



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 09:56 AM UTC

PDB ID : 6L08 / pdb_00006l08
Title : Crystal structure of Arabidopsis cytidine deaminase
Authors : Jia, W.; Xiao, W.; Lin, L.
Deposited on : 2019-09-26
Resolution : 3.00 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

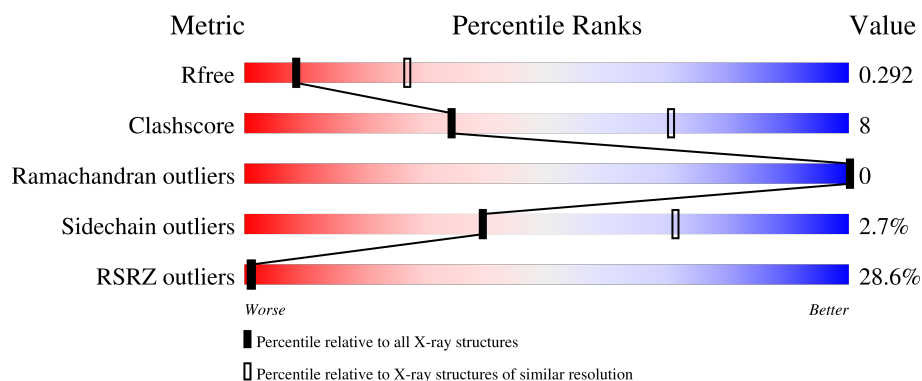
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	308	25% 74% 17% • 7%
1	B	308	28% 76% 15% • 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4183 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 Cytidine deaminase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	2	0
			2073	1310	361	395	7			
1	B	284	Total	C	N	O	S	0	1	0
			2062	1297	357	399	9			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP O65896
A	302	HIS	-	expression tag	UNP O65896
A	303	HIS	-	expression tag	UNP O65896
A	304	HIS	-	expression tag	UNP O65896
A	305	HIS	-	expression tag	UNP O65896
A	306	HIS	-	expression tag	UNP O65896
A	307	HIS	-	expression tag	UNP O65896
B	0	SER	-	expression tag	UNP O65896
B	302	HIS	-	expression tag	UNP O65896
B	303	HIS	-	expression tag	UNP O65896
B	304	HIS	-	expression tag	UNP O65896
B	305	HIS	-	expression tag	UNP O65896
B	306	HIS	-	expression tag	UNP O65896
B	307	HIS	-	expression tag	UNP O65896

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S) (labeled as "Ligand of Interest" by depositor).



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

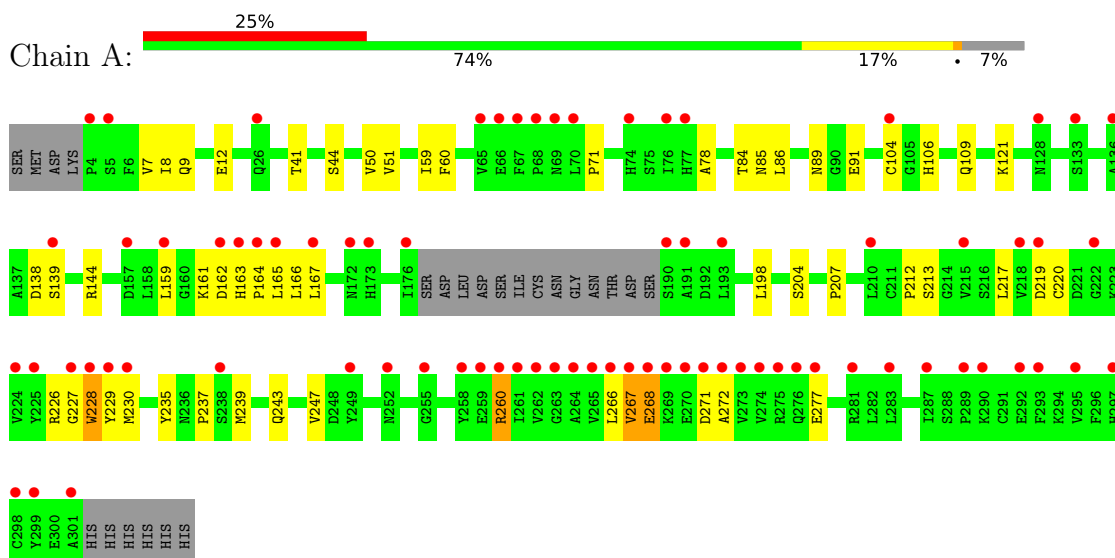
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total	O	0	0
			22	22		
3	B	21	Total	O	0	0
			21	21		

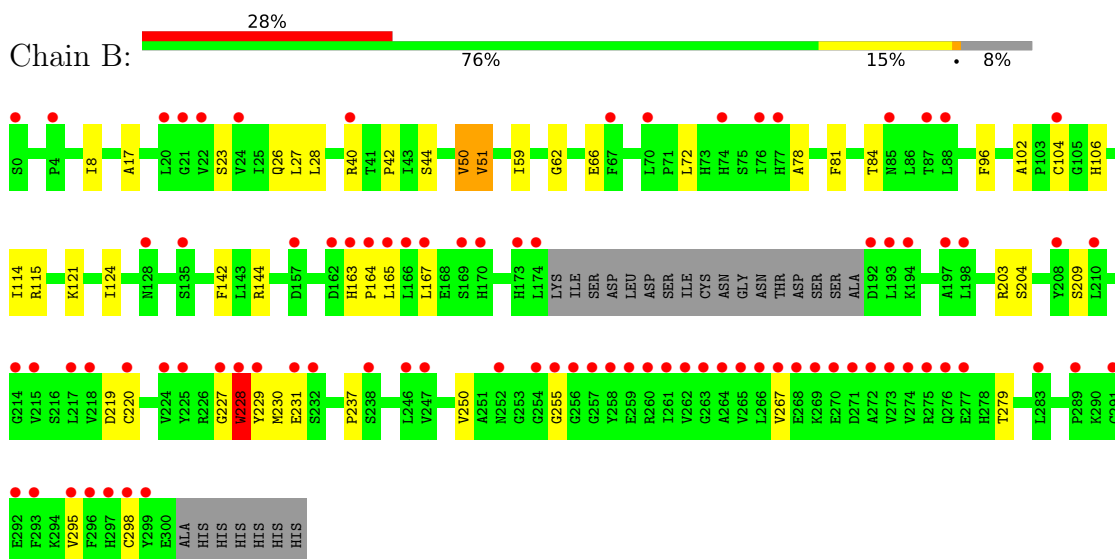
3 Residue-property plots [i](#)

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: Cytidine deaminase 1



• Molecule 1: Cytidine deaminase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	112.64Å 112.64Å 111.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.64 – 3.00 39.64 – 3.00	Depositor EDS
% Data completeness (in resolution range)	81.4 (39.64-3.00) 86.1 (39.64-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.8.1 _1168	Depositor
R, R_{free}	0.249 , 0.285 0.260 , 0.292	Depositor DCC
R_{free} test set	730 reflections (4.34%)	wwPDB-VP
Wilson B-factor (Å ²)	55.4	Xtriage
Anisotropy	0.381	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 76.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4183	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.36	0/2120	0.85	5/2896 (0.2%)
1	B	0.37	1/2109 (0.0%)	0.78	2/2881 (0.1%)
All	All	0.36	1/4229 (0.0%)	0.82	7/5777 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	228	TRP	CA-C	-7.64	1.45	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	272	ALA	N-CA-C	7.80	120.54	111.02
1	A	161	LYS	N-CA-C	7.69	122.00	111.39
1	B	227	GLY	CA-C-N	-7.24	112.61	122.16
1	B	227	GLY	C-N-CA	-7.24	112.61	122.16
1	A	271	ASP	N-CA-C	6.51	120.22	111.24
1	A	268	GLU	N-CA-C	5.85	118.09	110.43
1	A	162	ASP	N-CA-CB	-5.23	102.65	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2073	0	1948	35	0
1	B	2062	0	1930	32	0
2	B	5	0	0	0	0
3	A	22	0	0	0	0
3	B	21	0	0	0	0
All	All	4183	0	3878	61	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 (61) 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:84:THR:HG22	1:A:230:MET:HE1	1.72	0.72
1:B:23:SER:HB3	1:B:26:GLN:HG3	1.77	0.67
1:B:204:SER:HA	1:B:228:TRP:HD1	1.64	0.63
1:B:267:VAL:HG13	1:B:298:CYS:HB3	1.81	0.63
1:A:85:ASN:O	1:A:89:ASN:ND2	2.32	0.62
1:A:226:ARG:HD3	1:A:228:TRP:CH2	2.35	0.61
1:B:84:THR:HG22	1:B:230:MET:HE1	1.83	0.60
1:A:84:THR:HG21	1:A:239:MET:HE2	1.83	0.59
1:B:209:SER:OG	1:B:231:GLU:OE1	2.22	0.58
1:B:96:PHE:HE1	1:B:121:LYS:HD2	1.69	0.58
1:B:121:LYS:HB3	1:B:142:PHE:HB3	1.86	0.57
1:B:40:ARG:HH22	1:B:255:GLY:HA2	1.70	0.57
1:A:213:SER:HB2	1:A:229:TYR:HB2	1.88	0.56
1:B:42:PRO:HD2	1:B:66:GLU:HB2	1.88	0.56
1:A:164:PRO:HD2	1:A:167:LEU:HD22	1.86	0.56
1:A:237:PRO:HG3	1:B:106:HIS:HB2	1.88	0.55
1:A:267:VAL:HG12	1:A:268:GLU:H	1.72	0.55
1:B:204:SER:HA	1:B:228:TRP:CD1	2.41	0.55
1:B:164:PRO:HD2	1:B:167:LEU:HD22	1.89	0.54
1:B:203:ARG:O	1:B:228:TRP:CD1	2.61	0.54
1:B:8:ILE:HB	1:B:59:ILE:HB	1.89	0.53
1:A:159:LEU:HD21	1:A:207:PRO:HB2	1.90	0.53
1:A:227:GLY:C	1:A:228:TRP:HE3	2.18	0.52
1:A:267:VAL:HG21	1:A:277:GLU:HG2	1.91	0.52
1:A:138:ASP:OD1	1:A:139:SER:N	2.40	0.52
1:B:115:ARG:HH21	1:B:164:PRO:HD3	1.75	0.52
1:A:8:ILE:HB	1:A:59:ILE:HB	1.92	0.51
1:A:71:PRO:HG3	1:B:279:THR:HG21	1.93	0.51
1:B:219:ASP:OD1	1:B:220:CYS:N	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:HIS:HB2	1:B:237:PRO:HG3	1.95	0.49
1:B:62:GLY:HA2	1:B:81:PHE:HD2	1.78	0.48
1:A:50:VAL:HG12	1:A:78:ALA:HB1	1.95	0.48
1:B:114:ILE:HG12	1:B:165:LEU:HD21	1.96	0.47
1:A:165:LEU:HD11	1:A:230:MET:HB3	1.97	0.46
1:B:50:VAL:HG12	1:B:78:ALA:HB1	1.97	0.46
1:A:198:LEU:HA	1:A:266:LEU:HD22	1.97	0.46
1:A:219:ASP:OD1	1:A:260:ARG:NH1	2.48	0.46
1:A:228:TRP:N	1:A:228:TRP:CE3	2.83	0.46
1:A:41:THR:OG1	1:A:44:SER:OG	2.33	0.45
1:A:9:GLN:HB2	1:A:12:GLU:HG2	1.98	0.45
1:B:104:CYS:SG	1:B:106:HIS:HB3	2.57	0.45
1:B:163:HIS:HA	1:B:164:PRO:HD3	1.88	0.45
1:A:228:TRP:HE3	1:A:228:TRP:N	2.15	0.44
1:A:86:LEU:HB2	1:A:91:GLU:HB3	2.00	0.44
1:B:28:LEU:HD22	1:B:51:VAL:HG22	1.98	0.44
1:B:228:TRP:HB2	1:B:229:TYR:H	1.57	0.44
1:A:235:TYR:HB3	1:B:72:LEU:HB3	2.01	0.43
1:A:121:LYS:HA	1:A:144:ARG:HA	1.99	0.43
1:B:62:GLY:HA2	1:B:81:PHE:CD2	2.53	0.43
1:A:219:ASP:OD1	1:A:220:CYS:N	2.51	0.43
1:B:121:LYS:HA	1:B:144:ARG:HA	2.01	0.42
1:A:163:HIS:HA	1:A:164:PRO:HD3	1.83	0.42
1:A:260:ARG:HD3	1:A:260:ARG:HA	1.82	0.42
1:B:102:ALA:HA	1:B:124:ILE:HD13	2.02	0.42
1:A:204:SER:HB3	1:A:212:PRO:HB3	2.01	0.42
1:B:17:ALA:HB2	1:B:27:LEU:HD22	2.02	0.42
1:A:109:GLN:HG2	1:B:104:CYS:HB2	2.02	0.41
1:A:104:CYS:SG	1:A:106:HIS:HB3	2.61	0.41
1:A:235:TYR:OH	1:B:44:SER:HB3	2.21	0.41
1:A:243:GLN:O	1:A:247:VAL:HG13	2.21	0.41
1:A:7:VAL:HG12	1:A:60:PHE:CD1	2.56	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	283/308 (92%)	272 (96%)	11 (4%)	0	100	100
1	B	281/308 (91%)	264 (94%)	17 (6%)	0	100	100
All	All	564/616 (92%)	536 (95%)	28 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	207/257 (80%)	201 (97%)	6 (3%)	37	70
1	B	210/257 (82%)	205 (98%)	5 (2%)	43	73
All	All	417/514 (81%)	406 (97%)	11 (3%)	39	72

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	VAL
1	A	166	LEU
1	A	217	LEU
1	A	228	TRP
1	A	260	ARG
1	A	267	VAL
1	B	50	VAL
1	B	51	VAL
1	B	228	TRP
1	B	250	VAL
1	B	295	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	HIS
1	B	73	HIS
1	B	77	HIS
1	B	80	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	SO4	B	400	-	4,4,4	0.24	0	6,6,6	0.07	0

There are no bond length outliers.

There are no bond angle outliers.

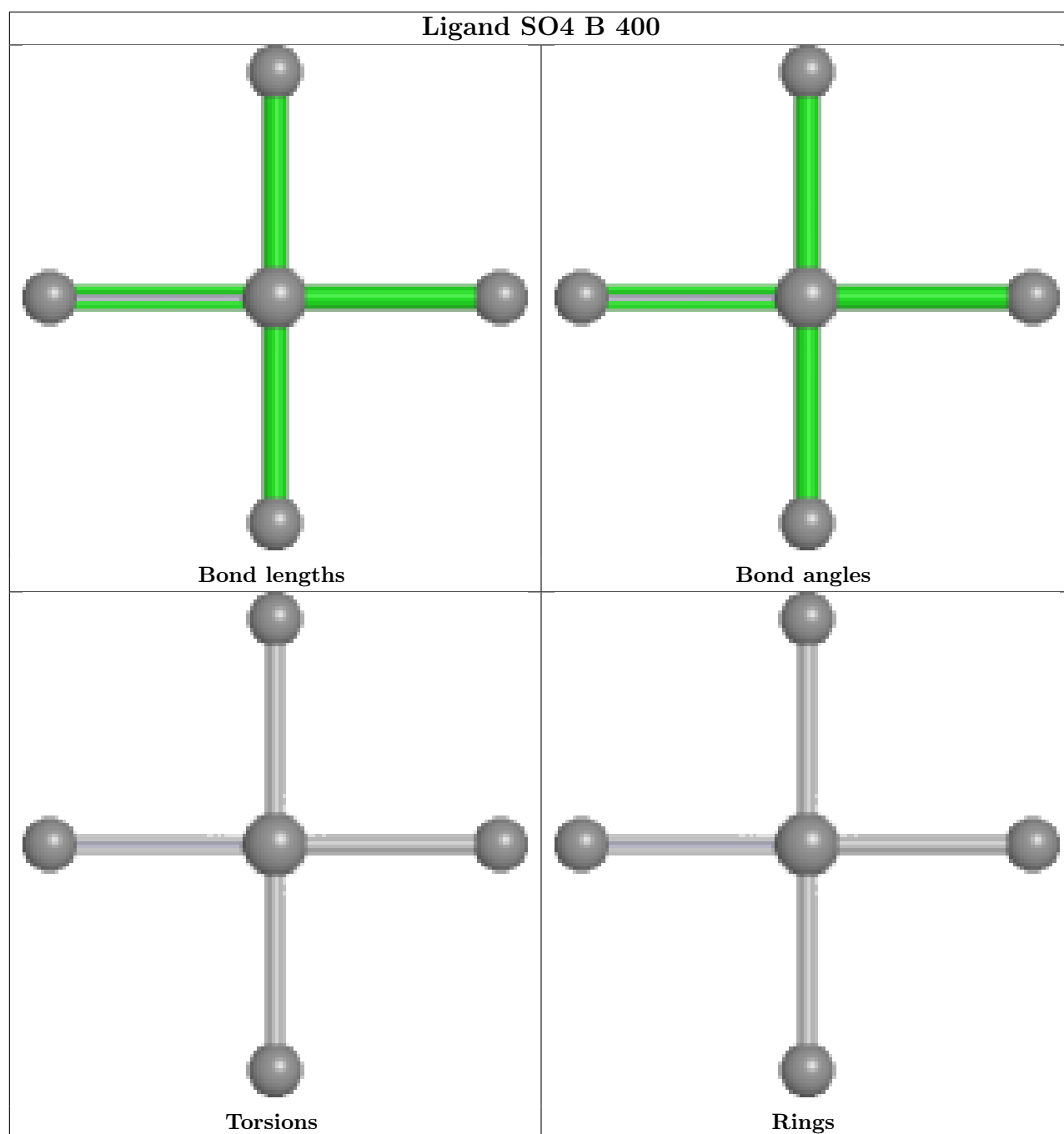
There are no chirality outliers.

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	285/308 (92%)	1.43	77 (27%)  	28, 87, 156, 268	2 (0%)
1	B	284/308 (92%)	1.46	86 (30%)  	38, 102, 180, 210	1 (0%)
All	All	569/616 (92%)	1.45	163 (28%)  	28, 94, 176, 268	3 (0%)

All (163) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	263	GLY	7.4
1	B	266	LEU	7.3
1	B	265	VAL	7.1
1	A	271	ASP	6.9
1	A	264	ALA	6.9
1	B	254	GLY	6.0
1	A	301	ALA	5.7
1	B	257	GLY	5.4
1	B	291	CYS	5.1
1	A	272	ALA	5.1
1	A	267	VAL	5.1
1	A	69	ASN	4.6
1	A	228	TRP	4.5
1	B	174	LEU	4.4
1	A	261	ILE	4.4
1	B	164	PRO	4.3
1	B	21	GLY	4.3
1	A	273	VAL	4.2
1	B	215	VAL	4.2
1	A	163	HIS	4.1
1	B	194	LYS	4.1
1	A	262	VAL	4.0
1	B	267	VAL	4.0
1	A	299	TYR	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	270	GLU	3.9
1	A	176	ILE	3.9
1	A	295	VAL	3.8
1	A	276	GLN	3.8
1	A	164	PRO	3.8
1	B	163	HIS	3.8
1	A	136	ALA	3.8
1	A	5	SER	3.7
1	A	225	TYR	3.7
1	A	238	SER	3.7
1	B	167	LEU	3.7
1	A	258	TYR	3.6
1	B	193	LEU	3.6
1	A	268	GLU	3.6
1	B	256	GLY	3.5
1	B	260	ARG	3.5
1	B	268	GLU	3.5
1	B	165	LEU	3.4
1	A	260	ARG	3.4
1	A	218	VAL	3.3
1	B	20	LEU	3.3
1	A	76	ILE	3.3
1	A	173	HIS	3.3
1	B	214	GLY	3.3
1	A	277	GLU	3.2
1	A	274	VAL	3.2
1	A	298	CYS	3.2
1	B	255	GLY	3.2
1	B	298	CYS	3.2
1	A	266	LEU	3.1
1	B	162	ASP	3.1
1	B	261	ILE	3.1
1	B	87	THR	3.1
1	A	77	HIS	3.1
1	A	190	SER	3.1
1	B	252	ASN	3.1
1	B	264	ALA	3.0
1	A	255	GLY	3.0
1	A	165	LEU	3.0
1	B	231	GLU	3.0
1	B	74	HIS	3.0
1	B	170	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	295	VAL	3.0
1	B	263	GLY	2.9
1	B	197	ALA	2.9
1	B	272	ALA	2.9
1	B	273	VAL	2.9
1	A	159	LEU	2.9
1	A	68	PRO	2.9
1	A	265	VAL	2.9
1	A	269	LYS	2.9
1	B	225	TYR	2.9
1	A	281[A]	ARG	2.9
1	B	198	LEU	2.8
1	B	210	LEU	2.8
1	A	283	LEU	2.8
1	B	275	ARG	2.8
1	A	133	SER	2.8
1	A	4	PRO	2.8
1	A	224	VAL	2.8
1	B	135	SER	2.8
1	A	219	ASP	2.7
1	A	259	GLU	2.7
1	A	252	ASN	2.7
1	A	66	GLU	2.7
1	B	169[A]	SER	2.7
1	B	262	VAL	2.7
1	B	166	LEU	2.7
1	A	67	PHE	2.7
1	B	22	VAL	2.7
1	A	230	MET	2.6
1	B	269	LYS	2.6
1	B	276	GLN	2.6
1	B	104	CYS	2.6
1	A	227	GLY	2.6
1	B	274	VAL	2.6
1	B	0	SER	2.6
1	B	229	TYR	2.6
1	A	292	GLU	2.6
1	B	277	GLU	2.6
1	B	85	ASN	2.6
1	A	275	ARG	2.5
1	A	297	HIS	2.5
1	A	289	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	70	LEU	2.5
1	B	88	LEU	2.5
1	A	104	CYS	2.5
1	A	162	ASP	2.5
1	B	297	HIS	2.5
1	A	290	LYS	2.5
1	A	26	GLN	2.5
1	B	224	VAL	2.5
1	A	293	PHE	2.5
1	B	289	PRO	2.4
1	B	67	PHE	2.4
1	B	293	PHE	2.4
1	B	246	LEU	2.4
1	B	270	GLU	2.4
1	A	128	ASN	2.4
1	A	172	ASN	2.4
1	B	238	SER	2.4
1	B	283	LEU	2.4
1	A	222	GLY	2.4
1	A	191	ALA	2.4
1	A	193	LEU	2.4
1	A	210	LEU	2.4
1	B	173	HIS	2.4
1	A	215	VAL	2.3
1	B	192	ASP	2.3
1	B	292	GLU	2.3
1	B	232	SER	2.3
1	B	247	VAL	2.3
1	B	228	TRP	2.3
1	B	258	TYR	2.3
1	B	24	VAL	2.3
1	B	296	PHE	2.2
1	B	128	ASN	2.2
1	A	157	ASP	2.2
1	B	40	ARG	2.2
1	B	218	VAL	2.2
1	A	139	SER	2.2
1	A	65	VAL	2.2
1	B	220	CYS	2.2
1	B	299	TYR	2.2
1	A	287	ILE	2.1
1	B	4	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	167	LEU	2.1
1	B	70	LEU	2.1
1	B	227	GLY	2.1
1	B	271	ASP	2.1
1	B	76	ILE	2.1
1	B	259	GLU	2.1
1	A	249	TYR	2.1
1	A	74	HIS	2.0
1	B	217	LEU	2.0
1	B	77	HIS	2.0
1	B	157	ASP	2.0
1	A	229	TYR	2.0
1	B	208	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

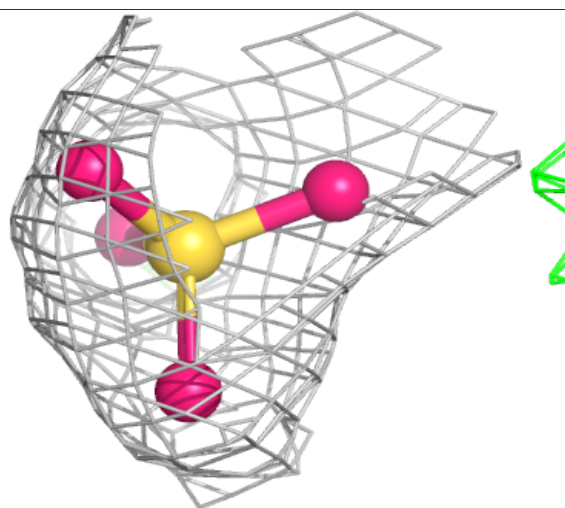
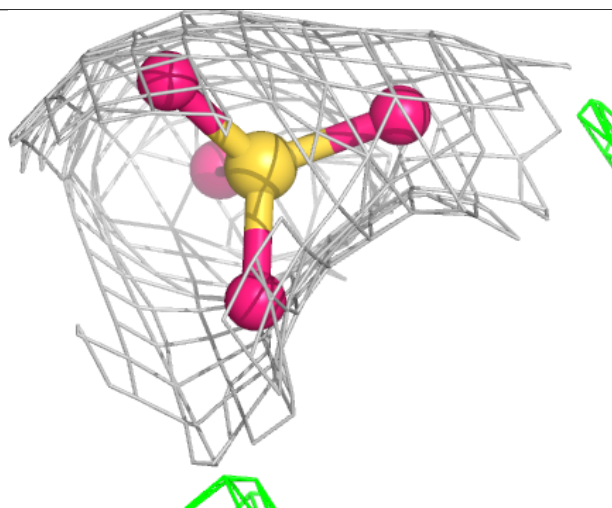
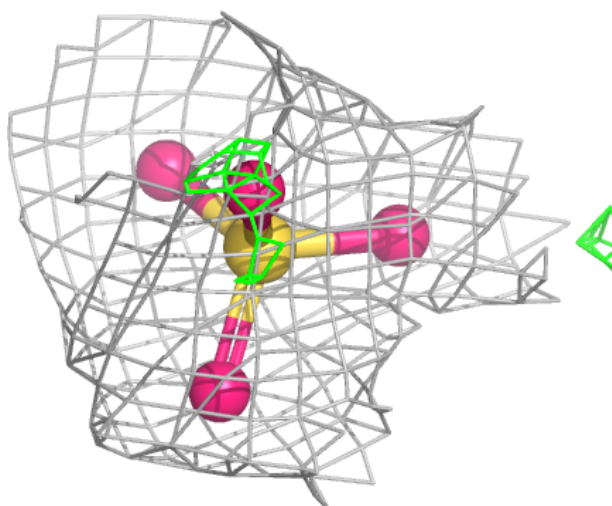
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	B	400	5/5	0.62	0.17	90,92,93,93	5

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 SO4 B 400:

$2mF_o-DF_c$ (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.