



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:03 PM UTC

PDB ID : 5A86 / pdb_00005a86
Title : Structure of pregnane X receptor in complex with a Sphingosine 1- Phosphate Receptor 1 Antagonist
Authors : Xue, Y.; Oster, L.
Deposited on : 2015-07-13
Resolution : 2.25 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

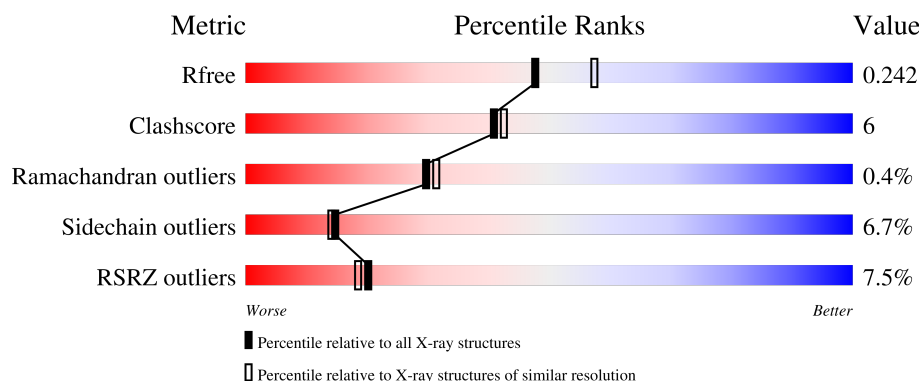
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	314	7% 70% 14% • 14%
1	B	314	6% 71% 14% • 13%
2	C	17	6% 71% 18% 12%
2	D	17	6% 100%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5081 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 NUCLEAR RECEPTOR SUBFAMILY 1 GROUP I MEMBER 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2213	1423	375	397	18			
1	B	274	Total	C	N	O	S	0	0	0
			2239	1439	382	400	18			

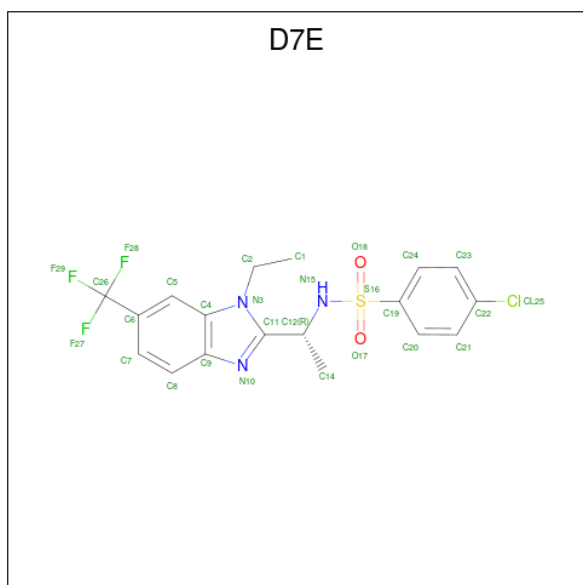
There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	119	MET	-	expression tag	UNP O75469
A	120	LYS	-	expression tag	UNP O75469
A	121	LYS	-	expression tag	UNP O75469
A	122	GLY	-	expression tag	UNP O75469
A	123	HIS	-	expression tag	UNP O75469
A	124	HIS	-	expression tag	UNP O75469
A	125	HIS	-	expression tag	UNP O75469
A	126	HIS	-	expression tag	UNP O75469
A	127	HIS	-	expression tag	UNP O75469
A	128	HIS	-	expression tag	UNP O75469
A	129	GLY	-	expression tag	UNP O75469
B	119	MET	-	expression tag	UNP O75469
B	120	LYS	-	expression tag	UNP O75469
B	121	LYS	-	expression tag	UNP O75469
B	122	GLY	-	expression tag	UNP O75469
B	123	HIS	-	expression tag	UNP O75469
B	124	HIS	-	expression tag	UNP O75469
B	125	HIS	-	expression tag	UNP O75469
B	126	HIS	-	expression tag	UNP O75469
B	127	HIS	-	expression tag	UNP O75469
B	128	HIS	-	expression tag	UNP O75469
B	129	GLY	-	expression tag	UNP O75469

- Molecule 2 is a protein called NUCLEAR RECEPTOR COACTIVATOR 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	0	0	0
			131	82	27	22			
2	D	17	Total	C	N	O	0	0	0
			141	87	29	25			

- Molecule 3 is 4-chloro-N-[(1R)-1-[1-ethyl-6-(trifluoromethyl)benzimidazol-2-yl]ethyl]benzene sulfonamide (CCD ID: D7E) (formula: $C_{18}H_{17}ClF_3N_3O_2S$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	S	0	0
			28	18	1	3	3	2	1		
3	B	1	Total	C	Cl	F	N	O	S	0	0
			28	18	1	3	3	2	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	125	Total	O	0	0
			125	125		
4	A	3	Total	O	0	0
			3	3		
4	A	1	Total	O	0	0
			1	1		
4	A	1	Total	O	0	0
			1	1		
4	A	1	Total	O	0	0
			1	1		

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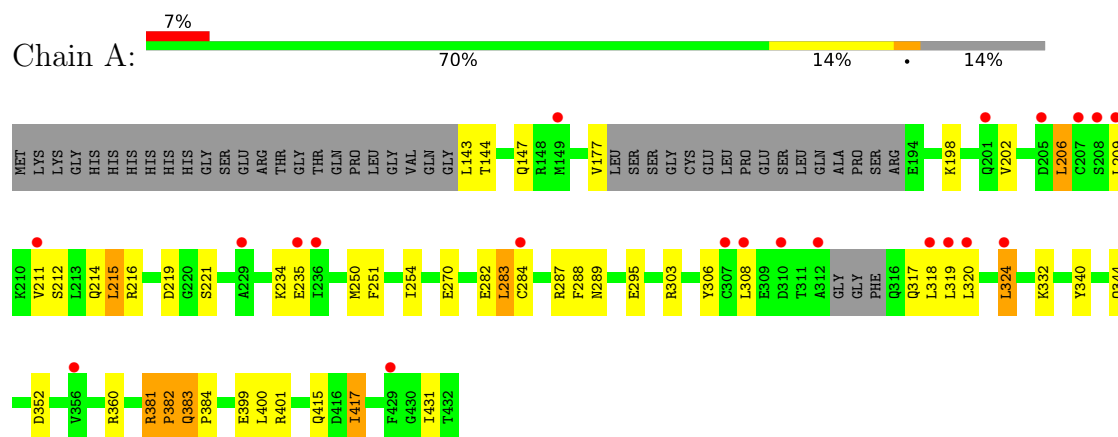
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	O 1	0	0
4	A	6	Total 6	O 6	0	0
4	B	133	Total 133	O 133	0	0
4	B	1	Total 1	O 1	0	0
4	B	1	Total 1	O 1	0	0
4	B	1	Total 1	O 1	0	0
4	B	1	Total 1	O 1	0	0
4	B	3	Total 3	O 3	0	0
4	B	5	Total 5	O 5	0	0
4	B	1	Total 1	O 1	0	0
4	B	3	Total 3	O 3	0	0
4	C	5	Total 5	O 5	0	0
4	D	2	Total 2	O 2	0	0
4	D	6	Total 6	O 6	0	0
4	D	1	Total 1	O 1	0	0

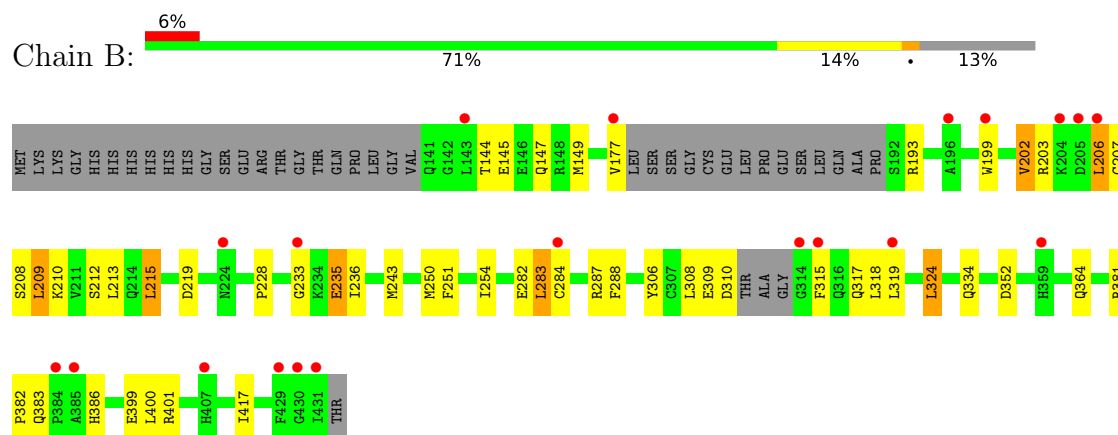
3 Residue-property plots [i](#)

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.

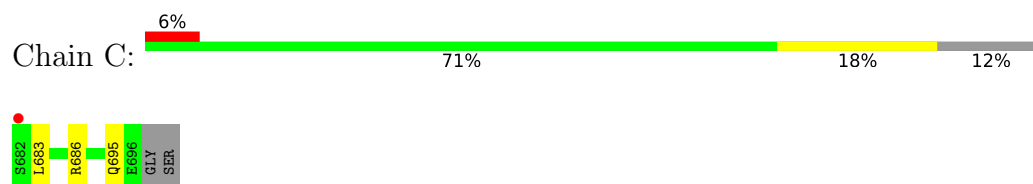
• Molecule 1: NUCLEAR RECEPTOR SUBFAMILY 1 GROUP I MEMBER 2



• Molecule 1: NUCLEAR RECEPTOR SUBFAMILY 1 GROUP I MEMBER 2



• Molecule 2: NUCLEAR RECEPTOR COACTIVATOR 1



• Molecule 2: NUCLEAR RECEPTOR COACTIVATOR 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.67Å 89.20Å 106.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.88 – 2.25 14.88 – 2.25	Depositor EDS
% Data completeness (in resolution range)	97.6 (14.88-2.25) 97.2 (14.88-2.25)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.24Å)	Xtriage
Refinement program	BUSTER 2.11.6	Depositor
R, R_{free}	0.231 , 0.256 0.222 , 0.242	Depositor DCC
R_{free} test set	1877 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 65.7	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	0.94	EDS
Total number of atoms	5081	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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: D7E

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	0.64	0/2261	1.04	1/3045 (0.0%)
1	B	0.61	0/2288	1.02	2/3078 (0.1%)
2	C	0.50	0/132	1.01	0/175
2	D	0.59	0/142	1.25	0/188
All	All	0.62	0/4823	1.03	3/6486 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	399	GLU	CB-CG-CD	6.15	123.05	112.60
1	A	212	SER	N-CA-C	-5.42	102.50	110.52
1	B	212	SER	N-CA-C	-5.08	102.99	110.52

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	2213	0	2222	39	0
1	B	2239	0	2244	29	0
2	C	131	0	139	1	0
2	D	141	0	147	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	28	0	17	2	0
3	B	28	0	17	3	0
4	A	138	0	0	0	0
4	B	149	0	0	1	0
4	C	5	0	0	0	0
4	D	9	0	0	0	0
All	All	5081	0	4786	62	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 6.

All (62) 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:383:GLN:HB3	1:A:384:PRO:CD	1.91	1.00
1:A:383:GLN:HB3	1:A:384:PRO:HD2	1.59	0.84
1:A:198:LYS:HB3	1:A:417:ILE:HD11	1.60	0.83
1:A:381:ARG:HH11	1:A:381:ARG:HG2	1.46	0.81
1:A:308:LEU:HD13	1:A:324:LEU:HD21	1.64	0.80
1:A:214:GLN:HE21	1:A:216:ARG:HH11	1.33	0.75
1:A:352:ASP:CG	1:A:401:ARG:HH12	1.98	0.71
1:A:177:VAL:HG21	1:B:215:LEU:HD13	1.74	0.69
2:C:683:LEU:HD12	2:C:686:ARG:HH21	1.60	0.67
1:A:383:GLN:HB3	1:A:384:PRO:HD3	1.77	0.66
1:B:199:TRP:HZ3	1:B:233:GLY:O	1.82	0.63
1:A:177:VAL:CG2	1:B:215:LEU:HD13	2.29	0.62
1:B:251:PHE:CD1	1:B:284:CYS:SG	2.94	0.60
1:A:198:LYS:CB	1:A:417:ILE:HD11	2.29	0.60
1:A:235:GLU:HB3	1:B:219:ASP:HB2	1.84	0.59
1:A:211:VAL:HG21	1:A:306:TYR:HB3	1.83	0.59
1:B:254:ILE:HD11	1:B:287:ARG:HD2	1.85	0.59
1:A:254:ILE:HD11	1:A:287:ARG:HD2	1.84	0.58
1:A:235:GLU:CB	1:B:219:ASP:HB2	2.35	0.57
1:B:144:THR:OG1	1:B:147:GLN:HG3	2.07	0.54
1:B:199:TRP:O	1:B:203:ARG:HG3	2.09	0.53
1:A:282:GLU:HG2	1:A:400:LEU:HG	1.90	0.53
1:A:144:THR:OG1	1:A:147:GLN:HG3	2.09	0.53
1:B:383:GLN:H	1:B:383:GLN:CD	2.16	0.53
1:A:214:GLN:HE21	1:A:216:ARG:NH1	2.02	0.52
1:B:282:GLU:HG2	1:B:400:LEU:HG	1.92	0.52
1:A:283:LEU:HD13	1:A:344:GLN:HB3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:LYS:HB3	1:A:417:ILE:CD1	2.37	0.51
1:A:381:ARG:HG2	1:A:381:ARG:NH1	2.17	0.51
1:A:215:LEU:HD13	1:B:177:VAL:CG2	2.42	0.50
1:B:383:GLN:HB2	1:B:386:HIS:CD2	2.46	0.49
1:B:318:LEU:HB3	1:B:324:LEU:HD12	1.95	0.49
1:A:289:ASN:ND2	1:A:340:TYR:OH	2.46	0.49
1:B:382:PRO:HG2	1:B:383:GLN:OE1	2.12	0.49
1:A:221:SER:HB3	1:B:228:PRO:HD3	1.94	0.48
1:A:318:LEU:HB3	1:A:324:LEU:HD22	1.96	0.47
1:B:207:CYS:HA	1:B:208:SER:HA	1.66	0.47
1:A:209:LEU:HA	3:A:1433:D7E:H11C	1.97	0.46
1:A:308:LEU:CD1	1:A:324:LEU:HD21	2.38	0.46
1:B:202:VAL:HG11	1:B:417:ILE:HD11	1.98	0.46
1:A:251:PHE:CD1	1:A:284:CYS:SG	3.10	0.45
1:B:243:MET:HG3	3:B:1432:D7E:C6	2.46	0.45
1:B:352:ASP:CG	1:B:401:ARG:HH12	2.24	0.45
1:A:317:GLN:O	1:A:320:LEU:HB2	2.17	0.45
1:A:206:LEU:HD23	1:A:209:LEU:HD12	1.99	0.44
1:B:306:TYR:CD2	3:B:1432:D7E:CL25	3.07	0.44
1:A:219:ASP:HB2	1:B:235:GLU:HB3	1.99	0.44
1:A:306:TYR:CD2	3:A:1433:D7E:CL25	3.08	0.43
1:B:251:PHE:HD1	1:B:284:CYS:SG	2.40	0.43
1:A:381:ARG:O	1:A:382:PRO:C	2.61	0.43
1:A:308:LEU:HD22	1:A:324:LEU:HD11	2.01	0.42
1:A:202:VAL:O	1:A:206:LEU:HD12	2.19	0.42
1:A:211:VAL:CG2	1:A:306:TYR:HB3	2.49	0.42
1:A:415:GLN:HG2	1:A:431:ILE:HD13	2.01	0.42
1:A:250:MET:HG3	1:A:288:PHE:CZ	2.55	0.42
1:A:381:ARG:HA	1:A:382:PRO:HD2	1.68	0.41
1:B:209:LEU:HB2	3:B:1432:D7E:H11C	2.03	0.41
1:B:254:ILE:HD12	1:B:283:LEU:HB3	2.02	0.41
1:B:250:MET:HG3	1:B:288:PHE:CZ	2.55	0.41
1:B:334:GLN:HA	4:B:2089:HOH:O	2.21	0.41
1:B:206:LEU:HD23	1:B:236:ILE:HB	2.03	0.40
1:B:213:LEU:HD23	1:B:213:LEU:C	2.46	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	265/314 (84%)	256 (97%)	7 (3%)	2 (1%)	16	14
1	B	268/314 (85%)	257 (96%)	11 (4%)	0	100	100
2	C	13/17 (76%)	12 (92%)	1 (8%)	0	100	100
2	D	15/17 (88%)	14 (93%)	1 (7%)	0	100	100
All	All	561/662 (85%)	539 (96%)	20 (4%)	2 (0%)	30	31

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	PRO
1	A	383	GLN

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	244/278 (88%)	229 (94%)	15 (6%)	17	17
1	B	246/278 (88%)	227 (92%)	19 (8%)	12	10
2	C	15/16 (94%)	14 (93%)	1 (7%)	15	14
2	D	16/16 (100%)	16 (100%)	0	100	100
All	All	521/588 (89%)	486 (93%)	35 (7%)	15	14

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	143	LEU
1	A	206	LEU
1	A	215	LEU
1	A	234	LYS
1	A	270	GLU
1	A	283	LEU
1	A	295	GLU
1	A	303	ARG
1	A	319	LEU
1	A	324	LEU
1	A	332	LYS
1	A	360	ARG
1	A	381	ARG
1	A	399	GLU
1	A	417	ILE
1	B	145	GLU
1	B	149	MET
1	B	193	ARG
1	B	202	VAL
1	B	206	LEU
1	B	209	LEU
1	B	210	LYS
1	B	215	LEU
1	B	235	GLU
1	B	283	LEU
1	B	308	LEU
1	B	309	GLU
1	B	310	ASP
1	B	315	PHE
1	B	317	GLN
1	B	319	LEU
1	B	324	LEU
1	B	364	GLN
1	B	381	ARG
2	C	695	GLN

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

Mol	Chain	Res	Type
1	A	168	HIS
1	A	214	GLN
1	A	224	ASN
1	A	272	GLN

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Mol	Chain	Res	Type
1	A	289	ASN
1	B	168	HIS
1	B	201	GLN
1	B	285	GLN
1	B	289	ASN
1	B	383	GLN
1	B	386	HIS

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 ⓘ

2 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	D7E	B	1432	-	30,30,30	1.16	2 (6%)	46,46,46	1.68	9 (19%)
3	D7E	A	1433	-	30,30,30	1.11	3 (10%)	46,46,46	1.63	7 (15%)

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	D7E	B	1432	-	-	4/23/23/23	0/3/3/3
3	D7E	A	1433	-	-	3/23/23/23	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1432	D7E	S16-N15	3.69	1.67	1.61
3	A	1433	D7E	S16-N15	3.11	1.66	1.61
3	B	1432	D7E	F29-C26	2.27	1.41	1.33
3	A	1433	D7E	F29-C26	2.10	1.40	1.33
3	A	1433	D7E	F27-C26	2.03	1.40	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1433	D7E	N3-C11-N10	-5.55	108.94	113.26
3	A	1433	D7E	C9-N10-C11	5.35	110.47	104.75
3	B	1432	D7E	C9-N10-C11	5.01	110.10	104.75
3	B	1432	D7E	N3-C11-N10	-4.12	110.06	113.26
3	B	1432	D7E	C4-C9-N10	-3.40	105.88	110.19
3	B	1432	D7E	C24-C19-S16	-2.70	116.78	119.76
3	A	1433	D7E	C4-C9-N10	-2.67	106.80	110.19
3	A	1433	D7E	C1-C2-N3	-2.66	102.23	111.57
3	B	1432	D7E	O18-S16-C19	-2.43	104.91	107.98
3	B	1432	D7E	C5-C4-C9	-2.28	119.74	122.47
3	B	1432	D7E	C8-C9-N10	2.24	134.56	128.21
3	A	1433	D7E	F29-C26-C6	-2.23	108.12	112.90
3	B	1432	D7E	C20-C19-S16	2.14	122.11	119.76
3	A	1433	D7E	C8-C9-N10	2.10	134.17	128.21
3	B	1432	D7E	O18-S16-O17	2.07	122.04	119.52
3	A	1433	D7E	C4-N3-C11	2.05	108.12	106.61

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1432	D7E	C11-C12-N15-S16
3	A	1433	D7E	N10-C11-C12-N15
3	B	1432	D7E	N3-C11-C12-C14
3	B	1432	D7E	C14-C12-N15-S16

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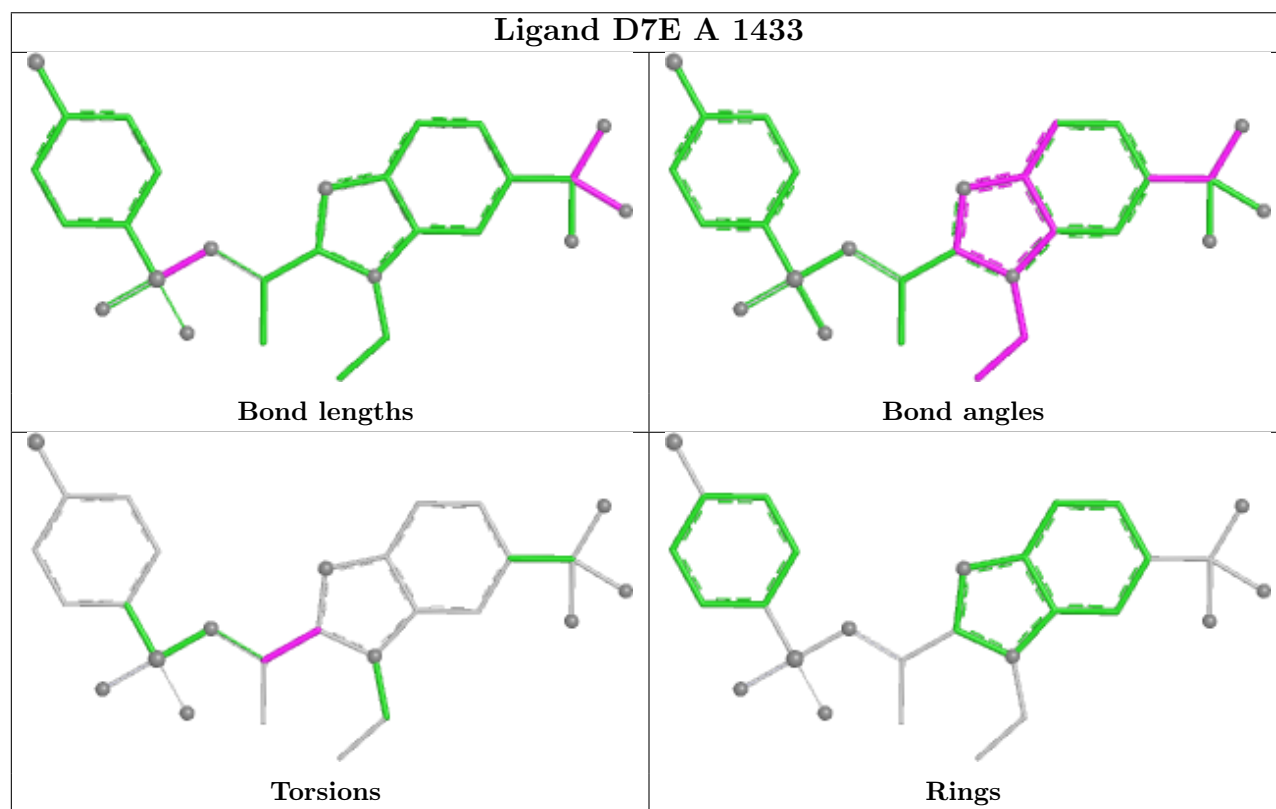
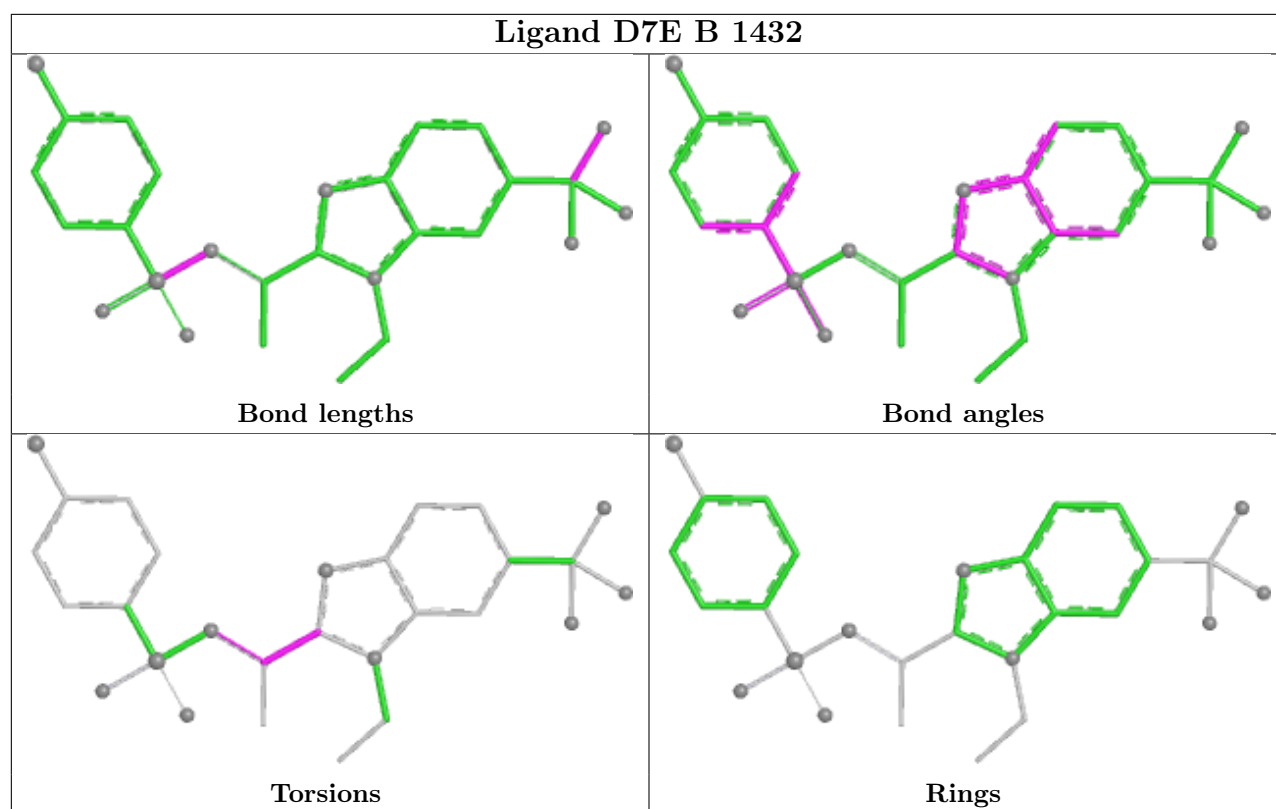
Mol	Chain	Res	Type	Atoms
3	A	1433	D7E	N3-C11-C12-C14
3	A	1433	D7E	N10-C11-C12-C14
3	B	1432	D7E	N10-C11-C12-C14

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1432	D7E	3	0
3	A	1433	D7E	2	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	271/314 (86%)	0.51	21 (7%) 19 17	26, 52, 106, 137	0
1	B	274/314 (87%)	0.31	20 (7%) 21 19	26, 49, 114, 145	0
2	C	15/17 (88%)	0.54	1 (6%) 24 22	47, 60, 92, 95	0
2	D	17/17 (100%)	0.57	1 (5%) 28 26	39, 55, 95, 105	0
All	All	577/662 (87%)	0.42	43 (7%) 20 18	26, 52, 110, 145	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	308	LEU	6.8
1	B	233	GLY	4.8
1	B	385	ALA	4.5
1	A	229	ALA	4.5
1	B	196	ALA	4.0
1	A	307	CYS	3.8
1	A	235	GLU	3.4
1	B	204	LYS	3.4
1	A	211	VAL	3.4
2	C	682	SER	3.2
1	A	207	CYS	3.1
1	A	319	LEU	3.1
1	A	324	LEU	3.1
1	A	429	PHE	3.0
1	A	209	LEU	2.9
1	A	320	LEU	2.8
1	A	236	ILE	2.7
1	B	314	GLY	2.7
1	A	310	ASP	2.7
1	B	430	GLY	2.7
1	B	429	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	356	VAL	2.5
1	A	312	ALA	2.5
1	B	284	CYS	2.4
1	A	318	LEU	2.4
1	A	149	MET	2.4
1	B	205	ASP	2.4
1	A	284	CYS	2.4
1	B	315	PHE	2.4
1	B	431	ILE	2.3
1	B	319	LEU	2.3
1	A	208	SER	2.3
2	D	683	LEU	2.3
1	B	359	HIS	2.3
1	B	384	PRO	2.2
1	A	201	GLN	2.2
1	B	199	TRP	2.2
1	B	206	LEU	2.2
1	B	177	VAL	2.2
1	B	143	LEU	2.1
1	B	224	ASN	2.1
1	B	407	HIS	2.0
1	A	205	ASP	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 [i](#)

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

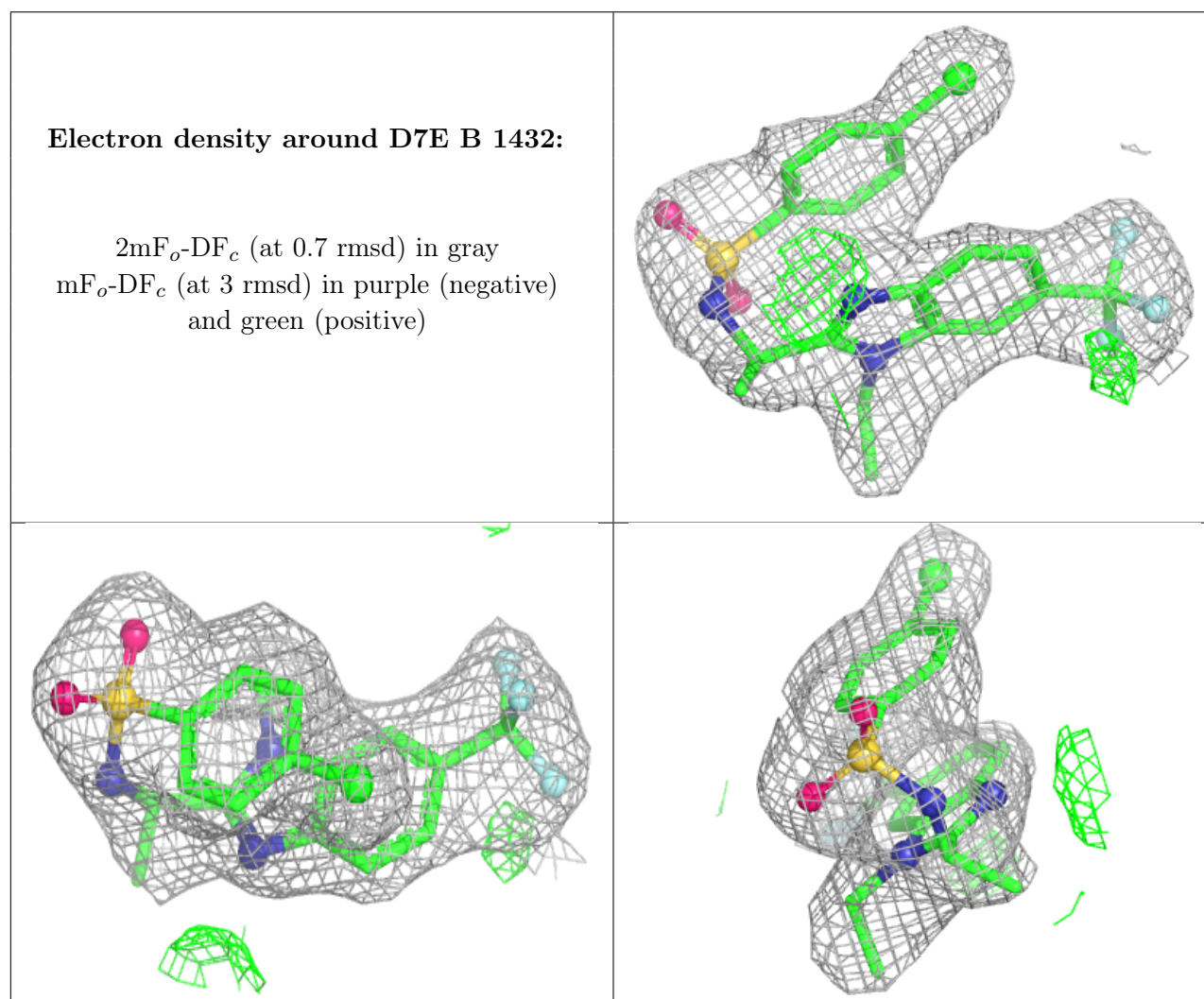
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	D7E	B	1432	28/28	0.91	0.09	59,71,73,74	0

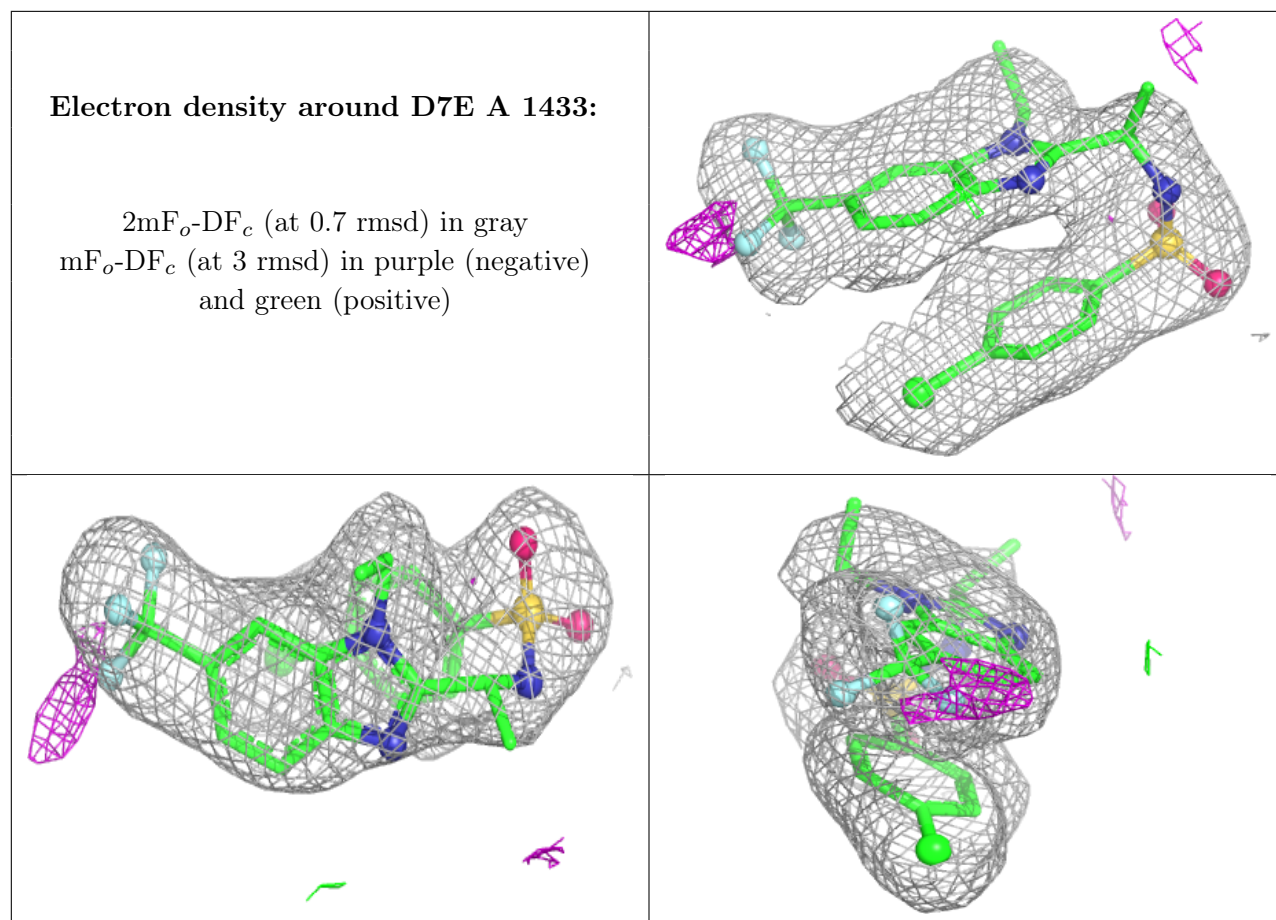
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	D7E	A	1433	28/28	0.93	0.09	49,61,73,75	0

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.