



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 05:33 PM BST

PDB ID : 3ZfZ
Title : Crystal structure of ceftaroline acyl-PBP2a from MRSA with non-covalently bound ceftaroline and muramic acid at allosteric site obtained by soaking
Authors : Otero, L.H.; Rojas-Altuve, A.; Hermoso, J.A.
Deposited on : 2012-12-13
Resolution : 2.25 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

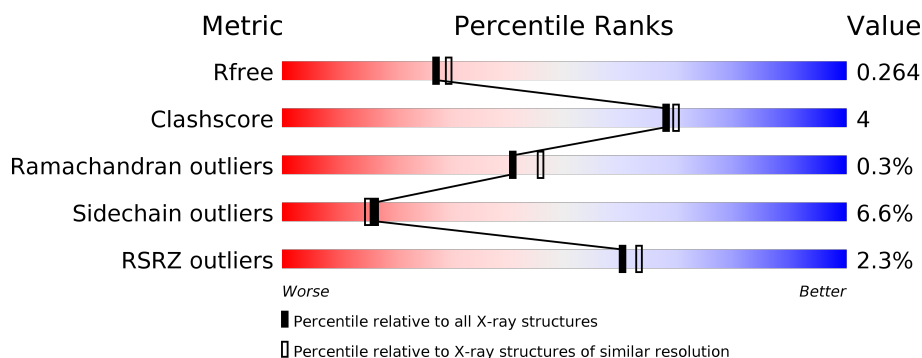
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 on 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	642	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>••</div> </div> </div>
1	B	642	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>•</div> </div> </div>

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
5	1W8	A	1669	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 11069 atoms, of which 32 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 PENICILLIN BINDING PROTEIN 2 PRIME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	635	Total	C	N	O	S	0	0	0
			5096	3213	860	1008	15			
1	B	642	Total	C	N	O	S	0	0	0
			5151	3246	870	1019	16			

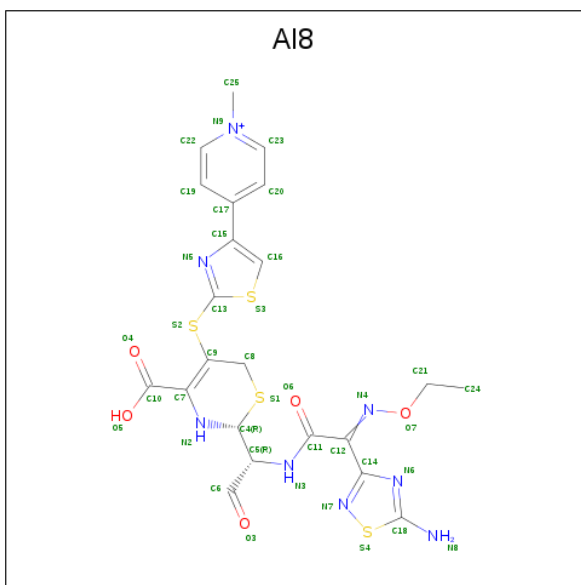
- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Cd	0	0
			3	3		
2	A	5	Total	Cd	0	0
			5	5		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

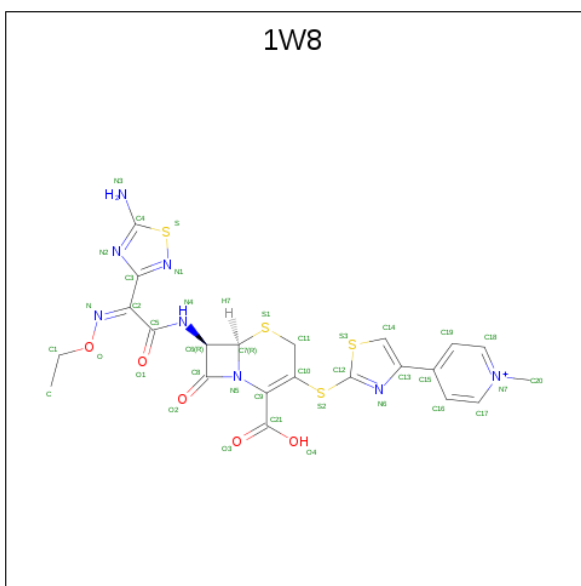
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		
3	A	2	Total	Cl	0	0
			2	2		

- Molecule 4 is Ceftaroline, bound form (three-letter code: AI8) (formula: C₂₂H₂₃N₈O₅S₄).



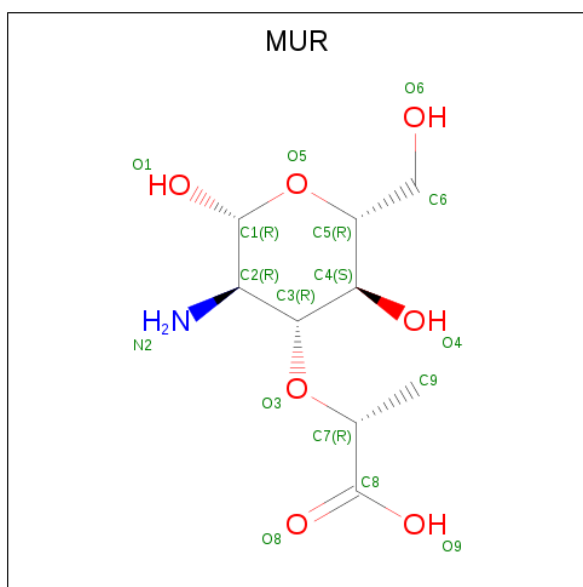
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			39	22	8	5	4		

- Molecule 5 is Ceftaroline (three-letter code: 1W8) (formula: $C_{22}H_{21}N_8O_5S_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			39	22	8	5	4		

- Molecule 6 is beta-muramic acid (three-letter code: MUR) (formula: $C_9H_{17}NO_7$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	H	N	O	0	0
			33	9	16	1	7		
6	B	1	Total	C	H	N	O	0	0
			33	9	16	1	7		

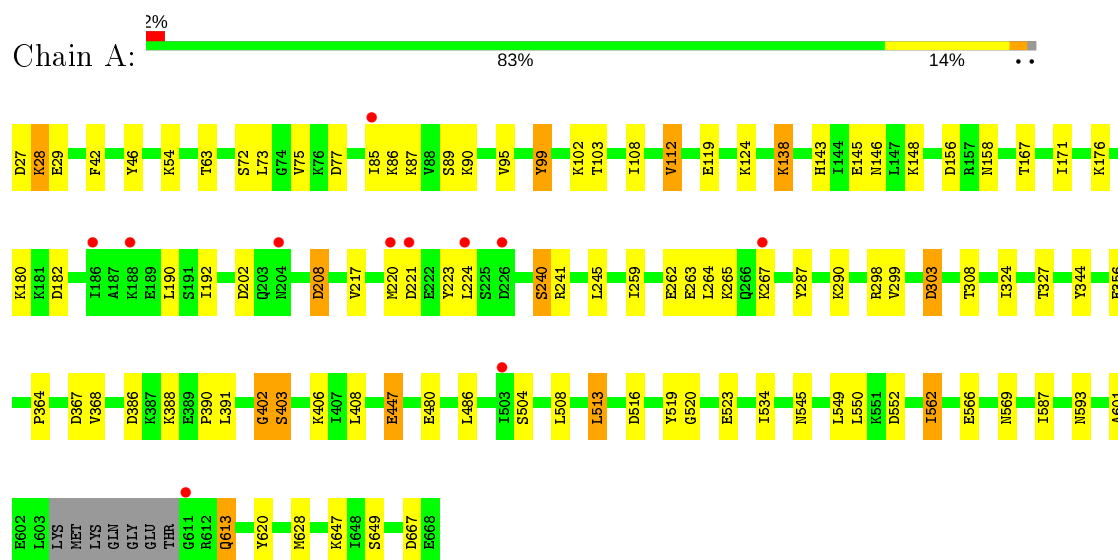
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	334	Total	O	0	0
			334	334		
7	B	332	Total	O	0	0
			332	332		

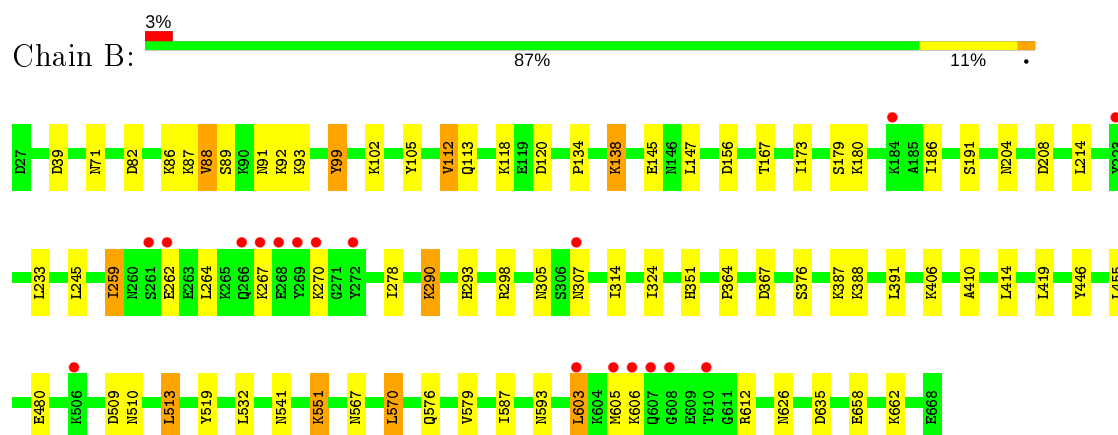
3 Residue-property plots [i](#)

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: PENICILLIN BINDING PROTEIN 2 PRIME



• Molecule 1: PENICILLIN BINDING PROTEIN 2 PRIME



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.52Å 101.48Å 187.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.76 – 2.25 63.08 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (68.76-2.25) 100.0 (63.08-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 2.25Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.194 , 0.249 0.209 , 0.264	Depositor DCC
R_{free} test set	5315 reflections (7.24%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.775	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11069	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 1W8, CD, MUR, AI8, CL

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	0.51	1/5181 (0.0%)	0.71	2/6967 (0.0%)
1	B	0.50	0/5237	0.70	0/7041
All	All	0.50	1/10418 (0.0%)	0.70	2/14008 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	403	SER	C-N	-8.41	1.14	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	402	GLY	C-N-CA	-6.46	105.54	121.70
1	A	403	SER	CB-CA-C	6.24	121.95	110.10

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	5096	0	5088	45	0
1	B	5151	0	5150	36	0
2	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	3	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	39	0	21	1	0
5	A	39	0	20	7	0
6	A	17	16	16	3	0
6	B	17	16	16	0	0
7	A	334	0	0	1	0
7	B	332	0	0	1	0
All	All	11037	32	10311	79	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 4.

All (79) 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)
5:A:1669:1W8:H18	1:B:145:GLU:HG3	1.62	0.82
5:A:1669:1W8:H201	1:B:145:GLU:HB3	1.72	0.72
1:A:138:LYS:H	1:A:138:LYS:HD2	1.56	0.71
1:A:601:ALA:O	1:A:613:GLN:HA	1.92	0.69
1:A:259:ILE:HG13	1:A:263:GLU:HG3	1.76	0.67
1:B:91:ASN:HB3	1:B:118:LYS:HB3	1.75	0.66
5:A:1669:1W8:C17	1:B:298:ARG:HH22	2.09	0.65
1:B:138:LYS:H	1:B:138:LYS:CD	2.09	0.64
1:A:402:GLY:O	1:A:520:GLY:HA3	1.96	0.64
1:B:138:LYS:H	1:B:138:LYS:HD2	1.62	0.63
1:A:447:GLU:HG3	4:A:1403:AI8:H20	1.80	0.62
1:B:88:VAL:HG13	1:B:89:SER:H	1.66	0.60
1:A:408:LEU:HD22	1:A:534:ILE:HG21	1.84	0.60
1:A:138:LYS:H	1:A:138:LYS:CD	2.13	0.60
1:A:99:TYR:HB2	1:A:112:VAL:HG22	1.85	0.58
1:A:220:MET:HG3	1:A:221:ASP:H	1.68	0.58
1:B:414:LEU:HD13	1:B:567:ASN:HB3	1.86	0.58
1:A:259:ILE:HG12	1:A:264:LEU:HG	1.85	0.58
1:B:99:TYR:HB2	1:B:112:VAL:HG22	1.87	0.57
1:B:173:ILE:HD12	1:B:214:LEU:HD11	1.86	0.56
1:A:158:ASN:HA	1:B:551:LYS:HE2	1.86	0.56
1:B:186:ILE:HD13	1:B:233:LEU:HD21	1.89	0.55
1:A:290:LYS:HB3	1:A:324:ILE:HD11	1.87	0.55
1:B:112:VAL:HG13	1:B:134:PRO:HB3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:364:PRO:HG3	1:B:388:LYS:HB3	1.91	0.53
1:A:587:ILE:HD12	1:A:649:SER:HB2	1.90	0.53
1:B:455:LEU:HD23	1:B:570:LEU:HD22	1.91	0.52
1:A:143:HIS:HB3	1:A:145:GLU:OE2	2.10	0.52
1:B:259:ILE:HG12	1:B:264:LEU:HG	1.92	0.51
5:A:1669:1W8:H202	1:B:147:LEU:HD23	1.91	0.51
1:A:27:ASP:N	1:A:28:LYS:HZ2	2.08	0.51
1:B:99:TYR:HB2	1:B:112:VAL:CG2	2.40	0.51
1:A:73:LEU:HD21	1:A:299:VAL:HG11	1.93	0.50
1:B:191:SER:HB3	1:B:376:SER:HB3	1.93	0.50
1:B:87:LYS:HA	1:B:93:LYS:HG2	1.93	0.49
1:A:146:ASN:HB2	5:A:1669:1W8:H111	1.93	0.49
1:A:217:VAL:HG21	1:A:224:LEU:HD13	1.94	0.48
1:A:72:SER:HB3	1:B:305:ASN:HB3	1.95	0.48
1:A:486:LEU:HD23	1:A:562:ILE:HG13	1.95	0.48
1:B:89:SER:HB2	1:B:92:LYS:HB3	1.96	0.48
1:A:240:SER:HA	6:A:1670:MUR:H2	1.96	0.47
1:A:85:ILE:HG23	1:A:95:VAL:HG22	1.97	0.47
1:A:182:ASP:HA	7:A:2092:HOH:O	2.15	0.47
1:B:510:ASN:HB3	1:B:513:LEU:HB2	1.97	0.47
1:A:190:LEU:HB2	1:A:192:ILE:HG12	1.96	0.47
1:A:176:LYS:HG3	1:A:208:ASP:HB3	1.97	0.46
1:A:303:ASP:HB2	1:A:308:THR:HG23	1.98	0.46
1:A:364:PRO:HG2	1:A:388:LYS:HB3	1.96	0.46
1:B:267:LYS:HA	1:B:270:LYS:HD2	1.97	0.46
1:A:513:LEU:HD21	1:A:523:GLU:HG3	1.98	0.46
1:A:158:ASN:HD21	1:A:667:ASP:HA	1.80	0.46
1:A:119:GLU:HB2	1:A:124:LYS:HE3	1.97	0.46
1:A:327:THR:OG1	1:A:549:LEU:HA	2.16	0.45
1:B:406:LYS:HE3	1:B:519:TYR:HB2	1.99	0.45
1:B:576:GLN:HE22	1:B:593:ASN:HD21	1.65	0.44
1:A:99:TYR:HB2	1:A:112:VAL:CG2	2.48	0.44
1:B:410:ALA:O	1:B:414:LEU:HG	2.18	0.44
1:A:241:ARG:H	6:A:1670:MUR:H92	1.82	0.43
1:A:480:GLU:HG3	1:A:508:LEU:HD12	2.00	0.43
1:B:105:TYR:O	1:B:314:ILE:HG21	2.19	0.43
1:B:579:VAL:HG13	1:B:587:ILE:HG23	2.00	0.43
1:B:658:GLU:HB2	1:B:662:LYS:HG2	2.00	0.43
1:A:403:SER:CB	1:A:406:LYS:HZ2	2.31	0.43
1:A:46:TYR:O	1:A:54:LYS:HD3	2.18	0.43
1:B:290:LYS:HB3	1:B:324:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:TYR:CZ	1:A:550:LEU:HD11	2.54	0.43
1:B:603:LEU:HD22	1:B:612:ARG:HD3	2.00	0.43
5:A:1669:1W8:H18	1:B:145:GLU:CG	2.42	0.42
1:A:327:THR:HB	1:A:356:GLU:HB3	2.00	0.42
1:A:386:ASP:HB3	1:A:390:PRO:HD2	2.00	0.42
5:A:1669:1W8:S2	1:B:307:ASN:O	2.77	0.42
1:A:241:ARG:HB2	6:A:1670:MUR:H4	2.01	0.42
1:B:293:HIS:CE1	7:B:2079:HOH:O	2.72	0.42
1:A:516:ASP:HA	1:A:519:TYR:CE1	2.55	0.42
1:A:620:TYR:HB3	1:A:628:MET:HG2	2.01	0.42
1:A:42:PHE:HB3	1:A:63:THR:HA	2.01	0.42
1:B:351:HIS:HD2	1:B:626:ASN:O	2.02	0.41
1:A:75:VAL:HG22	1:A:103:THR:HG22	2.03	0.41
1:A:504:SER:HA	1:A:523:GLU:HB2	2.03	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	631/642 (98%)	611 (97%)	19 (3%)	1 (0%)	47	55
1	B	640/642 (100%)	611 (96%)	26 (4%)	3 (0%)	29	29
All	All	1271/1284 (99%)	1222 (96%)	45 (4%)	4 (0%)	41	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	613	GLN
1	B	635	ASP
1	B	606	LYS
1	B	88	VAL

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	566/572 (99%)	526 (93%)	40 (7%)	14	13
1	B	572/572 (100%)	537 (94%)	35 (6%)	18	18
All	All	1138/1144 (100%)	1063 (93%)	75 (7%)	16	15

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	29	GLU
1	A	77	ASP
1	A	86	LYS
1	A	87	LYS
1	A	89	SER
1	A	90	LYS
1	A	99	TYR
1	A	102	LYS
1	A	108	ILE
1	A	112	VAL
1	A	138	LYS
1	A	148	LYS
1	A	156	ASP
1	A	167	THR
1	A	171	ILE
1	A	180	LYS
1	A	202	ASP
1	A	208	ASP
1	A	223	TYR
1	A	240	SER
1	A	245	LEU
1	A	262	GLU
1	A	265	LYS
1	A	267	LYS
1	A	298	ARG
1	A	303	ASP

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Mol	Chain	Res	Type
1	A	344	TYR
1	A	367	ASP
1	A	368	VAL
1	A	391	LEU
1	A	447	GLU
1	A	513	LEU
1	A	545	ASN
1	A	552	ASP
1	A	562	ILE
1	A	566	GLU
1	A	569	ASN
1	A	593	ASN
1	A	647	LYS
1	B	39	ASP
1	B	71	ASN
1	B	82	ASP
1	B	86	LYS
1	B	99	TYR
1	B	102	LYS
1	B	112	VAL
1	B	113	GLN
1	B	120	ASP
1	B	138	LYS
1	B	156	ASP
1	B	167	THR
1	B	179	SER
1	B	180	LYS
1	B	204	ASN
1	B	208	ASP
1	B	245	LEU
1	B	259	ILE
1	B	262	GLU
1	B	278	ILE
1	B	290	LYS
1	B	367	ASP
1	B	387	LYS
1	B	391	LEU
1	B	419	LEU
1	B	446	TYR
1	B	480	GLU
1	B	509	ASP
1	B	513	LEU

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Mol	Chain	Res	Type
1	B	532	LEU
1	B	541	ASN
1	B	551	LYS
1	B	570	LEU
1	B	603	LEU
1	B	605	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	A	203	GLN
1	A	381	ASN
1	A	593	ASN
1	B	266	GLN
1	B	293	HIS
1	B	593	ASN

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 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	AI8	A	1403	1	27,42,42	2.58	10 (37%)	26,58,58	3.07	11 (42%)
6	MUR	B	1669	-	14,17,17	1.48	2 (14%)	19,24,24	2.09	8 (42%)
6	MUR	A	1670	-	14,17,17	1.49	3 (21%)	19,24,24	2.32	8 (42%)
5	1W8	A	1669	-	31,43,43	0.56	1 (3%)	35,62,62	1.79	4 (11%)

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
4	AI8	A	1403	1	-	3/15/47/47	0/3/4/4
6	MUR	B	1669	-	-	1/6/30/30	0/1/1/1
6	MUR	A	1670	-	-	2/6/30/30	0/1/1/1
5	1W8	A	1669	-	-	4/15/57/57	0/5/5/5

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1403	AI8	O7-N4	-6.29	1.30	1.42
4	A	1403	AI8	C11-N3	5.89	1.45	1.34
4	A	1403	AI8	C7-N2	4.07	1.47	1.33
4	A	1403	AI8	C17-C15	3.92	1.55	1.48
4	A	1403	AI8	O7-C21	-3.64	1.37	1.44
4	A	1403	AI8	O6-C11	-3.39	1.17	1.23
4	A	1403	AI8	C18-N8	3.34	1.45	1.35
6	B	1669	MUR	O5-C5	-3.19	1.36	1.44
6	A	1670	MUR	O5-C5	-3.14	1.36	1.44
4	A	1403	AI8	C12-N4	-3.11	1.24	1.29
6	B	1669	MUR	O5-C1	-2.88	1.35	1.42
4	A	1403	AI8	C8-S1	-2.85	1.75	1.82
4	A	1403	AI8	C8-C9	2.44	1.55	1.51
6	A	1670	MUR	O5-C1	-2.13	1.37	1.42
5	A	1669	1W8	C10-S2	-2.04	1.73	1.77
6	A	1670	MUR	C4-C5	-2.02	1.48	1.53

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1403	AI8	C8-S1-C4	8.05	110.26	94.47
5	A	1669	1W8	C10-C11-S1	7.50	125.01	112.61
5	A	1669	1W8	C10-C9-N5	-6.31	111.90	121.61
4	A	1403	AI8	C12-C11-N3	5.99	124.32	114.38
4	A	1403	AI8	N8-C18-N6	4.65	129.21	123.19
4	A	1403	AI8	C16-C15-C17	-4.62	123.02	129.44
6	A	1670	MUR	C9-C7-C8	-4.59	107.42	113.35
4	A	1403	AI8	C9-C8-S1	-4.43	105.28	112.61
4	A	1403	AI8	C15-C16-S3	-4.02	106.85	111.79
6	A	1670	MUR	C6-C5-C4	3.90	122.14	113.00
6	A	1670	MUR	O3-C7-C8	3.74	121.73	111.85
4	A	1403	AI8	O6-C11-C12	-3.73	116.05	120.35
6	B	1669	MUR	C1-O5-C5	3.59	120.44	113.66
6	B	1669	MUR	C1-C2-C3	3.55	116.61	110.14
6	B	1669	MUR	C6-C5-C4	3.37	120.89	113.00
4	A	1403	AI8	N6-C14-N7	-3.18	116.35	122.26
6	B	1669	MUR	C9-C7-C8	-3.06	109.40	113.35
6	A	1670	MUR	C1-C2-C3	2.92	115.47	110.14
4	A	1403	AI8	C19-C17-C15	-2.72	116.98	121.28
6	B	1669	MUR	O1-C1-C2	2.72	114.57	108.96
6	A	1670	MUR	O5-C5-C4	-2.67	104.85	109.69
6	B	1669	MUR	O3-C7-C8	2.66	118.89	111.85
6	A	1670	MUR	O3-C7-C9	2.57	113.93	107.33
6	A	1670	MUR	O3-C3-C2	2.51	115.03	108.89
6	A	1670	MUR	O1-C1-C2	2.46	114.04	108.96
6	B	1669	MUR	O5-C5-C4	-2.44	105.27	109.69
4	A	1403	AI8	O7-C21-C24	2.39	115.63	109.44
4	A	1403	AI8	C9-C7-N2	-2.35	118.97	123.28
5	A	1669	1W8	C14-C13-C15	-2.30	126.24	129.44
6	B	1669	MUR	C4-C3-C2	2.24	115.80	111.37
5	A	1669	1W8	O-C1-C	-2.05	104.13	109.44

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1669	1W8	C-C1-O-N
5	A	1669	1W8	O1-C5-N4-C6
5	A	1669	1W8	C2-C5-N4-C6
6	A	1670	MUR	C2-C3-O3-C7
4	A	1403	AI8	C11-C12-N4-O7
6	B	1669	MUR	C8-C7-O3-C3
6	A	1670	MUR	C9-C7-O3-C3

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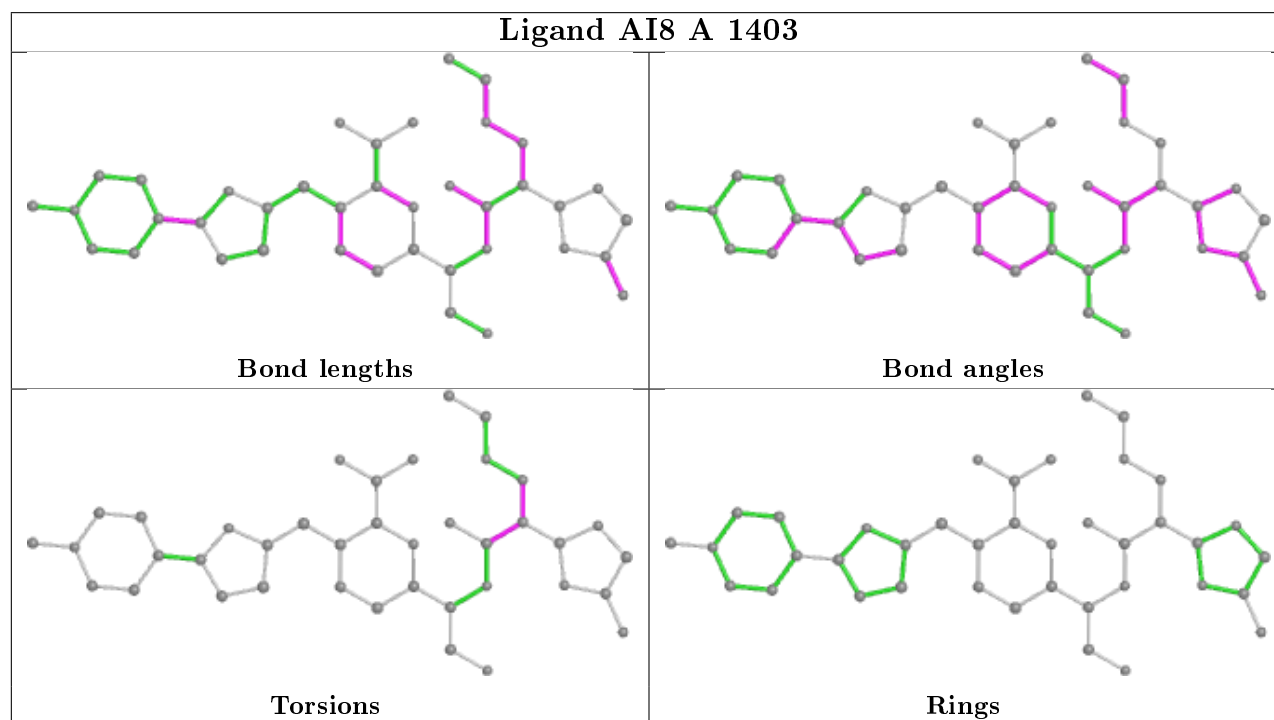
Mol	Chain	Res	Type	Atoms
4	A	1403	AI8	N3-C11-C12-N4
5	A	1669	1W8	C3-C2-C5-O1
4	A	1403	AI8	O6-C11-C12-N4

There are no ring outliers.

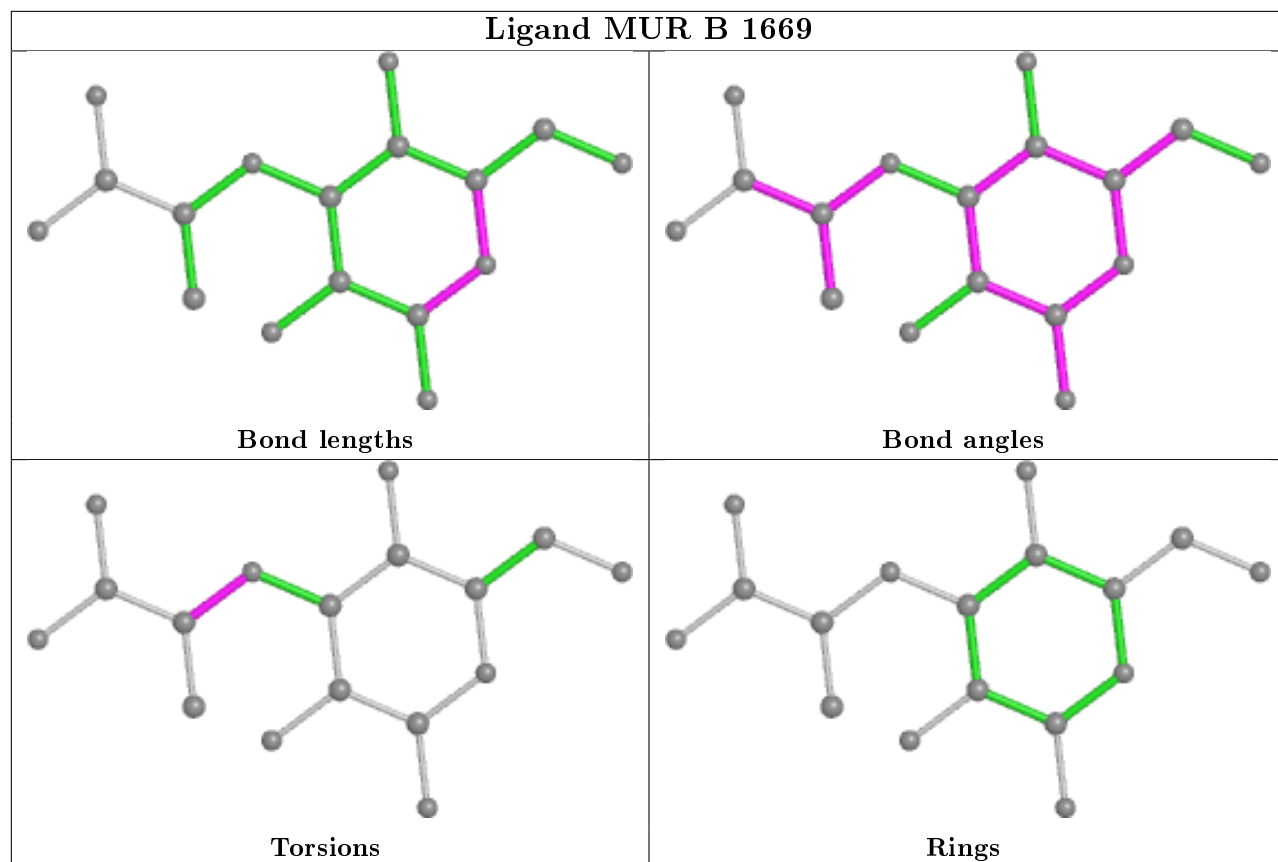
3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1403	AI8	1	0
6	A	1670	MUR	3	0
5	A	1669	1W8	7	0

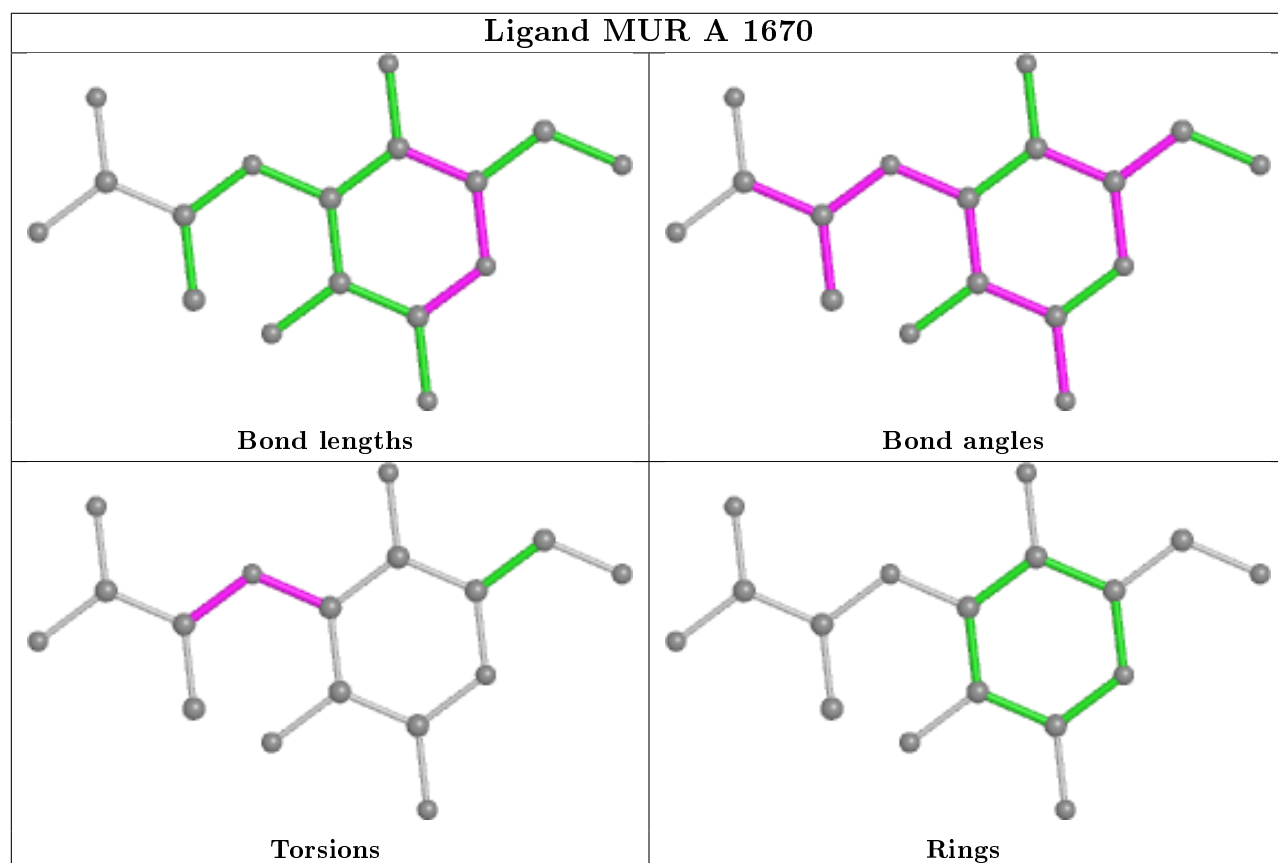
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

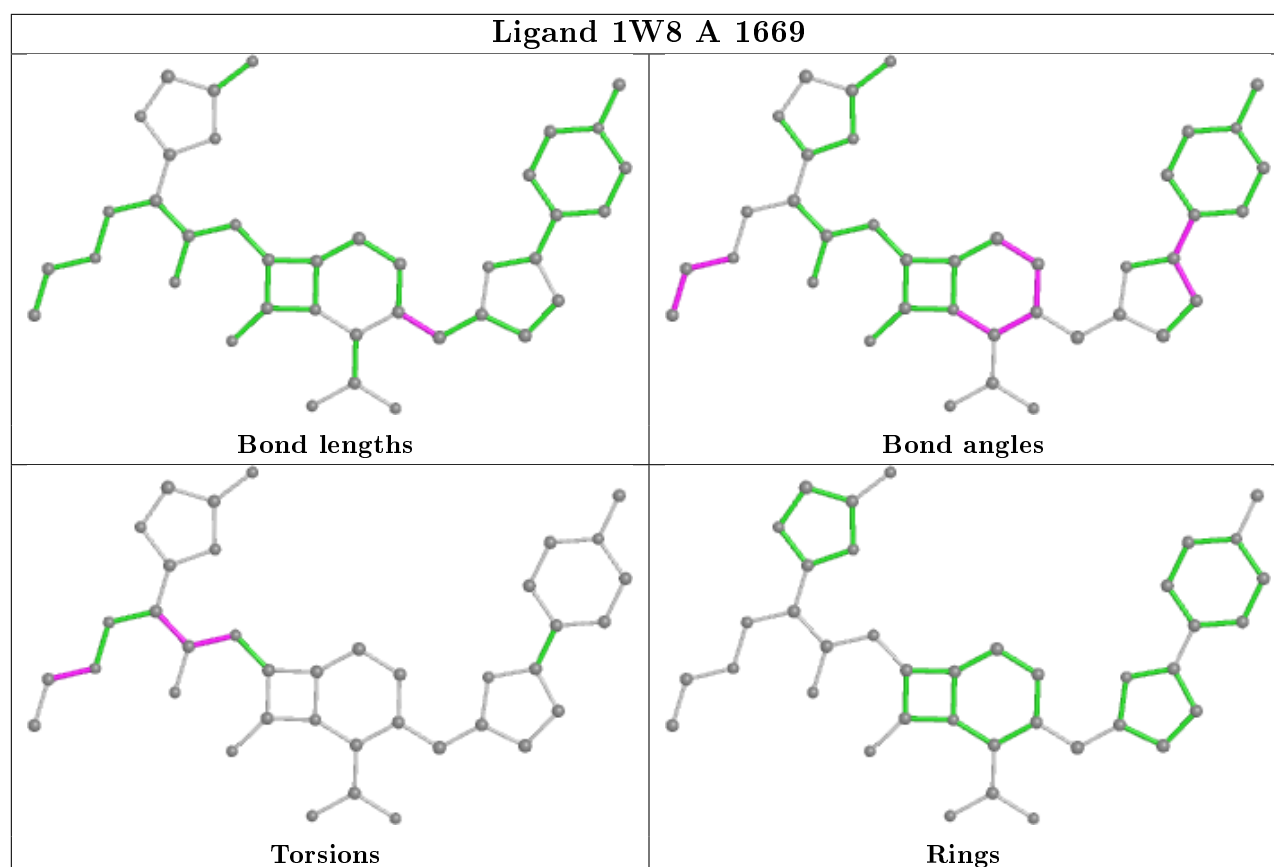


Ligand MUR B 1669



Ligand MUR A 1670





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	403:SER	C	404:THR	N	1.14

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	635/642 (98%)	0.05	11 (1%) 70 73	42, 62, 95, 129	0
1	B	642/642 (100%)	0.09	18 (2%) 53 55	43, 63, 104, 148	0
All	All	1277/1284 (99%)	0.07	29 (2%) 60 63	42, 63, 100, 148	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	261	SER	5.0
1	B	607	GLN	4.9
1	A	221	ASP	4.6
1	B	608	GLY	4.4
1	A	611	GLY	4.0
1	B	610	THR	3.9
1	B	605	MET	3.8
1	B	603	LEU	3.4
1	A	220	MET	3.3
1	B	223	TYR	3.2
1	B	506	LYS	3.1
1	B	272	TYR	3.1
1	B	268	GLU	3.1
1	A	226	ASP	3.1
1	B	606	LYS	3.0
1	A	503	ILE	3.0
1	B	184	LYS	2.8
1	B	262	GLU	2.8
1	B	269	TYR	2.7
1	B	266	GLN	2.6
1	B	270	LYS	2.5
1	A	188	LYS	2.5
1	A	224	LEU	2.4
1	B	267	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	85	ILE	2.2
1	A	186	ILE	2.2
1	A	267	LYS	2.1
1	A	204	ASN	2.1
1	B	307	ASN	2.1

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, 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.

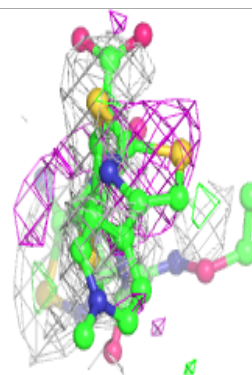
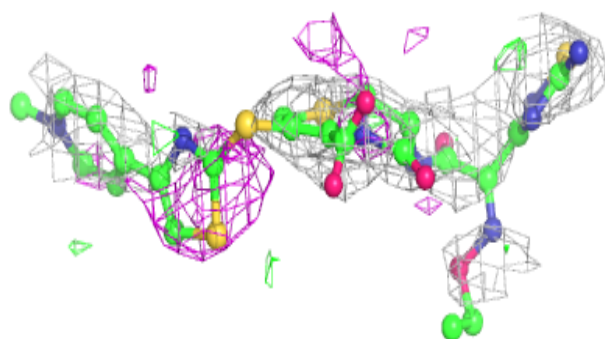
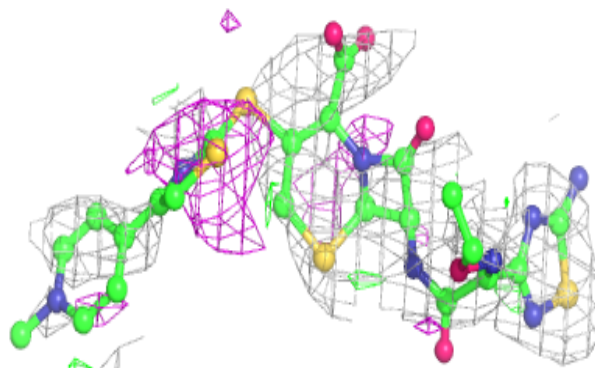
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CD	B	1006	1/1	-0.34	0.29	300,300,300,300	0
2	CD	A	1008	1/1	0.23	0.32	297,297,297,297	0
5	1W8	A	1669	39/39	0.54	0.47	130,134,136,136	0
6	MUR	B	1669	17/17	0.72	0.39	73,85,100,102	0
6	MUR	A	1670	17/17	0.84	0.30	37,51,68,73	0
4	AI8	A	1403	39/39	0.85	0.18	54,75,94,95	0
3	CL	B	1009	1/1	0.97	0.11	50,50,50,50	0
2	CD	A	1004	1/1	0.98	0.06	135,135,135,135	0
2	CD	B	1007	1/1	0.99	0.08	93,93,93,93	0
3	CL	A	1011	1/1	0.99	0.13	50,50,50,50	0
2	CD	A	1001	1/1	0.99	0.15	52,52,52,52	0
3	CL	A	1010	1/1	0.99	0.08	48,48,48,48	0
2	CD	B	1002	1/1	1.00	0.13	52,52,52,52	0
2	CD	A	1003	1/1	1.00	0.16	62,62,62,62	0
3	CL	B	1012	1/1	1.00	0.13	57,57,57,57	0
2	CD	A	1005	1/1	1.00	0.13	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.

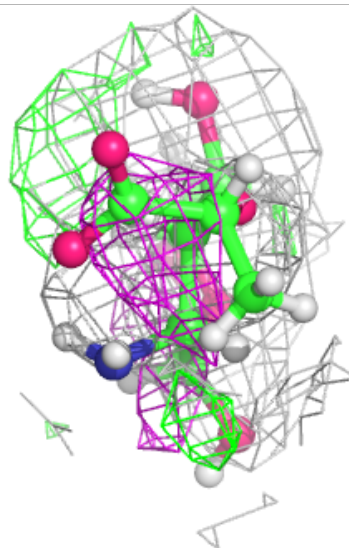
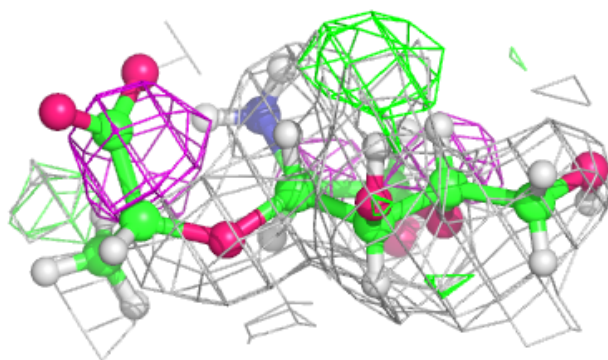
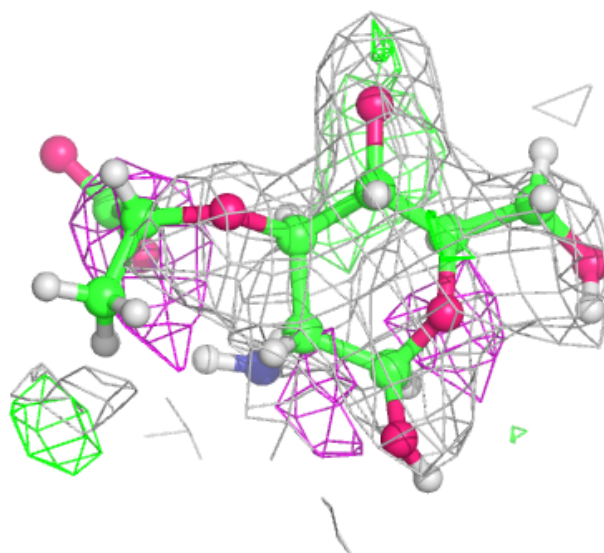
Electron density around 1W8 A 1669:

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



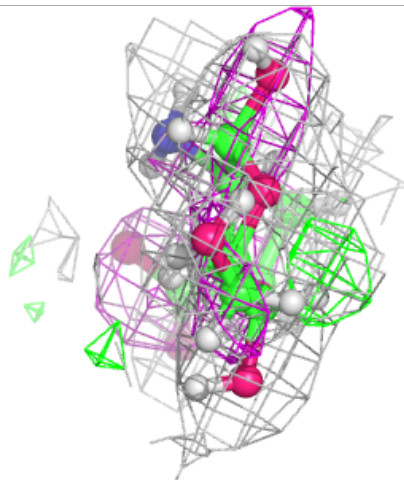
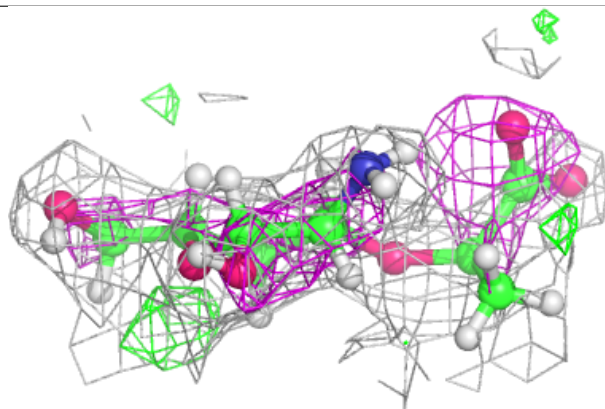
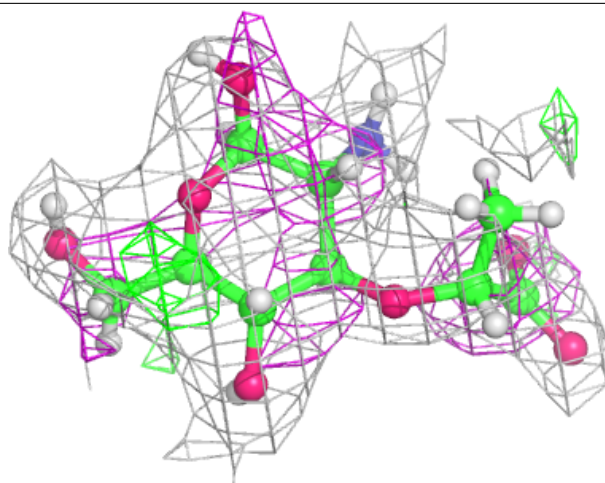
Electron density around MUR B 1669:

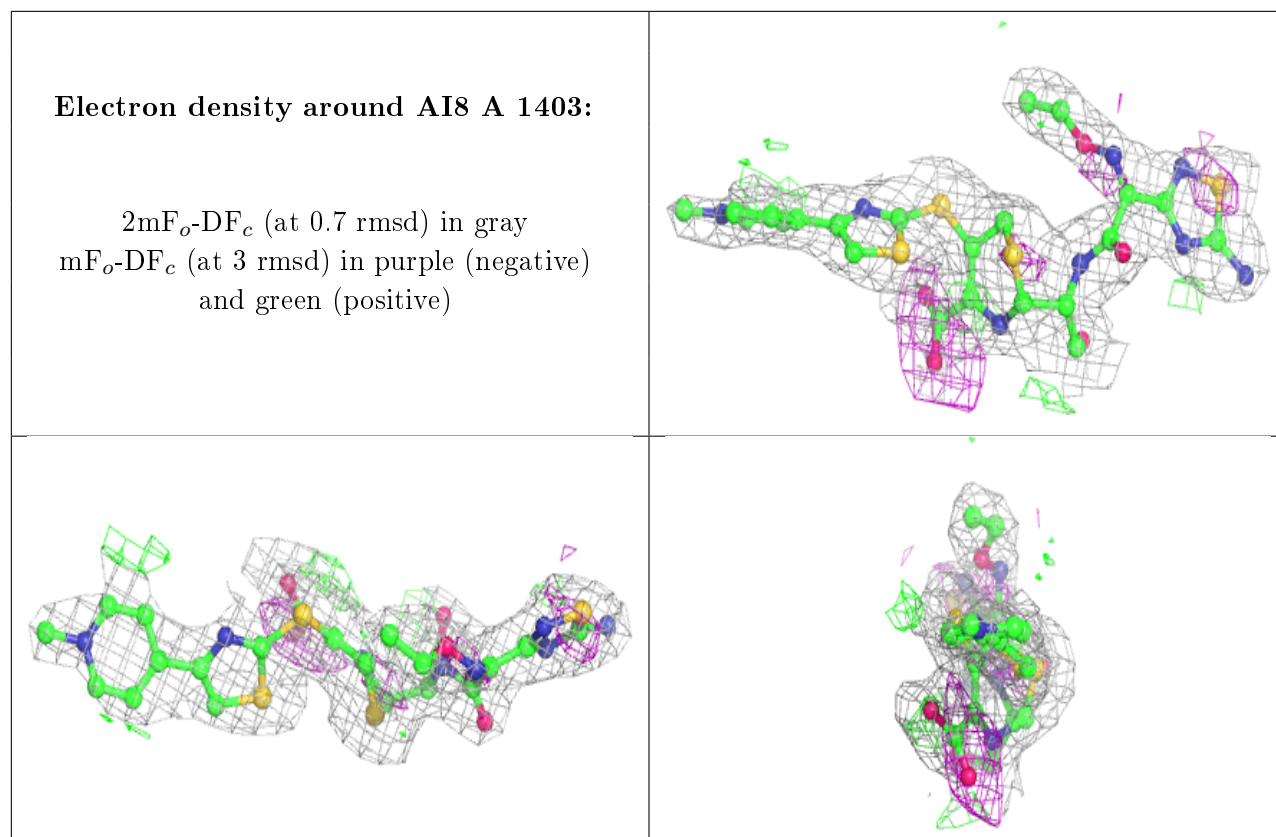
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MUR A 1670:

$2mF_o-DF_c$ (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.