



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 01:46 PM EDT

PDB ID : 3TGS / pdb_00003tgs
Title : Crystal structure of HIV-1 clade C strain C1086 gp120 core in complex with NBD-556
Authors : Kwon, Y.D.; Kwong, P.D.
Deposited on : 2011-08-17
Resolution : 2.70 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

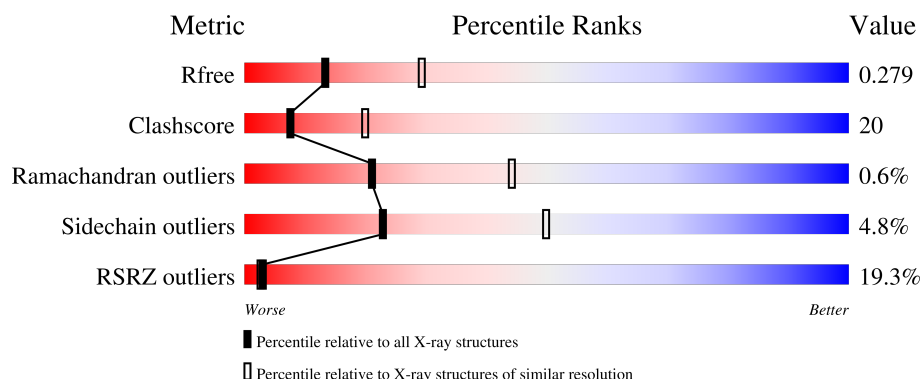
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	358	18% 59% 34% • •
1	B	358	20% 57% 37% • •

2 Entry composition [i](#)

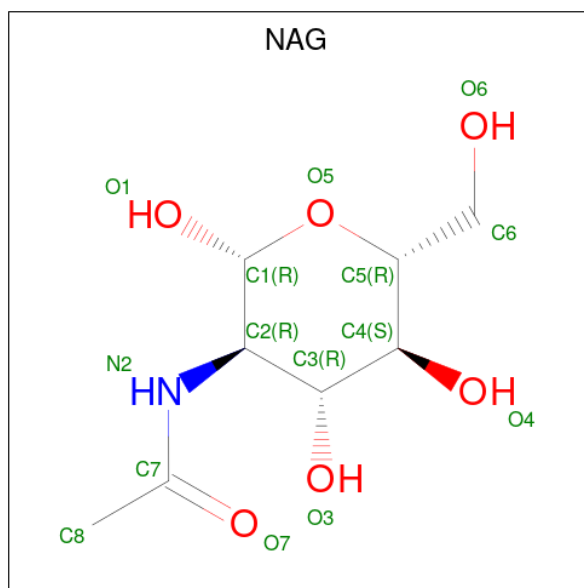
There are 4 unique types of molecules in this entry. The entry contains 5742 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HIV-1 clade C1086 gp120 core.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	345	Total	C	N	O	S	0	0	0
			2705	1692	473	520	20			
1	B	345	Total	C	N	O	S	0	0	0
			2705	1692	473	520	20			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



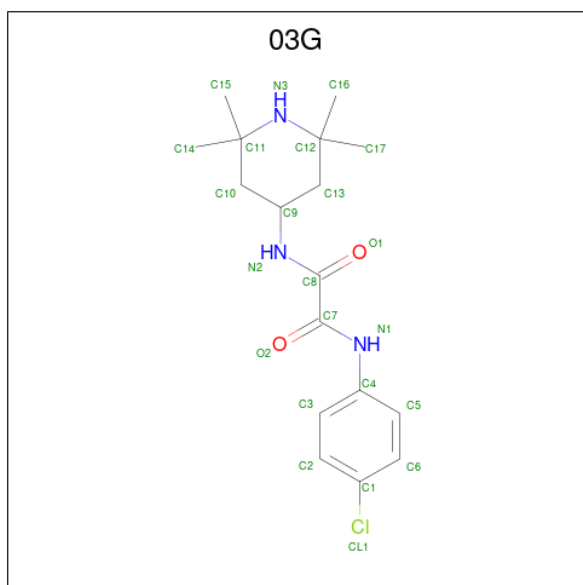
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is N-(4-chlorophenyl)-N'-(2,2,6,6-tetramethylpiperidin-4-yl)ethanediamide (CCD ID: 03G) (formula: $C_{17}H_{24}ClN_3O_2$).

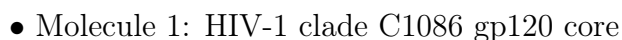


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0
			23	17	1	3	2	0
3	B	1	Total	C	Cl	N	O	0
			23	17	1	3	2	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	58	Total 58	O 58	0	0
4	B	46	Total 46	O 46	0	0

- Molecule 1: HIV-1 clade C1086 gp120 core



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	66.96Å 126.00Å 191.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.84 – 2.70 44.84 – 2.70	Depositor EDS
% Data completeness (in resolution range)	73.5 (44.84-2.70) 75.1 (44.84-2.70)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.62 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.239 , 0.295 0.229 , 0.279	Depositor DCC
R_{free} test set	861 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 60.8	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	0.88	EDS
Total number of atoms	5742	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, O3G

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	0.52	5/2762 (0.2%)	0.72	2/3746 (0.1%)
1	B	0.28	0/2762	0.71	2/3746 (0.1%)
All	All	0.42	5/5524 (0.1%)	0.72	4/7492 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	287	HIS	CG-ND1	-10.37	1.26	1.38
1	A	287	HIS	CD2-NE2	-7.48	1.29	1.37
1	A	290	GLU	C-O	-6.62	1.16	1.23
1	A	290	GLU	CA-C	-5.69	1.45	1.52
1	A	290	GLU	N-CA	-5.19	1.40	1.46

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	211	ASP	CA-C-N	5.49	125.49	119.89
1	B	211	ASP	C-N-CA	5.49	125.49	119.89
1	A	353	PHE	CA-C-N	5.17	124.62	119.24
1	A	353	PHE	C-N-CA	5.17	124.62	119.24

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	2705	0	2616	110	0
1	B	2705	0	2622	114	0
2	A	112	0	104	5	0
2	B	70	0	65	4	0
3	A	23	0	24	7	0
3	B	23	0	24	5	0
4	A	58	0	0	4	0
4	B	46	0	0	3	0
All	All	5742	0	5455	224	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 20.

All (224) 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:64:GLU:HG2	1:A:209:SER:HB3	1.49	0.94
1:A:289:ASN:OD1	1:A:290:GLU:HB2	1.75	0.86
1:A:347:GLU:HG2	4:A:653:HOH:O	1.86	0.75
1:A:104:MET:O	1:A:108:ILE:HG12	1.87	0.73
1:A:335:GLU:OE2	1:A:408:ARG:HB3	1.88	0.72
1:B:405:HIS:CG	1:B:406:THR:H	2.04	0.72
1:B:104:MET:O	1:B:108:ILE:HG12	1.91	0.71
1:A:277:LEU:HD12	2:A:501:NAG:H81	1.74	0.70
1:A:405:HIS:CG	1:A:406:THR:H	2.11	0.69
1:A:325:ASN:HB3	1:A:328:GLN:HB2	1.73	0.69
1:B:411:ASN:HA	4:B:623:HOH:O	1.94	0.67
1:B:279:ASN:HD21	1:B:281:ALA:HB3	1.60	0.66
1:B:395:TYR:HD1	1:B:404:ASN:H	1.43	0.65
1:B:476:ARG:O	1:B:480:ARG:HG3	1.97	0.64
1:A:105:HIS:HD1	1:A:479:TRP:HZ3	1.45	0.64
1:A:456:ARG:NH1	1:A:458:GLY:HA2	2.12	0.64
1:B:395:TYR:CD1	1:B:403:TYR:HA	2.34	0.63
1:A:93:PHE:HB2	1:A:233:PHE:CZ	2.34	0.63
1:B:226:LEU:HD13	1:B:489:VAL:HG11	1.82	0.62
1:A:61:TYR:CE1	1:B:214:PRO:HG2	2.35	0.61
1:B:357:LYS:HG2	1:B:464:ASP:HA	1.81	0.61
1:B:46:LYS:O	1:B:489:VAL:HG23	2.01	0.61
1:A:395:TYR:CD1	1:A:403:TYR:HA	2.35	0.60
1:A:213:ILE:O	1:A:253:PRO:HD3	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:242:VAL:HG12	1:B:243:SER:H	1.66	0.60
1:B:451:GLY:C	1:B:452:LEU:HD12	2.26	0.60
1:A:59:LYS:HG2	1:A:61:TYR:OH	2.01	0.60
1:A:207:LYS:HE2	1:A:436:ALA:HB3	1.85	0.59
1:B:335:GLU:OE2	1:B:408:ARG:HB3	2.03	0.59
1:A:102:GLU:O	1:A:106:GLU:HG2	2.03	0.59
1:B:105:HIS:HD1	1:B:479:TRP:HZ3	1.50	0.59
1:A:279:ASN:HD22	1:A:282:LYS:HG2	1.68	0.58
1:A:424:ILE:HD12	3:A:507:03G:H3	1.85	0.58
1:A:46:LYS:O	1:A:489:VAL:HG23	2.02	0.58
1:B:395:TYR:HD1	1:B:403:TYR:HA	1.65	0.58
1:A:69:TRP:HA	1:A:72:HIS:CD2	2.37	0.58
1:B:207:LYS:HE2	1:B:436:ALA:HB3	1.86	0.58
1:B:406:THR:HG23	2:B:506:NAG:O6	2.03	0.58
1:B:286:VAL:HB	1:B:452:LEU:HB2	1.86	0.58
1:B:465:THR:HG23	1:B:465:THR:O	2.04	0.57
1:A:424:ILE:CD1	3:A:507:03G:H3	2.34	0.57
1:A:215:LEU:H	1:A:251:ILE:H	1.52	0.57
1:A:286:VAL:HB	1:A:452:LEU:HB2	1.86	0.57
1:B:358:THR:OG1	1:B:396:ARG:HD2	2.05	0.57
1:B:406:THR:HG21	2:B:506:NAG:C1	2.34	0.57
1:A:59:LYS:HG2	1:A:61:TYR:CZ	2.39	0.56
1:A:339:ASN:HA	1:A:403:TYR:CZ	2.40	0.56
1:B:339:ASN:HA	1:B:403:TYR:CZ	2.41	0.56
1:B:121:LYS:HE3	1:B:426:MET:HE1	1.86	0.56
1:A:463:ASN:O	1:A:464:ASP:HB2	2.04	0.55
1:B:99:ASP:HA	1:B:102:GLU:OE1	2.07	0.55
1:A:451:GLY:C	1:A:452:LEU:HD12	2.32	0.55
1:A:476:ARG:O	1:A:480:ARG:HG3	2.07	0.54
1:B:484:TYR:CZ	1:B:485:LYS:HG3	2.43	0.54
1:B:78:ASP:HB2	1:B:81:PRO:HD3	1.89	0.54
1:B:370:GLU:HG3	1:B:384:TYR:CE2	2.42	0.54
1:A:342:LEU:HB2	1:A:403:TYR:HE2	1.73	0.54
1:A:280:ASN:ND2	1:A:457:ASP:O	2.37	0.54
1:A:105:HIS:ND1	1:A:479:TRP:HZ3	2.05	0.54
1:B:384:TYR:CE1	1:B:421:LYS:HB2	2.43	0.54
1:B:484:TYR:OH	1:B:485:LYS:HE3	2.08	0.53
1:A:69:TRP:HD1	1:A:114:GLU:OE1	1.91	0.53
1:A:78:ASP:N	1:A:79:PRO:HA	2.24	0.53
1:A:215:LEU:N	1:A:251:ILE:H	2.05	0.53
1:B:279:ASN:ND2	1:B:282:LYS:HG2	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:CYS:HB3	1:A:383:PHE:CE1	2.43	0.53
1:A:93:PHE:HB2	1:A:233:PHE:HZ	1.74	0.52
1:A:342:LEU:HB3	1:A:395:TYR:CE2	2.44	0.52
1:B:78:ASP:N	1:B:79:PRO:HA	2.24	0.52
1:A:226:LEU:HD13	1:A:489:VAL:HG11	1.90	0.52
1:B:405:HIS:CG	1:B:406:THR:N	2.72	0.52
1:B:335:GLU:HG2	1:B:414:ILE:HG13	1.91	0.52
1:B:102:GLU:O	1:B:106:GLU:HG2	2.10	0.52
1:B:279:ASN:HD22	1:B:282:LYS:HG2	1.74	0.51
1:A:358:THR:OG1	1:A:396:ARG:HD2	2.10	0.51
1:B:84:MET:HB2	1:B:244:THR:HB	1.93	0.51
1:A:242:VAL:HG12	1:A:243:SER:H	1.75	0.51
1:A:72:HIS:CE1	1:A:73:ALA:HB2	2.46	0.51
1:A:269:GLU:O	1:A:271:ILE:HD12	2.11	0.51
1:B:64:GLU:HB2	1:B:67:ASN:HD22	1.74	0.51
1:A:456:ARG:HH12	1:A:458:GLY:HA2	1.76	0.51
1:A:47:GLU:HG2	1:A:487:LYS:HE3	1.93	0.50
1:A:405:HIS:CG	1:A:406:THR:N	2.79	0.50
1:A:398:GLY:HA2	4:A:653:HOH:O	2.11	0.50
1:A:89:VAL:HG13	4:A:624:HOH:O	2.11	0.50
1:B:105:HIS:O	1:B:109:ILE:HG13	2.11	0.50
1:B:370:GLU:HG3	1:B:384:TYR:HE2	1.77	0.50
1:B:378:CYS:HB3	1:B:383:PHE:CE1	2.46	0.50
1:B:343:GLN:O	1:B:347:GLU:HG3	2.12	0.49
1:B:64:GLU:HG2	1:B:209:SER:HB3	1.93	0.49
1:A:242:VAL:HG12	1:A:243:SER:N	2.28	0.49
1:A:273:ARG:NH2	1:A:287:HIS:ND1	2.50	0.49
1:A:67:ASN:ND2	1:A:213:ILE:HD11	2.28	0.49
1:A:255:VAL:HG13	1:A:475:MET:HE3	1.94	0.49
1:B:242:VAL:HG12	1:B:243:SER:N	2.28	0.49
1:A:217:TYR:O	1:A:248:THR:HG23	2.13	0.48
1:A:368:ASP:HB3	1:A:370:GLU:OE1	2.13	0.48
1:A:395:TYR:HD1	1:A:404:ASN:H	1.59	0.48
1:B:338:TRP:CZ2	1:B:390:LEU:HG	2.48	0.48
1:B:216:HIS:HD2	1:B:248:THR:O	1.96	0.48
1:B:62:GLU:HG3	1:B:64:GLU:H	1.78	0.48
1:A:342:LEU:HB2	1:A:403:TYR:CE2	2.49	0.48
1:B:64:GLU:HB2	1:B:67:ASN:ND2	2.28	0.48
1:B:69:TRP:CB	1:B:111:LEU:HD13	2.43	0.48
1:B:339:ASN:HA	1:B:403:TYR:OH	2.13	0.47
1:A:340:ASN:O	1:A:344:LYS:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:SER:HA	1:A:383:PHE:O	2.14	0.47
1:B:280:ASN:HB2	1:B:456:ARG:O	2.14	0.47
1:A:270:ILE:HB	1:A:348:GLU:HG3	1.95	0.47
1:A:381:GLU:HG3	1:A:443:ILE:HD13	1.96	0.47
1:B:105:HIS:HA	1:B:479:TRP:CZ3	2.50	0.47
1:B:105:HIS:ND1	1:B:479:TRP:HZ3	2.11	0.47
1:B:349:LEU:O	1:B:353:PHE:HD2	1.97	0.47
1:A:249:HIS:O	1:A:251:ILE:HG13	2.14	0.47
1:B:69:TRP:HA	1:B:72:HIS:CD2	2.50	0.47
1:B:490:GLU:HG2	1:B:492:LYS:H	1.80	0.47
1:A:342:LEU:CB	1:A:403:TYR:HE2	2.27	0.47
1:B:388:SER:O	1:B:392:ASN:HB2	2.15	0.47
1:A:343:GLN:O	1:A:347:GLU:HG3	2.15	0.46
1:A:387:THR:HG23	1:A:387:THR:O	2.14	0.46
1:A:405:HIS:C	1:A:408:ARG:NH2	2.74	0.46
1:B:427:TRP:CE2	1:B:428:GLN:HG3	2.50	0.46
3:B:502:03G:H14	3:B:502:03G:H23	1.97	0.46
1:A:69:TRP:CD1	1:A:111:LEU:HA	2.50	0.46
1:A:104:MET:HA	1:A:217:TYR:OH	2.16	0.46
1:A:216:HIS:HD2	1:A:248:THR:O	1.98	0.46
1:B:45:TRP:HB3	1:B:489:VAL:HG21	1.98	0.46
1:B:384:TYR:OH	1:B:424:ILE:HB	2.16	0.46
1:A:69:TRP:CB	1:A:111:LEU:HD13	2.45	0.46
1:A:269:GLU:HB3	2:A:504:NAG:H61	1.98	0.46
1:B:269:GLU:HA	1:B:289:ASN:HD22	1.81	0.46
1:B:55:ALA:HA	1:B:75:VAL:O	2.16	0.45
1:B:395:TYR:HB2	4:B:619:HOH:O	2.17	0.45
1:A:213:ILE:HG23	1:A:214:PRO:HD2	1.98	0.45
1:A:338:TRP:CZ2	1:A:390:LEU:HG	2.52	0.45
1:A:456:ARG:HD2	1:A:466:GLU:OE1	2.17	0.45
1:B:259:LEU:HD13	1:B:449:ILE:HD13	1.98	0.45
1:A:55:ALA:HA	1:A:75:VAL:O	2.17	0.45
1:A:234:ASN:ND2	2:A:501:NAG:O7	2.50	0.45
1:B:67:ASN:ND2	1:B:213:ILE:HD11	2.31	0.45
1:A:427:TRP:HB3	3:A:507:03G:C7	2.46	0.44
1:B:375:SER:HA	1:B:383:PHE:O	2.16	0.44
1:B:255:VAL:HG13	1:B:475:MET:SD	2.56	0.44
1:B:363:PRO:HG3	1:B:388:SER:HA	1.99	0.44
1:B:364:SER:HB3	4:B:622:HOH:O	2.17	0.44
1:A:376:PHE:HE1	1:A:385:CYS:SG	2.40	0.44
1:B:69:TRP:HD1	1:B:114:GLU:OE1	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:424:ILE:HD12	3:B:502:03G:H4	1.99	0.44
1:B:329:ALA:HB3	1:B:418:CYS:HB2	1.98	0.44
1:A:62:GLU:HG3	1:A:64:GLU:H	1.82	0.44
1:B:104:MET:HA	1:B:217:TYR:OH	2.18	0.44
1:B:368:ASP:HB3	1:B:370:GLU:OE1	2.17	0.44
1:B:75:VAL:HG13	1:B:76:PRO:HD2	2.00	0.44
1:B:226:LEU:HD13	1:B:489:VAL:CG1	2.47	0.44
1:A:105:HIS:O	1:A:109:ILE:HG13	2.17	0.44
1:A:335:GLU:CD	1:A:408:ARG:HB3	2.41	0.44
1:A:370:GLU:HG2	3:A:507:03G:H5	1.83	0.44
1:B:396:ARG:H	1:B:399:THR:HG22	1.83	0.44
1:A:396:ARG:O	1:A:399:THR:HG22	2.17	0.43
1:B:65:VAL:HG13	1:B:208:VAL:HG21	2.00	0.43
1:B:78:ASP:CB	1:B:81:PRO:HD3	2.47	0.43
1:B:258:GLN:OE1	1:B:387:THR:HG21	2.18	0.43
1:A:208:VAL:HG22	1:A:209:SER:N	2.33	0.43
1:B:406:THR:HG21	2:B:506:NAG:O5	2.18	0.43
1:A:399:THR:HG23	1:A:403:TYR:N	2.33	0.43
1:B:325:ASN:HB3	1:B:328:GLN:HB2	1.99	0.43
1:B:342:LEU:CB	1:B:403:TYR:HE1	2.32	0.43
1:B:362:GLU:O	1:B:469:ARG:HA	2.18	0.43
1:A:78:ASP:HB2	1:A:80:ASN:N	2.33	0.43
1:A:108:ILE:HD12	1:A:253:PRO:HB3	2.01	0.43
1:A:299:PRO:O	1:A:300:ASN:C	2.60	0.43
1:B:270:ILE:HG23	1:B:287:HIS:O	2.19	0.43
1:A:407:GLY:C	1:A:408:ARG:HG3	2.43	0.43
1:B:52:LEU:HG	1:B:218:CYS:O	2.18	0.43
1:A:61:TYR:CD1	1:B:214:PRO:HG2	2.53	0.43
1:A:69:TRP:CG	1:A:111:LEU:HD13	2.53	0.43
1:B:392:ASN:CG	1:B:406:THR:HG21	2.44	0.43
1:A:405:HIS:C	1:A:408:ARG:HH22	2.26	0.43
1:B:279:ASN:ND2	1:B:281:ALA:HB3	2.32	0.43
1:B:360:LYS:HG2	1:B:394:THR:HB	2.00	0.43
1:A:70:ALA:O	1:A:74:CYS:N	2.52	0.42
1:B:396:ARG:O	1:B:399:THR:HG22	2.19	0.42
1:B:430:VAL:HG22	3:B:502:03G:H20	2.00	0.42
1:A:421:LYS:HB3	1:A:421:LYS:HE2	1.79	0.42
1:B:66:HIS:CE1	1:B:212:PRO:HA	2.54	0.42
1:B:69:TRP:HB3	1:B:111:LEU:HD13	2.01	0.42
1:B:59:LYS:HB3	1:B:61:TYR:CE2	2.55	0.42
1:B:229:ASN:OD1	1:B:243:SER:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:445:CYS:SG	2:B:503:NAG:H83	2.59	0.42
1:B:217:TYR:O	1:B:248:THR:HG23	2.19	0.42
1:A:289:ASN:OD1	1:A:289:ASN:C	2.61	0.42
1:B:474:ASP:HB2	3:B:502:03G:H15	2.02	0.42
2:A:504:NAG:O3	2:A:504:NAG:C7	2.68	0.41
1:B:78:ASP:HB2	1:B:80:ASN:N	2.35	0.41
1:A:78:ASP:HB2	1:A:81:PRO:HD3	2.01	0.41
1:A:64:GLU:HB2	1:A:67:ASN:HD22	1.85	0.41
1:A:395:TYR:HD1	1:A:403:TYR:HA	1.81	0.41
1:B:111:LEU:HD12	1:B:115:SER:OG	2.21	0.41
1:B:252:LYS:HA	1:B:253:PRO:HD3	1.80	0.41
1:A:289:ASN:OD1	1:A:290:GLU:N	2.52	0.41
1:A:83:GLU:HG2	1:A:245:VAL:CG1	2.50	0.41
1:B:427:TRP:HB3	3:B:502:03G:C7	2.51	0.41
1:A:339:ASN:HA	1:A:403:TYR:OH	2.20	0.41
1:B:396:ARG:HG2	1:B:397:ASN:N	2.36	0.41
1:B:369:LEU:HB3	1:B:384:TYR:CD2	2.56	0.41
1:A:219:ALA:HA	1:A:220:PRO:HD3	1.94	0.41
1:A:376:PHE:C	3:A:507:03G:CL1	3.00	0.41
1:B:118:PRO:HB3	1:B:433:ALA:HB1	2.03	0.41
1:B:212:PRO:HB3	1:B:253:PRO:HD2	2.03	0.41
1:B:341:THR:O	1:B:345:VAL:HG23	2.21	0.41
1:A:45:TRP:HB3	1:A:489:VAL:HG21	2.03	0.41
1:A:80:ASN:HB2	4:A:634:HOH:O	2.21	0.41
1:A:376:PHE:HA	3:A:507:03G:CL1	2.57	0.41
1:A:427:TRP:HA	3:A:507:03G:C13	2.51	0.41
1:A:66:HIS:CE1	1:A:212:PRO:HA	2.56	0.40
1:A:335:GLU:HB2	2:A:505:NAG:H62	2.04	0.40
1:B:429:GLU:HG2	1:B:430:VAL:N	2.35	0.40
1:A:396:ARG:HG2	1:A:397:ASN:N	2.35	0.40
1:B:47:GLU:HG2	1:B:487:LYS:HE3	2.03	0.40
1:B:69:TRP:HD1	1:B:114:GLU:CD	2.29	0.40
1:B:476:ARG:HB2	1:B:480:ARG:NH1	2.37	0.40
1:B:407:GLY:C	1:B:408:ARG:HG3	2.47	0.40
1:A:105:HIS:HA	1:A:479:TRP:CZ3	2.57	0.40
1:B:421:LYS:HB3	1:B:421:LYS:HE2	1.82	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	339/358 (95%)	305 (90%)	31 (9%)	3 (1%)	14	35
1	B	339/358 (95%)	306 (90%)	32 (9%)	1 (0%)	36	60
All	All	678/716 (95%)	611 (90%)	63 (9%)	4 (1%)	21	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	356	SER
1	A	115	SER
1	A	341	THR
1	B	407	GLY

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	304/312 (97%)	287 (94%)	17 (6%)	19	44
1	B	304/312 (97%)	292 (96%)	12 (4%)	28	57
All	All	608/624 (97%)	579 (95%)	29 (5%)	23	50

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	VAL
1	A	45	TRP

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Mol	Chain	Res	Type
1	A	85	VAL
1	A	90	THR
1	A	123	THR
1	A	226	LEU
1	A	277	LEU
1	A	290	GLU
1	A	387	THR
1	A	399	THR
1	A	406	THR
1	A	424	ILE
1	A	429	GLU
1	A	444	THR
1	A	445	CYS
1	A	467	THR
1	A	488	VAL
1	B	45	TRP
1	B	72	HIS
1	B	90	THR
1	B	111	LEU
1	B	202	THR
1	B	226	LEU
1	B	236	THR
1	B	245	VAL
1	B	336	SER
1	B	399	THR
1	B	406	THR
1	B	424	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
1	A	203	GLN
1	A	229	ASN
1	A	230	ASN
1	A	279	ASN
1	A	417	GLN
1	B	67	ASN
1	B	72	HIS
1	B	94	ASN
1	B	216	HIS
1	B	279	ASN

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Mol	Chain	Res	Type
1	B	425	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

15 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	NAG	A	502	1	14,14,15	0.30	0	17,19,21	0.76	0
2	NAG	B	503	1	14,14,15	0.28	0	17,19,21	0.56	0
2	NAG	A	505	1	14,14,15	0.29	0	17,19,21	0.56	0
2	NAG	A	504	1	14,14,15	0.29	0	17,19,21	0.57	0
2	NAG	A	503	1	14,14,15	0.34	0	17,19,21	0.57	0
2	NAG	A	506	1	14,14,15	0.37	0	17,19,21	1.02	1 (5%)
2	NAG	B	501	1	14,14,15	0.33	0	17,19,21	0.51	0
2	NAG	B	505	1	14,14,15	0.29	0	17,19,21	0.64	0
2	NAG	A	508	1	14,14,15	0.26	0	17,19,21	0.50	0
3	03G	A	507	-	24,24,24	2.79	7 (29%)	36,36,36	2.27	15 (41%)
2	NAG	A	509	1	14,14,15	0.27	0	17,19,21	0.84	1 (5%)
2	NAG	B	504	1	14,14,15	0.30	0	17,19,21	0.62	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	B	506	1	14,14,15	0.30	0	17,19,21	0.58	0
2	NAG	A	501	-	14,14,15	0.28	0	17,19,21	0.57	0
3	03G	B	502	-	24,24,24	2.72	7 (29%)	36,36,36	1.97	11 (30%)

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	NAG	A	502	1	-	2/6/23/26	0/1/1/1
2	NAG	B	503	1	-	2/6/23/26	0/1/1/1
2	NAG	A	505	1	-	4/6/23/26	0/1/1/1
2	NAG	A	504	1	-	5/6/23/26	0/1/1/1
2	NAG	A	503	1	-	2/6/23/26	0/1/1/1
2	NAG	A	506	1	-	2/6/23/26	0/1/1/1
2	NAG	B	501	1	-	2/6/23/26	0/1/1/1
2	NAG	B	505	1	-	3/6/23/26	0/1/1/1
2	NAG	A	508	1	-	3/6/23/26	0/1/1/1
3	03G	A	507	-	-	1/12/28/28	0/2/2/2
2	NAG	A	509	1	-	2/6/23/26	0/1/1/1
2	NAG	B	504	1	-	3/6/23/26	0/1/1/1
2	NAG	B	506	1	-	3/6/23/26	0/1/1/1
2	NAG	A	501	-	-	4/6/23/26	0/1/1/1
3	03G	B	502	-	-	2/12/28/28	0/2/2/2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	507	03G	C1-CL1	-10.75	1.49	1.74
3	B	502	03G	C1-CL1	-9.98	1.51	1.74
3	B	502	03G	C7-N1	4.00	1.43	1.35
3	A	507	03G	C7-N1	3.77	1.43	1.35
3	B	502	03G	C8-C7	3.76	1.61	1.53
3	A	507	03G	C13-C12	3.42	1.57	1.53
3	A	507	03G	C8-C7	3.21	1.60	1.53
3	B	502	03G	C8-N2	2.95	1.40	1.34
3	A	507	03G	C8-N2	2.85	1.40	1.34
3	B	502	03G	C4-N1	2.80	1.47	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	03G	C10-C11	2.66	1.56	1.53
3	A	507	03G	C4-N1	2.45	1.46	1.41
3	B	502	03G	O2-C7	-2.18	1.19	1.23
3	A	507	03G	O2-C7	-2.17	1.19	1.23

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	507	03G	C12-C13-C9	5.17	116.43	112.41
3	A	507	03G	C6-C1-C2	-5.07	114.96	121.24
3	B	502	03G	C6-C1-C2	-4.94	115.12	121.24
3	B	502	03G	C8-C7-N1	4.13	119.23	112.25
3	B	502	03G	C5-C6-C1	4.04	123.31	119.24
3	A	507	03G	C8-C7-N1	4.02	119.04	112.25
3	A	507	03G	C5-C6-C1	3.73	122.99	119.24
3	A	507	03G	C3-C2-C1	3.49	122.76	119.24
3	A	507	03G	C13-C12-N3	3.26	116.40	110.42
3	A	507	03G	C5-C4-C3	-3.21	114.78	119.04
3	A	507	03G	C6-C1-CL1	3.00	123.78	119.36
3	B	502	03G	C3-C2-C1	2.99	122.25	119.24
3	B	502	03G	O2-C7-C8	-2.94	116.45	121.24
3	B	502	03G	C5-C4-C3	-2.93	115.15	119.04
3	A	507	03G	O2-C7-C8	-2.83	116.63	121.24
3	B	502	03G	C15-C11-C10	-2.80	106.97	110.06
3	A	507	03G	C17-C12-C13	-2.79	106.98	110.06
3	B	502	03G	C6-C1-CL1	2.61	123.21	119.36
2	A	506	NAG	O5-C1-C2	-2.54	107.36	111.29
3	A	507	03G	C11-C10-C9	-2.48	110.49	112.41
3	B	502	03G	C10-C11-N3	2.42	114.85	110.42
2	A	509	NAG	O5-C1-C2	-2.41	107.57	111.29
3	A	507	03G	C4-N1-C7	-2.33	123.31	127.45
3	B	502	03G	C2-C3-C4	2.30	122.95	120.30
3	B	502	03G	C12-N3-C11	-2.26	116.14	118.98
3	A	507	03G	C12-N3-C11	-2.25	116.14	118.98
3	A	507	03G	C15-C11-C10	-2.22	107.60	110.06
3	A	507	03G	C2-C3-C4	2.17	122.80	120.30

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	A	502	NAG	C8-C7-N2-C2
2	A	502	NAG	O7-C7-N2-C2
2	A	504	NAG	C3-C2-N2-C7
2	A	504	NAG	C8-C7-N2-C2
2	A	504	NAG	O7-C7-N2-C2
2	A	505	NAG	C8-C7-N2-C2
2	A	508	NAG	C8-C7-N2-C2
2	A	508	NAG	O7-C7-N2-C2
2	A	509	NAG	C8-C7-N2-C2
2	A	509	NAG	O7-C7-N2-C2
2	B	501	NAG	C8-C7-N2-C2
2	B	501	NAG	O7-C7-N2-C2
2	B	503	NAG	C8-C7-N2-C2
2	B	503	NAG	O7-C7-N2-C2
2	B	505	NAG	C8-C7-N2-C2
2	B	505	NAG	O7-C7-N2-C2
2	B	506	NAG	C8-C7-N2-C2
2	B	506	NAG	O7-C7-N2-C2
2	A	501	NAG	C8-C7-N2-C2
2	A	501	NAG	O5-C5-C6-O6
2	A	505	NAG	O7-C7-N2-C2
2	A	505	NAG	O5-C5-C6-O6
2	A	501	NAG	C4-C5-C6-O6
2	A	505	NAG	C4-C5-C6-O6
2	A	503	NAG	C8-C7-N2-C2
2	A	503	NAG	O7-C7-N2-C2
2	A	506	NAG	C8-C7-N2-C2
2	B	504	NAG	C8-C7-N2-C2
2	B	504	NAG	O7-C7-N2-C2
2	A	506	NAG	O7-C7-N2-C2
2	A	504	NAG	O5-C5-C6-O6
3	A	507	03G	O1-C8-N2-C9
3	B	502	03G	C10-C9-N2-C8
2	B	504	NAG	O5-C5-C6-O6
2	B	505	NAG	C3-C2-N2-C7
2	A	508	NAG	O5-C5-C6-O6
2	B	506	NAG	C1-C2-N2-C7
2	A	504	NAG	C4-C5-C6-O6
3	B	502	03G	C13-C9-N2-C8

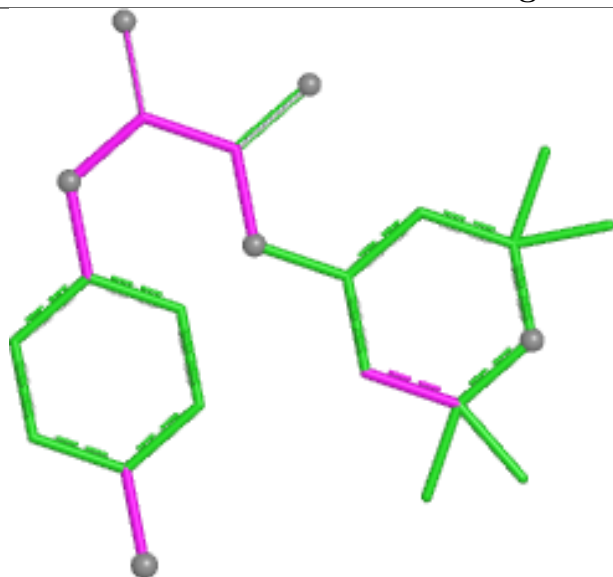
There are no ring outliers.

7 monomers are involved in 21 short contacts:

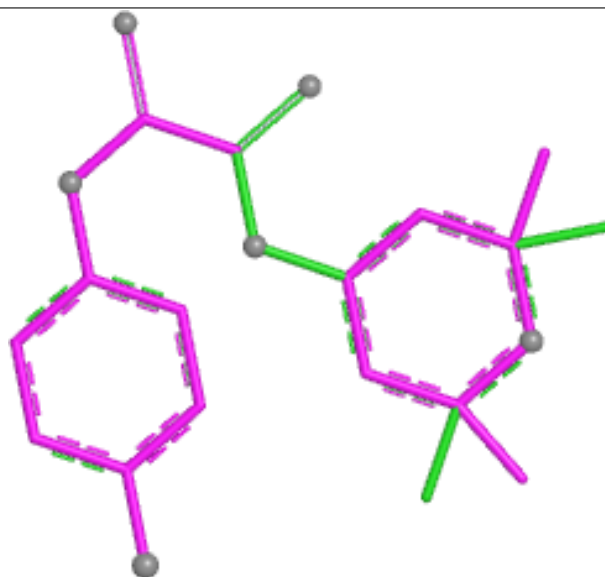
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	503	NAG	1	0
2	A	505	NAG	1	0
2	A	504	NAG	2	0
3	A	507	03G	7	0
2	B	506	NAG	3	0
2	A	501	NAG	2	0
3	B	502	03G	5	0

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.

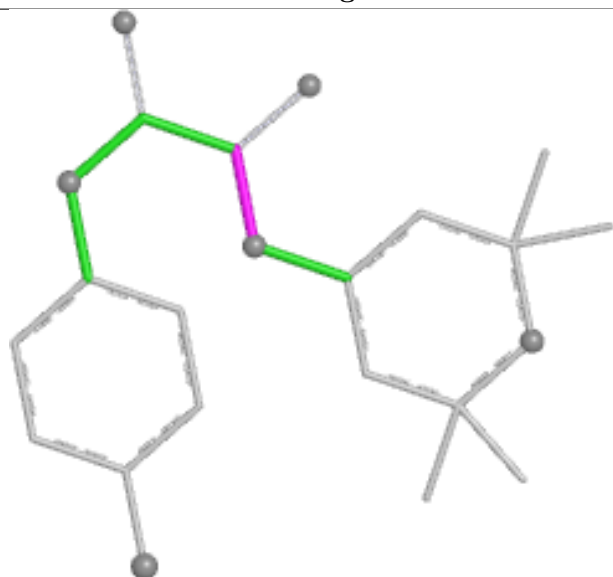
Ligand 03G A 507



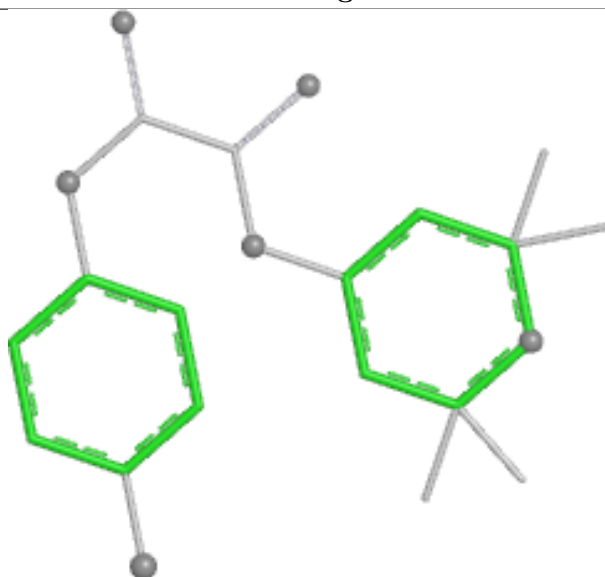
Bond lengths



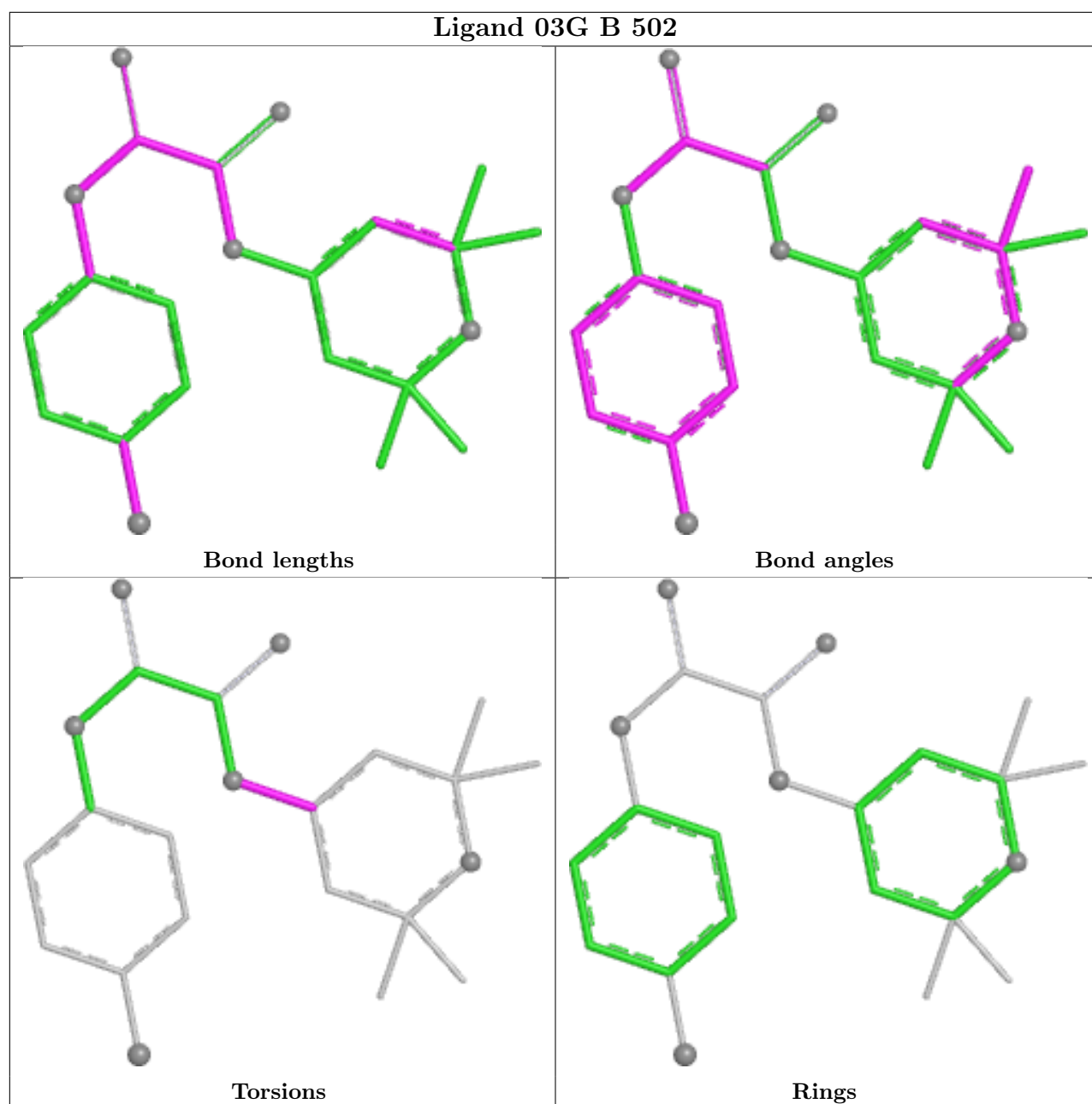
Bond angles



Torsions



Rings



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	345/358 (96%)	1.12	63 (18%) 3 3	20, 56, 148, 235	0
1	B	345/358 (96%)	1.22	70 (20%) 3 2	34, 73, 146, 252	0
All	All	690/716 (96%)	1.17	133 (19%) 3 3	20, 65, 148, 252	0

All (133) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	459	GLY	8.0
1	A	403	TYR	6.6
1	A	48	ALA	6.6
1	B	44	VAL	6.5
1	A	354	PRO	6.3
1	A	463	ASN	6.0
1	A	410	SER	5.9
1	A	459	GLY	5.9
1	A	408	ARG	5.4
1	B	45	TRP	5.3
1	B	463	ASN	5.1
1	B	224	ALA	5.0
1	B	296	CYS	4.9
1	A	79	PRO	4.8
1	A	64	GLU	4.7
1	B	272	ILE	4.5
1	A	44	VAL	4.5
1	A	409	SER	4.4
1	B	403	TYR	4.4
1	A	335	GLU	4.3
1	A	45	TRP	4.2
1	A	396	ARG	4.2
1	A	464	ASP	4.1
1	B	48	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	79	PRO	3.8
1	A	488	VAL	3.8
1	B	438	PRO	3.8
1	B	404	ASN	3.6
1	A	395	TYR	3.6
1	A	78	ASP	3.5
1	A	81	PRO	3.5
1	A	489	VAL	3.5
1	A	224	ALA	3.5
1	B	217	TYR	3.4
1	B	409	SER	3.4
1	A	47	GLU	3.4
1	B	86	LEU	3.4
1	A	46	LYS	3.4
1	B	492	LYS	3.4
1	B	81	PRO	3.4
1	A	50	THR	3.3
1	B	87	ALA	3.3
1	B	99	ASP	3.2
1	A	221	ALA	3.2
1	A	356	SER	3.2
1	B	80	ASN	3.1
1	A	491	ILE	3.1
1	B	407	GLY	3.1
1	B	458	GLY	3.1
1	A	458	GLY	3.1
1	A	80	ASN	3.0
1	A	399	THR	3.0
1	A	479	TRP	3.0
1	A	77	THR	3.0
1	A	246	GLN	3.0
1	A	240	ARG	3.0
1	A	245	VAL	3.0
1	B	82	GLN	2.9
1	A	327	ARG	2.9
1	B	225	ILE	2.9
1	A	407	GLY	2.9
1	A	300	ASN	2.8
1	A	208	VAL	2.8
1	A	84	MET	2.8
1	B	356	SER	2.8
1	B	464	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	492	LYS	2.8
1	B	410	SER	2.8
1	B	83	GLU	2.8
1	A	217	TYR	2.8
1	B	354	PRO	2.7
1	B	52	LEU	2.7
1	A	119	CYS	2.7
1	B	226	LEU	2.7
1	B	479	TRP	2.7
1	A	87	ALA	2.7
1	B	395	TYR	2.7
1	A	241	ASN	2.7
1	B	247	CYS	2.6
1	B	248	THR	2.6
1	B	231	LYS	2.6
1	B	241	ASN	2.6
1	A	465	THR	2.6
1	B	399	THR	2.6
1	B	77	THR	2.5
1	B	276	ASN	2.5
1	B	396	ARG	2.5
1	B	411	ASN	2.5
1	B	47	GLU	2.5
1	B	353	PHE	2.5
1	A	289	ASN	2.4
1	A	213	ILE	2.4
1	B	119	CYS	2.4
1	A	472	GLY	2.4
1	B	124	GLY	2.4
1	A	411	ASN	2.4
1	B	64	GLU	2.4
1	B	352	HIS	2.4
1	B	408	ARG	2.4
1	B	267	GLU	2.3
1	B	245	VAL	2.3
1	B	327	ARG	2.3
1	A	49	LYS	2.3
1	A	347	GLU	2.3
1	B	46	LYS	2.3
1	B	442	GLU	2.3
1	A	388	SER	2.3
1	B	365	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	398	GLY	2.3
1	B	89	VAL	2.3
1	A	225	ILE	2.2
1	B	440	GLU	2.2
1	B	198	GLY	2.2
1	A	89	VAL	2.2
1	A	55	ALA	2.2
1	A	88	ASN	2.2
1	A	269	GLU	2.2
1	A	398	GLY	2.2
1	B	85	VAL	2.2
1	B	58	ALA	2.2
1	B	300	ASN	2.1
1	B	339	ASN	2.1
1	B	232	THR	2.1
1	A	397	ASN	2.1
1	A	281	ALA	2.1
1	B	244	THR	2.1
1	B	325	ASN	2.1
1	B	341	THR	2.1
1	B	78	ASP	2.0
1	B	114	GLU	2.0
1	A	220	PRO	2.0
1	A	272	ILE	2.0
1	B	393	GLY	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 ⓘ

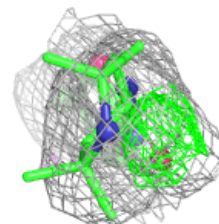
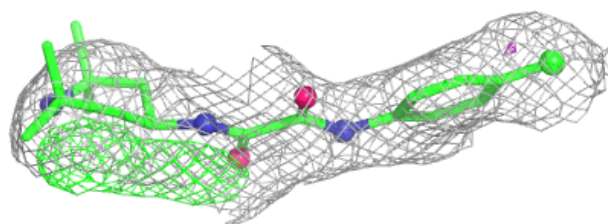
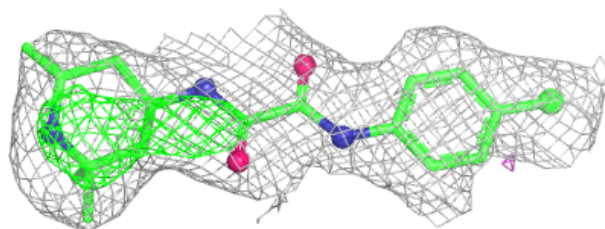
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	A	504	14/15	0.56	0.27	63,73,85,85	0
2	NAG	B	505	14/15	0.62	0.18	54,80,85,89	0
2	NAG	A	505	14/15	0.64	0.21	37,68,88,100	0
2	NAG	A	509	14/15	0.65	0.20	105,113,126,126	0
2	NAG	B	501	14/15	0.67	0.16	57,100,109,110	0
2	NAG	A	501	14/15	0.79	0.20	93,99,107,108	0
2	NAG	B	504	14/15	0.81	0.13	57,67,85,88	0
2	NAG	A	503	14/15	0.81	0.12	60,62,66,68	0
2	NAG	B	506	14/15	0.83	0.15	67,78,88,88	0
2	NAG	A	506	14/15	0.84	0.14	47,55,71,72	0
2	NAG	B	503	14/15	0.87	0.14	57,62,71,73	0
3	03G	B	502	23/23	0.88	0.18	30,50,77,112	0
3	03G	A	507	23/23	0.91	0.12	5,26,39,96	0
2	NAG	A	508	14/15	0.92	0.11	32,39,54,56	0
2	NAG	A	502	14/15	0.93	0.11	31,40,54,56	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.

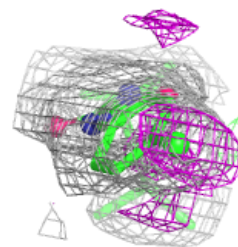
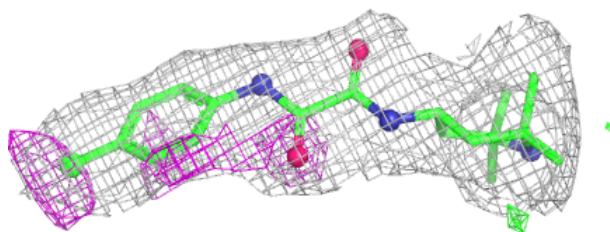
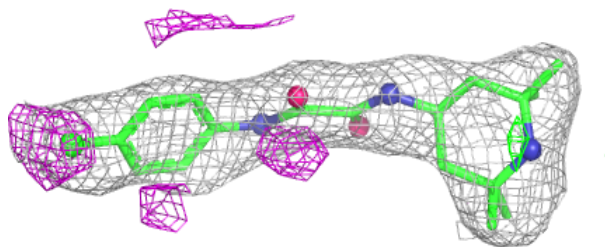
Electron density around 03G B 502:

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 03G A 507:

$2mF_o-DF_c$ (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.