



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:18 AM UTC

PDB ID : 3RZE / pdb_00003rze
Title : Structure of the human histamine H1 receptor in complex with doxepin
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Deposited on : 2011-05-11
Resolution : 3.10 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

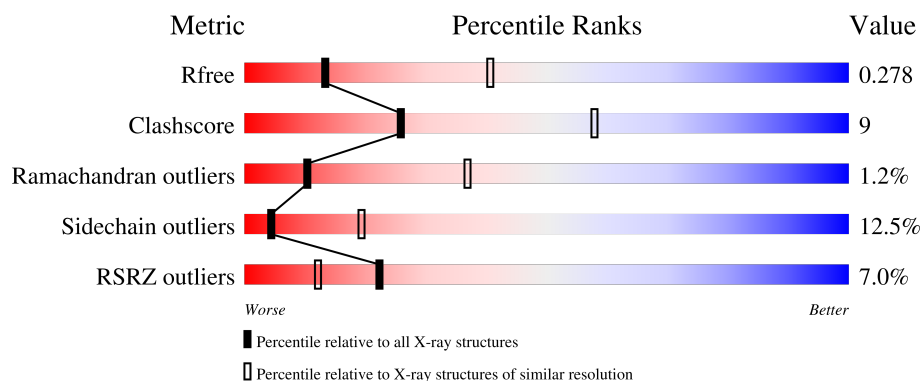
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	452	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3555 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.

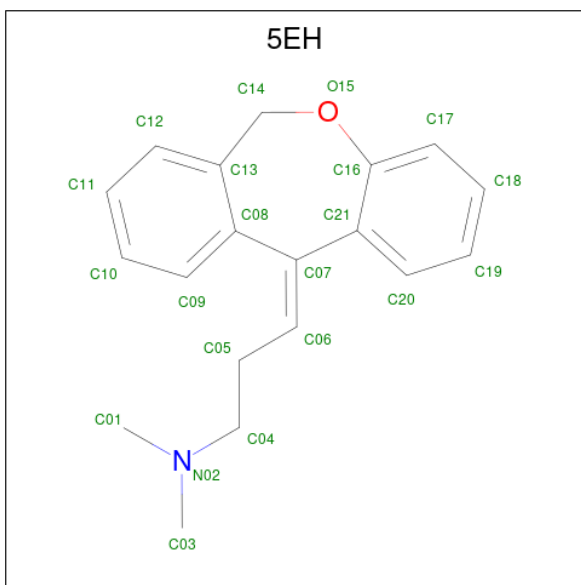
- Molecule 1 is a protein called Histamine H1 receptor, Lysozyme chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	0	0
			3481	2279	591	586	25			

There are 9 discrepancies between the modelled and reference sequences:

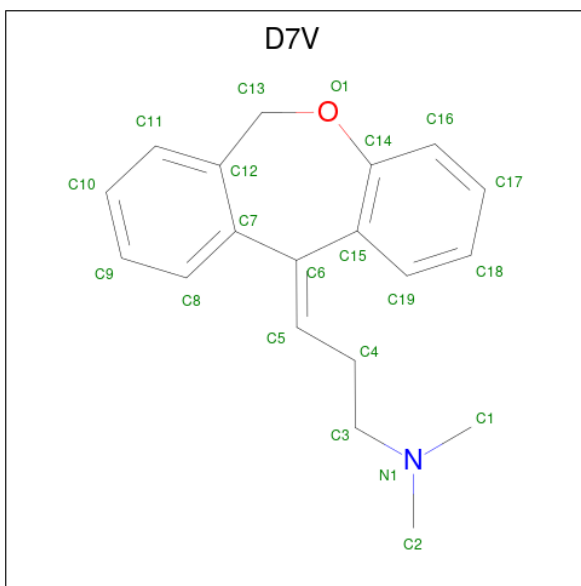
Chain	Residue	Modelled	Actual	Comment	Reference
A	1054	THR	CYS	engineered mutation	UNP P00720
A	1097	ALA	CYS	engineered mutation	UNP P00720
A	488	GLY	-	expression tag	UNP P35367
A	489	GLU	-	expression tag	UNP P35367
A	490	ASN	-	expression tag	UNP P35367
A	491	LEU	-	expression tag	UNP P35367
A	492	TYR	-	expression tag	UNP P35367
A	493	PHE	-	expression tag	UNP P35367
A	494	GLN	-	expression tag	UNP P35367

- Molecule 2 is (3E)-3-(dibenzo[b,e]oxepin-11(6H)-ylidene)-N,N-dimethylpropan-1-amine (CCD ID: 5EH) (formula: C₁₉H₂₁NO).



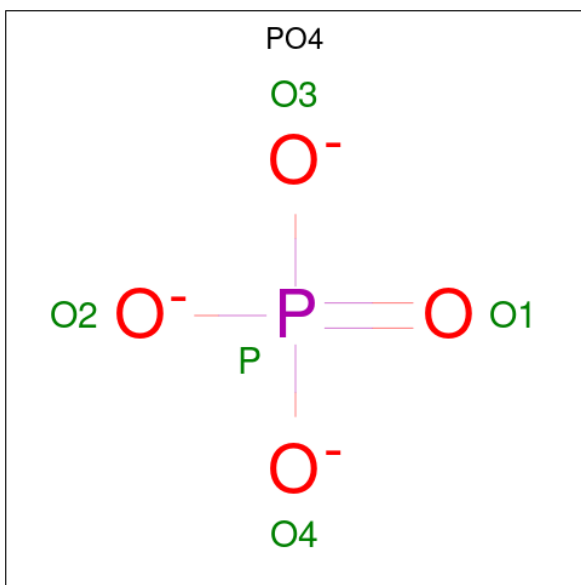
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	1
			21	19	1	1		

- Molecule 3 is (3Z)-3-(dibenzo[b,e]oxepin-11(6H)-ylidene)-N,N-dimethylpropan-1-amine (CCD ID: D7V) (formula: C₁₉H₂₁NO).



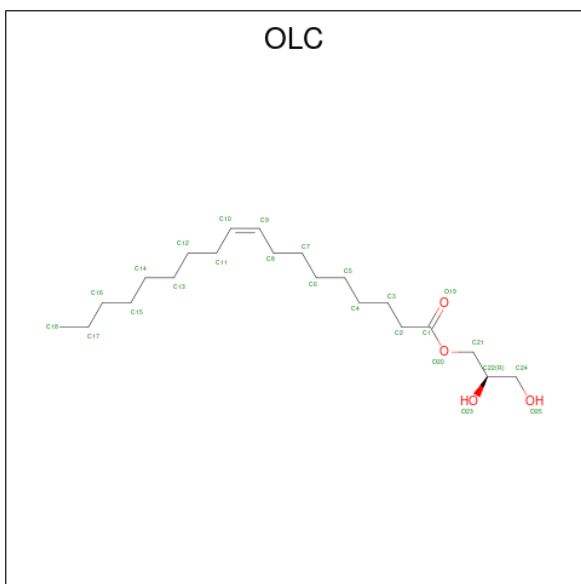
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			21	19	1	1		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 5	O 4	P 1	0	0
4	A	1	Total 5	O 4	P 1	0	0
4	A	1	Total 5	O 4	P 1	0	0

- Molecule 5 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLC) (formula: $C_{21}H_{40}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			15	11	4		

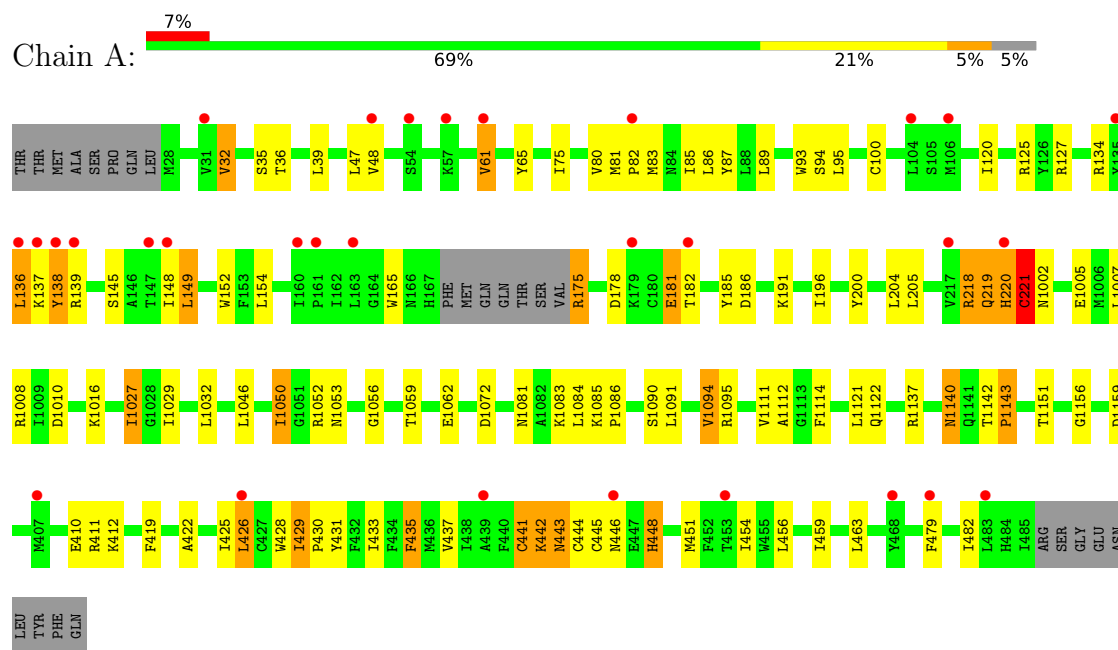
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	O	0	0
			2	2		

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: Histamine H1 receptor, Lysozyme chimera



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	88.14Å 88.14Å 331.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.50 – 3.10 34.50 – 3.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (34.50-3.10) 96.4 (34.50-3.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 3.12Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, R_{free}	0.214 , 0.249 0.231 , 0.278	Depositor DCC
R_{free} test set	797 reflections (6.64%)	wwPDB-VP
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.138	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 76.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	3555	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.79% 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: OLC, PO4, 5EH, D7V

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	0.87	3/3566 (0.1%)	1.46	27/4835 (0.6%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1002	ASN	C-O	-6.79	1.17	1.24
1	A	448	HIS	ND1-CE1	5.48	1.38	1.32
1	A	220	HIS	CD2-NE2	-5.20	1.32	1.37

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	437	VAL	N-CA-C	8.87	119.42	110.82
1	A	95	LEU	N-CA-C	8.79	120.79	111.03
1	A	448	HIS	CB-CG-CD2	-7.74	121.14	131.20
1	A	220	HIS	CB-CG-CD2	-7.03	122.06	131.20
1	A	138	TYR	N-CA-C	6.33	121.14	113.41
1	A	149	LEU	N-CA-C	-6.14	105.44	113.12
1	A	220	HIS	CB-CG-ND1	5.96	131.63	122.70
1	A	1142	THR	CB-CA-C	5.90	118.01	110.34
1	A	437	VAL	CB-CA-C	-5.86	104.37	112.04
1	A	437	VAL	CA-C-N	5.77	128.61	120.42
1	A	437	VAL	C-N-CA	5.77	128.61	120.42
1	A	1010	ASP	N-CA-C	5.68	118.68	111.69
1	A	1008	ARG	N-CA-C	-5.55	105.14	111.14
1	A	1072	ASP	N-CA-C	-5.52	105.16	111.07
1	A	444	CYS	N-CA-C	5.44	119.39	112.87
1	A	80	VAL	N-CA-C	-5.42	105.56	110.82
1	A	1002	ASN	N-CA-C	-5.36	100.54	107.73
1	A	1140	ASN	CA-CB-CG	5.36	117.95	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1005	GLU	CA-C-N	5.25	127.58	120.65
1	A	1005	GLU	C-N-CA	5.25	127.58	120.65
1	A	410	GLU	CA-C-N	5.08	127.40	120.54
1	A	410	GLU	C-N-CA	5.08	127.40	120.54
1	A	94	SER	CA-C-N	5.07	127.52	120.63
1	A	94	SER	C-N-CA	5.07	127.52	120.63
1	A	1121	LEU	CA-C-N	5.05	127.36	120.54
1	A	1121	LEU	C-N-CA	5.05	127.36	120.54
1	A	448	HIS	CB-CG-ND1	5.04	130.26	122.70

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	3481	0	3566	68	0
2	A	21	0	21	3	0
3	A	21	0	21	2	0
4	A	15	0	0	1	0
5	A	15	0	19	0	0
6	A	2	0	0	0	0
All	All	3555	0	3627	68	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 9.

All (68) 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:429:ILE:HG22	1:A:430:PRO:CD	1.80	1.11
1:A:429:ILE:HG22	1:A:430:PRO:HD3	1.14	1.07
1:A:442:LYS:O	1:A:443:ASN:HB2	1.52	1.07
1:A:145:SER:HA	1:A:148:ILE:HG12	1.55	0.86
1:A:429:ILE:CG2	1:A:430:PRO:CD	2.60	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:GLN:OE1	1:A:219:GLN:C	2.30	0.75
1:A:1081:ASN:HB3	1:A:1084:LEU:HB2	1.69	0.74
1:A:429:ILE:CG2	1:A:430:PRO:N	2.52	0.72
1:A:429:ILE:CG2	1:A:430:PRO:HD3	2.07	0.72
1:A:220:HIS:O	1:A:221:CYS:O	2.11	0.68
1:A:425:ILE:CG2	1:A:429:ILE:HD13	2.23	0.68
1:A:219:GLN:OE1	1:A:219:GLN:CA	2.42	0.67
1:A:425:ILE:O	1:A:429:ILE:HB	1.95	0.66
1:A:1094:VAL:HG21	1:A:1156:GLY:C	2.19	0.66
1:A:442:LYS:O	1:A:443:ASN:CB	2.36	0.64
1:A:145:SER:HA	1:A:148:ILE:CG1	2.29	0.63
1:A:136:LEU:O	1:A:137:LYS:HB2	1.98	0.63
1:A:220:HIS:NE2	1:A:411:ARG:NH1	2.48	0.61
1:A:441:CYS:O	1:A:442:LYS:CB	2.50	0.59
1:A:196:ILE:HA	1:A:200:TYR:HB2	1.85	0.59
1:A:425:ILE:HG23	1:A:429:ILE:HD13	1.85	0.58
1:A:1053:ASN:HD22	1:A:1053:ASN:C	2.12	0.58
1:A:429:ILE:HG23	1:A:430:PRO:N	2.20	0.57
1:A:185:TYR:HA	1:A:191:LYS:HE2	1.88	0.56
1:A:148:ILE:O	1:A:152:TRP:HD1	1.90	0.55
1:A:1143:PRO:HD2	4:A:1204:PO4:O4	2.06	0.55
1:A:175:ARG:N	1:A:185:TYR:HH	2.05	0.54
1:A:127:ARG:HH22	1:A:139:ARG:HD3	1.73	0.53
1:A:219:GLN:OE1	1:A:219:GLN:HA	2.09	0.53
1:A:1111:VAL:O	1:A:1114:PHE:HB2	2.08	0.53
1:A:425:ILE:HG23	1:A:429:ILE:CD1	2.38	0.53
1:A:65:TYR:HD1	1:A:120:ILE:HG23	1.75	0.51
1:A:1050:ILE:HG22	1:A:1052:ARG:HG2	1.92	0.51
1:A:134:ARG:C	1:A:136:LEU:H	2.20	0.49
1:A:1027:ILE:HG21	1:A:1046:LEU:HD13	1.94	0.49
1:A:431:TYR:HE2	2:A:1200[A]:5EH:H09	1.77	0.49
1:A:1083:LYS:HD3	1:A:1112:ALA:O	2.13	0.48
1:A:219:GLN:OE1	1:A:219:GLN:O	2.30	0.48
1:A:61:VAL:HG11	1:A:139:ARG:HG3	1.95	0.48
1:A:429:ILE:N	1:A:430:PRO:HD2	2.29	0.48
1:A:428:TRP:HH2	2:A:1200[A]:5EH:H01	1.79	0.48
1:A:1091:LEU:HD22	1:A:1095:ARG:HB3	1.97	0.47
1:A:425:ILE:HG22	1:A:429:ILE:HD13	1.97	0.47
1:A:220:HIS:CD2	1:A:411:ARG:NH1	2.84	0.46
1:A:36:THR:HA	1:A:39:LEU:HD12	1.96	0.46
1:A:428:TRP:CZ3	2:A:1200[A]:5EH:H20	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:TYR:CZ	1:A:93:TRP:HB3	2.52	0.45
1:A:220:HIS:CD2	1:A:220:HIS:N	2.83	0.44
1:A:218:ARG:O	1:A:218:ARG:HG2	2.16	0.44
1:A:1016:LYS:HA	1:A:1056:GLY:O	2.16	0.44
1:A:1059:THR:OG1	1:A:1062:GLU:HG3	2.18	0.44
1:A:82:PRO:HA	1:A:85:ILE:HD12	2.01	0.42
1:A:435:PHE:CZ	3:A:1201[B]:D7V:H18	2.55	0.42
1:A:182:THR:HG22	1:A:185:TYR:HB3	2.01	0.42
1:A:65:TYR:CD1	1:A:120:ILE:HG23	2.55	0.42
1:A:456:LEU:HA	1:A:459:ILE:HD12	2.02	0.42
1:A:127:ARG:HH11	1:A:134:ARG:HG3	1.84	0.42
1:A:422:ALA:O	1:A:426:LEU:HB2	2.20	0.42
1:A:1094:VAL:HG21	1:A:1156:GLY:O	2.19	0.41
1:A:431:TYR:OH	3:A:1201[B]:D7V:H3	2.20	0.41
1:A:48:VAL:HG22	1:A:479:PHE:HZ	1.85	0.41
1:A:1094:VAL:CG2	1:A:1095:ARG:N	2.81	0.41
1:A:83:MET:HA	1:A:86:LEU:HD12	2.01	0.41
1:A:1084:LEU:O	1:A:1085:LYS:C	2.63	0.41
1:A:1086:PRO:HB2	1:A:1122:GLN:OE1	2.19	0.41
1:A:32:VAL:HA	1:A:35:SER:HB2	2.03	0.41
1:A:148:ILE:O	1:A:152:TRP:CD1	2.72	0.41
1:A:61:VAL:HG13	1:A:139:ARG:NH1	2.36	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	424/452 (94%)	401 (95%)	18 (4%)	5 (1%)	10 37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	CYS
1	A	443	ASN
1	A	442	LYS
1	A	181	GLU
1	A	1143	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	375/399 (94%)	328 (88%)	47 (12%)	4 19

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	VAL
1	A	47	LEU
1	A	61	VAL
1	A	75	ILE
1	A	81	MET
1	A	89	LEU
1	A	100	CYS
1	A	125	ARG
1	A	136	LEU
1	A	138	TYR
1	A	149	LEU
1	A	154	LEU
1	A	165	TRP
1	A	175	ARG
1	A	178	ASP
1	A	181	GLU
1	A	186	ASP
1	A	204	LEU
1	A	205	LEU
1	A	218	ARG
1	A	219	GLN
1	A	221	CYS

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Mol	Chain	Res	Type
1	A	1007	LEU
1	A	1027	ILE
1	A	1029	ILE
1	A	1032	LEU
1	A	1050	ILE
1	A	1090	SER
1	A	1094	VAL
1	A	1137	ARG
1	A	1140	ASN
1	A	1151	THR
1	A	1159	ASP
1	A	412	LYS
1	A	419	PHE
1	A	426	LEU
1	A	429	ILE
1	A	433	ILE
1	A	435	PHE
1	A	441	CYS
1	A	445	CYS
1	A	446	ASN
1	A	448	HIS
1	A	451	MET
1	A	454	ILE
1	A	463	LEU
1	A	482	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1053	ASN
1	A	1055	ASN
1	A	1081	ASN
1	A	448	HIS
1	A	474	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	D7V	A	1201[B]	-	23,23,23	0.72	0	30,31,31	1.21	2 (6%)
2	5EH	A	1200[A]	-	23,23,23	0.68	0	30,31,31	1.20	2 (6%)
4	PO4	A	1204	-	4,4,4	2.58	1 (25%)	6,6,6	1.24	1 (16%)
4	PO4	A	1203	-	4,4,4	2.15	1 (25%)	6,6,6	0.60	0
4	PO4	A	1202	-	4,4,4	2.06	1 (25%)	6,6,6	0.59	0
5	OLC	A	1205	-	14,14,24	0.67	0	15,15,25	0.91	1 (6%)

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	D7V	A	1201[B]	-	-	3/6/19/19	0/3/3/3
2	5EH	A	1200[A]	-	-	3/6/19/19	0/3/3/3
5	OLC	A	1205	-	-	4/14/14/24	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1204	PO4	P-O1	4.57	1.61	1.50
4	A	1203	PO4	P-O1	3.52	1.58	1.50
4	A	1202	PO4	P-O1	3.38	1.58	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1200[A]	5EH	C14-O15-C16	4.63	125.58	117.68
3	A	1201[B]	D7V	C13-O1-C14	4.50	125.35	117.68
4	A	1204	PO4	O4-P-O3	2.26	114.94	107.91
5	A	1205	OLC	C21-O20-C1	2.15	124.98	117.12
3	A	1201[B]	D7V	O1-C13-C12	-2.11	108.60	111.73
2	A	1200[A]	5EH	C05-C06-C07	-2.01	123.66	127.28

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1200[A]	5EH	N02-C04-C05-C06
3	A	1201[B]	D7V	N1-C3-C4-C5
2	A	1200[A]	5EH	C05-C04-N02-C01
3	A	1201[B]	D7V	C4-C3-N1-C1
2	A	1200[A]	5EH	C05-C04-N02-C03
3	A	1201[B]	D7V	C4-C3-N1-C2
5	A	1205	OLC	C2-C1-O20-C21
5	A	1205	OLC	O19-C1-O20-C21
5	A	1205	OLC	C2-C3-C4-C5
5	A	1205	OLC	C5-C6-C7-C8

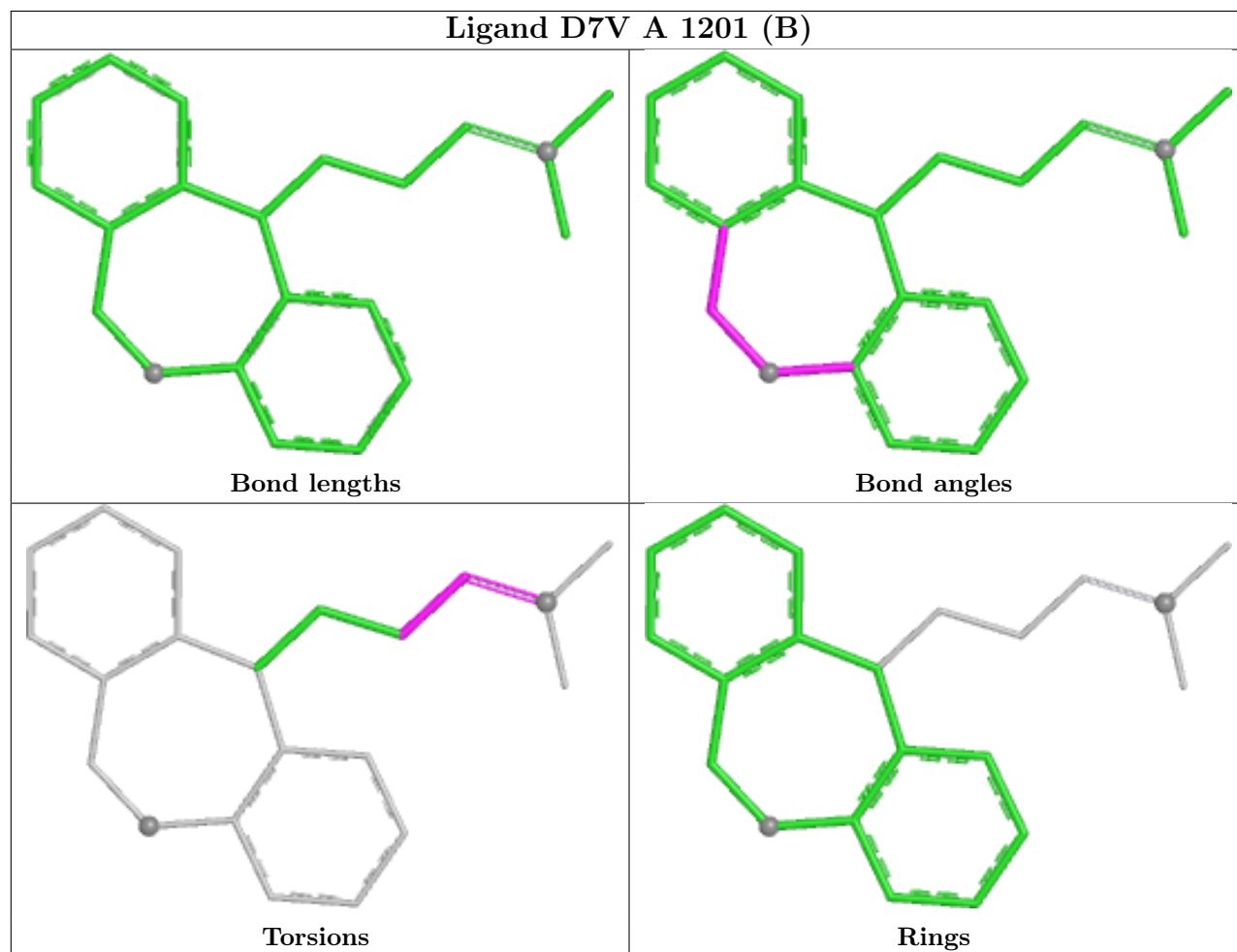
There are no ring outliers.

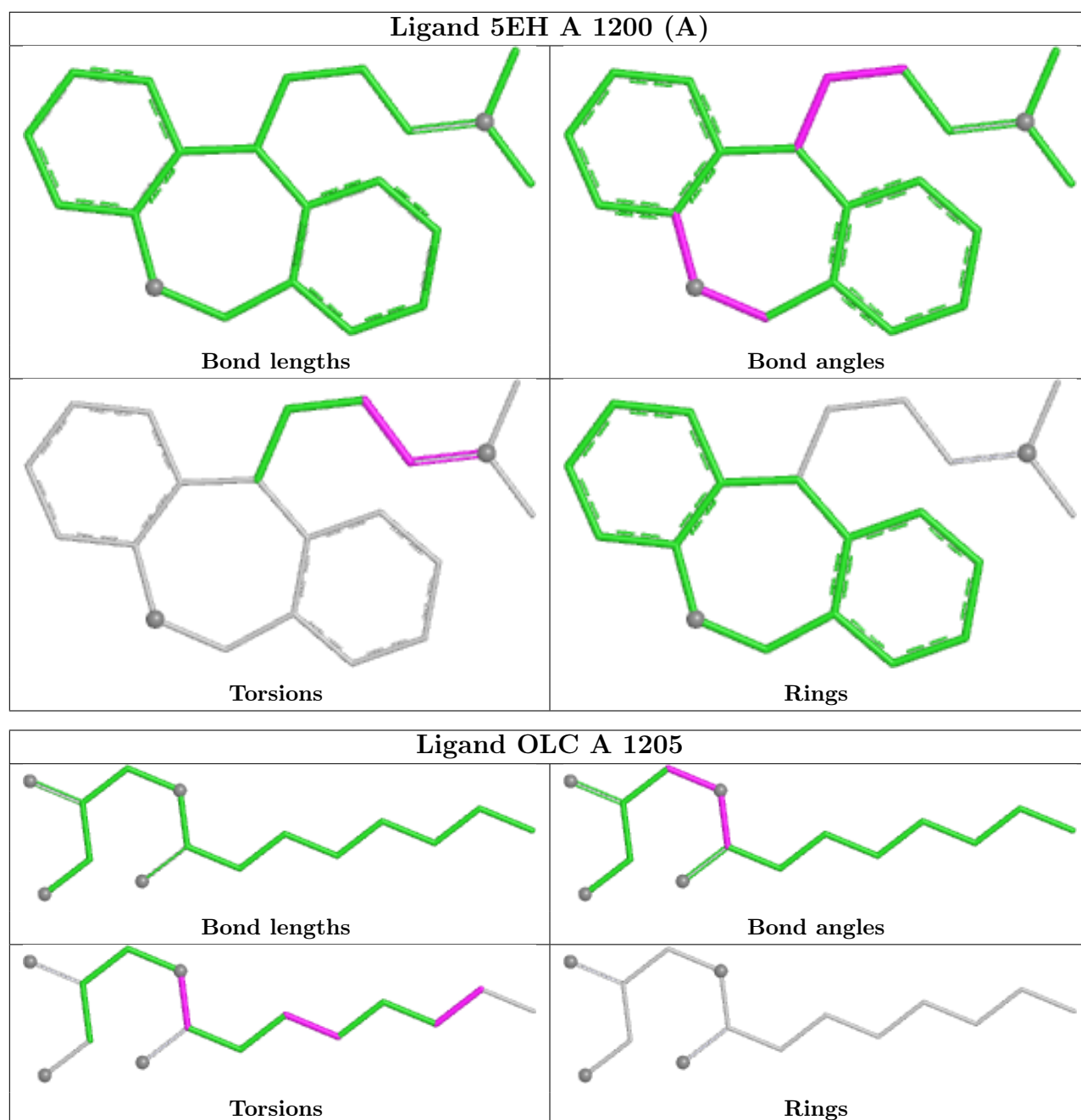
3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1201[B]	D7V	2	0
2	A	1200[A]	5EH	3	0
4	A	1204	PO4	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 ⓘ

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	428/452 (94%)	0.46	30 (7%) 22 12	10, 93, 162, 179	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	106	MET	3.3
1	A	160	ILE	3.1
1	A	137	LYS	3.0
1	A	138	TYR	2.9
1	A	136	LEU	2.8
1	A	217	VAL	2.8
1	A	479	PHE	2.7
1	A	407	MET	2.6
1	A	135	TYR	2.6
1	A	139	ARG	2.6
1	A	439	ALA	2.6
1	A	483	LEU	2.5
1	A	453	THR	2.4
1	A	161	PRO	2.4
1	A	148	ILE	2.4
1	A	182	THR	2.4
1	A	61	VAL	2.4
1	A	446	ASN	2.3
1	A	220	HIS	2.3
1	A	468	TYR	2.3
1	A	163	LEU	2.3
1	A	179	LYS	2.2
1	A	31	VAL	2.2
1	A	82	PRO	2.2
1	A	48	VAL	2.1
1	A	426	LEU	2.1
1	A	54	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	57	LYS	2.1
1	A	104	LEU	2.1
1	A	147	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

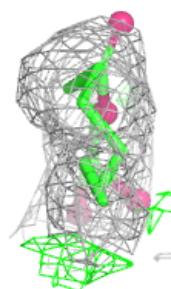
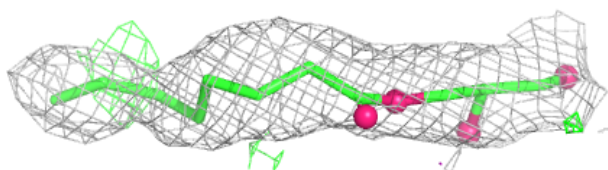
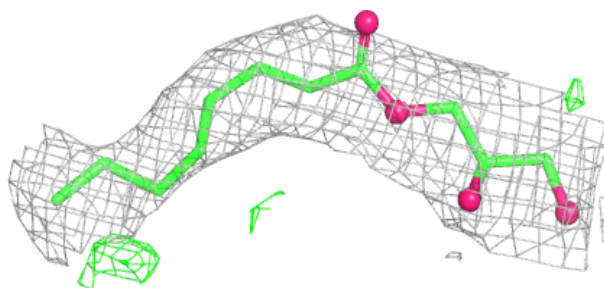
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PO4	A	1202	5/5	0.69	0.17	172,178,178,178	0
4	PO4	A	1203	5/5	0.81	0.20	143,149,150,150	0
5	OLC	A	1205	15/25	0.90	0.18	10,39,208,209	0
3	D7V	A	1201[B]	21/21	0.91	0.14	50,105,151,159	21
2	5EH	A	1200[A]	21/21	0.92	0.15	23,123,153,165	21
4	PO4	A	1204	5/5	0.93	0.15	69,74,75,76	0

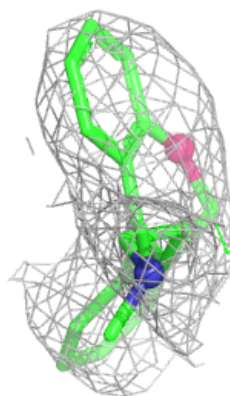
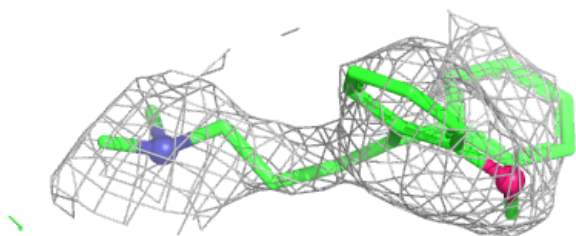
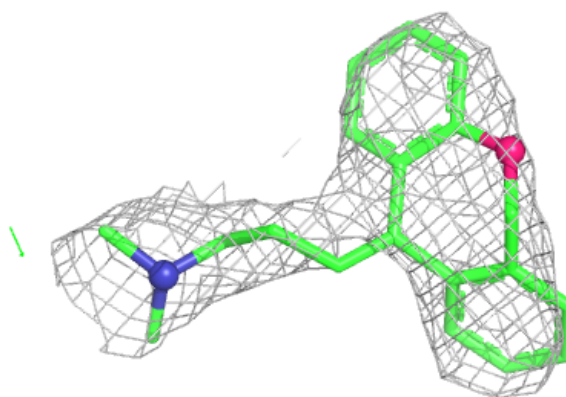
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OLC A 1205:

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

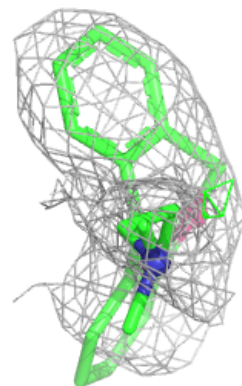
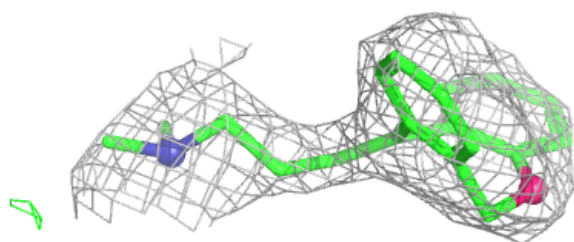
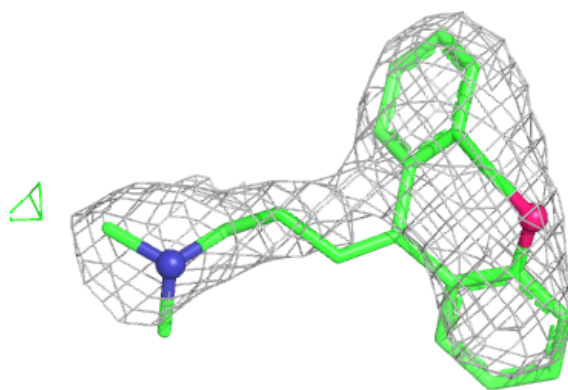
**Electron density around D7V A 1201 (B):**

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



Electron density around 5EH A 1200 (A):

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



6.5 Other polymers ⓘ

There are no such residues in this entry.