



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:24 PM EDT

PDB ID : 3P0G / pdb_00003p0g
Title : Structure of a nanobody-stabilized active state of the beta2 adrenoceptor
Authors : Rasmussen, S.G.F.; Choi, H.-J.; Fung, J.J.; Pardon, E.; Casarosa, P.; Chae, P.S.; DeVree, B.T.; Rosenbaum, D.M.; Thian, F.S.; Kobilka, T.S.; Schnapp, A.; Konetzki, I.; Sunahara, R.K.; Gellman, S.H.; Pautsch, A.; Steyaert, J.; Weis, W.I.; Kobilka, B.K.
Deposited on : 2010-09-28
Resolution : 3.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references](#) ①) 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

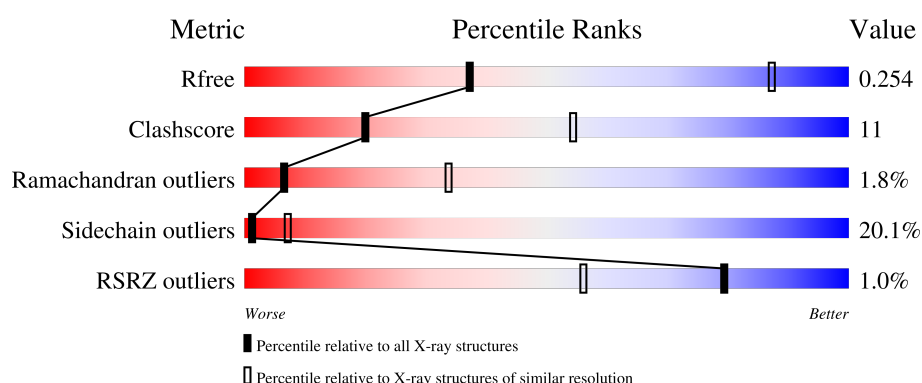
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	501	% 35% 16% 5% 43%
2	B	126	% 54% 35% 6% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3238 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 Beta-2 adrenergic receptor, Lysozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2290	1519	366	386	19			

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	ASP	-	expression tag	UNP P07550
A	-6	TYR	-	expression tag	UNP P07550
A	-5	LYS	-	expression tag	UNP P07550
A	-4	ASP	-	expression tag	UNP P07550
A	-3	ASP	-	expression tag	UNP P07550
A	-2	ASP	-	expression tag	UNP P07550
A	-1	ASP	-	expression tag	UNP P07550
A	0	ALA	-	expression tag	UNP P07550
A	16	ARG	GLY	SEE REMARK 999	UNP P07550
A	27	GLN	GLU	SEE REMARK 999	UNP P07550
A	187	GLU	ASN	engineered mutation	UNP P07550
A	229D	THR	CYS	conflict	UNP P00720
A	230U	ALA	CYS	conflict	UNP P00720

- Molecule 2 is a protein called Camelid Antibody Fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	121	Total	C	N	O	S	0	0	0
			921	573	163	181	4			

There are 6 discrepancies between the modelled and reference sequences:

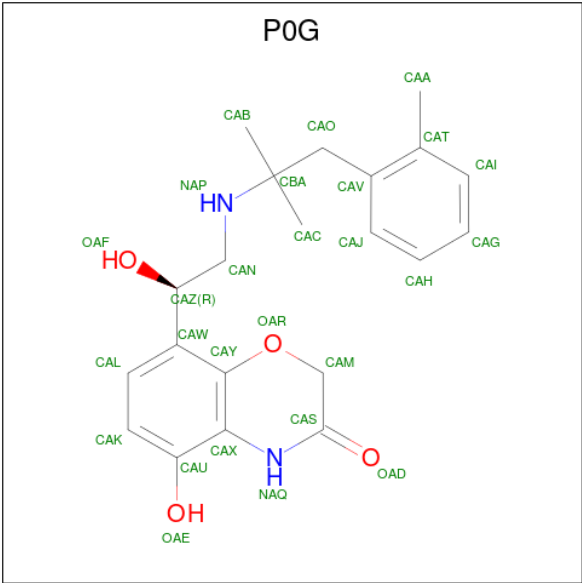
Chain	Residue	Modelled	Actual	Comment	Reference
B	121	HIS	-	expression tag	PDB 3P0G
B	122	HIS	-	expression tag	PDB 3P0G
B	123	HIS	-	expression tag	PDB 3P0G

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Chain	Residue	Modelled	Actual	Comment	Reference
B	124	HIS	-	expression tag	PDB 3P0G
B	125	HIS	-	expression tag	PDB 3P0G
B	126	HIS	-	expression tag	PDB 3P0G

- Molecule 3 is 8-[(1R)-2-{[1,1-dimethyl-2-(2-methylphenyl)ethyl]amino}-1-hydroxyethyl]-5-hydroxy-2H-1,4-benzoxazin-3(4H)-one (CCD ID: P0G) (formula: C₂₁H₂₆N₂O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	27	21	2	4	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	236.69Å 45.66Å 71.38Å 90.00° 102.34° 90.00°	Depositor
Resolution (Å)	37.00 – 3.50 37.00 – 3.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (37.00-3.50) 94.7 (37.00-3.50)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 3.48Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.8.0, BUSTER 2.8.0	Depositor
R, R_{free}	0.235 , 0.308 0.255 , 0.254	Depositor DCC
R_{free} test set	714 reflections (7.43%)	wwPDB-VP
Wilson B-factor (Å ²)	74.4	Xtriage
Anisotropy	0.847	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	3238	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.42% 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: P0G

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	0.89	0/2348	1.57	23/3197 (0.7%)
2	B	0.77	0/940	1.34	9/1274 (0.7%)
All	All	0.86	0/3288	1.51	32/4471 (0.7%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	ILE	N-CA-C	6.82	117.58	110.62
2	B	38	ARG	N-CA-C	6.54	119.17	108.52
1	A	149	LYS	N-CA-C	-6.33	104.26	112.68
2	B	101	GLY	CA-C-N	6.15	132.76	121.70
2	B	101	GLY	C-N-CA	6.15	132.76	121.70
1	A	129	VAL	CA-C-N	6.00	128.24	120.44
1	A	129	VAL	C-N-CA	6.00	128.24	120.44
1	A	112	ILE	CA-C-N	5.93	128.54	120.54
1	A	112	ILE	C-N-CA	5.93	128.54	120.54
2	B	67	PHE	CA-CB-CG	5.78	119.58	113.80
1	A	273	LYS	CA-C-N	5.70	128.17	120.65
1	A	273	LYS	C-N-CA	5.70	128.17	120.65
1	A	273	LYS	N-CA-C	-5.63	105.14	111.28
1	A	301	ASN	CA-CB-CG	5.59	118.19	112.60
1	A	332	PHE	N-CA-C	-5.52	105.18	111.14
2	B	97	VAL	N-CA-CB	5.50	118.73	111.25
2	B	29	PHE	CA-C-N	5.45	128.34	120.38
2	B	29	PHE	C-N-CA	5.45	128.34	120.38
2	B	103	VAL	N-CA-C	5.39	120.55	109.34
1	A	87	VAL	N-CA-CB	5.33	114.60	110.45
1	A	36	MET	CA-C-N	5.21	126.69	120.13
1	A	36	MET	C-N-CA	5.21	126.69	120.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	42	LEU	CA-C-N	5.19	127.29	120.60
1	A	42	LEU	C-N-CA	5.19	127.29	120.60
1	A	155	LEU	CA-C-N	5.19	127.54	120.54
1	A	155	LEU	C-N-CA	5.19	127.54	120.54
1	A	209	TYR	N-CA-C	5.18	116.92	111.28
1	A	140	LYS	CA-C-N	5.12	127.90	120.79
1	A	140	LYS	C-N-CA	5.12	127.90	120.79
2	B	78	VAL	N-CA-C	5.07	115.26	108.17
1	A	169	ILE	CA-C-N	5.05	127.04	120.28
1	A	169	ILE	C-N-CA	5.05	127.04	120.28

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	2290	0	2334	51	0
2	B	921	0	875	25	0
3	A	27	0	25	8	0
All	All	3238	0	3234	74	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 11.

All (74) 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:109:TRP:HZ3	3:A:366:P0G:HAAA	1.33	0.93
1:A:293:ASN:HD21	3:A:366:P0G:HAMA	1.46	0.79
1:A:309:ILE:HG12	3:A:366:P0G:HAH	1.71	0.73
1:A:168:PRO:HB2	1:A:199:TYR:HE1	1.56	0.70
1:A:72:ILE:HD12	1:A:326:TYR:HB3	1.74	0.68
1:A:109:TRP:CZ3	3:A:366:P0G:HAAA	2.24	0.67
1:A:71:PHE:HD1	1:A:154:ILE:HD11	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:GLN:HG2	1:A:302:LEU:HD13	1.80	0.64
2:B:59:TYR:HE1	2:B:69:ILE:H	1.45	0.62
1:A:333:ARG:O	1:A:337:GLN:HG2	1.99	0.61
1:A:135:ILE:HG12	2:B:100:TYR:HB3	1.81	0.61
2:B:90:THR:HG23	2:B:117:THR:HA	1.83	0.60
1:A:131:ARG:HD2	2:B:103:VAL:O	2.02	0.59
2:B:2:VAL:HB	2:B:25:SER:O	2.02	0.58
1:A:125:CYS:HB2	1:A:215:MET:HG3	1.85	0.58
1:A:163:LEU:O	1:A:168:PRO:HD3	2.03	0.57
2:B:34:MET:HB2	2:B:78:VAL:HG11	1.88	0.55
2:B:47:LEU:HD11	2:B:50:ALA:HB2	1.90	0.54
1:A:168:PRO:HB2	1:A:199:TYR:CE1	2.39	0.53
1:A:117:VAL:O	1:A:121:ILE:HG12	2.08	0.53
2:B:101:GLY:HA2	2:B:103:VAL:N	2.24	0.53
1:A:193:PHE:HE2	3:A:366:P0G:HAM	1.73	0.52
2:B:51:ILE:HG23	2:B:71:ARG:HD2	1.92	0.52
1:A:109:TRP:HA	1:A:112:ILE:HD12	1.93	0.51
2:B:59:TYR:HB2	2:B:64:LYS:HG2	1.92	0.51
1:A:193:PHE:CE2	3:A:366:P0G:HAM	2.46	0.51
2:B:38:ARG:HG2	2:B:48:VAL:HG22	1.92	0.51
1:A:167:LEU:O	1:A:171:MET:HB2	2.11	0.50
1:A:329:SER:HB3	1:A:332:PHE:HD1	1.77	0.50
1:A:58:ILE:O	1:A:65:GLN:NE2	2.44	0.50
1:A:71:PHE:HB3	1:A:127:ILE:HG13	1.93	0.50
1:A:269:HIS:HA	1:A:272:LEU:HD23	1.95	0.49
1:A:321:PHE:HA	1:A:324:LEU:HD12	1.95	0.48
1:A:121:ILE:HD11	1:A:286:TRP:CD1	2.49	0.48
1:A:195:THR:HG21	1:A:199:TYR:CD2	2.49	0.48
2:B:38:ARG:HH12	2:B:63:VAL:HG11	1.78	0.48
2:B:40:ALA:HB3	2:B:43:LYS:HB2	1.96	0.48
1:A:174:TYR:HD1	1:A:195:THR:CG2	2.26	0.48
1:A:25:THR:H	1:A:28:ARG:HG2	1.79	0.48
2:B:85:LEU:HB3	2:B:118:VAL:HG21	1.96	0.47
1:A:89:PHE:HE1	1:A:105:TRP:NE1	2.12	0.47
1:A:293:ASN:ND2	3:A:366:P0G:HAMA	2.23	0.47
1:A:174:TYR:HD1	1:A:195:THR:HG23	1.80	0.47
1:A:274:THR:HA	1:A:277:ILE:HD12	1.96	0.47
2:B:101:GLY:HA2	2:B:102:ALA:C	2.40	0.47
2:B:20:LEU:HG	2:B:82:MET:HE1	1.96	0.47
2:B:76:ASN:HD22	2:B:76:ASN:HA	1.63	0.46
1:A:221:ARG:HA	1:A:224:GLN:HG3	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:47:LEU:O	2:B:60:ALA:HB2	2.17	0.45
1:A:89:PHE:CE1	1:A:105:TRP:NE1	2.82	0.45
2:B:59:TYR:CE1	2:B:69:ILE:HG22	2.52	0.45
2:B:97:VAL:HG22	2:B:109:TYR:HB2	1.99	0.44
1:A:268:GLU:O	1:A:272:LEU:HB3	2.18	0.44
1:A:292:VAL:HG21	1:A:311:LEU:CD1	2.48	0.44
1:A:267:LYS:HD3	1:A:267:LYS:H	1.83	0.44
1:A:40:MET:HA	1:A:43:ILE:HD12	1.99	0.44
1:A:218:VAL:O	1:A:222:VAL:HG23	2.18	0.44
1:A:166:PHE:O	1:A:170:GLN:HB2	2.19	0.43
1:A:69:ASN:HA	1:A:72:ILE:HG12	2.00	0.43
1:A:175:ARG:HD2	1:A:185:TYR:CZ	2.54	0.42
2:B:33:THR:HG22	2:B:52:HIS:HA	2.00	0.42
1:A:87:VAL:HG22	1:A:316:TYR:CD1	2.54	0.42
2:B:11:LEU:HA	2:B:117:THR:O	2.19	0.42
2:B:66:ARG:HH22	2:B:89:ASP:CG	2.26	0.42
1:A:330:PRO:HD3	2:B:105:TYR:CE1	2.55	0.41
1:A:50:GLY:O	1:A:54:VAL:HG23	2.19	0.41
1:A:295:VAL:HA	1:A:298:ILE:HD12	2.01	0.41
1:A:83:GLY:HA2	1:A:87:VAL:HB	2.03	0.41
1:A:163:LEU:HD12	1:A:163:LEU:HA	1.92	0.41
1:A:71:PHE:CD1	1:A:154:ILE:HD11	2.48	0.41
3:A:366:P0G:HANA	3:A:366:P0G:HABB	1.86	0.41
1:A:121:ILE:HD12	1:A:282:PHE:HE1	1.86	0.41
1:A:86:VAL:HA	1:A:109:TRP:HE1	1.86	0.40
2:B:19:ARG:HB2	2:B:81:GLN:OE1	2.21	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	280/501 (56%)	252 (90%)	25 (9%)	3 (1%)	11	43
2	B	119/126 (94%)	101 (85%)	14 (12%)	4 (3%)	3	23
All	All	399/627 (64%)	353 (88%)	39 (10%)	7 (2%)	6	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	298	ILE
2	B	61	ASN
1	A	92	ALA
2	B	41	PRO
2	B	54	GLY
2	B	85	LEU
1	A	93	HIS

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	252/429 (59%)	197 (78%)	55 (22%)	1	6
2	B	96/101 (95%)	81 (84%)	15 (16%)	2	15
All	All	348/530 (66%)	278 (80%)	70 (20%)	1	7

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ARG
1	A	29	ASP
1	A	36	MET
1	A	41	SER
1	A	44	VAL
1	A	45	LEU
1	A	52	VAL
1	A	53	LEU
1	A	65	GLN

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Mol	Chain	Res	Type
1	A	73	THR
1	A	86	VAL
1	A	100	THR
1	A	115	LEU
1	A	122	GLU
1	A	125	CYS
1	A	131	ARG
1	A	132	TYR
1	A	133	PHE
1	A	140	LYS
1	A	142	GLN
1	A	144	LEU
1	A	146	THR
1	A	151	ARG
1	A	152	VAL
1	A	159	ILE
1	A	163	LEU
1	A	167	LEU
1	A	170	GLN
1	A	171	MET
1	A	174	TYR
1	A	177	THR
1	A	188	GLU
1	A	195	THR
1	A	203	SER
1	A	218	VAL
1	A	221	ARG
1	A	224	GLN
1	A	266	LEU
1	A	267	LYS
1	A	272	LEU
1	A	275	LEU
1	A	284	LEU
1	A	291	ILE
1	A	292	VAL
1	A	297	VAL
1	A	298	ILE
1	A	299	GLN
1	A	300	ASP
1	A	302	LEU
1	A	303	ILE
1	A	312	ASN

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Mol	Chain	Res	Type
1	A	317	VAL
1	A	329	SER
1	A	333	ARG
1	A	334	ILE
2	B	3	GLN
2	B	11	LEU
2	B	28	ILE
2	B	39	GLN
2	B	46	GLU
2	B	76	ASN
2	B	78	VAL
2	B	88	GLU
2	B	96	ASN
2	B	97	VAL
2	B	98	LYS
2	B	106	GLU
2	B	114	THR
2	B	117	THR
2	B	120	SER

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	178	HIS
1	A	293	ASN
1	A	301	ASN
2	B	39	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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$
3	P0G	A	366	-	27,29,29	1.61	3 (11%)	34,42,42	1.51	7 (20%)

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	P0G	A	366	-	-	5/15/24/24	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	366	P0G	CAW-CAZ	4.94	1.56	1.52
3	A	366	P0G	CAS-NAQ	4.32	1.40	1.35
3	A	366	P0G	CAO-CAV	3.10	1.55	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	366	P0G	CAN-NAP-CBA	-5.12	110.86	116.35
3	A	366	P0G	OAR-CAY-CAW	2.74	120.60	116.76
3	A	366	P0G	OAR-CAM-CAS	2.47	121.10	114.81
3	A	366	P0G	CAO-CAV-CAJ	-2.29	116.04	119.90
3	A	366	P0G	CAY-CAX-NAQ	2.27	121.41	118.86
3	A	366	P0G	CAC-CBA-CAO	2.10	112.69	109.96
3	A	366	P0G	CAI-CAT-CAV	2.05	121.20	118.59

There are no chirality outliers.

All (5) torsion outliers are listed below:

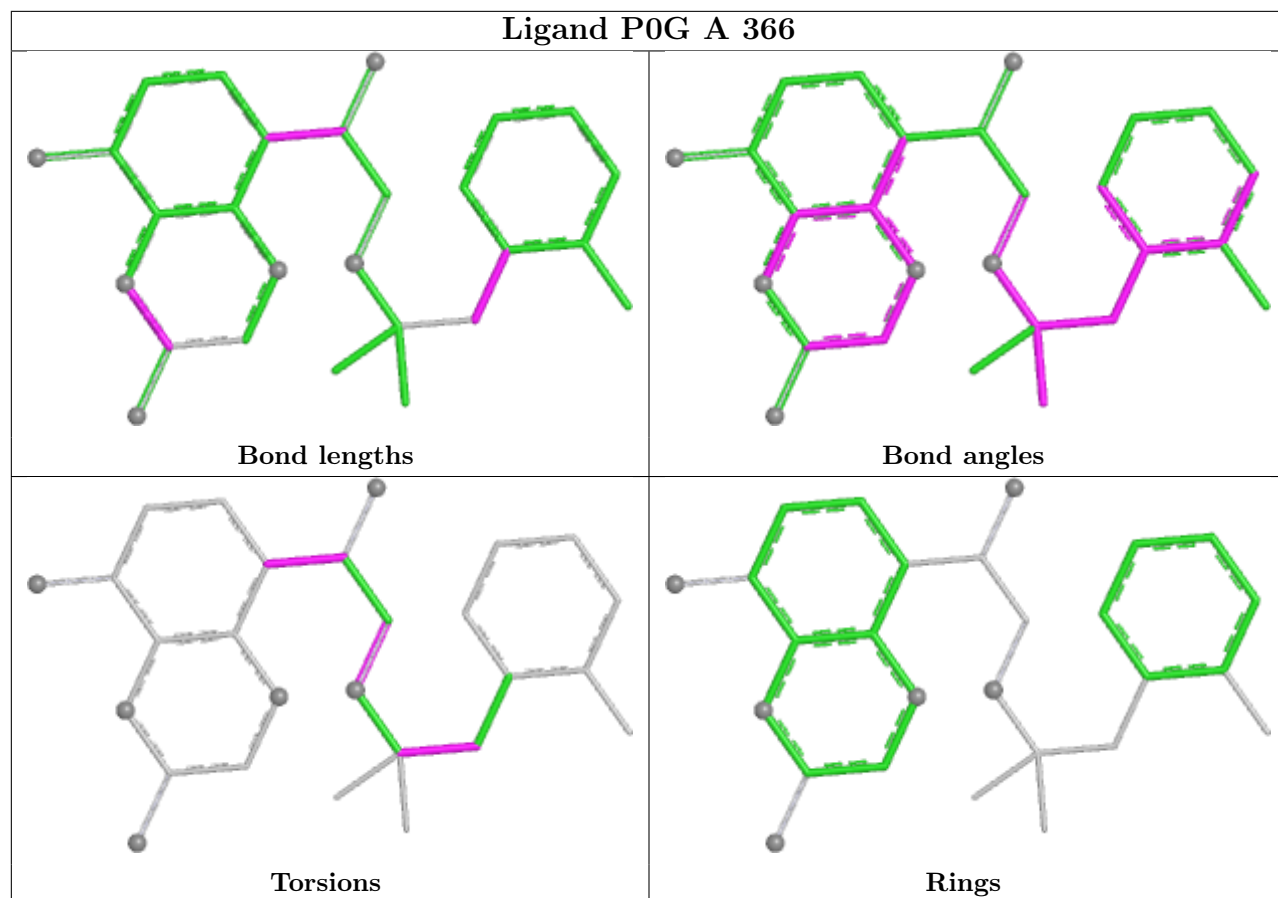
Mol	Chain	Res	Type	Atoms
3	A	366	P0G	CAV-CAO-CBA-CAB
3	A	366	P0G	CAV-CAO-CBA-NAP
3	A	366	P0G	CAV-CAO-CBA-CAC
3	A	366	P0G	CAL-CAW-CAZ-OAF
3	A	366	P0G	CAZ-CAN-NAP-CBA

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	366	P0G	8	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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, 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	284/501 (56%)	0.12	3 (1%)	78 54	44, 67, 114, 185	0
2	B	121/126 (96%)	0.49	1 (0%)	82 60	76, 93, 111, 118	0
All	All	405/627 (64%)	0.23	4 (0%)	79 56	44, 79, 113, 185	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	269	HIS	2.6
1	A	270	LYS	2.2
2	B	2	VAL	2.1
1	A	306	GLU	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](#)

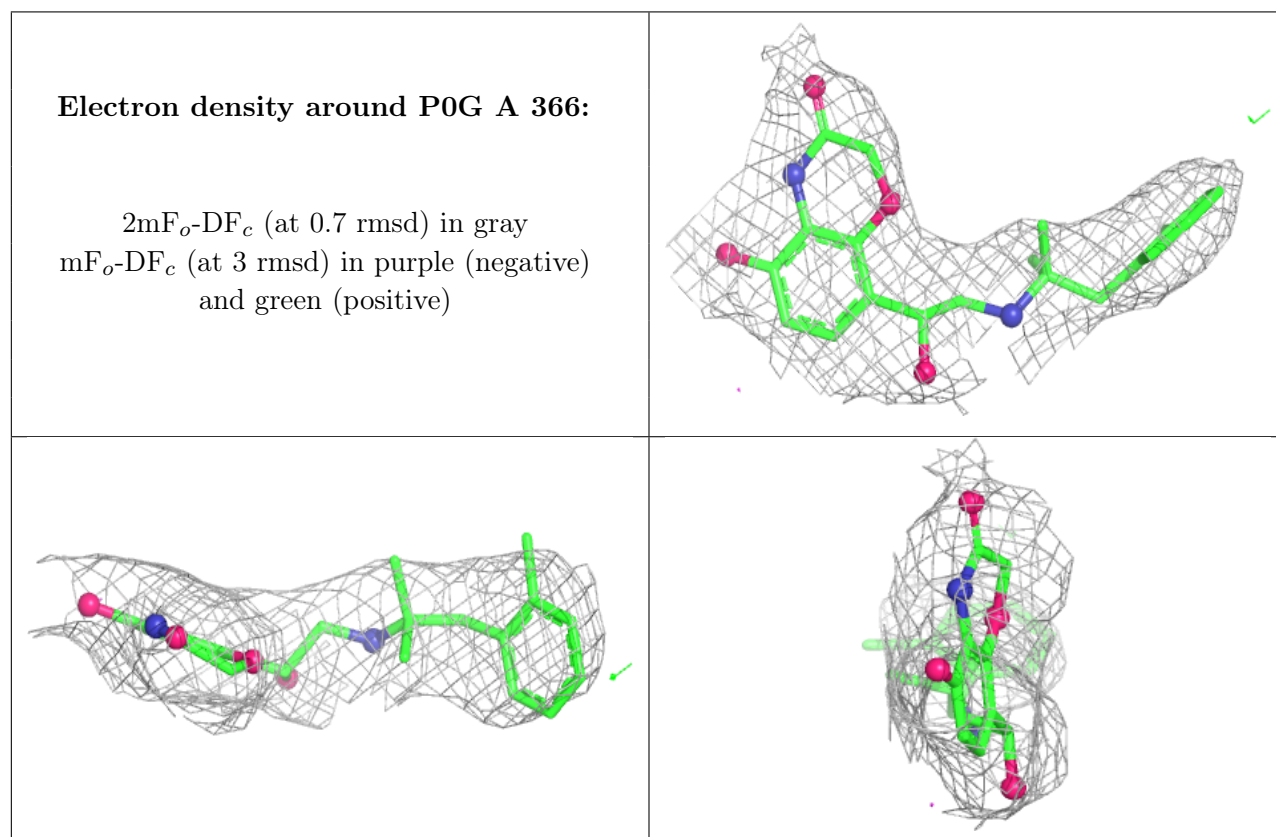
There are no oligosaccharides in this entry.

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(Å ²)	Q < 0.9
3	P0G	A	366	27/27	0.94	0.10	57,57,57,57	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 [i](#)

There are no such residues in this entry.