



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:07 PM GMT

PDB ID : 1NN2
Title : THREE-DIMENSIONAL STRUCTURE OF THE NEURAMINIDASE OF INFLUENZA VIRUS A(SLASH)TOKYO(SLASH)3(SLASH)67 AT 2.2 ANGSTROMS RESOLUTION
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Deposited on : 1991-03-28
Resolution : 2.20 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

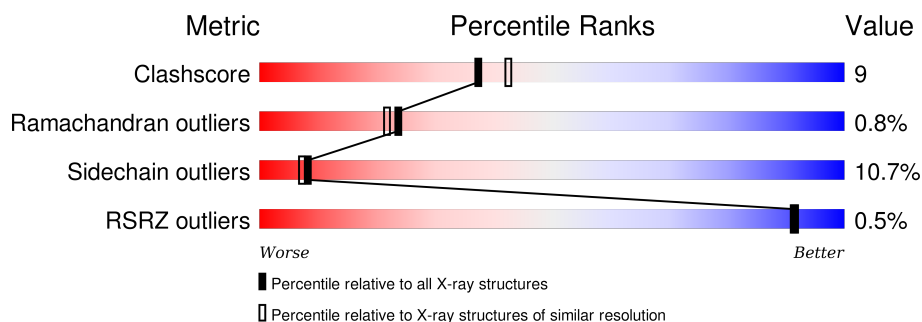
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	388	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NGK	A	477(F)	-	-	-	X
4	NAG	A	479(A)	-	-	-	X
5	NAG	A	485(A)	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 4429 atoms, of which 1100 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 NEURAMINIDASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	388	Total	C	H	N	O	S	0	0	0
			3746	1866	724	545	588	23			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	ASP	ASN	CONFLICT	UNP P06820

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	2	Total	C	H	N	O	28	0
			55	16	27	2	10		

- Molecule 3 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	7	Total	C	H	N	O	S	69	0
			174	50	82	4	37	1		

- Molecule 4 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	6	Total	C	H	N	O	0	0
			139	40	67	2	30		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	0	0
			28	8	14	1	5		
5	A	1	Total	C	H	N	O	28	0
			28	8	14	1	5		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

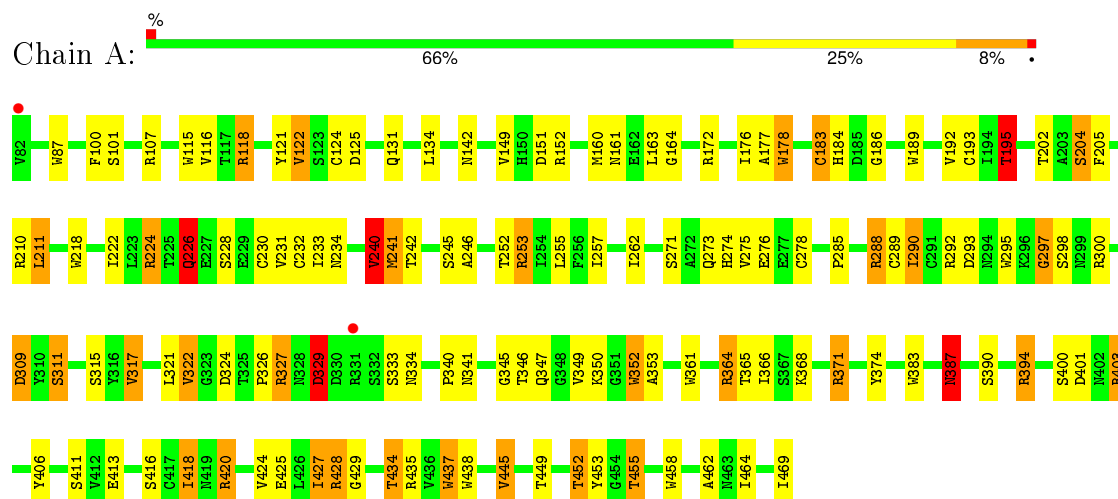
- Molecule 7 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	86	Total	H	O	0	0
			258	172	86		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: NEURAMINIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	139.60Å 139.60Å 191.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.20 9.99 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.20) 63.9 (9.99-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.210 , (Not available) 0.194 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.585	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.4	EDS
Estimated twinning fraction	0.043 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k 0.047 for -1/2*h-1/2*k+1/2*l,-1/2*h-1/2*k-1/2*l,h-k	Xtriage
L-test for twinning ¹	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	0 of 30245 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4429	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: BMA, NAG, CA, NGK, FUL, MAN

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	0.99	3/3092 (0.1%)	1.91	104/4194 (2.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	240	VAL	CA-CB	5.83	1.67	1.54
1	A	122	VAL	CA-CB	5.56	1.66	1.54
1	A	231	VAL	CA-CB	5.20	1.65	1.54

All (104) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	428	ARG	NE-CZ-NH1	12.53	126.56	120.30
1	A	124	CYS	N-CA-CB	-12.16	88.71	110.60
1	A	300	ARG	NE-CZ-NH1	11.14	125.87	120.30
1	A	361	TRP	CD1-CG-CD2	10.44	114.66	106.30
1	A	224	ARG	NE-CZ-NH1	10.42	125.51	120.30
1	A	224	ARG	NE-CZ-NH2	-10.31	115.14	120.30
1	A	253	ARG	NE-CZ-NH1	9.64	125.12	120.30
1	A	300	ARG	NE-CZ-NH2	-9.47	115.57	120.30
1	A	160	MET	CG-SD-CE	9.39	115.23	100.20
1	A	241	MET	CG-SD-CE	-9.34	85.26	100.20
1	A	297	GLY	CA-C-N	-9.26	96.83	117.20
1	A	383	TRP	CD1-CG-CD2	9.20	113.66	106.30
1	A	428	ARG	NE-CZ-NH2	-9.07	115.77	120.30
1	A	437	TRP	CD1-CG-CD2	8.74	113.29	106.30
1	A	327	ARG	NE-CZ-NH1	8.59	124.59	120.30
1	A	172	ARG	NE-CZ-NH2	-8.51	116.05	120.30
1	A	189	TRP	CD1-CG-CD2	8.41	113.03	106.30
1	A	403	ARG	NE-CZ-NH1	8.16	124.38	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	TRP	CD1-CG-CD2	8.06	112.75	106.30
1	A	107	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	A	178	TRP	CE2-CD2-CG	-8.01	100.90	107.30
1	A	371	ARG	NE-CZ-NH1	7.98	124.29	120.30
1	A	364	ARG	NE-CZ-NH1	7.95	124.28	120.30
1	A	455	THR	N-CA-CB	-7.90	95.29	110.30
1	A	361	TRP	CE2-CD2-CG	-7.85	101.02	107.30
1	A	293	ASP	CA-C-N	-7.84	99.96	117.20
1	A	371	ARG	NE-CZ-NH2	-7.80	116.40	120.30
1	A	420	ARG	NE-CZ-NH1	7.64	124.12	120.30
1	A	383	TRP	CE2-CD2-CG	-7.62	101.20	107.30
1	A	115	TRP	CD1-CG-CD2	7.61	112.39	106.30
1	A	87	TRP	CD1-CG-CD2	7.59	112.37	106.30
1	A	152	ARG	NE-CZ-NH1	7.57	124.08	120.30
1	A	189	TRP	CE2-CD2-CG	-7.53	101.28	107.30
1	A	178	TRP	CB-CG-CD1	-7.48	117.27	127.00
1	A	437	TRP	CE2-CD2-CG	-7.39	101.39	107.30
1	A	178	TRP	CG-CD2-CE3	7.34	140.51	133.90
1	A	124	CYS	CB-CA-C	7.33	125.06	110.40
1	A	87	TRP	CE2-CD2-CG	-7.25	101.50	107.30
1	A	438	TRP	CE2-CD2-CG	-7.23	101.52	107.30
1	A	434	THR	N-CA-CB	-7.22	96.58	110.30
1	A	292	ARG	NE-CZ-NH1	7.16	123.88	120.30
1	A	438	TRP	CD1-CG-CD2	7.13	112.01	106.30
1	A	458	TRP	CE2-CD2-CG	-7.12	101.61	107.30
1	A	361	TRP	CG-CD1-NE1	-7.09	103.01	110.10
1	A	295	TRP	CD1-CG-CD2	7.02	111.92	106.30
1	A	297	GLY	O-C-N	6.94	133.80	122.70
1	A	218	TRP	CD1-CG-CD2	6.93	111.84	106.30
1	A	210	ARG	NE-CZ-NH1	6.92	123.76	120.30
1	A	218	TRP	CE2-CD2-CG	-6.87	101.81	107.30
1	A	172	ARG	CG-CD-NE	-6.85	97.42	111.80
1	A	115	TRP	CE2-CD2-CG	-6.68	101.95	107.30
1	A	226	GLN	CA-CB-CG	6.58	127.86	113.40
1	A	458	TRP	CD1-CG-CD2	6.56	111.55	106.30
1	A	253	ARG	NE-CZ-NH2	-6.40	117.10	120.30
1	A	352	TRP	CD1-CG-CD2	6.39	111.41	106.30
1	A	352	TRP	CE2-CD2-CG	-6.28	102.28	107.30
1	A	406	TYR	CB-CG-CD2	-6.21	117.28	121.00
1	A	394	ARG	CG-CD-NE	-6.19	98.80	111.80
1	A	347	GLN	CA-C-N	-6.17	103.85	116.20
1	A	295	TRP	CE2-CD2-CG	-6.15	102.38	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	420	ARG	NE-CZ-NH2	-6.12	117.24	120.30
1	A	383	TRP	CG-CD2-CE3	6.04	139.33	133.90
1	A	107	ARG	NE-CZ-NH2	-6.03	117.28	120.30
1	A	210	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	383	TRP	CB-CG-CD1	-5.95	119.26	127.00
1	A	383	TRP	CG-CD1-NE1	-5.94	104.16	110.10
1	A	288	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	A	290	ILE	N-CA-C	-5.92	95.03	111.00
1	A	438	TRP	CB-CG-CD1	-5.90	119.33	127.00
1	A	121	TYR	CA-CB-CG	5.88	124.57	113.40
1	A	317	VAL	CG1-CB-CG2	-5.86	101.53	110.90
1	A	403	ARG	CB-CG-CD	-5.85	96.38	111.60
1	A	87	TRP	CB-CG-CD1	-5.71	119.57	127.00
1	A	435	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	195	THR	OG1-CB-CG2	5.62	122.94	110.00
1	A	183	CYS	CA-CB-SG	5.56	124.01	114.00
1	A	230	CYS	O-C-N	-5.55	113.81	122.70
1	A	435	ARG	NE-CZ-NH1	5.55	123.07	120.30
1	A	458	TRP	CG-CD2-CE3	5.50	138.85	133.90
1	A	329	ASP	CA-CB-CG	5.49	125.49	113.40
1	A	87	TRP	CG-CD2-CE3	5.49	138.84	133.90
1	A	295	TRP	CG-CD1-NE1	-5.41	104.69	110.10
1	A	458	TRP	CB-CG-CD1	-5.40	119.98	127.00
1	A	276	GLU	N-CA-CB	-5.40	100.89	110.60
1	A	276	GLU	CA-CB-CG	5.33	125.13	113.40
1	A	437	TRP	CG-CD1-NE1	-5.32	104.78	110.10
1	A	118	ARG	CB-CG-CD	5.31	125.41	111.60
1	A	121	TYR	CB-CG-CD2	-5.31	117.81	121.00
1	A	192	VAL	N-CA-C	-5.29	96.72	111.00
1	A	178	TRP	CA-CB-CG	5.28	123.73	113.70
1	A	118	ARG	CA-CB-CG	-5.28	101.79	113.40
1	A	364	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	A	416	SER	N-CA-CB	-5.23	102.66	110.50
1	A	230	CYS	CA-C-N	5.22	128.68	117.20
1	A	293	ASP	CA-C-O	5.18	130.98	120.10
1	A	309	ASP	CB-CA-C	-5.13	100.13	110.40
1	A	211	LEU	N-CA-C	-5.13	97.15	111.00
1	A	438	TRP	CG-CD2-CE3	5.11	138.50	133.90
1	A	210	ARG	CA-CB-CG	5.10	124.62	113.40
1	A	116	VAL	CG1-CB-CG2	-5.07	102.78	110.90
1	A	292	ARG	CA-CB-CG	5.05	124.52	113.40
1	A	387	ASN	N-CA-C	5.04	124.62	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	346	THR	CA-C-N	-5.03	106.13	117.20
1	A	449	THR	CA-CB-CG2	5.01	119.42	112.40

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	3022	724	2853	53	0
2	A	28	27	25	0	0
3	A	92	82	74	0	0
4	A	72	67	61	0	0
5	A	28	28	26	4	0
6	A	1	0	0	0	0
7	A	86	172	0	1	1
All	All	3329	1100	3039	54	1

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 9.

All (54) 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:233:ILE:HG22	1:A:234:ASN:HD22	1.46	0.78
1:A:233:ILE:HG22	1:A:234:ASN:ND2	2.03	0.74
1:A:334:ASN:HA	1:A:387:ASN:HD21	1.53	0.73
1:A:322:VAL:HG12	1:A:327:ARG:HG3	1.71	0.72
1:A:101:SER:HB3	1:A:445:VAL:HG13	1.77	0.67
1:A:226:GLN:HG3	1:A:278:CYS:O	1.95	0.67
1:A:177:ALA:HB2	1:A:193:CYS:HB3	1.76	0.66
1:A:317:VAL:HG23	7:A:551:HOH:O	1.96	0.65
1:A:228:SER:HB3	1:A:350:LYS:HE2	1.79	0.65
1:A:273:GLN:HG3	1:A:340:PRO:HG3	1.79	0.64
1:A:242:THR:HG21	1:A:275:VAL:O	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:PRO:HB2	5:A:485(A):NAG:H82	1.83	0.60
1:A:234:ASN:ND2	5:A:485(A):NAG:C1	2.65	0.60
1:A:309:ASP:HB2	1:A:311:SER:OG	2.02	0.59
1:A:329:ASP:O	1:A:333:SER:HB3	2.01	0.59
1:A:184:HIS:CD2	1:A:186:GLY:H	2.21	0.59
1:A:334:ASN:HA	1:A:387:ASN:ND2	2.19	0.58
1:A:176:ILE:HG22	1:A:195:THR:HG21	1.90	0.53
1:A:245:SER:O	1:A:274:HIS:HE1	1.91	0.53
5:A:485(A):NAG:C1	5:A:485(A):NAG:O7	2.58	0.53
1:A:298:SER:HB2	1:A:341:ASN:HD21	1.74	0.52
1:A:326:PRO:HA	1:A:368:LYS:O	2.09	0.51
1:A:365:THR:HG21	1:A:371:ARG:HA	1.92	0.51
1:A:349:VAL:CG2	1:A:371:ARG:HE	2.25	0.50
1:A:349:VAL:HG23	1:A:371:ARG:NE	2.27	0.50
1:A:131:GLN:NE2	1:A:164:GLY:H	2.11	0.48
1:A:428:ARG:NH2	1:A:462:ALA:O	2.47	0.48
1:A:437:TRP:HD1	1:A:469:ILE:HG22	1.79	0.47
1:A:321:LEU:HD23	1:A:321:LEU:HA	1.76	0.47
1:A:403:ARG:NH2	1:A:429:GLY:O	2.48	0.47
1:A:204:SER:HB3	1:A:211:LEU:HD11	1.97	0.46
1:A:242:THR:HG22	1:A:252:THR:HG23	1.98	0.46
1:A:183:CYS:SG	1:A:232:CYS:SG	3.14	0.45
1:A:240:VAL:HG21	1:A:278:CYS:SG	2.57	0.45
1:A:425:GLU:HG2	1:A:427:ILE:HG22	1.99	0.45
1:A:184:HIS:HD2	1:A:186:GLY:H	1.63	0.45
1:A:349:VAL:HG23	1:A:371:ARG:HE	1.82	0.45
1:A:131:GLN:HE21	1:A:163:LEU:HD12	1.82	0.44
1:A:298:SER:O	1:A:322:VAL:HG13	2.19	0.43
1:A:418:ILE:HD11	1:A:420:ARG:NH1	2.34	0.43
1:A:100:PHE:HB3	1:A:445:VAL:HG22	2.00	0.42
1:A:228:SER:HB3	1:A:350:LYS:CE	2.47	0.42
1:A:278:CYS:HB3	1:A:289:CYS:HB3	2.01	0.42
1:A:246:ALA:O	1:A:274:HIS:NE2	2.53	0.42
1:A:205:PHE:CE1	1:A:262:ILE:HD11	2.55	0.42
1:A:352:TRP:HE1	1:A:374:TYR:HH	1.68	0.42
1:A:352:TRP:NE1	1:A:374:TYR:OH	2.53	0.42
1:A:241:MET:HE3	1:A:255:LEU:HG	2.01	0.41
1:A:290:ILE:HG12	1:A:353:ALA:HB3	2.02	0.41
1:A:411:SER:HB3	1:A:418:ILE:CD1	2.50	0.41
1:A:366:ILE:HG21	1:A:400:SER:HB3	2.03	0.41
1:A:285:PRO:CB	5:A:485(A):NAG:H82	2.50	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:GLY:O	1:A:345:GLY:O	2.39	0.40
1:A:452:THR:HG22	1:A:453:TYR:H	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:490:HOH:H1	7:A:490:HOH:H1[16_665]	0.83	0.77

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	386/388 (100%)	353 (92%)	30 (8%)	3 (1%)	24 22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	329	ASP
1	A	222	ILE
1	A	322	VAL

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	338/338 (100%)	302 (89%)	36 (11%)	8 7

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	ARG
1	A	122	VAL
1	A	125	ASP
1	A	134	LEU
1	A	142	ASN
1	A	149	VAL
1	A	151	ASP
1	A	161	ASN
1	A	178	TRP
1	A	195	THR
1	A	202	THR
1	A	204	SER
1	A	224	ARG
1	A	226	GLN
1	A	240	VAL
1	A	253	ARG
1	A	257	ILE
1	A	271	SER
1	A	288	ARG
1	A	311	SER
1	A	315	SER
1	A	324	ASP
1	A	364	ARG
1	A	387	ASN
1	A	390	SER
1	A	394	ARG
1	A	401	ASP
1	A	413	GLU
1	A	418	ILE
1	A	424	VAL
1	A	427	ILE
1	A	434	THR
1	A	445	VAL
1	A	452	THR
1	A	455	THR
1	A	464	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	GLN
1	A	161	ASN
1	A	226	GLN
1	A	234	ASN
1	A	274	HIS
1	A	334	ASN
1	A	387	ASN
1	A	419	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](#)

15 carbohydrates 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	470(A)	1,2	14,14,15	0.65	0	15,19,21	1.00	1 (6%)
2	NAG	A	471(B)	2	14,14,15	0.66	0	15,19,21	1.39	2 (13%)
3	NAG	A	472(A)	1,3	14,14,15	1.02	1 (7%)	15,19,21	1.96	5 (33%)
3	NAG	A	473(B)	3	14,14,15	1.02	1 (7%)	15,19,21	1.36	2 (13%)
3	BMA	A	474(C)	3	11,11,12	1.00	0	14,15,17	1.52	3 (21%)
3	MAN	A	475(D)	3	11,11,12	0.87	0	14,15,17	1.50	2 (14%)
3	NAG	A	476(E)	3	14,14,15	1.57	3 (21%)	15,19,21	2.87	4 (26%)
3	NGK	A	477(F)	3	18,18,19	1.58	2 (11%)	21,26,28	2.36	6 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	FUL	A	478(I)	3	10,10,11	3.35	4 (40%)	14,14,16	5.52	9 (64%)
4	NAG	A	479(A)	1,4	14,14,15	0.86	1 (7%)	15,19,21	1.21	3 (20%)
4	NAG	A	480(B)	4	14,14,15	1.10	1 (7%)	15,19,21	1.44	4 (26%)
4	BMA	A	481(C)	4	11,11,12	0.81	0	14,15,17	1.17	1 (7%)
4	MAN	A	482(D)	4	11,11,12	0.57	0	14,15,17	1.36	2 (14%)
4	MAN	A	483(E)	4	11,11,12	0.82	0	14,15,17	1.26	1 (7%)
4	MAN	A	484(F)	4	11,11,12	0.68	0	14,15,17	0.99	0

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	470(A)	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	471(B)	2	-	0/6/23/26	0/1/1/1
3	NAG	A	472(A)	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	473(B)	3	-	0/6/23/26	0/1/1/1
3	BMA	A	474(C)	3	-	0/2/19/22	0/1/1/1
3	MAN	A	475(D)	3	-	0/2/19/22	0/1/1/1
3	NAG	A	476(E)	3	-	0/6/23/26	0/1/1/1
3	NGK	A	477(F)	3	-	0/11/28/31	0/1/1/1
3	FUL	A	478(I)	3	-	0/0/17/20	0/1/1/1
4	NAG	A	479(A)	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	480(B)	4	-	0/6/23/26	0/1/1/1
4	BMA	A	481(C)	4	-	0/2/19/22	0/1/1/1
4	MAN	A	482(D)	4	-	0/2/19/22	0/1/1/1
4	MAN	A	483(E)	4	-	0/2/19/22	0/1/1/1
4	MAN	A	484(F)	4	-	0/2/19/22	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	477(F)	NGK	O4-S	-4.44	1.42	1.57
4	A	480(B)	NAG	C1-C2	-3.29	1.47	1.52
4	A	479(A)	NAG	O4-C4	2.05	1.47	1.43
3	A	473(B)	NAG	C4-C5	2.18	1.57	1.53
3	A	476(E)	NAG	O4-C4	2.29	1.48	1.43
3	A	476(E)	NAG	C4-C5	2.44	1.58	1.53
3	A	478(I)	FUL	C6-C5	2.48	1.57	1.51
3	A	472(A)	NAG	C1-C2	2.48	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	478(I)	FUL	C1-C2	2.66	1.58	1.52
3	A	477(F)	NGK	C1-C2	3.61	1.57	1.52
3	A	476(E)	NAG	C1-C2	3.86	1.57	1.52
3	A	478(I)	FUL	C4-C3	4.81	1.65	1.52
3	A	478(I)	FUL	C4-C5	8.40	1.70	1.52

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	478(I)	FUL	O2-C2-C3	-4.48	101.12	110.12
3	A	477(F)	NGK	O7-C7-N	-3.69	114.34	121.86
3	A	476(E)	NAG	C3-C4-C5	-3.57	103.98	110.20
3	A	472(A)	NAG	C4-C3-C2	-3.28	106.13	111.23
3	A	478(I)	FUL	C2-C3-C4	-2.96	106.02	111.04
3	A	476(E)	NAG	C4-C3-C2	-2.70	107.03	111.23
4	A	480(B)	NAG	C3-C2-N2	-2.41	104.78	110.56
4	A	481(C)	BMA	C1-C2-C3	-2.36	106.75	109.54
4	A	479(A)	NAG	O7-C7-C8	-2.33	117.78	122.06
3	A	474(C)	BMA	C2-C3-C4	-2.27	107.18	111.04
4	A	482(D)	MAN	C3-C4-C5	-2.26	106.25	110.20
2	A	470(A)	NAG	C3-C2-N2	-2.21	105.27	110.56
3	A	472(A)	NAG	O7-C7-N2	-2.19	117.40	121.86
3	A	472(A)	NAG	O4-C4-C3	-2.14	105.51	110.34
4	A	480(B)	NAG	O4-C4-C3	-2.13	105.55	110.34
3	A	477(F)	NGK	C4-O4-S	2.08	122.73	118.77
4	A	480(B)	NAG	C8-C7-N2	2.09	120.10	116.11
3	A	473(B)	NAG	C8-C7-N2	2.12	120.16	116.11
3	A	475(D)	MAN	C3-C4-C5	2.13	113.92	110.20
4	A	480(B)	NAG	C1-O5-C5	2.13	114.96	112.25
3	A	477(F)	NGK	O4-S-O3S	2.25	114.16	106.86
4	A	479(A)	NAG	C8-C7-N2	2.25	120.42	116.11
3	A	477(F)	NGK	C2-N-C7	2.30	125.99	123.04
3	A	474(C)	BMA	C3-C4-C5	2.47	114.50	110.20
3	A	473(B)	NAG	C1-O5-C5	2.56	115.50	112.25
4	A	482(D)	MAN	C1-O5-C5	2.63	115.58	112.25
4	A	479(A)	NAG	C1-O5-C5	2.63	115.58	112.25
2	A	471(B)	NAG	C6-C5-C4	2.74	119.77	113.02
3	A	474(C)	BMA	C1-O5-C5	2.87	115.89	112.25
3	A	476(E)	NAG	O4-C4-C5	3.17	117.63	109.24
3	A	472(A)	NAG	C8-C7-N2	3.42	122.64	116.11
4	A	483(E)	MAN	C1-O5-C5	3.65	116.88	112.25
3	A	472(A)	NAG	C1-O5-C5	3.65	116.88	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	478(I)	FUL	C1-O5-C5	3.73	118.14	112.38
2	A	471(B)	NAG	C1-O5-C5	3.77	117.03	112.25
3	A	478(I)	FUL	O4-C4-C3	3.83	118.95	110.34
3	A	475(D)	MAN	C1-O5-C5	3.87	117.16	112.25
3	A	478(I)	FUL	C3-C4-C5	4.44	117.21	109.72
3	A	477(F)	NGK	C8-C7-N	5.78	127.16	116.11
3	A	477(F)	NGK	C1-O5-C5	6.27	120.20	112.25
3	A	478(I)	FUL	O2-C2-C1	7.56	124.36	109.21
3	A	478(I)	FUL	C1-C2-C3	7.95	118.95	109.54
3	A	476(E)	NAG	C1-O5-C5	9.29	124.04	112.25
3	A	478(I)	FUL	C6-C5-C4	9.63	132.03	113.08
3	A	478(I)	FUL	O3-C3-C2	11.44	130.67	110.00

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	A	485(A)	-	14,14,15	0.84	0	15,19,21	1.32	3 (20%)
5	NAG	A	486(B)	-	14,14,15	0.54	0	15,19,21	1.05	1 (6%)

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	485(A)	-	-	0/6/23/26	0/1/1/1
5	NAG	A	486(B)	-	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
5	A	485(A)	NAG	O7-C7-C8	-2.21	118.00	122.06
5	A	485(A)	NAG	C8-C7-N2	2.49	120.87	116.11
5	A	486(B)	NAG	O3-C3-C2	2.55	114.16	109.11
5	A	485(A)	NAG	C6-C5-C4	2.56	119.32	113.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	485(A)	NAG	4	0

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	388/388 (100%)	-0.68	2 (0%) 91 91	10, 25, 42, 60	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	82	VAL	3.2
1	A	331	ARG	2.9

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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	NGK	A	477(F)	18/19	0.69	0.36	11.01	0,93,99,100	0
4	NAG	A	479(A)	14/15	0.86	0.15	2.31	0,0,38,40	0
4	NAG	A	480(B)	14/15	0.89	0.13	0.65	0,0,36,38	0
3	BMA	A	474(C)	11/12	-	-	-	0,0,80,81	21
3	NAG	A	472(A)	14/15	0.86	0.14	-	0,66,74,82	0
3	NAG	A	476(E)	14/15	-	-	-	0,0,92,94	27
3	FUL	A	478(I)	10/11	0.66	0.45	-	0,0,87,89	0
3	NAG	A	473(B)	14/15	0.61	0.28	-	0,0,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	MAN	A	483(E)	11/12	0.87	0.32	-	0,0,61,65	0
3	MAN	A	475(D)	11/12	-	-	-	0,0,84,87	21
4	MAN	A	484(F)	11/12	0.77	0.33	-	0,0,61,63	0
4	MAN	A	482(D)	11/12	0.86	0.12	-	0,0,51,53	0
2	NAG	A	470(A)	14/15	0.80	0.24	-	0,0,54,55	0
2	NAG	A	471(B)	14/15	-	-	-	0,0,55,55	28
4	BMA	A	481(C)	11/12	0.92	0.12	-	0,41,49,54	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
5	NAG	A	485(A)	14/15	0.72	0.24	2.50	0,0,78,79	0
6	CA	A	1	1/1	0.96	0.08	-0.74	45,45,45,45	0
5	NAG	A	486(B)	14/15	-	-	-	0,0,92,92	28

6.5 Other polymers [i](#)

There are no such residues in this entry.