



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:07 AM UTC

PDB ID : 3H73 / pdb_00003h73
Title : Crystal structure of Streptococcus pneumoniae D39 neuraminidase A precursor (NanA) in complex with DANA
Authors : Hsiao, Y.-S.; Tong, L.
Deposited on : 2009-04-24
Resolution : 1.70 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

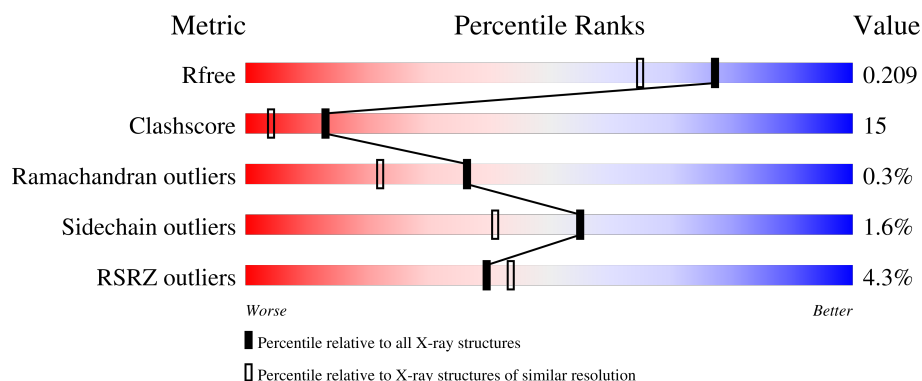
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	477	5% 75% 23% ..
1	B	477	4% 73% 25% .

2 Entry composition [i](#)

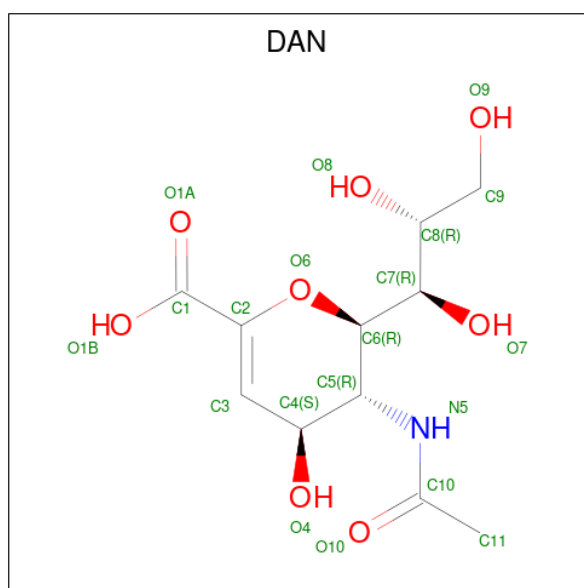
There are 3 unique types of molecules in this entry. The entry contains 9097 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Sialidase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	474	Total	C	N	O	S	0	0	0
			3747	2347	661	727	12			
1	B	477	Total	C	N	O	S	0	0	0
			3771	2363	664	732	12			

- Molecule 2 is 2-DEOXY-2,3-DEHYDRO-N-ACETYL-NEURAMINIC ACID (CCD ID: DAN) (formula: $C_{11}H_{17}NO_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			20	11	1	8		
2	B	1	Total	C	N	O	0	0
			20	11	1	8		

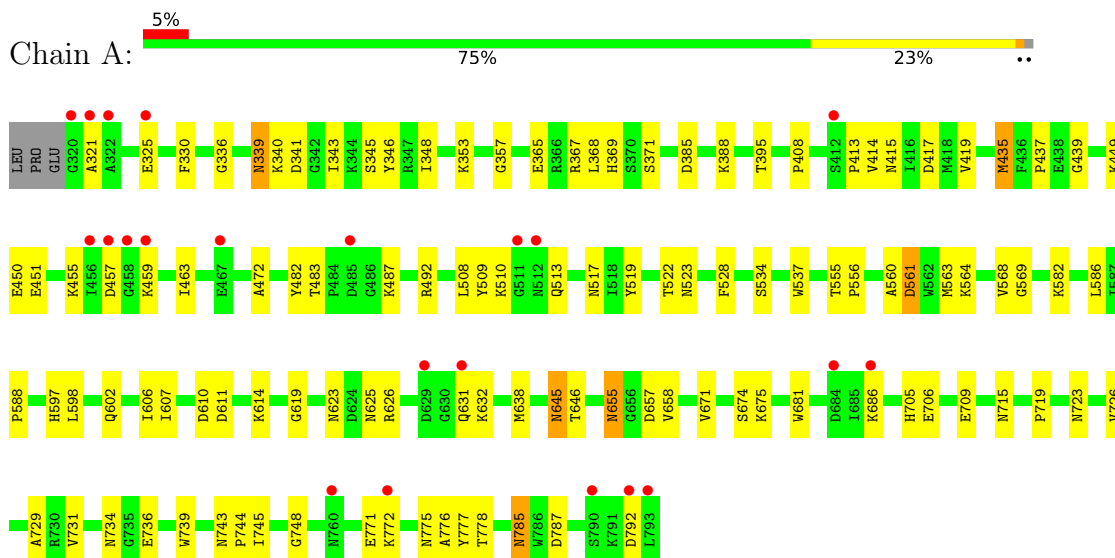
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	752	Total 752	O 752	0	0
3	B	787	Total 787	O 787	0	0

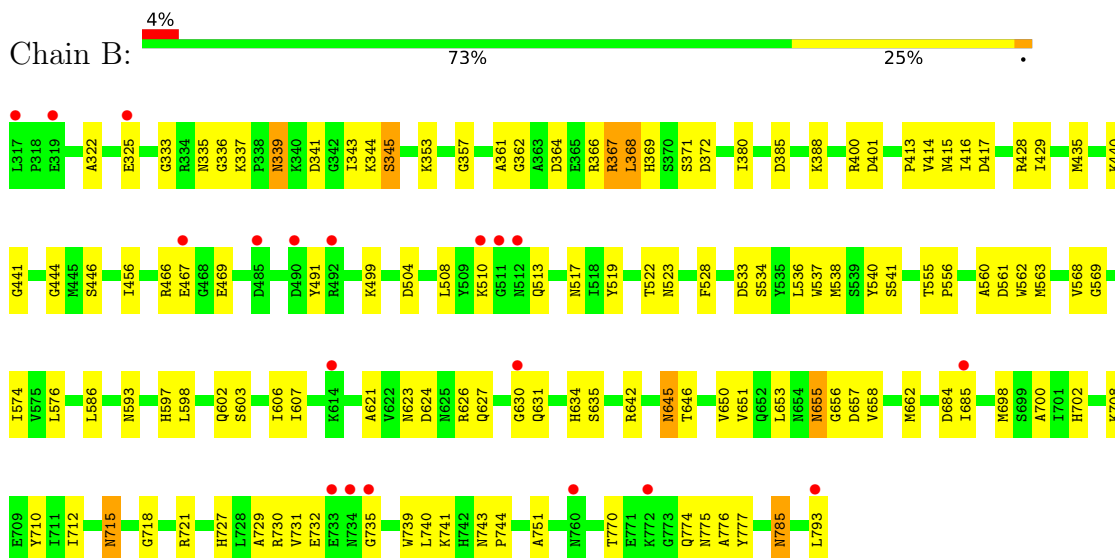
3 Residue-property plots

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: Sialidase A



• Molecule 1: Sialidase A



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	158.34Å 47.40Å 137.32Å 90.00° 116.45° 90.00°	Depositor
Resolution (Å)	27.34 – 1.70 27.34 – 1.70	Depositor EDS
% Data completeness (in resolution range)	92.6 (27.34-1.70) 92.6 (27.34-1.70)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.51 (at 1.70Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.178 , 0.209 0.178 , 0.209	Depositor DCC
R_{free} test set	7082 reflections (7.49%)	wwPDB-VP
Wilson B-factor (Å ²)	10.4	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.1	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	0.95	EDS
Total number of atoms	9097	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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: DAN

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	0.44	1/3828 (0.0%)	0.89	13/5169 (0.3%)
1	B	0.39	0/3853	0.84	12/5204 (0.2%)
All	All	0.41	1/7681 (0.0%)	0.86	25/10373 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	435	MET	CG-SD	-5.07	1.68	1.80

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	569	GLY	CA-C-N	7.42	127.47	119.90
1	A	569	GLY	C-N-CA	7.42	127.47	119.90
1	A	414	VAL	N-CA-C	7.18	118.81	108.48
1	B	367	ARG	N-CA-C	7.16	118.73	111.07
1	A	367	ARG	N-CA-C	6.82	118.37	111.07
1	A	569	GLY	N-CA-C	-6.77	98.53	112.34
1	A	528	PHE	N-CA-C	6.70	120.33	109.40
1	B	569	GLY	N-CA-C	-6.66	98.75	112.34
1	B	368	LEU	N-CA-C	6.50	120.42	112.23
1	B	414	VAL	N-CA-C	6.45	117.88	108.53
1	A	345	SER	N-CA-C	5.95	118.45	109.23
1	A	619	GLY	N-CA-C	-5.83	103.67	112.60
1	A	368	LEU	N-CA-C	5.79	119.52	112.23
1	A	437	PRO	N-CA-C	-5.74	103.18	111.22
1	A	607	ILE	N-CA-C	-5.54	100.34	108.99
1	A	413	PRO	N-CA-C	-5.53	102.42	111.21
1	A	415	ASN	N-CA-C	-5.53	100.39	109.40
1	B	528	PHE	N-CA-C	5.52	118.55	109.72

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	607	ILE	N-CA-C	-5.41	100.53	108.85
1	B	415	ASN	N-CA-C	-5.37	100.65	109.40
1	B	569	GLY	CA-C-N	5.35	126.53	119.84
1	B	569	GLY	C-N-CA	5.35	126.53	119.84
1	B	662	MET	N-CA-C	5.28	117.56	109.07
1	B	345	SER	N-CA-C	5.24	117.15	109.24
1	B	743	ASN	N-CA-C	5.22	116.28	109.64

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	3747	0	3664	102	0
1	B	3771	0	3688	121	0
2	A	20	0	16	0	0
2	B	20	0	16	0	0
3	A	752	0	0	45	0
3	B	787	0	0	54	0
All	All	9097	0	7384	222	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 15.

All (222) 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:681:TRP:HA	3:A:1173:HOH:O	1.54	1.05
1:B:428:ARG:HG2	3:B:1136:HOH:O	1.59	1.01
1:A:419:VAL:HG21	3:A:1140:HOH:O	1.61	1.00
1:A:517:ASN:HD22	1:A:519:TYR:H	1.11	0.94
1:B:336:GLY:H	1:B:775:ASN:HD21	1.15	0.93
1:B:517:ASN:HD22	1:B:519:TYR:H	1.16	0.91
1:A:686:LYS:HB3	3:A:1135:HOH:O	1.75	0.87

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:LYS:HG3	3:A:1118:HOH:O	1.73	0.86
1:A:671:VAL:HB	3:A:1135:HOH:O	1.77	0.85
1:B:541:SER:HB3	3:B:848:HOH:O	1.76	0.84
1:B:698:MET:HA	3:B:993:HOH:O	1.78	0.83
1:A:336:GLY:H	1:A:775:ASN:HD21	1.24	0.81
1:A:482:TYR:HB3	3:A:1263:HOH:O	1.81	0.81
1:A:395:THR:HG23	3:A:1200:HOH:O	1.81	0.80
1:B:593:ASN:HB3	3:B:1104:HOH:O	1.81	0.79
1:B:635:SER:HB3	3:B:1135:HOH:O	1.83	0.78
1:B:366:ARG:HD3	3:B:1199:HOH:O	1.83	0.77
1:B:603:SER:HB2	3:B:1497:HOH:O	1.85	0.77
1:A:674:SER:HA	3:A:1173:HOH:O	1.88	0.74
1:A:568:VAL:HG22	3:A:1159:HOH:O	1.86	0.73
1:A:560:ALA:H	1:A:563:MET:HE3	1.54	0.73
1:B:634:HIS:HA	3:B:1207:HOH:O	1.88	0.72
1:A:451:GLU:HA	3:A:1266:HOH:O	1.88	0.72
1:A:341:ASP:HB2	3:A:1249:HOH:O	1.90	0.72
1:B:429:ILE:HB	3:B:848:HOH:O	1.90	0.72
1:B:774:GLN:HG3	3:B:1210:HOH:O	1.88	0.72
1:B:635:SER:N	3:B:1207:HOH:O	2.23	0.72
1:A:517:ASN:ND2	1:A:519:TYR:H	1.86	0.71
3:A:1276:HOH:O	1:B:561:ASP:HB2	1.90	0.70
1:B:372:ASP:HB2	3:B:1103:HOH:O	1.91	0.69
1:B:655:ASN:ND2	1:B:657:ASP:H	1.90	0.69
1:B:624:ASP:HA	3:B:1207:HOH:O	1.93	0.69
1:A:582:LYS:HB3	3:A:1137:HOH:O	1.93	0.68
1:B:562:TRP:HZ3	3:B:1104:HOH:O	1.76	0.68
1:A:602:GLN:HB3	1:A:646:THR:HG22	1.76	0.67
1:B:751:ALA:HB2	3:B:1115:HOH:O	1.95	0.67
1:B:517:ASN:ND2	1:B:519:TYR:H	1.91	0.66
1:B:435:MET:HE2	1:B:537:TRP:CZ2	2.30	0.66
1:A:655:ASN:ND2	1:A:657:ASP:H	1.94	0.65
1:B:685:ILE:HA	3:B:980:HOH:O	1.95	0.65
1:A:343:ILE:HD12	3:A:1249:HOH:O	1.96	0.65
1:A:560:ALA:CB	1:A:563:MET:HE3	2.27	0.64
1:B:339:ASN:ND2	1:B:343:ILE:H	1.96	0.64
1:A:340:LYS:HE3	3:A:1307:HOH:O	1.97	0.64
1:B:345:SER:HB3	3:B:1164:HOH:O	1.97	0.63
1:A:674:SER:CA	3:A:1173:HOH:O	2.45	0.63
1:B:513:GLN:HB3	3:B:1156:HOH:O	1.98	0.63
1:A:339:ASN:ND2	1:A:343:ILE:H	1.97	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:ILE:HB	3:A:1249:HOH:O	2.01	0.61
1:B:627:GLN:NE2	1:B:630:GLY:H	1.99	0.60
1:B:385:ASP:CG	1:B:388:LYS:HB3	2.27	0.60
1:B:650:VAL:HG23	3:B:1051:HOH:O	2.02	0.60
1:B:339:ASN:C	1:B:339:ASN:HD22	2.09	0.60
1:B:499:LYS:HD2	1:B:504:ASP:HB3	1.83	0.60
1:B:623:ASN:HD21	1:B:645:ASN:HD21	1.48	0.60
1:B:623:ASN:C	3:B:1207:HOH:O	2.44	0.60
1:B:655:ASN:C	1:B:655:ASN:HD22	2.09	0.59
1:A:339:ASN:C	1:A:339:ASN:HD22	2.10	0.59
1:B:367:ARG:HH21	1:B:368:LEU:HD21	1.67	0.59
1:B:655:ASN:HD22	1:B:656:GLY:N	2.01	0.58
1:A:417:ASP:O	3:A:1159:HOH:O	2.16	0.58
1:A:588:PRO:HD2	3:A:1140:HOH:O	2.03	0.58
1:B:538:MET:HE1	3:B:1521:HOH:O	2.03	0.58
1:A:771:GLU:O	1:A:772:LYS:HG2	2.04	0.58
1:A:655:ASN:HD21	1:A:657:ASP:HB2	1.70	0.57
1:B:560:ALA:HB3	1:B:563:MET:HG3	1.87	0.57
1:A:586:LEU:HD22	1:A:606:ILE:HD13	1.87	0.57
1:B:623:ASN:HB2	3:B:1135:HOH:O	2.04	0.56
1:B:715:ASN:HA	3:B:993:HOH:O	2.05	0.56
1:B:598:LEU:HB3	3:B:941:HOH:O	2.06	0.56
1:B:339:ASN:ND2	1:B:341:ASP:H	2.03	0.56
1:B:623:ASN:HD21	1:B:626:ARG:HH11	1.54	0.56
1:B:642:ARG:HD2	3:B:1006:HOH:O	2.06	0.56
1:A:785:ASN:C	1:A:785:ASN:HD22	2.14	0.55
1:A:561:ASP:HB2	1:B:562:TRP:HD1	1.71	0.55
1:A:339:ASN:ND2	1:A:341:ASP:H	2.04	0.55
1:B:621:ALA:HB3	3:B:1135:HOH:O	2.06	0.55
1:A:560:ALA:H	1:A:563:MET:CE	2.19	0.54
1:A:792:ASP:HA	3:A:1442:HOH:O	2.07	0.54
1:A:408:PRO:HG3	3:A:1362:HOH:O	2.07	0.54
1:B:623:ASN:ND2	1:B:626:ARG:HH11	2.06	0.54
1:B:576:LEU:HA	3:B:1011:HOH:O	2.08	0.54
1:B:574:ILE:HD12	3:B:1011:HOH:O	2.07	0.53
1:B:336:GLY:N	1:B:775:ASN:HD21	1.97	0.53
1:B:540:TYR:HB2	3:B:1136:HOH:O	2.08	0.53
1:A:510:LYS:HB3	3:A:1196:HOH:O	2.08	0.53
1:A:564:LYS:HG3	3:A:1235:HOH:O	2.09	0.52
1:A:439:GLY:HA2	3:A:1213:HOH:O	2.08	0.52
1:B:336:GLY:H	1:B:775:ASN:ND2	1.96	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:ALA:H	1:B:785:ASN:HD21	1.57	0.52
1:A:339:ASN:ND2	1:A:339:ASN:C	2.66	0.52
1:A:623:ASN:HD21	1:A:645:ASN:HD21	1.58	0.52
1:B:586:LEU:HD22	1:B:606:ILE:HD13	1.92	0.51
1:A:748:GLY:N	3:A:1154:HOH:O	2.43	0.51
1:A:449:LYS:HE3	3:A:1266:HOH:O	2.10	0.51
1:B:325:GLU:HG3	3:B:1286:HOH:O	2.10	0.51
1:A:417:ASP:HB2	3:A:1159:HOH:O	2.10	0.51
1:B:602:GLN:HB3	1:B:646:THR:HG22	1.93	0.51
1:A:602:GLN:CB	1:A:646:THR:HG22	2.41	0.51
1:A:729:ALA:HB2	1:A:739:TRP:CE3	2.45	0.51
1:A:625:ASN:HB2	3:B:1390:HOH:O	2.11	0.51
1:B:335:ASN:HB2	1:B:337:LYS:HE2	1.93	0.50
1:B:623:ASN:ND2	1:B:645:ASN:HD21	2.09	0.50
1:A:487:LYS:N	3:A:1263:HOH:O	2.43	0.50
1:A:435:MET:HG3	1:A:537:TRP:CE2	2.47	0.50
1:B:401:ASP:HB3	3:B:270:HOH:O	2.10	0.50
1:B:653:LEU:HB2	1:B:655:ASN:ND2	2.27	0.50
1:A:534:SER:HA	3:A:992:HOH:O	2.12	0.49
1:B:729:ALA:HB2	1:B:739:TRP:CE3	2.47	0.49
1:B:401:ASP:HB3	3:B:881:HOH:O	2.12	0.49
1:B:651:VAL:HG13	1:B:702:HIS:HB2	1.94	0.49
1:B:708:LYS:HB2	1:B:710:TYR:CE2	2.48	0.49
1:B:413:PRO:HG2	3:B:236:HOH:O	2.12	0.49
1:B:536:LEU:HD11	1:B:568:VAL:HG13	1.94	0.49
1:A:385:ASP:CG	1:A:388:LYS:HB2	2.38	0.49
1:B:364:ASP:OD2	1:B:416:ILE:HG13	2.12	0.49
1:B:655:ASN:ND2	1:B:655:ASN:C	2.70	0.49
1:A:657:ASP:OD1	1:A:675:LYS:HD3	2.13	0.49
1:B:446:SER:HB3	3:B:827:HOH:O	2.12	0.48
1:B:456:ILE:HA	3:B:1528:HOH:O	2.13	0.48
1:A:611:ASP:CG	1:A:614:LYS:HB3	2.39	0.48
1:A:638:MET:HB2	3:A:1067:HOH:O	2.12	0.48
1:A:658:VAL:HB	1:A:674:SER:HB3	1.96	0.48
1:B:367:ARG:HG3	3:B:1352:HOH:O	2.12	0.48
1:A:597:HIS:CG	1:A:598:LEU:H	2.32	0.48
1:A:455:LYS:HB3	3:A:1103:HOH:O	2.14	0.47
1:B:435:MET:HE2	1:B:537:TRP:HZ2	1.74	0.47
1:B:634:HIS:CA	3:B:1207:HOH:O	2.53	0.47
1:A:631:GLN:HB2	3:A:1473:HOH:O	2.14	0.47
1:A:457:ASP:O	1:A:459:LYS:HG2	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:ASN:ND2	1:B:339:ASN:C	2.72	0.47
1:B:555:THR:N	1:B:556:PRO:HD2	2.30	0.47
1:A:555:THR:N	1:A:556:PRO:HD2	2.30	0.47
1:B:602:GLN:NE2	3:B:941:HOH:O	2.47	0.47
1:B:441:GLY:N	3:B:1201:HOH:O	2.47	0.47
1:B:444:GLY:HA3	3:B:1201:HOH:O	2.14	0.47
1:B:744:PRO:HD2	1:B:793:LEU:HD12	1.97	0.47
1:A:560:ALA:HB3	1:A:563:MET:HE3	1.96	0.47
1:B:429:ILE:C	3:B:1136:HOH:O	2.57	0.47
1:A:513:GLN:HB3	3:A:1196:HOH:O	2.13	0.46
1:A:623:ASN:ND2	1:A:626:ARG:HH11	2.12	0.46
1:A:330:PHE:HB3	1:A:346:TYR:CG	2.51	0.46
1:A:614:LYS:HD2	3:A:1146:HOH:O	2.16	0.46
1:B:576:LEU:HD23	3:B:1011:HOH:O	2.16	0.46
1:A:623:ASN:HD21	1:A:626:ARG:HH11	1.62	0.46
1:B:435:MET:HG3	1:B:537:TRP:CZ2	2.50	0.46
1:B:740:LEU:O	1:B:741:LYS:HD2	2.16	0.46
1:B:322:ALA:H	1:B:785:ASN:ND2	2.13	0.46
1:A:435:MET:HE2	1:A:537:TRP:CZ2	2.50	0.46
1:B:730:ARG:HB2	1:B:740:LEU:HD11	1.98	0.46
1:A:674:SER:HB2	3:A:1173:HOH:O	2.15	0.46
1:A:336:GLY:H	1:A:775:ASN:ND2	2.04	0.46
1:B:369:HIS:CE1	1:B:371:SER:HB2	2.51	0.46
1:A:450:GLU:N	3:A:1118:HOH:O	2.40	0.45
1:A:625:ASN:OD1	1:A:632:LYS:HD2	2.16	0.45
1:A:560:ALA:N	1:A:563:MET:HE3	2.29	0.45
1:B:606:ILE:C	1:B:606:ILE:HD12	2.41	0.45
1:B:715:ASN:HD22	1:B:715:ASN:N	2.14	0.45
1:A:369:HIS:CE1	1:A:371:SER:HB2	2.51	0.45
1:A:610:ASP:N	3:A:940:HOH:O	2.49	0.45
1:A:655:ASN:C	1:A:655:ASN:HD22	2.24	0.45
1:B:491:TYR:CE2	1:B:510:LYS:HG2	2.51	0.45
1:B:770:THR:HA	1:B:774:GLN:HE21	1.81	0.45
1:A:674:SER:CB	3:A:1173:HOH:O	2.63	0.45
1:A:743:ASN:HA	1:A:744:PRO:HD3	1.85	0.45
1:A:483:THR:N	3:A:1263:HOH:O	2.50	0.44
1:B:700:ALA:HA	1:B:712:ILE:O	2.17	0.44
1:B:715:ASN:ND2	1:B:727:HIS:NE2	2.66	0.44
1:A:435:MET:HG3	1:A:537:TRP:CZ2	2.53	0.44
1:A:723:ASN:N	3:A:1154:HOH:O	2.50	0.44
1:B:602:GLN:CB	1:B:646:THR:HG22	2.48	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:ILE:HG21	1:A:417:ASP:HA	2.00	0.44
1:A:726:VAL:HG23	3:A:1022:HOH:O	2.17	0.44
1:B:586:LEU:HG	3:B:1109:HOH:O	2.17	0.44
1:A:343:ILE:HD13	1:A:365:GLU:CG	2.48	0.44
1:B:385:ASP:OD2	1:B:388:LYS:HB3	2.17	0.44
1:B:730:ARG:HD3	1:B:731:VAL:N	2.32	0.44
1:A:655:ASN:ND2	1:A:655:ASN:C	2.74	0.43
1:A:745:ILE:HB	3:A:1022:HOH:O	2.17	0.43
1:B:362:GLY:HA2	1:B:380:ILE:HG22	2.00	0.43
1:B:785:ASN:C	1:B:785:ASN:HD22	2.26	0.43
1:A:343:ILE:HD13	1:A:365:GLU:HG3	2.00	0.43
1:B:466:ARG:HD3	1:B:469:GLU:OE2	2.19	0.43
1:B:353:LYS:HE3	1:B:357:GLY:HA2	2.00	0.43
1:B:718:GLY:N	3:B:1066:HOH:O	2.52	0.43
1:A:744:PRO:HB2	3:A:1090:HOH:O	2.19	0.43
1:B:333:GLY:HA3	1:B:337:LYS:O	2.19	0.43
1:B:631:GLN:HB3	3:B:1048:HOH:O	2.18	0.42
1:B:658:VAL:HG12	3:B:1051:HOH:O	2.17	0.42
1:A:606:ILE:C	1:A:606:ILE:HD12	2.44	0.42
1:A:597:HIS:CG	1:A:598:LEU:N	2.86	0.42
1:B:721:ARG:HB3	3:B:1115:HOH:O	2.18	0.42
1:A:321:ALA:HA	1:A:787:ASP:HB3	2.01	0.42
1:A:561:ASP:HB2	3:A:1276:HOH:O	2.19	0.42
1:B:339:ASN:HB2	3:B:926:HOH:O	2.19	0.42
1:B:400:ARG:HD2	3:B:1084:HOH:O	2.18	0.42
1:A:734:ASN:CG	1:A:736:GLU:HG2	2.45	0.42
1:B:533:ASP:CG	1:B:534:SER:H	2.28	0.42
1:A:492:ARG:HG2	1:A:509:TYR:HB2	2.02	0.42
1:B:597:HIS:CG	1:B:598:LEU:H	2.37	0.42
1:A:463:ILE:HG23	1:A:472:ALA:HB1	2.02	0.41
1:B:440:LYS:C	3:B:1201:HOH:O	2.63	0.41
1:A:709:GLU:HG2	1:A:731:VAL:HB	2.02	0.41
1:A:776:ALA:HA	1:A:777:TYR:HA	1.78	0.41
1:B:435:MET:HG3	1:B:537:TRP:CE2	2.56	0.41
1:B:651:VAL:CG1	1:B:702:HIS:HB2	2.51	0.41
1:B:730:ARG:HD3	1:B:731:VAL:C	2.45	0.41
1:B:344:LYS:CB	3:B:1164:HOH:O	2.68	0.41
1:A:705:HIS:CE1	1:A:706:GLU:HG3	2.55	0.41
1:B:776:ALA:HA	1:B:777:TYR:HA	1.71	0.41
1:A:353:LYS:HE3	1:A:357:GLY:HA2	2.02	0.41
1:B:621:ALA:CB	3:B:1135:HOH:O	2.67	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:522:THR:O	1:A:523:ASN:C	2.64	0.41
1:B:344:LYS:HD2	1:B:368:LEU:O	2.21	0.41
1:B:508:LEU:HD13	1:B:508:LEU:C	2.46	0.41
1:B:730:ARG:HD2	1:B:732:GLU:CD	2.46	0.41
1:A:508:LEU:C	1:A:508:LEU:HD13	2.46	0.41
1:A:623:ASN:ND2	1:A:645:ASN:HD21	2.18	0.41
1:B:735:GLY:HA2	3:B:985:HOH:O	2.21	0.40
1:A:325:GLU:HG2	3:A:1108:HOH:O	2.21	0.40
1:B:522:THR:O	1:B:523:ASN:C	2.62	0.40
1:B:361:ALA:O	1:B:380:ILE:HA	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	472/477 (99%)	442 (94%)	29 (6%)	1 (0%)	43	28
1	B	475/477 (100%)	446 (94%)	27 (6%)	2 (0%)	30	16
All	All	947/954 (99%)	888 (94%)	56 (6%)	3 (0%)	36	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	417	ASP
1	B	684	ASP
1	A	719	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	404/407 (99%)	397 (98%)	7 (2%)	53	38
1	B	407/407 (100%)	401 (98%)	6 (2%)	57	43
All	All	811/814 (100%)	798 (98%)	13 (2%)	55	41

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	339	ASN
1	A	561	ASP
1	A	645	ASN
1	A	655	ASN
1	A	715	ASN
1	A	778	THR
1	A	785	ASN
1	B	339	ASN
1	B	467	GLU
1	B	645	ASN
1	B	655	ASN
1	B	715	ASN
1	B	785	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	339	ASN
1	A	415	ASN
1	A	478	ASN
1	A	517	ASN
1	A	523	ASN
1	A	594	ASN
1	A	623	ASN
1	A	655	ASN
1	A	715	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	753	ASN
1	A	774	GLN
1	A	775	ASN
1	A	785	ASN
1	B	339	ASN
1	B	386	ASN
1	B	517	ASN
1	B	523	ASN
1	B	623	ASN
1	B	627	GLN
1	B	655	ASN
1	B	715	ASN
1	B	753	ASN
1	B	774	GLN
1	B	775	ASN
1	B	785	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](#)

2 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	DAN	A	900	-	20,20,20	3.03	4 (20%)	24,28,28	1.69	2 (8%)
2	DAN	B	901	-	20,20,20	3.05	4 (20%)	24,28,28	1.59	3 (12%)

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	DAN	A	900	-	-	0/18/34/34	0/1/1/1
2	DAN	B	901	-	-	0/18/34/34	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	DAN	C3-C2	11.41	1.53	1.33
2	A	900	DAN	C3-C2	11.18	1.52	1.33
2	B	901	DAN	O10-C10	5.99	1.36	1.23
2	A	900	DAN	O10-C10	5.97	1.36	1.23
2	A	900	DAN	C4-C3	2.55	1.53	1.50
2	B	901	DAN	C4-C3	2.36	1.53	1.50
2	A	900	DAN	O1B-C1	-2.14	1.24	1.30
2	B	901	DAN	C2-C1	2.01	1.53	1.48

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	DAN	C4-C3-C2	-6.44	110.55	121.45
2	B	901	DAN	C4-C3-C2	-5.25	112.57	121.45
2	B	901	DAN	O6-C2-C1	3.16	118.32	112.02
2	A	900	DAN	O6-C2-C1	2.98	117.97	112.02
2	B	901	DAN	C3-C2-C1	-2.16	118.89	123.56

There are no chirality outliers.

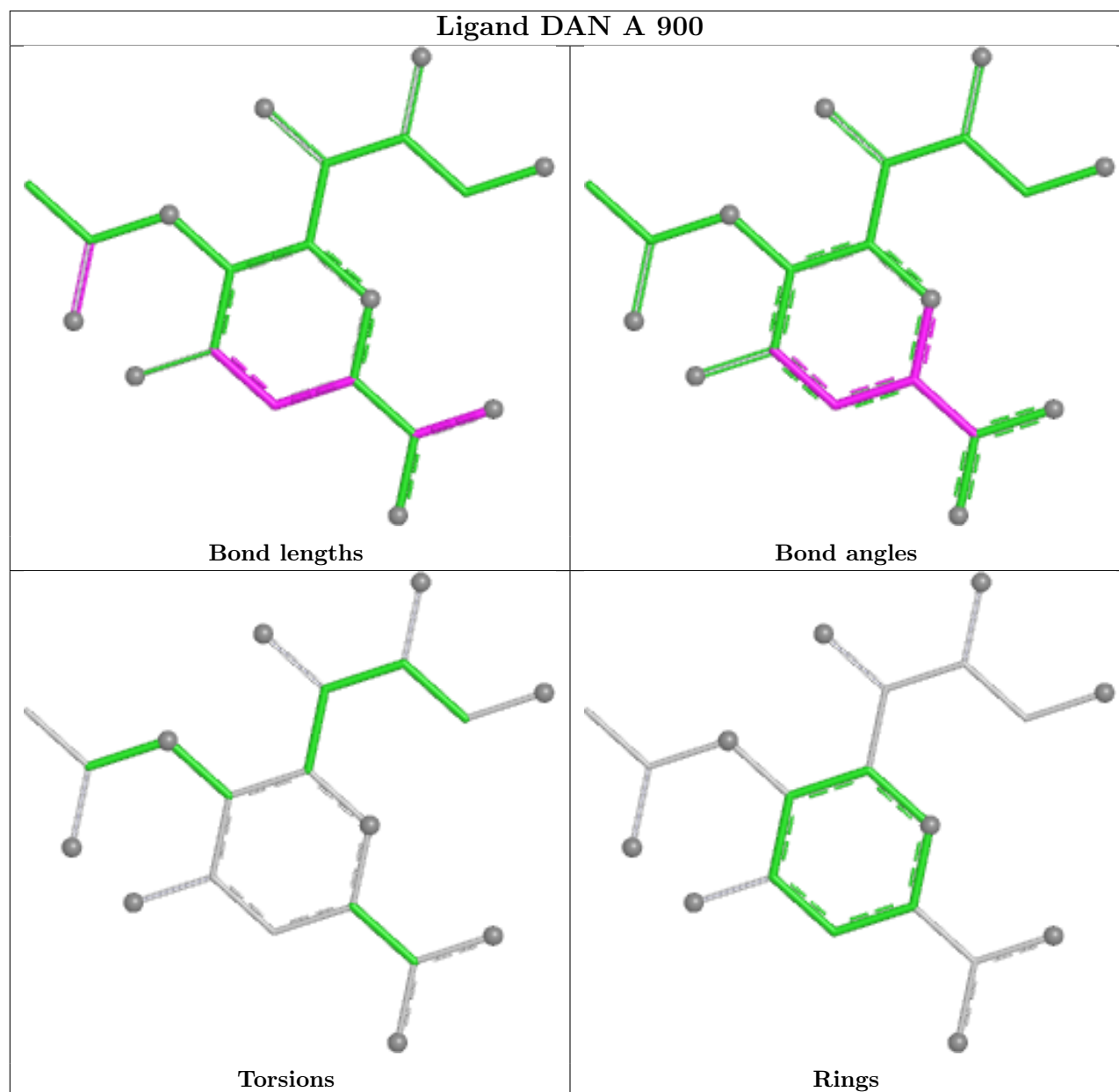
There are no torsion outliers.

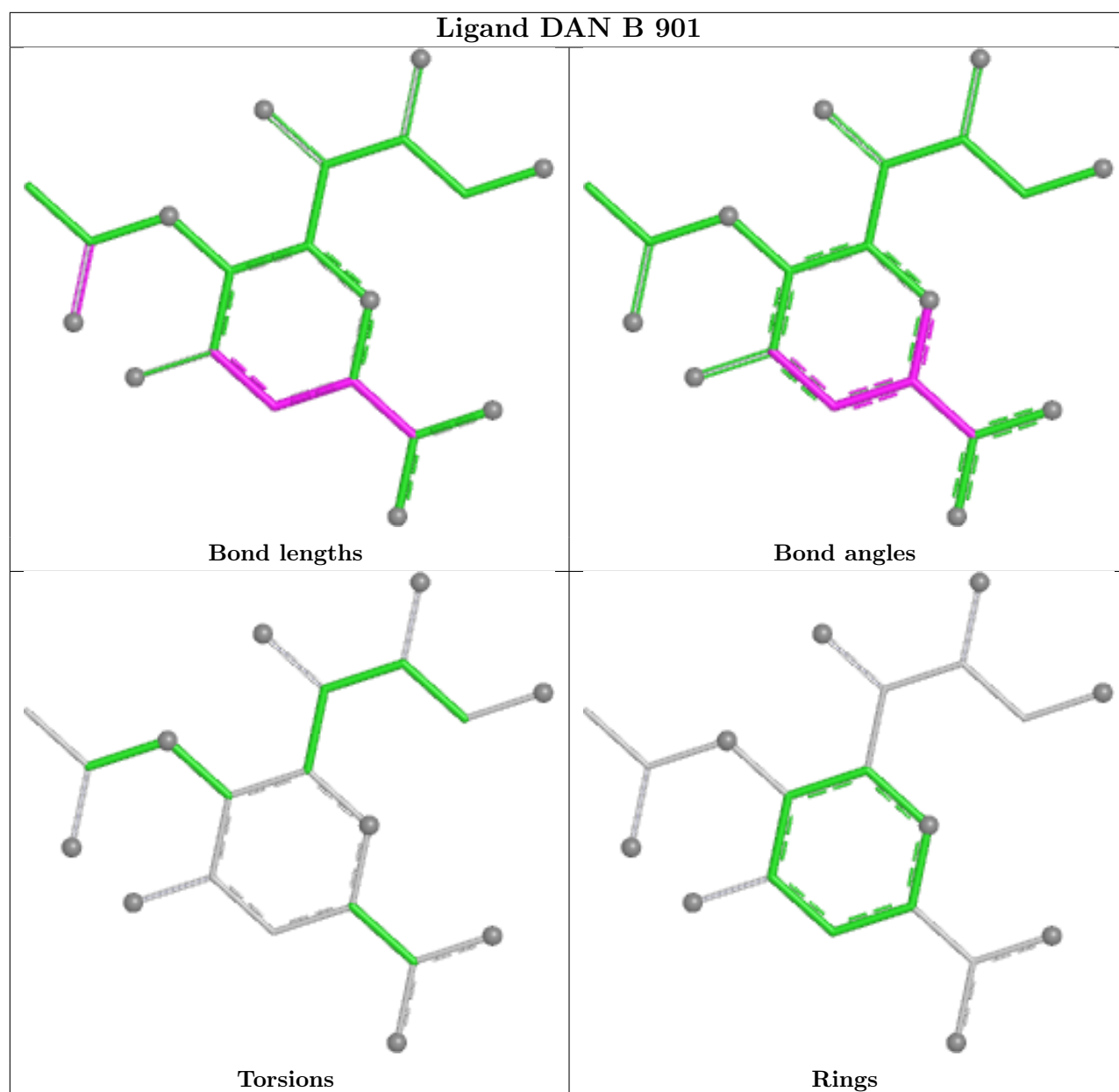
There are no ring outliers.

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	474/477 (99%)	-0.05	22 (4%) 37 41	3, 9, 24, 44	0
1	B	477/477 (100%)	0.03	19 (3%) 42 46	4, 10, 24, 35	0
All	All	951/954 (99%)	-0.01	41 (4%) 40 44	3, 9, 24, 44	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	321	ALA	9.7
1	A	793	LEU	6.3
1	B	793	LEU	4.4
1	A	320	GLY	4.4
1	B	511	GLY	4.2
1	B	317	LEU	3.8
1	A	684	ASP	3.7
1	A	458	GLY	3.7
1	B	630	GLY	3.6
1	A	631	GLN	3.4
1	A	790	SER	3.4
1	A	512	ASN	3.2
1	A	467	GLU	3.2
1	A	792	ASP	3.2
1	B	319	GLU	3.1
1	A	511	GLY	3.0
1	B	735	GLY	3.0
1	A	322	ALA	3.0
1	B	760	ASN	2.9
1	B	485	ASP	2.8
1	B	492	ARG	2.7
1	A	456	ILE	2.7
1	B	685	ILE	2.7
1	B	733	GLU	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	512	ASN	2.6
1	B	467	GLU	2.6
1	A	457	ASP	2.6
1	A	485	ASP	2.5
1	A	459	LYS	2.5
1	A	772	LYS	2.5
1	A	686	LYS	2.3
1	B	325	GLU	2.2
1	B	614	LYS	2.2
1	A	325	GLU	2.2
1	A	412	SER	2.2
1	B	490	ASP	2.1
1	A	760	ASN	2.1
1	B	734	ASN	2.1
1	A	629	ASP	2.1
1	B	510	LYS	2.1
1	B	772	LYS	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 oligosaccharides in this entry.

6.4 Ligands ⓘ

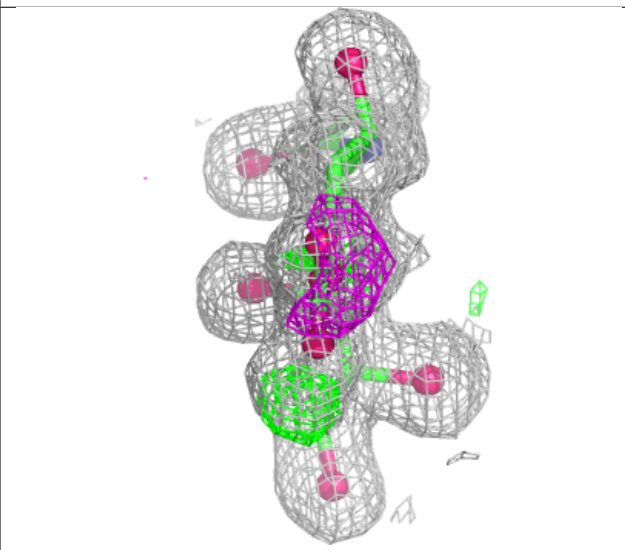
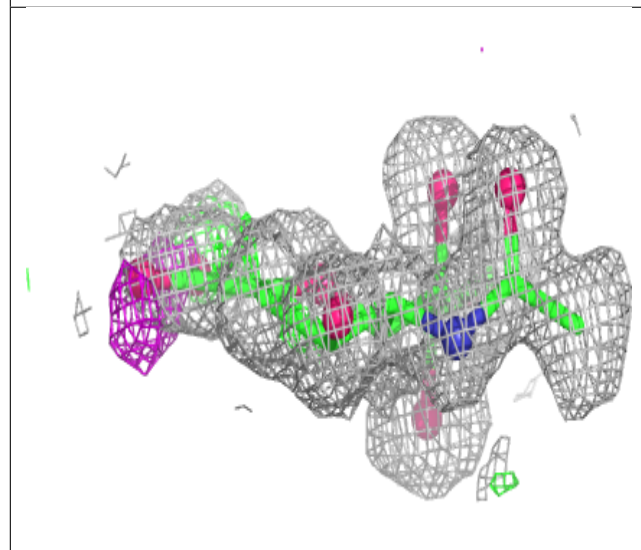
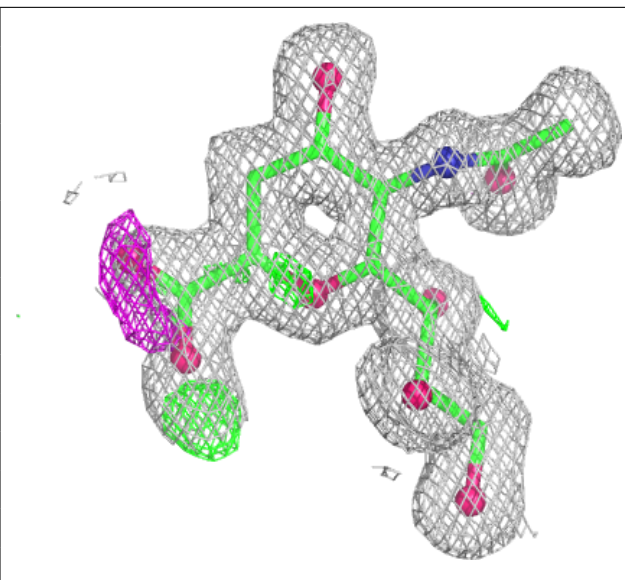
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	DAN	B	901	20/20	0.91	0.08	5,8,20,20	0
2	DAN	A	900	20/20	0.94	0.07	4,6,17,18	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.

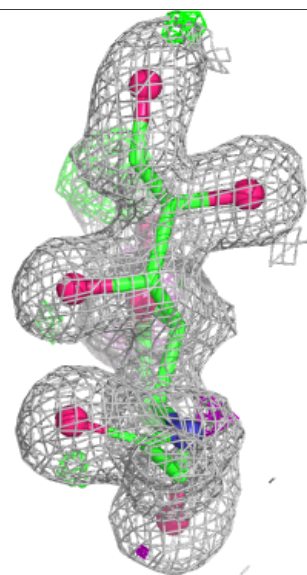
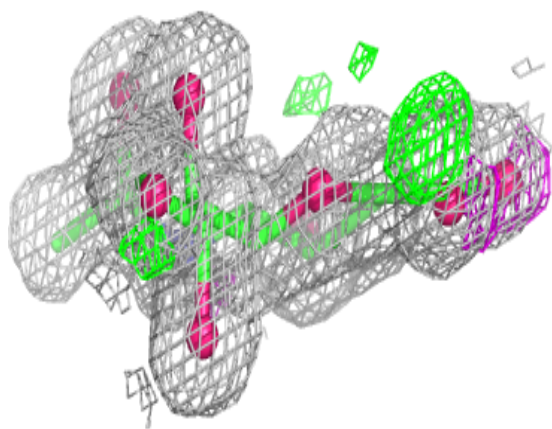
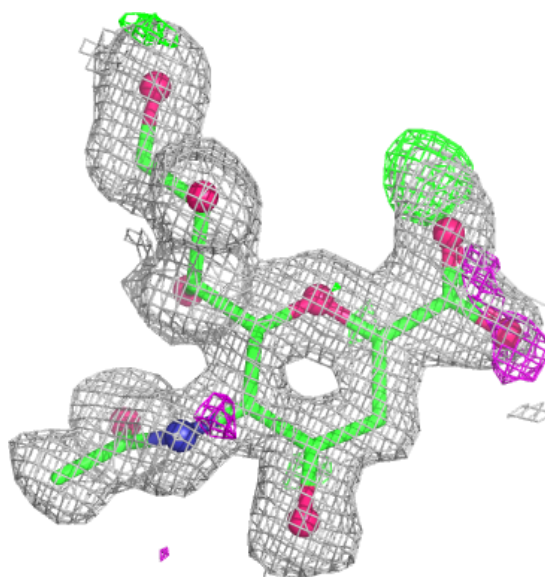
Electron density around DAN B 901:

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



Electron density around DAN A 900:

$2mF_o-DF_c$ (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.