



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 11, 2023 – 04:35 PM EDT

PDB ID : 8EC4
Title : XFEL structure of Mycobacterium tuberculosis beta lactamase microcrystals mixed with sulbactam for 240ms
Authors : Malla, T.N.; Schmidt, M.
Deposited on : 2022-09-01
Resolution : 2.35 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

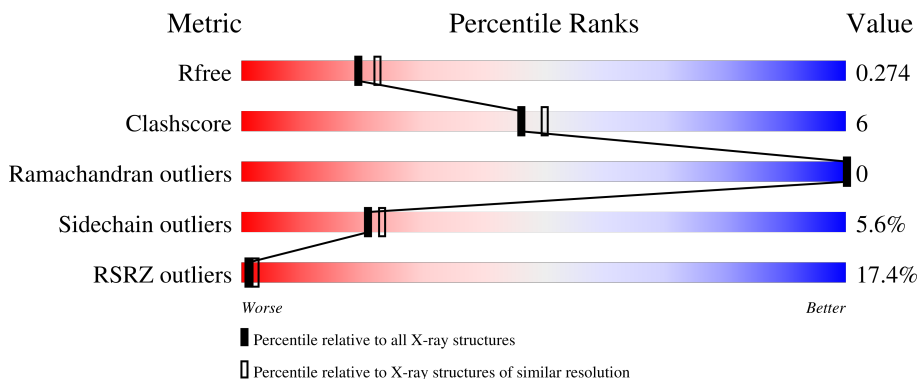
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	267	 19% 84% 15%
1	B	267	 13% 85% 13%
1	C	267	 24% 73% 23%
1	D	267	 13% 87% 12%

2 Entry composition [i](#)

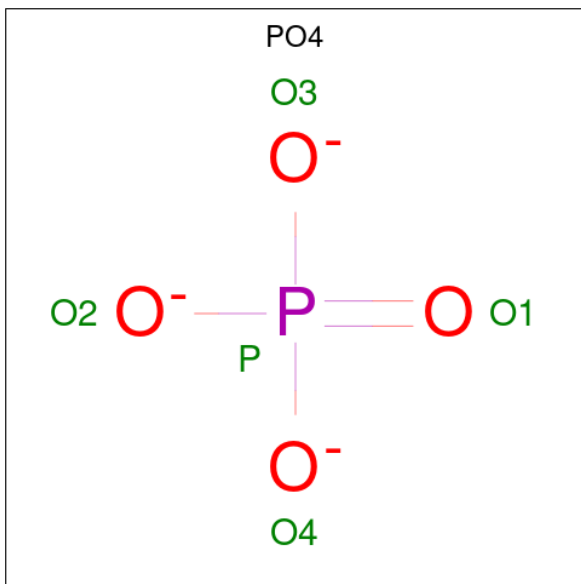
There are 4 unique types of molecules in this entry. The entry contains 8329 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	265	Total 1988	C 1243	N 352	O 387	S 6	0	0	0
1	B	265	Total 1988	C 1243	N 352	O 387	S 6	0	0	0
1	C	265	Total 1988	C 1243	N 352	O 387	S 6	0	0	0
1	D	265	Total 1988	C 1243	N 352	O 387	S 6	0	0	0

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



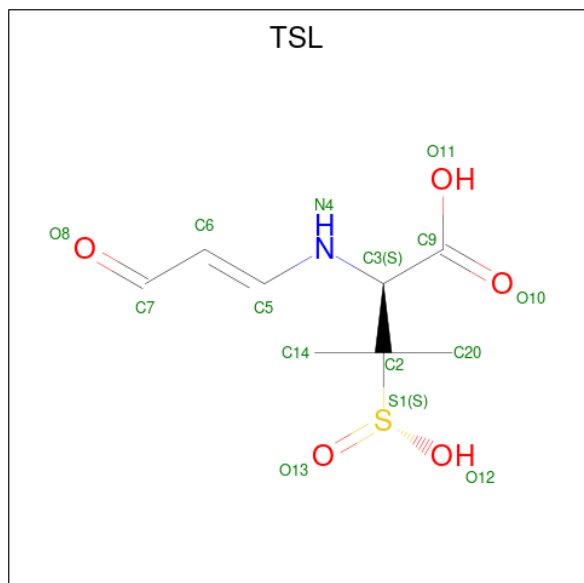
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	Total 5	O 4	P 1	0	0
2	B	1	Total 5	O 4	P 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is TRANS-ENAMINE INTERMEDIATE OF SULBACTAM (three-letter code: TSL) (formula: $C_8H_{13}NO_5S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	S	0	0
			15	8	1	5	1		
3	C	1	Total	C	N	O	S	0	0
			15	8	1	5	1		
3	D	1	Total	C	N	O	S	0	0
			15	8	1	5	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	88	Total	O	0	0
			88	88		
4	B	39	Total	O	0	0
			39	39		
4	C	118	Total	O	0	0
			118	118		

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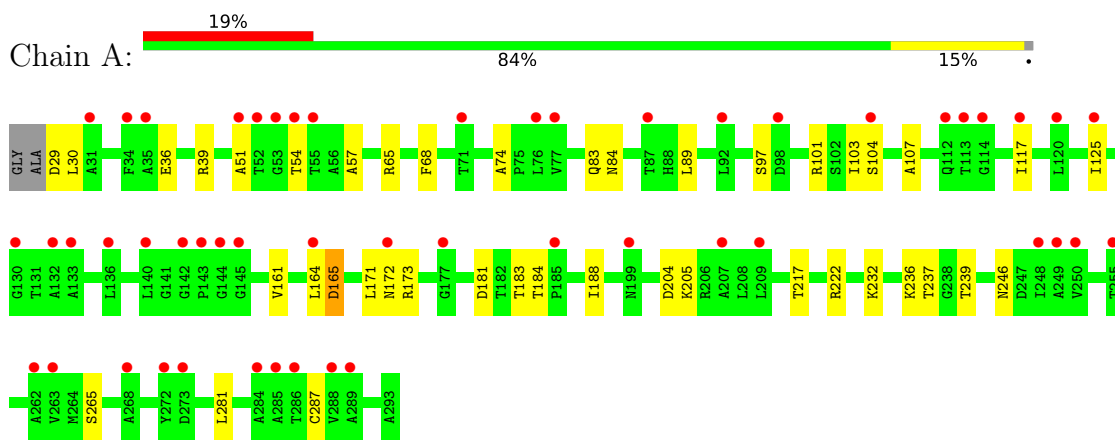
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	52	Total	O	0	0
			52	52		

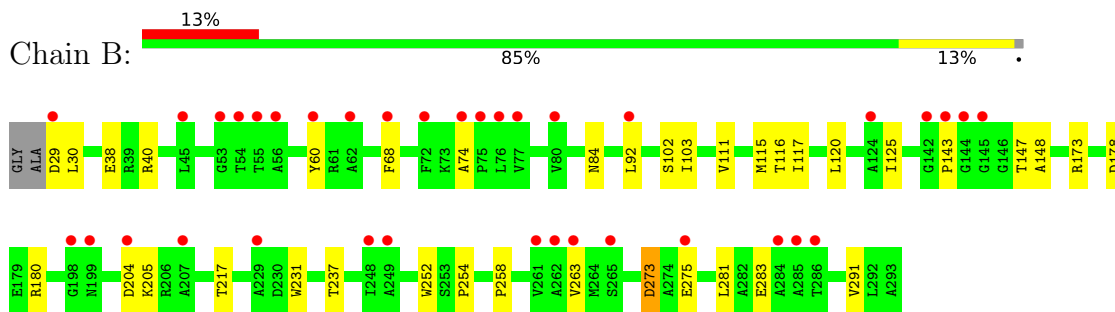
3 Residue-property plots [i](#)

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.

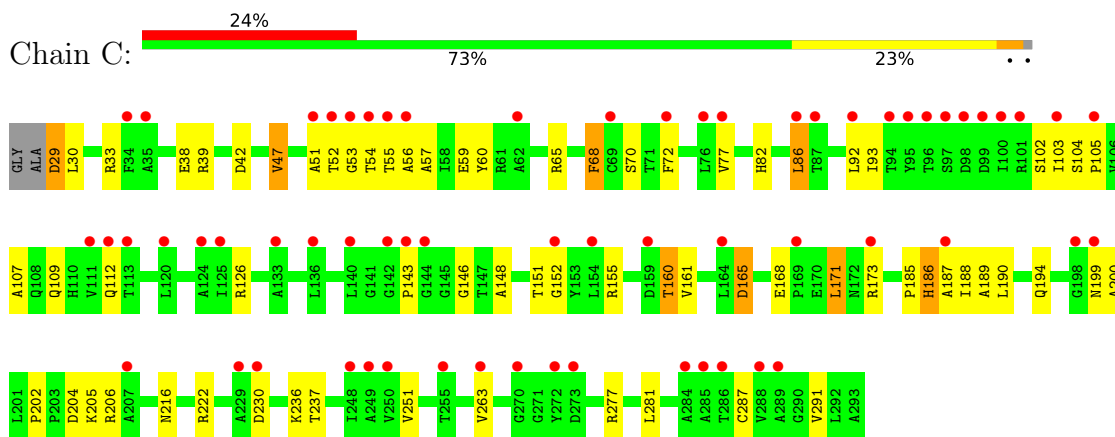
- Molecule 1: Beta-lactamase



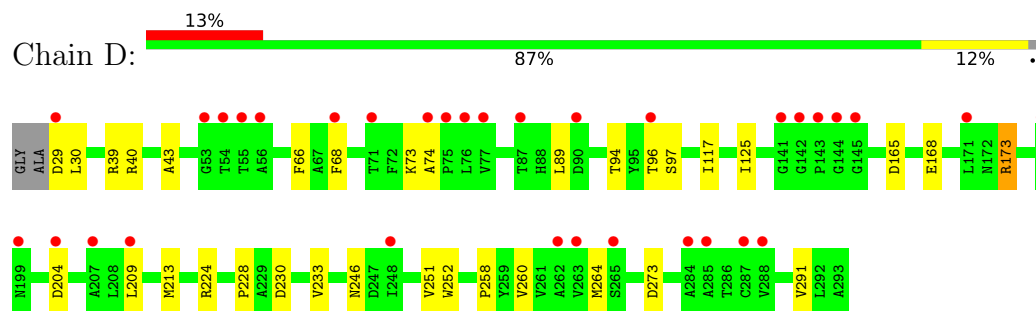
- Molecule 1: Beta-lactamase



- Molecule 1: Beta-lactamase



● Molecule 1: Beta-lactamase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.82Å 97.89Å 113.17Å 90.00° 109.34° 90.00°	Depositor
Resolution (Å)	22.42 – 2.35 22.42 – 2.35	Depositor EDS
% Data completeness (in resolution range)	78.5 (22.42-2.35) 78.5 (22.42-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 2.36Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.241 , 0.275 0.241 , 0.274	Depositor DCC
R_{free} test set	2696 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	58.0	Xtrriage
Anisotropy	0.094	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 46.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.071 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8329	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.40 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.7667e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: TSL, PO4

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	0.24	0/2027	0.50	0/2767
1	B	0.24	0/2027	0.49	0/2767
1	C	0.24	0/2027	0.51	0/2767
1	D	0.23	0/2027	0.50	0/2767
All	All	0.24	0/8108	0.50	0/11068

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 [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	1988	0	1955	24	0
1	B	1988	0	1955	17	0
1	C	1988	0	1955	45	0
1	D	1988	0	1955	17	0
2	A	5	0	0	0	0
2	B	5	0	0	1	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	15	0	10	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	15	0	11	0	0
3	C	15	0	10	0	0
3	D	15	0	11	0	0
4	A	88	0	0	8	0
4	B	39	0	0	2	0
4	C	118	0	0	13	0
4	D	52	0	0	3	0
All	All	8329	0	7862	100	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 6.

All (100) 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:C:287:CYS:O	4:C:402:HOH:O	2.10	0.70
1:C:187:ALA:O	4:C:401:HOH:O	2.10	0.70
1:A:183:THR:OG1	4:A:401:HOH:O	2.09	0.69
1:A:173:ARG:NH2	3:A:302:TSL:O12	2.27	0.68
1:C:185:PRO:O	4:C:403:HOH:O	2.13	0.66
1:D:233:VAL:HG22	1:D:251:VAL:HG12	1.80	0.64
1:A:222:ARG:HD3	1:A:281:LEU:HD12	1.81	0.62
1:B:29:ASP:N	4:B:404:HOH:O	2.32	0.62
1:C:202:PRO:HG2	1:C:205:LYS:HG3	1.80	0.62
1:B:117:ILE:HA	1:B:120:LEU:HD12	1.83	0.60
1:C:33:ARG:NH2	4:C:413:HOH:O	2.33	0.60
1:C:199:ASN:HA	1:C:206:ARG:HH12	1.67	0.60
1:D:230:ASP:N	1:D:230:ASP:OD1	2.35	0.60
1:C:65:ARG:NH2	1:C:161:VAL:O	2.36	0.59
1:A:184:THR:O	4:A:401:HOH:O	2.16	0.59
1:D:43:ALA:N	4:D:401:HOH:O	2.34	0.58
1:C:29:ASP:N	4:C:415:HOH:O	2.36	0.58
1:C:86:LEU:H	1:C:86:LEU:HD23	1.69	0.58
1:A:39:ARG:NH1	4:A:405:HOH:O	2.29	0.58
1:A:103:ILE:HD11	1:B:111:VAL:HG21	1.86	0.58
1:A:30:LEU:HD11	1:A:287:CYS:HB3	1.86	0.57
1:A:65:ARG:NH2	1:A:161:VAL:O	2.37	0.57
1:C:82:HIS:ND1	4:C:412:HOH:O	2.32	0.57
1:D:29:ASP:N	4:D:408:HOH:O	2.37	0.56
1:A:204:ASP:OD1	1:A:205:LYS:N	2.38	0.56
1:D:30:LEU:HD22	1:D:291:VAL:HG21	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:224:ARG:NH2	1:D:228:PRO:O	2.38	0.56
1:D:74:ALA:HA	1:D:125:ILE:HD11	1.87	0.55
1:B:84:ASN:O	1:B:205:LYS:NZ	2.40	0.55
1:C:51:ALA:HB2	1:C:57:ALA:HB2	1.88	0.55
1:C:155:ARG:N	4:C:408:HOH:O	2.36	0.54
1:A:172:ASN:ND2	4:A:413:HOH:O	2.40	0.54
1:A:104:SER:HB3	1:A:107:ALA:HB3	1.89	0.54
1:B:147:THR:OG1	4:B:401:HOH:O	2.17	0.53
1:A:171:LEU:H	1:A:171:LEU:HD23	1.73	0.53
1:A:89:LEU:HD23	1:A:117:ILE:HG22	1.89	0.53
1:D:192:LEU:HD23	1:D:260:VAL:HG13	1.91	0.53
1:C:103:ILE:HG22	1:D:97:SER:HB3	1.91	0.52
1:C:109:GLN:HB3	1:D:173:ARG:HH12	1.74	0.52
1:C:189:ALA:N	4:C:403:HOH:O	2.43	0.52
1:B:102:SER:OG	1:B:103:ILE:N	2.43	0.52
1:B:38:GLU:OE1	1:B:60:TYR:OH	2.26	0.51
1:C:47:VAL:HG13	1:C:263:VAL:HG22	1.92	0.51
1:D:73:LYS:NZ	1:D:168:GLU:OE2	2.42	0.51
1:A:239:THR:O	3:A:302:TSL:H5	2.09	0.51
1:C:104:SER:HB3	1:C:107:ALA:HB3	1.92	0.51
1:A:83:GLN:NE2	4:A:415:HOH:O	2.45	0.50
1:C:146:GLY:HA3	4:C:407:HOH:O	2.11	0.50
1:C:222:ARG:HD3	1:C:281:LEU:HD12	1.93	0.50
1:C:236:LYS:NZ	1:C:237:THR:O	2.45	0.50
1:C:39:ARG:NH2	1:C:42:ASP:OD1	2.45	0.49
1:C:38:GLU:OE1	1:C:60:TYR:OH	2.20	0.49
1:C:168:GLU:HG2	1:C:171:LEU:HD21	1.94	0.49
1:D:39:ARG:NH2	4:D:411:HOH:O	2.44	0.49
1:B:40:ARG:NH1	1:B:283:GLU:OE1	2.46	0.48
1:C:188:ILE:HG13	1:C:188:ILE:O	2.13	0.48
1:C:291:VAL:HG22	4:C:402:HOH:O	2.13	0.48
1:C:155:ARG:HD3	1:C:160:THR:HA	1.94	0.48
1:C:55:THR:OG1	1:C:56:ALA:N	2.46	0.48
1:C:52:THR:OG1	1:C:53:GLY:N	2.46	0.48
1:A:236:LYS:NZ	1:A:237:THR:O	2.47	0.47
1:C:68:PHE:HE1	1:C:188:ILE:HD13	1.79	0.47
1:D:89:LEU:HA	1:D:117:ILE:HB	1.97	0.46
1:B:30:LEU:HD22	1:B:291:VAL:HG21	1.98	0.46
1:B:143:PRO:HD2	1:B:148:ALA:HB3	1.98	0.46
1:C:186:HIS:O	1:C:186:HIS:ND1	2.44	0.46
1:C:186:HIS:HE1	1:C:190:LEU:HD11	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:66:PHE:HB2	1:D:264:MET:HE2	1.96	0.46
1:C:230:ASP:OD1	1:C:230:ASP:N	2.36	0.45
1:D:209:LEU:O	1:D:213:MET:HG3	2.17	0.45
1:B:237:THR:OG1	2:B:301:PO4:O4	2.35	0.45
1:C:55:THR:HG23	4:C:405:HOH:O	2.17	0.45
1:B:74:ALA:HA	1:B:125:ILE:HD11	1.99	0.44
1:A:101:ARG:HA	1:A:101:ARG:HD2	1.83	0.44
1:A:165:ASP:N	1:A:181:ASP:OD1	2.47	0.44
1:C:105:PRO:O	1:C:109:GLN:NE2	2.41	0.44
1:B:92:LEU:HD12	1:B:116:THR:HG22	2.00	0.44
1:B:273:ASP:OD2	1:B:273:ASP:N	2.51	0.44
1:C:92:LEU:HD13	1:C:93:ILE:N	2.32	0.43
1:A:83:GLN:NE2	4:A:422:HOH:O	2.51	0.43
1:A:74:ALA:HA	1:A:125:ILE:HD11	2.01	0.43
1:C:200:ALA:H	1:C:206:ARG:HH22	1.66	0.43
1:D:252:TRP:NE1	1:D:258:PRO:HB3	2.34	0.43
1:D:204:ASP:OD1	1:D:204:ASP:N	2.47	0.43
1:A:164:LEU:HD11	1:A:171:LEU:HD13	2.01	0.42
1:C:30:LEU:HD23	1:C:30:LEU:HA	1.81	0.42
1:C:165:ASP:N	1:C:165:ASP:OD1	2.52	0.42
1:C:194:GLN:O	1:C:206:ARG:NH2	2.52	0.42
1:C:200:ALA:H	1:C:206:ARG:NH2	2.18	0.42
1:A:84:ASN:ND2	4:A:408:HOH:O	2.33	0.42
1:A:188:ILE:N	4:A:401:HOH:O	2.52	0.42
1:B:231:TRP:CE2	1:B:254:PRO:HD3	2.54	0.42
1:C:148:ALA:O	1:C:152:GLY:N	2.51	0.42
1:C:102:SER:N	4:C:409:HOH:O	2.46	0.41
1:C:143:PRO:O	4:C:406:HOH:O	2.22	0.41
1:C:126:ARG:HA	1:C:216:ASN:HD22	1.84	0.41
1:B:263:VAL:HG12	1:B:281:LEU:HD22	2.02	0.41
1:B:252:TRP:NE1	1:B:258:PRO:HB3	2.36	0.40
1:C:173:ARG:HE	1:C:173:ARG:HB3	1.60	0.40
1:A:51:ALA:HB2	1:A:57:ALA: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	263/267 (98%)	256 (97%)	7 (3%)	0	100	100
1	B	263/267 (98%)	252 (96%)	11 (4%)	0	100	100
1	C	263/267 (98%)	252 (96%)	11 (4%)	0	100	100
1	D	263/267 (98%)	253 (96%)	10 (4%)	0	100	100
All	All	1052/1068 (98%)	1013 (96%)	39 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	202/202 (100%)	192 (95%)	10 (5%)	24	28
1	B	202/202 (100%)	193 (96%)	9 (4%)	27	33
1	C	202/202 (100%)	184 (91%)	18 (9%)	9	8
1	D	202/202 (100%)	194 (96%)	8 (4%)	31	39
All	All	808/808 (100%)	763 (94%)	45 (6%)	21	23

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	36	GLU

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Mol	Chain	Res	Type
1	A	54	THR
1	A	68	PHE
1	A	97	SER
1	A	165	ASP
1	A	217	THR
1	A	232	LYS
1	A	246	ASN
1	A	265	SER
1	B	68	PHE
1	B	115	MET
1	B	173	ARG
1	B	178	ASP
1	B	180	ARG
1	B	204	ASP
1	B	217	THR
1	B	273	ASP
1	B	275	GLU
1	C	29	ASP
1	C	47	VAL
1	C	54	THR
1	C	59	GLU
1	C	68	PHE
1	C	70	SER
1	C	72	PHE
1	C	77	VAL
1	C	86	LEU
1	C	112	GLN
1	C	151	THR
1	C	160	THR
1	C	165	ASP
1	C	171	LEU
1	C	186	HIS
1	C	204	ASP
1	C	251	VAL
1	C	277	ARG
1	D	40	ARG
1	D	68	PHE
1	D	94	THR
1	D	96	THR
1	D	165	ASP
1	D	173	ARG
1	D	246	ASN

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Mol	Chain	Res	Type
1	D	273	ASP

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

Mol	Chain	Res	Type
1	B	112	GLN
1	C	216	ASN
1	D	194	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	PO4	A	301	-	4,4,4	0.92	0	6,6,6	0.44	0
2	PO4	B	301	-	4,4,4	0.91	0	6,6,6	0.46	0
2	PO4	C	301	-	4,4,4	0.94	0	6,6,6	0.40	0
3	TSL	A	302	1	9,14,14	1.66	2 (22%)	9,19,19	3.54	1 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	TSL	D	302	1	9,14,14	1.66	2 (22%)	9,19,19	3.35	2 (22%)
3	TSL	C	302	1	9,14,14	1.70	2 (22%)	9,19,19	3.60	2 (22%)
2	PO4	D	301	-	4,4,4	0.91	0	6,6,6	0.42	0
3	TSL	B	302	1	9,14,14	1.66	2 (22%)	9,19,19	3.51	1 (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
3	TSL	A	302	1	-	8/17/21/21	-
3	TSL	B	302	1	-	9/17/21/21	-
3	TSL	C	302	1	-	3/17/21/21	-
3	TSL	D	302	1	-	8/17/21/21	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	TSL	C6-C7	3.32	1.54	1.44
3	C	302	TSL	C6-C7	3.25	1.54	1.44
3	A	302	TSL	C6-C7	3.25	1.54	1.44
3	D	302	TSL	C6-C7	3.10	1.53	1.44
3	D	302	TSL	C5-C6	-2.29	1.33	1.37
3	A	302	TSL	C5-C6	-2.17	1.33	1.37
3	C	302	TSL	C5-C6	-2.02	1.33	1.37
3	B	302	TSL	C5-C6	-2.02	1.33	1.37

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	302	TSL	C5-C6-C7	10.29	152.84	121.80
3	A	302	TSL	C5-C6-C7	10.16	152.47	121.80
3	B	302	TSL	C5-C6-C7	10.09	152.25	121.80
3	D	302	TSL	C5-C6-C7	9.52	150.54	121.80
3	C	302	TSL	O11-C9-O10	-2.25	118.98	124.09
3	D	302	TSL	O8-C7-C6	-2.11	118.48	125.67

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	302	TSL	C14-C2-C3-C9
3	A	302	TSL	C14-C2-C3-N4
3	A	302	TSL	S1-C2-C3-C9
3	A	302	TSL	S1-C2-C3-N4
3	A	302	TSL	C20-C2-C3-C9
3	A	302	TSL	C20-C2-C3-N4
3	A	302	TSL	C6-C5-N4-C3
3	A	302	TSL	C5-C6-C7-O8
3	B	302	TSL	C14-C2-C3-C9
3	B	302	TSL	C14-C2-C3-N4
3	B	302	TSL	S1-C2-C3-C9
3	B	302	TSL	S1-C2-C3-N4
3	B	302	TSL	C20-C2-C3-C9
3	B	302	TSL	C20-C2-C3-N4
3	B	302	TSL	C6-C5-N4-C3
3	B	302	TSL	C5-C6-C7-O8
3	C	302	TSL	C6-C5-N4-C3
3	C	302	TSL	C5-C6-C7-O8
3	D	302	TSL	C14-C2-C3-C9
3	D	302	TSL	C14-C2-C3-N4
3	D	302	TSL	S1-C2-C3-C9
3	D	302	TSL	S1-C2-C3-N4
3	D	302	TSL	C20-C2-C3-C9
3	D	302	TSL	C20-C2-C3-N4
3	D	302	TSL	C6-C5-N4-C3
3	D	302	TSL	C5-C6-C7-O8
3	B	302	TSL	C2-C3-N4-C5
3	C	302	TSL	C20-C2-C3-N4

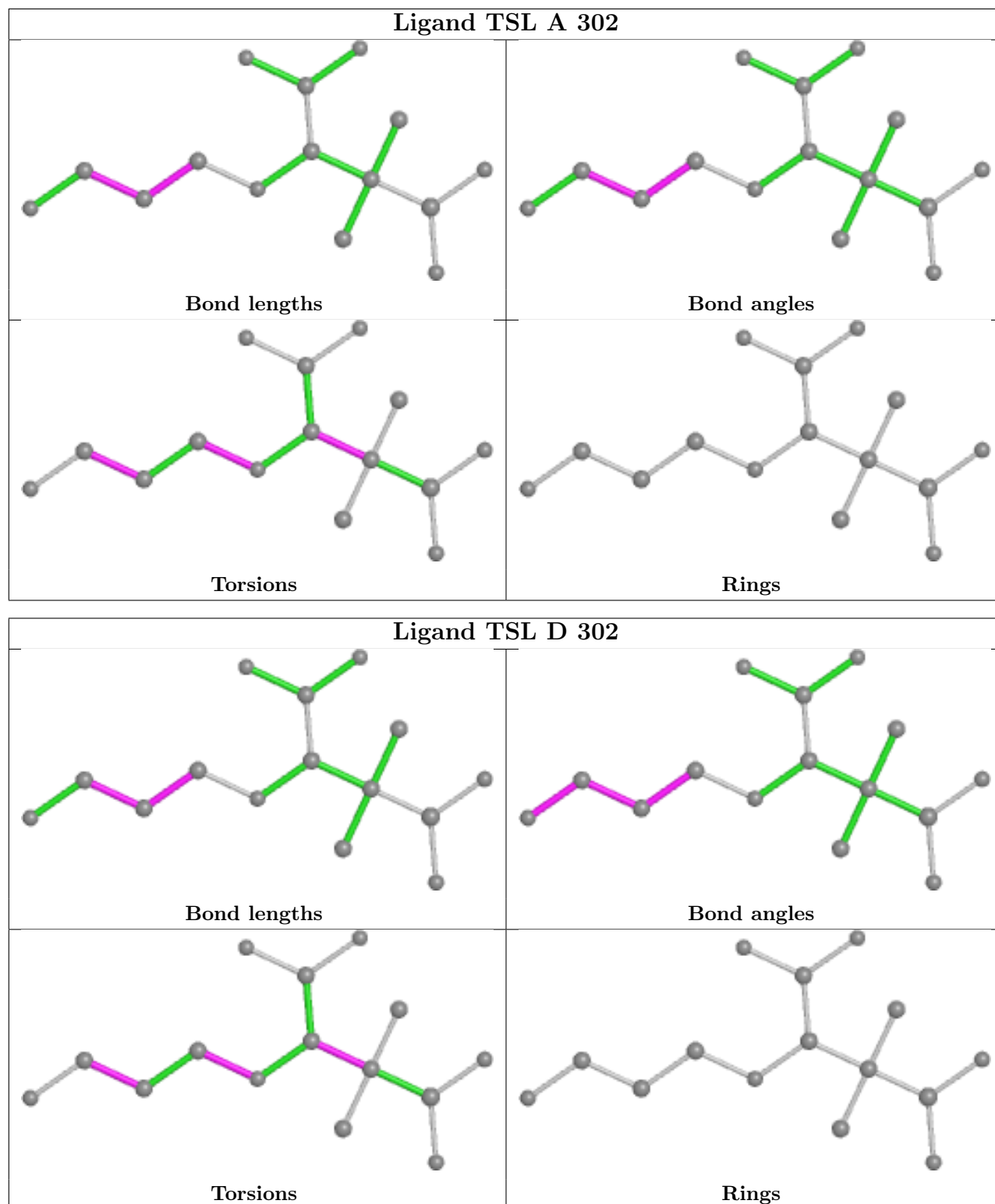
There are no ring outliers.

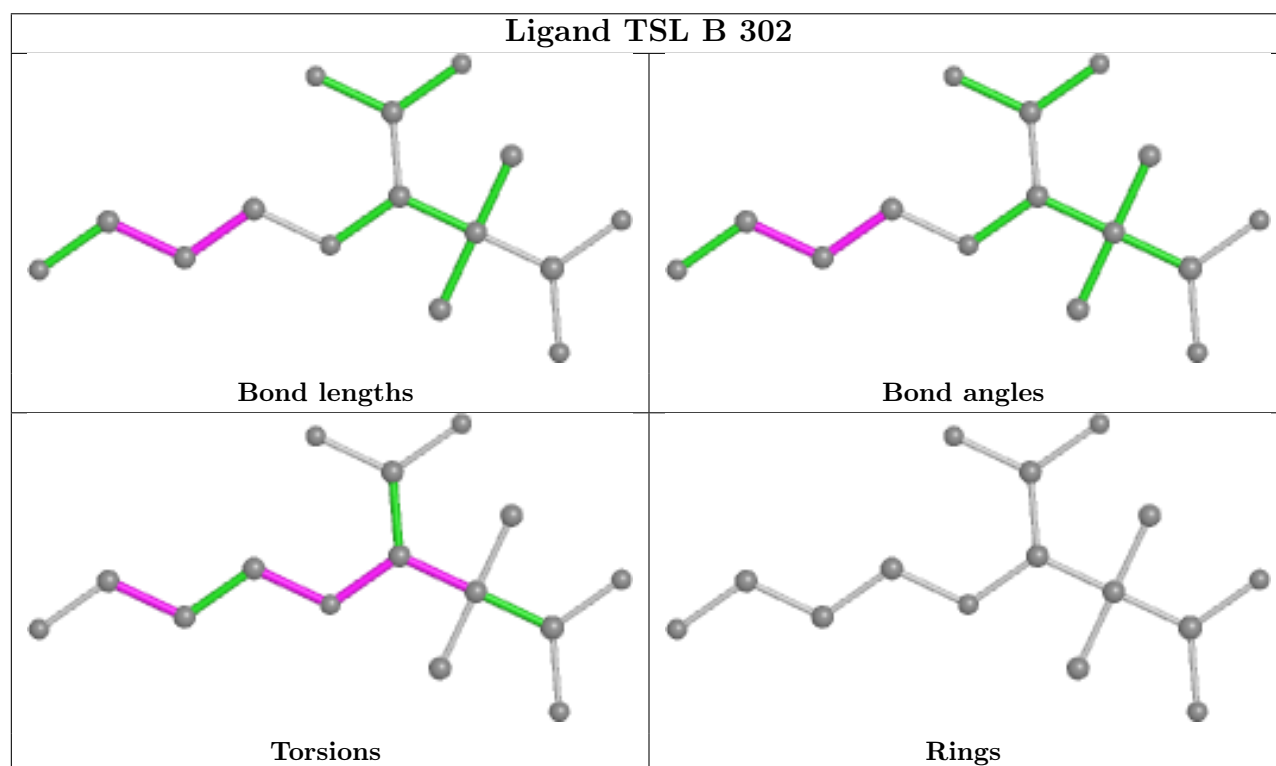
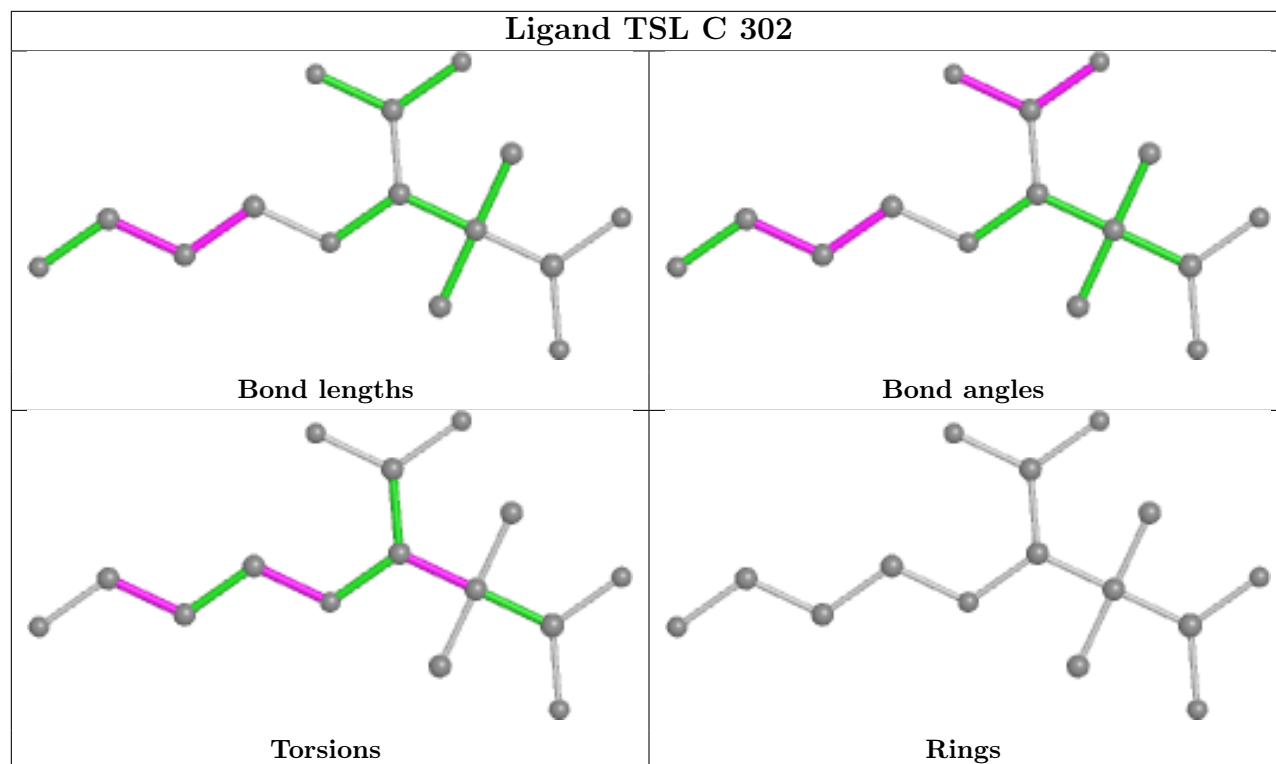
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	PO4	1	0
3	A	302	TSL	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 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

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	265/267 (99%)	1.14	51 (19%) 1 2	45, 58, 75, 92	0
1	B	265/267 (99%)	0.89	36 (13%) 3 4	49, 58, 73, 93	0
1	C	265/267 (99%)	1.32	63 (23%) 0 1	45, 61, 80, 93	0
1	D	265/267 (99%)	0.81	34 (12%) 3 6	48, 58, 74, 84	0
All	All	1060/1068 (99%)	1.04	184 (17%) 1 2	45, 59, 77, 93	0

All (184) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	144	GLY	7.7
1	A	53	GLY	6.6
1	A	142	GLY	6.4
1	C	87	THR	5.8
1	B	143	PRO	5.8
1	C	52	THR	5.8
1	C	53	GLY	5.5
1	C	55	THR	5.4
1	A	144	GLY	5.2
1	A	143	PRO	5.2
1	C	94	THR	5.1
1	A	114	GLY	5.0
1	D	53	GLY	4.8
1	C	96	THR	4.8
1	B	54	THR	4.7
1	B	263	VAL	4.4
1	C	285	ALA	4.4
1	A	87	THR	4.3
1	B	77	VAL	4.3
1	D	77	VAL	4.2
1	A	263	VAL	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	144	GLY	4.1
1	C	92	LEU	4.1
1	C	140	LEU	3.9
1	B	55	THR	3.9
1	C	263	VAL	3.9
1	D	143	PRO	3.9
1	B	144	GLY	3.9
1	C	143	PRO	3.8
1	A	268	ALA	3.7
1	B	76	LEU	3.7
1	D	76	LEU	3.7
1	A	55	THR	3.6
1	C	207	ALA	3.6
1	C	272	TYR	3.6
1	C	289	ALA	3.6
1	C	136	LEU	3.6
1	C	54	THR	3.6
1	B	60	TYR	3.5
1	B	261	VAL	3.5
1	C	249	ALA	3.5
1	C	56	ALA	3.4
1	C	198	GLY	3.4
1	B	248	ILE	3.4
1	A	285	ALA	3.4
1	A	98	ASP	3.4
1	D	141	GLY	3.4
1	D	288	VAL	3.3
1	D	90	ASP	3.3
1	C	164	LEU	3.3
1	C	270	GLY	3.3
1	C	284	ALA	3.3
1	A	92	LEU	3.3
1	B	53	GLY	3.3
1	A	272	TYR	3.3
1	B	75	PRO	3.3
1	B	74	ALA	3.3
1	C	173	ARG	3.2
1	C	113	THR	3.2
1	D	56	ALA	3.2
1	B	142	GLY	3.2
1	D	142	GLY	3.2
1	C	112	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	54	THR	3.2
1	A	199	ASN	3.1
1	C	97	SER	3.1
1	C	99	ASP	3.1
1	C	199	ASN	3.0
1	A	177	GLY	3.0
1	A	112	GLN	3.0
1	C	34	PHE	3.0
1	C	288	VAL	3.0
1	D	75	PRO	2.9
1	C	62	ALA	2.9
1	C	95	TYR	2.9
1	A	76	LEU	2.9
1	D	68	PHE	2.9
1	D	263	VAL	2.9
1	C	35	ALA	2.9
1	D	207	ALA	2.8
1	C	98	ASP	2.8
1	A	34	PHE	2.8
1	B	199	ASN	2.8
1	B	229	ALA	2.8
1	A	54	THR	2.7
1	A	35	ALA	2.7
1	D	262	ALA	2.7
1	C	255	THR	2.7
1	A	164	LEU	2.7
1	A	51	ALA	2.7
1	A	249	ALA	2.7
1	A	136	LEU	2.7
1	A	250	VAL	2.7
1	C	286	THR	2.7
1	A	71	THR	2.6
1	D	171	LEU	2.6
1	C	120	LEU	2.6
1	C	187	ALA	2.6
1	C	101	ARG	2.6
1	A	77	VAL	2.6
1	B	68	PHE	2.6
1	B	198	GLY	2.6
1	C	229	ALA	2.6
1	C	142	GLY	2.5
1	C	51	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	248	ILE	2.5
1	A	262	ALA	2.5
1	B	29	ASP	2.5
1	B	207	ALA	2.5
1	D	199	ASN	2.5
1	A	132	ALA	2.4
1	B	56	ALA	2.4
1	B	286	THR	2.4
1	C	125	ILE	2.4
1	D	209	LEU	2.4
1	A	288	VAL	2.4
1	D	248	ILE	2.4
1	C	77	VAL	2.4
1	C	230	ASP	2.4
1	D	145	GLY	2.4
1	A	284	ALA	2.4
1	D	87	THR	2.4
1	A	273	ASP	2.3
1	C	248	ILE	2.3
1	D	55	THR	2.3
1	A	209	LEU	2.3
1	D	284	ALA	2.3
1	A	52	THR	2.3
1	A	120	LEU	2.3
1	B	284	ALA	2.3
1	B	124	ALA	2.3
1	B	249	ALA	2.3
1	C	273	ASP	2.3
1	A	113	THR	2.3
1	A	255	THR	2.3
1	D	198	GLY	2.3
1	B	265	SER	2.3
1	B	285	ALA	2.3
1	C	124	ALA	2.3
1	B	275	GLU	2.2
1	C	72	PHE	2.2
1	C	76	LEU	2.2
1	C	111	VAL	2.2
1	A	31	ALA	2.2
1	A	289	ALA	2.2
1	C	86	LEU	2.2
1	A	145	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	207	ALA	2.2
1	D	96	THR	2.2
1	D	177	GLY	2.2
1	B	62	ALA	2.2
1	B	262	ALA	2.2
1	C	152	GLY	2.2
1	C	250	VAL	2.2
1	C	69	CYS	2.2
1	D	29	ASP	2.2
1	C	100	ILE	2.1
1	D	287	CYS	2.1
1	C	169	PRO	2.1
1	B	72	PHE	2.1
1	C	105	PRO	2.1
1	A	286	THR	2.1
1	D	71	THR	2.1
1	B	204	ASP	2.1
1	B	92	LEU	2.1
1	D	74	ALA	2.1
1	D	285	ALA	2.1
1	B	145	GLY	2.1
1	C	103	ILE	2.1
1	D	265	SER	2.1
1	B	45	LEU	2.1
1	A	133	ALA	2.1
1	A	185	PRO	2.1
1	A	140	LEU	2.0
1	A	172	ASN	2.0
1	D	204	ASP	2.0
1	A	125	ILE	2.0
1	C	154	LEU	2.0
1	B	80	VAL	2.0
1	A	117	ILE	2.0
1	A	130	GLY	2.0
1	A	104	SER	2.0
1	C	159	ASP	2.0
1	C	133	ALA	2.0

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

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

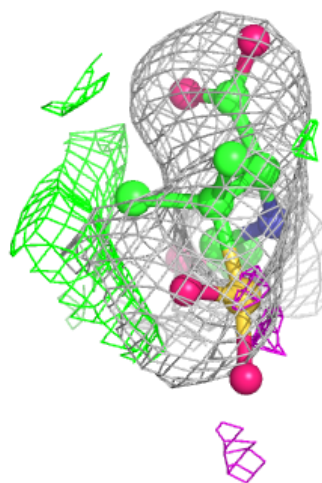
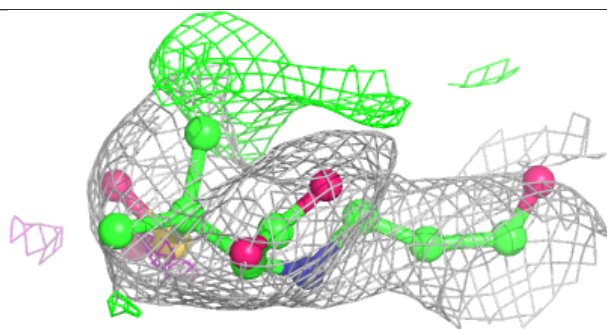
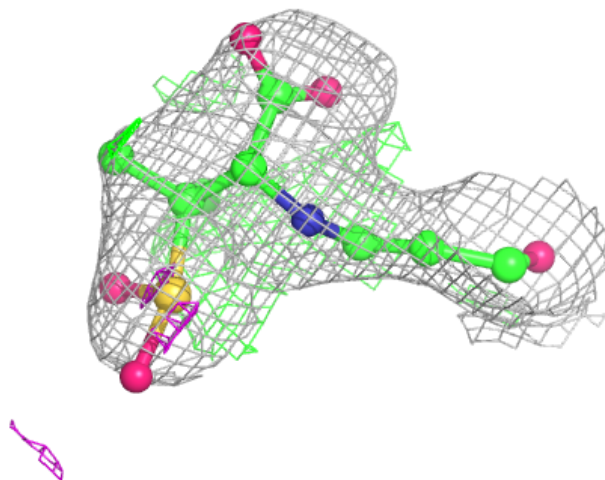
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	TSL	C	302	15/15	0.77	0.30	58,68,81,84	15
3	TSL	B	302	15/15	0.85	0.19	61,72,80,89	0
3	TSL	D	302	15/15	0.85	0.22	55,65,78,80	0
3	TSL	A	302	15/15	0.88	0.21	58,65,78,80	15
2	PO4	B	301	5/5	0.97	0.12	55,56,59,60	0
2	PO4	C	301	5/5	0.98	0.15	52,57,62,63	0
2	PO4	D	301	5/5	0.98	0.15	53,55,60,62	0
2	PO4	A	301	5/5	0.98	0.16	51,53,57,60	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.

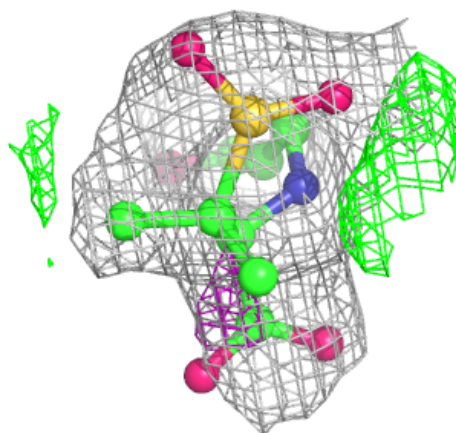
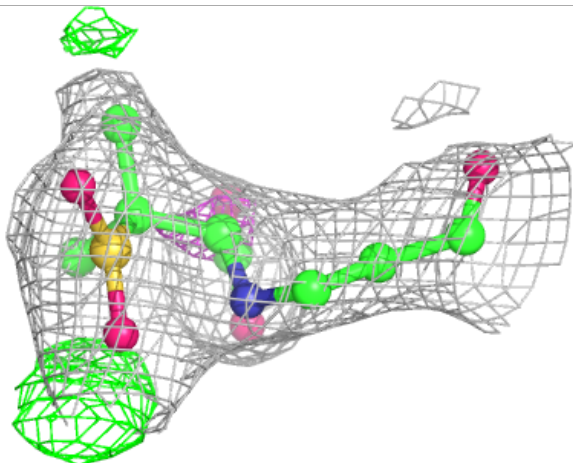
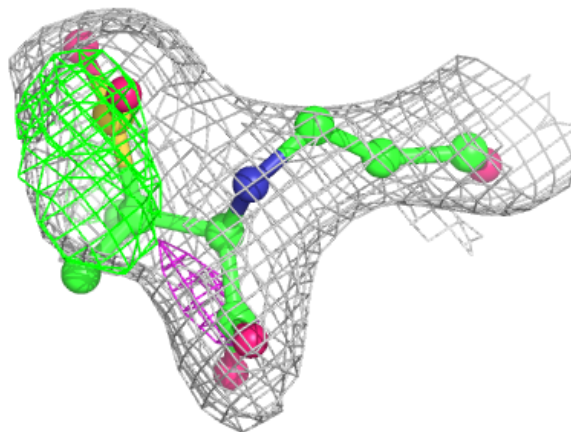
Electron density around TSL C 302:

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



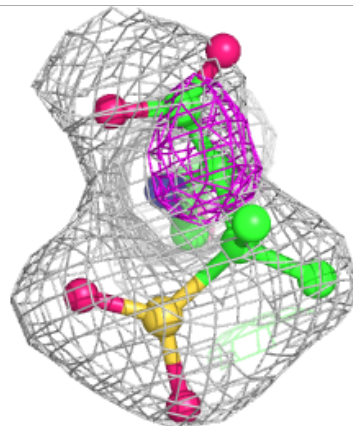
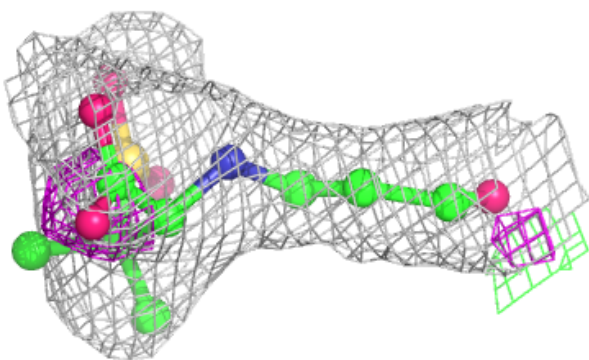
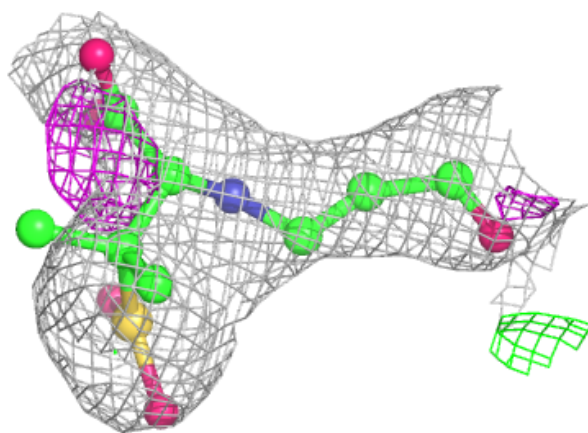
Electron density around TSL B 302:

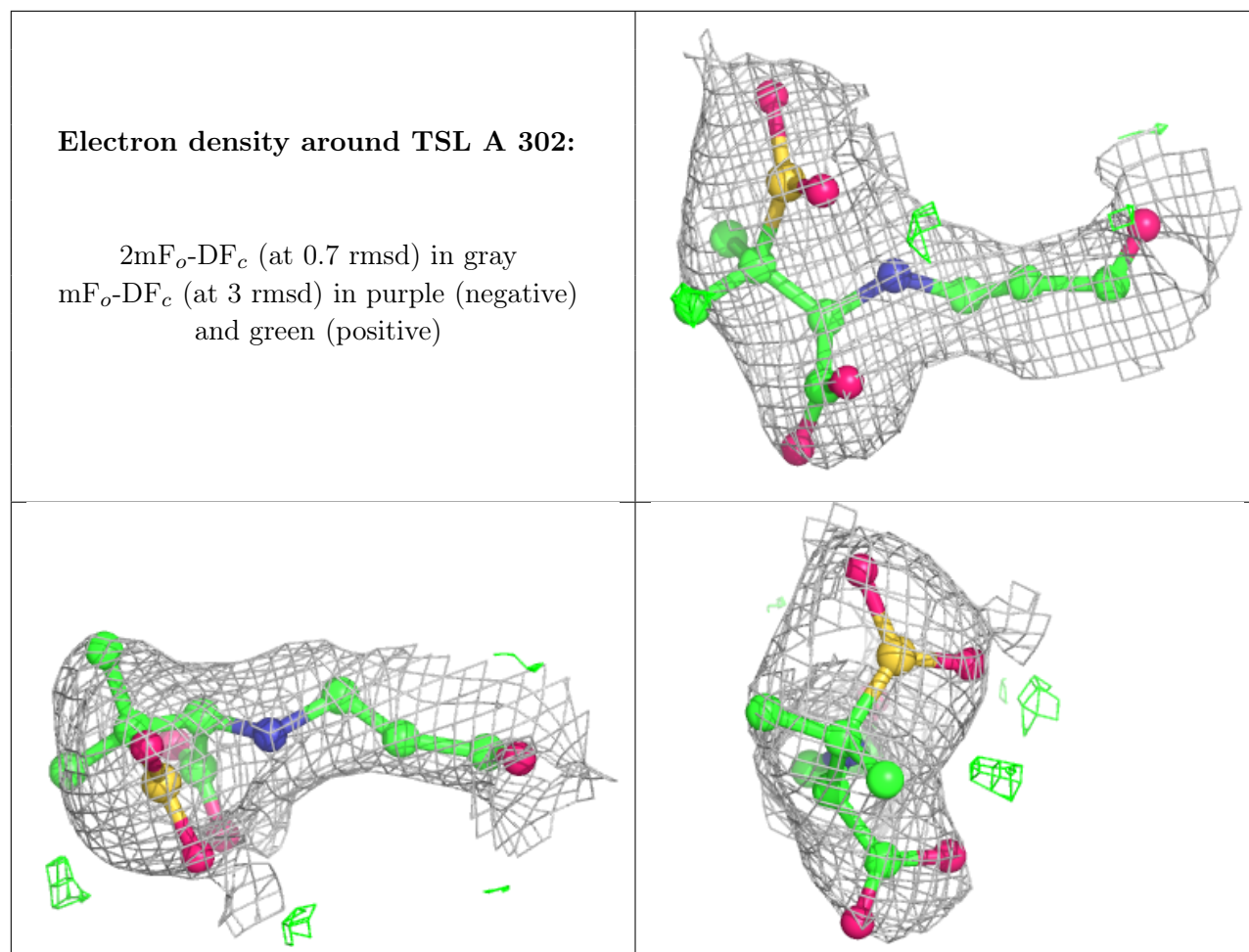
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around TSL D 302:

$2mF_o-DF_c$ (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.