



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 06:16 pm BST

PDB ID : 3B1T
Title : Crystal structure of human peptidylarginine deiminase 4 in complex with o-Cl-amidine
Authors : Causey, C.P.; Jones, J.E.; Slack, J.L.; Kamei, D.; Jones Jr, L.E.; Subramanian, V.; Knuckley, B.; Ebrahimi, P.; Chumanevich, A.A.; Luo, Y.; Hashimoto, H.; Shimizu, T.; Sato, M.; Hofseth, L.J.; Thompson, P.R.
Deposited on : 2011-07-13
Resolution : 2.50 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 : 2.11
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) : 2.11

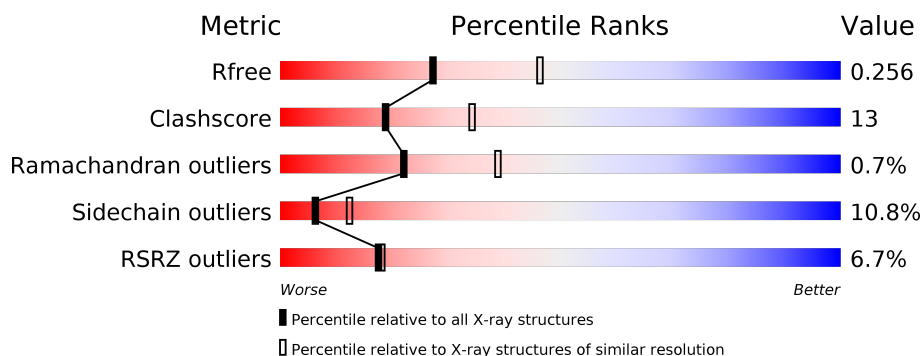
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	671	<div> <div>6%</div> <div>67%</div> <div>21%</div> <div>•</div> <div>7%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5109 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.

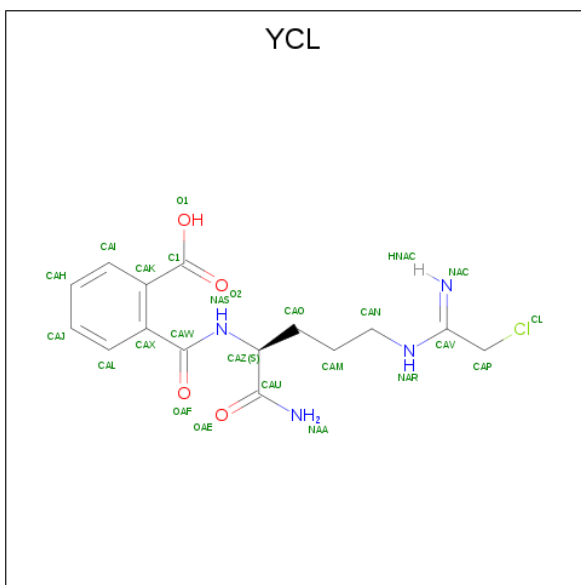
- Molecule 1 is a protein called Protein-arginine deiminase type-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	624	4912	3135	823	919	35	0	0	0

There are 11 discrepancies between the modelled and reference sequences:

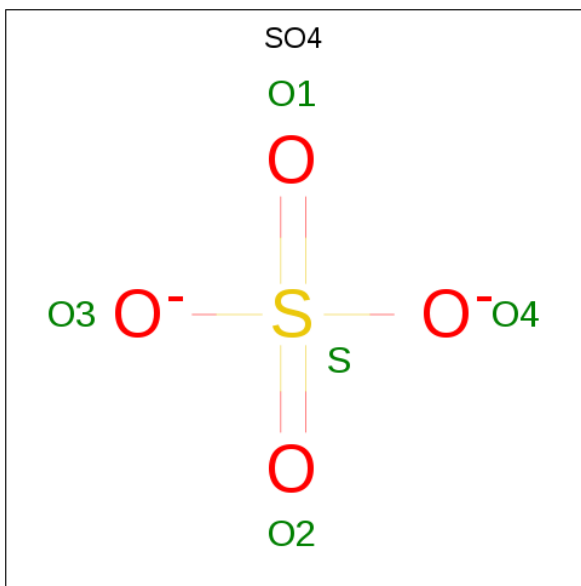
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	EXPRESSION TAG	UNP Q9UM07
A	-6	PRO	-	EXPRESSION TAG	UNP Q9UM07
A	-5	LEU	-	EXPRESSION TAG	UNP Q9UM07
A	-4	GLY	-	EXPRESSION TAG	UNP Q9UM07
A	-3	SER	-	EXPRESSION TAG	UNP Q9UM07
A	-2	PRO	-	EXPRESSION TAG	UNP Q9UM07
A	-1	GLU	-	EXPRESSION TAG	UNP Q9UM07
A	0	PHE	-	EXPRESSION TAG	UNP Q9UM07
A	55	SER	GLY	SEE REMARK 999	UNP Q9UM07
A	82	ALA	VAL	SEE REMARK 999	UNP Q9UM07
A	112	ALA	GLY	SEE REMARK 999	UNP Q9UM07

- Molecule 2 is 2-{{[(2S)-1-amino-5-{{[(1Z)-2-chloroethanimidoyl]amino}-1-oxopentan-2-yl]carbonyl]benzoic acid (three-letter code: YCL) (formula: C₁₅H₁₉ClN₄O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			23	15	4	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total 5	Ca 5	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	154	Total 154	O 154	0	0

- Molecule 1: Protein-arginine deiminase type-4



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.79Å 60.85Å 114.30Å 90.00° 123.99° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 19.96 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.50) 94.8 (19.96-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 2.50Å)	Xtriage
Refinement program	REFMAC 5.6.0085	Depositor
R, R_{free}	0.211 , 0.263 0.210 , 0.256	Depositor DCC
R_{free} test set	1367 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.232	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5109	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: CA, SO4, YCL

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	0.65	0/5030	0.74	2/6826 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	412	VAL	CB-CA-C	-6.66	98.74	111.40
1	A	427	ARG	NE-CZ-NH2	-5.87	117.36	120.30

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	4912	0	4861	129	0
2	A	23	0	15	2	0
3	A	15	0	0	0	0
4	A	5	0	0	0	0
5	A	154	0	0	5	0
All	All	5109	0	4876	129	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 13.

All (129) 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:334:LEU:HD13	1:A:335:THR:N	1.52	1.24
1:A:334:LEU:HD13	1:A:334:LEU:C	1.75	1.05
1:A:367:VAL:HG21	1:A:384:VAL:HG11	1.41	1.01
1:A:137:ARG:HD2	1:A:288:SER:HB2	1.43	1.00
1:A:50:VAL:HG22	1:A:78:LEU:HD12	1.51	0.91
1:A:334:LEU:CD1	1:A:334:LEU:C	2.37	0.91
1:A:367:VAL:HG21	1:A:384:VAL:CG1	2.06	0.85
1:A:313:ILE:HG23	1:A:316:ASN:HB3	1.59	0.83
1:A:188:LYS:CE	1:A:244:LYS:HE2	2.08	0.83
1:A:41:SER:HG	1:A:67:THR:N	1.77	0.82
1:A:317:GLU:O	1:A:321:LYS:HG3	1.83	0.79
1:A:38:THR:OG1	1:A:39:SER:N	2.14	0.78
1:A:47:GLY:HA3	1:A:90:GLN:NE2	2.00	0.77
1:A:67:THR:N	1:A:97:TYR:HH	1.83	0.77
1:A:203:VAL:HG22	1:A:229:LEU:HD13	1.68	0.75
1:A:188:LYS:NZ	1:A:244:LYS:HE2	2.02	0.74
1:A:47:GLY:CA	1:A:90:GLN:NE2	2.51	0.73
1:A:313:ILE:HG23	1:A:316:ASN:CB	2.18	0.72
1:A:166:CYS:HB2	1:A:254:LEU:HD22	1.71	0.72
1:A:317:GLU:O	1:A:321:LYS:CG	2.38	0.71
1:A:203:VAL:CG2	1:A:229:LEU:HD13	2.20	0.71
1:A:354:ILE:HD12	1:A:367:VAL:HG22	1.72	0.70
1:A:47:GLY:H	1:A:90:GLN:HE22	1.38	0.69
1:A:484:ARG:NH1	1:A:563:ASP:OD1	2.27	0.68
1:A:47:GLY:N	1:A:90:GLN:HE22	1.91	0.68
1:A:50:VAL:HG22	1:A:78:LEU:CD1	2.23	0.68
1:A:521:LYS:HG3	1:A:522:LYS:N	2.10	0.67
1:A:272:LEU:CD2	1:A:283:VAL:HG22	2.25	0.66
1:A:47:GLY:CA	1:A:90:GLN:HE22	2.09	0.66
1:A:29:ILE:HD11	1:A:42:ILE:HD11	1.79	0.65
1:A:367:VAL:CG2	1:A:384:VAL:HG11	2.22	0.64
1:A:123:ARG:HD3	1:A:659:TRP:CD1	2.33	0.63
1:A:188:LYS:HZ2	1:A:244:LYS:HE2	1.61	0.63
1:A:313:ILE:HG12	1:A:314:PHE:N	2.13	0.63
1:A:313:ILE:HG22	1:A:316:ASN:ND2	2.14	0.62
1:A:334:LEU:HD11	1:A:336:ILE:HG13	1.83	0.61
1:A:34:PRO:HG2	1:A:104:VAL:HG21	1.82	0.61
1:A:139:TRP:CD1	1:A:147:GLY:HA3	2.35	0.61
1:A:483:ASP:OD1	1:A:483:ASP:N	2.34	0.60
1:A:151:LEU:HD21	1:A:359:ALA:HB2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:ILE:CG2	1:A:316:ASN:CG	2.71	0.59
1:A:212:ARG:HD3	1:A:227:VAL:HG11	1.85	0.58
1:A:334:LEU:HD13	1:A:335:THR:CA	2.31	0.58
1:A:21:LEU:HD23	1:A:114:GLU:HG2	1.86	0.56
1:A:354:ILE:CD1	1:A:367:VAL:HG22	2.35	0.56
1:A:276:ASN:CB	1:A:279:LEU:HD12	2.35	0.56
1:A:313:ILE:HG23	1:A:316:ASN:CG	2.27	0.55
1:A:318:ASP:HA	1:A:321:LYS:HG3	1.89	0.55
1:A:188:LYS:HE3	1:A:244:LYS:HE2	1.85	0.55
1:A:10:THR:HG22	1:A:32:SER:CB	2.37	0.54
1:A:29:ILE:HG23	1:A:96:TYR:HE1	1.73	0.54
1:A:432:ASP:OD2	1:A:433:SER:N	2.38	0.54
1:A:212:ARG:HD3	1:A:227:VAL:CG1	2.39	0.53
1:A:425:LEU:HD12	1:A:456:VAL:HG13	1.91	0.53
1:A:312:SER:HA	1:A:320:LEU:HD11	1.91	0.52
1:A:5:THR:CG2	5:A:768:HOH:O	2.57	0.52
1:A:108:LEU:HG	1:A:110:LEU:HD13	1.90	0.52
1:A:152:VAL:HG12	1:A:154:CYS:HB3	1.91	0.52
1:A:71:ASP:O	1:A:74:VAL:HG12	2.09	0.52
1:A:313:ILE:HG22	1:A:316:ASN:CG	2.31	0.51
1:A:497:CYS:HB3	1:A:570:LEU:HD13	1.93	0.51
1:A:101:THR:HG22	1:A:102:PRO:O	2.12	0.50
1:A:347:TRP:CE2	2:A:800:YCL:HAO	2.46	0.50
1:A:159:LEU:O	1:A:159:LEU:HD13	2.12	0.50
1:A:45:SER:HB2	1:A:46:PRO:CD	2.41	0.49
1:A:313:ILE:HG12	1:A:314:PHE:H	1.77	0.49
1:A:427:ARG:HD2	1:A:458:ALA:O	2.13	0.49
1:A:203:VAL:HG22	1:A:229:LEU:CD1	2.40	0.49
1:A:51:ASP:HB2	1:A:77:THR:HG23	1.94	0.49
1:A:188:LYS:NZ	1:A:244:LYS:CE	2.74	0.49
1:A:317:GLU:O	1:A:321:LYS:HG2	2.12	0.49
1:A:621:GLU:N	1:A:622:PRO:CD	2.77	0.48
1:A:171:VAL:HG23	1:A:214:PHE:CE2	2.48	0.48
1:A:5:THR:HG23	5:A:768:HOH:O	2.14	0.47
1:A:150:LEU:HD21	1:A:251:VAL:HG12	1.95	0.47
1:A:594:LYS:NZ	5:A:741:HOH:O	2.47	0.47
1:A:467:LEU:HD13	1:A:474:GLU:HB2	1.96	0.47
1:A:228:VAL:O	1:A:229:LEU:HD23	2.15	0.46
1:A:10:THR:HG22	1:A:32:SER:HB3	1.97	0.46
1:A:195:THR:HG22	1:A:196:ASN:ND2	2.31	0.46
1:A:347:TRP:CD2	2:A:800:YCL:HAO	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:SER:OG	1:A:338:PRO:HA	2.15	0.46
1:A:52:ILE:HG21	1:A:68:TRP:CH2	2.51	0.46
1:A:200:VAL:HG23	1:A:236:HIS:O	2.17	0.45
1:A:367:VAL:HG23	1:A:389:PHE:CE2	2.52	0.45
1:A:21:LEU:HD23	1:A:114:GLU:CG	2.46	0.45
1:A:333:LYS:HG2	1:A:334:LEU:N	2.32	0.45
1:A:47:GLY:HA3	1:A:90:GLN:HE22	1.73	0.45
1:A:188:LYS:HE3	1:A:244:LYS:CE	2.45	0.45
1:A:345:ASP:HB2	1:A:375:GLY:HA3	1.99	0.45
1:A:313:ILE:O	1:A:314:PHE:O	2.35	0.45
1:A:396:PRO:HG2	1:A:401:ILE:CD1	2.47	0.45
1:A:367:VAL:CG2	1:A:384:VAL:CG1	2.87	0.45
1:A:513:LEU:HD22	1:A:523:GLN:HB3	1.99	0.44
1:A:333:LYS:CG	1:A:334:LEU:N	2.80	0.44
1:A:13:GLN:HG2	1:A:14:PRO:HD2	1.98	0.44
1:A:276:ASN:HB2	1:A:279:LEU:HD12	1.98	0.44
1:A:108:LEU:HG	1:A:110:LEU:CD1	2.47	0.44
1:A:42:ILE:O	1:A:42:ILE:HG22	2.17	0.44
1:A:473:ASP:HB2	5:A:694:HOH:O	2.17	0.44
1:A:150:LEU:HD21	1:A:251:VAL:CG1	2.48	0.43
1:A:313:ILE:CG1	1:A:314:PHE:N	2.80	0.43
1:A:416:VAL:HG22	5:A:739:HOH:O	2.17	0.43
1:A:45:SER:CB	1:A:46:PRO:CD	2.96	0.43
1:A:493:SER:CB	1:A:566:ASP:HB3	2.48	0.43
1:A:171:VAL:HG23	1:A:214:PHE:CD2	2.53	0.43
1:A:96:TYR:HB2	1:A:104:VAL:CG2	2.49	0.43
1:A:34:PRO:CG	1:A:104:VAL:HG21	2.47	0.43
1:A:156:ARG:HB2	1:A:385:MET:HE2	2.00	0.43
1:A:198:THR:HG23	1:A:274:THR:HG21	2.01	0.42
1:A:203:VAL:HG23	1:A:229:LEU:HD13	1.99	0.42
1:A:188:LYS:HZ2	1:A:244:LYS:CE	2.32	0.42
1:A:631:ASN:C	1:A:631:ASN:OD1	2.57	0.42
1:A:74:VAL:HG13	1:A:74:VAL:O	2.20	0.42
1:A:457:GLN:O	1:A:458:ALA:C	2.57	0.42
1:A:583:PHE:O	1:A:584:PRO:C	2.57	0.42
1:A:514:LEU:HD23	1:A:603:GLY:HA2	2.02	0.41
1:A:80:MET:HE3	1:A:80:MET:HB2	1.97	0.41
1:A:273:ASP:OD2	1:A:275:SER:OG	2.27	0.41
1:A:484:ARG:HH12	1:A:563:ASP:CG	2.23	0.41
1:A:40:PHE:HD1	1:A:42:ILE:HD13	1.86	0.41
1:A:238:LEU:HD12	1:A:269:ILE:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:ILE:CG2	1:A:316:ASN:ND2	2.83	0.41
1:A:596:LEU:HD11	1:A:625:LEU:HG	2.03	0.41
1:A:114:GLU:O	1:A:187:THR:HA	2.21	0.41
1:A:623:LEU:HB2	1:A:625:LEU:HD22	2.03	0.41
1:A:137:ARG:O	1:A:290:VAL:HG21	2.21	0.40
1:A:497:CYS:SG	1:A:570:LEU:HD22	2.61	0.40
1:A:633:PHE:HA	1:A:637:HIS:HB3	2.04	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	612/671 (91%)	582 (95%)	26 (4%)	4 (1%)	22 39

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	314	PHE
1	A	585	ASN
1	A	458	ALA
1	A	45	SER

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	555/594 (93%)	495 (89%)	60 (11%)	6 12

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LEU
1	A	23	THR
1	A	24	LEU
1	A	38	THR
1	A	41	SER
1	A	42	ILE
1	A	43	ASN
1	A	67	THR
1	A	68	TRP
1	A	70	LEU
1	A	71	ASP
1	A	75	GLU
1	A	78	LEU
1	A	79	THR
1	A	80	MET
1	A	86	SER
1	A	91	LYS
1	A	95	SER
1	A	110	LEU
1	A	116	SER
1	A	117	LEU
1	A	159	LEU
1	A	169	ASP
1	A	182	LEU
1	A	185	LEU
1	A	198	THR
1	A	201	LEU
1	A	203	VAL
1	A	206	SER
1	A	259	THR
1	A	278	GLU
1	A	302	THR
1	A	312	SER
1	A	316	ASN
1	A	322	SER
1	A	328	MET
1	A	334	LEU
1	A	344	ASP

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Mol	Chain	Res	Type
1	A	364	LEU
1	A	372	ARG
1	A	374	ARG
1	A	376	LEU
1	A	389	PHE
1	A	398	THR
1	A	412	VAL
1	A	416	VAL
1	A	429	LEU
1	A	479	VAL
1	A	489	LEU
1	A	495	ARG
1	A	500	LEU
1	A	514	LEU
1	A	522	LYS
1	A	523	GLN
1	A	538	HIS
1	A	570	LEU
1	A	592	LEU
1	A	609	ARG
1	A	625	LEU
1	A	640	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	90	GLN
1	A	215	GLN
1	A	303	GLN
1	A	342	ASN
1	A	448	GLN
1	A	505	GLN
1	A	523	GLN
1	A	538	HIS

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 5 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	$\# Z > 2$
3	SO4	A	905	-	4,4,4	0.23	0	6,6,6	0.19	0
3	SO4	A	906	-	4,4,4	0.16	0	6,6,6	0.18	0
2	YCL	A	800	1	21,23,24	1.05	1 (4%)	23,30,31	1.43	2 (8%)
3	SO4	A	907	-	4,4,4	0.19	0	6,6,6	0.15	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	YCL	A	800	1	-	4/18/23/25	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	800	YCL	CAK-C1	3.87	1.51	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	YCL	CAO-CAZ-NAS	-3.34	104.12	110.88
2	A	800	YCL	CAP-CAV-NAR	2.62	117.60	114.24

There are no chirality outliers.

All (4) torsion outliers are listed below:

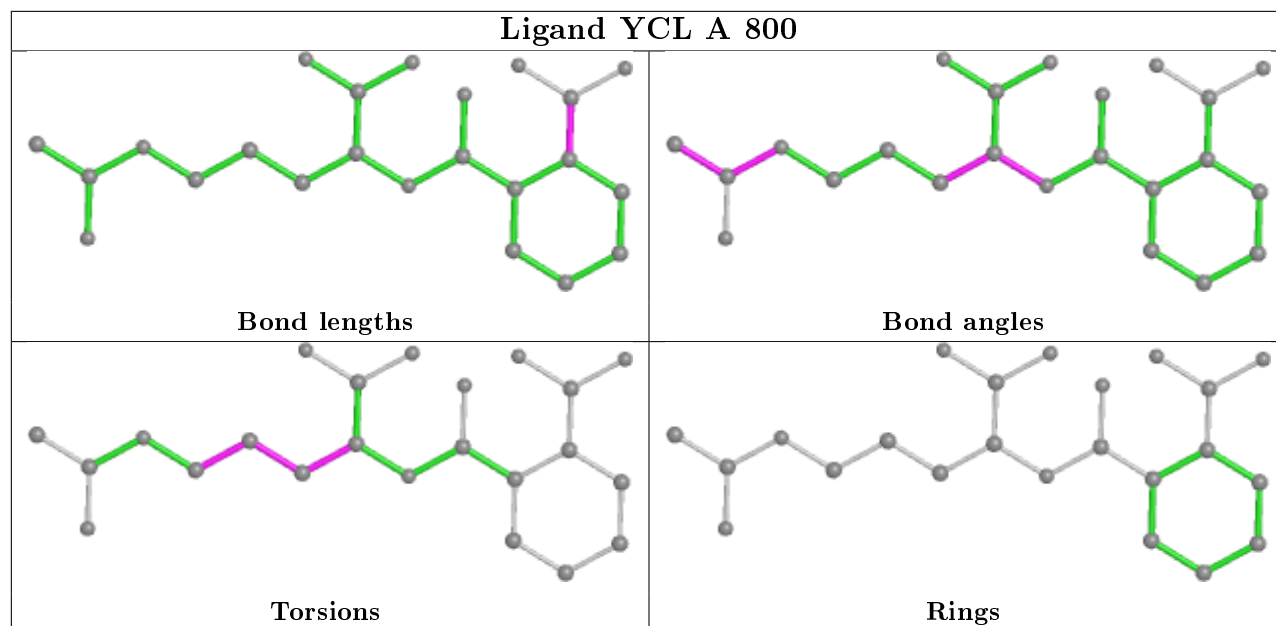
Mol	Chain	Res	Type	Atoms
2	A	800	YCL	CAM-CAO-CAZ-NAS
2	A	800	YCL	CAM-CAO-CAZ-CAU
2	A	800	YCL	CAO-CAM-CAN-NAR
2	A	800	YCL	CAN-CAM-CAO-CAZ

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	800	YCL	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 [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	624/671 (92%)	0.22	42 (6%) 17 18	42, 72, 115, 142	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	ARG	6.7
1	A	217	THR	5.9
1	A	314	PHE	5.6
1	A	100	LYS	5.5
1	A	328	MET	4.9
1	A	170	GLU	4.5
1	A	103	PRO	4.0
1	A	172	LEU	3.6
1	A	163	ALA	3.4
1	A	315	GLU	3.4
1	A	522	LYS	3.4
1	A	244	LYS	3.2
1	A	162	SER	3.2
1	A	102	PRO	3.2
1	A	213	VAL	3.2
1	A	521	LYS	2.9
1	A	161	SER	2.9
1	A	398	THR	2.8
1	A	276	ASN	2.8
1	A	136	GLN	2.8
1	A	104	VAL	2.7
1	A	340	GLU	2.7
1	A	101	THR	2.6
1	A	646	GLY	2.6
1	A	75	GLU	2.6
1	A	97	TYR	2.5
1	A	339	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	34	PRO	2.4
1	A	400	GLY	2.4
1	A	638	ILE	2.3
1	A	4	GLY	2.3
1	A	99	PRO	2.3
1	A	159	LEU	2.3
1	A	124	THR	2.2
1	A	576	PHE	2.2
1	A	634	PHE	2.2
1	A	13	GLN	2.1
1	A	438	ASN	2.1
1	A	633	PHE	2.1
1	A	24	LEU	2.1
1	A	125	GLY	2.1
1	A	639	ARG	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 carbohydrates 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.

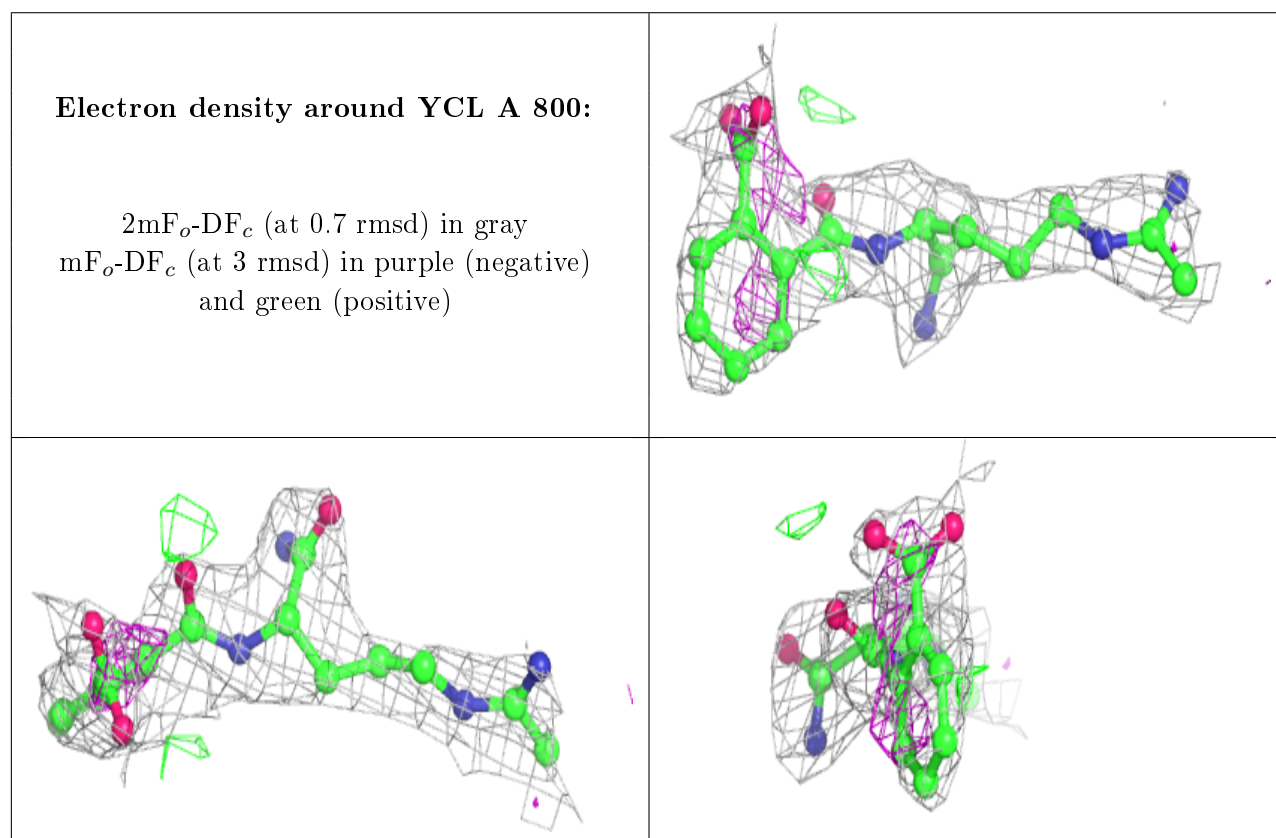
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	YCL	A	800	23/24	0.83	0.30	76,78,81,81	0
4	CA	A	904	1/1	0.88	0.11	67,67,67,67	0
3	SO4	A	907	5/5	0.90	0.31	85,85,86,86	0
3	SO4	A	906	5/5	0.91	0.40	102,102,103,103	0
4	CA	A	902	1/1	0.94	0.17	63,63,63,63	0
3	SO4	A	905	5/5	0.96	0.20	88,88,88,89	0
4	CA	A	901	1/1	0.97	0.12	56,56,56,56	0
4	CA	A	900	1/1	0.98	0.04	52,52,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CA	A	903	1/1	0.99	0.22	65,65,65,65	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.