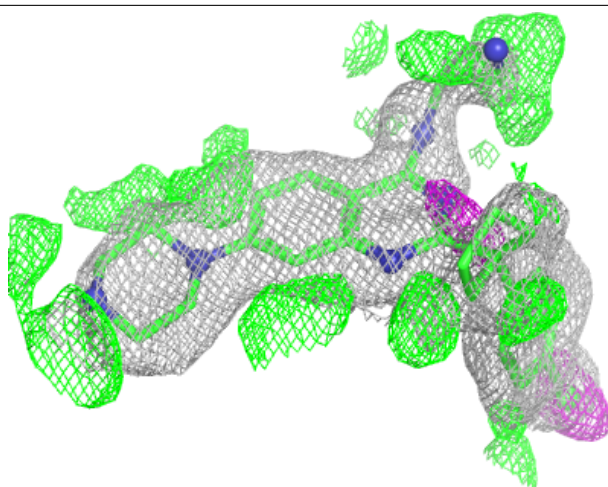
Electron density around VY3 A 1510 (A):

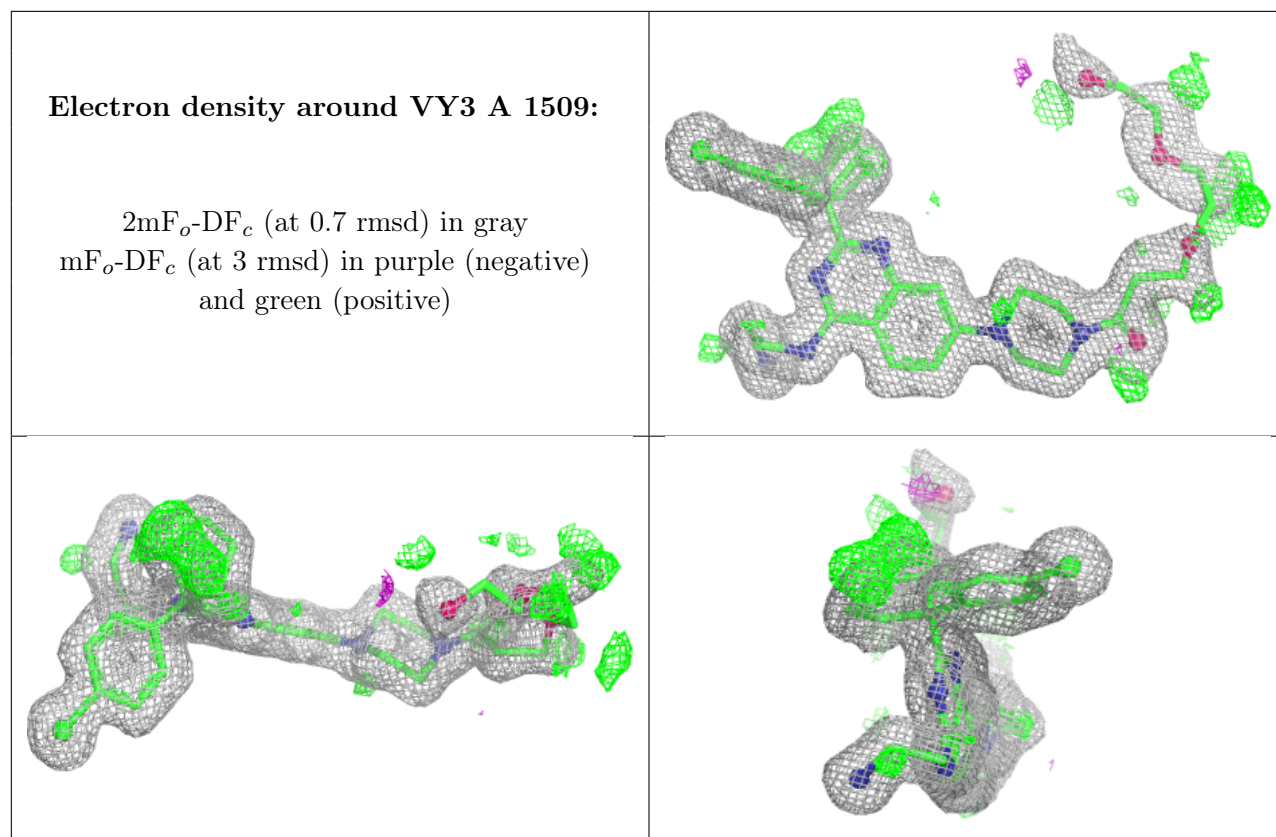
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around VY3 A 1510 (B):

$2mF_o-DF_c$ (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.