



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 03:13 pm GMT

PDB ID : 2H9Y
Title : Crystal structure of mouse acetylcholinesterase complexed with m-(N,N,N-trimethylammonio)trifluoroacetophenone
Authors : Bourne, Y.; Radic, Z.; Sulzenbacher, G.; Kim, E.; Taylor, P.; Marchot, P.
Deposited on : 2006-06-12
Resolution : 2.40 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

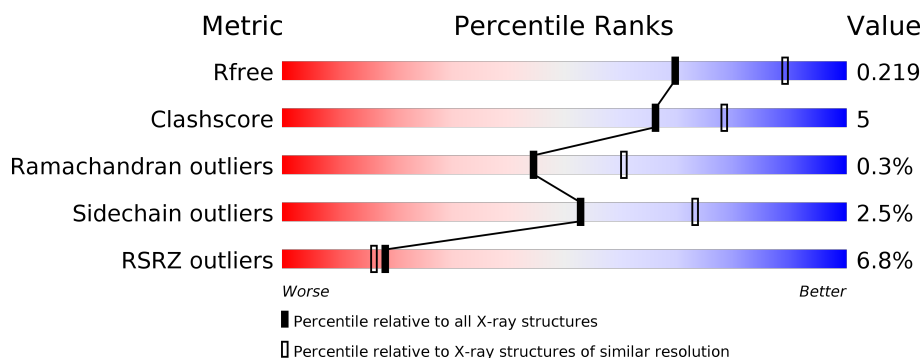
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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	543	<div> <div>6%</div> <div>89%</div> <div>9%</div> <div>..</div> </div>
1	B	543	<div> <div>8%</div> <div>87%</div> <div>10%</div> <div>..</div> </div>

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
2	NAG	A	544	-	-	-	X
2	FUC	A	545	X	-	-	-
5	P6G	A	901	-	-	-	X
6	PG4	A	902	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 8911 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.

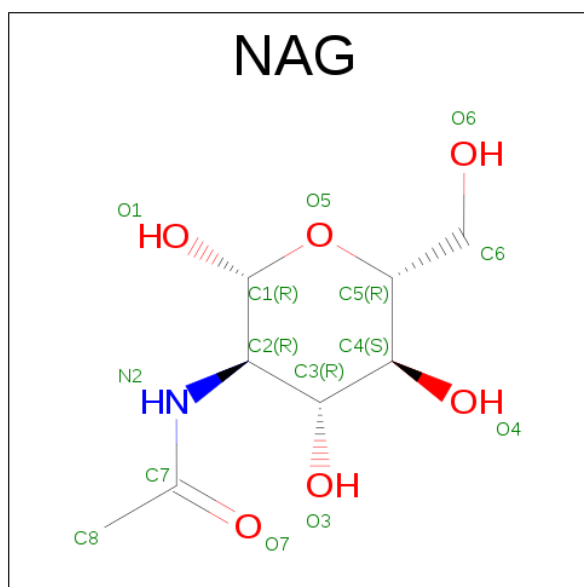
- Molecule 1 is a protein called Acetylcholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	536	Total	C	N	O	S	0	1	0
			4190	2687	729	760	14			
1	B	534	Total	C	N	O	S	0	3	0
			4183	2685	723	761	14			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			24	14	1	9		

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



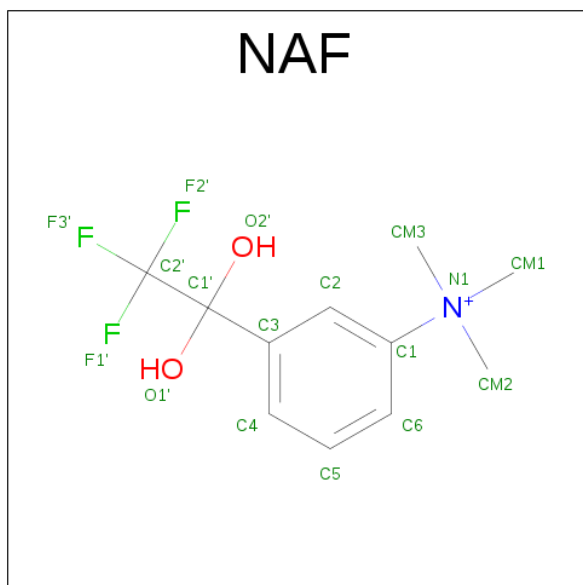
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

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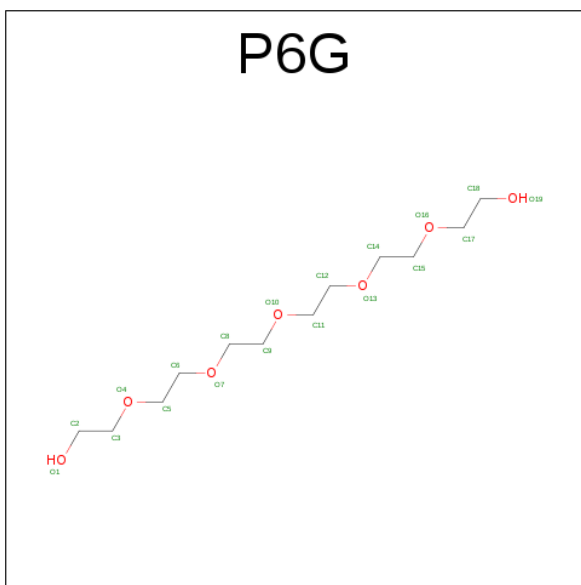
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is M-(N,N,N-TRIMETHYLAMMONIO)-2,2,2-TRIFLUORO-1,1-DIHYDROXY ETHYLBENZENE (three-letter code: NAF) (formula: $C_{11}H_{15}F_3NO_2$).



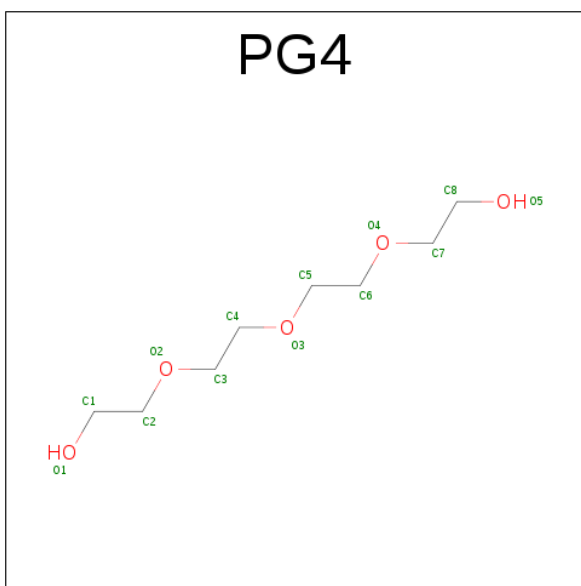
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	F	N	O	0
			16	11	3	1	1	
4	B	1	Total	C	F	N	O	0
			16	11	3	1	1	

- Molecule 5 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $C_{12}H_{26}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			19	12	7		

- Molecule 6 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			13	8	5		

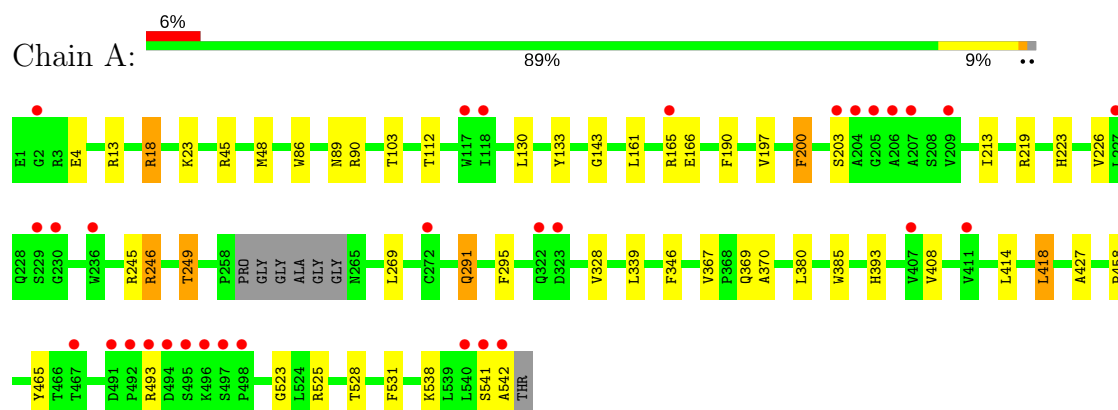
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	237	Total 237	O 237	0	0
7	B	185	Total 185	O 185	0	0

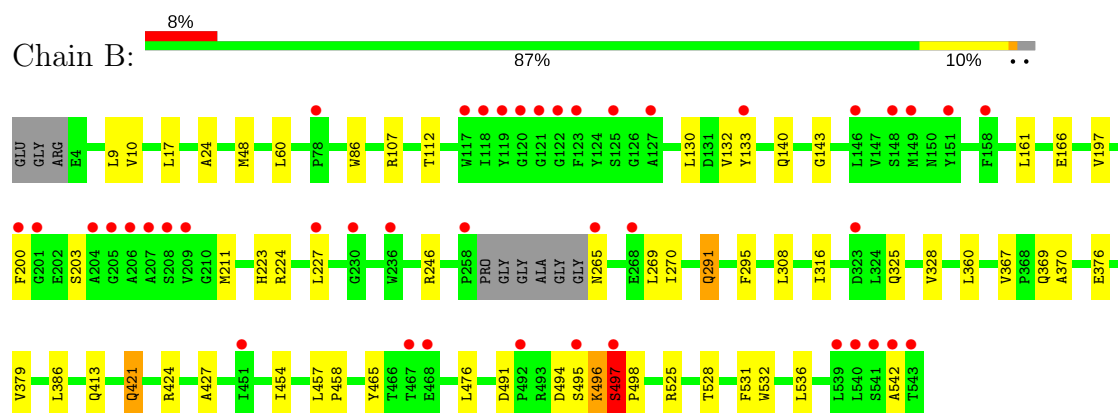
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Acetylcholinesterase



• Molecule 1: Acetylcholinesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.01Å 111.39Å 226.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 39.36 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.00-2.40) 91.0 (39.36-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.91 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.186 , 0.214 0.200 , 0.219	Depositor DCC
R_{free} test set	1564 reflections (2.22%)	DCC
Wilson B-factor (Å ²)	42.2	Xtriage
Anisotropy	0.801	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 52.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8911	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.72% 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: PG4, NAF, NAG, P6G, FUC

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	0.54	0/4319	0.65	3/5901 (0.1%)
1	B	0.48	0/4319	0.62	0/5901
All	All	0.51	0/8638	0.64	3/11802 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
2	A	1	0
All	All	1	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	A	246	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	18	ARG	NE-CZ-NH1	5.00	122.80	120.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	545	FUC	C1

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	496	LYS	Peptide

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	4190	0	4079	39	0
1	B	4183	0	4068	38	0
2	A	24	0	22	0	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
4	A	16	0	14	4	0
4	B	16	0	14	4	0
5	A	19	0	26	0	0
6	A	13	0	18	0	0
7	A	237	0	0	2	0
7	B	185	0	0	1	0
All	All	8911	0	8267	75	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 5.

All (75) 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:B:161:LEU:HD11	1:B:269:LEU:HD22	1.64	0.80
1:A:380:LEU:HD23	1:A:385:TRP:HZ2	1.53	0.73
1:A:291:GLN:HE22	1:A:369:GLN:NE2	1.90	0.68
1:A:245:ARG:O	1:A:249:THR:CG2	2.42	0.67
1:B:166:GLU:HG2	1:B:270:ILE:HD13	1.80	0.64
1:A:380:LEU:HD23	1:A:385:TRP:CZ2	2.34	0.63
1:B:86:TRP:CD2	4:B:952:NAF:HM11	2.34	0.62
1:B:227:LEU:HB2	1:B:328:VAL:HG12	1.81	0.62
1:B:203:SER:HG	4:B:952:NAF:C1'	2.10	0.62
1:A:291:GLN:NE2	1:A:369:GLN:NE2	2.48	0.61
1:B:112:THR:HG21	1:B:143:GLY:O	2.01	0.61
1:A:245:ARG:O	1:A:249:THR:HG22	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:TRP:CD2	4:A:951:NAF:HM11	2.36	0.60
1:A:291:GLN:NE2	1:A:369:GLN:HE21	1.98	0.60
1:A:291:GLN:HE22	1:A:369:GLN:HE21	1.48	0.60
1:B:316:ILE:O	1:B:421:GLN:NE2	2.35	0.60
1:A:4:GLU:OE2	1:A:18:ARG:HD3	2.02	0.59
1:B:367:VAL:HG12	1:B:370:ALA:HB2	1.84	0.58
1:B:497:SER:HB2	1:B:498:PRO:O	2.06	0.56
1:A:203:SER:OG	4:A:951:NAF:C2'	2.53	0.56
1:A:414:LEU:HG	1:A:418:LEU:HD22	1.88	0.55
1:A:203:SER:HG	4:A:951:NAF:C1'	2.13	0.55
1:B:197:VAL:H	1:B:223:HIS:HD2	1.55	0.55
1:A:203:SER:CB	4:A:951:NAF:C1'	2.84	0.55
1:A:245:ARG:O	1:A:249:THR:HG23	2.06	0.55
1:B:291:GLN:HE22	1:B:369:GLN:NE2	2.05	0.54
1:B:24:ALA:HB3	1:B:140:GLN:HG3	1.90	0.54
1:B:203:SER:OG	4:B:952:NAF:C2'	2.55	0.53
1:B:17:LEU:HD23	1:B:60:LEU:HB3	1.91	0.53
1:A:197:VAL:H	1:A:223:HIS:HD2	1.56	0.53
1:A:112:THR:HG21	1:A:143:GLY:O	2.08	0.52
1:A:103:THR:HG21	1:A:190:PHE:HB3	1.91	0.52
1:A:328:VAL:O	1:A:427:ALA:HA	2.08	0.52
1:A:538:LYS:HD3	1:B:376[A]:GLU:HG2	1.91	0.52
1:B:161:LEU:HD11	1:B:269:LEU:CD2	2.39	0.51
1:B:265:ASN:ND2	7:B:1104:HOH:O	2.37	0.50
1:B:161:LEU:CD1	1:B:269:LEU:HD22	2.38	0.49
1:A:408:VAL:HG11	1:A:525:ARG:HG3	1.95	0.49
1:B:497:SER:HB2	1:B:498:PRO:C	2.33	0.49
1:B:211:MET:HG2	1:B:308:LEU:HD21	1.94	0.48
1:B:197:VAL:H	1:B:223:HIS:CD2	2.31	0.48
1:A:197:VAL:H	1:A:223:HIS:CD2	2.31	0.48
1:B:10:VAL:HG22	1:B:107:ARG:NH1	2.28	0.48
1:A:130:LEU:HD12	1:A:133:TYR:CE2	2.48	0.48
1:B:360:LEU:HD22	1:B:379:VAL:HG21	1.95	0.48
1:A:458:PRO:HA	1:A:465:TYR:CD2	2.49	0.47
1:B:491:ASP:HB3	1:B:494:ASP:HB3	1.95	0.47
1:B:166:GLU:HG2	1:B:270:ILE:CD1	2.45	0.46
1:B:130:LEU:HD12	1:B:133:TYR:CE2	2.50	0.46
1:A:48:MET:HE1	1:A:166:GLU:HA	1.98	0.44
1:B:458:PRO:HA	1:B:465:TYR:CD2	2.52	0.44
1:A:523:GLY:HA3	1:B:386:LEU:HD21	1.98	0.44
1:A:541:SER:O	1:A:542:ALA:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:SER:CB	4:B:952:NAF:C1'	2.89	0.44
1:A:291:GLN:HB2	1:A:291:GLN:HE21	1.66	0.43
1:B:457:LEU:N	1:B:458:PRO:CD	2.81	0.43
1:A:89:ASN:O	1:A:90:ARG:NH1	2.48	0.43
1:A:200:PHE:CB	1:A:226:VAL:HB	2.49	0.43
1:A:339:LEU:HD13	1:A:346:PHE:CE2	2.54	0.43
1:B:328:VAL:O	1:B:427:ALA:HA	2.19	0.43
1:B:224:ARG:HG2	1:B:325:GLN:HB2	2.00	0.42
1:A:48:MET:HE1	1:A:165:ARG:O	2.19	0.42
1:A:393:HIS:CD2	7:A:995:HOH:O	2.73	0.42
1:A:45:ARG:HG3	7:A:1164:HOH:O	2.19	0.42
1:B:367:VAL:CG1	1:B:370:ALA:HB2	2.50	0.42
1:B:454:ILE:HD13	1:B:476:LEU:HB3	2.01	0.42
1:B:532:TRP:CE3	1:B:536:LEU:HD12	2.55	0.41
1:A:213:ILE:O	1:A:219:ARG:HD3	2.19	0.41
1:B:291:GLN:HE21	1:B:291:GLN:HB2	1.66	0.41
1:A:161:LEU:HD11	1:A:269:LEU:HD22	2.02	0.41
1:B:161:LEU:HD12	1:B:270:ILE:CG1	2.51	0.41
1:A:528:THR:O	1:A:531:PHE:HB3	2.21	0.41
1:A:367:VAL:HG12	1:A:370:ALA:HB2	2.03	0.41
1:A:269:LEU:C	1:A:269:LEU:HD23	2.42	0.40
1:B:528:THR:O	1:B:531:PHE:HB3	2.21	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	533/543 (98%)	511 (96%)	22 (4%)	0	100	100
1	B	533/543 (98%)	513 (96%)	17 (3%)	3 (1%)	28	41
All	All	1066/1086 (98%)	1024 (96%)	39 (4%)	3 (0%)	44	60

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	497	SER
1	B	542	ALA
1	B	495	SER

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	441/443 (100%)	432 (98%)	9 (2%)	60	79
1	B	442/443 (100%)	429 (97%)	13 (3%)	48	68
All	All	883/886 (100%)	861 (98%)	22 (2%)	53	73

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

Mol	Chain	Res	Type
1	A	13	ARG
1	A	23	LYS
1	A	200	PHE
1	A	246	ARG
1	A	249	THR
1	A	291	GLN
1	A	295	PHE
1	A	418	LEU
1	A	493	ARG
1	B	9	LEU
1	B	48	MET
1	B	132	VAL
1	B	200	PHE
1	B	246	ARG
1	B	291	GLN
1	B	295	PHE
1	B	413	GLN
1	B	421	GLN
1	B	424	ARG
1	B	496	LYS

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Mol	Chain	Res	Type
1	B	497	SER
1	B	525	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	HIS
1	A	287	HIS
1	A	291	GLN
1	B	223	HIS
1	B	291	GLN
1	B	421	GLN
1	B	499	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 ⓘ

2 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	544	1,2	14,14,15	0.59	0	15,19,21	1.61	3 (20%)
2	FUC	A	545	2	9,10,11	0.66	0	13,14,16	1.25	2 (15%)

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	544	1,2	-	0/6/23/26	0/1/1/1
2	FUC	A	545	2	1/1/4/5	0/0/17/20	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	544	NAG	C2-N2-C7	-2.47	119.34	122.94
2	A	545	FUC	C6-C5-C4	-2.18	109.18	113.07
2	A	544	NAG	O5-C1-C2	-2.02	108.66	111.47
2	A	545	FUC	O5-C1-C2	2.01	113.94	110.79
2	A	544	NAG	C1-O5-C5	4.29	118.09	112.17

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	545	FUC	C1

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry ⓘ

6 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	NAG	A	701	1	14,14,15	0.54	0	15,19,21	0.88	0
5	P6G	A	901	-	18,18,18	2.19	6 (33%)	17,17,17	1.22	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	PG4	A	902	-	12,12,12	0.69	0	11,11,11	0.73	0
4	NAF	A	951	1	16,16,17	1.34	2 (12%)	24,25,28	1.29	3 (12%)
3	NAG	B	601	1	14,14,15	0.64	0	15,19,21	1.67	2 (13%)
4	NAF	B	952	1	16,16,17	1.43	2 (12%)	24,25,28	1.57	4 (16%)

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	NAG	A	701	1	-	0/6/23/26	0/1/1/1
5	P6G	A	901	-	-	0/16/16/16	0/0/0/0
6	PG4	A	902	-	-	0/10/10/10	0/0/0/0
4	NAF	A	951	1	-	0/16/16/21	0/1/1/1
3	NAG	B	601	1	-	0/6/23/26	0/1/1/1
4	NAF	B	952	1	-	0/16/16/21	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	952	NAF	C1-N1	-3.35	1.42	1.50
4	A	951	NAF	C1-N1	-3.33	1.42	1.50
4	B	952	NAF	C3-C1'	-3.27	1.48	1.51
4	A	951	NAF	C3-C1'	-2.91	1.48	1.51
5	A	901	P6G	O7-C6	2.83	1.54	1.42
5	A	901	P6G	O19-C18	3.35	1.59	1.42
5	A	901	P6G	O4-C3	3.38	1.56	1.42
5	A	901	P6G	O16-C15	3.52	1.57	1.42
5	A	901	P6G	O13-C12	3.81	1.58	1.42
5	A	901	P6G	O10-C9	4.48	1.61	1.42

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	601	NAG	O5-C1-C2	-3.03	107.26	111.47
4	A	951	NAF	F2'-C2'-C1'	-2.65	105.94	112.28
4	B	952	NAF	C5-C4-C3	-2.36	117.73	120.64
4	B	952	NAF	F1'-C2'-C1'	-2.29	106.80	112.28
4	A	951	NAF	F2'-C2'-F1'	2.85	113.17	106.85
4	A	951	NAF	C4-C3-C1'	2.98	124.15	120.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	901	P6G	O1-C2-C3	3.44	131.66	111.89
4	B	952	NAF	F2'-C2'-F1'	3.80	115.28	106.85
4	B	952	NAF	C4-C3-C1'	4.41	125.77	120.77
3	B	601	NAG	C1-O5-C5	4.82	118.81	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	951	NAF	4	0
4	B	952	NAF	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 ⓘ

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	536/543 (98%)	0.22	31 (5%)	24 22	46, 61, 85, 110	0
1	B	534/543 (98%)	0.41	42 (7%)	13 12	50, 67, 89, 116	0
All	All	1070/1086 (98%)	0.31	73 (6%)	18 16	46, 64, 88, 116	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	495	SER	8.7
1	B	543	THR	6.7
1	B	542	ALA	5.8
1	A	497	SER	5.6
1	B	497	SER	5.5
1	B	540	LEU	5.3
1	B	118	ILE	5.0
1	B	541	SER	4.9
1	B	495	SER	4.8
1	A	494	ASP	4.5
1	B	204	ALA	4.2
1	A	496	LYS	4.1
1	B	117	TRP	4.1
1	A	542	ALA	3.8
1	A	493	ARG	3.8
1	B	268	GLU	3.7
1	B	451	ILE	3.6
1	A	498	PRO	3.5
1	A	272	CYS	3.2
1	A	204	ALA	3.1
1	B	539	LEU	3.1
1	A	492	PRO	3.0
1	B	123	PHE	3.0
1	B	121	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	206	ALA	2.8
1	A	491	ASP	2.8
1	B	205	GLY	2.8
1	A	209	VAL	2.8
1	B	209	VAL	2.8
1	A	117	TRP	2.8
1	B	265	ASN	2.8
1	A	118	ILE	2.8
1	B	206	ALA	2.8
1	A	323	ASP	2.8
1	A	236	TRP	2.7
1	A	205	GLY	2.7
1	A	322	GLN	2.7
1	A	467	THR	2.6
1	B	467	THR	2.6
1	A	2	GLY	2.6
1	A	227	LEU	2.6
1	A	165	ARG	2.6
1	B	133	TYR	2.6
1	B	208	SER	2.6
1	A	230	GLY	2.6
1	B	120	GLY	2.5
1	B	207	ALA	2.5
1	B	119	TYR	2.5
1	B	227	LEU	2.5
1	B	230	GLY	2.4
1	B	125	SER	2.4
1	B	323	ASP	2.4
1	A	540	LEU	2.4
1	B	258	PRO	2.3
1	A	203	SER	2.3
1	B	200	PHE	2.3
1	B	127	ALA	2.3
1	B	149	MET	2.3
1	A	541	SER	2.2
1	A	411	VAL	2.2
1	A	207	ALA	2.2
1	B	78	PRO	2.2
1	A	229	SER	2.2
1	B	151	TYR	2.1
1	A	407	VAL	2.1
1	B	468	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	201	GLY	2.1
1	B	492	PRO	2.1
1	B	148	SER	2.1
1	B	122	GLY	2.1
1	B	236	TRP	2.1
1	B	158	PHE	2.0
1	B	146	LEU	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](#)

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
2	NAG	A	544	14/15	0.86	0.41	3.76	95,99,101,105	0
2	FUC	A	545	10/11	0.85	0.60	-	108,109,111,111	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(Å ²)	Q<0.9
6	PG4	A	902	13/13	0.71	0.25	2.57	87,93,99,99	0
5	P6G	A	901	19/19	0.89	0.20	2.40	76,82,86,87	0
4	NAF	A	951	16/17	0.96	0.20	-0.29	51,55,56,56	0
4	NAF	B	952	16/17	0.96	0.17	-1.36	57,60,61,64	0
3	NAG	B	601	14/15	0.58	0.45	-	96,99,102,102	0
3	NAG	A	701	14/15	0.75	0.41	-	98,101,103,104	0

6.5 Other polymers [i](#)

There are no such residues in this entry.