



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:59 AM UTC

PDB ID : 6UJP / pdb_00006ujp
Title : P-glycoprotein mutant-F979A and C952A-with BDE100
Authors : Aller, S.G.; Le, C.A.
Deposited on : 2019-10-03
Resolution : 3.98 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 ⓘ](#)) 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

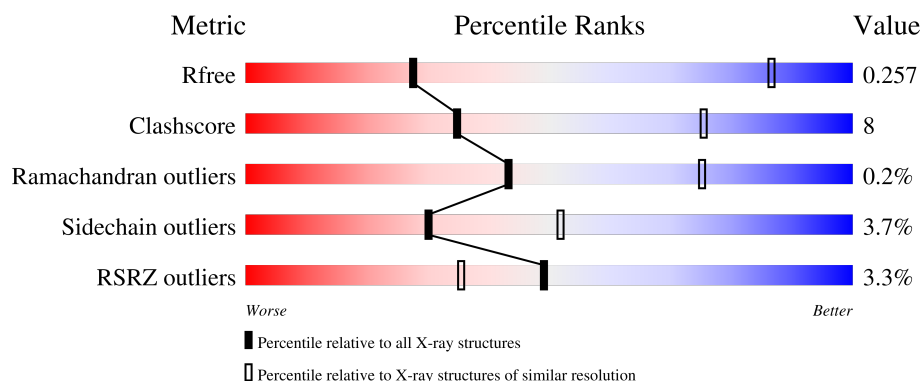
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.98 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	1016 (4.18-3.78)
Clashscore	190562	1007 (4.16-3.80)
Ramachandran outliers	187476	1000 (4.18-3.78)
Sidechain outliers	187428	1131 (4.20-3.76)
RSRZ outliers	180081	1016 (4.18-3.78)

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	1282	3% 73% 19% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9175 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.

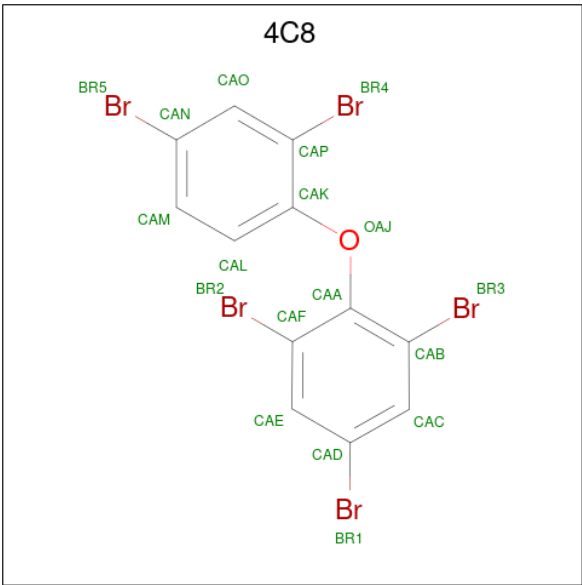
- Molecule 1 is a protein called ATP-dependent translocase ABCB1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1182	Total	C	N	O	S	0	0	0
			9157	5887	1553	1680	37			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	83	GLN	ASN	engineered mutation	UNP P21447
A	87	GLN	ASN	engineered mutation	UNP P21447
A	90	GLN	ASN	engineered mutation	UNP P21447
A	952	ALA	CYS	engineered mutation	UNP P21447
A	979	ALA	PHE	engineered mutation	UNP P21447
A	1277	HIS	-	expression tag	UNP P21447
A	1278	HIS	-	expression tag	UNP P21447
A	1279	HIS	-	expression tag	UNP P21447
A	1280	HIS	-	expression tag	UNP P21447
A	1281	HIS	-	expression tag	UNP P21447
A	1282	HIS	-	expression tag	UNP P21447

- Molecule 2 is 2,4-dibromophenyl 2,4,6-tribromophenyl ether (CCD ID: 4C8) (formula: C₁₂H₅Br₅O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	Br	C	O		
2	A	1	18	5	12	1	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.05Å 138.13Å 188.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.59 – 3.98 29.59 – 3.98	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.59-3.98) 91.8 (29.59-3.98)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.66 (at 3.86Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.239 , 0.263 0.239 , 0.257	Depositor DCC
R_{free} test set	1996 reflections (8.54%)	wwPDB-VP
Wilson B-factor (Å ²)	169.0	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 100.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9175	wwPDB-VP
Average B, all atoms (Å ²)	197.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 4C8

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	0.16	0/9325	0.36	1/12605 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	428	GLY	N-CA-C	-5.31	107.73	114.16

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	9157	0	9346	139	0
2	A	18	0	0	0	0
All	All	9175	0	9346	139	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 8.

All (139) 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:156:ILE:HD11	1:A:904:VAL:HG11	1.69	0.73
1:A:388:LEU:HD11	1:A:547:ILE:HD12	1.70	0.72
1:A:960:VAL:HG13	1:A:965:MET:HG2	1.70	0.72
1:A:160:ASP:OD1	1:A:901:ARG:NH2	2.26	0.67
1:A:1196:ASP:HA	1:A:1226:ILE:HG12	1.76	0.67
1:A:817:ASP:HA	1:A:1000:SER:HB3	1.82	0.62
1:A:1050:GLN:NE2	1:A:1244:ASN:O	2.32	0.62
1:A:846:SER:HB3	1:A:973:VAL:HG13	1.83	0.61
1:A:421:LEU:HB2	1:A:581:ILE:HG13	1.82	0.60
1:A:1063:ALA:HB3	1:A:1239:ILE:HA	1.84	0.60
1:A:256:ALA:HB2	1:A:1117:ILE:HG12	1.83	0.60
1:A:384:ILE:HD13	1:A:465:ILE:HD12	1.84	0.58
1:A:1050:GLN:H	1:A:1245:GLY:HA3	1.68	0.58
1:A:711:ILE:HD12	1:A:833:PHE:HD2	1.68	0.58
1:A:1120:ASP:HA	1:A:1165:VAL:O	2.03	0.58
1:A:446:MET:HE2	1:A:448:SER:HB2	1.87	0.57
1:A:388:LEU:HB2	1:A:413:VAL:HG13	1.86	0.57
1:A:1204:THR:OG1	1:A:1205:GLU:N	2.34	0.57
1:A:436:MET:HE1	1:A:449:ILE:HB	1.87	0.56
1:A:801:ASP:OD2	1:A:1083:TYR:OH	2.23	0.56
1:A:1173:SER:OG	1:A:1176:GLN:OE1	2.24	0.55
1:A:1253:HIS:ND1	1:A:1263:TYR:OH	2.32	0.55
1:A:603:VAL:HG12	1:A:604:GLU:HG3	1.88	0.55
1:A:711:ILE:HD12	1:A:833:PHE:CD2	2.41	0.55
1:A:1157:LEU:HD13	1:A:1163:THR:HG21	1.87	0.55
1:A:78:PHE:HE1	1:A:961:THR:HG21	1.71	0.55
1:A:235:PHE:O	1:A:287:LYS:NZ	2.40	0.55
1:A:721:GLN:OE1	1:A:834:GLN:NE2	2.37	0.54
1:A:221:LEU:HD22	1:A:302:ILE:HG12	1.89	0.54
1:A:1037:VAL:HG22	1:A:1051:GLY:H	1.73	0.54
1:A:1207:GLU:OE1	1:A:1229:ARG:NH2	2.41	0.53
1:A:466:ILE:HG12	1:A:547:ILE:HB	1.89	0.53
1:A:217:ILE:HD12	1:A:305:SER:HB3	1.90	0.53
1:A:944:MET:O	1:A:948:SER:OG	2.18	0.53
1:A:1123:ILE:HD13	1:A:1154:ILE:HD12	1.89	0.53
1:A:211:THR:HG22	1:A:334:VAL:HG21	1.91	0.53
1:A:479:THR:HA	1:A:518:THR:O	2.09	0.53
1:A:805:ASN:HB3	1:A:810:LEU:HD21	1.91	0.52
1:A:1202:LEU:HB3	1:A:1207:GLU:HG2	1.91	0.52
1:A:707:PHE:O	1:A:711:ILE:HG12	2.09	0.52
1:A:418:THR:HG21	1:A:573:ARG:HE	1.75	0.52
1:A:922:ILE:HB	1:A:923:PRO:HD3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1096:GLU:HG2	1:A:1098:LYS:H	1.74	0.51
1:A:1067:SER:HA	1:A:1266:MET:HE1	1.91	0.51
1:A:851:TRP:C	1:A:853:LEU:H	2.18	0.51
1:A:1197:GLU:HA	1:A:1227:ALA:HA	1.93	0.51
1:A:1202:LEU:HD23	1:A:1207:GLU:HA	1.93	0.51
1:A:705:PRO:C	1:A:707:PHE:H	2.19	0.51
1:A:189:PHE:CE1	1:A:348:ILE:HD11	2.47	0.50
1:A:481:ALA:O	1:A:485:ARG:HB2	2.12	0.50
1:A:462:LEU:O	1:A:465:ILE:HG12	2.11	0.50
1:A:719:GLY:HA2	1:A:841:THR:OG1	2.11	0.50
1:A:1050:GLN:O	1:A:1246:LYS:HG3	2.12	0.50
1:A:900:PHE:O	1:A:903:VAL:HG12	2.12	0.50
1:A:550:LEU:HB2	1:A:580:VAL:HG23	1.95	0.49
1:A:160:ASP:HB2	1:A:398:PRO:HG2	1.94	0.49
1:A:1239:ILE:HG21	1:A:1263:TYR:CE1	2.48	0.49
1:A:1154:ILE:HA	1:A:1157:LEU:HD12	1.95	0.49
1:A:1076:VAL:HG13	1:A:1194:LEU:HD13	1.95	0.48
1:A:1124:ALA:HB1	1:A:1141:ILE:HD12	1.95	0.48
1:A:1218:ARG:C	1:A:1220:GLY:H	2.21	0.48
1:A:74:MET:HG2	1:A:110:TYR:CD2	2.48	0.48
1:A:235:PHE:HB3	1:A:287:LYS:HE2	1.95	0.48
1:A:200:PHE:HE1	1:A:215:LEU:HD11	1.79	0.48
1:A:1122:SER:HB3	1:A:1125:GLU:HG2	1.95	0.48
1:A:477:ALA:HA	1:A:520:VAL:O	2.14	0.48
1:A:1079:LEU:HD23	1:A:1194:LEU:HD21	1.95	0.48
1:A:144:ARG:HD3	1:A:175:VAL:HG11	1.96	0.47
1:A:206:ARG:HH11	1:A:326:GLN:HG2	1.78	0.47
1:A:544:ASN:O	1:A:544:ASN:ND2	2.48	0.47
1:A:360:GLU:O	1:A:364:ILE:HG12	2.15	0.47
1:A:1092:LEU:HG	1:A:1093:ASP:H	1.80	0.47
1:A:38:MET:SD	1:A:355:ARG:HG2	2.55	0.47
1:A:938:PHE:O	1:A:942:GLN:HG2	2.15	0.47
1:A:867:ALA:HA	1:A:870:VAL:HG12	1.97	0.46
1:A:423:GLY:N	1:A:429:LYS:HD3	2.30	0.46
1:A:65:PRO:HB2	1:A:202:ILE:HD12	1.98	0.46
1:A:1001:ALA:O	1:A:1005:ILE:HD12	2.15	0.46
1:A:964:LEU:HD12	1:A:964:LEU:H	1.80	0.46
1:A:709:VAL:O	1:A:713:CYS:HB2	2.16	0.46
1:A:1199:THR:HG21	1:A:1211:GLN:HG2	1.98	0.46
1:A:573:ARG:C	1:A:575:GLY:H	2.25	0.45
1:A:388:LEU:HB2	1:A:413:VAL:CG1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1154:ILE:O	1:A:1160:LYS:HA	2.16	0.45
1:A:263:PHE:HA	1:A:1188:ARG:HH22	1.82	0.45
1:A:851:TRP:O	1:A:853:LEU:N	2.47	0.45
1:A:856:LEU:O	1:A:860:ILE:HG12	2.17	0.45
1:A:1021:GLY:HA3	1:A:1101:ASN:HB2	1.99	0.45
1:A:762:SER:HA	1:A:765:THR:HG22	1.98	0.45
1:A:857:LEU:HD12	1:A:973:VAL:HG12	1.99	0.45
1:A:1028:GLU:HB2	1:A:1058:LYS:HD2	1.99	0.45
1:A:833:PHE:HA	1:A:836:ILE:HG12	1.98	0.45
1:A:171:LEU:HD21	1:A:916:TYR:OH	2.17	0.44
1:A:88:SER:OG	1:A:89:THR:N	2.50	0.44
1:A:1186:LEU:HA	1:A:1190:PRO:HD3	2.00	0.44
1:A:313:GLY:O	1:A:317:VAL:HG23	2.17	0.44
1:A:784:LEU:HD12	1:A:1004:ILE:HD13	2.00	0.44
1:A:436:MET:HE3	1:A:466:ILE:HD11	2.00	0.43
1:A:1144:ALA:CB	1:A:1187:VAL:HG23	2.48	0.43
1:A:724:PHE:CE1	1:A:758:LEU:HB3	2.54	0.43
1:A:893:ALA:HB2	1:A:916:TYR:CE1	2.53	0.43
1:A:696:ILE:HB	1:A:1005:ILE:HD11	2.01	0.43
1:A:966:THR:HB	1:A:968:GLU:HG2	2.00	0.43
1:A:1150:ILE:HD13	1:A:1179:ARG:HH11	1.84	0.43
1:A:824:ALA:HA	1:A:828:ARG:HD3	2.00	0.43
1:A:1150:ILE:CD1	1:A:1179:ARG:HH11	2.32	0.43
1:A:178:ILE:HA	1:A:354:ALA:HB1	2.01	0.43
1:A:1070:CYS:O	1:A:1074:THR:N	2.36	0.43
1:A:1150:ILE:HB	1:A:1179:ARG:HB3	2.01	0.42
1:A:1234:GLN:HG2	1:A:1253:HIS:CE1	2.54	0.42
1:A:83:GLN:O	1:A:87:GLN:HB2	2.19	0.42
1:A:585:LEU:HD12	1:A:588:VAL:HG11	2.01	0.42
1:A:200:PHE:CE1	1:A:215:LEU:HD11	2.53	0.42
1:A:200:PHE:CE1	1:A:334:VAL:HG13	2.54	0.42
1:A:1063:ALA:HB3	1:A:1239:ILE:HG13	2.01	0.42
1:A:1113:SER:O	1:A:1178:GLN:NE2	2.51	0.42
1:A:1068:SER:OG	1:A:1069:GLY:N	2.52	0.42
1:A:113:TYR:O	1:A:117:ILE:HG12	2.19	0.42
1:A:1176:GLN:OE1	1:A:1176:GLN:N	2.47	0.42
1:A:151:ILE:HD12	1:A:167:LEU:HD21	2.01	0.42
1:A:608:HIS:O	1:A:612:MET:HG2	2.19	0.42
1:A:228:TRP:HZ2	1:A:295:MET:HB2	1.85	0.42
1:A:257:ILE:HG12	1:A:800:PHE:CE2	2.55	0.42
1:A:698:LYS:O	1:A:701:SER:OG	2.30	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:GLU:HB2	1:A:287:LYS:HZ1	1.84	0.41
1:A:41:TYR:HE2	1:A:362:PHE:HE1	1.69	0.41
1:A:33:VAL:O	1:A:355:ARG:NH1	2.53	0.41
1:A:40:ARG:HD2	1:A:41:TYR:CZ	2.56	0.41
1:A:147:PHE:O	1:A:151:ILE:HG12	2.20	0.41
1:A:474:VAL:HG13	1:A:902:THR:HG21	2.02	0.41
1:A:434:GLN:HG3	1:A:439:LEU:HD12	2.02	0.41
1:A:1064:LEU:HD12	1:A:1064:LEU:HA	1.93	0.41
1:A:238:LYS:O	1:A:242:ALA:N	2.53	0.41
1:A:806:THR:HG23	1:A:809:ALA:H	1.86	0.41
1:A:1181:ALA:HA	1:A:1184:ARG:HG2	2.03	0.41
1:A:864:ILE:HG22	1:A:944:MET:HB3	2.02	0.40
1:A:317:VAL:HG13	1:A:322:TYR:O	2.20	0.40
1:A:471:GLN:HB2	1:A:552:GLU:HB2	2.04	0.40
1:A:170:ARG:O	1:A:174:ASP:HB2	2.22	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	1178/1282 (92%)	1128 (96%)	48 (4%)	2 (0%)	43 75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1204	THR
1	A	852	GLN

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	971/1061 (92%)	935 (96%)	36 (4%)	30	52

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	LEU
1	A	144	ARG
1	A	260	VAL
1	A	261	ILE
1	A	270	LEU
1	A	372	ASP
1	A	406	LEU
1	A	409	LEU
1	A	413	VAL
1	A	434	GLN
1	A	447	VAL
1	A	465	ILE
1	A	474	VAL
1	A	496	ILE
1	A	518	THR
1	A	593	VAL
1	A	761	ILE
1	A	795	GLN
1	A	897	ILE
1	A	911	LYS
1	A	930	LYS
1	A	966	THR
1	A	1005	ILE
1	A	1030	ASN
1	A	1032	GLN
1	A	1054	LEU
1	A	1090	VAL
1	A	1093	ASP
1	A	1102	VAL
1	A	1155	ASP

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Mol	Chain	Res	Type
1	A	1172	LEU
1	A	1182	ILE
1	A	1187	VAL
1	A	1225	VAL
1	A	1233	ILE
1	A	1247	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	153	ASN
1	A	154	GLN
1	A	452	GLN
1	A	504	ASN
1	A	544	ASN
1	A	769	GLN
1	A	805	ASN
1	A	942	GLN
1	A	1211	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 [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
2	4C8	A	1301	-	19,19,19	1.08	2 (10%)	27,27,27	0.95	1 (3%)

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	4C8	A	1301	-	-	0/4/4/4	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1301	4C8	BR1-CAD	2.18	1.94	1.90
2	A	1301	4C8	BR3-CAB	2.09	1.94	1.89

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1301	4C8	CAK-OAJ-CAA	-2.43	112.72	117.94

There are no chirality outliers.

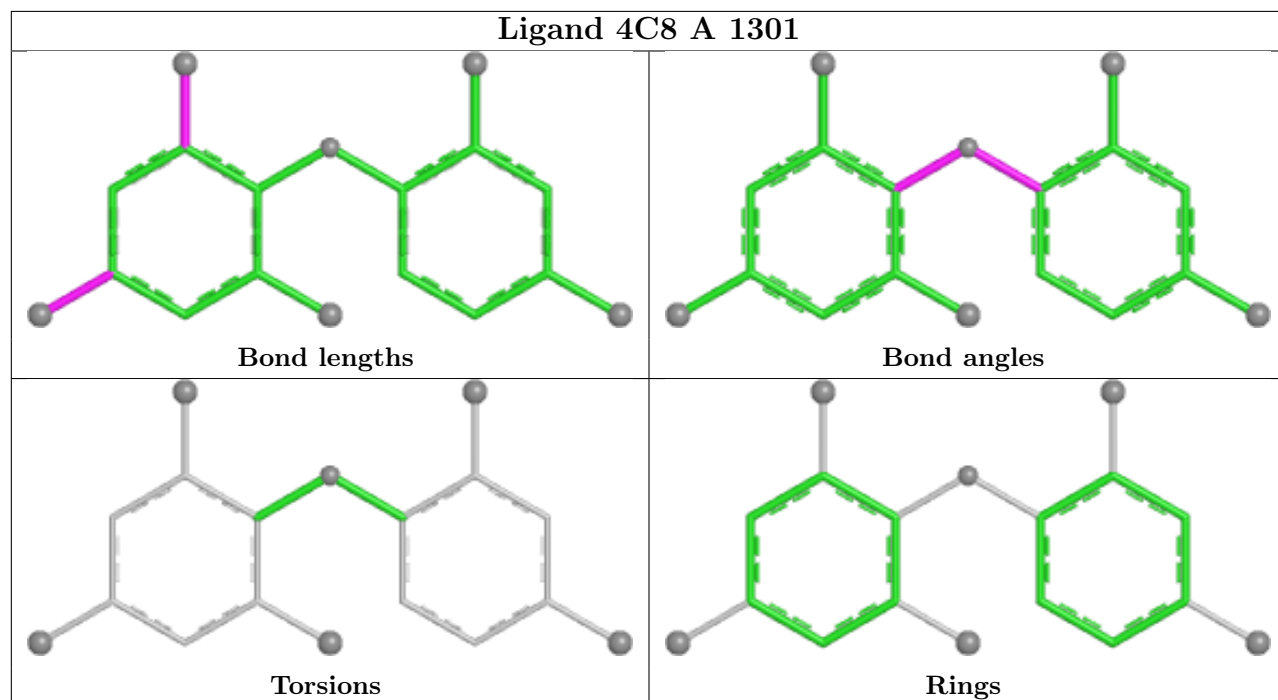
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	1182/1282 (92%)	0.02	39 (3%)	49 35	133, 192, 262, 319	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	321	GLU	4.9
1	A	605	GLN	4.7
1	A	391	LYS	4.5
1	A	410	ASN	4.0
1	A	477	ALA	3.9
1	A	49	TYR	3.4
1	A	974	PHE	3.4
1	A	732	VAL	3.3
1	A	460	ARG	3.1
1	A	978	VAL	3.0
1	A	364	ILE	3.0
1	A	1059	GLY	2.8
1	A	324	ILE	2.7
1	A	519	LEU	2.7
1	A	1223	CYS	2.6
1	A	734	VAL	2.6
1	A	705	PRO	2.5
1	A	502	GLU	2.5
1	A	1225	VAL	2.5
1	A	322	TYR	2.4
1	A	389	GLU	2.4
1	A	422	VAL	2.4
1	A	45	LEU	2.4
1	A	343	GLN	2.3
1	A	62	VAL	2.3
1	A	323	SER	2.3
1	A	964	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	348	ILE	2.2
1	A	967	PHE	2.2
1	A	707	PHE	2.2
1	A	412	LYS	2.2
1	A	327	VAL	2.2
1	A	971	LEU	2.1
1	A	700	ASN	2.1
1	A	71	PHE	2.1
1	A	154	GLN	2.1
1	A	184	ASP	2.1
1	A	153	ASN	2.1
1	A	776	ALA	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](#)

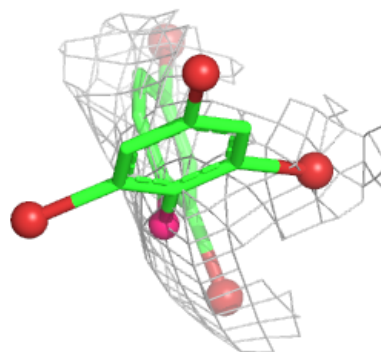
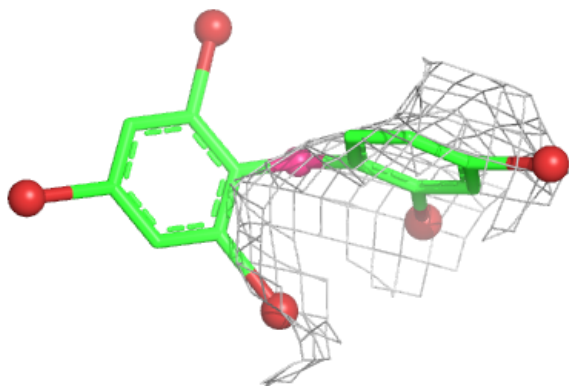
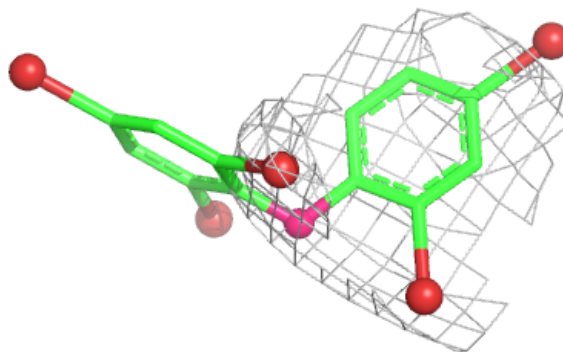
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	4C8	A	1301	18/18	0.72	0.14	203,216,236,239	18

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 4C8 A 1301:

$2mF_o-DF_c$ (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.