



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:33 AM UTC

PDB ID : 8HV9 / pdb_00008hv9
Title : Crystal structure of EGFR_TMX in complex with covalently bound fragment 12
Authors : Dokurno, P.
Deposited on : 2022-12-26
Resolution : 2.50 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

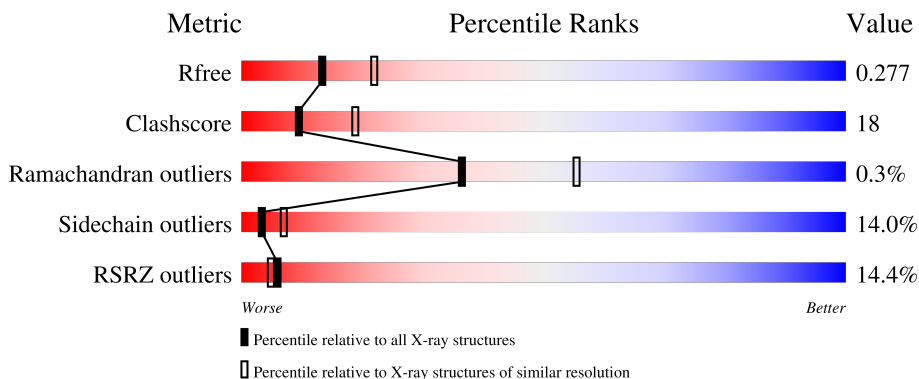
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	331	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2564 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.

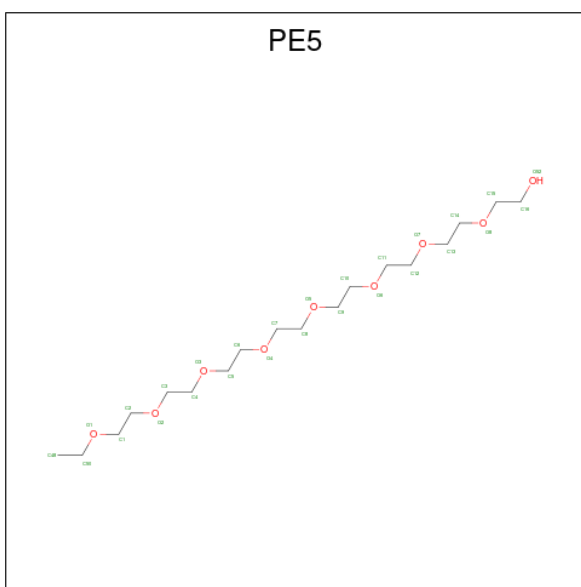
- Molecule 1 is a protein called Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	1	0
			2409	1546	408	438	17			

There are 9 discrepancies between the modelled and reference sequences:

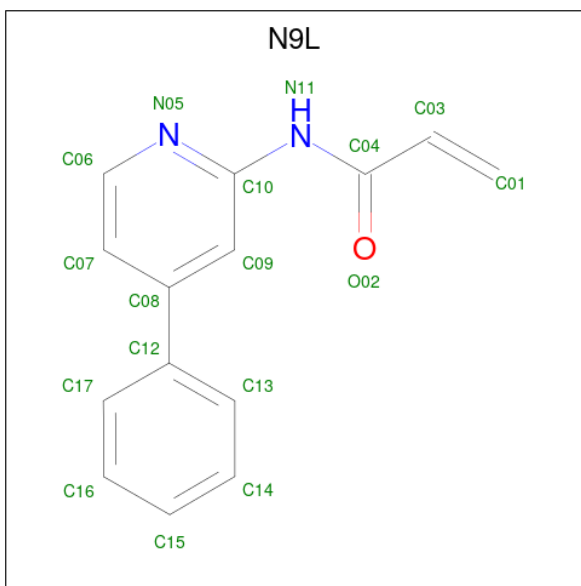
Chain	Residue	Modelled	Actual	Comment	Reference
A	694	GLY	-	expression tag	UNP P00533
A	790	MET	THR	variant	UNP P00533
A	797	SER	CYS	variant	UNP P00533
A	858	ARG	LEU	variant	UNP P00533
A	865	ALA	GLU	engineered mutation	UNP P00533
A	866	ALA	GLU	engineered mutation	UNP P00533
A	867	ALA	LYS	engineered mutation	UNP P00533
A	1023	ASN	-	expression tag	UNP P00533
A	1024	SER	-	expression tag	UNP P00533

- Molecule 2 is 3,6,9,12,15,18,21,24-OCTAOXAHEXACOSAN-1-OL (CCD ID: PE5) (formula: C₁₈H₃₈O₉).



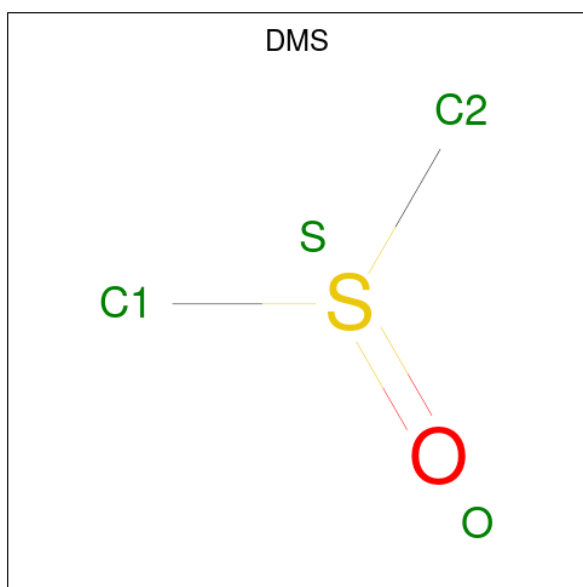
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			26	18	8		

- Molecule 3 is {N}-(4-phenylpyridin-2-yl)prop-2-enamide (CCD ID: N9L) (formula: $C_{14}H_{12}N_2O$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	14	2	1		

- Molecule 4 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			4	2	1	1		

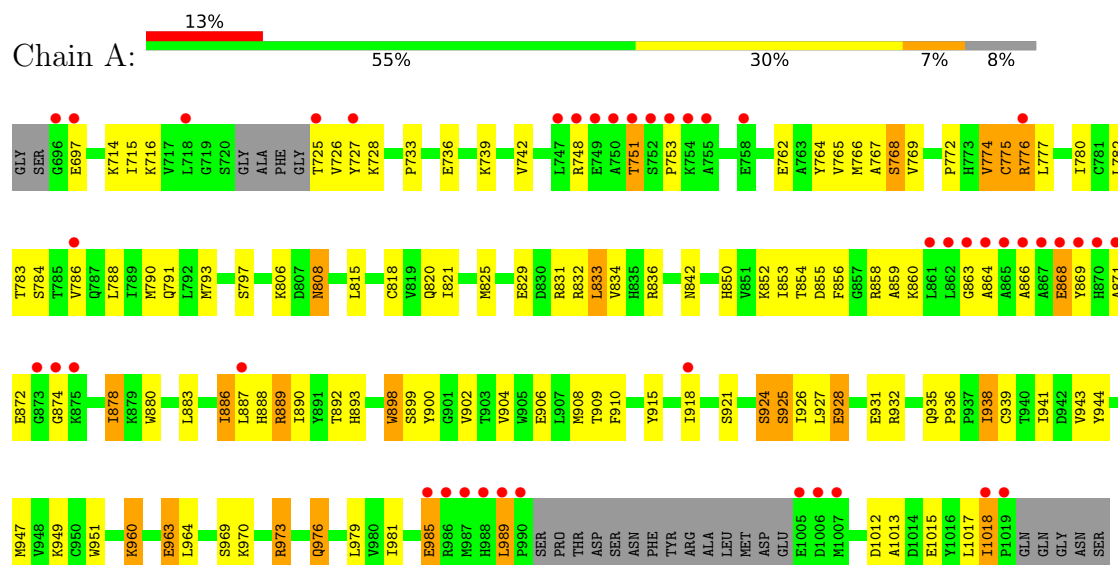
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	108	Total	O	0	0
			108	108		

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: Epidermal growth factor receptor



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	144.44Å 144.44Å 144.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.0 (20.00-2.50) 94.0 (20.00-2.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.253 , 0.274 0.256 , 0.277	Depositor DCC
R_{free} test set	823 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.54$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.008 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2564	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: N9L, PE5, DMS

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	1.05	1/2463 (0.0%)	1.47	9/3340 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	989	LEU	N-CA	5.73	1.50	1.45

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	874	GLY	CA-C-O	-5.60	117.14	122.13
1	A	774	VAL	CA-C-N	-5.25	114.31	122.09
1	A	774	VAL	C-N-CA	-5.25	114.31	122.09
1	A	963	GLU	N-CA-C	-5.24	104.50	112.04
1	A	775	CYS	CA-C-N	5.23	128.28	120.95
1	A	775	CYS	C-N-CA	5.23	128.28	120.95
1	A	775	CYS	N-CA-CB	5.18	118.62	110.23
1	A	898	TRP	N-CA-C	-5.14	105.57	111.07
1	A	938	ILE	N-CA-C	-5.00	106.49	113.00

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	2409	0	2413	88	0
2	A	26	0	35	1	0
3	A	17	0	0	1	0
4	A	4	0	6	0	0
5	A	108	0	0	5	0
All	All	2564	0	2454	88	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 18.

All (88) 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:772:PRO:O	1:A:852:LYS:HD3	1.72	0.90
1:A:697:GLU:HG2	1:A:864:ALA:HB2	1.59	0.84
1:A:829:GLU:HG3	1:A:893:HIS:CD2	2.21	0.76
1:A:768:SER:O	1:A:831:ARG:NH2	2.19	0.75
1:A:883:LEU:HG	1:A:887:LEU:HD12	1.72	0.71
1:A:825:MET:SD	1:A:853:ILE:HD13	2.32	0.70
1:A:863:GLY:H	1:A:866:ALA:HB3	1.59	0.68
1:A:938:ILE:HD12	1:A:981:ILE:HD11	1.74	0.68
1:A:931:GLU:OE1	5:A:1201:HOH:O	2.12	0.68
1:A:860:LYS:NZ	1:A:869:TYR:OH	2.27	0.67
1:A:924:SER:O	1:A:928:GLU:HG2	1.94	0.67
1:A:924:SER:O	1:A:928:GLU:CG	2.44	0.65
1:A:791:GLN:OE1	1:A:852:LYS:NZ	2.29	0.65
1:A:860:LYS:HG3	1:A:869:TYR:CZ	2.31	0.65
1:A:793:MET:CE	1:A:852:LYS:HG3	2.28	0.64
1:A:764:TYR:O	1:A:768:SER:OG	2.15	0.63
1:A:766:MET:HB3	1:A:777:LEU:HB2	1.80	0.63
1:A:1015:GLU:OE2	5:A:1202:HOH:O	2.15	0.63
1:A:985:GLU:H	1:A:985:GLU:CD	2.06	0.63
1:A:820:GLN:HG2	2:A:1101:PE5:H142	1.82	0.61
1:A:726:VAL:HG21	3:A:1102:N9L:C16	2.32	0.60
1:A:793:MET:HE1	1:A:852:LYS:HG3	1.83	0.60
1:A:829:GLU:HG3	1:A:893:HIS:CG	2.37	0.59
1:A:791:GLN:NE2	1:A:1012:ASP:OD2	2.36	0.59
1:A:697:GLU:CG	1:A:864:ALA:HB2	2.34	0.57
1:A:860:LYS:HG3	1:A:869:TYR:CE2	2.39	0.57
1:A:935:GLN:HA	1:A:944:TYR:CD1	2.39	0.57
1:A:733:PRO:HB2	1:A:736:GLU:HB2	1.86	0.56
1:A:855:ASP:OD2	5:A:1203:HOH:O	2.18	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:943:VAL:HG12	1:A:947:MET:HE2	1.90	0.53
1:A:842:ASN:O	1:A:854:THR:HG22	2.09	0.53
1:A:832:ARG:NH1	1:A:863:GLY:O	2.41	0.52
1:A:883:LEU:HG	1:A:887:LEU:CD1	2.39	0.52
1:A:836:ARG:NH1	1:A:858:ARG:O	2.43	0.52
1:A:808:ASN:HB3	1:A:989:LEU:HD21	1.92	0.52
1:A:872:GLU:HA	1:A:889:ARG:NH2	2.25	0.51
1:A:973[B]:ARG:NH1	5:A:1209:HOH:O	2.43	0.51
1:A:815:LEU:HG	1:A:979:LEU:HD12	1.93	0.51
1:A:932:ARG:HE	1:A:951:TRP:HB3	1.74	0.51
1:A:836:ARG:HH11	1:A:860:LYS:HG2	1.77	0.50
1:A:765:VAL:HG11	1:A:856:PHE:HB3	1.93	0.50
1:A:780:ILE:HG22	1:A:788:LEU:HD12	1.92	0.50
1:A:970:LYS:O	1:A:973[A]:ARG:HB2	2.12	0.49
1:A:925:SER:HA	1:A:928:GLU:HG3	1.95	0.49
1:A:943:VAL:HG12	1:A:947:MET:CE	2.43	0.49
1:A:960:LYS:O	1:A:963:GLU:HB2	2.12	0.49
1:A:767:ALA:HB2	1:A:777:LEU:HD23	1.95	0.49
1:A:908:MET:HG3	1:A:939:CYS:SG	2.54	0.48
1:A:714:LYS:HD3	1:A:727:TYR:CG	2.49	0.48
1:A:806:LYS:HB3	1:A:910:PHE:HB3	1.95	0.48
1:A:924:SER:O	1:A:928:GLU:HG3	2.12	0.48
1:A:939:CYS:SG	1:A:947:MET:HE1	2.53	0.48
1:A:772:PRO:O	1:A:852:LYS:CD	2.55	0.48
1:A:900:TYR:CE1	1:A:964:LEU:HD13	2.49	0.48
1:A:833:LEU:HD21	1:A:859:ALA:HB1	1.96	0.47
1:A:878:ILE:HG23	1:A:886:ILE:HD13	1.96	0.47
1:A:915:TYR:C	1:A:918:ILE:HD13	2.40	0.47
1:A:842:ASN:C	1:A:854:THR:HG22	2.40	0.47
1:A:791:GLN:HG3	1:A:1013:ALA:HB3	1.97	0.46
1:A:790:MET:HB2	1:A:790:MET:HE3	1.64	0.45
1:A:793:MET:HE2	1:A:852:LYS:HG3	1.99	0.45
1:A:898:TRP:C	1:A:898:TRP:CD1	2.95	0.45
1:A:909:THR:HG22	1:A:936:PRO:HB3	1.99	0.45
1:A:850:HIS:CD2	1:A:850:HIS:C	2.93	0.45
1:A:860:LYS:HD3	1:A:860:LYS:HA	1.79	0.45
1:A:868:GLU:CB	1:A:890:ILE:HG23	2.47	0.45
1:A:869:TYR:CE2	1:A:871:ALA:HB2	2.52	0.44
1:A:1018:ILE:HD13	1:A:1018:ILE:HA	1.86	0.43
1:A:751:THR:O	1:A:753:PRO:HD3	2.18	0.43
1:A:926:ILE:HG23	1:A:931:GLU:HB3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:976:GLN:HE21	1:A:976:GLN:HB3	1.60	0.43
1:A:818:CYS:SG	1:A:904:VAL:HG22	2.58	0.43
1:A:880:TRP:O	1:A:899:SER:OG	2.30	0.42
1:A:815:LEU:HG	1:A:979:LEU:CD1	2.50	0.42
1:A:868:GLU:HB3	1:A:892:THR:CG2	2.50	0.41
1:A:944:TYR:HA	1:A:947:MET:HE3	2.01	0.41
1:A:833:LEU:CD2	1:A:859:ALA:HB1	2.50	0.41
1:A:762:GLU:HG2	1:A:766:MET:HE2	2.02	0.41
1:A:860:LYS:HG3	1:A:869:TYR:CE1	2.55	0.41
1:A:834:VAL:HG12	1:A:836:ARG:CG	2.51	0.41
1:A:766:MET:O	1:A:769:VAL:HG22	2.21	0.41
1:A:898:TRP:CZ2	1:A:927:LEU:HD13	2.55	0.41
1:A:821:ILE:O	1:A:853:ILE:HD11	2.21	0.41
1:A:832:ARG:HB3	1:A:832:ARG:HH11	1.86	0.41
1:A:898:TRP:CE3	1:A:951:TRP:HA	2.56	0.41
1:A:902:VAL:O	1:A:906:GLU:HG3	2.21	0.40
1:A:776:ARG:HG3	5:A:1292:HOH:O	2.20	0.40
1:A:935:GLN:HA	1:A:936:PRO:HD3	1.99	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	301/331 (91%)	287 (95%)	13 (4%)	1 (0%)	36 55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	748	ARG

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	258/287 (90%)	221 (86%)	37 (14%)	3 6

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	715	ILE
1	A	716	LYS
1	A	725	THR
1	A	728	LYS
1	A	739	LYS
1	A	742	VAL
1	A	751	THR
1	A	768	SER
1	A	774	VAL
1	A	775	CYS
1	A	776	ARG
1	A	782	LEU
1	A	783	THR
1	A	784	SER
1	A	786	VAL
1	A	797	SER
1	A	808	ASN
1	A	833	LEU
1	A	868	GLU
1	A	878	ILE
1	A	886	ILE
1	A	888	HIS
1	A	889	ARG
1	A	921	SER
1	A	924	SER
1	A	925	SER
1	A	928	GLU
1	A	941	ILE
1	A	949	LYS
1	A	960	LYS

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Mol	Chain	Res	Type
1	A	969	SER
1	A	973[A]	ARG
1	A	973[B]	ARG
1	A	976	GLN
1	A	985	GLU
1	A	1017	LEU
1	A	1018	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	756	ASN
1	A	787	GLN
1	A	805	HIS
1	A	808	ASN
1	A	849	GLN

5.3.3 RNA [i](#)

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](#)

3 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PE5	A	1101	-	25,25,26	0.23	0	24,24,25	0.13	0
3	N9L	A	1102	1	18,18,18	1.80	2 (11%)	23,23,23	2.29	6 (26%)
4	DMS	A	1103	-	3,3,3	0.29	0	3,3,3	0.14	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
2	PE5	A	1101	-	-	12/23/23/24	-
3	N9L	A	1102	1	-	2/10/10/10	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1102	N9L	C01-C03	5.00	1.54	1.30
3	A	1102	N9L	C03-C04	4.78	1.55	1.48

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1102	N9L	C01-C03-C04	-7.34	114.15	122.22
3	A	1102	N9L	C06-N05-C10	3.81	122.65	117.21
3	A	1102	N9L	C09-C08-C12	-3.56	114.87	120.84
3	A	1102	N9L	C09-C10-N05	-3.27	118.63	122.92
3	A	1102	N9L	C07-C06-N05	-3.01	120.28	123.97
3	A	1102	N9L	C07-C08-C12	2.42	125.44	121.25

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1102	N9L	C01-C03-C04-O02
2	A	1101	PE5	O2-C3-C4-O3
2	A	1101	PE5	O4-C7-C8-O5
2	A	1101	PE5	O1-C1-C2-O2
2	A	1101	PE5	C48-C50-O1-C1
2	A	1101	PE5	C9-C10-O6-C11
2	A	1101	PE5	C14-C13-O7-C12

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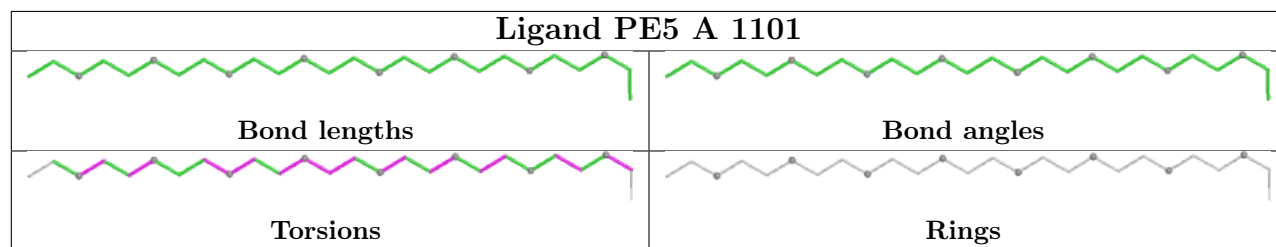
Mol	Chain	Res	Type	Atoms
2	A	1101	PE5	C10-C9-O5-C8
2	A	1101	PE5	C13-C14-O8-C15
2	A	1101	PE5	C7-C8-O5-C9
3	A	1102	N9L	C01-C03-C04-N11
2	A	1101	PE5	C6-C5-O3-C4
2	A	1101	PE5	C5-C6-O4-C7
2	A	1101	PE5	C12-C11-O6-C10

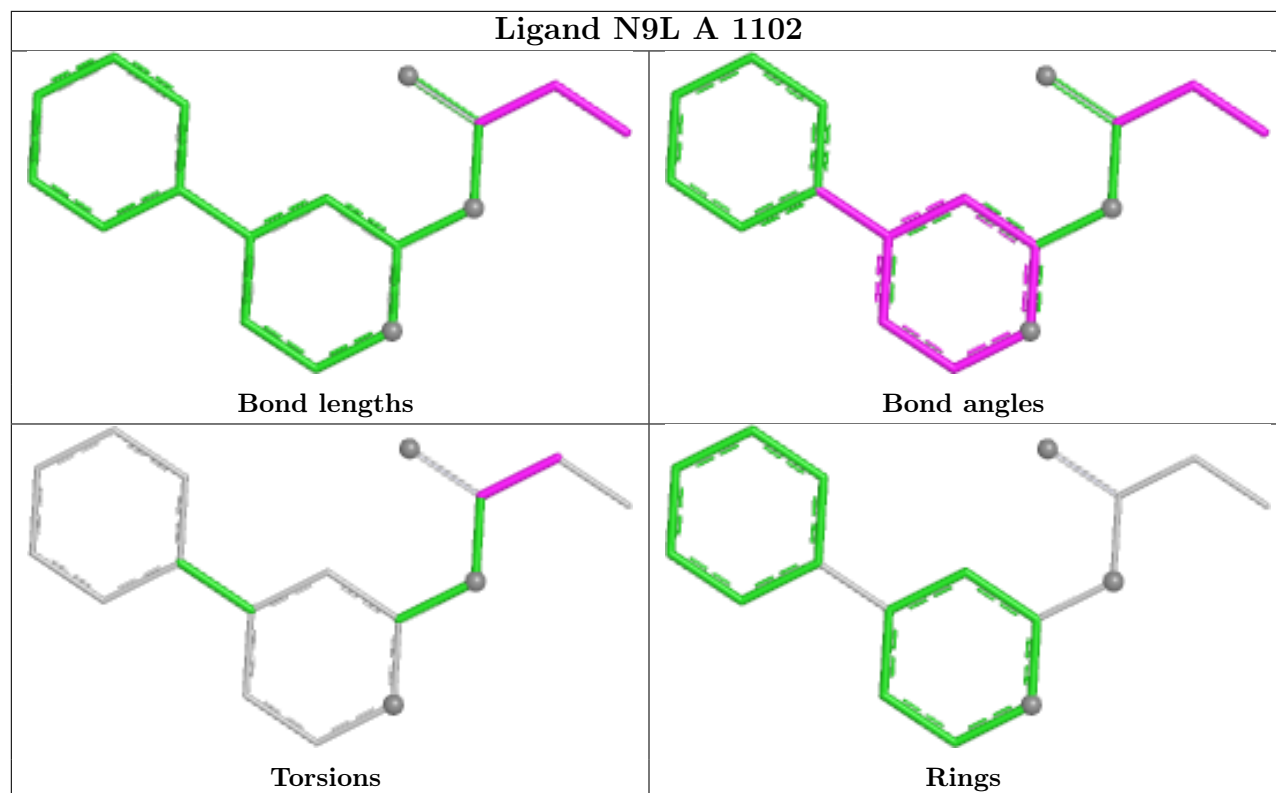
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1101	PE5	1	0
3	A	1102	N9L	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	306/331 (92%)	1.02	44 (14%) 6 5	16, 36, 103, 156	1 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	867	ALA	7.0
1	A	861	LEU	7.0
1	A	864	ALA	6.5
1	A	869	TYR	6.2
1	A	862	LEU	6.1
1	A	751	THR	5.8
1	A	866	ALA	5.8
1	A	750	ALA	5.4
1	A	990	PRO	5.2
1	A	865	ALA	5.1
1	A	870	HIS	5.1
1	A	696	GLY	5.0
1	A	749	GLU	4.7
1	A	1005	GLU	4.7
1	A	989	LEU	4.4
1	A	871	ALA	4.1
1	A	863	GLY	4.0
1	A	868	GLU	3.8
1	A	776	ARG	3.8
1	A	874	GLY	3.7
1	A	753	PRO	3.6
1	A	875	LYS	3.6
1	A	748	ARG	3.6
1	A	747	LEU	3.5
1	A	1007	MET	3.4
1	A	873	GLY	3.1
1	A	1018	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	987	MET	2.8
1	A	986	ARG	2.8
1	A	758	GLU	2.8
1	A	988	HIS	2.8
1	A	985	GLU	2.7
1	A	755	ALA	2.7
1	A	752	SER	2.6
1	A	786	VAL	2.5
1	A	754	LYS	2.5
1	A	918	ILE	2.4
1	A	1006	ASP	2.4
1	A	1019	PRO	2.3
1	A	718	LEU	2.2
1	A	887	LEU	2.1
1	A	725	THR	2.1
1	A	697	GLU	2.1
1	A	727	TYR	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.

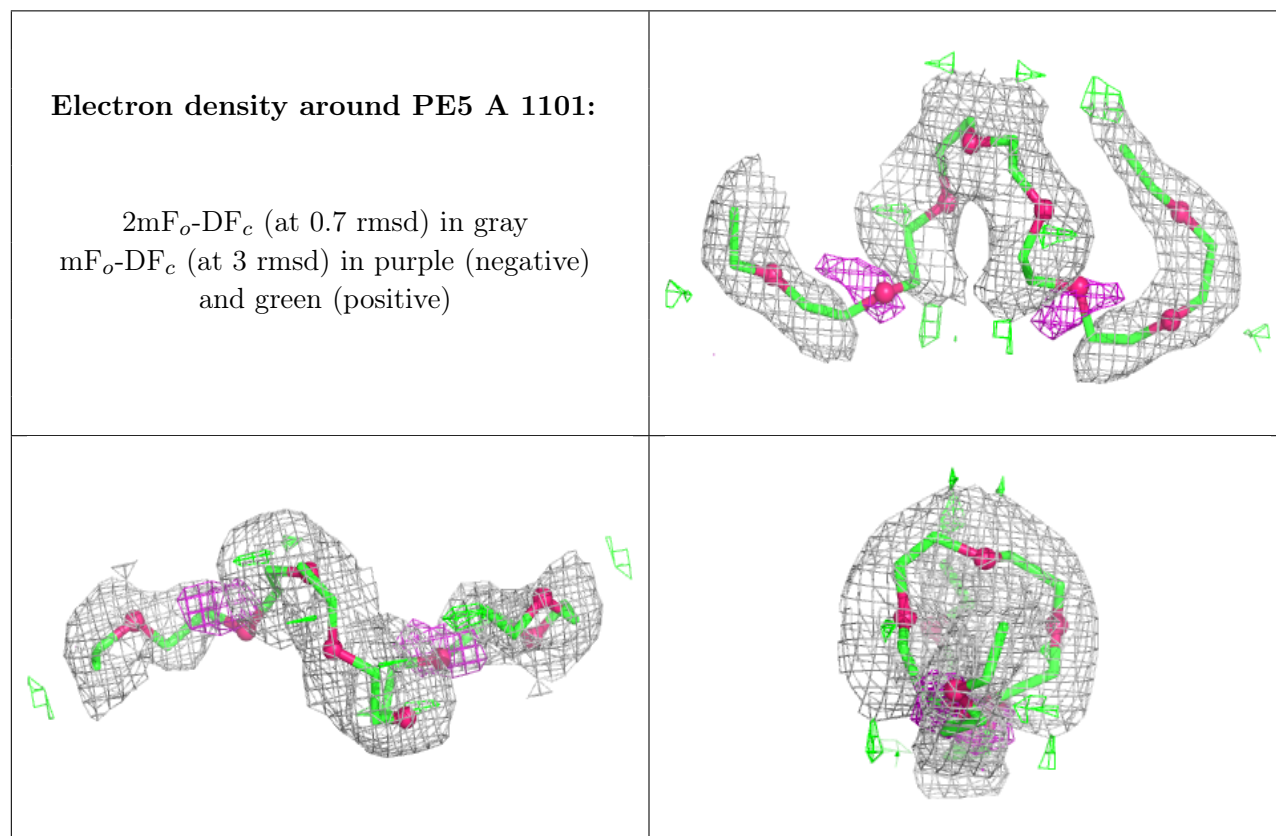
6.4 Ligands [i](#)

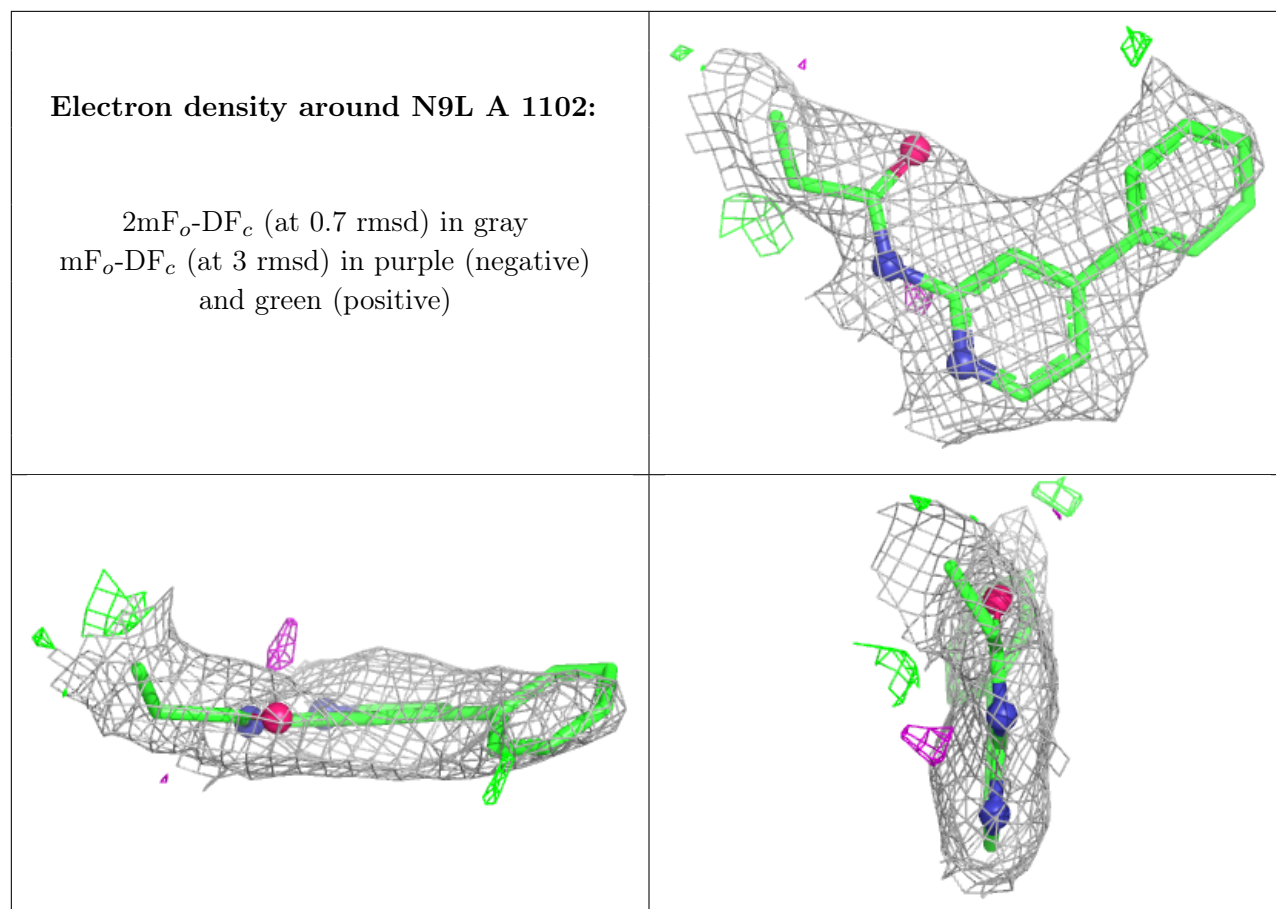
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	A	1103	4/4	0.77	0.21	70,72,75,75	0
2	PE5	A	1101	26/27	0.78	0.26	24,45,112,115	26
3	N9L	A	1102	17/17	0.81	0.18	44,57,74,76	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.





6.5 Other polymers ⓘ

There are no such residues in this entry.