



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2022 – 02:10 PM JST

PDB ID : 7X5G  
Title : Nrf2 (A510Y)-MafG heterodimer bound with CsMBE2  
Authors : Sengoku, T.; Shiina, M.; Suzuki, K.; Hamada, K.; Sato, K.; Uchiyama, A.;  
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Deposited on : 2022-03-04  
Resolution : 2.30 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](mailto: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>.

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

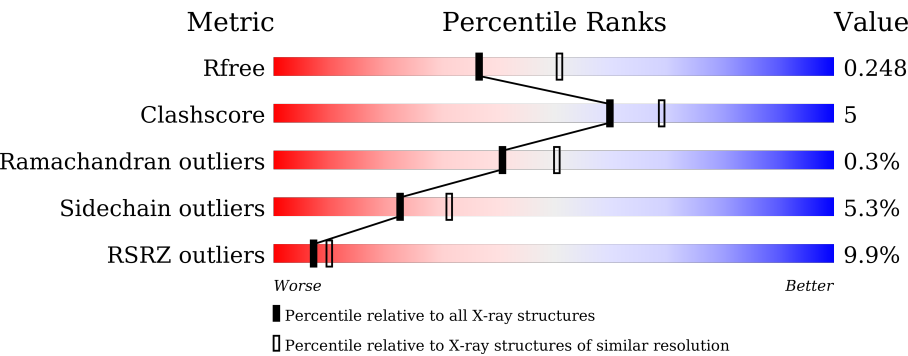
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.31.2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:  
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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.



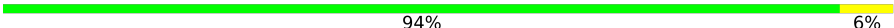

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	104	<div><div>9%</div><div><div></div><div></div><div></div><div></div></div><div>77%17%• 5%</div></div>
1	E	104	<div><div>16%</div><div><div></div><div></div><div></div><div></div></div><div>74%17%9%</div></div>
2	B	113	<div><div>%</div><div><div></div><div></div><div></div><div></div></div><div>86%6%• 7%</div></div>
2	F	113	<div><div>17%</div><div><div></div><div></div><div></div><div></div></div><div>72%19%• 8%</div></div>
3	C	16	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>81%19%</div></div>
3	G	16	<div><div></div><div><div></div><div></div><div></div><div></div></div><div>81%19%</div></div>

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Mol	Chain	Length	Quality of chain
4	D	16	 94%6%
4	H	16	 69%31%

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4634 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 Transcription factor MafG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	99	Total	C	N	O	S	0	0	0
			781	480	144	154	3			
1	E	95	Total	C	N	O	S	0	0	0
			747	455	141	148	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	initiating methionine	UNP O15525
E	20	MET	-	initiating methionine	UNP O15525

- Molecule 2 is a protein called Nuclear factor erythroid 2-related factor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	105	Total	C	N	O	S	0	0	0
			871	547	161	160	3			
2	F	104	Total	C	N	O	S	0	0	0
			814	507	148	156	3			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	448	GLY	-	expression tag	UNP Q16236
B	449	PRO	-	expression tag	UNP Q16236
B	450	HIS	-	expression tag	UNP Q16236
B	451	MET	-	expression tag	UNP Q16236
B	510	TYR	ALA	engineered mutation	UNP Q16236
F	448	GLY	-	expression tag	UNP Q16236
F	449	PRO	-	expression tag	UNP Q16236
F	450	HIS	-	expression tag	UNP Q16236
F	451	MET	-	expression tag	UNP Q16236
F	510	TYR	ALA	engineered mutation	UNP Q16236

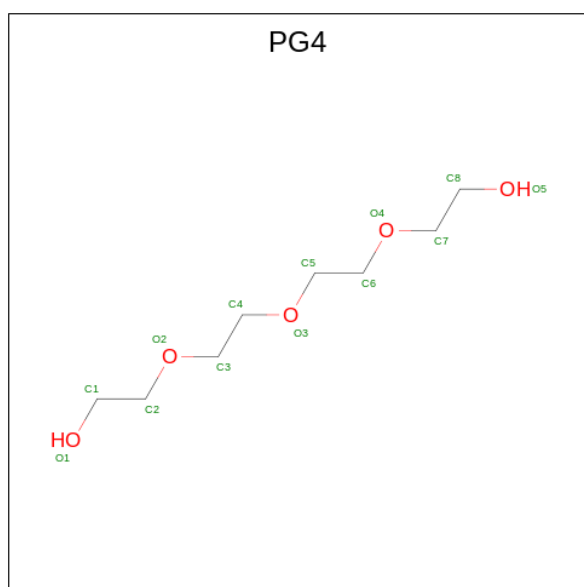
- Molecule 3 is a DNA chain called DNA (5'-D(\*GP\*CP\*TP\*GP\*CP\*TP\*GP\*AP\*GP\*TP\*CP\*AP\*CP\*TP\*GP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	16	Total	C	N	O	P	0	0	0
			325	156	57	97	15			
3	G	16	Total	C	N	O	P	0	0	0
			325	156	57	97	15			

- Molecule 4 is a DNA chain called DNA (5'-D(\*CP\*AP\*CP\*AP\*GP\*TP\*GP\*AP\*CP\*TP\*CP\*AP\*GP\*CP\*AP\*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	16	Total	C	N	O	P	0	0	0
			325	155	64	91	15			
4	H	16	Total	C	N	O	P	0	0	0
			325	155	64	91	15			

- Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			13	8	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	12	Total	O	0	0
			12	12		

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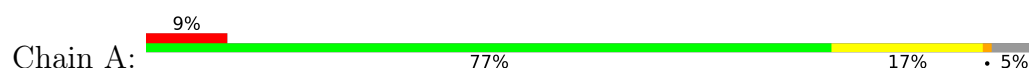
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	29	Total 29	O 29	0	0
6	C	21	Total 21	O 21	0	0
6	D	16	Total 16	O 16	0	0
6	E	8	Total 8	O 8	0	0
6	F	9	Total 9	O 9	0	0
6	G	4	Total 4	O 4	0	0
6	H	9	Total 9	O 9	0	0

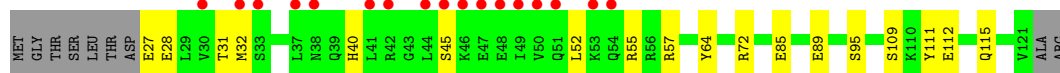
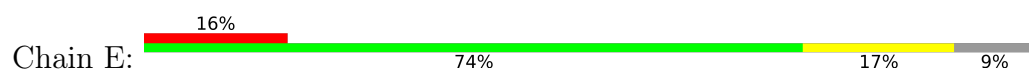
### 3 Residue-property plots [i](#)

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.

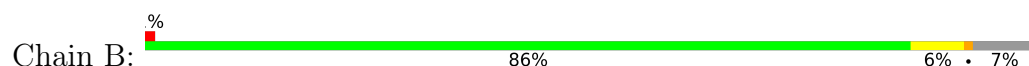
- Molecule 1: Transcription factor MafG



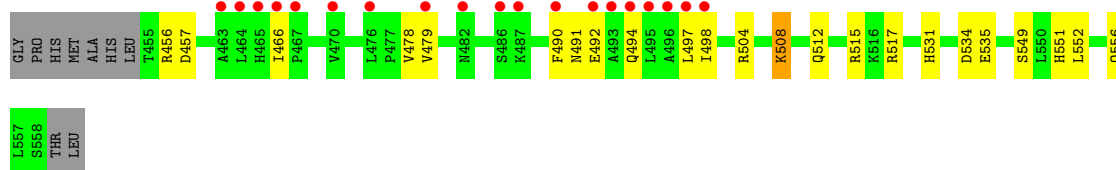
- Molecule 1: Transcription factor MafG



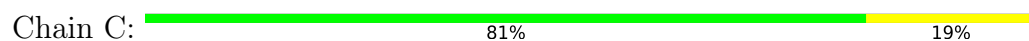
- Molecule 2: Nuclear factor erythroid 2-related factor 2



- Molecule 2: Nuclear factor erythroid 2-related factor 2

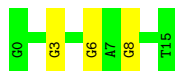
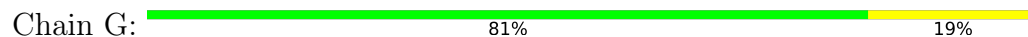


- Molecule 3: DNA (5'-D(\*GP\*CP\*TP\*GP\*CP\*TP\*GP\*AP\*GP\*TP\*CP\*AP\*CP\*TP\*GP\*T)-3')





- Molecule 3: DNA (5'-D(\*GP\*CP\*TP\*GP\*CP\*TP\*GP\*AP\*GP\*TP\*CP\*AP\*CP\*TP\*GP\*T)-3')



- Molecule 4: DNA (5'-D(\*CP\*AP\*CP\*AP\*GP\*TP\*GP\*AP\*CP\*TP\*CP\*AP\*GP\*CP\*AP\*G)-3')



- Molecule 4: DNA (5'-D(\*CP\*AP\*CP\*AP\*GP\*TP\*GP\*AP\*CP\*TP\*CP\*AP\*GP\*CP\*AP\*G)-3')





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	246.02Å 55.02Å 86.15Å 90.00° 97.30° 90.00°	Depositor
Resolution (Å)	33.89 – 2.30 33.89 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.0 (33.89-2.30) 98.0 (33.89-2.30)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.37 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.219 , 0.250 0.217 , 0.248	Depositor DCC
$R_{free}$ test set	5070 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	53.3	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4634	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.54% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: PG4

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	0.22	0/784	0.45	0/1050
1	E	0.24	0/750	0.47	0/1000
2	B	0.23	0/880	0.43	0/1174
2	F	0.23	0/821	0.43	0/1102
3	C	0.48	0/363	0.91	0/559
3	G	0.51	0/363	0.93	0/559
4	D	0.49	0/365	0.79	0/561
4	H	0.51	0/365	0.79	0/561
All	All	0.34	0/4691	0.62	0/6566

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	781	0	793	10	0
1	E	747	0	743	12	0
2	B	871	0	910	4	0
2	F	814	0	804	10	0
3	C	325	0	183	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	325	0	183	4	0
4	D	325	0	180	1	0
4	H	325	0	180	3	0
5	C	13	0	18	1	0
6	A	12	0	0	0	0
6	B	29	0	0	0	0
6	C	21	0	0	0	0
6	D	16	0	0	0	0
6	E	8	0	0	0	0
6	F	9	0	0	0	0
6	G	4	0	0	0	0
6	H	9	0	0	0	0
All	All	4634	0	3994	39	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 5.

All (39) 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)
3:C:6:DG:H2''	3:C:7:DA:H5''	1.63	0.81
1:A:41:LEU:HA	1:A:44:LEU:HD12	1.71	0.70
1:A:41:LEU:HD22	1:A:49:ILE:HG12	1.75	0.67
1:A:26:ASP:OD1	1:A:55:ARG:NH2	2.31	0.63
1:E:72:ARG:NH2	3:G:6:DG:OP2	2.35	0.60
2:F:552:LEU:O	2:F:556:GLN:HG3	2.01	0.59
2:F:478:VAL:HG21	4:H:1:DA:H5''	1.85	0.58
1:E:28:GLU:OE2	1:E:28:GLU:N	2.30	0.54
1:A:116:THR:HA	1:A:119:ARG:HD2	1.91	0.52
2:B:541:LYS:O	2:B:545:GLU:HG3	2.09	0.52
1:A:50:VAL:O	1:A:54:GLN:HG3	2.10	0.51
4:H:13:DC:H2'	4:H:14:DA:C8	2.46	0.51
1:A:45:SER:N	1:A:48:GLU:OE2	2.27	0.51
1:A:70:VAL:O	1:A:74:THR:HG23	2.11	0.51
2:F:466:ILE:HG12	2:F:498:ILE:HD11	1.93	0.51
2:B:553:LEU:O	2:B:557:LEU:HG	2.12	0.50
3:C:12:DC:H1'	5:C:101:PG4:H12	1.92	0.50
1:E:64:TYR:OH	3:G:3:DG:OP2	2.19	0.49
1:E:57:ARG:NH2	3:G:3:DG:O6	2.44	0.49
2:F:491:ASN:OD1	2:F:492:GLU:N	2.47	0.48
2:F:531:HIS:O	2:F:535:GLU:HG3	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:518:LYS:NZ	4:D:6:DG:OP2	2.33	0.46
1:E:112:GLU:HA	1:E:115:GLN:NE2	2.31	0.46
1:E:112:GLU:HA	1:E:115:GLN:HE21	1.81	0.45
1:E:111:TYR:O	1:E:115:GLN:HG3	2.16	0.45
1:E:28:GLU:O	1:E:32:MET:HG3	2.18	0.44
1:A:112:GLU:O	1:A:116:THR:HG23	2.18	0.44
2:F:508:LYS:O	2:F:512:GLN:HG3	2.17	0.43
2:F:457:ASP:OD1	2:F:504:ARG:NH2	2.49	0.43
1:A:107:LEU:HG	2:B:553:LEU:HD13	2.01	0.43
4:H:4:DG:H2'	4:H:5:DT:C6	2.53	0.43
2:F:494:GLN:O	2:F:497:LEU:N	2.52	0.42
1:A:47:GLU:O	1:A:51:GLN:HG3	2.19	0.42
2:F:552:LEU:HD12	2:F:552:LEU:HA	1.83	0.41
1:E:27:GLU:O	1:E:31:THR:HG22	2.20	0.41
1:E:40:HIS:O	1:E:40:HIS:ND1	2.45	0.41
1:E:85:GLU:O	1:E:89:GLU:HG3	2.20	0.41
2:F:515:ARG:NH2	3:G:8:DG:O6	2.49	0.41
1:E:52:LEU:HD12	1:E:52:LEU:HA	1.83	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	97/104 (93%)	94 (97%)	3 (3%)	0	100	100
1	E	93/104 (89%)	91 (98%)	2 (2%)	0	100	100
2	B	103/113 (91%)	100 (97%)	2 (2%)	1 (1%)	15	17
2	F	102/113 (90%)	101 (99%)	1 (1%)	0	100	100
All	All	395/434 (91%)	386 (98%)	8 (2%)	1 (0%)	41	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	557	LEU

### 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	83/92 (90%)	80 (96%)	3 (4%)	35	49
1	E	77/92 (84%)	73 (95%)	4 (5%)	23	32
2	B	97/104 (93%)	94 (97%)	3 (3%)	40	55
2	F	85/104 (82%)	77 (91%)	8 (9%)	8	10
All	All	342/392 (87%)	324 (95%)	18 (5%)	22	31

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ARG
1	A	95	SER
1	A	109	SER
2	B	456	ARG
2	B	489	GLN
2	B	549	SER
1	E	45	SER
1	E	55	ARG
1	E	95	SER
1	E	109	SER
2	F	456	ARG
2	F	479	VAL
2	F	490	PHE
2	F	508	LYS
2	F	517	ARG
2	F	534	ASP
2	F	549	SER
2	F	551	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
2	B	521	ASN
1	E	115	GLN
2	F	527	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z  > 2$
5	PG4	C	101	-	12,12,12	0.11	0	11,11,11	0.66	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PG4	C	101	-	-	5/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	101	PG4	O2-C3-C4-O3
5	C	101	PG4	C1-C2-O2-C3
5	C	101	PG4	C4-C3-O2-C2
5	C	101	PG4	O4-C7-C8-O5
5	C	101	PG4	C5-C6-O4-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	101	PG4	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	99/104 (95%)	0.67	9 (9%) 9 12	54, 76, 124, 144	0
1	E	95/104 (91%)	0.86	17 (17%) 1 1	46, 82, 150, 171	0
2	B	105/113 (92%)	0.31	1 (0%) 82 86	39, 61, 103, 124	0
2	F	104/113 (92%)	0.97	19 (18%) 1 1	47, 92, 153, 165	0
3	C	16/16 (100%)	0.27	0 100 100	49, 67, 90, 95	0
3	G	16/16 (100%)	0.07	0 100 100	63, 83, 98, 103	0
4	D	16/16 (100%)	0.09	0 100 100	52, 64, 98, 102	0
4	H	16/16 (100%)	-0.05	0 100 100	63, 81, 98, 105	0
All	All	467/498 (93%)	0.62	46 (9%) 7 10	39, 75, 142, 171	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	493	ALA	8.0
2	F	463	ALA	5.2
1	E	49	ILE	5.0
1	E	47	GLU	4.9
1	E	54	GLN	4.4
1	A	122	ALA	4.4
2	F	490	PHE	4.3
1	E	50	VAL	4.2
1	E	46	LYS	4.1
1	A	31	THR	3.9
2	F	464	LEU	3.9
2	F	497	LEU	3.9
1	E	41	LEU	3.6
1	E	37	LEU	3.4
1	E	42	ARG	3.4
1	E	32	MET	3.4

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Mol	Chain	Res	Type	RSRZ
2	F	465	HIS	3.3
1	A	44	LEU	3.1
2	F	496	ALA	2.9
1	E	30	VAL	2.9
1	A	25	THR	2.9
1	E	38	ASN	2.8
2	F	482	ASN	2.8
2	F	492	GLU	2.7
1	E	53	LYS	2.6
2	F	467	PRO	2.6
2	F	498	ILE	2.6
1	A	43	GLY	2.5
1	A	41	LEU	2.5
2	B	548	LYS	2.5
1	A	32	MET	2.4
2	F	487	LYS	2.4
1	A	47	GLU	2.3
2	F	466	ILE	2.3
2	F	494	GLN	2.3
2	F	479	VAL	2.3
1	E	33	SER	2.3
1	E	44	LEU	2.2
1	E	48	GLU	2.2
2	F	495	LEU	2.2
2	F	470	VAL	2.2
2	F	486	SER	2.2
1	E	51	GLN	2.0
1	E	45	SER	2.0
2	F	476	LEU	2.0
1	A	74	THR	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 monosaccharides 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
5	PG4	C	101	13/13	0.67	0.27	67,86,94,101	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.