



Full wwPDB X-ray Structure Validation Report

May 14, 2020 – 09:11 pm BST

PDB ID : 6EMX
Title : X-ray structure of the N15'C mutant of GLIC in complex with bromoform
Authors : Fourati, Z.; Delarue, M.
Deposited on : 2017-10-03
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the  symbol.

The following versions of software and data (see [references](#) ) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

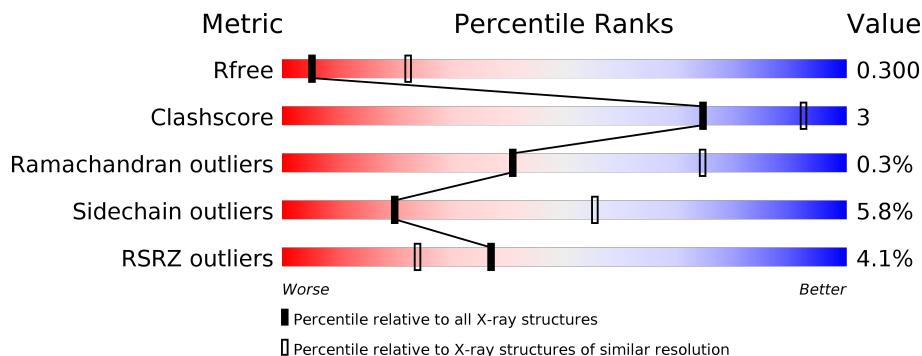
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	327	 2% 81% 14% 5%
1	B	327	 3% 82% 13% 5%
1	C	327	 5% 82% 13% 5%
1	D	327	 4% 81% 14% 5%
1	E	327	 5% 83% 12% 5%

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	MBR	E	404	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12685 atoms, of which 6 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 Proton-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	311	2523	1663	403	452	5	0	0	0
1	B	311	2523	1663	403	452	5	0	0	0
1	C	311	2523	1663	403	452	5	0	0	0
1	D	311	2523	1663	403	452	5	0	0	0
1	E	311	2523	1663	403	452	5	0	0	0

There are 55 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	239	CYS	ASN	engineered mutation	UNP Q7NDN8
A	318	GLY	-	expression tag	UNP Q7NDN8
A	319	TYR	-	expression tag	UNP Q7NDN8
A	320	PRO	-	expression tag	UNP Q7NDN8
A	321	TYR	-	expression tag	UNP Q7NDN8
A	322	ASP	-	expression tag	UNP Q7NDN8
A	323	VAL	-	expression tag	UNP Q7NDN8
A	324	PRO	-	expression tag	UNP Q7NDN8
A	325	ASP	-	expression tag	UNP Q7NDN8
A	326	TYR	-	expression tag	UNP Q7NDN8
A	327	ALA	-	expression tag	UNP Q7NDN8
B	239	CYS	ASN	engineered mutation	UNP Q7NDN8
B	318	GLY	-	expression tag	UNP Q7NDN8
B	319	TYR	-	expression tag	UNP Q7NDN8
B	320	PRO	-	expression tag	UNP Q7NDN8
B	321	TYR	-	expression tag	UNP Q7NDN8
B	322	ASP	-	expression tag	UNP Q7NDN8
B	323	VAL	-	expression tag	UNP Q7NDN8
B	324	PRO	-	expression tag	UNP Q7NDN8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	325	ASP	-	expression tag	UNP Q7NDN8
B	326	TYR	-	expression tag	UNP Q7NDN8
B	327	ALA	-	expression tag	UNP Q7NDN8
C	239	CYS	ASN	engineered mutation	UNP Q7NDN8
C	318	GLY	-	expression tag	UNP Q7NDN8
C	319	TYR	-	expression tag	UNP Q7NDN8
C	320	PRO	-	expression tag	UNP Q7NDN8
C	321	TYR	-	expression tag	UNP Q7NDN8
C	322	ASP	-	expression tag	UNP Q7NDN8
C	323	VAL	-	expression tag	UNP Q7NDN8
C	324	PRO	-	expression tag	UNP Q7NDN8
C	325	ASP	-	expression tag	UNP Q7NDN8
C	326	TYR	-	expression tag	UNP Q7NDN8
C	327	ALA	-	expression tag	UNP Q7NDN8
D	239	CYS	ASN	engineered mutation	UNP Q7NDN8
D	318	GLY	-	expression tag	UNP Q7NDN8
D	319	TYR	-	expression tag	UNP Q7NDN8
D	320	PRO	-	expression tag	UNP Q7NDN8
D	321	TYR	-	expression tag	UNP Q7NDN8
D	322	ASP	-	expression tag	UNP Q7NDN8
D	323	VAL	-	expression tag	UNP Q7NDN8
D	324	PRO	-	expression tag	UNP Q7NDN8
D	325	ASP	-	expression tag	UNP Q7NDN8
D	326	TYR	-	expression tag	UNP Q7NDN8
D	327	ALA	-	expression tag	UNP Q7NDN8
E	239	CYS	ASN	engineered mutation	UNP Q7NDN8
E	318	GLY	-	expression tag	UNP Q7NDN8
E	319	TYR	-	expression tag	UNP Q7NDN8
E	320	PRO	-	expression tag	UNP Q7NDN8
E	321	TYR	-	expression tag	UNP Q7NDN8
E	322	ASP	-	expression tag	UNP Q7NDN8
E	323	VAL	-	expression tag	UNP Q7NDN8
E	324	PRO	-	expression tag	UNP Q7NDN8
E	325	ASP	-	expression tag	UNP Q7NDN8
E	326	TYR	-	expression tag	UNP Q7NDN8
E	327	ALA	-	expression tag	UNP Q7NDN8

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Cl 1 1	0	0

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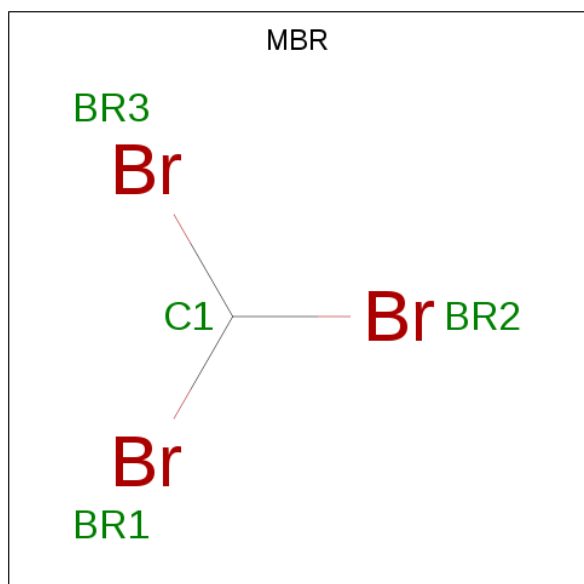
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	D	1	Total Cl 1 1	0	0
2	E	1	Total Cl 1 1	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Na 1 1	0	0
3	A	2	Total Na 2 2	0	0
3	D	1	Total Na 1 1	0	0
3	C	1	Total Na 1 1	0	0
3	E	1	Total Na 1 1	0	0

- Molecule 4 is TRIBROMOMETHANE (three-letter code: MBR) (formula: CHBr_3).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Br C H 5 3 1 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	Br	C	H		
4	B	1	5	3	1	1	0	0
4	C	1	5	3	1	1	0	0
4	D	1	5	3	1	1	0	0
4	E	1	5	3	1	1	0	0
4	E	1	5	3	1	1	0	0

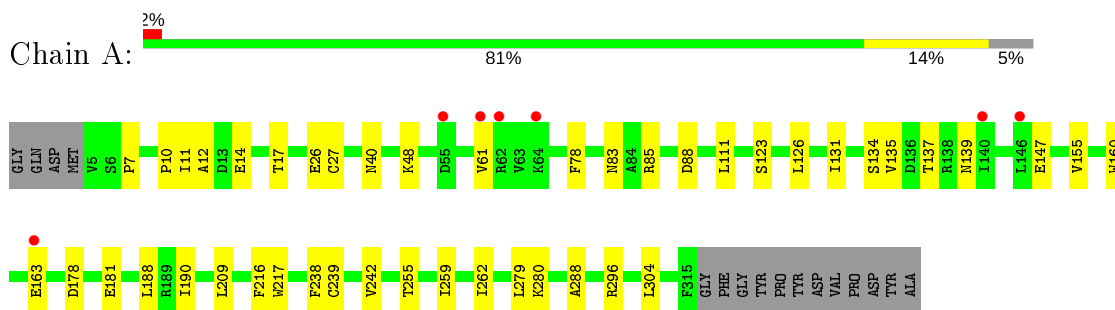
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	5	Total	O	0	0
			5	5		
5	B	5	Total	O	0	0
			5	5		
5	C	4	Total	O	0	0
			4	4		
5	D	5	Total	O	0	0
			5	5		
5	E	11	Total	O	0	0
			11	11		

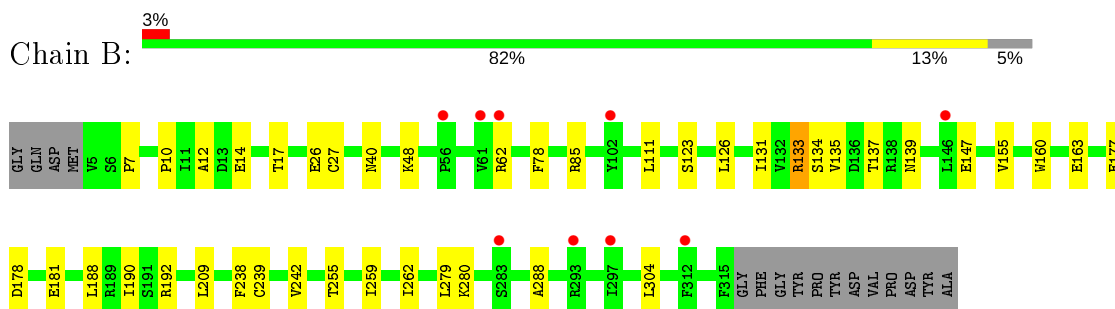
3 Residue-property plots [i](#)

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

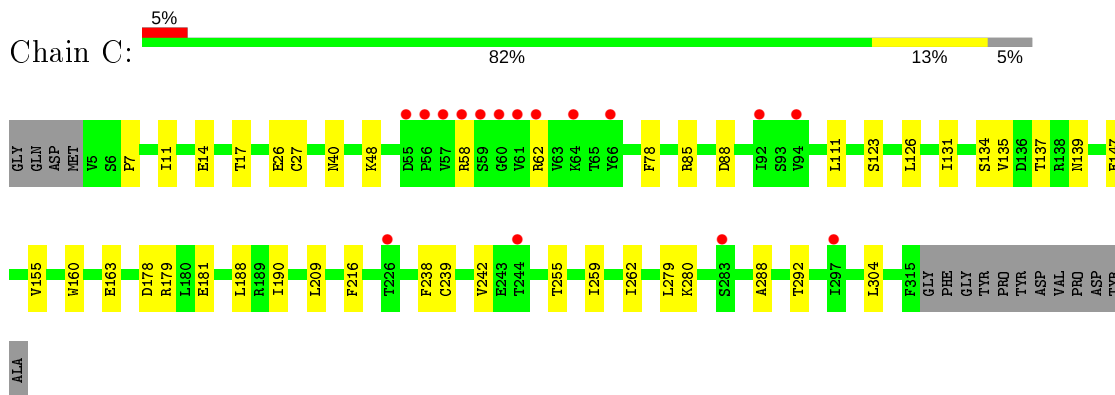
- Molecule 1: Proton-gated ion channel



- Molecule 1: Proton-gated ion channel

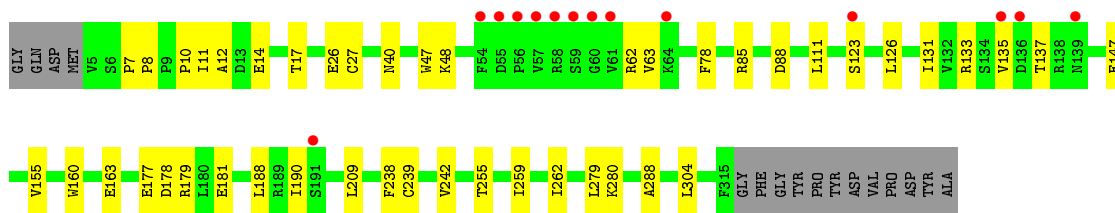


- Molecule 1: Proton-gated ion channel

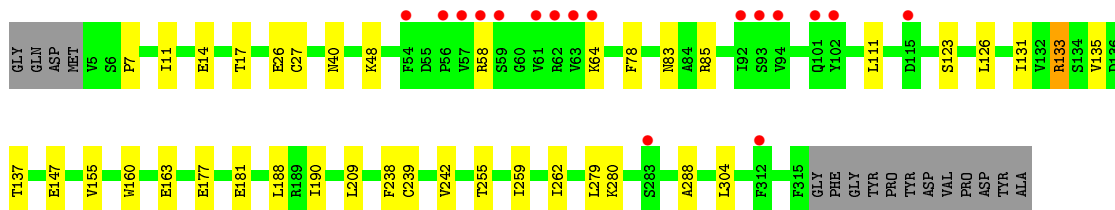
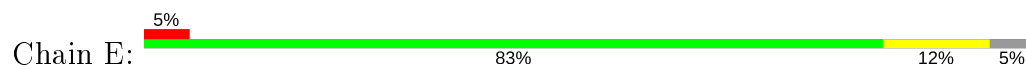


- Molecule 1: Proton-gated ion channel





- Molecule 1: Proton-gated ion channel



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	179.79Å 133.32Å 158.77Å 90.00° 101.68° 90.00°	Depositor
Resolution (Å)	28.62 – 3.20 28.53 – 3.20	Depositor EDS
% Data completeness (in resolution range)	98.8 (28.62-3.20) 98.8 (28.53-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 3.17Å)	Xtrriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.246 , 0.268 0.275 , 0.300	Depositor DCC
R_{free} test set	2914 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	103.6	Xtrriage
Anisotropy	0.332	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 66.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	12685	wwPDB-VP
Average B, all atoms (Å ²)	147.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MBR, CL

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	0.43	0/2591	0.61	0/3542
1	B	0.43	0/2591	0.63	0/3542
1	C	0.43	0/2591	0.63	0/3542
1	D	0.43	0/2591	0.61	0/3542
1	E	0.43	0/2591	0.62	0/3542
All	All	0.43	0/12955	0.62	0/17710

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2523	0	2544	20	0
1	B	2523	0	2544	16	0
1	C	2523	0	2544	14	0
1	D	2523	0	2544	18	0
1	E	2523	0	2544	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
4	A	4	1	0	1	0
4	B	4	1	0	1	0
4	C	4	1	0	1	0
4	D	4	1	0	1	0
4	E	8	2	0	0	0
5	A	5	0	0	5	0
5	B	5	0	0	0	0
5	C	4	0	0	0	0
5	D	5	0	0	3	0
5	E	11	0	0	0	0
All	All	12679	6	12720	81	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 3.

All (81) 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:296:ARG:HG2	5:A:501:HOH:O	1.44	1.13
1:D:47:TRP:HD1	5:D:501:HOH:O	1.71	0.72
1:B:78:PHE:HE2	1:B:85:ARG:HD3	1.65	0.62
1:A:78:PHE:HE2	1:A:85:ARG:HD3	1.64	0.62
1:E:78:PHE:HE2	1:E:85:ARG:HD3	1.65	0.62
1:D:78:PHE:HE2	1:D:85:ARG:HD3	1.65	0.60
1:B:26:GLU:HB2	1:B:40:ASN:HB3	1.85	0.58
1:C:26:GLU:HB2	1:C:40:ASN:HB3	1.86	0.58
1:A:134:SER:HB3	1:A:139:ASN:HA	1.84	0.58
1:A:78:PHE:CE2	1:A:85:ARG:HD3	2.39	0.58
1:E:26:GLU:HB2	1:E:40:ASN:HB3	1.86	0.57
1:C:78:PHE:HE2	1:C:85:ARG:HD3	1.68	0.57
1:D:26:GLU:HB2	1:D:40:ASN:HB3	1.85	0.57
1:E:78:PHE:CE2	1:E:85:ARG:HD3	2.41	0.56
1:A:26:GLU:HB2	1:A:40:ASN:HB3	1.87	0.56
1:B:78:PHE:CE2	1:B:85:ARG:HD3	2.41	0.56
1:B:7:PRO:HG3	1:B:135:VAL:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:78:PHE:CE2	1:D:85:ARG:HD3	2.41	0.55
1:A:216:PHE:CB	5:A:501:HOH:O	2.56	0.54
1:A:216:PHE:HB2	5:A:501:HOH:O	2.06	0.54
1:A:131:ILE:HG12	1:A:181:GLU:HG2	1.90	0.54
1:C:78:PHE:CE2	1:C:85:ARG:HD3	2.44	0.53
1:A:7:PRO:HG3	1:A:135:VAL:HG21	1.92	0.52
1:D:133:ARG:HH12	1:D:177:GLU:HB2	1.75	0.52
1:D:8:PRO:HG3	5:D:501:HOH:O	2.11	0.51
1:B:239:CYS:SG	1:B:259:ILE:HG23	2.50	0.51
1:B:131:ILE:HG12	1:B:181:GLU:HG2	1.92	0.50
1:D:7:PRO:HG3	1:D:135:VAL:HG21	1.92	0.50
1:E:7:PRO:HG3	1:E:135:VAL:HG21	1.94	0.50
1:A:126:LEU:HD12	1:A:188:LEU:HD23	1.93	0.50
1:C:131:ILE:HG12	1:C:181:GLU:HG2	1.94	0.49
1:E:131:ILE:HG12	1:E:181:GLU:HG2	1.93	0.49
1:D:239:CYS:SG	1:D:259:ILE:HG23	2.53	0.49
1:C:7:PRO:HG3	1:C:135:VAL:HG21	1.93	0.49
1:D:126:LEU:HD12	1:D:188:LEU:HD23	1.94	0.49
1:E:239:CYS:SG	1:E:259:ILE:HG23	2.53	0.48
1:D:131:ILE:HG12	1:D:181:GLU:HG2	1.95	0.48
1:E:126:LEU:HD12	1:E:188:LEU:HD23	1.95	0.48
1:B:126:LEU:HD12	1:B:188:LEU:HD23	1.94	0.47
1:C:126:LEU:HD12	1:C:188:LEU:HD23	1.96	0.47
1:A:239:CYS:SG	1:A:259:ILE:HG23	2.55	0.47
1:C:239:CYS:SG	1:C:259:ILE:HG23	2.54	0.47
1:C:238:PHE:O	1:C:242:VAL:HG23	2.15	0.47
1:A:217:TRP:CE3	5:A:501:HOH:O	2.68	0.47
1:E:238:PHE:O	1:E:242:VAL:HG23	2.14	0.47
1:C:134:SER:HB3	1:C:139:ASN:HA	1.98	0.47
1:B:10:PRO:HB2	1:B:12:ALA:O	2.16	0.46
1:C:255:THR:HG23	4:C:402:MBR:BR1	2.71	0.46
1:A:10:PRO:HB2	1:A:12:ALA:O	2.15	0.46
1:A:216:PHE:C	5:A:501:HOH:O	2.54	0.46
1:A:238:PHE:O	1:A:242:VAL:HG23	2.16	0.46
1:D:238:PHE:O	1:D:242:VAL:HG23	2.16	0.45
1:B:160:TRP:CE3	1:B:190:ILE:HD12	2.51	0.45
1:D:10:PRO:HB2	1:D:12:ALA:O	2.17	0.45
1:B:238:PHE:O	1:B:242:VAL:HG23	2.16	0.45
1:E:160:TRP:CE3	1:E:190:ILE:HD12	2.53	0.44
1:B:134:SER:HB3	1:B:139:ASN:HA	1.99	0.44
1:A:279:LEU:HB2	1:A:288:ALA:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:TRP:CE3	1:A:190:ILE:HD12	2.52	0.43
1:A:242:VAL:HG11	1:A:255:THR:HG21	2.00	0.43
1:C:160:TRP:CE3	1:C:190:ILE:HD12	2.52	0.43
1:D:160:TRP:CE3	1:D:190:ILE:HD12	2.53	0.43
1:B:133:ARG:HH12	1:B:177:GLU:HB2	1.82	0.43
1:B:209:LEU:HD13	1:B:262:ILE:HG23	2.01	0.43
1:B:279:LEU:HB2	1:B:288:ALA:HB2	1.99	0.43
1:E:209:LEU:HD13	1:E:262:ILE:HG23	2.00	0.43
1:D:242:VAL:HG11	1:D:255:THR:HG21	2.00	0.43
1:D:279:LEU:HB2	1:D:288:ALA:HB2	2.00	0.43
1:A:255:THR:HG23	4:A:404:MBR:BR1	2.74	0.42
1:D:47:TRP:CD1	5:D:501:HOH:O	2.56	0.42
1:E:133:ARG:HH12	1:E:177:GLU:HB2	1.84	0.42
1:D:255:THR:HG23	4:D:403:MBR:BR1	2.74	0.42
1:C:279:LEU:HB2	1:C:288:ALA:HB2	2.01	0.42
1:E:242:VAL:HG11	1:E:255:THR:HG21	2.00	0.42
1:E:279:LEU:HB2	1:E:288:ALA:HB2	2.00	0.42
1:B:255:THR:HG23	4:B:403:MBR:BR1	2.75	0.41
1:B:242:VAL:HG11	1:B:255:THR:HG21	2.01	0.41
1:D:209:LEU:HD13	1:D:262:ILE:HG23	2.02	0.41
1:C:209:LEU:HD13	1:C:262:ILE:HG23	2.01	0.41
1:A:209:LEU:HD13	1:A:262:ILE:HG23	2.02	0.41
1:C:216:PHE:O	1:C:292:THR:HG22	2.22	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	309/327 (94%)	298 (96%)	10 (3%)	1 (0%)	41 74
1	B	309/327 (94%)	297 (96%)	11 (4%)	1 (0%)	41 74

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	309/327 (94%)	297 (96%)	11 (4%)	1 (0%)	41	74
1	D	309/327 (94%)	297 (96%)	11 (4%)	1 (0%)	41	74
1	E	309/327 (94%)	298 (96%)	10 (3%)	1 (0%)	41	74
All	All	1545/1635 (94%)	1487 (96%)	53 (3%)	5 (0%)	41	74

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	155	VAL
1	D	155	VAL
1	E	155	VAL
1	B	155	VAL
1	A	155	VAL

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	280/292 (96%)	264 (94%)	16 (6%)	20	56
1	B	280/292 (96%)	265 (95%)	15 (5%)	22	58
1	C	280/292 (96%)	263 (94%)	17 (6%)	18	54
1	D	280/292 (96%)	263 (94%)	17 (6%)	18	54
1	E	280/292 (96%)	264 (94%)	16 (6%)	20	56
All	All	1400/1460 (96%)	1319 (94%)	81 (6%)	20	55

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	14	GLU
1	A	17	THR
1	A	27	CYS
1	A	48	LYS

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Mol	Chain	Res	Type
1	A	61	VAL
1	A	83	ASN
1	A	88	ASP
1	A	111	LEU
1	A	123	SER
1	A	137	THR
1	A	147	GLU
1	A	163	GLU
1	A	178	ASP
1	A	280	LYS
1	A	304	LEU
1	B	14	GLU
1	B	17	THR
1	B	27	CYS
1	B	48	LYS
1	B	62	ARG
1	B	111	LEU
1	B	123	SER
1	B	133	ARG
1	B	137	THR
1	B	147	GLU
1	B	163	GLU
1	B	178	ASP
1	B	192	ARG
1	B	280	LYS
1	B	304	LEU
1	C	11	ILE
1	C	14	GLU
1	C	17	THR
1	C	27	CYS
1	C	48	LYS
1	C	58	ARG
1	C	62	ARG
1	C	88	ASP
1	C	111	LEU
1	C	123	SER
1	C	137	THR
1	C	147	GLU
1	C	163	GLU
1	C	178	ASP
1	C	179	ARG
1	C	280	LYS

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Mol	Chain	Res	Type
1	C	304	LEU
1	D	11	ILE
1	D	14	GLU
1	D	17	THR
1	D	27	CYS
1	D	48	LYS
1	D	62	ARG
1	D	63	VAL
1	D	88	ASP
1	D	111	LEU
1	D	123	SER
1	D	137	THR
1	D	147	GLU
1	D	163	GLU
1	D	178	ASP
1	D	179	ARG
1	D	280	LYS
1	D	304	LEU
1	E	11	ILE
1	E	14	GLU
1	E	17	THR
1	E	27	CYS
1	E	48	LYS
1	E	58	ARG
1	E	64	LYS
1	E	83	ASN
1	E	111	LEU
1	E	123	SER
1	E	133	ARG
1	E	137	THR
1	E	147	GLU
1	E	163	GLU
1	E	280	LYS
1	E	304	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	139	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 10 are monoatomic - leaving 6 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	$\# Z > 2$
4	MBR	D	403	-	3,3,3	0.05	0	3,3,3	0.14	0
4	MBR	B	403	-	3,3,3	0.08	0	3,3,3	0.14	0
4	MBR	C	402	-	3,3,3	0.09	0	3,3,3	0.18	0
4	MBR	A	404	-	3,3,3	0.06	0	3,3,3	0.15	0
4	MBR	E	404	-	3,3,3	0.08	0	3,3,3	0.10	0
4	MBR	E	403	-	3,3,3	0.06	0	3,3,3	0.06	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

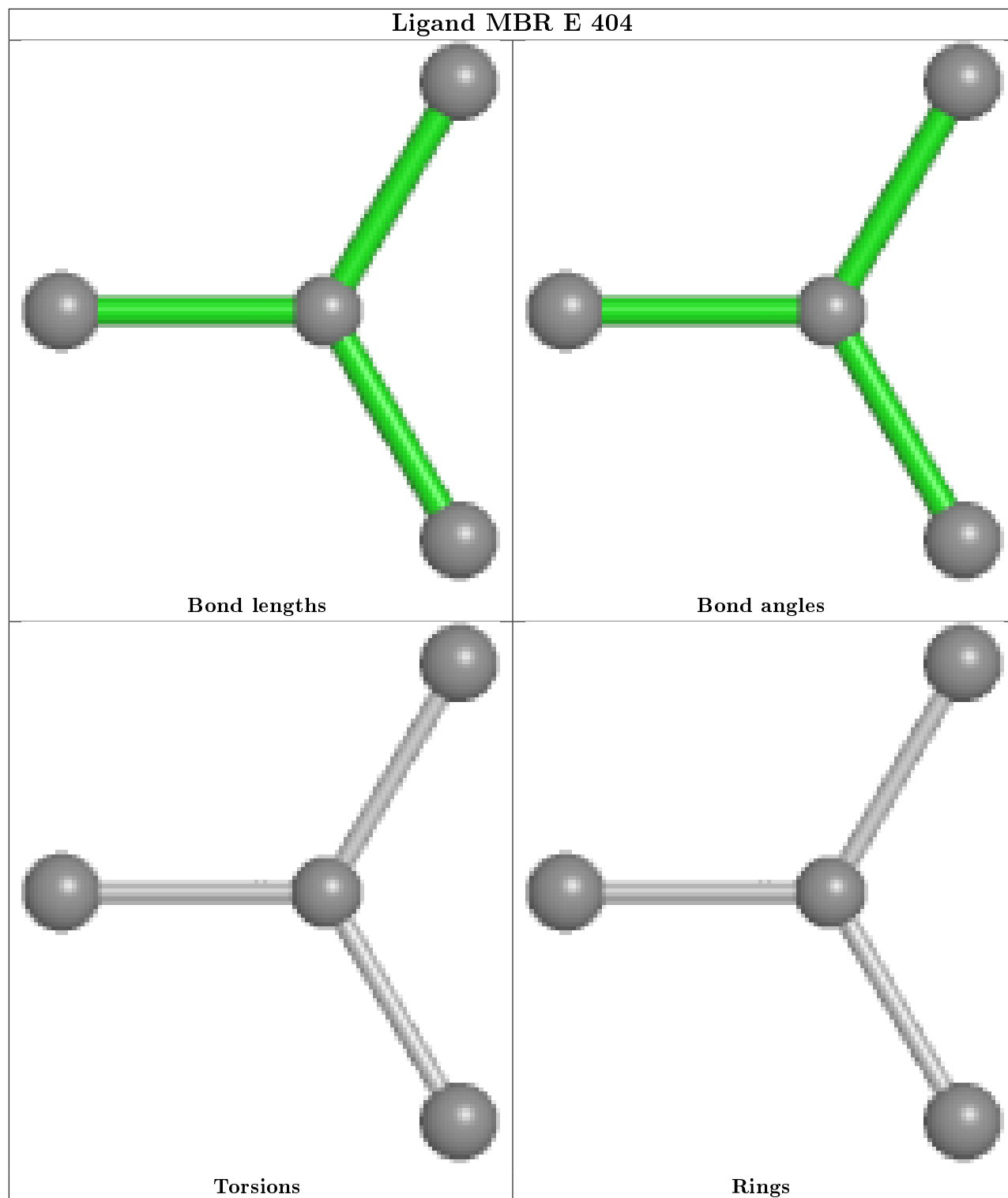
There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	403	MBR	1	0
4	B	403	MBR	1	0
4	C	402	MBR	1	0
4	A	404	MBR	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

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	311/327 (95%)	-0.18	7 (2%) 60 47	110, 139, 199, 243	0
1	B	311/327 (95%)	-0.12	9 (2%) 51 36	95, 134, 177, 216	0
1	C	311/327 (95%)	-0.10	16 (5%) 28 16	117, 145, 205, 231	0
1	D	311/327 (95%)	-0.06	14 (4%) 33 21	119, 149, 205, 253	0
1	E	311/327 (95%)	-0.05	17 (5%) 25 14	112, 145, 202, 237	0
All	All	1555/1635 (95%)	-0.10	63 (4%) 37 24	95, 142, 199, 253	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	58	ARG	6.4
1	C	64	LYS	5.5
1	C	61	VAL	5.4
1	B	62	ARG	5.2
1	E	64	LYS	5.2
1	C	283	SER	5.1
1	D	61	VAL	4.9
1	E	283	SER	4.9
1	D	57	VAL	4.8
1	D	60	GLY	4.6
1	E	58	ARG	4.4
1	C	66	TYR	4.4
1	E	62	ARG	4.4
1	C	57	VAL	4.3
1	C	56	PRO	4.3
1	D	56	PRO	4.1
1	E	61	VAL	4.0
1	D	59	SER	4.0
1	E	115	ASP	3.9
1	D	54	PHE	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	60	GLY	3.7
1	E	56	PRO	3.6
1	D	55	ASP	3.6
1	E	93	SER	3.6
1	E	94	VAL	3.4
1	B	61	VAL	3.4
1	D	64	LYS	3.3
1	D	123	SER	3.3
1	E	57	VAL	3.1
1	B	283	SER	3.1
1	C	92	ILE	3.0
1	D	139	ASN	3.0
1	D	136	ASP	2.9
1	D	191	SER	2.9
1	A	62	ARG	2.9
1	A	163	GLU	2.8
1	E	102	TYR	2.8
1	E	63	VAL	2.8
1	E	92	ILE	2.6
1	C	244	THR	2.6
1	B	312	PHE	2.5
1	A	61	VAL	2.5
1	E	312	PHE	2.5
1	D	58	ARG	2.5
1	B	56	PRO	2.5
1	E	101	GLN	2.4
1	D	135	VAL	2.4
1	C	59	SER	2.4
1	E	54	PHE	2.3
1	B	146	LEU	2.2
1	B	293	ARG	2.2
1	C	62	ARG	2.2
1	C	94	VAL	2.2
1	A	146	LEU	2.1
1	C	297	ILE	2.1
1	A	55	ASP	2.1
1	A	140	ILE	2.1
1	C	55	ASP	2.1
1	C	226	THR	2.1
1	B	297	ILE	2.1
1	E	59	SER	2.1
1	A	64	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	102	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](#)

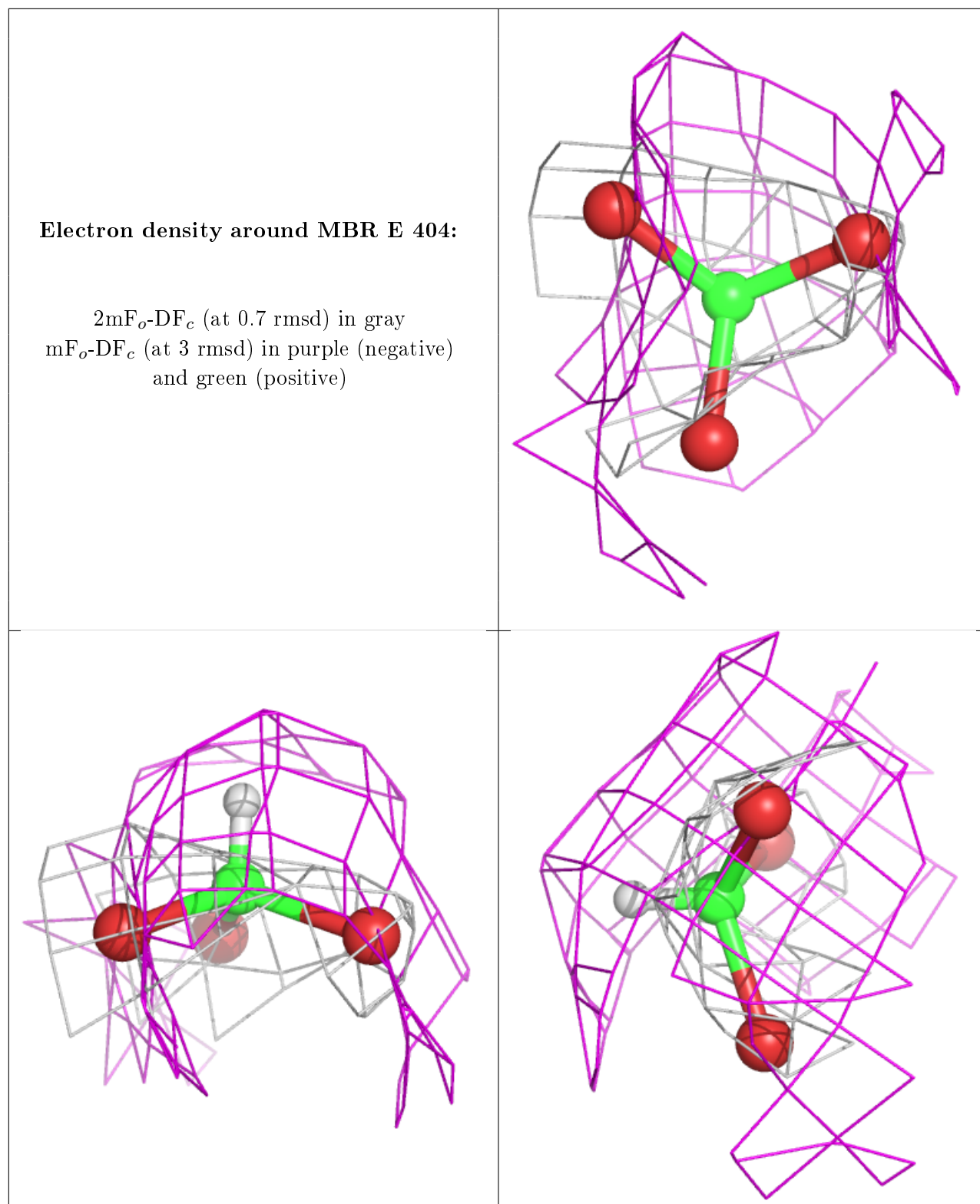
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NA	B	402	1/1	0.65	0.32	152,152,152,152	0
4	MBR	E	404	4/4	0.70	0.48	288,289,290,292	0
4	MBR	E	403	4/4	0.72	0.22	282,285,286,286	0
3	NA	E	402	1/1	0.73	0.34	146,146,146,146	0
2	CL	D	401	1/1	0.81	0.27	126,126,126,126	0
2	CL	B	401	1/1	0.81	0.25	119,119,119,119	0
2	CL	E	401	1/1	0.81	0.32	126,126,126,126	0
3	NA	A	403	1/1	0.84	1.99	132,132,132,132	0
3	NA	A	402	1/1	0.84	0.55	145,145,145,145	0
3	NA	C	401	1/1	0.85	0.19	130,130,130,130	0
3	NA	D	402	1/1	0.87	0.17	133,133,133,133	0
4	MBR	D	403	4/4	0.89	0.19	289,289,289,289	0
2	CL	A	401	1/1	0.89	0.16	102,102,102,102	0
4	MBR	B	403	4/4	0.92	0.28	251,252,252,253	0
4	MBR	A	404	4/4	0.92	0.55	252,254,255,255	0
4	MBR	C	402	4/4	0.93	0.31	254,256,256,258	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.