



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 08:54 PM UTC

PDB ID : 2XVW / pdb_00002xvw
Title : Human serum albumin complexed with dansyl-L-arginine and myristic acid
Authors : Ghuman, J.; Curry, S.
Deposited on : 2010-10-28
Resolution : 2.65 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

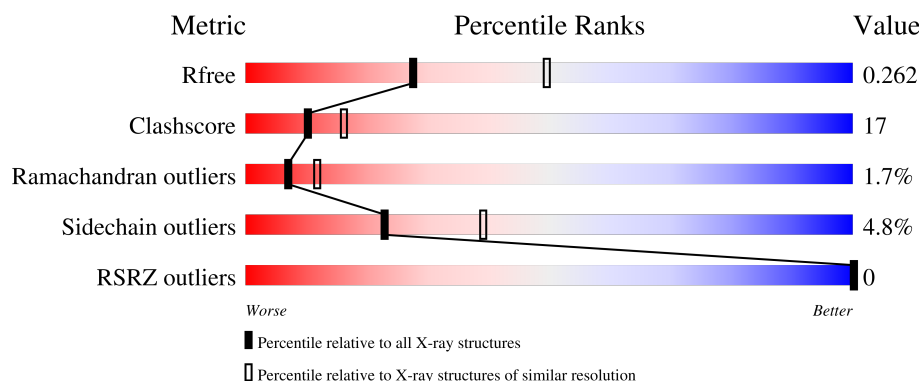
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	585	 64% 31% . .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4630 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.

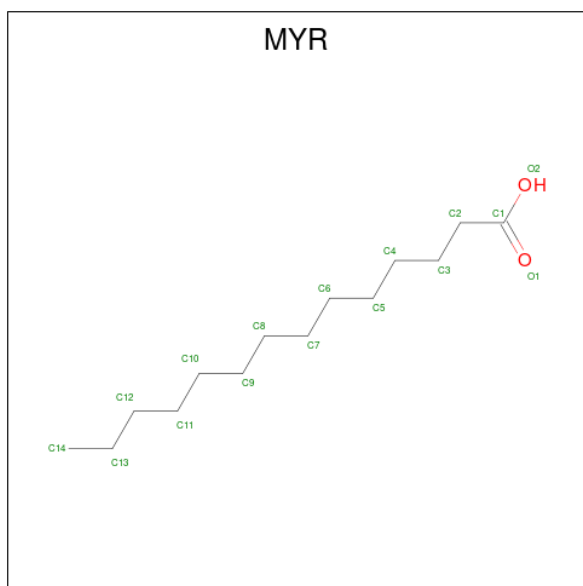
- Molecule 1 is a protein called SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	582	Total	C	N	O	S	0	0	0
			4465	2831	746	847	41			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	521	GLU	ARG	conflict	UNP P02768

- Molecule 2 is MYRISTIC ACID (CCD ID: MYR) (formula: C₁₄H₂₈O₂).



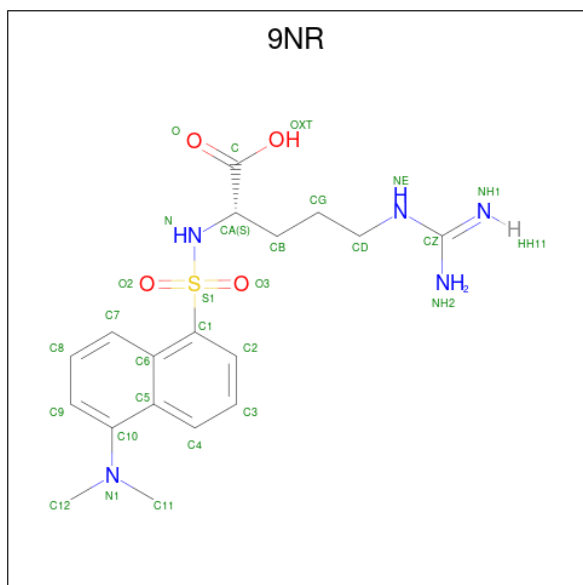
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			14	12	2		
2	A	1	Total	C	O	0	0
			14	12	2		
2	A	1	Total	C	O	0	0
			13	11	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	11	2		
2	A	1	Total	C	O	0	0
			16	14	2		
2	A	1	Total	C	O	0	0
			16	14	2		

- Molecule 3 is DANSYL-L-ARGININE (CCD ID: 9NR) (formula: $C_{18}H_{25}N_5O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			28	18	5	4	1		
3	A	1	Total	C	N	O	S	12	0
			28	18	5	4	1		

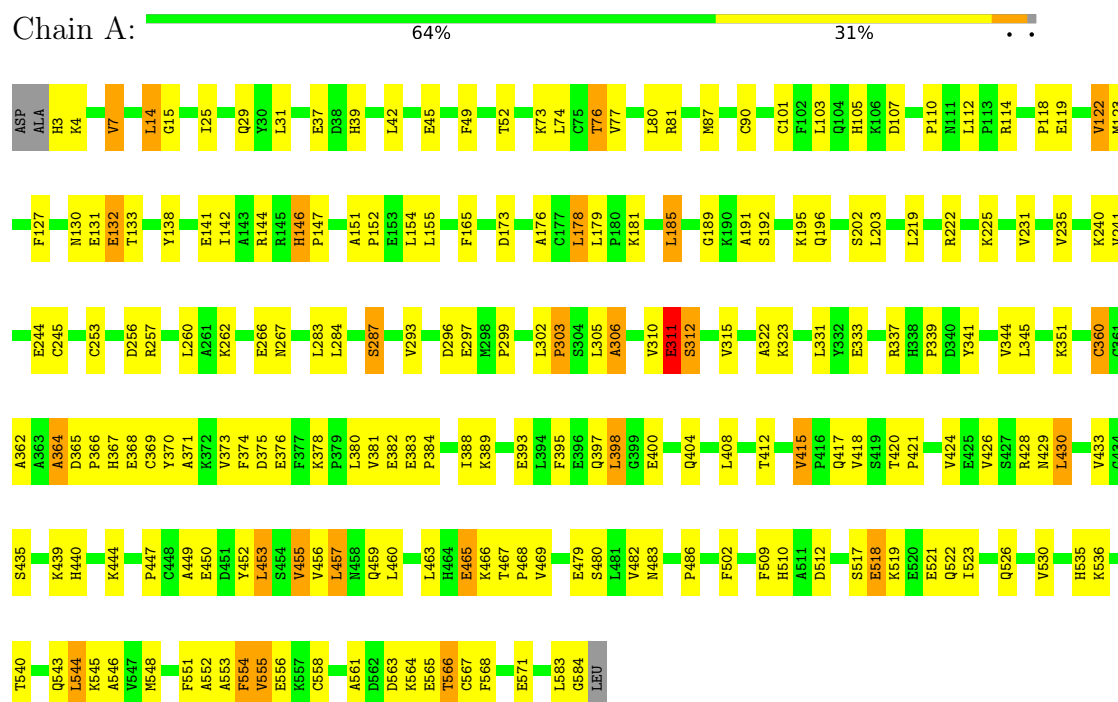
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	23	Total	O	0	0
			23	23		

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: SERUM ALBUMIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	187.08Å 38.76Å 95.73Å 90.00° 105.64° 90.00°	Depositor
Resolution (Å)	32.57 – 2.65 32.57 – 2.65	Depositor EDS
% Data completeness (in resolution range)	96.7 (32.57-2.65) 96.7 (32.57-2.65)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.64Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.191 , 0.252 0.207 , 0.262	Depositor DCC
R_{free} test set	913 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	57.6	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4630	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.89% 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: MYR, 9NR

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	0.48	0/4555	0.95	18/6177 (0.3%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	360	CYS	N-CA-C	9.30	121.02	111.07
1	A	14	LEU	N-CA-C	7.11	118.67	111.07
1	A	178	LEU	N-CA-C	6.50	118.25	111.03
1	A	339	PRO	N-CA-C	-6.36	105.91	113.86
1	A	415	VAL	CA-C-N	6.19	125.81	119.56
1	A	415	VAL	C-N-CA	6.19	125.81	119.56
1	A	536	LYS	CA-C-N	6.07	126.18	119.87
1	A	536	LYS	C-N-CA	6.07	126.18	119.87
1	A	132	GLU	N-CA-C	5.96	117.45	111.07
1	A	15	GLY	N-CA-C	-5.53	106.09	112.29
1	A	146	HIS	CA-C-N	5.40	126.59	119.84
1	A	146	HIS	C-N-CA	5.40	126.59	119.84
1	A	323	LYS	N-CA-C	5.32	117.51	111.02
1	A	465	GLU	N-CA-C	-5.24	105.46	111.07
1	A	412	THR	N-CA-C	-5.14	105.67	111.28
1	A	112	LEU	CA-C-N	5.04	124.80	119.76
1	A	112	LEU	C-N-CA	5.04	124.80	119.76
1	A	154	LEU	N-CA-C	-5.00	105.91	111.36

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	4465	0	4206	153	0
2	A	86	0	130	3	0
3	A	56	0	46	3	0
4	A	23	0	0	1	0
All	All	4630	0	4382	153	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 17.

All (153) 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:480:SER:HB3	1:A:483:ASN:HD22	1.23	0.98
1:A:424:VAL:O	1:A:428:ARG:HG3	1.70	0.91
1:A:305:LEU:HD22	1:A:337:ARG:HH11	1.36	0.88
1:A:257:ARG:HH21	1:A:287:SER:HB3	1.36	0.88
1:A:118:PRO:HD2	1:A:123:MET:CE	2.08	0.84
1:A:37:GLU:H	1:A:37:GLU:CD	1.94	0.75
1:A:479:GLU:HG3	1:A:480:SER:H	1.53	0.74
1:A:305:LEU:HD22	1:A:337:ARG:NH1	2.03	0.73
1:A:222:ARG:HD3	1:A:293:VAL:HG12	1.72	0.72
1:A:364:ALA:O	1:A:366:PRO:HD3	1.89	0.72
1:A:118:PRO:HD2	1:A:123:MET:HE1	1.70	0.71
1:A:510:HIS:HB2	1:A:512:ASP:OD1	1.92	0.69
1:A:118:PRO:HD2	1:A:123:MET:HE3	1.73	0.69
1:A:119:GLU:HB3	1:A:122:VAL:HG13	1.75	0.68
1:A:449:ALA:O	1:A:453:LEU:HB2	1.95	0.66
1:A:389:LYS:O	1:A:393:GLU:HG3	1.95	0.66
1:A:305:LEU:CD2	1:A:337:ARG:HH11	2.08	0.66
1:A:222:ARG:NH1	1:A:293:VAL:O	2.29	0.65
1:A:479:GLU:HG3	1:A:480:SER:N	2.10	0.65
1:A:563:ASP:CG	1:A:566:THR:HB	2.21	0.64
1:A:262:LYS:O	1:A:266:GLU:HG3	1.97	0.64
1:A:467:THR:O	1:A:469:VAL:HG23	1.97	0.64
1:A:551:PHE:O	1:A:554:PHE:HB3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LYS:O	1:A:76:THR:HB	1.98	0.62
1:A:367:HIS:O	1:A:371:ALA:HB2	1.99	0.61
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.84	0.60
1:A:81:ARG:HB2	1:A:81:ARG:NH1	2.16	0.60
1:A:558:CYS:HB3	1:A:568:PHE:CE1	2.36	0.59
1:A:39:HIS:HA	1:A:42:LEU:HD12	1.85	0.59
1:A:365:ASP:OD2	1:A:368:GLU:HB2	2.03	0.59
1:A:260:LEU:HD23	3:A:2001:9NR:H122	1.84	0.59
1:A:420:THR:O	1:A:424:VAL:HG23	2.02	0.59
1:A:191:ALA:O	1:A:195:LYS:HG3	2.04	0.58
1:A:306:ALA:HA	1:A:310:VAL:HG23	1.84	0.58
1:A:77:VAL:O	1:A:80:LEU:HB2	2.04	0.57
1:A:518:GLU:HA	1:A:518:GLU:OE1	2.04	0.57
1:A:267:ASN:HD22	1:A:267:ASN:N	2.02	0.57
1:A:561:ALA:H	1:A:564:LYS:HZ2	1.53	0.57
1:A:296:ASP:OD1	1:A:297:GLU:N	2.39	0.56
1:A:305:LEU:HD11	1:A:333:GLU:HB3	1.87	0.56
1:A:519:LYS:O	1:A:523:ILE:HG13	2.05	0.56
1:A:310:VAL:O	1:A:370:TYR:HE1	1.88	0.56
1:A:566:THR:HG22	1:A:567:CYS:N	2.21	0.56
1:A:426:VAL:HG21	1:A:460:LEU:HD13	1.88	0.55
1:A:565:GLU:HA	1:A:565:GLU:OE1	2.06	0.55
1:A:408:LEU:HD21	1:A:526:GLN:HB3	1.88	0.55
1:A:417:GLN:HB3	1:A:469:VAL:HG12	1.88	0.55
1:A:222:ARG:HD3	1:A:293:VAL:CG1	2.34	0.55
1:A:420:THR:HG23	1:A:530:VAL:HB	1.87	0.55
1:A:417:GLN:HB3	1:A:469:VAL:CG1	2.37	0.55
1:A:311:GLU:O	1:A:312:SER:O	2.25	0.54
1:A:452:TYR:O	1:A:455:VAL:HG13	2.08	0.54
1:A:480:SER:HB3	1:A:483:ASN:ND2	2.08	0.54
1:A:517:SER:O	1:A:521:GLU:HG3	2.08	0.53
1:A:430:LEU:HD21	2:A:1004:MYR:H61	1.89	0.53
1:A:231:VAL:O	1:A:235:VAL:HG23	2.09	0.53
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.39	0.52
1:A:553:ALA:O	1:A:554:PHE:C	2.51	0.52
1:A:561:ALA:N	1:A:564:LYS:NZ	2.57	0.52
1:A:311:GLU:O	1:A:312:SER:C	2.53	0.52
1:A:420:THR:N	1:A:421:PRO:HD2	2.25	0.52
1:A:551:PHE:CD2	2:A:1005:MYR:H52	2.44	0.52
1:A:522:GLN:O	1:A:526:GLN:HG3	2.10	0.52
1:A:101:CYS:O	1:A:105:HIS:HD2	1.94	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:VAL:HG12	1:A:450:GLU:OE2	2.10	0.51
1:A:344:VAL:HG23	1:A:482:VAL:HG13	1.92	0.51
1:A:7:VAL:HG22	4:A:2006:HOH:O	2.09	0.51
1:A:123:MET:HG2	1:A:165:PHE:CE1	2.46	0.51
1:A:283:LEU:C	1:A:283:LEU:HD13	2.36	0.51
1:A:459:GLN:O	1:A:463:LEU:HG	2.11	0.50
1:A:563:ASP:OD1	1:A:566:THR:HB	2.12	0.49
1:A:240:LYS:HE2	1:A:244:GLU:OE2	2.12	0.49
1:A:37:GLU:CD	1:A:37:GLU:N	2.69	0.49
1:A:127:PHE:CE1	1:A:131:GLU:HG3	2.48	0.49
1:A:540:THR:OG1	1:A:543:GLN:HG3	2.13	0.49
1:A:543:GLN:O	1:A:546:ALA:HB3	2.13	0.49
1:A:267:ASN:N	1:A:267:ASN:ND2	2.61	0.48
1:A:299:PRO:HB2	1:A:302:LEU:HG	1.94	0.48
1:A:123:MET:HG2	1:A:165:PHE:HE1	1.77	0.48
1:A:266:GLU:C	1:A:267:ASN:HD22	2.21	0.48
1:A:398:LEU:N	1:A:398:LEU:HD23	2.28	0.48
1:A:502:PHE:HB2	1:A:535:HIS:CE1	2.48	0.48
1:A:25:ILE:O	1:A:29:GLN:HG3	2.13	0.48
1:A:107:ASP:OD2	1:A:110:PRO:HA	2.14	0.47
1:A:165:PHE:CE1	1:A:178:LEU:HD21	2.50	0.47
1:A:465:GLU:O	1:A:468:PRO:HD3	2.13	0.47
1:A:49:PHE:O	1:A:52:THR:HB	2.14	0.47
1:A:181:LYS:O	1:A:185:LEU:HD22	2.13	0.47
1:A:141:GLU:HA	1:A:144:ARG:HD3	1.97	0.47
1:A:219:LEU:HD12	1:A:235:VAL:HG22	1.96	0.47
1:A:395:PHE:CE2	1:A:435:SER:HB3	2.50	0.47
1:A:155:LEU:HD23	1:A:284:LEU:HD21	1.96	0.47
1:A:119:GLU:HB3	1:A:122:VAL:CG1	2.44	0.47
1:A:545:LYS:HA	1:A:548:MET:HE3	1.97	0.47
1:A:555:VAL:O	1:A:556:GLU:C	2.57	0.47
1:A:509:PHE:O	1:A:568:PHE:HB3	2.15	0.46
1:A:439:LYS:HD3	1:A:439:LYS:HA	1.65	0.46
1:A:341:TYR:CG	1:A:345:LEU:HD23	2.50	0.46
1:A:469:VAL:HG12	1:A:469:VAL:O	2.14	0.46
1:A:554:PHE:HE1	1:A:571:GLU:HB2	1.80	0.45
1:A:551:PHE:O	1:A:552:ALA:C	2.58	0.45
1:A:429:ASN:O	1:A:433:VAL:HG23	2.16	0.45
1:A:87:MET:O	1:A:90:CYS:HB2	2.16	0.45
1:A:225:LYS:HG2	1:A:299:PRO:HG3	1.99	0.45
1:A:483:ASN:C	1:A:486:PRO:HD2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:VAL:O	1:A:384:PRO:HG2	2.16	0.45
1:A:415:VAL:HG12	1:A:418:VAL:HG23	1.98	0.45
1:A:305:LEU:CD2	1:A:337:ARG:NH1	2.76	0.45
1:A:305:LEU:CD1	1:A:333:GLU:HB3	2.47	0.45
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.99	0.45
1:A:373:VAL:C	1:A:375:ASP:N	2.75	0.44
1:A:245:CYS:HA	1:A:253:CYS:HB2	1.99	0.44
1:A:400:GLU:O	1:A:404:GLN:HG3	2.17	0.44
1:A:440:HIS:N	1:A:440:HIS:CD2	2.84	0.44
1:A:395:PHE:O	1:A:398:LEU:O	2.36	0.44
1:A:417:GLN:CB	1:A:469:VAL:HG12	2.48	0.44
1:A:373:VAL:HG13	1:A:374:PHE:N	2.32	0.44
1:A:114:ARG:O	1:A:114:ARG:HG3	2.16	0.43
1:A:360:CYS:SG	1:A:369:CYS:C	3.02	0.43
1:A:203:LEU:O	1:A:203:LEU:HD23	2.17	0.43
1:A:583:LEU:O	1:A:584:GLY:C	2.61	0.43
1:A:452:TYR:O	1:A:456:VAL:HG23	2.18	0.43
1:A:241:VAL:HG22	1:A:256:ASP:HB3	2.01	0.42
1:A:558:CYS:HB3	1:A:568:PHE:CD1	2.54	0.42
1:A:138:TYR:O	1:A:142:ILE:HG12	2.20	0.42
1:A:518:GLU:O	1:A:519:LYS:C	2.61	0.42
1:A:544:LEU:O	1:A:545:LYS:C	2.62	0.42
1:A:189:GLY:HA3	3:A:2002:9NR:N1	2.35	0.42
1:A:373:VAL:O	1:A:376:GLU:N	2.48	0.42
1:A:173:ASP:OD2	1:A:176:ALA:HB2	2.20	0.42
1:A:360:CYS:C	1:A:362:ALA:H	2.28	0.42
1:A:517:SER:O	1:A:518:GLU:C	2.61	0.42
1:A:130:ASN:OD1	1:A:133:THR:HG23	2.20	0.42
1:A:370:TYR:CD1	1:A:370:TYR:C	2.98	0.42
1:A:566:THR:CG2	1:A:567:CYS:N	2.82	0.42
1:A:444:LYS:C	1:A:447:PRO:HD2	2.44	0.41
1:A:101:CYS:O	1:A:105:HIS:CD2	2.73	0.41
1:A:561:ALA:H	1:A:564:LYS:NZ	2.14	0.41
1:A:3:HIS:HB3	1:A:4:LYS:H	1.44	0.41
1:A:31:LEU:HD13	1:A:74:LEU:HD11	2.03	0.41
1:A:378:LYS:O	1:A:382:GLU:HG3	2.21	0.41
1:A:222:ARG:NH2	3:A:2001:9NR:OXT	2.53	0.41
1:A:415:VAL:CG1	1:A:418:VAL:HG23	2.50	0.41
1:A:45:GLU:OE1	1:A:73:LYS:NZ	2.39	0.41
1:A:457:LEU:HD12	1:A:457:LEU:HA	1.93	0.41
1:A:192:SER:O	1:A:196:GLN:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:LEU:HA	1:A:303:PRO:HD3	1.92	0.41
1:A:388:ILE:HD13	2:A:1003:MYR:H72	2.03	0.41
1:A:395:PHE:CD1	1:A:395:PHE:C	2.99	0.41
1:A:466:LYS:C	1:A:468:PRO:HD3	2.46	0.41
1:A:146:HIS:N	1:A:147:PRO:HD3	2.36	0.41
1:A:305:LEU:HD13	1:A:337:ARG:NH1	2.35	0.40
1:A:262:LYS:HE3	1:A:262:LYS:HB2	1.94	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	580/585 (99%)	524 (90%)	46 (8%)	10 (2%)	7 12

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	312	SER
1	A	306	ALA
1	A	311	GLU
1	A	397	GLN
1	A	555	VAL
1	A	322	ALA
1	A	554	PHE
1	A	364	ALA
1	A	303	PRO
1	A	315	VAL

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	457/511 (89%)	435 (95%)	22 (5%)	23	39

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	14	LEU
1	A	76	THR
1	A	103	LEU
1	A	122	VAL
1	A	132	GLU
1	A	179	LEU
1	A	185	LEU
1	A	202	SER
1	A	287	SER
1	A	311	GLU
1	A	331	LEU
1	A	351	LYS
1	A	380	LEU
1	A	398	LEU
1	A	430	LEU
1	A	453	LEU
1	A	455	VAL
1	A	457	LEU
1	A	518	GLU
1	A	544	LEU
1	A	566	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	39	HIS
1	A	105	HIS

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Mol	Chain	Res	Type
1	A	221	GLN
1	A	267	ASN
1	A	318	ASN
1	A	397	GLN
1	A	404	GLN
1	A	459	GLN
1	A	483	ASN
1	A	510	HIS

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](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	9NR	A	2002	-	29,29,29	1.90	11 (37%)	40,41,41	1.36	6 (15%)
2	MYR	A	1004	-	12,12,15	0.73	0	12,12,15	1.02	1 (8%)
2	MYR	A	1005	-	15,15,15	0.61	0	15,15,15	1.00	1 (6%)
2	MYR	A	1006	-	15,15,15	0.64	0	15,15,15	1.02	1 (6%)
2	MYR	A	1001	-	13,13,15	0.70	0	13,13,15	1.01	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	9NR	A	2001	-	29,29,29	1.68	8 (27%)	40,41,41	0.97	2 (5%)
2	MYR	A	1003	-	12,12,15	0.67	0	12,12,15	1.09	1 (8%)
2	MYR	A	1002	-	13,13,15	0.74	0	13,13,15	1.01	1 (7%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	9NR	A	2002	-	-	9/26/26/26	0/2/2/2
2	MYR	A	1004	-	-	7/10/10/13	-
2	MYR	A	1005	-	-	9/13/13/13	-
2	MYR	A	1006	-	-	11/13/13/13	-
2	MYR	A	1001	-	-	10/11/11/13	-
3	9NR	A	2001	-	-	4/26/26/26	0/2/2/2
2	MYR	A	1003	-	-	6/10/10/13	-
2	MYR	A	1002	-	-	7/11/11/13	-

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2002	9NR	C1-S1	3.71	1.81	1.77
3	A	2001	9NR	C1-C6	3.64	1.49	1.43
3	A	2001	9NR	O-C	3.54	1.32	1.22
3	A	2002	9NR	C1-C6	3.36	1.49	1.43
3	A	2001	9NR	OXT-C	-2.99	1.21	1.30
3	A	2002	9NR	C2-C1	2.89	1.41	1.37
3	A	2001	9NR	C6-C5	2.81	1.48	1.43
3	A	2002	9NR	C3-C2	2.80	1.43	1.38
3	A	2002	9NR	O-C	2.79	1.30	1.22
3	A	2002	9NR	C6-C5	2.70	1.48	1.43
3	A	2002	9NR	C10-C5	2.56	1.48	1.42
3	A	2002	9NR	OXT-C	-2.54	1.22	1.30
3	A	2001	9NR	C12-N1	2.54	1.51	1.45
3	A	2001	9NR	C10-C5	2.49	1.48	1.42
3	A	2001	9NR	C8-C9	2.40	1.43	1.38
3	A	2002	9NR	C12-N1	2.28	1.50	1.45
3	A	2002	9NR	CB-CA	2.25	1.58	1.53
3	A	2002	9NR	C8-C9	2.24	1.42	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2001	9NR	C3-C2	2.11	1.42	1.38

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2002	9NR	O3-S1-O2	-3.80	114.90	119.52
3	A	2002	9NR	CA-N-S1	3.57	129.13	121.30
3	A	2001	9NR	C12-N1-C11	-3.20	105.90	116.18
3	A	2002	9NR	C12-N1-C11	-3.12	106.16	116.18
3	A	2002	9NR	CB-CA-N	3.03	115.67	110.21
3	A	2002	9NR	O-C-CA	-2.46	114.31	122.26
2	A	1005	MYR	C4-C3-C2	-2.45	104.14	113.13
2	A	1002	MYR	C4-C3-C2	-2.39	104.35	113.13
2	A	1006	MYR	C4-C3-C2	-2.31	104.64	113.13
3	A	2001	9NR	O-C-CA	-2.28	114.90	122.26
2	A	1001	MYR	C4-C3-C2	-2.22	104.99	113.13
2	A	1004	MYR	C4-C3-C2	-2.08	105.50	113.13
2	A	1003	MYR	C4-C3-C2	-2.03	105.67	113.13
3	A	2002	9NR	CG-CB-CA	2.01	119.98	113.80

There are no chirality outliers.

All (63) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2002	9NR	CB-CA-N-S1
3	A	2001	9NR	CA-N-S1-O3
3	A	2002	9NR	NE-CD-CG-CB
3	A	2001	9NR	CA-N-S1-C1
2	A	1005	MYR	C1-C2-C3-C4
2	A	1001	MYR	C1-C2-C3-C4
2	A	1001	MYR	C6-C7-C8-C9
2	A	1006	MYR	C4-C5-C6-C7
2	A	1004	MYR	C1-C2-C3-C4
2	A	1002	MYR	C11-C10-C9-C8
2	A	1003	MYR	C5-C6-C7-C8
2	A	1004	MYR	C2-C3-C4-C5
2	A	1006	MYR	C1-C2-C3-C4
2	A	1005	MYR	C9-C10-C11-C12
2	A	1005	MYR	C10-C11-C12-C13
2	A	1006	MYR	C5-C6-C7-C8
2	A	1006	MYR	C10-C11-C12-C13
2	A	1002	MYR	C2-C3-C4-C5

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Mol	Chain	Res	Type	Atoms
2	A	1006	MYR	C2-C3-C4-C5
2	A	1001	MYR	C3-C4-C5-C6
2	A	1005	MYR	C7-C8-C9-C10
3	A	2002	9NR	OXT-C-CA-N
2	A	1004	MYR	C11-C10-C9-C8
2	A	1004	MYR	C4-C5-C6-C7
2	A	1005	MYR	C5-C6-C7-C8
3	A	2002	9NR	CA-CB-CG-CD
2	A	1001	MYR	C2-C3-C4-C5
3	A	2002	9NR	C-CA-CB-CG
2	A	1003	MYR	C4-C5-C6-C7
2	A	1003	MYR	C1-C2-C3-C4
2	A	1006	MYR	C11-C10-C9-C8
3	A	2001	9NR	CA-CB-CG-CD
2	A	1006	MYR	C7-C8-C9-C10
2	A	1002	MYR	C4-C5-C6-C7
2	A	1004	MYR	C7-C8-C9-C10
2	A	1005	MYR	C2-C3-C4-C5
2	A	1001	MYR	C5-C6-C7-C8
2	A	1005	MYR	C6-C7-C8-C9
2	A	1002	MYR	C6-C7-C8-C9
2	A	1005	MYR	C4-C5-C6-C7
2	A	1002	MYR	C7-C8-C9-C10
2	A	1001	MYR	C7-C8-C9-C10
3	A	2002	9NR	O-C-CA-N
2	A	1003	MYR	C11-C10-C9-C8
3	A	2002	9NR	CA-N-S1-O2
2	A	1001	MYR	C4-C5-C6-C7
3	A	2001	9NR	NE-CD-CG-CB
2	A	1004	MYR	O2-C1-C2-C3
2	A	1002	MYR	O1-C1-C2-C3
2	A	1003	MYR	O1-C1-C2-C3
2	A	1004	MYR	O1-C1-C2-C3
2	A	1003	MYR	O2-C1-C2-C3
2	A	1006	MYR	C9-C10-C11-C12
2	A	1002	MYR	O2-C1-C2-C3
2	A	1006	MYR	C11-C12-C13-C14
2	A	1001	MYR	C9-C10-C11-C12
2	A	1006	MYR	O2-C1-C2-C3
2	A	1006	MYR	O1-C1-C2-C3
3	A	2002	9NR	N-CA-CB-CG
2	A	1001	MYR	O1-C1-C2-C3

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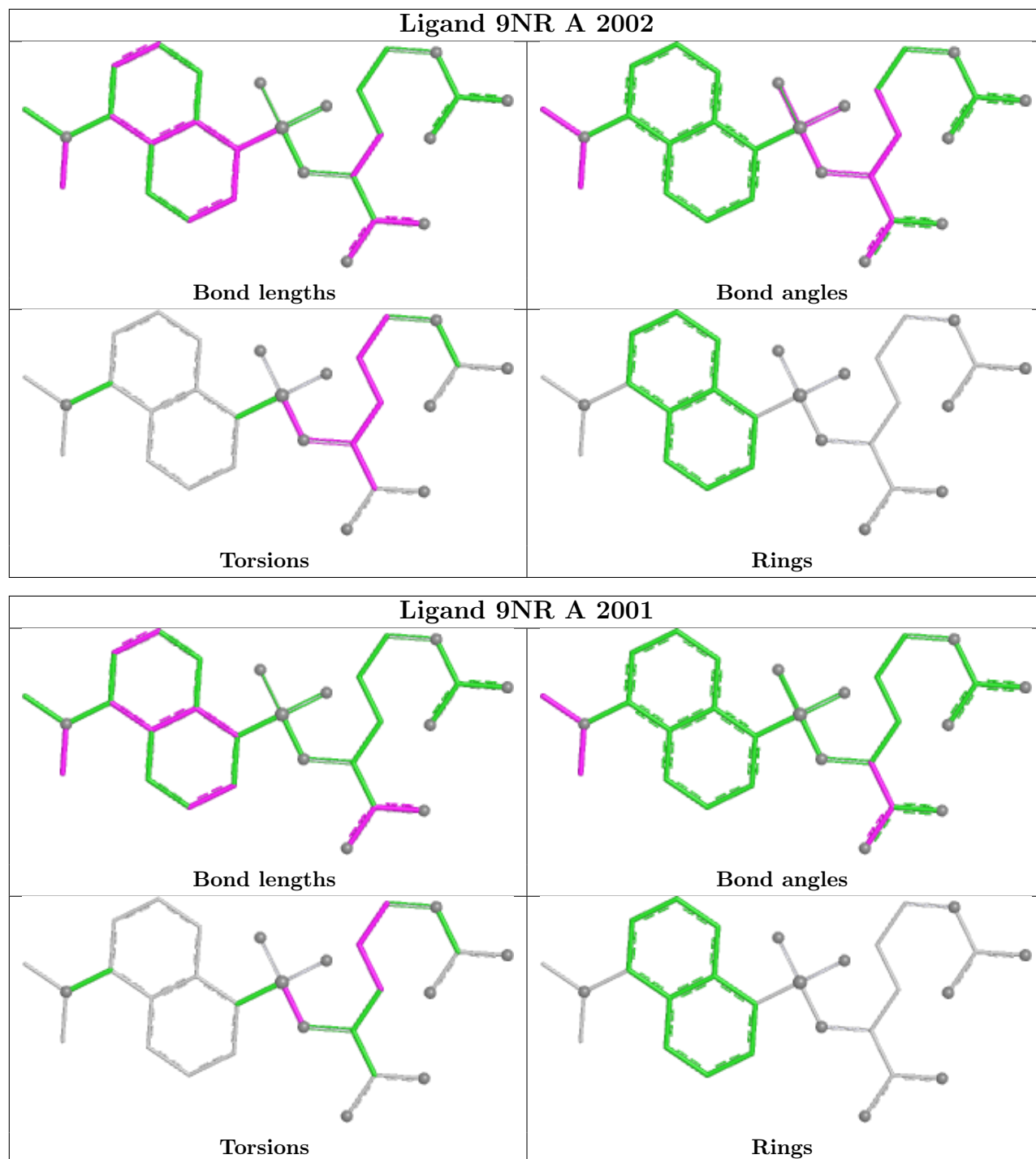
Mol	Chain	Res	Type	Atoms
2	A	1001	MYR	O2-C1-C2-C3
3	A	2002	9NR	CA-N-S1-O3
2	A	1005	MYR	O2-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2002	9NR	1	0
2	A	1004	MYR	1	0
2	A	1005	MYR	1	0
3	A	2001	9NR	2	0
2	A	1003	MYR	1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	582/585 (99%)	-0.50	0 100 100	30, 54, 80, 98	0

There are no RSRZ outliers to report.

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.

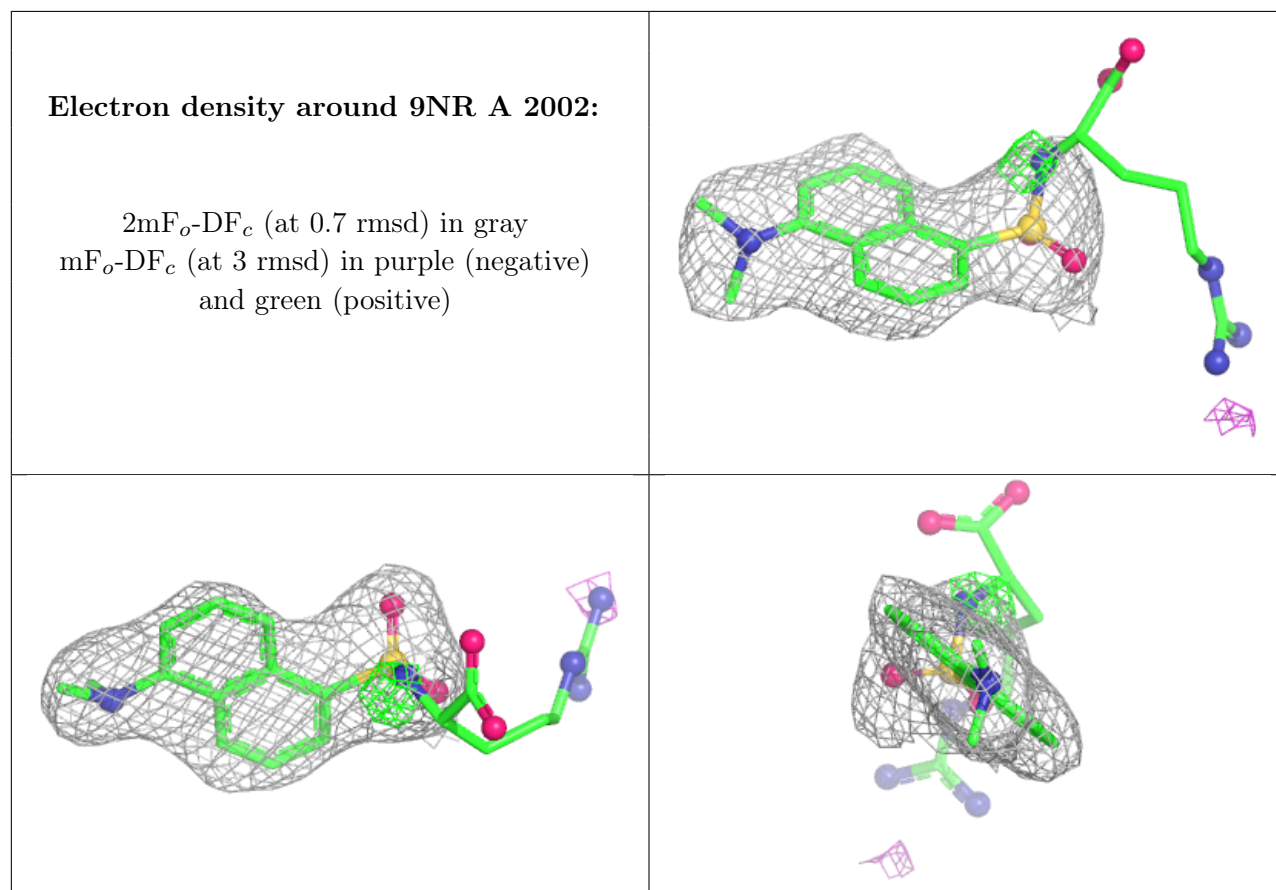
6.4 Ligands [i](#)

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	MYR	A	1002	14/16	0.89	0.13	57,59,61,62	0
2	MYR	A	1006	16/16	0.92	0.13	55,63,71,73	0
2	MYR	A	1004	13/16	0.93	0.09	51,54,57,58	0
3	9NR	A	2002	28/28	0.93	0.09	64,69,74,75	12
2	MYR	A	1003	13/16	0.94	0.11	40,46,57,58	0
3	9NR	A	2001	28/28	0.94	0.09	59,70,87,88	0
2	MYR	A	1001	14/16	0.94	0.09	44,52,56,56	0
2	MYR	A	1005	16/16	0.95	0.12	60,61,74,74	0

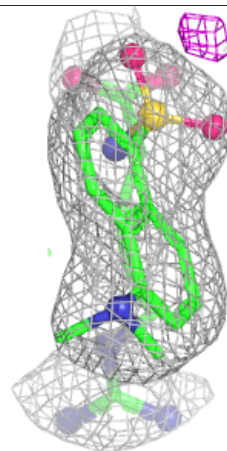
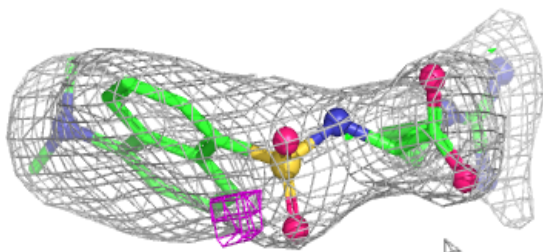
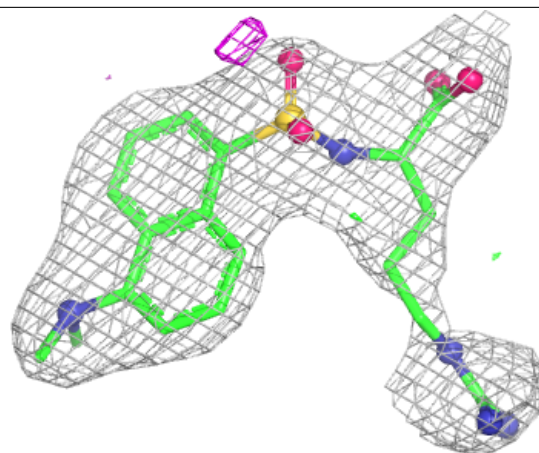
The following is a graphical depiction of the model fit to experimental electron density of all

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 9NR A 2001:

$2mF_o-DF_c$ (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.