Full wwPDB X-ray Structure Validation Report (i)

Apr 25, 2022 – 04:19 pm BST

PDB ID	:	7QB1
Title	:	PPARg in complex with inhibitor
Authors	:	Petersen, J.
Deposited on	:	2021-11-17
Resolution	:	2.20 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 (i) symbol.

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

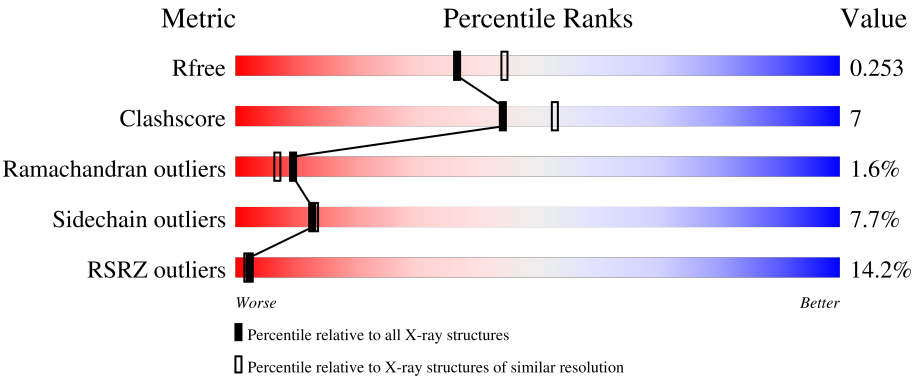
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.28
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.28

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 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 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	AAA	287	10% 72% 17% • 9%
1	BBB	287	16% 71% 16% • 9%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4381 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.

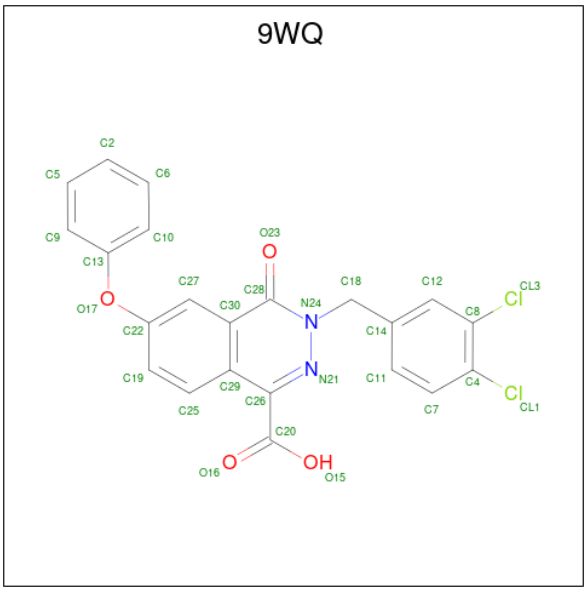
- Molecule 1 is a protein called Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	261	Total	C	N	O	S	0	9	0
			2147	1381	353	403	10			
1	BBB	260	Total	C	N	O	S	0	0	0
			2086	1348	341	388	9			

There are 8 discrepancies between the modelled and reference sequences:

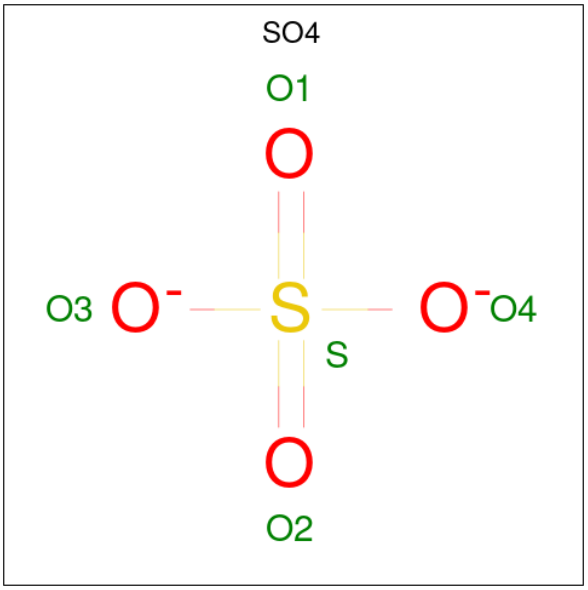
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	191	GLY	-	expression tag	UNP P37231
AAA	192	SER	-	expression tag	UNP P37231
AAA	193	HIS	-	expression tag	UNP P37231
AAA	194	MET	-	expression tag	UNP P37231
BBB	191	GLY	-	expression tag	UNP P37231
BBB	192	SER	-	expression tag	UNP P37231
BBB	193	HIS	-	expression tag	UNP P37231
BBB	194	MET	-	expression tag	UNP P37231

- Molecule 2 is 3-[(3,4-dichlorophenyl)methyl]-4-oxidanylidene-6-phenoxy-phthalazine-1-carboxylic acid (three-letter code: 9WQ) (formula: $C_{22}H_{14}Cl_2N_2O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	AAA	1	Total	C	Cl	N	O	0	0
			30	22	2	2	4		
2	BBB	1	Total	C	Cl	N	O	0	0
			30	22	2	2	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	AAA	1	Total	O	S	0	0
			5	4	1		
3	AAA	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

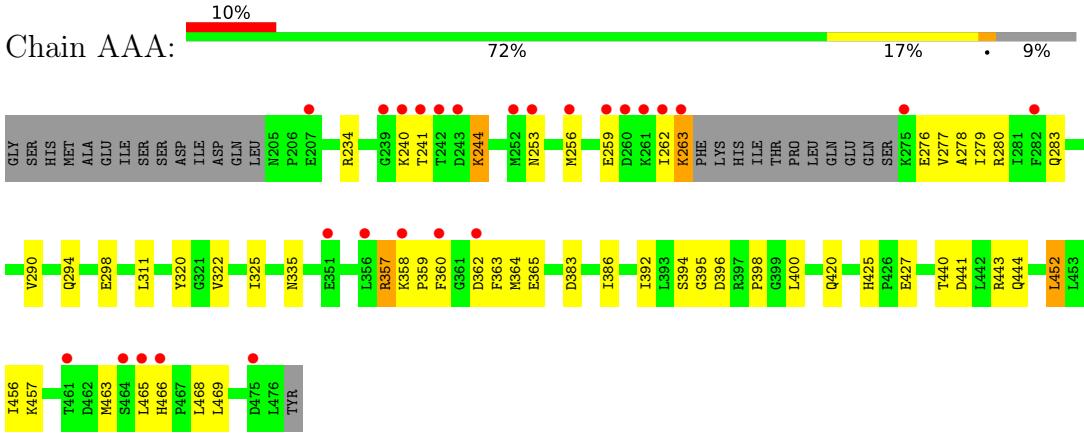
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	AAA	46	Total O 46 46	0	0
4	BBB	32	Total O 32 32	0	0

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.

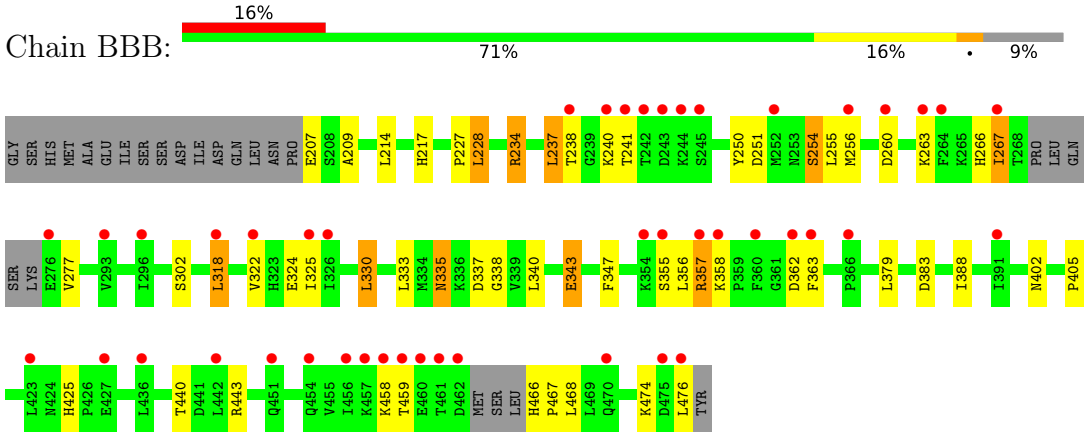
• Molecule 1: Peroxisome proliferator-activated receptor gamma

Chain AAA:



• Molecule 1: Peroxisome proliferator-activated receptor gamma

Chain BBB:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	93.37Å 60.66Å 118.25Å 90.00° 102.93° 90.00°	Depositor
Resolution (Å)	45.98 – 2.20 45.98 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (45.98-2.20) 99.1 (45.98-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.19 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.199 , 0.249 0.204 , 0.253	Depositor DCC
R_{free} test set	1626 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	42.1	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	4381	wwPDB-VP
Average B, all atoms (Å^2)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical 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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 9WQ

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	AAA	0.79	0/2183	0.88	0/2939
1	BBB	0.77	0/2120	0.84	0/2854
All	All	0.78	0/4303	0.86	0/5793

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

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	AAA	2147	0	2214	29	0
1	BBB	2086	0	2147	30	0
2	AAA	30	0	0	0	0
2	BBB	30	0	0	2	0
3	AAA	10	0	0	0	0
4	AAA	46	0	0	1	0
4	BBB	32	0	0	1	0
All	All	4381	0	4361	57	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 7.

All (57) 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:AAA:325:ILE:HD11	1:AAA:392:ILE:HG13	1.60	0.83
1:AAA:240[B]:LYS:HG3	1:AAA:241[B]:THR:HG23	1.65	0.78
1:AAA:259:GLU:OE1	1:AAA:280:ARG:NH2	2.18	0.77
1:AAA:244[A]:LYS:H	1:AAA:244[A]:LYS:HD2	1.55	0.72
1:BBB:383:ASP:OD2	1:BBB:425:HIS:HE1	1.78	0.66
1:AAA:443[B]:ARG:HG3	1:BBB:440:THR:CG2	2.27	0.64
1:AAA:443[B]:ARG:HG3	1:BBB:440:THR:HG23	1.80	0.62
1:BBB:251:ASP:HB3	1:BBB:254:SER:OG	2.02	0.60
1:AAA:383:ASP:OD2	1:AAA:425:HIS:HE1	1.85	0.59
1:BBB:325:ILE:HD12	1:BBB:388:ILE:HG23	1.85	0.58
1:BBB:266:HIS:O	1:BBB:267:ILE:HB	2.03	0.57
1:BBB:250:TYR:CD2	1:BBB:251:ASP:HB2	2.41	0.56
1:AAA:395:GLY:HA2	1:AAA:400:LEU:CD1	2.37	0.55
1:BBB:217:HIS:HE1	1:BBB:302:SER:O	1.89	0.55
1:BBB:255:LEU:HD11	2:BBB:501:9WQ:C2	2.38	0.54
1:BBB:227:PRO:HB2	1:BBB:343:GLU:OE2	2.08	0.54
1:BBB:255:LEU:HD23	1:BBB:277:VAL:CG2	2.39	0.53
1:AAA:320:TYR:CZ	1:AAA:398:PRO:HG2	2.43	0.53
1:AAA:364:MET:HA	1:AAA:364:MET:HE2	1.90	0.53
1:AAA:253:ASN:HA	1:AAA:256:MET:CE	2.39	0.52
1:AAA:286:GLN:HE22	1:AAA:465:LEU:HD12	1.76	0.51
1:BBB:234:ARG:HG3	4:BBB:603:HOH:O	2.11	0.51
1:AAA:452:LEU:O	1:AAA:456:ILE:HD13	2.11	0.50
1:AAA:253:ASN:HA	1:AAA:256:MET:HE2	1.93	0.49
1:BBB:335:ASN:HD22	1:BBB:337:ASP:H	1.61	0.49
1:AAA:290:VAL:HG21	1:AAA:466:HIS:CD2	2.47	0.49
1:AAA:363:PHE:CD1	1:AAA:452:LEU:HD13	2.48	0.49
1:AAA:394:SER:HB3	1:AAA:396:ASP:OD1	2.14	0.48
1:BBB:335:ASN:ND2	1:BBB:337:ASP:H	2.11	0.47
1:BBB:362:ASP:OD1	1:BBB:362:ASP:N	2.47	0.47
1:AAA:279:ILE:HD13	1:AAA:360:PHE:CZ	2.51	0.46
1:AAA:276:GLU:OE2	1:AAA:357:ARG:NE	2.49	0.46
1:BBB:402:ASN:O	1:BBB:405:PRO:HD2	2.16	0.46
1:AAA:262:ILE:O	1:AAA:263:LYS:HG3	2.15	0.46
1:AAA:240[B]:LYS:NZ	1:AAA:240[B]:LYS:HB3	2.31	0.45
1:BBB:255:LEU:HD23	1:BBB:277:VAL:HG23	2.00	0.44
1:BBB:330:LEU:HD23	1:BBB:330:LEU:HA	1.84	0.44
1:AAA:456:ILE:HG21	1:AAA:463:MET:HE1	1.99	0.44
1:AAA:234:ARG:NH2	4:AAA:605:HOH:O	2.50	0.43
1:BBB:324:GLU:OE2	1:BBB:443:ARG:HD3	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:466:HIS:N	1:BBB:467:PRO:CD	2.81	0.43
1:BBB:338:GLY:HA3	1:BBB:347:PHE:CZ	2.54	0.43
1:AAA:358:LYS:HB2	1:AAA:359:PRO:HD3	2.01	0.43
1:BBB:255:LEU:CD1	2:BBB:501:9WQ:C2	2.97	0.43
1:BBB:468:LEU:HD12	1:BBB:468:LEU:HA	1.89	0.42
1:BBB:207:GLU:HG3	1:BBB:209:ALA:HB3	2.02	0.42
1:AAA:277:VAL:HG13	1:AAA:278:ALA:N	2.35	0.41
1:AAA:325:ILE:HD11	1:AAA:392:ILE:CG1	2.40	0.41
1:BBB:318:LEU:HD12	1:BBB:318:LEU:HA	1.87	0.41
1:BBB:228:LEU:HD13	1:BBB:333:LEU:CD2	2.50	0.41
1:BBB:237:LEU:HB3	1:BBB:238:THR:H	1.69	0.41
1:BBB:266:HIS:O	1:BBB:267:ILE:CB	2.67	0.41
1:AAA:290:VAL:O	1:AAA:294:GLN:HG3	2.20	0.41
1:AAA:386:ILE:HD12	1:AAA:420:GLN:HG2	2.02	0.41
1:BBB:318:LEU:O	1:BBB:322:VAL:HG13	2.22	0.40
1:BBB:356:LEU:O	1:BBB:357:ARG:O	2.39	0.40
1:AAA:365:GLU:OE1	1:AAA:365:GLU:HA	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	AAA	266/287 (93%)	258 (97%)	7 (3%)	1 (0%)	34 37
1	BBB	254/287 (88%)	238 (94%)	9 (4%)	7 (3%)	5 2
All	All	520/574 (91%)	496 (95%)	16 (3%)	8 (2%)	9 8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	241	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	BBB	357	ARG
1	AAA	357	ARG
1	BBB	267	ILE
1	BBB	458	LYS
1	BBB	237	LEU
1	BBB	459	THR
1	BBB	474	LYS

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	AAA	240/258 (93%)	222 (92%)	18 (8%)	13	14
1	BBB	233/258 (90%)	214 (92%)	19 (8%)	11	11
All	All	473/516 (92%)	436 (92%)	37 (8%)	13	13

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	244[A]	LYS
1	AAA	244[B]	LYS
1	AAA	263	LYS
1	AAA	283	GLN
1	AAA	286	GLN
1	AAA	298	GLU
1	AAA	311	LEU
1	AAA	322	VAL
1	AAA	335	ASN
1	AAA	362	ASP
1	AAA	427	GLU
1	AAA	440	THR
1	AAA	441	ASP
1	AAA	444	GLN
1	AAA	452	LEU
1	AAA	457	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	AAA	468	LEU
1	AAA	469	LEU
1	BBB	214	LEU
1	BBB	228	LEU
1	BBB	234	ARG
1	BBB	240	LYS
1	BBB	254	SER
1	BBB	256	MET
1	BBB	260	ASP
1	BBB	263	LYS
1	BBB	318	LEU
1	BBB	330	LEU
1	BBB	335	ASN
1	BBB	340	LEU
1	BBB	343	GLU
1	BBB	355	SER
1	BBB	358	LYS
1	BBB	363	PHE
1	BBB	379	LEU
1	BBB	412	ASN
1	BBB	476	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	9WQ	BBB	501	-	28,33,33	1.43	4 (14%)	37,47,47	2.02	11 (29%)
3	SO4	AAA	503	-	4,4,4	0.31	0	6,6,6	0.19	0
2	9WQ	AAA	501	-	28,33,33	1.65	4 (14%)	37,47,47	1.91	10 (27%)
3	SO4	AAA	502	-	4,4,4	0.35	0	6,6,6	0.25	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
2	9WQ	BBB	501	-	-	1/7/12/12	0/4/4/4
2	9WQ	AAA	501	-	-	1/7/12/12	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	AAA	501	9WQ	C26-C29	5.17	1.51	1.43
2	AAA	501	9WQ	C30-C28	3.81	1.48	1.41
2	BBB	501	9WQ	C26-C29	3.77	1.49	1.43
2	AAA	501	9WQ	C8-CL3	2.73	1.80	1.73
2	BBB	501	9WQ	C30-C28	2.68	1.46	1.41
2	BBB	501	9WQ	C19-C22	2.14	1.42	1.38
2	BBB	501	9WQ	C11-C7	2.09	1.42	1.38
2	AAA	501	9WQ	C26-N21	2.02	1.35	1.33

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	AAA	501	9WQ	C14-C18-N24	-6.57	102.86	112.48
2	BBB	501	9WQ	C27-C30-C29	5.55	123.22	118.50
2	BBB	501	9WQ	C8-C4-CL1	-4.11	110.82	120.85
2	AAA	501	9WQ	C12-C8-C4	-3.66	115.23	120.21
2	BBB	501	9WQ	C7-C4-CL1	3.55	125.52	118.41

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	BBB	501	9WQ	C12-C8-CL3	3.52	124.17	118.49
2	AAA	501	9WQ	C8-C12-C14	3.07	122.55	120.46
2	BBB	501	9WQ	C11-C7-C4	-2.95	115.74	120.00
2	AAA	501	9WQ	C27-C30-C29	2.82	120.90	118.50
2	AAA	501	9WQ	C26-N21-N24	2.79	124.76	118.13
2	BBB	501	9WQ	C22-C27-C30	-2.70	117.02	120.48
2	BBB	501	9WQ	C6-C10-C13	2.68	123.25	118.96
2	AAA	501	9WQ	C30-C28-N24	2.59	121.18	116.87
2	BBB	501	9WQ	C18-C14-C12	-2.58	116.25	120.25
2	AAA	501	9WQ	C22-C27-C30	-2.38	117.43	120.48
2	AAA	501	9WQ	C12-C8-CL3	2.19	122.02	118.49
2	AAA	501	9WQ	C18-N24-N21	2.19	121.36	115.46
2	BBB	501	9WQ	C30-C28-N24	2.13	120.41	116.87
2	BBB	501	9WQ	C19-C25-C29	-2.03	118.31	121.13
2	BBB	501	9WQ	C18-N24-N21	2.03	120.94	115.46
2	AAA	501	9WQ	C19-C25-C29	-2.02	118.32	121.13

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	BBB	501	9WQ	C14-C18-N24-C28
2	AAA	501	9WQ	C14-C18-N24-C28

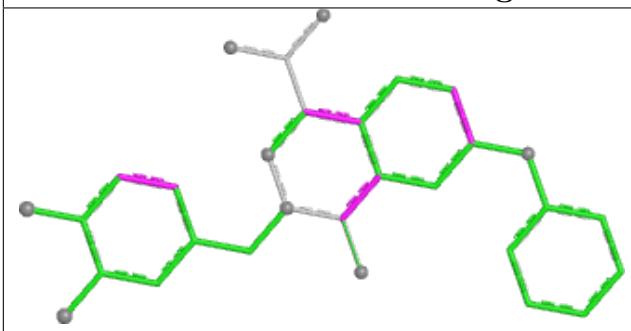
There are no ring outliers.

1 monomer is involved in 2 short contacts:

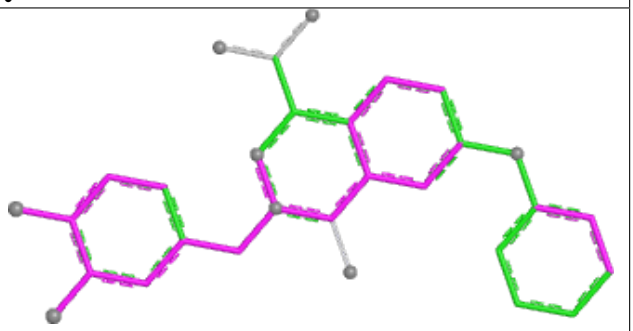
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	BBB	501	9WQ	2	0

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

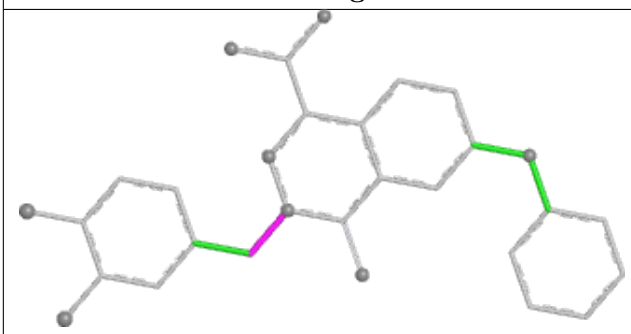
Ligand 9WQ BBB 501



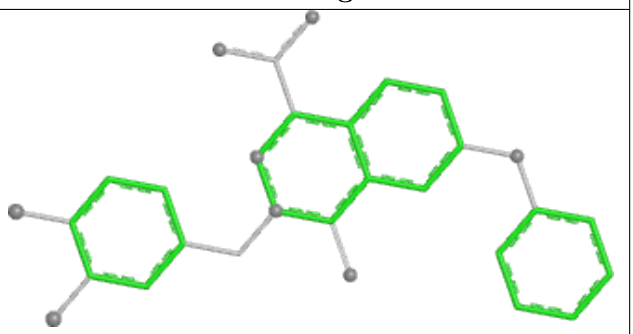
Bond lengths



Bond angles

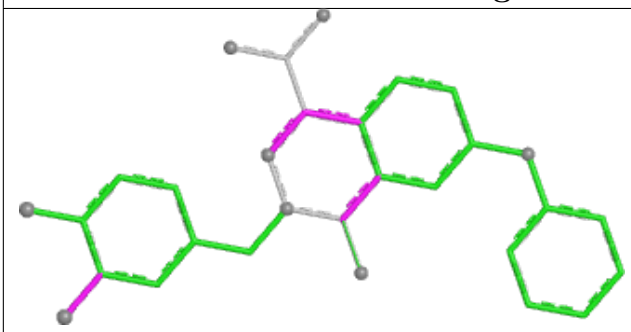


Torsions

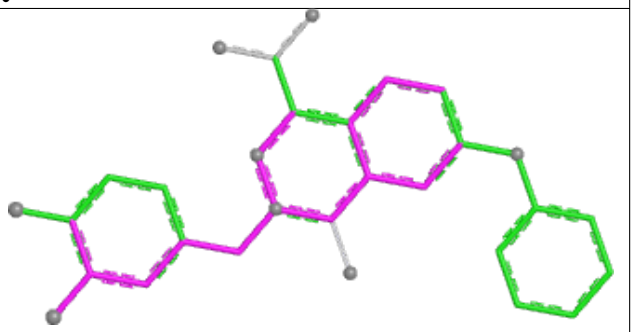


Rings

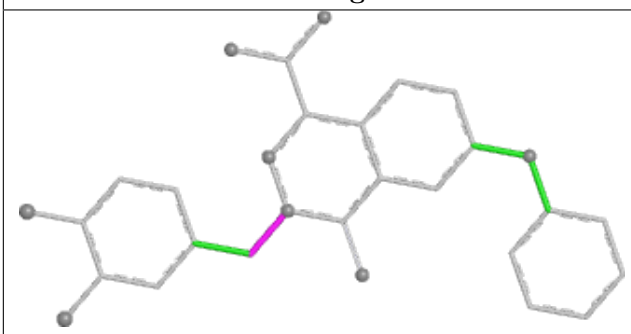
Ligand 9WQ AAA 501



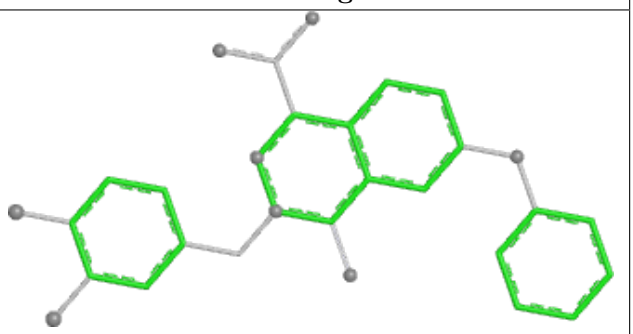
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

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^{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>2			OWAB(Å^2)	Q<0.9
1	AAA	261/287 (90%)	0.85	29 (11%)	5	4	40, 54, 87, 128	0
1	BBB	260/287 (90%)	1.08	45 (17%)	1	1	40, 59, 125, 156	0
All	All	521/574 (90%)	0.96	74 (14%)	2	2	40, 56, 104, 156	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	238	THR	9.2
1	BBB	462	ASP	8.2
1	AAA	242[A]	THR	6.5
1	BBB	459	THR	6.5
1	BBB	470	GLN	6.4
1	AAA	465	LEU	5.9
1	BBB	240	LYS	5.8
1	AAA	241[A]	THR	5.7
1	BBB	241	THR	5.5
1	BBB	476	LEU	5.5
1	BBB	243	ASP	5.5
1	BBB	456	ILE	5.2
1	AAA	256	MET	5.2
1	AAA	464	SER	5.1
1	AAA	475	ASP	5.0
1	AAA	263	LYS	4.8
1	AAA	362	ASP	4.7
1	BBB	242	THR	4.6
1	AAA	260	ASP	4.5
1	BBB	358	LYS	4.4
1	BBB	363	PHE	4.3
1	BBB	461	THR	4.0
1	BBB	454	GLN	4.0
1	BBB	360	PHE	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BBB	458	LYS	3.9
1	AAA	275	LYS	3.8
1	BBB	475	ASP	3.8
1	AAA	282	PHE	3.7
1	BBB	276	GLU	3.6
1	BBB	256	MET	3.6
1	BBB	260	ASP	3.6
1	AAA	262	ILE	3.5
1	AAA	261	LYS	3.5
1	BBB	460	GLU	3.4
1	AAA	252	MET	3.3
1	AAA	240[A]	LYS	3.2
1	AAA	239[A]	GLY	3.2
1	BBB	357	ARG	3.1
1	BBB	322	VAL	3.1
1	BBB	427	GLU	3.0
1	BBB	355	SER	3.0
1	AAA	358	LYS	3.0
1	BBB	325	ILE	3.0
1	AAA	207	GLU	2.9
1	BBB	264	PHE	2.9
1	BBB	318	LEU	2.9
1	AAA	455	VAL	2.9
1	BBB	423	LEU	2.9
1	BBB	326	ILE	2.8
1	BBB	293	VAL	2.8
1	BBB	267	ILE	2.8
1	BBB	244	LYS	2.8
1	BBB	436	LEU	2.7
1	BBB	457	LYS	2.7
1	BBB	451	GLN	2.6
1	AAA	287	PHE	2.6
1	BBB	366	PRO	2.5
1	BBB	362	ASP	2.5
1	BBB	252	MET	2.5
1	BBB	354	LYS	2.3
1	AAA	461	THR	2.3
1	BBB	391	ILE	2.3
1	AAA	259	GLU	2.3
1	BBB	296	ILE	2.3
1	BBB	263	LYS	2.2
1	AAA	356	LEU	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	AAA	351	GLU	2.1
1	BBB	245	SER	2.1
1	AAA	454	GLN	2.1
1	AAA	466	HIS	2.0
1	BBB	442	LEU	2.0
1	AAA	253	ASN	2.0
1	AAA	360	PHE	2.0
1	AAA	243[A]	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

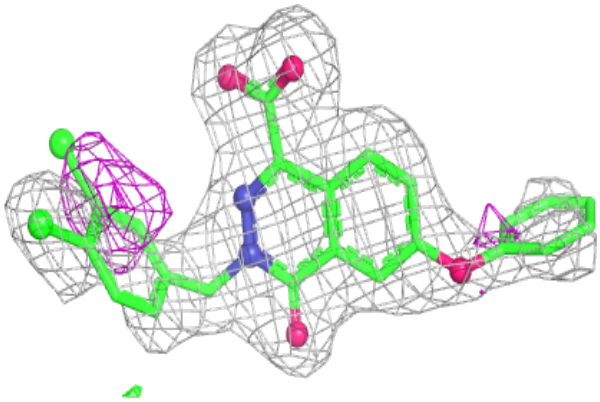
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

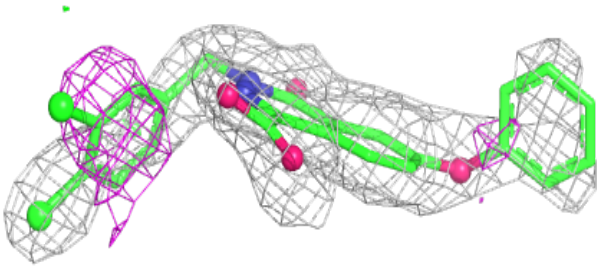
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	SO4	AAA	502	5/5	0.76	0.21	96,98,117,121	0
2	9WQ	AAA	501	30/30	0.79	0.33	52,66,91,106	0
3	SO4	AAA	503	5/5	0.84	0.37	85,96,101,103	0
2	9WQ	BBB	501	30/30	0.87	0.15	49,60,81,92	0

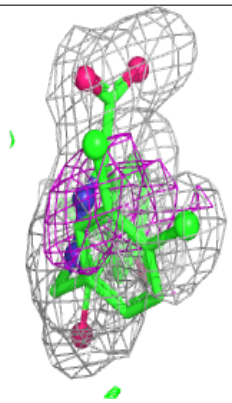
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 9WQ AAA 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

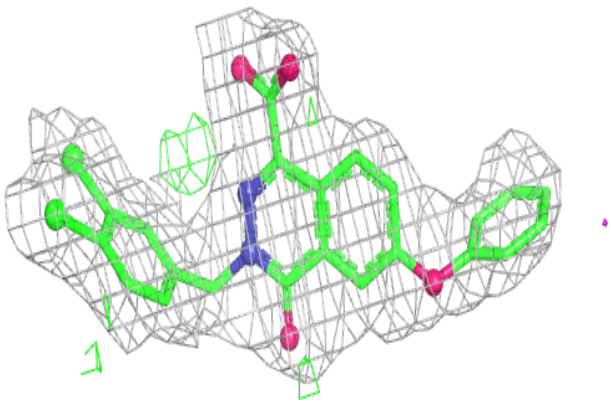


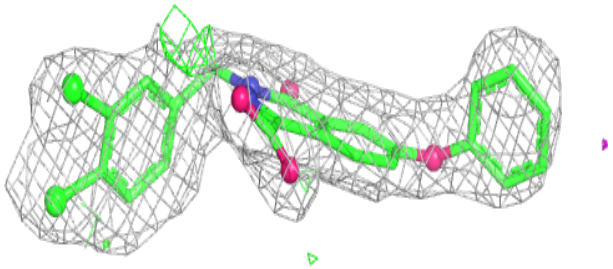


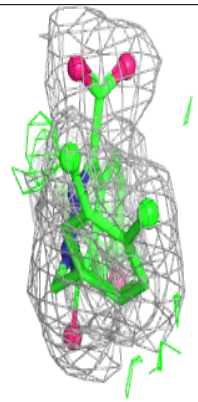


Electron density around 9WQ BBB 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)







6.5 Other polymers (i)

There are no such residues in this entry.