

wwPDB X-ray Structure Validation Summary Report (i)

Nov 16, 2023 – 10:39 AM JST

PDB ID : 6LLF

Title: Biphenyl-2,2',3-triol-soaked resting complex of Oxy and Fd in carbazole

1,9a-dioxygenase

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Deposited on : 2019-12-23

Resolution : 1.93 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) 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 : 2.36

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

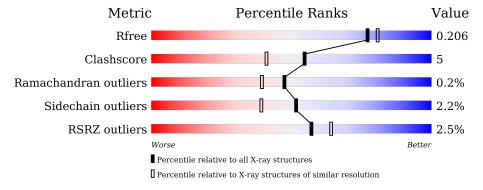
Validation Pipeline (wwPDB-VP) : 2.36

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 1.93 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	Similar resolution
	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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 <=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	392	86%		13%	
1	В	392	89%		10%	
1	С	392	89%		10%	
2	D	115	83%	7%	10%	_
2	Е	115	80%	2%	8%	
2	F	115	81%	% •	• 10%	6



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
6	EDO	С	406	-	-	X	-
7	GOL	С	408	-	-	X	-



2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 13321 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 Terminal oxygenase component of carbazole.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace	
1	Λ	388	Total	С	N	О	S	0	4	0	
1	A	300	3147	2010	537	587	13	U	4		
1	В	389	Total	С	N	О	S	0	6	0	
1	Б	369	3168	2025	543	587	13	U			
1	С	388	Total	С	N	О	S	0	0	0	
1		300	3174	2029	543	589	13		0		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	385	LEU	-	expression tag	UNP Q84II6
A	386	GLU	-	expression tag	UNP Q84II6
A	387	HIS	-	expression tag	UNP Q84II6
A	388	HIS	-	expression tag	UNP Q84II6
A	389	HIS	-	expression tag	UNP Q84II6
A	390	HIS	-	expression tag	UNP Q84II6
A	391	HIS	-	expression tag	UNP Q84II6
A	392	HIS	-	expression tag	UNP Q84II6
В	385	LEU	-	expression tag	UNP Q84II6
В	386	GLU	-	expression tag	UNP Q84II6
В	387	HIS	-	expression tag	UNP Q84II6
В	388	HIS	-	expression tag	UNP Q84II6
В	389	HIS	-	expression tag	UNP Q84II6
В	390	HIS	-	expression tag	UNP Q84II6
В	391	HIS	-	expression tag	UNP Q84II6
В	392	HIS	-	expression tag	UNP Q84II6
С	385	LEU	-	expression tag	UNP Q84II6
С	386	GLU	-	expression tag	UNP Q84II6
С	387	HIS	-	expression tag	UNP Q84II6
С	388	HIS	-	expression tag	UNP Q84II6
С	389	HIS	-	expression tag	UNP Q84II6
С	390	HIS	-	expression tag	UNP Q84II6
С	391	HIS	_	expression tag	UNP Q84II6

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Chain	Residue	Modelled	Actual	Comment	Reference
С	392	HIS	-	expression tag	UNP Q84II6

• Molecule 2 is a protein called Ferredoxin CarAc.

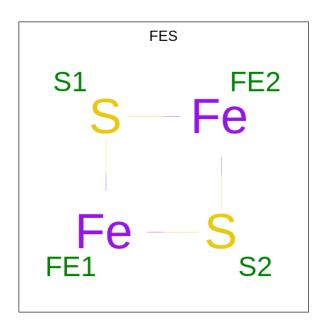
Mol	Chain	Residues		Atoms			ZeroOcc	AltConf	Trace		
2	D	104	Total	С	N	О	S	0	1	0	
2	D	104	771	485	129	150	7	0	1		
2	E	106	Total	С	N	О	S	0	1	0	
	ינו	100	788	496	132	153	7	0			
2	F	104	Total	С	N	О	S	0	9	0	
	Г	104	780	490	130	153	7				

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	108	LEU	-	expression tag	UNP Q8GI16
D	109	GLU	-	expression tag	UNP Q8GI16
D	110	HIS	-	expression tag	UNP Q8GI16
D	111	HIS	-	expression tag	UNP Q8GI16
D	112	HIS	-	expression tag	UNP Q8GI16
D	113	HIS	-	expression tag	UNP Q8GI16
D	114	HIS	-	expression tag	UNP Q8GI16
D	115	HIS	-	expression tag	UNP Q8GI16
Е	108	LEU	-	expression tag	UNP Q8GI16
Е	109	GLU	-	expression tag	UNP Q8GI16
Е	110	HIS	-	expression tag	UNP Q8GI16
Е	111	HIS	-	expression tag	UNP Q8GI16
Е	112	HIS	-	expression tag	UNP Q8GI16
Е	113	HIS	-	expression tag	UNP Q8GI16
Е	114	HIS	-	expression tag	UNP Q8GI16
Е	115	HIS	-	expression tag	UNP Q8GI16
F	108	LEU	-	expression tag	UNP Q8GI16
F	109	GLU	-	expression tag	UNP Q8GI16
F	110	HIS	-	expression tag	UNP Q8GI16
F	111	HIS	-	expression tag	UNP Q8GI16
F	112	HIS	-	expression tag	UNP Q8GI16
F	113	HIS	-	expression tag	UNP Q8GI16
F	114	HIS	-	expression tag	UNP Q8GI16
F	115	HIS	-	expression tag	UNP Q8GI16

• Molecule 3 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe₂S₂).





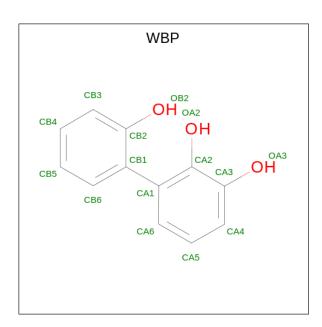
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe S 4 2 2	0	0
3	В	1	Total Fe S 4 2 2	0	0
3	С	1	Total Fe S 4 2 2	0	0
3	D	1	Total Fe S 4 2 2	0	0
3	E	1	Total Fe S 4 2 2	0	0
3	F	1	Total Fe S 4 2 2	0	0

• Molecule 4 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	${f Atoms}$	ZeroOcc	AltConf
4	A	1	Total Fe 1 1	0	0
4	В	1	Total Fe 1 1	0	0
4	С	1	Total Fe 1 1	0	0

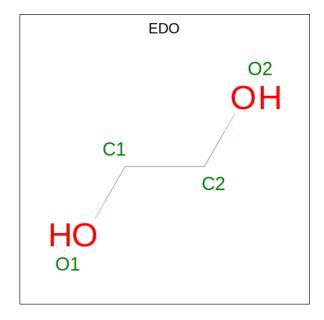
• Molecule 5 is 3-(2-hydroxyphenyl)benzene-1,2-diol (three-letter code: WBP) (formula: $C_{12}H_{10}O_3$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
5	Δ	1	Total C		0	1	
	Λ	1	30 24	6	U		
5	В	1	Total C	О	0	0	
	Б	1	15 12	3	0		
5	C	1	Total C	О	0	0	
		1	15 12	3		U	

 \bullet Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $\mathrm{C_2H_6O_2}).$



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

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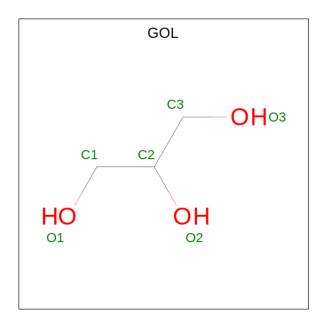


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Mol	Chain	$oxed{ ext{Residues} }$	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	В	1	Total C O 4 2 2	0	0
6	В	1	Total C O 4 2 2	0	0
6	В	1	Total C O 4 2 2	0	0
6	В	1	Total C O 4 2 2	0	0
6	С	1	Total C O 4 2 2	0	0
6	С	1	Total C O 4 2 2	0	0
6	С	1	Total C O 4 2 2	0	0
6	С	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	Е	1	Total C O 4 2 2	0	0
6	F	1	Total C O 4 2 2	0	0

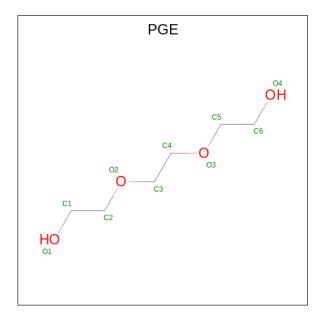
 \bullet Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: $\mathrm{C_3H_8O_3}).$





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 6 3 3	0	0
7	A	1	Total C O 6 3 3	0	0
7	В	1	Total C O 6 3 3	0	0
7	С	1	Total C O 6 3 3	0	0

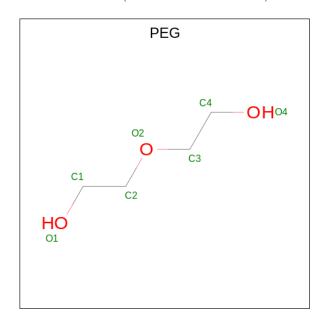
 \bullet Molecule 8 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $\mathrm{C_6H_{14}O_4}).$





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 10 6 4	0	0

 $\bullet \ \ Molecule \ 9 \ is \ DI(HYDROXYETHYL)ETHER \ (three-letter \ code: \ PEG) \ (formula: \ C_4H_{10}O_3).$



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	В	1	Total 7	C 4	O 3	0	0

• Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	371	Total O 371 371	0	0
10	В	378	Total O 378 378	0	0
10	С	376	Total O 376 376	0	0
10	D	32	Total O 32 32	0	0
10	E	68	Total O 68 68	0	0
10	F	72	Total O 72 72	0	0

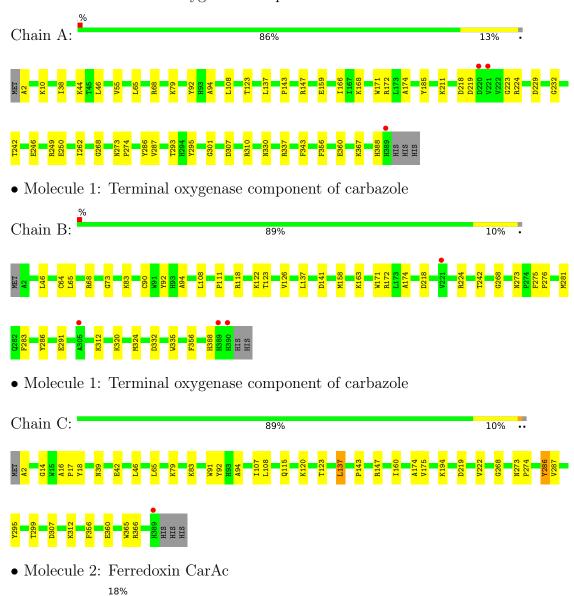


Chain D:

3 Residue-property plots (i)

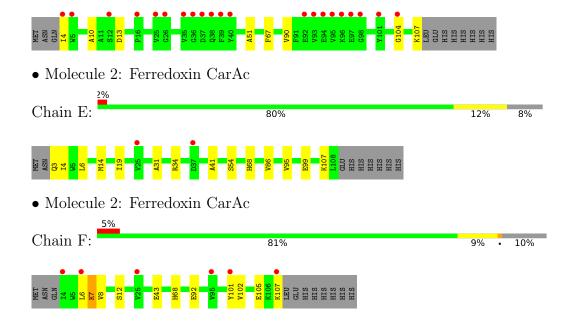
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Terminal oxygenase component of carbazole



83%







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	98.29Å 89.68Å 105.02Å	Depositor
a, b, c, α , β , γ	90.00° 104.08° 90.00°	Depositor
Resolution (Å)	44.88 - 1.93	Depositor
rtesolution (A)	44.84 - 1.93	EDS
% Data completeness	96.3 (44.88-1.93)	Depositor
(in resolution range)	96.4 (44.84-1.93)	EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.68 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
P. P.	0.158 , 0.200	Depositor
R, R_{free}	0.167 , 0.206	DCC
R_{free} test set	6252 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 45.2	EDS
L-test for twinning ²	$ < L >=0.50, < L^2>=0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13321	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.51% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: FE2, PEG, GOL, WBP, PGE, FES, EDO

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	Bo	nd lengths	Bond angles		
MIOI	Cham	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.74	$1/3246 \ (0.0\%)$	0.86	$1/4406 \ (0.0\%)$	
1	В	0.72	0/3276	0.84	0/4447	
1	С	0.74	$1/3285 \ (0.0\%)$	0.85	0/4455	
2	D	0.75	0/790	0.86	0/1074	
2	Е	0.77	0/807	0.88	0/1097	
2	F	0.80	1/799~(0.1%)	0.88	0/1086	
All	All	0.74	3/12203 (0.0%)	0.85	1/16565 (0.0%)	

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
2	F	43	GLU	CD-OE1	5.56	1.31	1.25
1	A	360	GLU	CD-OE1	5.31	1.31	1.25
1	С	360	GLU	CD-OE1	5.14	1.31	1.25

All (1) bond angle outliers are listed below:

N	/Iol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
	1	Α	224	ARG	CG-CD-NE	-5.58	100.08	111.80

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	3147	0	3049	32	0
1	В	3168	0	3071	29	0
1	С	3174	0	3088	27	0
2	D	771	0	750	4	0
2	Е	788	0	769	11	0
2	F	780	0	755	8	0
3	A	4	0	0	0	0
3	В	4	0	0	0	0
3	С	4	0	0	0	0
3	D	4	0	0	0	0
3	E	4	0	0	1	0
3	F	4	0	0	1	0
4	A	1	0	0	0	0
4	В	1	0	0	0	0
4	С	1	0	0	0	0
5	A	30	0	0	3	0
5	В	15	0	0	2	0
5	С	15	0	0	3	0
6	A	24	0	36	4	0
6	В	16	0	24	6	0
6	С	16	0	24	6	0
6	D	4	0	6	1	0
6	\mathbf{E}	4	0	6	0	0
6	F	4	0	6	0	0
7	A	12	0	16	1	0
7	В	6	0	8	0	0
7	С	6	0	8	5	0
8	A	10	0	14	3	0
9	В	7	0	10	2	0
10	A	371	0	0	10	0
10	В	378	0	0	11	0
10	С	376	0	0	7	0
10	D	32	0	0	1	0
10	Ε	68	0	0	1	0
10	F	72	0	0	2	0
All	All	13321	0	11640	122	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.

The worst 5 of 122 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
5:C:403:WBP:CA2	10:C:501:HOH:O	2.08	1.01
1:B:68:ARG:HH11	6:B:405:EDO:H11	1.37	0.87
1:A:218:ASP:OD1	1:A:218:ASP:O	1.90	0.87
1:A:79:LYS:HE3	10:A:832:HOH:O	1.82	0.78
6:B:407:EDO:H21	10:B:714:HOH:O	1.82	0.77

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	Perce	ntiles
1	A	390/392 (100%)	370 (95%)	19 (5%)	1 (0%)	41	32
1	В	393/392 (100%)	375 (95%)	17 (4%)	1 (0%)	41	32
1	С	394/392 (100%)	373 (95%)	20 (5%)	1 (0%)	41	32
2	D	103/115 (90%)	99 (96%)	4 (4%)	0	100	100
2	E	105/115 (91%)	103 (98%)	2 (2%)	0	100	100
2	F	104/115 (90%)	100 (96%)	4 (4%)	0	100	100
All	All	1489/1521 (98%)	1420 (95%)	66 (4%)	3 (0%)	47	39

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	268	GLY
1	A	268	GLY
1	С	268	GLY

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	339/339 (100%)	332 (98%)	7 (2%)	53	41	
1	В	342/339 (101%)	336 (98%)	6 (2%)	59	47	
1	С	343/339 (101%)	334 (97%)	9 (3%)	46	32	
2	D	83/93 (89%)	83 (100%)	0	100	100	
2	E	85/93 (91%)	83 (98%)	2 (2%)	49	36	
2	F	84/93 (90%)	81 (96%)	3 (4%)	35	20	
All	All	1276/1296 (98%)	1249 (98%)	27 (2%)	52	41	

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	92	TYR
1	С	222	VAL
2	F	6	LEU
1	С	219	ASP
1	С	286	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	В	234	HIS
1	В	388	HIS
2	F	38	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)

Of 36 ligands modelled in this entry, 3 are monoatomic - leaving 33 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).

Mal	Т	Clasia.	Dag	T ! 1-	Во	ond leng	$_{ m ths}$	В	ond ang	gles
Mol	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	GOL	A	407	-	5,5,5	0.10	0	5,5,5	0.37	0
6	EDO	Е	202	_	3,3,3	0.18	0	2,2,2	0.36	0
6	EDO	A	404	-	3,3,3	0.29	0	2,2,2	0.67	0
3	FES	Е	201	2	0,4,4	-	-	-		
7	GOL	С	408	-	5,5,5	0.21	0	5,5,5	0.64	0
8	PGE	A	408	-	9,9,9	0.34	0	8,8,8	0.20	0
6	EDO	A	406	-	3,3,3	0.25	0	2,2,2	0.31	0
6	EDO	A	405	-	3,3,3	0.11	0	2,2,2	0.12	0
3	FES	В	401	1	0,4,4	-	-	-		
7	GOL	A	412	-	5,5,5	0.12	0	5,5,5	0.31	0
6	EDO	A	410	-	3,3,3	0.13	0	2,2,2	0.27	0
5	WBP	С	403	-	16,16,16	2.70	3 (18%)	22,22,22	0.95	1 (4%)
6	EDO	С	407	-	3,3,3	0.16	0	2,2,2	0.27	0
6	EDO	С	406	-	3,3,3	0.66	0	2,2,2	0.50	0
6	EDO	F	202	-	3,3,3	0.12	0	2,2,2	0.24	0
6	EDO	С	404	-	3,3,3	0.21	0	2,2,2	0.24	0
6	EDO	С	405	_	3,3,3	0.42	0	2,2,2	0.59	0
6	EDO	В	409	-	3,3,3	0.26	0	2,2,2	0.65	0
9	PEG	В	404	-	6,6,6	0.19	0	5,5,5	0.43	0
6	EDO	В	407	-	3,3,3	0.39	0	2,2,2	0.36	0
6	EDO	В	405	-	3,3,3	0.35	0	2,2,2	0.08	0
6	EDO	В	408	-	3,3,3	0.06	0	2,2,2	0.40	0
6	EDO	A	409	-	3,3,3	0.13	0	2,2,2	0.31	0
5	WBP	A	403[B]	-	16,16,16	2.38	3 (18%)	22,22,22	1.00	1 (4%)
6	EDO	A	411	-	3,3,3	0.18	0	2,2,2	0.20	0
3	FES	A	401	1	0,4,4	-	-	-		
7	GOL	В	406	-	5,5,5	0.10	0	5,5,5	0.22	0



Mol	Т	Chain	Dog	Res Link		ond leng	ths	В	ond ang	les	
MIOI	Type	Chain	nes	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	WBP	В	403	-	16,16,16	2.78	3 (18%)	22,22,22	0.84	0	
6	EDO	D	202	-	3,3,3	0.36	0	2,2,2	0.22	0	
3	FES	D	201	2	0,4,4	-	-	-			
5	WBP	A	403[A]	-	16,16,16	2.40	3 (18%)	22,22,22	1.19	3 (13%)	
3	FES	F	201	2	0,4,4	-	-	-			
3	FES	С	401	1	0,4,4	-	-	-			

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
7	GOL	A	407	-	-	2/4/4/4	-
6	EDO	Е	202	-	-	1/1/1/1	-
6	EDO	A	404	-	-	1/1/1/1	-
7	GOL	С	408	-	-	0/4/4/4	-
3	FES	Е	201	2	-	-	0/1/1/1
8	PGE	A	408	-	-	4/7/7/7	-
6	EDO	A	406	-	-	1/1/1/1	-
6	EDO	A	405	-	-	0/1/1/1	-
3	FES	В	401	1	-	-	0/1/1/1
7	GOL	A	412	-	-	1/4/4/4	-
6	EDO	A	410	-	-	0/1/1/1	-
5	WBP	С	403	-	-	0/4/4/4	0/2/2/2
6	EDO	С	407	-	-	0/1/1/1	-
6	EDO	С	406	-	-	0/1/1/1	-
6	EDO	F	202	-	-	0/1/1/1	-
6	EDO	С	404	-	-	0/1/1/1	-
6	EDO	С	405	-	-	0/1/1/1	-
6	EDO	В	409	-	-	1/1/1/1	-
9	PEG	В	404	-	-	1/4/4/4	-
6	EDO	В	407	-	-	1/1/1/1	-
6	EDO	В	405	-	-	1/1/1/1	-
6	EDO	В	408	-	-	0/1/1/1	-
6	EDO	A	409	_	-	0/1/1/1	-
5	WBP	A	403[B]	-	-	1/4/4/4	0/2/2/2
6	EDO	A	411	-	-	1/1/1/1	-
7	GOL	В	406	-	-	4/4/4/4	-
3	FES	A	401	1	-	-	0/1/1/1
6	EDO	D	202	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	WBP	В	403	_	=	0/4/4/4	0/2/2/2
3	FES	D	201	2	-	-	0/1/1/1
5	WBP	A	403[A]	_	-	0/4/4/4	0/2/2/2
3	FES	F	201	2	-	-	0/1/1/1
3	FES	С	401	1	-	-	0/1/1/1

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	Ideal(Å)
5	В	403	WBP	CA3-CA2	8.75	1.50	1.40
5	С	403	WBP	CA3-CA2	7.72	1.49	1.40
5	A	403[A]	WBP	CA3-CA2	7.31	1.48	1.40
5	A	403[B]	WBP	CA3-CA2	6.57	1.47	1.40
5	С	403	WBP	CA1-CA2	6.37	1.51	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\mathrm{Ideal}(^{o})$
5	A	403[A]	WBP	CB6-CB1-CA1	3.12	125.18	118.74
5	A	403[A]	WBP	CA1-CB1-CB2	-2.69	117.00	123.14
5	A	403[B]	WBP	CA1-CA2-CA3	-2.38	118.51	120.03
5	A	403[A]	WBP	CA4-CA3-CA2	-2.18	117.83	120.06
5	С	403	WBP	CA1-CA2-CA3	-2.02	118.73	120.03

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	407	GOL	C1-C2-C3-O3
8	A	408	PGE	O3-C5-C6-O4
7	A	412	GOL	C1-C2-C3-O3
7	В	406	GOL	O1-C1-C2-C3
7	В	406	GOL	C1-C2-C3-O3

There are no ring outliers.

18 monomers are involved in 38 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	407	GOL	1	0
6	A	404	EDO	1	0
3	Е	201	FES	1	0

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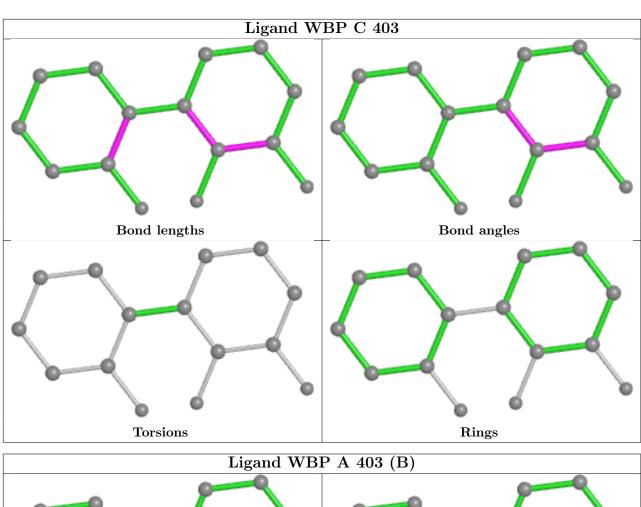


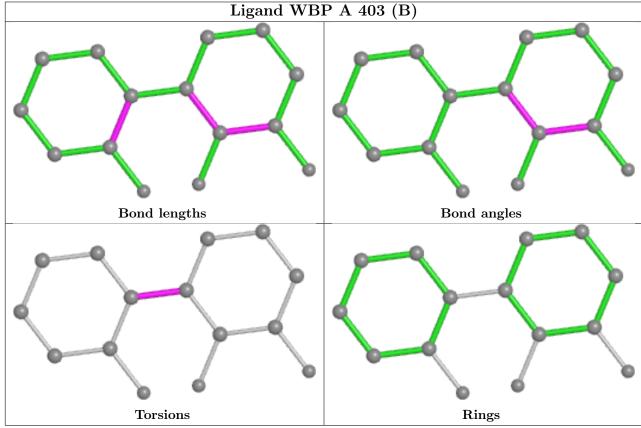
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	С	408	GOL	5	0
8	A	408	PGE	3	0
6	A	406	EDO	3	0
5	С	403	WBP	3	0
6	С	406	EDO	4	0
6	С	405	EDO	2	0
6	В	409	EDO	1	0
9	В	404	PEG	2	0
6	В	407	EDO	2	0
6	В	405	EDO	3	0
5	A	403[B]	WBP	1	0
5	В	403	WBP	2	0
6	D	202	EDO	1	0
5	A	403[A]	WBP	2	0
3	F	201	FES	1	0

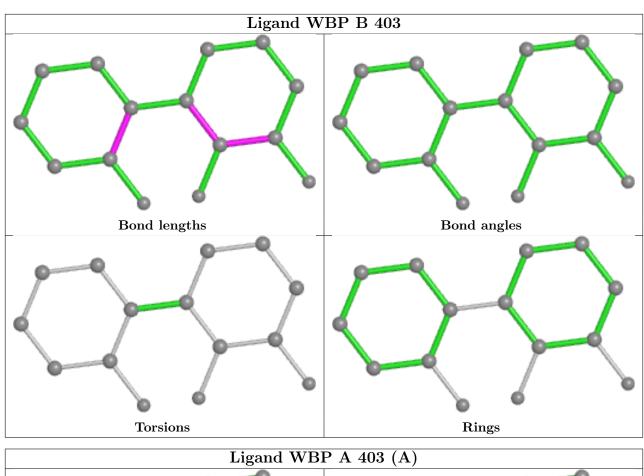
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

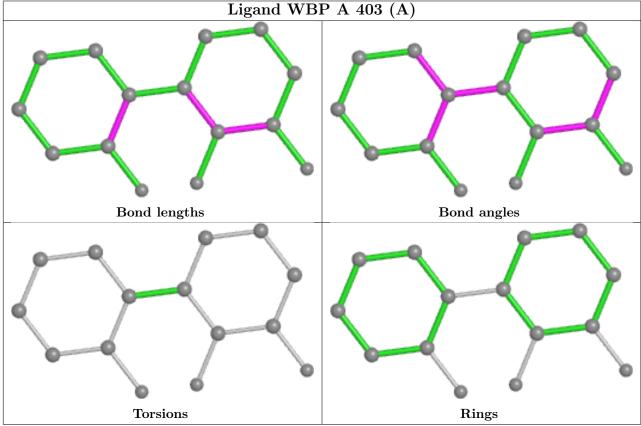














5.7 Other polymers (i)

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 (i)

6.1 Protein, DNA and RNA chains (i)

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^{th} 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>	$\# \mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q<0.9
1	A	388/392 (98%)	-0.42	3 (0%) 86 89	10, 17, 41, 58	0
1	В	389/392 (99%)	-0.36	4 (1%) 82 86	11, 21, 47, 78	0
1	С	388/392 (98%)	-0.54	1 (0%) 94 96	12, 20, 41, 65	0
2	D	104/115 (90%)	0.80	21 (20%) 1 1	15, 38, 74, 79	0
2	E	106/115 (92%)	-0.00	2 (1%) 66 72	16, 29, 50, 64	0
2	F	104/115 (90%)	0.07	6 (5%) 23 29	15, 29, 52, 75	0
All	All	1479/1521 (97%)	-0.29	37 (2%) 57 64	10, 21, 50, 79	0

The worst 5 of 37 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	93	VAL	4.9
2	D	4	ILE	4.5
2	D	38	GLN	4.4
2	D	98	GLY	4.2
2	D	37	ASP	3.9

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.



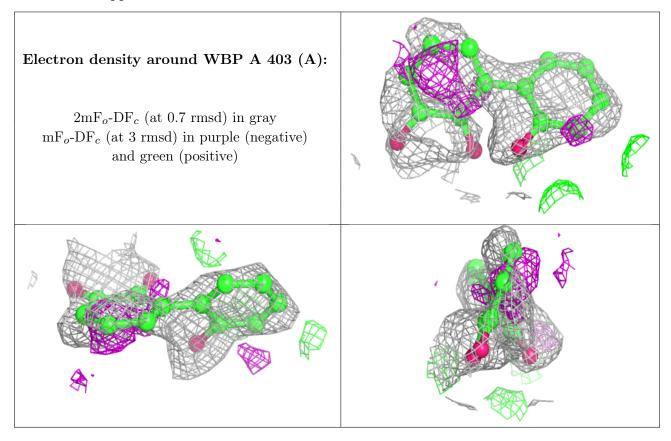
6.4 Ligands (i)

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^{th} 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.

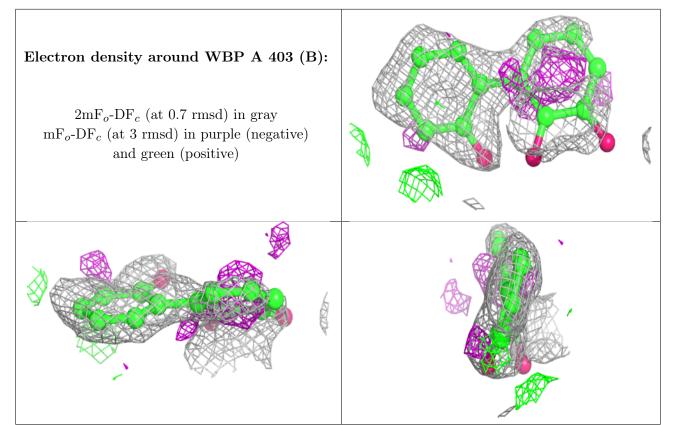
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q<0.9
6	EDO	A	405	4/4	0.81	0.14	45,47,47,49	0
6	EDO	A	409	4/4	0.81	0.20	54,58,60,63	0
8	PGE	A	408	10/10	0.82	0.17	28,40,47,49	0
6	EDO	Е	202	4/4	0.84	0.12	47,57,57,59	0
6	EDO	В	405	4/4	0.85	0.16	38,38,40,45	0
5	WBP	A	403[A]	15/15	0.85	0.29	41,48,50,51	15
5	WBP	A	403[B]	15/15	0.85	0.29	28,31,33,34	15
6	EDO	D	202	4/4	0.86	0.28	35,42,45,46	0
5	WBP	В	403	15/15	0.86	0.20	46,52,63,63	0
7	GOL	В	406	6/6	0.86	0.18	53,54,55,58	0
6	EDO	С	407	4/4	0.86	0.21	49,52,54,54	0
6	EDO	В	408	4/4	0.87	0.30	49,52,54,56	0
6	EDO	A	411	4/4	0.87	0.12	48,49,51,51	0
5	WBP	С	403	15/15	0.88	0.20	39,44,50,60	0
9	PEG	В	404	7/7	0.88	0.14	47,50,53,54	0
6	EDO	A	410	4/4	0.89	0.14	41,49,49,50	0
6	EDO	В	409	4/4	0.89	0.21	47,51,55,56	0
7	GOL	A	412	6/6	0.90	0.15	50,51,52,53	0
6	EDO	A	404	4/4	0.90	0.27	36,37,39,40	0
6	EDO	F	202	4/4	0.91	0.24	58,58,60,63	0
6	EDO	A	406	4/4	0.92	0.20	40,41,43,45	0
7	GOL	A	407	6/6	0.92	0.15	23,46,49,51	0
7	GOL	С	408	6/6	0.93	0.15	27,41,45,54	0
6	EDO	С	404	4/4	0.95	0.10	30,40,40,43	0
6	EDO	В	407	4/4	0.96	0.21	31,34,35,39	0
6	EDO	С	405	4/4	0.97	0.13	31,31,32,33	0
6	EDO	С	406	4/4	0.97	0.15	24,26,27,29	0
4	FE2	В	402	1/1	0.99	0.04	32,32,32,32	0
3	FES	С	401	4/4	0.99	0.07	16,16,18,18	0
4	FE2	С	402	1/1	1.00	0.02	24,24,24,24	0
3	FES	A	401	4/4	1.00	0.09	12,12,12,13	0
3	FES	D	201	4/4	1.00	0.05	16,16,17,19	0
3	FES	Е	201	4/4	1.00	0.07	17,17,18,18	0
3	FES	F	201	4/4	1.00	0.05	13,14,14,16	0
4	FE2	A	402	1/1	1.00	0.03	23,23,23,23	0
3	FES	В	401	4/4	1.00	0.09	11,11,12,12	0



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

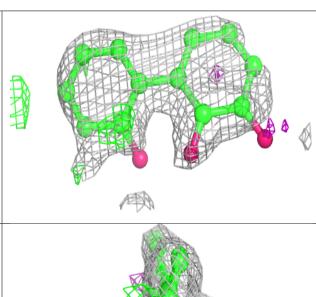


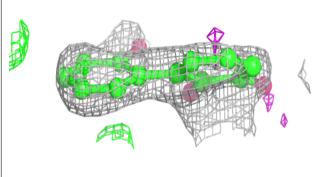


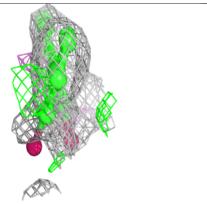


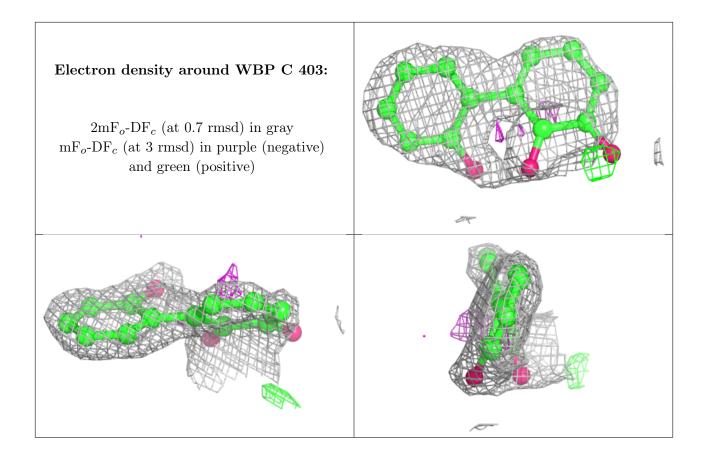
Electron density around WBP B 403:

 $2 {\rm mF}_o\text{-}{\rm DF}_c$ (at 0.7 rmsd) in gray ${\rm mF}_o\text{-}{\rm DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)









6.5 Other polymers (i)

There are no such residues in this entry.

