



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 15, 2026 – 12:46 AM UTC

PDB ID : 5A3Q / pdb_00005a3q
Title : Crystal structure of the (SR) Calcium ATPase E2-vanadate complex bound to thapsigargin and TNP-AMPPCP
Authors : Clausen, J.D.; Bublitz, M.; Arnou, B.; Olesen, C.; Andersen, J.P.; Moller, J.V.; Nissen, P.
Deposited on : 2015-06-02
Resolution : 3.05 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

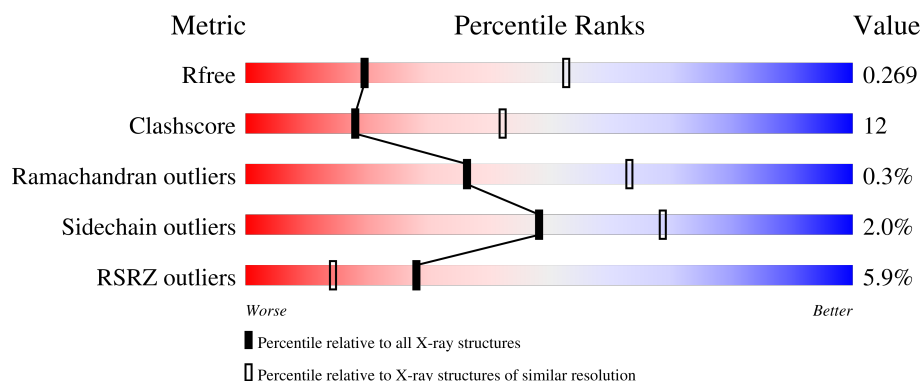
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	995	6% 73% 26% .

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7775 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.

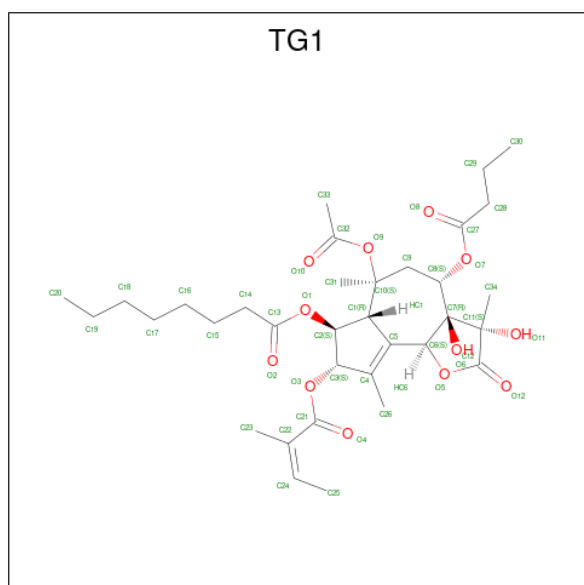
- Molecule 1 is a protein called SARCOPLASMIC/ENDOPLASMIC RETICULUM CALCIUM ATPASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	995	7667	4874	1286	1450	57	0	0	1

There is a discrepancy between the modelled and reference sequences:

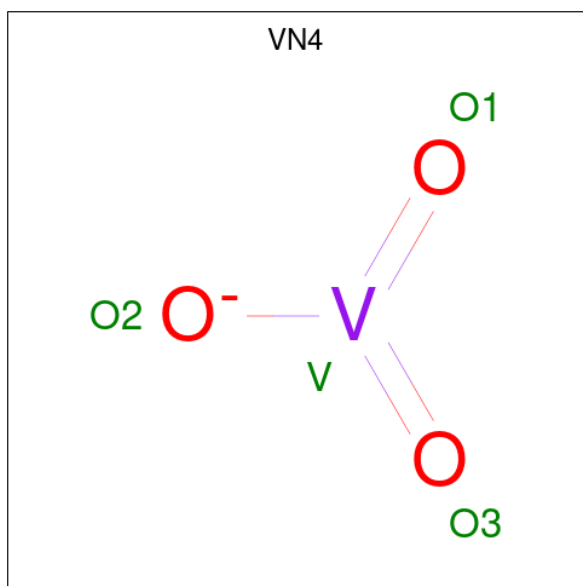
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ACE	-	acetylation	UNP P04191

- Molecule 2 is OCTANOIC ACID [3S-[3ALPHA, 3ABETA, 4ALPHA, 6BETA, 6ABETA, 7BETA, 8ALPHA(Z), 9BALPHA]]-6-(ACETYLOXY)-2,3,-3A,4,5,6,6A,7,8,9B-DECAHYDRO-3,3A-DIHYDROXY-3,6,9-TRIMETHYL-8-[(2-METHYL-1-OXO-2-BUTENYL)OXY]-2-OXO-4-(1-OXOBUTOXY)-AZULENO[4,5-B]FURAN-7-YL ESTER (CCD ID: TG1) (formula: C₃₄H₅₀O₁₂).



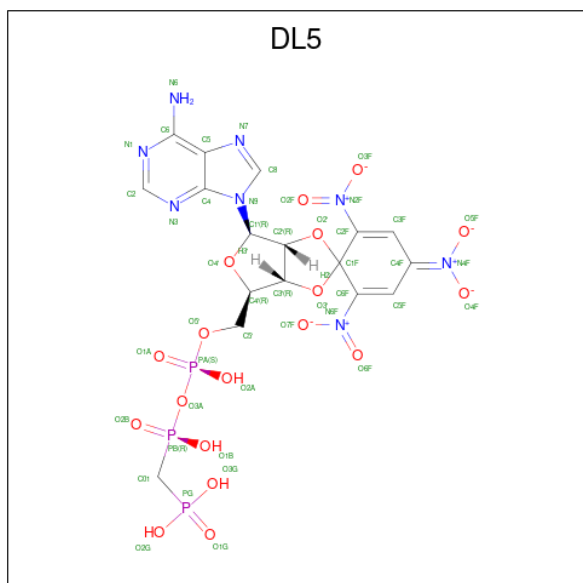
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			46	34	12		

- Molecule 3 is oxido(dioxo)vanadium (CCD ID: VN4) (formula: O_3V).



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

- Molecule 4 is Spiro(2,4,6-trinitrobenzene[1,2a]-O2',O3'-methylene-adenosine (beta,gamma-methylene)triphosphate (CCD ID: DL5) (formula: $\text{C}_{17}\text{H}_{18}\text{N}_8\text{O}_{18}\text{P}_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			46	17	8	18	3		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		

- Molecule 6 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	K	0	0
			1	1		

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Cl	0	0
			1	1		

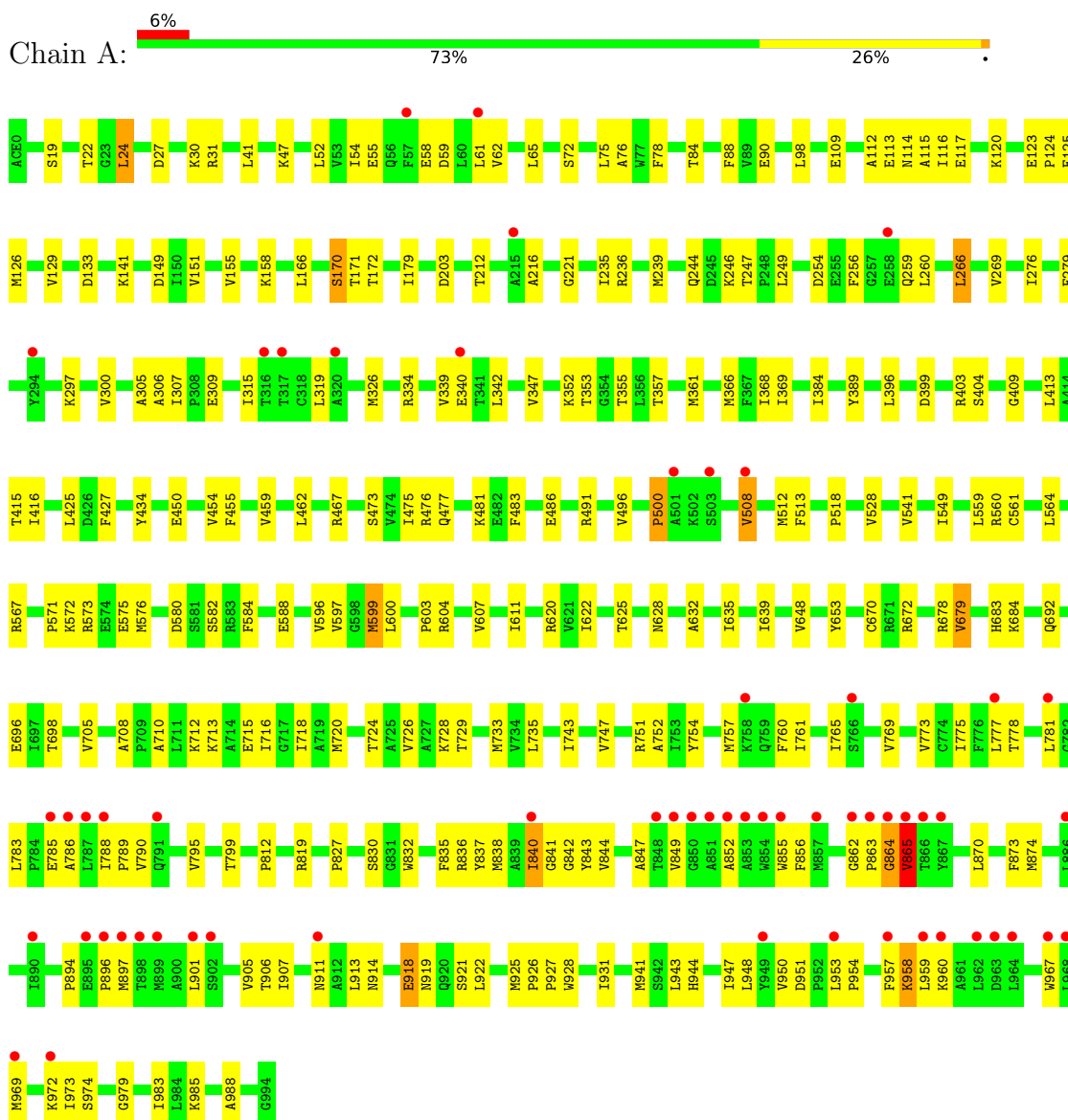
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	8	Total	O	0	0
			8	8		

3 Residue-property plots

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: SARCOPLASMIC/ENDOPLASMIC RETICULUM CALCIUM ATPASE 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	86.43Å 118.78Å 141.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	73.80 – 3.05 73.80 – 3.05	Depositor EDS
% Data completeness (in resolution range)	99.9 (73.80-3.05) 99.9 (73.80-3.05)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 3.07Å)	Xtriage
Refinement program	PHENIX (1.10_2152: ???)	Depositor
R, R_{free}	0.200 , 0.258 0.208 , 0.269	Depositor DCC
R_{free} test set	1448 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	82.7	Xtriage
Anisotropy	0.235	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 83.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7775	wwPDB-VP
Average B, all atoms (Å ²)	114.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% 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: TG1, K, DL5, ACE, CL, MG, VN4

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	0.23	0/7803	0.47	2/10575 (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	864	GLY	N-CA-C	15.21	133.52	114.66
1	A	865	VAL	N-CA-CB	10.22	128.88	111.50

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	7667	0	7754	185	0
2	A	46	0	50	11	0
3	A	4	0	0	0	0
4	A	46	0	0	3	0
5	A	2	0	0	0	0
6	A	1	0	0	0	0
7	A	1	0	0	0	0
8	A	8	0	0	0	0
All	All	7775	0	7804	190	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 12.

All (190) 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:862:GLY:C	1:A:863:PRO:CA	2.44	0.91
1:A:862:GLY:C	1:A:863:PRO:CD	2.48	0.87
1:A:894:PRO:HB3	1:A:959:LEU:H	1.41	0.84
1:A:603:PRO:HB3	1:A:639:ILE:HD11	1.70	0.74
1:A:326:MET:HE1	1:A:339:VAL:HG22	1.71	0.71
1:A:355:THR:HG23	1:A:720:MET:HE3	1.73	0.71
1:A:864:GLY:O	1:A:865:VAL:HG23	1.92	0.70
1:A:155:VAL:HG12	1:A:216:ALA:HA	1.74	0.70
1:A:239:MET:HE3	1:A:708:ALA:HB1	1.74	0.69
1:A:113:GLU:HB3	1:A:729:THR:HG23	1.73	0.69
1:A:864:GLY:O	1:A:865:VAL:CG2	2.42	0.68
1:A:396:LEU:HD13	1:A:399:ASP:HA	1.75	0.68
1:A:837:TYR:HB2	2:A:1000:TG1:H331	1.74	0.68
1:A:837:TYR:HA	1:A:840:ILE:HG22	1.75	0.68
1:A:256:PHE:CE1	2:A:1000:TG1:H231	2.28	0.68
1:A:62:VAL:HG23	1:A:98:LEU:HD22	1.74	0.68
1:A:179:ILE:H	1:A:179:ILE:HD12	1.58	0.67
1:A:773:VAL:HG11	1:A:842:GLY:HA2	1.77	0.67
1:A:827:PRO:HG2	1:A:830:SER:HB3	1.76	0.67
1:A:799:THR:HG21	1:A:905:VAL:HG22	1.76	0.67
1:A:203:ASP:OD1	1:A:678:ARG:NH2	2.21	0.66
1:A:947:ILE:HD12	1:A:953:LEU:HD13	1.78	0.66
1:A:389:TYR:HB3	1:A:425:LEU:HD11	1.81	0.63
1:A:366:MET:HG2	1:A:384:ILE:HD11	1.81	0.62
1:A:473:SER:HA	1:A:476:ARG:HD2	1.80	0.62
1:A:260:LEU:HD11	1:A:306:ALA:HB1	1.82	0.62
1:A:481:LYS:HG3	1:A:496:VAL:HG13	1.81	0.62
1:A:705:VAL:HG22	1:A:726:VAL:HG11	1.82	0.61
1:A:786:ALA:HB2	1:A:852:ALA:HB2	1.82	0.61
1:A:837:TYR:CB	2:A:1000:TG1:H331	2.29	0.61
1:A:786:ALA:HA	1:A:897:MET:HA	1.83	0.61
1:A:783:LEU:HD21	1:A:870:LEU:HG	1.81	0.61
1:A:125:GLU:HG3	1:A:126:MET:HG2	1.84	0.60
1:A:836:ARG:HH22	1:A:918:GLU:HA	1.66	0.60
1:A:256:PHE:HE1	2:A:1000:TG1:H231	1.67	0.59
1:A:235:ILE:O	1:A:239:MET:HG3	2.03	0.59
1:A:24:LEU:HD22	1:A:149:ASP:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:ARG:NH2	1:A:622:ILE:HD11	2.19	0.58
1:A:919:ASN:HA	1:A:985:LYS:HG3	1.85	0.58
1:A:59:ASP:HB3	1:A:62:VAL:HG12	1.86	0.57
1:A:52:LEU:HD11	1:A:109:GLU:HG2	1.85	0.57
1:A:560:ARG:NH2	4:A:1002:DL5:O1B	2.31	0.57
1:A:500:PRO:HG2	1:A:508:VAL:HB	1.87	0.56
1:A:604:ARG:HB2	1:A:607:VAL:HG23	1.88	0.56
1:A:778:THR:HB	1:A:783:LEU:O	2.05	0.56
1:A:78:PHE:O	1:A:84:THR:HG21	2.06	0.56
1:A:653:TYR:OH	1:A:672:ARG:NH2	2.31	0.56
1:A:179:ILE:HD13	1:A:212:THR:HA	1.86	0.55
1:A:862:GLY:CA	1:A:863:PRO:CD	2.85	0.55
1:A:415:THR:HA	1:A:475:ILE:HD13	1.89	0.55
1:A:491:ARG:NH2	1:A:584:PHE:HD2	2.04	0.55
1:A:957:PHE:O	1:A:958:LYS:HB2	2.06	0.55
1:A:266:LEU:HA	1:A:269:VAL:HG22	1.89	0.55
2:A:1000:TG1:H233	2:A:1000:TG1:C31	2.37	0.54
1:A:969:MET:HE3	1:A:972:LYS:HD2	1.90	0.54
1:A:65:LEU:HD11	1:A:307:ILE:HG21	1.89	0.54
1:A:512:MET:HB2	1:A:567:ARG:HB3	1.90	0.54
1:A:75:LEU:HD13	1:A:297:LYS:HZ3	1.73	0.53
1:A:843:TYR:CE2	1:A:973:ILE:HG23	2.43	0.53
1:A:315:ILE:HG22	1:A:760:PHE:HE1	1.73	0.52
1:A:855:TRP:CD1	1:A:896:PRO:HG3	2.45	0.52
1:A:960:LYS:HE2	1:A:967:TRP:CH2	2.44	0.52
2:A:1000:TG1:H233	2:A:1000:TG1:H311	1.92	0.52
2:A:1000:TG1:H232	2:A:1000:TG1:H262	1.92	0.52
1:A:61:LEU:HB2	1:A:307:ILE:HD12	1.92	0.52
1:A:894:PRO:HB3	1:A:959:LEU:N	2.17	0.51
1:A:486:GLU:O	1:A:491:ARG:NH2	2.43	0.51
1:A:943:LEU:O	1:A:947:ILE:HB	2.11	0.51
1:A:115:ALA:HB1	1:A:239:MET:CE	2.41	0.51
1:A:239:MET:CE	1:A:708:ALA:HB1	2.41	0.51
1:A:625:THR:HG21	1:A:632:ALA:HB2	1.94	0.50
1:A:777:LEU:HB2	1:A:849:VAL:HG21	1.93	0.50
1:A:361:MET:HG3	1:A:599:MET:HG2	1.94	0.50
1:A:894:PRO:HA	1:A:958:LYS:HD3	1.94	0.50
1:A:944:HIS:O	1:A:947:ILE:HG22	2.11	0.50
1:A:450:GLU:OE1	1:A:467:ARG:NH1	2.43	0.50
1:A:926:PRO:HB3	1:A:928:TRP:CE2	2.47	0.50
1:A:115:ALA:HB1	1:A:239:MET:HE2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:894:PRO:HA	1:A:958:LYS:HB3	1.94	0.50
1:A:560:ARG:HG3	4:A:1002:DL5:N1	2.26	0.50
1:A:835:PHE:HA	1:A:838:MET:HE2	1.94	0.49
1:A:979:GLY:O	1:A:983:ILE:HG12	2.13	0.49
1:A:795:VAL:HG13	1:A:799:THR:HB	1.95	0.49
2:A:1000:TG1:O4	2:A:1000:TG1:H161	2.13	0.49
1:A:483:PHE:HE2	1:A:573:ARG:HG3	1.76	0.49
1:A:113:GLU:HG3	1:A:334:ARG:HH22	1.78	0.48
1:A:894:PRO:HB3	1:A:958:LYS:HB3	1.95	0.48
1:A:580:ASP:OD1	1:A:582:SER:HB3	2.13	0.48
1:A:628:ASN:ND2	1:A:678:ARG:HH11	2.12	0.48
1:A:931:ILE:HD12	1:A:931:ILE:H	1.79	0.48
1:A:249:LEU:HD23	1:A:757:MET:HE2	1.95	0.48
1:A:76:ALA:HB1	1:A:88:PHE:CD2	2.48	0.48
1:A:832:TRP:CD1	1:A:988:ALA:HB2	2.49	0.48
1:A:785:GLU:HG2	1:A:856:PHE:HZ	1.77	0.47
1:A:47:LYS:HE3	1:A:112:ALA:HB3	1.96	0.47
1:A:922:LEU:HD23	1:A:927:PRO:HG3	1.96	0.47
1:A:954:PRO:O	1:A:958:LYS:N	2.43	0.47
1:A:847:ALA:HB2	1:A:973:ILE:HG21	1.96	0.47
1:A:246:LYS:HE3	1:A:254:ASP:OD2	2.14	0.47
1:A:757:MET:HA	1:A:760:PHE:CE2	2.49	0.47
1:A:559:LEU:HD22	1:A:600:LEU:HB2	1.97	0.46
1:A:628:ASN:HD21	1:A:678:ARG:HH11	1.62	0.46
1:A:75:LEU:HD11	1:A:300:VAL:HG11	1.98	0.46
1:A:561:CYS:HA	1:A:597:VAL:O	2.15	0.46
1:A:347:VAL:HB	1:A:698:THR:HG22	1.98	0.46
1:A:416:ILE:HG23	1:A:513:PHE:HB3	1.98	0.46
1:A:518:PRO:HB3	1:A:549:ILE:HD13	1.96	0.46
1:A:906:THR:HG22	1:A:941:MET:HE1	1.96	0.46
1:A:769:VAL:O	1:A:773:VAL:HG23	2.16	0.46
1:A:315:ILE:HG22	1:A:760:PHE:CE1	2.51	0.45
1:A:692:GLN:NE2	1:A:715:GLU:OE1	2.47	0.45
1:A:913:LEU:HB3	1:A:922:LEU:HD21	1.97	0.45
1:A:950:VAL:O	1:A:954:PRO:HD3	2.17	0.45
1:A:840:ILE:O	1:A:844:VAL:HG13	2.16	0.45
1:A:572:LYS:HB2	1:A:575:GLU:HG2	1.99	0.45
1:A:735:LEU:HD11	1:A:743:ILE:HG13	1.99	0.45
1:A:90:GLU:OE2	1:A:789:PRO:HB2	2.17	0.45
1:A:353:THR:HA	1:A:357:THR:OG1	2.16	0.45
1:A:113:GLU:CB	1:A:729:THR:HG23	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:862:GLY:O	1:A:863:PRO:C	2.59	0.45
1:A:352:LYS:HD2	1:A:635:ILE:HG13	1.98	0.44
1:A:692:GLN:OE1	1:A:713:LYS:HE2	2.17	0.44
1:A:894:PRO:CA	1:A:958:LYS:HB3	2.48	0.44
1:A:795:VAL:HG21	1:A:901:LEU:HD11	1.99	0.44
1:A:276:ILE:HA	1:A:279:PHE:CD2	2.52	0.44
1:A:247:THR:HG22	1:A:249:LEU:H	1.82	0.44
1:A:907:ILE:HD11	1:A:973:ILE:HG22	2.00	0.44
1:A:427:PHE:HA	1:A:434:TYR:HA	2.00	0.44
1:A:769:VAL:HG12	1:A:841:GLY:HA3	2.00	0.44
1:A:41:LEU:O	1:A:120:LYS:NZ	2.50	0.44
1:A:786:ALA:H	1:A:897:MET:HG2	1.83	0.44
1:A:249:LEU:HD13	1:A:754:TYR:HE1	1.83	0.43
1:A:259:GLN:HB3	2:A:1000:TG1:C26	2.48	0.43
1:A:620:ARG:NH2	1:A:670:CYS:O	2.42	0.43
1:A:114:ASN:ND2	1:A:117:GLU:HB2	2.34	0.43
1:A:342:LEU:HD12	1:A:716:ILE:HD13	2.00	0.43
1:A:713:LYS:HA	1:A:713:LYS:HD2	1.84	0.43
1:A:769:VAL:HG13	2:A:1000:TG1:H151	2.00	0.43
1:A:777:LEU:O	1:A:781:LEU:HG	2.19	0.43
2:A:1000:TG1:O10	2:A:1000:TG1:H313	2.19	0.43
1:A:129:VAL:HG12	1:A:151:VAL:HG22	1.99	0.43
1:A:696:GLU:O	1:A:698:THR:HG23	2.19	0.43
1:A:862:GLY:C	1:A:863:PRO:C	2.85	0.43
1:A:761:ILE:O	1:A:765:ILE:HG12	2.18	0.43
1:A:133:ASP:OD1	1:A:133:ASP:N	2.52	0.42
1:A:788:ILE:HG22	1:A:790:VAL:H	1.84	0.42
1:A:576:MET:HE2	1:A:576:MET:HB3	1.89	0.42
1:A:571:PRO:HG2	1:A:576:MET:SD	2.59	0.42
1:A:473:SER:O	1:A:477:GLN:HG3	2.20	0.42
1:A:560:ARG:HG3	4:A:1002:DL5:C6	2.49	0.42
1:A:752:ALA:HB1	1:A:812:PRO:HB3	2.02	0.42
1:A:368:ILE:HD12	1:A:409:GLY:HA3	2.02	0.42
1:A:718:ILE:HG12	1:A:733:MET:HE2	2.02	0.42
1:A:819:ARG:HE	1:A:819:ARG:HB3	1.68	0.42
1:A:116:ILE:HG23	1:A:236:ARG:HG3	2.02	0.42
1:A:483:PHE:CE2	1:A:573:ARG:HG3	2.54	0.42
1:A:491:ARG:NH1	1:A:588:GLU:OE2	2.53	0.41
1:A:19:SER:HB3	1:A:22:THR:OG1	2.19	0.41
1:A:27:ASP:O	1:A:31:ARG:HG3	2.19	0.41
1:A:256:PHE:CZ	1:A:765:ILE:HD12	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:SER:OG	1:A:300:VAL:HG21	2.19	0.41
1:A:369:ILE:HG13	1:A:528:VAL:CG1	2.49	0.41
1:A:549:ILE:HD11	1:A:596:VAL:HG21	2.01	0.41
1:A:715:GLU:C	1:A:716:ILE:HG13	2.45	0.41
1:A:114:ASN:HD22	1:A:117:GLU:HB2	1.86	0.41
1:A:911:ASN:HA	1:A:914:ASN:HB2	2.02	0.41
1:A:948:LEU:HB3	1:A:960:LYS:HB2	2.01	0.41
1:A:54:ILE:O	1:A:58:GLU:HG3	2.20	0.41
1:A:611:ILE:HG13	1:A:639:ILE:HG23	2.02	0.41
1:A:648:VAL:HG12	1:A:648:VAL:O	2.20	0.41
1:A:459:VAL:HA	1:A:462:LEU:HG	2.02	0.41
1:A:684:LYS:HB2	1:A:710:ALA:HB2	2.02	0.41
1:A:921:SER:O	1:A:925:MET:HB2	2.20	0.41
1:A:454:VAL:HG23	1:A:455:PHE:CD2	2.56	0.41
1:A:712:LYS:HE3	1:A:712:LYS:HB3	1.96	0.41
1:A:305:ALA:HB2	1:A:775:ILE:HD12	2.01	0.41
1:A:724:THR:O	1:A:728:LYS:HG3	2.21	0.41
1:A:747:VAL:O	1:A:751:ARG:HB2	2.21	0.41
1:A:837:TYR:O	1:A:841:GLY:N	2.54	0.41
1:A:124:PRO:HG3	1:A:158:LYS:O	2.20	0.41
1:A:166:LEU:HG	1:A:221:GLY:HA2	2.03	0.41
1:A:170:SER:HB2	1:A:172:THR:O	2.21	0.41
1:A:403:ARG:HD3	1:A:404:SER:H	1.86	0.40
1:A:413:LEU:HD11	1:A:564:LEU:HD13	2.04	0.40
1:A:907:ILE:HG12	1:A:974:SER:HA	2.03	0.40
1:A:873:PHE:CE1	1:A:874:MET:HG3	2.57	0.40
1:A:416:ILE:HG12	1:A:513:PHE:CD1	2.56	0.40
1:A:528:VAL:HG21	1:A:541:VAL:HG11	2.03	0.40
1:A:679:VAL:HB	1:A:683:HIS:HB2	2.03	0.40
1:A:951:ASP:O	1:A:954:PRO:HD2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	989/995 (99%)	934 (94%)	52 (5%)	3 (0%)	36 63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	865	VAL
1	A	958	LYS
1	A	500	PRO

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	838/840 (100%)	821 (98%)	17 (2%)	48 69

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	30	LYS
1	A	55	GLU
1	A	123	GLU
1	A	141	LYS
1	A	170	SER
1	A	171	THR
1	A	244	GLN
1	A	266	LEU
1	A	309	GLU
1	A	319	LEU
1	A	340	GLU
1	A	508	VAL
1	A	599	MET
1	A	679	VAL
1	A	840	ILE

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Mol	Chain	Res	Type
1	A	918	GLU

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

Mol	Chain	Res	Type
1	A	114	ASN
1	A	456	ASN
1	A	510	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	DL5	A	1002	5	41,50,50	1.53	6 (14%)	54,80,80	1.19	6 (11%)
2	TG1	A	1000	-	44,48,48	0.83	0	47,72,72	1.22	7 (14%)
3	VN4	A	1001	5,8,1	0,3,3	-	-	-		

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	DL5	A	1002	5	-	11/23/80/80	0/5/5/5
2	TG1	A	1000	-	-	11/33/99/99	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	DL5	C5F-C6F	5.05	1.51	1.37
4	A	1002	DL5	PA-O3A	3.63	1.63	1.59
4	A	1002	DL5	PA-O1A	2.37	1.59	1.50
4	A	1002	DL5	PG-O3G	-2.16	1.50	1.55
4	A	1002	DL5	C5F-C4F	-2.12	1.32	1.40
4	A	1002	DL5	O7F-N6F	-2.04	1.21	1.35

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1002	DL5	O2'-C1F-C6F	3.46	115.20	110.47
2	A	1000	TG1	C7-C6-C5	3.37	124.25	115.41
4	A	1002	DL5	O2'-C1F-C2F	-2.78	106.67	110.47
4	A	1002	DL5	O4'-C1'-N9	-2.56	103.17	108.09
2	A	1000	TG1	C10-C1-C5	2.49	118.46	115.07
4	A	1002	DL5	O3G-PG-C01	2.33	112.04	106.40
2	A	1000	TG1	O1-C13-O2	-2.29	118.34	123.70
4	A	1002	DL5	O2G-PG-O1G	-2.26	106.53	112.39
2	A	1000	TG1	C10-O9-C32	2.21	126.53	121.36
2	A	1000	TG1	C6-O5-C12	-2.17	107.80	110.80
2	A	1000	TG1	O5-C6-C5	2.12	114.31	111.17
4	A	1002	DL5	C4'-O4'-C1'	2.11	114.13	109.47
2	A	1000	TG1	C11-C7-C8	2.06	117.19	114.44

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1000	TG1	C14-C13-O1-C2
2	A	1000	TG1	C2-C3-O3-C21
2	A	1000	TG1	C22-C21-O3-C3

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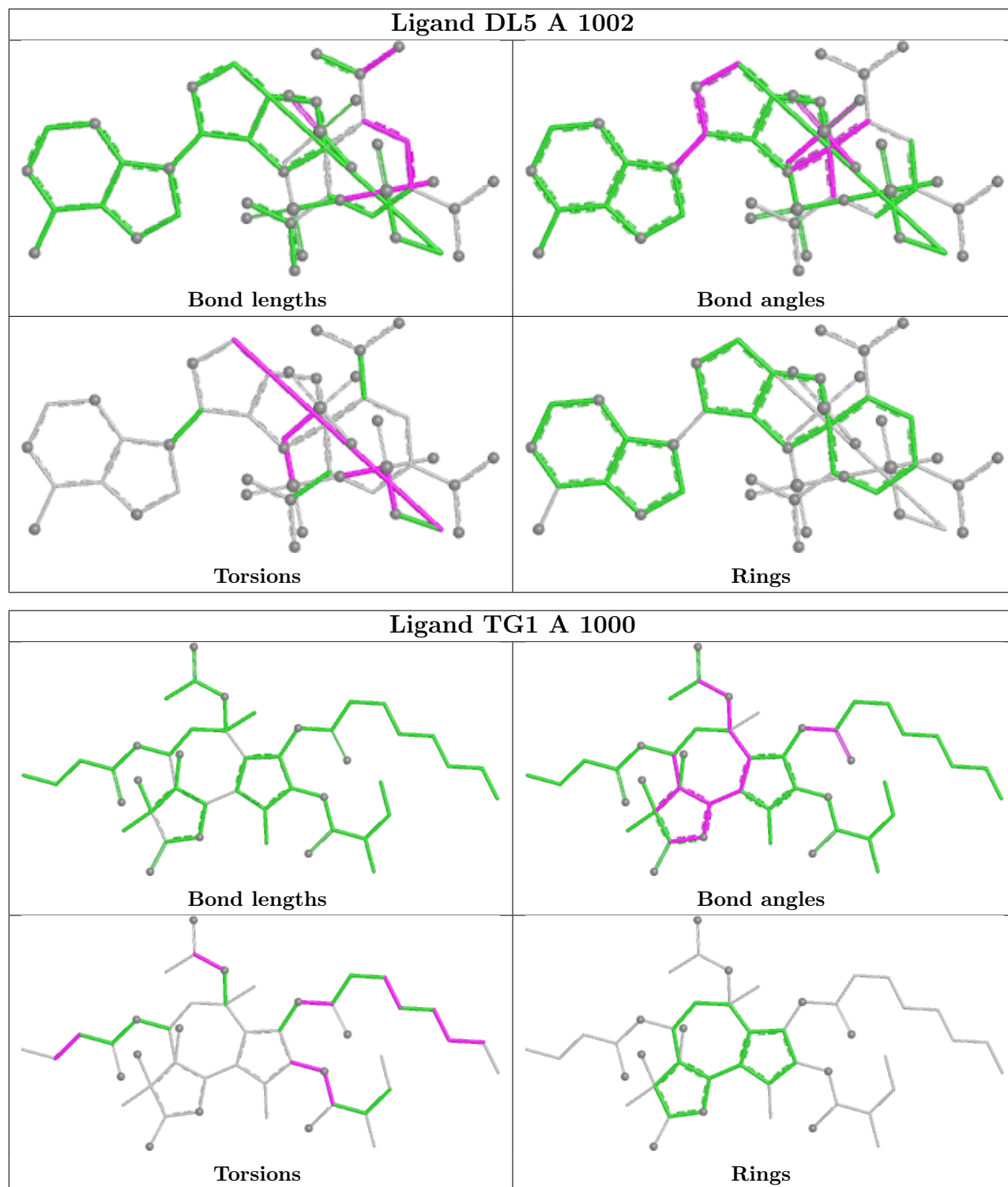
Mol	Chain	Res	Type	Atoms
4	A	1002	DL5	C5'-O5'-PA-O2A
4	A	1002	DL5	C5'-O5'-PA-O3A
4	A	1002	DL5	PG-C01-PB-O1B
4	A	1002	DL5	PG-C01-PB-O2B
4	A	1002	DL5	PG-C01-PB-O3A
4	A	1002	DL5	PB-C01-PG-O1G
2	A	1000	TG1	O2-C13-O1-C2
2	A	1000	TG1	O4-C21-O3-C3
2	A	1000	TG1	C33-C32-O9-C10
2	A	1000	TG1	O10-C32-O9-C10
4	A	1002	DL5	O4'-C4'-C5'-O5'
2	A	1000	TG1	C16-C17-C18-C19
2	A	1000	TG1	C14-C15-C16-C17
2	A	1000	TG1	C17-C18-C19-C20
2	A	1000	TG1	C27-C28-C29-C30
4	A	1002	DL5	C5'-O5'-PA-O1A
4	A	1002	DL5	PB-C01-PG-O3G
4	A	1002	DL5	PB-O3A-PA-O2A
4	A	1002	DL5	C3'-C4'-C5'-O5'

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1002	DL5	3	0
2	A	1000	TG1	11	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 ⓘ

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	994/995 (99%)	0.13	59 (5%) 28 14	44, 92, 225, 324	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	786	ALA	9.5
1	A	850	GLY	7.3
1	A	863	PRO	7.1
1	A	851	ALA	6.8
1	A	866	THR	5.5
1	A	785	GLU	5.4
1	A	899	MET	4.7
1	A	781	LEU	4.6
1	A	787	LEU	4.5
1	A	898	THR	4.3
1	A	791	GLN	4.3
1	A	901	LEU	4.0
1	A	963	ASP	3.9
1	A	849	VAL	3.9
1	A	897	MET	3.7
1	A	854	TRP	3.7
1	A	777	LEU	3.7
1	A	788	ILE	3.7
1	A	967	TRP	3.6
1	A	57	PHE	3.5
1	A	972	LYS	3.5
1	A	857	MET	3.5
1	A	867	TYR	3.4
1	A	501	ALA	3.4
1	A	902	SER	3.3
1	A	886	LEU	3.3
1	A	964	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	317	THR	3.2
1	A	960	LYS	3.2
1	A	758	LYS	3.1
1	A	969	MET	3.1
1	A	316	THR	3.1
1	A	896	PRO	3.1
1	A	959	LEU	3.0
1	A	890	ILE	3.0
1	A	852	ALA	2.9
1	A	968	LEU	2.8
1	A	957	PHE	2.8
1	A	848	THR	2.8
1	A	962	LEU	2.7
1	A	865	VAL	2.7
1	A	840	ILE	2.6
1	A	503	SER	2.6
1	A	949	TYR	2.6
1	A	864	GLY	2.5
1	A	61	LEU	2.5
1	A	862	GLY	2.5
1	A	853	ALA	2.5
1	A	508	VAL	2.5
1	A	215	ALA	2.4
1	A	895	GLU	2.4
1	A	766	SER	2.4
1	A	340	GLU	2.3
1	A	320	ALA	2.3
1	A	911	ASN	2.3
1	A	953	LEU	2.2
1	A	294	TYR	2.1
1	A	258	GLU	2.1
1	A	855	TRP	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

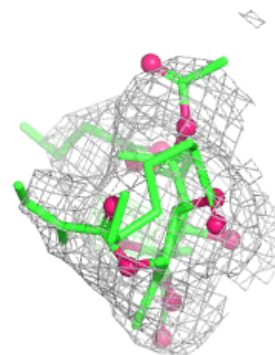
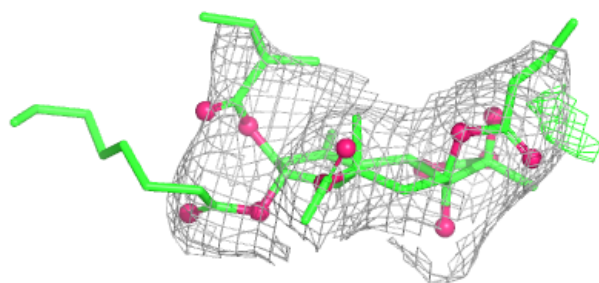
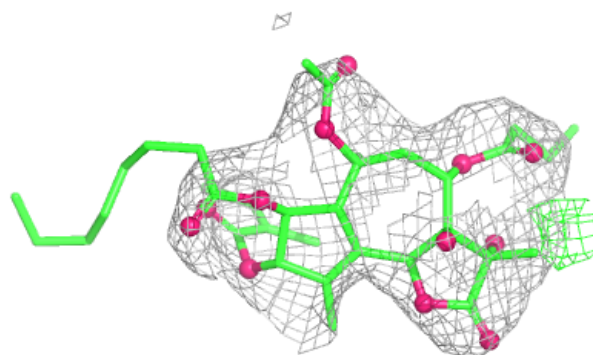
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

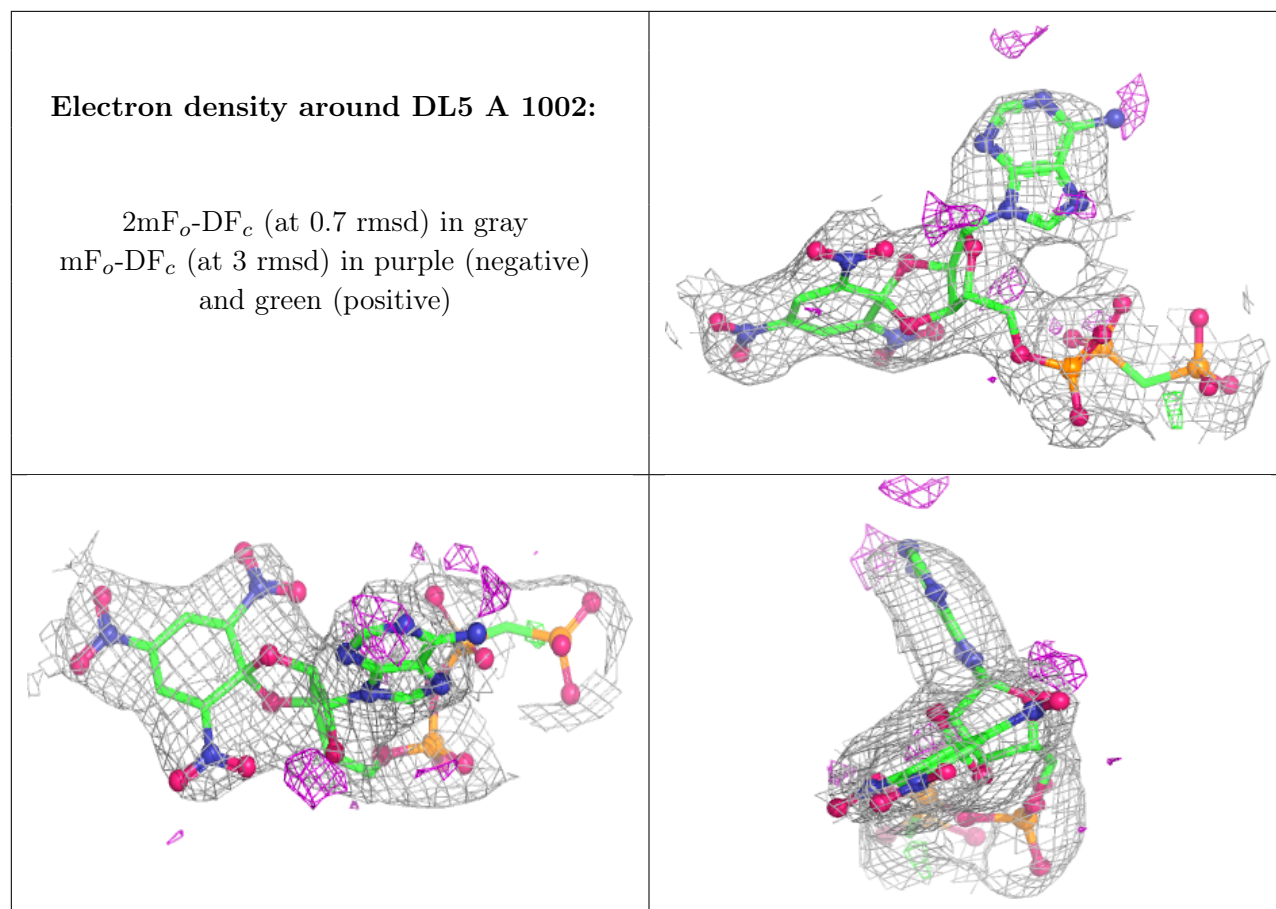
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	TG1	A	1000	46/46	0.89	0.13	111,146,158,171	0
4	DL5	A	1002	46/46	0.92	0.11	93,114,141,160	0
5	MG	A	1004	1/1	0.92	0.09	103,103,103,103	0
7	CL	A	1009	1/1	0.97	0.09	65,65,65,65	0
6	K	A	1008	1/1	0.99	0.04	73,73,73,73	0
3	VN4	A	1001	4/4	0.99	0.09	52,57,63,73	0
5	MG	A	1003	1/1	1.00	0.02	36,36,36,36	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 TG1 A 1000:

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 ⓘ

There are no such residues in this entry.