



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 08:05 PM GMT

PDB ID : 4XII
Title : X-ray structure of human butyrylcholinesterase in complex with N-((1-(2,3-dihydro-1H-inden-2-yl)piperidin-3-yl)methyl)-8-hydroxy-N-(2-methoxyethyl)-5-nitroquinoline-7-carboxamide
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Deposited on : 2015-01-07
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
<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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

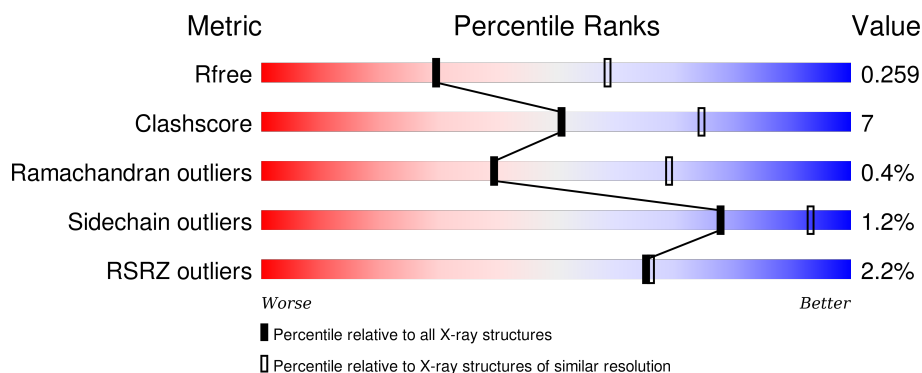
1 Overall quality at a glance

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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (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 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	544	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 85%, yellow 85%, yellow 97%, grey 97%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> % 85% 12% . </div> </div>
1	B	544	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, orange 3%, orange 82%, yellow 82%, yellow 96%, grey 96%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> 3% 82% 14% .. </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
3	NAG	A	1002	X	-	-	-
3	NAG	A	1003	-	-	X	-
3	NAG	A	1011	X	-	-	-
3	NAG	A	1015	X	-	-	-
3	NAG	B	1008	X	-	-	-
4	FUL	A	1005	X	-	-	-
4	FUL	A	1017	-	-	-	X
4	FUL	B	1006	-	-	-	X

2 Entry composition [i](#)

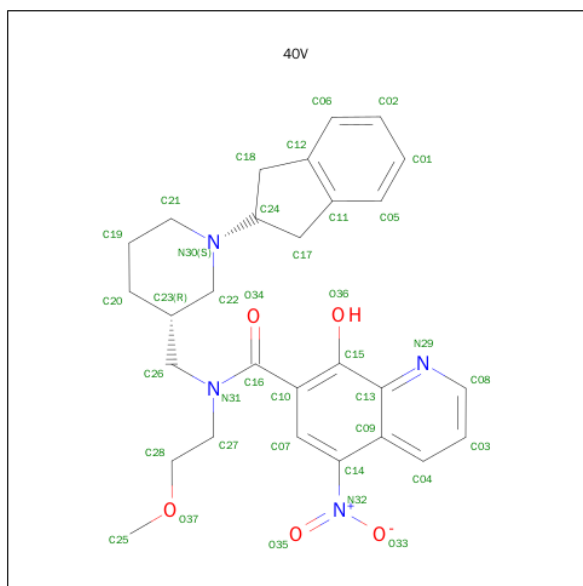
There are 8 unique types of molecules in this entry. The entry contains 8849 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	528	Total	C	N	O	S	0	2	1
			4195	2706	710	764	15			
1	B	525	Total	C	N	O	S	0	1	1
			4168	2693	701	759	15			

- Molecule 2 is N-([(3R)-1-(2,3-dihydro-1H-inden-2-yl)piperidin-3-yl]methyl)-8-hydroxy-N-(2-methoxyethyl)-5-nitroquinoline-7-carboxamide (three-letter code: 40V) (formula: $C_{28}H_{32}N_4O_5$).





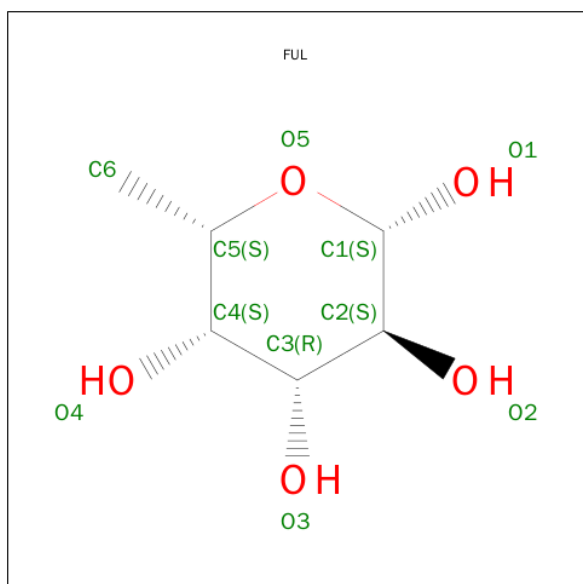
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1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

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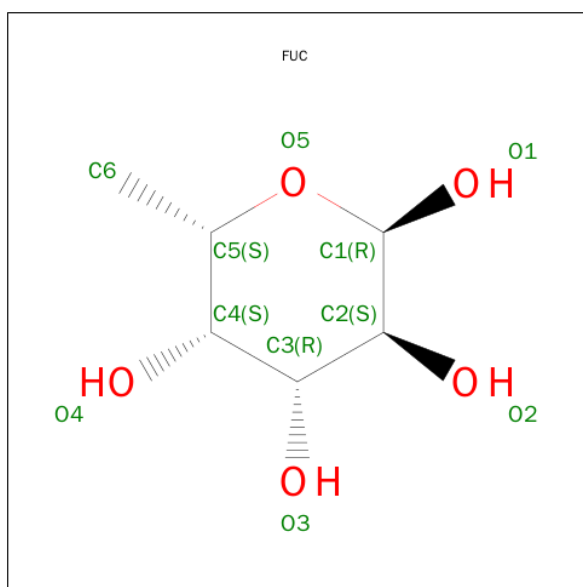
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is BETA-L-FUCOSE (three-letter code: FUL) (formula: $C_6H_{12}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	A	1	Total	C	O	0	0
			10	6	4		
4	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 5 is ALPHA-L-FUCOSE (three-letter code: FUC) (formula: $C_6H_{12}O_5$).

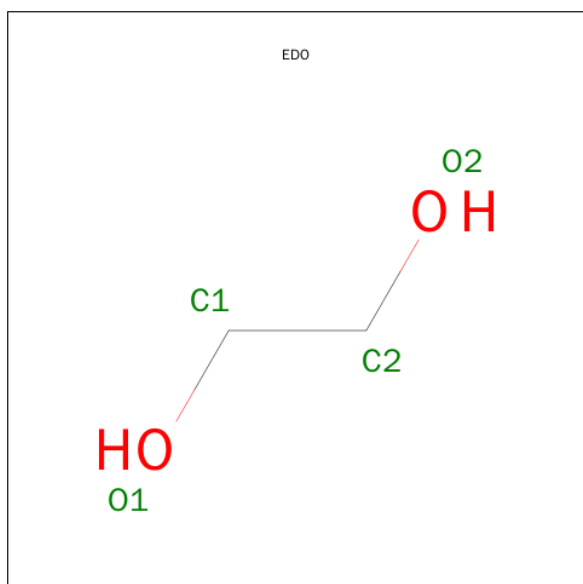


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Cl	0	0
			1	1		
6	A	3	Total	Cl	0	0
			3	3		

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		

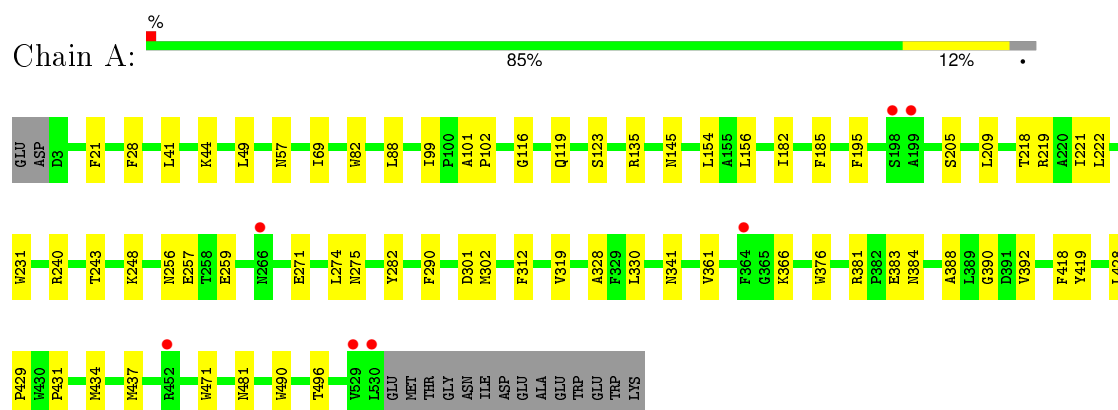
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	35	Total	O	0	0
			35	35		
8	B	27	Total	O	0	0
			27	27		

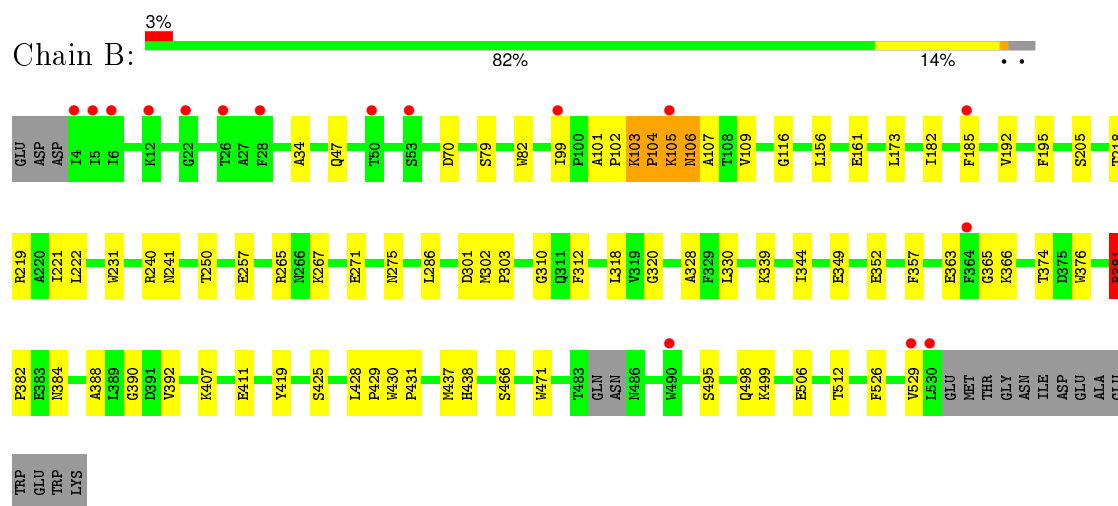
3 Residue-property plots [i](#)

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



• Molecule 1: Cholinesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.04Å 79.76Å 230.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.19 – 2.70 46.76 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (46.19-2.70) 99.0 (46.76-2.70)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.206 , 0.257 0.214 , 0.259	Depositor DCC
R_{free} test set	1509 reflections (3.96%)	DCC
Wilson B-factor (Å ²)	76.8	Xtriage
Anisotropy	0.420	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.9	EDS
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 39583 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8849	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 3.23% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, CL, EDO, 4OV, FUC, FUL

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	0.37	0/4317	0.54	0/5864
1	B	0.35	0/4286	0.54	3/5821 (0.1%)
All	All	0.36	0/8603	0.54	3/11685 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	381	ARG	C-N-CD	-8.37	102.19	120.60
1	B	381	ARG	C-N-CA	6.28	148.38	122.00
1	B	103	LYS	C-N-CD	5.07	139.04	128.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4195	0	4083	42	0
1	B	4168	0	4049	60	0
2	A	37	0	32	4	0
2	B	37	0	32	5	0
3	A	168	0	149	17	0
3	B	112	0	103	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	30	0	30	7	0
4	B	10	0	10	0	0
5	A	10	0	10	3	0
6	A	3	0	0	1	0
6	B	1	0	0	0	0
7	A	8	0	12	0	0
7	B	8	0	12	0	0
8	A	35	0	0	1	0
8	B	27	0	0	0	0
All	All	8849	0	8522	119	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 7.

All (119) 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:57:ASN:HD21	3:A:1003:NAG:C1	1.01	1.55
1:B:241:ASN:HD21	3:B:1005:NAG:C1	1.14	1.52
1:B:106:ASN:OD1	3:B:1004:NAG:C1	1.64	1.44
3:A:1008:NAG:O6	4:A:1010:FUL:C1	1.69	1.40
1:A:57:ASN:ND2	3:A:1003:NAG:C1	1.82	1.40
1:B:241:ASN:ND2	3:B:1005:NAG:C1	1.88	1.37
3:A:1015:NAG:O6	4:A:1017:FUL:C1	1.88	1.20
3:A:1003:NAG:O4	3:A:1004:NAG:H62	1.56	1.01
3:A:1015:NAG:HO6	4:A:1017:FUL:C1	1.70	1.01
3:A:1008:NAG:C6	4:A:1010:FUL:C1	2.45	0.95
3:A:1008:NAG:HO6	4:A:1010:FUL:C1	1.85	0.88
1:A:116:GLY:O	8:A:1101:HOH:O	1.94	0.84
1:B:241:ASN:CG	3:B:1005:NAG:C1	2.47	0.82
1:B:240:ARG:NH2	1:B:257:GLU:OE2	2.13	0.81
3:A:1011:NAG:O6	5:A:1012:FUC:O5	2.02	0.77
1:A:156:LEU:HD13	1:A:257:GLU:HB3	1.74	0.70
3:A:1008:NAG:O6	4:A:1010:FUL:C2	2.41	0.68
1:B:101:ALA:HA	1:B:102:PRO:C	2.17	0.65
3:A:1011:NAG:C6	5:A:1012:FUC:O5	2.44	0.65
1:A:381:ARG:O	1:A:384:ASN:ND2	2.31	0.64
1:B:241:ASN:OD1	3:B:1005:NAG:C1	2.47	0.63
3:A:1003:NAG:O4	3:A:1004:NAG:C6	2.28	0.62
1:A:123:SER:OG	1:A:145:ASN:OD1	2.10	0.59
3:A:1011:NAG:H62	5:A:1012:FUC:O5	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:256:ASN:OD1	1:A:259:GLU:N	2.37	0.57
1:A:154:LEU:HD11	1:A:243:THR:HG23	1.85	0.57
1:A:271:GLU:O	1:A:275:ASN:ND2	2.36	0.57
1:B:428:LEU:HD23	1:B:430:TRP:H	1.70	0.57
1:B:330:LEU:HD11	1:B:390:GLY:HA2	1.86	0.57
1:A:381:ARG:HH11	1:A:383:GLU:HB2	1.71	0.56
1:B:376:TRP:CH2	1:B:384:ASN:HB3	2.41	0.56
1:B:526:PHE:O	1:B:529:VAL:HG12	2.05	0.55
1:B:99:ILE:HD11	1:B:185:PHE:HB3	1.89	0.55
1:B:407:LYS:O	1:B:411:GLU:HG2	2.06	0.55
3:A:1008:NAG:O6	4:A:1010:FUL:O2	2.25	0.54
1:A:481:ASN:HA	6:A:1019:CL:CL	2.45	0.54
1:B:116:GLY:HA2	2:B:1001:40V:O34	2.07	0.53
1:B:376:TRP:CZ3	1:B:384:ASN:HB3	2.45	0.52
1:A:82:TRP:CE2	2:A:1001:40V:H13	2.44	0.52
3:A:1003:NAG:O3	3:A:1004:NAG:C1	2.58	0.51
1:A:240:ARG:NH1	1:A:257:GLU:OE2	2.36	0.51
1:A:381:ARG:HB3	1:A:384:ASN:ND2	2.26	0.51
1:B:105:LYS:CD	1:B:105:LYS:N	2.73	0.51
1:A:428:LEU:HD12	1:A:429:PRO:HD2	1.91	0.51
1:B:156:LEU:HD13	1:B:257:GLU:HB3	1.92	0.50
1:A:376:TRP:CZ3	1:A:384:ASN:HB3	2.46	0.50
1:B:438:HIS:CD2	2:B:1001:40V:H28	2.46	0.50
1:B:34:ALA:HB2	1:B:173:LEU:HD23	1.94	0.49
1:B:271:GLU:O	1:B:275:ASN:ND2	2.36	0.49
1:B:82:TRP:CE2	2:B:1001:40V:H13	2.48	0.49
1:A:69:ILE:HD11	1:A:88:LEU:HD11	1.94	0.49
1:A:205:SER:HB3	1:A:222:LEU:HD21	1.95	0.49
1:B:388:ALA:O	1:B:392:VAL:HG23	2.13	0.48
1:B:363:GLU:HA	1:B:366:LYS:HE3	1.94	0.48
3:A:1003:NAG:O3	3:A:1004:NAG:H2	2.14	0.48
1:B:103:LYS:HG3	1:B:104:PRO:HD2	1.96	0.48
1:A:431:PRO:HG2	1:A:434:MET:HG3	1.96	0.47
1:B:374:THR:HG22	1:B:376:TRP:CZ2	2.50	0.46
1:B:349:GLU:HA	1:B:352:GLU:OE1	2.14	0.46
1:B:104:PRO:C	1:B:105:LYS:HD3	2.35	0.46
1:B:320:GLY:HA3	1:B:419:TYR:CD2	2.50	0.46
1:B:381:ARG:HB2	1:B:384:ASN:ND2	2.31	0.46
1:A:361:VAL:O	1:A:366:LYS:NZ	2.44	0.46
1:B:105:LYS:HD3	1:B:105:LYS:N	2.31	0.46
1:A:49:LEU:HA	1:A:49:LEU:HD12	1.78	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:ASP:OD1	1:B:302:MET:N	2.45	0.45
1:A:419:TYR:HB3	1:A:490:TRP:CZ2	2.51	0.45
1:B:250:THR:HB	1:B:267:LYS:HE2	1.98	0.45
1:B:218:THR:HG22	1:B:219:ARG:HG3	1.99	0.45
1:A:44:LYS:HB3	1:A:44:LYS:HE3	1.85	0.45
1:B:428:LEU:HA	1:B:429:PRO:HD3	1.74	0.44
1:A:388:ALA:O	1:A:392:VAL:HG23	2.17	0.44
1:A:209:LEU:HD23	1:A:312:PHE:HB3	2.00	0.44
1:A:328:ALA:HB2	1:A:437:MET:HE3	2.00	0.44
1:B:328:ALA:HB2	1:B:437:MET:HE3	1.99	0.44
1:A:330:LEU:HD11	1:A:390:GLY:HA2	1.99	0.43
1:A:182:ILE:HA	1:A:182:ILE:HD12	1.79	0.43
3:A:1003:NAG:O3	3:A:1004:NAG:C2	2.67	0.43
1:B:381:ARG:HB3	1:B:382:PRO:CA	2.47	0.43
1:A:428:LEU:HA	1:A:429:PRO:HD3	1.82	0.43
1:B:310:GLY:HA2	1:B:312:PHE:CE2	2.53	0.43
1:A:21:PHE:O	1:A:135:ARG:NH2	2.51	0.43
1:A:119:GLN:NE2	1:A:290:PHE:O	2.47	0.43
1:B:101:ALA:HA	1:B:102:PRO:O	2.18	0.43
1:B:231:TRP:CE2	2:B:1001:40V:H21	2.54	0.43
1:A:28:PHE:HE1	1:A:99:ILE:HD12	1.83	0.43
1:B:104:PRO:CG	1:B:107:ALA:HB2	2.49	0.43
1:B:109:VAL:HB	1:B:192:VAL:HG22	2.00	0.43
1:A:218:THR:HG22	1:A:219:ARG:HG3	2.01	0.43
1:B:320:GLY:HA3	1:B:419:TYR:CE2	2.53	0.43
1:B:195:PHE:CB	1:B:221:ILE:HB	2.49	0.43
1:A:301:ASP:OD1	1:A:302:MET:N	2.51	0.42
1:B:106:ASN:HD22	1:B:106:ASN:HA	1.59	0.42
2:B:1001:40V:H5	2:B:1001:40V:H23	1.61	0.42
1:B:495:SER:O	1:B:498:GLN:NE2	2.40	0.42
1:A:82:TRP:CD2	2:A:1001:40V:H13	2.55	0.42
1:A:231:TRP:CE2	2:A:1001:40V:H21	2.55	0.42
1:B:103:LYS:HG3	1:B:104:PRO:CD	2.49	0.42
1:A:101:ALA:HA	1:A:102:PRO:C	2.39	0.42
1:B:499:LYS:HG2	1:B:512:THR:HG22	2.02	0.41
1:B:161:GLU:OE1	1:B:265:ARG:NH1	2.32	0.41
1:B:286:LEU:HD12	1:B:357:PHE:CE1	2.55	0.41
1:B:425:SER:HB3	1:B:428:LEU:HB2	2.02	0.41
1:B:109:VAL:HG21	1:B:182:ILE:HG12	2.03	0.41
1:B:99:ILE:HD11	1:B:185:PHE:HD2	1.85	0.41
1:B:365:GLY:CA	1:B:529:VAL:HG21	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:365:GLY:HA3	1:B:529:VAL:HG21	2.03	0.41
1:B:221:ILE:HG23	1:B:318:LEU:HB3	2.03	0.41
1:A:41:LEU:HA	1:A:41:LEU:HD23	1.80	0.41
1:A:195:PHE:CB	1:A:221:ILE:HB	2.51	0.41
1:B:344:ILE:HD13	1:B:382:PRO:HG2	2.03	0.41
2:A:1001:40V:H7	2:A:1001:40V:C07	2.51	0.40
1:B:302:MET:HA	1:B:303:PRO:HD3	1.90	0.40
1:A:99:ILE:HD11	1:A:185:PHE:CD2	2.56	0.40
1:B:339:LYS:HA	1:B:431:PRO:HG3	2.02	0.40
1:B:205:SER:HB3	1:B:222:LEU:HD21	2.02	0.40
1:B:70:ASP:OD2	1:B:79:SER:HB2	2.21	0.40
1:A:381:ARG:HB3	1:A:384:ASN:HD21	1.86	0.40
1:A:319:VAL:O	1:A:418:PHE:HA	2.22	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	528/544 (97%)	503 (95%)	24 (4%)	1 (0%)	52	80
1	B	522/544 (96%)	496 (95%)	23 (4%)	3 (1%)	30	59
All	All	1050/1088 (96%)	999 (95%)	47 (4%)	4 (0%)	39	69

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	381	ARG
1	B	506	GLU
1	A	282	TYR
1	B	104	PRO

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	450/467 (96%)	445 (99%)	5 (1%)	80	94
1	B	445/467 (95%)	439 (99%)	6 (1%)	76	92
All	All	895/934 (96%)	884 (99%)	11 (1%)	78	93

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	248	LYS
1	A	274	LEU
1	A	341	ASN
1	A	471	TRP
1	A	496	THR
1	B	47	GLN
1	B	105	LYS
1	B	106	ASN
1	B	381	ARG
1	B	466	SER
1	B	471	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	384	ASN
1	B	241	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 35 ligands modelled in this entry, 4 are monoatomic - leaving 31 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	$\# Z > 2$
2	40V	A	1001	-	38,41,41	5.44	26 (68%)	48,58,58	1.59	10 (20%)
3	NAG	A	1002	1	14,14,15	0.27	0	15,19,21	0.57	0
3	NAG	A	1003	3,4	14,14,15	0.39	0	15,19,21	0.54	0
3	NAG	A	1004	3	14,14,15	0.39	0	15,19,21	1.13	2 (13%)
4	FUL	A	1005	3	10,10,11	0.30	0	13,14,16	0.77	0
3	NAG	A	1006	1,3	14,14,15	0.24	0	15,19,21	0.37	0
3	NAG	A	1007	3	14,14,15	0.20	0	15,19,21	0.46	0
3	NAG	A	1008	1,3	14,14,15	0.55	0	15,19,21	0.57	0
3	NAG	A	1009	3	14,14,15	0.25	0	15,19,21	0.37	0
4	FUL	A	1010	-	10,10,11	0.30	0	13,14,16	0.79	0
3	NAG	A	1011	1	14,14,15	0.34	0	15,19,21	0.78	0
5	FUC	A	1012	-	10,10,11	0.52	0	13,14,16	1.38	2 (15%)
3	NAG	A	1013	1,3	14,14,15	0.38	0	15,19,21	0.66	1 (6%)
3	NAG	A	1014	3	14,14,15	0.48	0	15,19,21	0.37	0
3	NAG	A	1015	1,3	14,14,15	0.46	0	15,19,21	0.38	0
3	NAG	A	1016	3	14,14,15	0.24	0	15,19,21	0.43	0
4	FUL	A	1017	-	10,10,11	0.31	0	13,14,16	0.77	0
7	EDO	A	1021	-	3,3,3	0.50	0	2,2,2	0.38	0
7	EDO	A	1022	-	3,3,3	0.51	0	2,2,2	0.37	0
2	40V	B	1001	-	38,41,41	5.46	27 (71%)	48,58,58	1.60	10 (20%)
3	NAG	B	1002	1	14,14,15	0.28	0	15,19,21	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	1003	1	14,14,15	0.28	0	15,19,21	0.54	0
3	NAG	B	1004	-	14,14,15	0.35	0	15,19,21	0.81	0
3	NAG	B	1005	-	14,14,15	0.35	0	15,19,21	0.76	1 (6%)
4	FUL	B	1006	-	10,10,11	0.28	0	13,14,16	0.78	0
3	NAG	B	1007	1	14,14,15	0.28	0	15,19,21	0.53	0
3	NAG	B	1008	1	14,14,15	0.35	0	15,19,21	0.81	0
3	NAG	B	1009	1	14,14,15	0.27	0	15,19,21	0.53	0
3	NAG	B	1010	1	14,14,15	0.28	0	15,19,21	0.54	0
7	EDO	B	1012	-	3,3,3	0.46	0	2,2,2	0.42	0
7	EDO	B	1013	-	3,3,3	0.54	0	2,2,2	0.43	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	40V	A	1001	-	-	0/24/42/42	0/5/5/5
3	NAG	A	1002	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	1003	3,4	-	0/6/23/26	0/1/1/1
3	NAG	A	1004	3	-	0/6/23/26	0/1/1/1
4	FUL	A	1005	3	1/1/4/5	0/0/17/20	0/1/1/1
3	NAG	A	1006	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1007	3	-	0/6/23/26	0/1/1/1
3	NAG	A	1008	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1009	3	-	0/6/23/26	0/1/1/1
4	FUL	A	1010	-	-	0/0/17/20	0/1/1/1
3	NAG	A	1011	1	1/1/5/7	0/6/23/26	0/1/1/1
5	FUC	A	1012	-	-	0/0/17/20	0/1/1/1
3	NAG	A	1013	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1014	3	-	0/6/23/26	0/1/1/1
3	NAG	A	1015	1,3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	1016	3	-	0/6/23/26	0/1/1/1
4	FUL	A	1017	-	-	0/0/17/20	0/1/1/1
7	EDO	A	1021	-	-	0/1/1/1	0/0/0/0
7	EDO	A	1022	-	-	0/1/1/1	0/0/0/0
2	40V	B	1001	-	-	0/24/42/42	0/5/5/5
3	NAG	B	1002	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1003	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1004	-	-	0/6/23/26	0/1/1/1
3	NAG	B	1005	-	-	0/6/23/26	0/1/1/1
4	FUL	B	1006	-	-	0/0/17/20	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1007	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1008	1	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	B	1009	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1010	1	-	0/6/23/26	0/1/1/1
7	EDO	B	1012	-	-	0/1/1/1	0/0/0/0
7	EDO	B	1013	-	-	0/1/1/1	0/0/0/0

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	40V	C24-N30	-10.06	1.22	1.48
2	B	1001	40V	C24-N30	-10.04	1.22	1.48
2	A	1001	40V	C19-C20	-4.07	1.42	1.53
2	B	1001	40V	C19-C20	-3.70	1.43	1.53
2	A	1001	40V	C20-C23	-3.06	1.43	1.52
2	B	1001	40V	C20-C23	-2.99	1.43	1.52
2	B	1001	40V	C22-N30	-2.54	1.43	1.47
2	A	1001	40V	C26-N31	-2.52	1.42	1.47
2	B	1001	40V	C21-N30	-2.47	1.42	1.47
2	B	1001	40V	C26-N31	-2.35	1.43	1.47
2	A	1001	40V	C22-N30	-2.19	1.43	1.47
2	B	1001	40V	O36-C15	2.09	1.42	1.35
2	A	1001	40V	O36-C15	2.18	1.43	1.35
2	A	1001	40V	C18-C24	2.29	1.58	1.54
2	B	1001	40V	C18-C24	2.58	1.58	1.54
2	A	1001	40V	C04-C09	2.68	1.47	1.42
2	A	1001	40V	C17-C24	2.89	1.59	1.54
2	B	1001	40V	C17-C24	2.96	1.59	1.54
2	B	1001	40V	C04-C09	3.01	1.48	1.42
2	B	1001	40V	C09-C13	3.10	1.50	1.42
2	A	1001	40V	C09-C13	3.32	1.51	1.42
2	A	1001	40V	C10-C16	4.52	1.57	1.50
2	B	1001	40V	C10-C16	4.69	1.58	1.50
2	B	1001	40V	C03-C08	5.08	1.52	1.37
2	A	1001	40V	C03-C08	5.08	1.52	1.37
2	A	1001	40V	C15-C13	5.17	1.51	1.42
2	B	1001	40V	C15-C13	5.17	1.51	1.42
2	A	1001	40V	C16-N31	6.07	1.48	1.34
2	B	1001	40V	C16-N31	6.22	1.48	1.34
2	A	1001	40V	C03-C04	6.34	1.51	1.36
2	B	1001	40V	C02-C01	6.48	1.54	1.38
2	B	1001	40V	C03-C04	6.49	1.51	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	40V	C02-C01	6.52	1.54	1.38
2	A	1001	40V	C12-C11	6.70	1.52	1.39
2	A	1001	40V	C08-N29	6.78	1.46	1.32
2	A	1001	40V	C07-C10	6.82	1.51	1.39
2	B	1001	40V	C12-C11	6.89	1.52	1.39
2	A	1001	40V	C14-C09	6.93	1.61	1.43
2	B	1001	40V	C08-N29	6.99	1.46	1.32
2	B	1001	40V	C14-C09	7.01	1.61	1.43
2	B	1001	40V	C07-C10	7.03	1.51	1.39
2	B	1001	40V	C01-C05	7.43	1.53	1.38
2	A	1001	40V	C01-C05	7.52	1.53	1.38
2	B	1001	40V	C06-C12	7.76	1.53	1.39
2	A	1001	40V	C06-C12	7.84	1.53	1.39
2	A	1001	40V	C02-C06	8.23	1.54	1.38
2	B	1001	40V	C02-C06	8.36	1.55	1.38
2	B	1001	40V	C05-C11	9.03	1.55	1.39
2	A	1001	40V	C05-C11	9.12	1.55	1.39
2	B	1001	40V	C13-N29	10.08	1.50	1.37
2	A	1001	40V	C13-N29	11.14	1.51	1.37
2	A	1001	40V	C10-C15	12.91	1.58	1.39
2	B	1001	40V	C10-C15	13.65	1.59	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	40V	C17-C11-C12	-3.22	108.26	110.65
2	A	1001	40V	C03-C08-N29	-3.10	119.01	123.94
2	B	1001	40V	C17-C11-C12	-3.01	108.42	110.65
2	A	1001	40V	C28-C27-N31	-2.86	106.96	113.39
2	B	1001	40V	C28-C27-N31	-2.86	106.97	113.39
2	A	1001	40V	C21-N30-C24	-2.83	107.36	112.63
2	A	1001	40V	C07-C14-C09	-2.77	118.78	122.81
2	B	1001	40V	C03-C08-N29	-2.64	119.75	123.94
2	A	1001	40V	C18-C12-C11	-2.64	108.70	110.65
2	B	1001	40V	C18-C12-C11	-2.57	108.75	110.65
2	B	1001	40V	C19-C21-N30	-2.47	106.97	111.13
2	B	1001	40V	C07-C14-C09	-2.29	119.47	122.81
3	B	1005	NAG	C3-C4-C5	2.01	113.80	110.23
3	A	1013	NAG	C1-O5-C5	2.13	115.27	112.14
2	B	1001	40V	C21-N30-C24	2.26	116.84	112.63
3	A	1004	NAG	O5-C5-C4	2.29	113.93	110.13
5	A	1012	FUC	C3-C4-C5	2.43	113.28	109.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	40V	C20-C23-C22	2.44	111.67	108.59
3	A	1004	NAG	C4-C3-C2	2.51	115.24	111.34
2	B	1001	40V	C20-C23-C22	2.78	112.10	108.59
2	A	1001	40V	C07-C10-C15	2.92	120.77	118.82
2	A	1001	40V	C21-N30-C22	3.23	111.20	108.13
5	A	1012	FUC	C1-C2-C3	3.43	113.71	109.55
2	B	1001	40V	C27-N31-C26	3.80	122.79	117.22
2	A	1001	40V	C27-N31-C26	4.04	123.14	117.22
2	B	1001	40V	C21-N30-C22	4.57	112.48	108.13

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	1005	FUL	C1
3	A	1002	NAG	C1
3	A	1015	NAG	C1
3	B	1008	NAG	C1
3	A	1011	NAG	C1

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	40V	4	0
3	A	1003	NAG	7	0
3	A	1004	NAG	5	0
3	A	1008	NAG	5	0
4	A	1010	FUL	5	0
3	A	1011	NAG	3	0
5	A	1012	FUC	3	0
3	A	1015	NAG	2	0
4	A	1017	FUL	2	0
2	B	1001	40V	5	0
3	B	1004	NAG	1	0
3	B	1005	NAG	4	0

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	528/544 (97%)	-0.04	7 (1%) 79 79	41, 66, 94, 116	0
1	B	525/544 (96%)	0.18	16 (3%) 54 54	51, 73, 102, 125	0
All	All	1053/1088 (96%)	0.07	23 (2%) 65 66	41, 70, 99, 125	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	530	LEU	11.7
1	B	530	LEU	6.1
1	A	529	VAL	5.4
1	B	490	TRP	3.4
1	A	452	ARG	3.2
1	B	28	PHE	3.1
1	B	4	ILE	3.1
1	B	6	ILE	3.1
1	B	185	PHE	3.0
1	B	22	GLY	3.0
1	B	105	LYS	2.8
1	B	12	LYS	2.8
1	B	5	ILE	2.8
1	B	53	SER	2.5
1	B	50	THR	2.5
1	B	529	VAL	2.4
1	A	266	ASN	2.4
1	B	26	THR	2.4
1	B	99	ILE	2.3
1	B	364[A]	PHE	2.2
1	A	199	ALA	2.1
1	A	198	SER	2.0
1	A	364[A]	PHE	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 carbohydrates 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. 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
4	FUL	B	1006	10/11	0.89	0.55	9.98	96,106,112,116	0
4	FUL	A	1017	10/11	0.86	0.32	6.30	91,104,116,119	0
4	FUL	A	1010	10/11	0.91	0.24	2.00	91,111,114,114	0
3	NAG	B	1010	14/15	0.79	0.22	1.25	108,121,128,128	0
2	40V	A	1001	37/37	0.94	0.25	1.11	51,64,79,86	0
3	NAG	A	1014	14/15	0.80	0.22	1.00	78,92,103,104	0
7	EDO	A	1021	4/4	0.81	0.20	0.65	84,85,89,91	0
2	40V	B	1001	37/37	0.94	0.22	0.62	53,64,74,83	0
3	NAG	A	1011	14/15	0.80	0.21	0.31	97,104,115,116	0
3	NAG	B	1008	14/15	0.90	0.18	-0.89	89,97,103,108	0
6	CL	A	1019	1/1	0.90	0.11	-1.11	80,80,80,80	0
3	NAG	A	1013	14/15	0.94	0.13	-1.61	78,86,90,94	0
7	EDO	B	1012	4/4	0.95	0.10	-2.07	62,71,79,81	0
3	NAG	A	1015	14/15	0.83	0.18	-	109,124,136,145	0
7	EDO	B	1013	4/4	0.94	0.30	-	51,56,66,66	0
3	NAG	B	1007	14/15	0.68	0.28	-	112,128,131,136	0
3	NAG	B	1005	14/15	0.82	0.24	-	105,112,122,124	0
3	NAG	B	1004	14/15	0.87	0.31	-	107,113,118,121	0
4	FUL	A	1005	10/11	0.84	0.25	-	99,115,118,121	0
6	CL	A	1020	1/1	0.74	0.30	-	102,102,102,102	0
3	NAG	A	1016	14/15	0.79	0.35	-	138,144,150,152	0
7	EDO	A	1022	4/4	0.90	0.09	-	59,70,77,79	0
3	NAG	B	1002	14/15	0.66	0.20	-	118,129,138,142	0
6	CL	B	1011	1/1	0.90	0.13	-	76,76,76,76	0
5	FUC	A	1012	10/11	0.79	0.51	-	122,128,135,136	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NAG	A	1009	14/15	0.77	0.37	-	136,147,154,158	0
3	NAG	B	1009	14/15	0.84	0.25	-	109,117,122,123	0
3	NAG	A	1004	14/15	0.77	0.31	-	126,140,147,147	0
3	NAG	A	1003	14/15	0.79	0.17	-	101,112,135,135	0
3	NAG	B	1003	14/15	0.69	0.24	-	98,112,118,129	0
6	CL	A	1018	1/1	0.89	0.56	-	99,99,99,99	0
3	NAG	A	1002	14/15	0.93	0.20	-	90,103,112,113	0
3	NAG	A	1006	14/15	0.92	0.19	-	77,88,99,101	0
3	NAG	A	1008	14/15	0.86	0.17	-	118,122,136,147	0
3	NAG	A	1007	14/15	0.86	0.22	-	93,110,116,118	0

6.5 Other polymers [i](#)

There are no such residues in this entry.