



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:33 PM UTC

PDB ID : 8GV5 / pdb_00008gv5
Title : Crystal structure of PN-SIA28 in complex with influenza hemagglutinin A/s
wine/Guangdong/104/2013 (H1N1)
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Deposited on : 2022-09-14
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

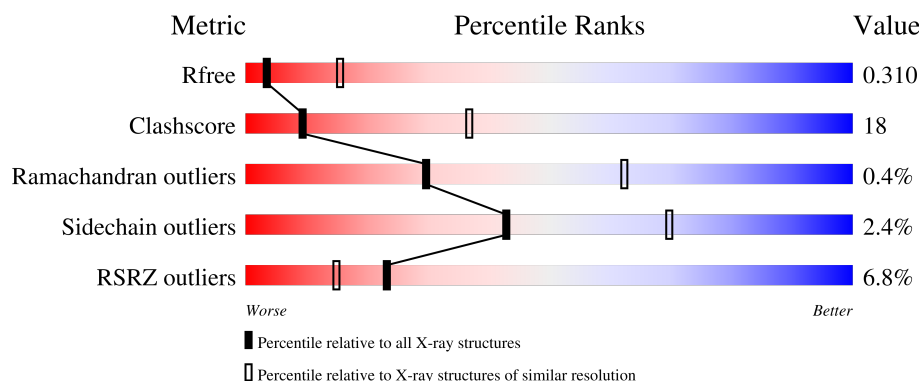
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	330	6% 62% 33% ..
2	B	173	10% 69% 22% 9%
3	C	125	5% 74% 24% ..
4	D	108	4% 69% 30% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5503 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 Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2470	1556	427	475	12			

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	157	Total	C	N	O	S	0	0	0
			1267	796	217	247	7			

- Molecule 3 is a protein called PN-SIA28 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	124	Total	C	N	O	S	0	0	0
			951	606	162	179	4			

- Molecule 4 is a protein called PN-SIA28 Light Chain.

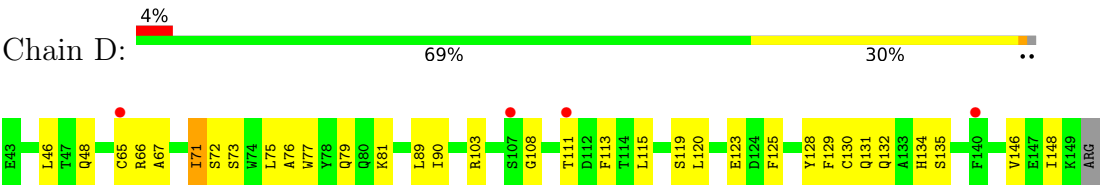
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	107	Total	C	N	O	S	0	0	0
			801	510	133	156	2			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



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

● Molecule 4: PN-SIA28 Light Chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	85.80Å 85.80Å 231.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.63 – 3.20 45.63 – 3.20	Depositor EDS
% Data completeness (in resolution range)	73.2 (45.63-3.20) 73.2 (45.63-3.20)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 3.19Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.267 , 0.308 0.266 , 0.310	Depositor DCC
R_{free} test set	654 reflections (4.11%)	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.353	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.084 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	5503	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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

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	0/2531	0.78	5/3450 (0.1%)
2	B	0.19	0/1294	0.48	0/1741
3	C	0.12	0/975	0.38	0/1323
4	D	0.28	0/821	0.51	0/1115
All	All	0.33	0/5621	0.63	5/7629 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	214	ILE	CA-C-N	-9.45	112.52	121.97
1	A	214	ILE	C-N-CA	-9.45	112.52	121.97
1	A	74	ASP	N-CA-C	-6.37	105.82	113.97
1	A	229	TYR	CB-CA-C	-5.05	102.69	110.78
1	A	72	THR	CB-CA-C	-5.02	103.00	110.88

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	2470	0	2327	118	0
2	B	1267	0	1186	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	951	0	924	24	0
4	D	801	0	788	32	0
5	A	14	0	13	0	0
All	All	5503	0	5238	193	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 18.

All (193) 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)
4:D:71:ILE:HD11	4:D:113:PHE:CZ	1.59	1.35
2:B:452:GLN:HE22	2:B:482:GLY:HA2	1.32	0.95
1:A:293:GLN:HE22	1:A:299:THR:HG21	1.32	0.94
2:B:457:ALA:HA	2:B:466:GLU:O	1.69	0.93
4:D:71:ILE:HD11	4:D:113:PHE:HZ	1.07	0.86
1:A:180:HIS:CE1	1:A:182:PRO:HG3	2.12	0.85
1:A:171:LYS:HD3	1:A:258:LYS:CB	2.07	0.84
3:C:43:GLN:H	3:C:46:ARG:HH21	1.23	0.83
1:A:31:ASN:HB2	1:A:313:ARG:HH21	1.43	0.82
1:A:92:PRO:HB2	1:A:226:ARG:HD3	1.60	0.82
1:A:16:VAL:HG21	1:A:315:ALA:HB2	1.62	0.82
3:C:71:PRO:HD3	3:C:122:ALA:HA	1.59	0.82
4:D:71:ILE:CD1	4:D:113:PHE:CZ	2.55	0.80
4:D:71:ILE:CD1	4:D:113:PHE:HZ	1.91	0.78
1:A:219:LYS:HA	1:A:223:GLN:O	1.86	0.76
3:C:139:LEU:HD21	4:D:135:SER:HA	1.70	0.74
1:A:41:LEU:HG	1:A:270:ALA:HB3	1.70	0.73
1:A:180:HIS:HE1	1:A:182:PRO:CA	2.02	0.73
1:A:29:SER:OG	1:A:313:ARG:NH1	2.24	0.71
1:A:44:LEU:C	1:A:44:LEU:HD12	2.16	0.71
4:D:71:ILE:HD11	4:D:113:PHE:CE2	2.23	0.70
1:A:180:HIS:CE1	1:A:182:PRO:CG	2.74	0.70
1:A:180:HIS:CE1	1:A:182:PRO:CA	2.76	0.68
4:D:65:CYS:O	4:D:113:PHE:HB2	1.94	0.68
1:A:2:THR:HG22	2:B:466:GLU:HG2	1.76	0.68
1:A:44:LEU:HD12	1:A:44:LEU:O	1.94	0.67
2:B:395:LYS:HG3	2:B:408:ASN:ND2	2.10	0.67
1:A:192:TYR:HE2	1:A:247:ASN:HA	1.60	0.67
4:D:71:ILE:C	4:D:71:ILE:HD13	2.20	0.66
1:A:293:GLN:NE2	1:A:299:THR:HG21	2.06	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:GLN:NE2	1:A:281:THR:O	2.29	0.65
1:A:281:THR:HG22	1:A:299:THR:HG22	1.78	0.65
1:A:103:GLU:O	1:A:106:SER:HB2	1.97	0.64
3:C:50:LEU:HD22	3:C:124:TYR:HD2	1.63	0.64
1:A:92:PRO:HG2	1:A:220:VAL:HB	1.80	0.64
1:A:16:VAL:HG12	1:A:313:ARG:HB3	1.79	0.64
2:B:395:LYS:HG3	2:B:408:ASN:HD21	1.63	0.63
1:A:133:THR:HG23	1:A:136:CYS:H	1.62	0.63
2:B:352:HIS:HD2	2:B:361:TYR:CE1	2.16	0.63
1:A:180:HIS:HE1	1:A:182:PRO:CB	2.12	0.63
3:C:35:VAL:O	3:C:52:CYS:HA	1.99	0.63
1:A:217:ARG:HB3	1:A:226:ARG:NH2	2.13	0.62
3:C:81:VAL:HG23	3:C:88:LYS:HG2	1.80	0.62
4:D:46:LEU:HD23	4:D:67:ALA:HB2	1.81	0.62
4:D:66:ARG:NH2	4:D:67:ALA:O	2.34	0.61
1:A:180:HIS:CE1	1:A:182:PRO:HA	2.34	0.61
2:B:352:HIS:HD2	2:B:361:TYR:HE1	1.48	0.60
4:D:76:ALA:HA	4:D:90:ILE:O	2.00	0.60
1:A:150:TRP:CD1	1:A:192:TYR:HH	2.19	0.60
1:A:226:ARG:HG3	1:A:226:ARG:NH1	2.16	0.59
3:C:142:LEU:HD13	4:D:131:GLN:HE22	1.67	0.59
1:A:242:PHE:HE2	1:A:251:PRO:HG2	1.67	0.59
4:D:48:GLN:HE22	4:D:129:PHE:HA	1.68	0.59
4:D:79:GLN:HB2	4:D:89:LEU:HD11	1.85	0.59
4:D:79:GLN:HG3	4:D:128:TYR:HE1	1.68	0.58
1:A:102:ARG:NE	2:B:396:GLU:OE1	2.37	0.58
1:A:263:SER:HA	2:B:390:PHE:CE1	2.39	0.58
4:D:75:LEU:HD11	4:D:130:CYS:HB2	1.86	0.57
4:D:108:GLY:HA3	4:D:113:PHE:HA	1.86	0.57
1:A:226:ARG:CG	1:A:226:ARG:HH11	2.17	0.56
1:A:77:SER:O	1:A:263:SER:CB	2.53	0.56
1:A:31:ASN:HB2	1:A:313:ARG:NH2	2.16	0.56
1:A:133:THR:HG1	1:A:223:GLN:NE2	2.03	0.56
2:B:409:LYS:NZ	2:B:413:ASP:OD2	2.38	0.56
1:A:4:CYS:HB2	2:B:352:HIS:HB3	1.88	0.55
1:A:180:HIS:HE1	1:A:182:PRO:HA	1.70	0.55
1:A:25:THR:HG22	1:A:320:ASN:HB3	1.87	0.55
1:A:188:GLN:HE22	1:A:247:ASN:HD21	1.54	0.55
4:D:77:TRP:CE2	4:D:115:LEU:HB2	2.41	0.55
1:A:180:HIS:HB2	1:A:249:ILE:HD11	1.88	0.55
1:A:293:GLN:HG2	1:A:304:PRO:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:266:MET:HE1	1:A:280:GLN:HG3	1.89	0.55
1:A:180:HIS:HE1	1:A:182:PRO:CG	2.18	0.55
2:B:355:ASN:ND2	2:B:473:ASN:HA	2.21	0.55
1:A:180:HIS:O	1:A:182:PRO:HD3	2.07	0.54
1:A:188:GLN:NE2	1:A:247:ASN:HD21	2.05	0.54
1:A:40:LYS:HG3	1:A:273:HIS:HB2	1.88	0.54
4:D:66:ARG:NE	4:D:111:THR:O	2.41	0.54
1:A:95:PHE:CE1	1:A:101:LEU:HD22	2.43	0.54
1:A:150:TRP:NE1	1:A:192:TYR:OH	2.40	0.53
1:A:126:HIS:CE1	1:A:161:LEU:HB3	2.43	0.53
1:A:133:THR:OG1	1:A:223:GLN:NE2	2.41	0.53
1:A:301:GLY:HA2	2:B:390:PHE:CE2	2.44	0.53
4:D:76:ALA:HB3	4:D:131:GLN:HB3	1.90	0.53
1:A:171:LYS:HB3	1:A:257:LEU:O	2.09	0.53
2:B:453:LEU:HD13	2:B:467:PHE:HE1	1.72	0.53
1:A:51:GLN:NE2	1:A:272:VAL:HG23	2.24	0.53
1:A:180:HIS:ND1	1:A:182:PRO:HG3	2.22	0.53
3:C:94:VAL:HG12	3:C:97:ARG:HH21	1.72	0.53
1:A:68:ASP:OD1	1:A:146:ARG:NE	2.42	0.53
1:A:115:GLU:OE2	1:A:118:PRO:HA	2.10	0.52
1:A:98:TYR:CE2	1:A:102:ARG:HD2	2.45	0.52
1:A:126:HIS:HE1	1:A:161:LEU:HB3	1.74	0.52
1:A:312:LEU:HD22	2:B:427:ILE:HD12	1.91	0.52
1:A:153:GLN:HB3	1:A:192:TYR:HA	1.92	0.52
1:A:301:GLY:HA2	2:B:390:PHE:CD2	2.46	0.51
1:A:104:GLN:C	1:A:106:SER:H	2.20	0.50
1:A:223:GLN:OE1	1:A:223:GLN:HA	2.11	0.50
1:A:226:ARG:NH1	1:A:226:ARG:CG	2.73	0.50
4:D:132:GLN:OE1	4:D:134:HIS:N	2.44	0.50
4:D:48:GLN:NE2	4:D:130:CYS:H	2.10	0.50
1:A:291:PRO:HG3	2:B:383:ILE:HA	1.92	0.50
1:A:1:ASP:HB3	2:B:467:PHE:HB2	1.94	0.49
2:B:425:LEU:O	2:B:429:LEU:HG	2.11	0.49
4:D:81:LYS:NZ	4:D:123:GLU:O	2.41	0.49
1:A:43:SER:HB2	1:A:48:SER:HA	1.94	0.49
1:A:263:SER:HA	2:B:390:PHE:HE1	1.75	0.49
1:A:133:THR:O	1:A:142:LYS:HA	2.13	0.49
1:A:226:ARG:HG3	1:A:226:ARG:HH11	1.77	0.49
2:B:344:MET:HE3	2:B:363:ALA:HA	1.94	0.49
4:D:71:ILE:HG12	4:D:71:ILE:O	2.12	0.49
1:A:180:HIS:CE1	1:A:182:PRO:N	2.81	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:344:MET:HE1	2:B:350:GLY:HA3	1.94	0.48
2:B:481:ASN:O	2:B:481:ASN:ND2	2.45	0.48
2:B:450:LYS:HE2	2:B:459:GLU:OE2	2.13	0.48
1:A:60:TRP:CD1	1:A:70:LEU:HB3	2.49	0.48
1:A:319:ARG:HE	2:B:435:LEU:HD13	1.78	0.48
1:A:135:ALA:HB2	1:A:223:GLN:HE21	1.79	0.48
1:A:208:GLN:HG2	1:A:209:ARG:O	2.14	0.48
3:C:62:TYR:CD2	3:C:128:ARG:HD2	2.49	0.48
1:A:79:ILE:HB	1:A:265:MET:HG3	1.96	0.47
4:D:75:LEU:HD22	4:D:113:PHE:CD2	2.49	0.47
1:A:174:LEU:O	1:A:254:ALA:HB1	2.14	0.47
1:A:44:LEU:C	1:A:44:LEU:CD1	2.85	0.47
1:A:32:LEU:O	1:A:292:PHE:HB2	2.15	0.47
1:A:227:ILE:CD1	1:A:249:ILE:HG13	2.45	0.47
1:A:227:ILE:HD13	1:A:249:ILE:HG13	1.96	0.46
1:A:40:LYS:O	1:A:42:CYS:N	2.48	0.46
1:A:180:HIS:ND1	1:A:180:HIS:C	2.72	0.46
3:C:69:GLN:HB2	3:C:75:LEU:HD23	1.97	0.46
1:A:114:PHE:O	1:A:253:HIS:HB2	2.16	0.45
1:A:65:PRO:HA	1:A:146:ARG:HE	1.82	0.45
1:A:77:SER:O	1:A:263:SER:OG	2.33	0.45
4:D:125:PHE:CE2	4:D:148:ILE:HG12	2.52	0.45
1:A:28:HIS:ND1	3:C:133:PHE:O	2.50	0.45
1:A:112:GLU:O	1:A:113:ARG:C	2.59	0.45
2:B:452:GLN:HE22	2:B:482:GLY:CA	2.17	0.45
1:A:185:TYR:HA	1:A:188:GLN:HB3	1.99	0.44
3:C:140:ASN:OD1	3:C:141:GLY:N	2.48	0.44
1:A:105:LEU:HD13	1:A:105:LEU:HA	1.77	0.44
1:A:192:TYR:CE2	1:A:247:ASN:HA	2.47	0.44
1:A:274:ASN:C	1:A:274:ASN:HD22	2.25	0.44
3:C:50:LEU:HD13	3:C:113:MET:HE2	1.98	0.44
1:A:133:THR:HG1	1:A:223:GLN:HE21	1.65	0.44
1:A:266:MET:HE3	1:A:300:ILE:HD12	2.00	0.44
4:D:75:LEU:HD12	4:D:131:GLN:O	2.18	0.44
1:A:126:HIS:CD2	1:A:159:PRO:HG2	2.53	0.44
1:A:28:HIS:CD2	3:C:133:PHE:HB3	2.53	0.44
1:A:112:GLU:O	1:A:114:PHE:HB3	2.18	0.44
3:C:113:MET:SD	3:C:124:TYR:HE2	2.41	0.44
1:A:43:SER:HA	1:A:49:PRO:HD3	2.00	0.43
1:A:133:THR:CG2	1:A:136:CYS:H	2.28	0.43
1:A:212:PRO:CB	1:A:247:ASN:HD22	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:347:GLY:HA2	3:C:135:ILE:HD12	1.99	0.43
2:B:355:ASN:HD22	2:B:473:ASN:HA	1.82	0.43
1:A:219:LYS:HG2	1:A:223:GLN:C	2.44	0.43
3:C:82:SER:O	3:C:102:ARG:NH1	2.51	0.43
2:B:348:TRP:CZ3	2:B:438:HIS:HE1	2.36	0.43
1:A:219:LYS:CA	1:A:223:GLN:O	2.64	0.43
1:A:274:ASN:C	1:A:274:ASN:ND2	2.77	0.43
3:C:34:LEU:HB3	3:C:52:CYS:SG	2.58	0.43
1:A:150:TRP:CD1	1:A:192:TYR:OH	2.71	0.43
1:A:289:ASN:OD1	1:A:290:LEU:N	2.49	0.43
4:D:75:LEU:HA	4:D:131:GLN:O	2.19	0.43
1:A:181:HIS:ND1	1:A:213:GLU:HB3	2.34	0.42
4:D:66:ARG:NH2	4:D:111:THR:HG23	2.34	0.42
1:A:112:GLU:HB3	1:A:256:ALA:HB3	2.01	0.42
1:A:119:LYS:HB2	1:A:252:TRP:NE1	2.35	0.42
2:B:365:GLN:O	2:B:369:GLN:HG2	2.19	0.42
3:C:50:LEU:HD22	3:C:124:TYR:CD2	2.50	0.42
3:C:34:LEU:HD23	3:C:52:CYS:SG	2.59	0.42
4:D:103:ARG:HD2	4:D:119:SER:O	2.19	0.42
1:A:91:TYR:CE2	1:A:227:ILE:HG13	2.54	0.42
3:C:62:TYR:CE2	3:C:128:ARG:HD2	2.55	0.42
2:B:355:ASN:C	2:B:357:HIS:H	2.28	0.41
2:B:455:ASN:HB2	2:B:487:PRO:O	2.20	0.41
2:B:481:ASN:HD22	2:B:481:ASN:C	2.29	0.41
1:A:115:GLU:HG3	1:A:252:TRP:O	2.20	0.41
3:C:61:SER:HB3	3:C:136:TYR:CD2	2.55	0.41
2:B:397:PHE:CE1	2:B:404:MET:HE3	2.56	0.41
1:A:191:LEU:O	1:A:192:TYR:HD1	2.02	0.41
1:A:108:VAL:HG22	1:A:109:SER:H	1.85	0.41
1:A:177:TRP:HZ3	1:A:232:THR:HG22	1.86	0.41
1:A:198:TYR:HA	1:A:212:PRO:HD3	2.03	0.41
3:C:42:VAL:HG11	3:C:116:LEU:HD13	2.02	0.41
3:C:132:ILE:HD12	3:C:133:PHE:N	2.36	0.41
1:A:212:PRO:HG2	1:A:247:ASN:HD22	1.86	0.41
1:A:291:PRO:O	1:A:292:PHE:CG	2.74	0.41
2:B:355:ASN:N	2:B:476:MET:HE3	2.35	0.41
1:A:187:ASP:O	1:A:191:LEU:HB2	2.22	0.40
4:D:120:LEU:HD11	4:D:146:VAL:HG11	2.02	0.40
1:A:217:ARG:HB3	1:A:226:ARG:HH21	1.86	0.40
4:D:77:TRP:CE3	4:D:115:LEU:HD22	2.56	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	317/330 (96%)	272 (86%)	43 (14%)	2 (1%)	21	56
2	B	155/173 (90%)	147 (95%)	8 (5%)	0	100	100
3	C	122/125 (98%)	116 (95%)	6 (5%)	0	100	100
4	D	105/108 (97%)	98 (93%)	6 (6%)	1 (1%)	12	45
All	All	699/736 (95%)	633 (91%)	63 (9%)	3 (0%)	30	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	108	VAL
1	A	38	ASN
4	D	72	SER

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	265/292 (91%)	254 (96%)	11 (4%)	26	60
2	B	134/150 (89%)	134 (100%)	0	100	100
3	C	101/102 (99%)	100 (99%)	1 (1%)	68	80
4	D	90/91 (99%)	88 (98%)	2 (2%)	45	71
All	All	590/635 (93%)	576 (98%)	14 (2%)	43	70

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	THR
1	A	90	CYS
1	A	105	LEU
1	A	107	THR
1	A	111	PHE
1	A	132	THR
1	A	133	THR
1	A	226	ARG
1	A	261	SER
1	A	274	ASN
1	A	293	GLN
3	C	142	LEU
4	D	71	ILE
4	D	73	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	HIS
1	A	87	ASN
1	A	180	HIS
1	A	181	HIS
1	A	189	GLN
1	A	195	ASN
1	A	223	GLN
1	A	247	ASN
1	A	274	ASN
1	A	293	GLN
2	B	352	HIS
2	B	354	GLN
2	B	355	ASN
2	B	365	GLN
2	B	452	GLN
2	B	456	ASN
2	B	469	HIS
2	B	481	ASN
3	C	69	GLN
3	C	114	ASN
4	D	48	GLN

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 ⓘ

1 ligand is 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	$\# Z > 2$
5	NAG	A	401	1	14,14,15	0.43	0	17,19,21	1.15	2 (11%)

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
5	NAG	A	401	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	401	NAG	C8-C7-N2	2.35	120.02	116.12
5	A	401	NAG	C2-N2-C7	-2.16	120.01	122.90

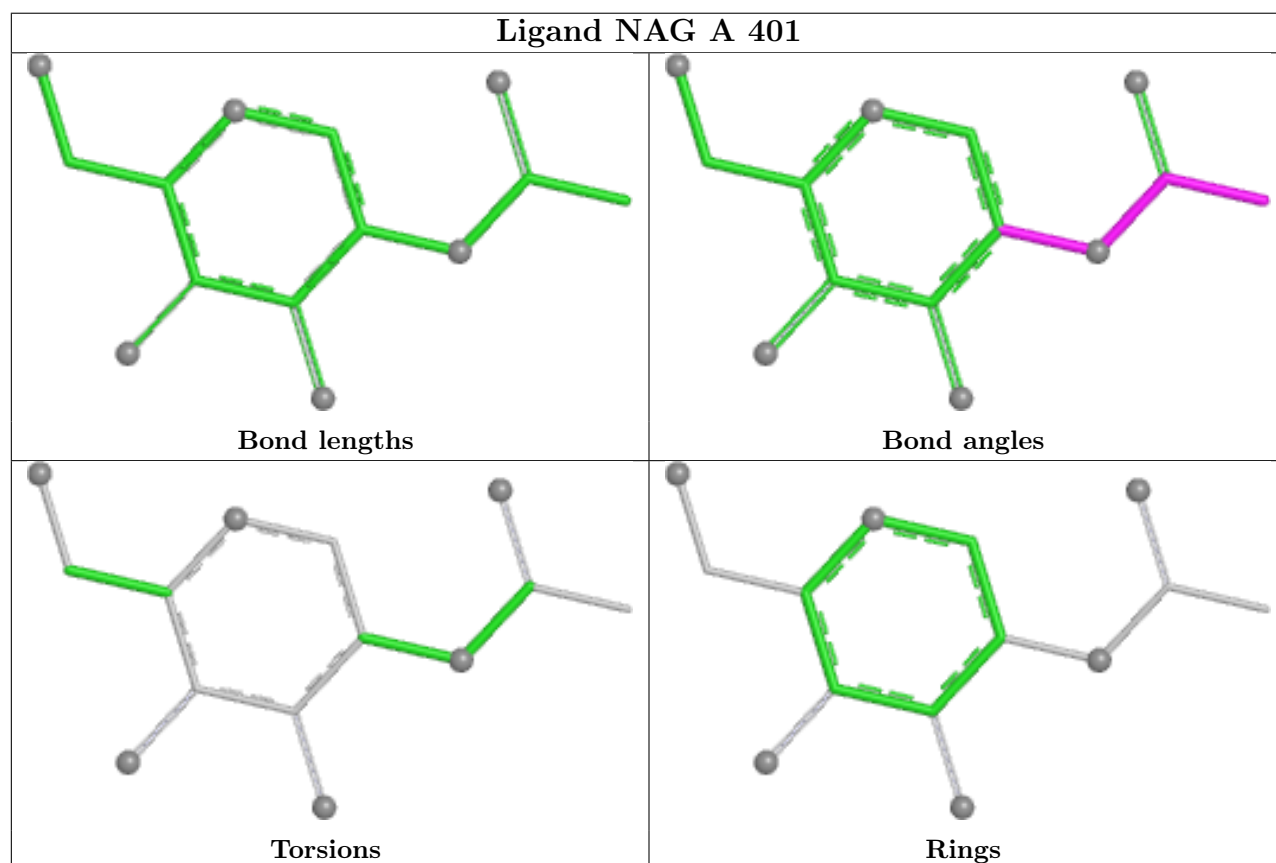
There are no chirality outliers.

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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	321/330 (97%)	0.43	20 (6%) 26 17	10, 63, 119, 147	0
2	B	157/173 (90%)	0.61	18 (11%) 9 7	7, 49, 108, 128	0
3	C	124/125 (99%)	0.67	6 (4%) 35 22	59, 111, 147, 167	0
4	D	107/108 (99%)	0.62	4 (3%) 45 28	58, 103, 146, 156	0
All	All	709/736 (96%)	0.54	48 (6%) 23 15	7, 76, 137, 167	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	440	PHE	4.6
1	A	125	ASN	4.4
1	A	186	ARG	4.4
2	B	486	TYR	4.1
3	C	154	SER	4.0
2	B	334	ALA	3.5
1	A	141	ALA	3.4
4	D	65	CYS	3.2
2	B	461	GLY	3.2
2	B	451	SER	3.1
1	A	155	GLU	3.1
2	B	454	ARG	3.1
1	A	207	TYR	3.1
2	B	333	ILE	3.0
1	A	108	VAL	3.0
2	B	432	GLU	3.0
1	A	139	SER	2.9
1	A	156	ASN	2.9
1	A	127	ASP	2.9
1	A	157	SER	2.8
3	C	132	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	140	GLY	2.8
2	B	450	LYS	2.7
2	B	465	PHE	2.7
3	C	48	LEU	2.7
4	D	111	THR	2.7
2	B	459	GLU	2.6
1	A	262	SER	2.6
3	C	112	GLN	2.6
3	C	131	ALA	2.5
1	A	191	LEU	2.5
4	D	140	PHE	2.4
2	B	453	LEU	2.4
2	B	462	ASN	2.3
1	A	158	TYR	2.3
1	A	161	LEU	2.3
2	B	376	ARG	2.3
2	B	437	TYR	2.2
1	A	154	LYS	2.2
1	A	153	GLN	2.2
3	C	60	SER	2.1
1	A	163	ASN	2.1
1	A	123	TRP	2.1
2	B	464	CYS	2.1
4	D	107	SER	2.0
1	A	192	TYR	2.0
2	B	336	PHE	2.0
2	B	460	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

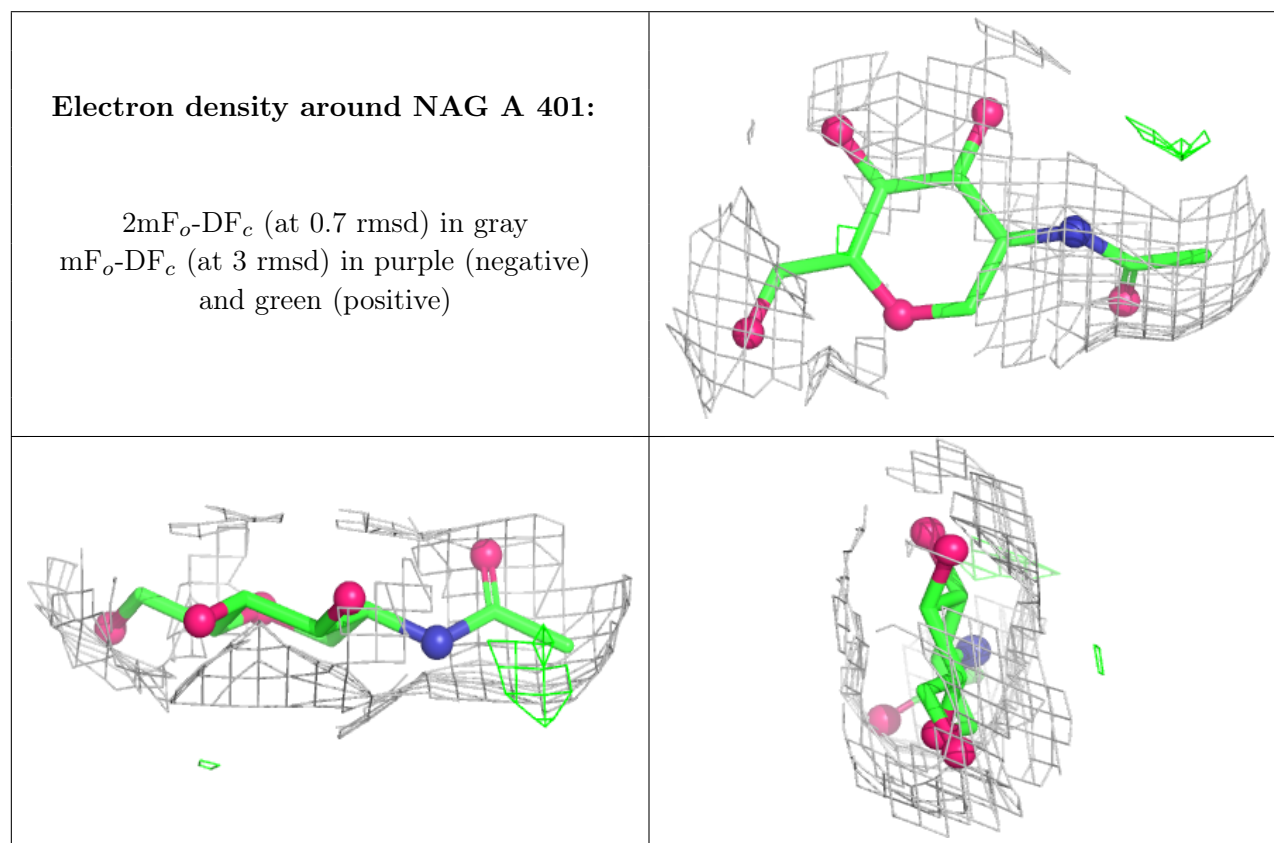
6.4 Ligands [i](#)

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
5	NAG	A	401	14/15	0.60	0.11	24,43,49,54	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.