



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:09 AM UTC

PDB ID : 7R2F / pdb_00007r2f
Title : Structure of tabun inhibited acetylcholinesterase in complex with 2-((hydroxymimino)methyl)-1-(5-(4-methyl-3-nitrobenzamido)pentyl)pyridinium
Authors : Forsgren, N.; Lindgren, C.; Edvinsson, L.; Linusson, A.; Ekstrom, F.
Deposited on : 2022-02-04
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

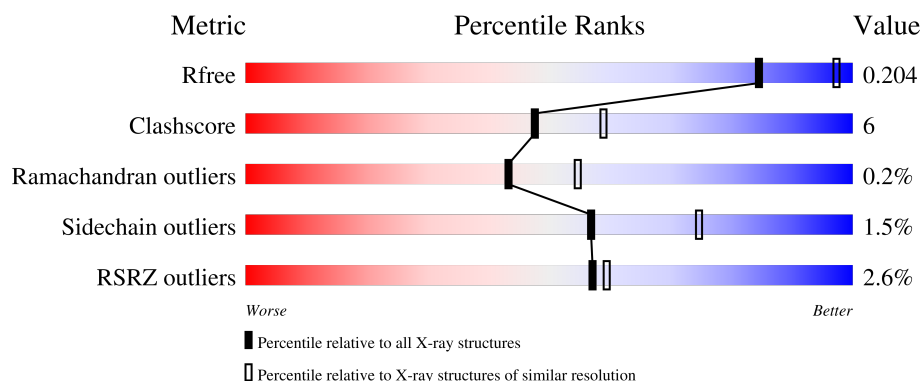
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	543	2% 84% 14% ..
1	B	543	3% 83% 14% ..

2 Entry composition [i](#)

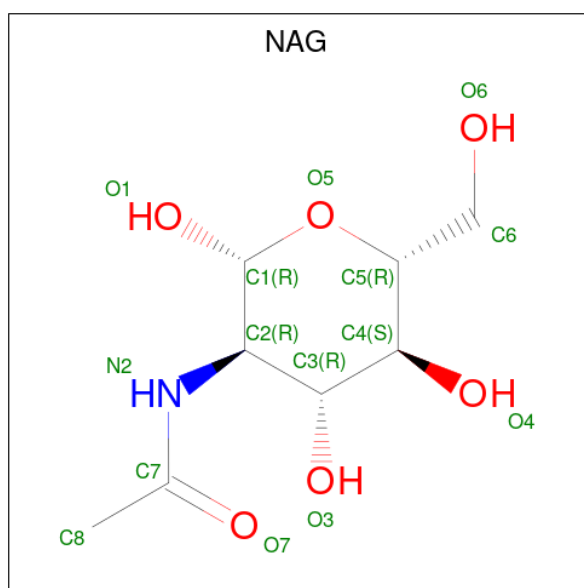
There are 6 unique types of molecules in this entry. The entry contains 9476 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

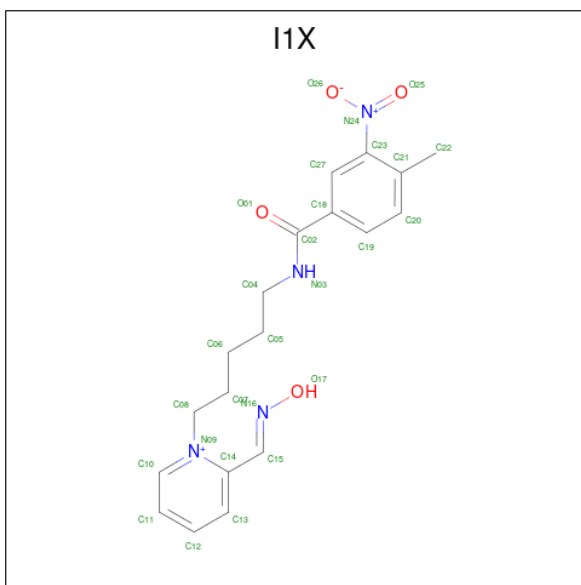
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	536	Total	C	N	O	P	S	0	42	0
			4528	2931	781	800	1	15			
1	B	533	Total	C	N	O	P	S	0	42	0
			4502	2914	771	799	1	17			

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



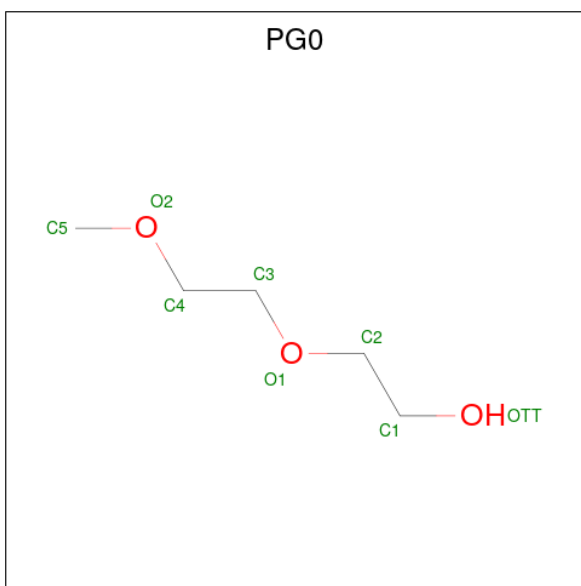
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is 4-methyl-3-nitro- {N}-[(2 {E},4 {E})-5-[2-[(oxidanylamino)methyl]pyridin-1-yl]penta-2,4-dienyl]benzamide (CCD ID: I1X) (formula: $C_{19}H_{23}N_4O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			16	11	2	3		
3	B	1	Total	C	N	O	0	0
			18	13	2	3		

- Molecule 4 is 2-(2-METHOXYETHOXY)ETHANOL (CCD ID: PG0) (formula: C₅H₁₂O₃).



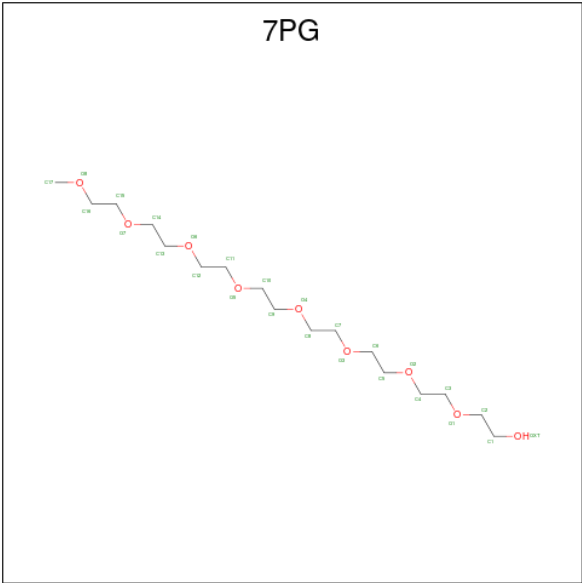
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	5	3		
4	A	1	Total	C	O	0	0
			7	5	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	5	3		
4	A	1	Total	C	O	0	0
			6	4	2		
4	A	1	Total	C	O	0	0
			8	5	3		
4	A	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		
4	B	1	Total	C	O	0	0
			8	5	3		

- Molecule 5 is 2,5,8,11,14,17,20,23-OCTAOXAPENTACOSAN-25-OL (CCD ID: 7PG) (formula: C₁₇H₃₆O₉).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			23	15	8		

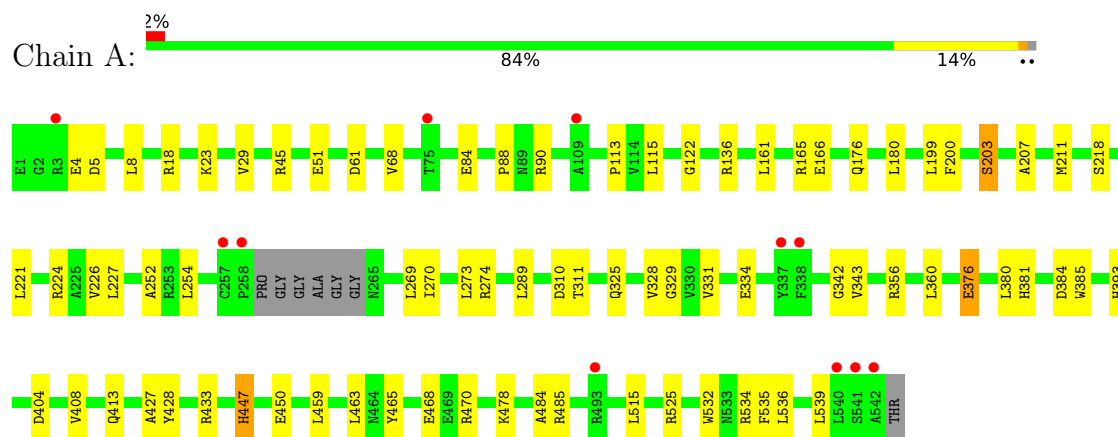
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	151	Total	O	0	0
			151	151		
6	B	101	Total	O	0	0
			101	101		

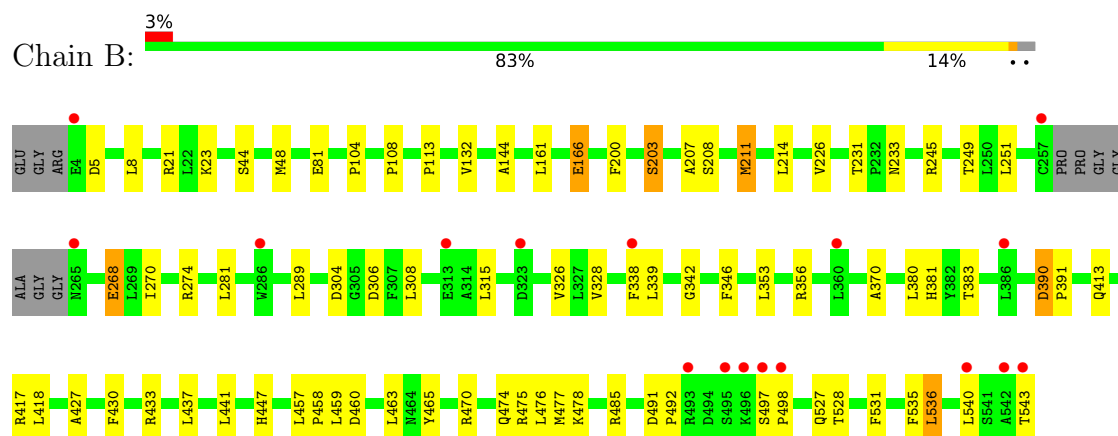
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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	78.83Å 110.75Å 227.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.14 – 2.30 46.14 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.4 (46.14-2.30) 95.6 (46.14-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.172 , 0.205 0.176 , 0.204	Depositor DCC
R_{free} test set	1692 reflections (1.69%)	wwPDB-VP
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.862	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9476	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: I1X, NAG, 7PG, SUN, PG0

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	0.47	0/4642	0.76	1/6343 (0.0%)
1	B	0.44	0/4614	0.78	2/6305 (0.0%)
All	All	0.46	0/9256	0.77	3/12648 (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	A	343	VAL	N-CA-C	5.53	112.59	107.56
1	B	390	ASP	CA-C-N	5.11	124.56	119.24
1	B	390	ASP	C-N-CA	5.11	124.56	119.24

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	4528	0	4481	52	0
1	B	4502	0	4458	53	0
2	A	28	0	26	0	0
3	A	16	0	0	0	0
3	B	18	0	0	0	0
4	A	45	0	64	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	64	0	96	4	0
5	B	23	0	29	4	0
6	A	151	0	0	0	0
6	B	101	0	0	2	0
All	All	9476	0	9154	102	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 6.

All (102) 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:113:PRO:HG2	4:B:605:PG0:H52	1.58	0.85
1:A:413:GLN:HG3	4:A:605:PG0:H31	1.66	0.77
1:B:338:PHE:HE1	1:B:447[B]:HIS:HE2	1.31	0.77
1:B:214:LEU:HB3	1:B:315[B]:LEU:HD23	1.71	0.71
1:B:268:GLU:N	1:B:268:GLU:OE1	2.27	0.68
1:B:132[B]:VAL:HA	1:B:457[B]:LEU:HD11	1.79	0.64
1:B:356[A]:ARG:NH1	1:B:383:THR:OG1	2.32	0.63
1:A:68[B]:VAL:HG21	1:A:88:PRO:HB3	1.82	0.61
1:A:404:ASP:OD1	1:A:525:ARG:NH1	2.34	0.61
1:B:380[B]:LEU:HB3	5:B:610:7PG:H91	1.82	0.61
1:B:207:ALA:O	1:B:211[A]:MET:HG2	2.02	0.59
1:A:45:ARG:NE	1:A:51:GLU:OE2	2.36	0.58
1:A:227[B]:LEU:HB2	1:A:328[B]:VAL:HG13	1.84	0.58
1:B:326[B]:VAL:HG11	1:B:418[B]:LEU:HD23	1.88	0.56
1:B:353[A]:LEU:HB3	1:B:391:PRO:HB2	1.88	0.56
1:B:211[A]:MET:HE2	1:B:308:LEU:HD21	1.88	0.56
1:B:381:HIS:CE1	5:B:610:7PG:H132	2.41	0.55
1:B:459[B]:LEU:HD11	1:B:477[B]:MET:HE1	1.90	0.54
1:A:463[B]:LEU:HD23	1:A:465:TYR:HE2	1.73	0.54
1:B:200:PHE:CB	1:B:226:VAL:HB	2.38	0.54
1:A:207:ALA:O	1:A:211:MET:HG2	2.08	0.53
1:B:245:ARG:O	1:B:249:THR:HG23	2.09	0.53
1:A:218:SER:HA	1:A:221[B]:LEU:HD23	1.90	0.53
1:A:115[A]:LEU:HD21	1:A:484:ALA:HB2	1.91	0.52
1:A:433[B]:ARG:H	4:A:607:PG0:H51	1.74	0.52
1:A:270[A]:ILE:HG22	1:A:274[A]:ARG:HH11	1.75	0.51
1:B:144:ALA:HA	4:B:605:PG0:H53	1.92	0.51
1:B:338:PHE:HE1	1:B:447[B]:HIS:NE2	2.04	0.51
1:B:23:LYS:NZ	6:B:704:HOH:O	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:GLU:OE2	1:A:18[B]:ARG:NE	2.44	0.51
1:A:534[B]:ARG:HB3	1:B:380[B]:LEU:HD21	1.92	0.50
1:A:180[B]:LEU:HD21	1:A:199[B]:LEU:HD21	1.93	0.50
1:B:161[B]:LEU:HD12	1:B:270:ILE:HD11	1.93	0.50
1:B:459[B]:LEU:HD23	1:B:470:ARG:HG2	1.92	0.49
1:A:113:PRO:HG2	1:A:485:ARG:HG2	1.94	0.49
1:A:165:ARG:HH21	1:A:166:GLU:HG3	1.78	0.49
1:B:413:GLN:HG3	4:B:603:PG0:H12	1.95	0.49
1:B:430:PHE:HE2	1:B:476[B]:LEU:HD21	1.77	0.48
1:A:122:GLY:HA2	1:A:203:SUN:H1C3	1.95	0.48
1:A:161:LEU:HD11	1:A:269:LEU:HD22	1.96	0.48
1:B:328:VAL:O	1:B:427:ALA:HA	2.14	0.48
1:B:200:PHE:HB2	1:B:226:VAL:HB	1.96	0.47
1:B:370:ALA:HA	1:B:540:LEU:HD21	1.96	0.47
1:B:458:PRO:HA	1:B:465:TYR:CD2	2.50	0.47
1:B:433:ARG:CZ	1:B:437[A]:LEU:HD23	2.44	0.47
1:A:328[B]:VAL:O	1:A:427:ALA:HA	2.15	0.47
1:A:18[A]:ARG:NH2	1:A:61:ASP:OD2	2.39	0.46
1:A:532:TRP:CE3	1:A:536[B]:LEU:HD12	2.51	0.46
1:B:485:ARG:HG2	4:B:605:PG0:H12	1.98	0.46
1:A:376:GLU:O	1:A:380[A]:LEU:HG	2.16	0.46
1:B:251[B]:LEU:HD11	1:B:281[B]:LEU:HG	1.98	0.45
1:A:224:ARG:HG2	1:A:325:GLN:HB2	1.96	0.45
1:A:450:GLU:OE1	1:A:450:GLU:N	2.42	0.45
1:B:304:ASP:OD2	1:B:306:ASP:HB3	2.16	0.45
1:B:339[B]:LEU:HD23	1:B:346:PHE:CE2	2.52	0.45
1:A:270[A]:ILE:HD13	1:A:273[A]:LEU:HD12	1.99	0.45
1:B:540:LEU:O	1:B:543:THR:HG22	2.16	0.45
1:A:200:PHE:CB	1:A:226:VAL:HB	2.47	0.45
1:B:417:ARG:HD3	1:B:417:ARG:HA	1.72	0.44
1:B:531:PHE:CZ	1:B:536:LEU:HD11	2.52	0.44
1:A:468:GLU:OE1	1:A:468:GLU:N	2.50	0.44
1:B:48:MET:SD	1:B:166:GLU:HA	2.57	0.44
1:B:497:SER:HB2	1:B:498:PRO:HA	2.00	0.44
1:A:534[B]:ARG:NH2	4:A:605:PG0:H41	2.33	0.44
1:B:459[B]:LEU:HD21	1:B:474:GLN:HG3	2.00	0.44
1:A:535:PHE:CE2	1:A:539:LEU:HD22	2.53	0.44
1:A:165:ARG:NH2	1:A:166:GLU:HG3	2.33	0.43
1:B:491:ASP:HA	1:B:492:PRO:HD3	1.83	0.43
1:A:29:VAL:HG21	1:A:136:ARG:HB2	1.99	0.43
1:A:331:VAL:HG22	1:A:334:GLU:OE2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:VAL:HG11	1:A:525:ARG:HG3	2.00	0.43
1:A:536[B]:LEU:HA	1:A:536[B]:LEU:HD23	1.66	0.42
1:B:5:ASP:HB3	1:B:8[B]:LEU:HD22	2.00	0.42
1:B:208:SER:HA	1:B:211[A]:MET:HG3	2.02	0.42
1:B:231:THR:HB	1:B:233:ASN:OD1	2.19	0.42
1:A:459[B]:LEU:HD23	1:A:470:ARG:HG3	2.02	0.42
1:A:381:HIS:ND1	5:B:610:7PG:H12	2.35	0.42
1:A:356:ARG:O	1:A:360[A]:LEU:HG	2.19	0.42
1:A:200:PHE:HB2	1:A:226:VAL:HB	2.02	0.42
1:A:310:ASP:OD1	1:A:311:THR:N	2.44	0.42
1:A:515[A]:LEU:HG	4:A:607:PG0:H31	2.02	0.42
1:B:289[B]:LEU:HD12	1:B:289[B]:LEU:HA	1.81	0.41
1:A:176:GLN:O	1:A:180[A]:LEU:HG	2.19	0.41
1:B:21:ARG:NH1	6:B:702:HOH:O	2.53	0.41
1:B:528:THR:O	1:B:531:PHE:HB3	2.20	0.41
1:A:5:ASP:HB3	1:A:8:LEU:HD12	2.02	0.41
1:A:176:GLN:O	1:A:180[B]:LEU:HD13	2.20	0.41
1:A:329:GLY:HA3	1:A:428:TYR:CZ	2.55	0.41
1:A:252:ALA:HA	1:A:273[A]:LEU:HD21	2.01	0.41
1:A:289[B]:LEU:HD12	1:A:289[B]:LEU:HA	1.89	0.41
1:B:203:SUN:H2C2	1:B:338:PHE:CZ	2.56	0.41
1:B:104:PRO:HG2	1:B:108:PRO:HG3	2.02	0.41
1:B:460:ASP:HB3	1:B:463[B]:LEU:HD22	2.03	0.41
1:A:254[B]:LEU:HD12	1:A:254[B]:LEU:HA	1.90	0.41
1:A:384:ASP:HB2	1:A:393:HIS:CE1	2.56	0.41
1:B:527:GLN:HB3	5:B:610:7PG:H22	2.02	0.41
1:A:203:SUN:OG	1:A:447[B]:HIS:NE2	2.54	0.40
1:A:68[A]:VAL:HG23	1:A:90:ARG:HB2	2.02	0.40
1:A:380[A]:LEU:HD12	1:B:535:PHE:HB2	2.02	0.40
1:B:44:SER:HA	1:B:274:ARG:HD2	2.03	0.40
1:A:385:TRP:HD1	1:B:527:GLN:OE1	2.04	0.40
1:B:390:ASP:HA	1:B:391:PRO:HD2	1.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	573/543 (106%)	553 (96%)	19 (3%)	1 (0%)	43	55
1	B	570/543 (105%)	554 (97%)	15 (3%)	1 (0%)	43	55
All	All	1143/1086 (105%)	1107 (97%)	34 (3%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	342	GLY
1	A	342	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	478/442 (108%)	472 (99%)	6 (1%)	61	77
1	B	477/442 (108%)	468 (98%)	9 (2%)	50	69
All	All	955/884 (108%)	940 (98%)	15 (2%)	57	73

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LYS
1	A	84	GLU
1	A	376	GLU
1	A	447[A]	HIS
1	A	447[B]	HIS
1	A	478	LYS
1	B	81	GLU
1	B	166	GLU
1	B	211[A]	MET
1	B	211[B]	MET

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Mol	Chain	Res	Type
1	B	268	GLU
1	B	441	LEU
1	B	475	ARG
1	B	478	LYS
1	B	536	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	393	HIS
1	A	499	GLN
1	B	317	ASN
1	B	421	GLN
1	B	464	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	SUN	B	203	1	12,13,14	1.47	3 (25%)	10,17,19	2.61	3 (30%)
1	SUN	A	203	1	12,13,14	1.56	3 (25%)	10,17,19	2.27	3 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SUN	B	203	1	-	4/16/18/20	-
1	SUN	A	203	1	-	3/16/18/20	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	203	SUN	P1-O1	2.87	1.50	1.46
1	B	203	SUN	P1-OG	2.77	1.68	1.57
1	A	203	SUN	OG-CB	-2.77	1.34	1.44
1	B	203	SUN	P1-O1	2.07	1.49	1.46
1	A	203	SUN	P1-O2	2.06	1.65	1.57
1	B	203	SUN	P1-O2	2.04	1.65	1.57

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	203	SUN	OG-CB-CA	6.60	114.57	108.14
1	A	203	SUN	OG-CB-CA	5.78	113.77	108.14
1	B	203	SUN	OG-P1-O1	-2.89	106.50	115.71
1	B	203	SUN	O2-P1-OG	2.55	108.64	100.62
1	A	203	SUN	O2-P1-OG	2.33	107.94	100.62
1	A	203	SUN	OG-P1-O1	-2.22	108.63	115.71

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	203	SUN	N-CA-CB-OG
1	A	203	SUN	C1-N1-P1-OG
1	B	203	SUN	C1-N1-P1-O2
1	B	203	SUN	N-CA-CB-OG
1	B	203	SUN	C3-O2-P1-N1
1	A	203	SUN	C-CA-CB-OG
1	B	203	SUN	C-CA-CB-OG

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	203	SUN	1	0
1	A	203	SUN	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

19 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
4	PG0	B	605	-	7,7,7	0.63	0	6,6,6	1.03	1 (16%)
4	PG0	B	606	-	7,7,7	0.59	0	6,6,6	0.87	0
4	PG0	B	609	-	7,7,7	0.57	0	6,6,6	0.89	0
4	PG0	B	607	-	7,7,7	0.59	0	6,6,6	0.73	0
2	NAG	A	601	1	14,14,15	0.52	0	17,19,21	0.64	0
2	NAG	A	602	1	14,14,15	0.40	0	17,19,21	0.55	0
4	PG0	A	606	-	7,7,7	0.61	0	6,6,6	0.80	0
3	I1X	A	603	-	16,16,28	2.17	3 (18%)	17,21,36	1.12	1 (5%)
4	PG0	A	608	-	7,7,7	0.56	0	6,6,6	0.85	0
4	PG0	A	609	-	7,7,7	0.61	0	6,6,6	0.78	0
4	PG0	A	607	-	5,5,7	0.52	0	4,4,6	0.91	0
4	PG0	A	605	-	6,6,7	0.53	0	5,5,6	1.11	1 (20%)
5	7PG	B	610	-	22,22,25	0.65	0	21,21,24	0.94	0
4	PG0	B	602	-	7,7,7	0.58	0	6,6,6	0.79	0
4	PG0	B	604	-	7,7,7	0.62	0	6,6,6	0.87	0
3	I1X	B	601	-	18,18,28	2.07	2 (11%)	19,23,36	1.05	0
4	PG0	B	603	-	7,7,7	0.57	0	6,6,6	0.81	0
4	PG0	B	608	-	7,7,7	0.61	0	6,6,6	0.73	0
4	PG0	A	604	-	7,7,7	0.60	0	6,6,6	0.75	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
4	PG0	B	605	-	-	3/5/5/5	-
4	PG0	B	606	-	-	3/5/5/5	-
4	PG0	B	609	-	-	1/5/5/5	-
4	PG0	B	607	-	-	3/5/5/5	-
2	NAG	A	601	1	-	2/6/23/26	0/1/1/1
2	NAG	A	602	1	-	2/6/23/26	0/1/1/1
4	PG0	A	606	-	-	5/5/5/5	-
3	I1X	A	603	-	-	0/10/12/20	0/1/1/2
4	PG0	A	608	-	-	3/5/5/5	-
4	PG0	A	609	-	-	2/5/5/5	-
4	PG0	A	607	-	-	2/3/3/5	-
4	PG0	A	605	-	-	2/4/4/5	-
5	7PG	B	610	-	-	14/20/20/23	-
4	PG0	B	602	-	-	4/5/5/5	-
4	PG0	B	604	-	-	4/5/5/5	-
3	I1X	B	601	-	-	5/12/14/20	0/1/1/2
4	PG0	B	603	-	-	2/5/5/5	-
4	PG0	B	608	-	-	4/5/5/5	-
4	PG0	A	604	-	-	3/5/5/5	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	601	I1X	C02-N03	6.87	1.48	1.33
3	A	603	I1X	C02-N03	6.63	1.48	1.33
3	A	603	I1X	O25-N24	4.61	1.30	1.22
3	B	601	I1X	O25-N24	4.58	1.30	1.22
3	A	603	I1X	O26-N24	-2.01	1.21	1.35

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	603	I1X	O25-N24-C23	2.38	123.10	119.03
4	A	605	PG0	C2-O1-C3	2.07	120.34	113.06
4	B	605	PG0	C2-O1-C3	2.03	122.14	113.26

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	I1X	C21-C23-N24-O25
3	B	601	I1X	C27-C23-N24-O25
2	A	601	NAG	O5-C5-C6-O6
2	A	601	NAG	C4-C5-C6-O6
4	A	607	PG0	O1-C3-C4-O2
4	A	604	PG0	O1-C3-C4-O2
4	A	606	PG0	O1-C3-C4-O2
3	B	601	I1X	N03-C04-C05-C06
4	A	608	PG0	O1-C3-C4-O2
2	A	602	NAG	O5-C5-C6-O6
5	B	610	7PG	O3-C7-C8-O4
4	B	607	PG0	O1-C3-C4-O2
5	B	610	7PG	O5-C10-C9-O4
4	B	607	PG0	OTT-C1-C2-O1
4	B	608	PG0	OTT-C1-C2-O1
4	B	606	PG0	OTT-C1-C2-O1
4	A	609	PG0	O1-C3-C4-O2
4	B	605	PG0	OTT-C1-C2-O1
4	B	604	PG0	O1-C3-C4-O2
4	A	606	PG0	OTT-C1-C2-O1
4	A	608	PG0	OTT-C1-C2-O1
4	B	602	PG0	OTT-C1-C2-O1
4	B	604	PG0	OTT-C1-C2-O1
4	B	606	PG0	O1-C3-C4-O2
5	B	610	7PG	OXT-C1-C2-O1
4	B	605	PG0	O1-C3-C4-O2
5	B	610	7PG	O5-C11-C12-O6
5	B	610	7PG	O1-C3-C4-O2
3	B	601	I1X	C04-C05-C06-C07
5	B	610	7PG	C12-C11-O5-C10
4	A	606	PG0	C1-C2-O1-C3
4	B	604	PG0	C4-C3-O1-C2
5	B	610	7PG	C4-C3-O1-C2
4	A	604	PG0	C4-C3-O1-C2
5	B	610	7PG	C10-C9-O4-C8
4	B	602	PG0	C1-C2-O1-C3
4	B	605	PG0	C3-C4-O2-C5
4	B	602	PG0	C4-C3-O1-C2
4	B	608	PG0	C3-C4-O2-C5
5	B	610	7PG	C9-C10-O5-C11
4	B	608	PG0	O1-C3-C4-O2
4	A	606	PG0	C3-C4-O2-C5
4	B	603	PG0	C3-C4-O2-C5

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Mol	Chain	Res	Type	Atoms
4	B	604	PG0	C1-C2-O1-C3
4	A	607	PG0	C4-C3-O1-C2
4	A	606	PG0	C4-C3-O1-C2
4	A	608	PG0	C3-C4-O2-C5
5	B	610	7PG	C3-C4-O2-C5
5	B	610	7PG	O2-C5-C6-O3
3	B	601	I1X	C05-C06-C07-C08
4	A	609	PG0	C1-C2-O1-C3
4	B	608	PG0	C4-C3-O1-C2
4	B	607	PG0	C1-C2-O1-C3
4	B	602	PG0	O1-C3-C4-O2
2	A	602	NAG	C4-C5-C6-O6
4	A	605	PG0	C4-C3-O1-C2
5	B	610	7PG	C1-C2-O1-C3
4	A	605	PG0	O1-C3-C4-O2
4	B	609	PG0	O1-C3-C4-O2
4	B	603	PG0	O1-C3-C4-O2
4	B	606	PG0	C1-C2-O1-C3
5	B	610	7PG	O6-C13-C14-O7
5	B	610	7PG	C14-C13-O6-C12
4	A	604	PG0	OTT-C1-C2-O1

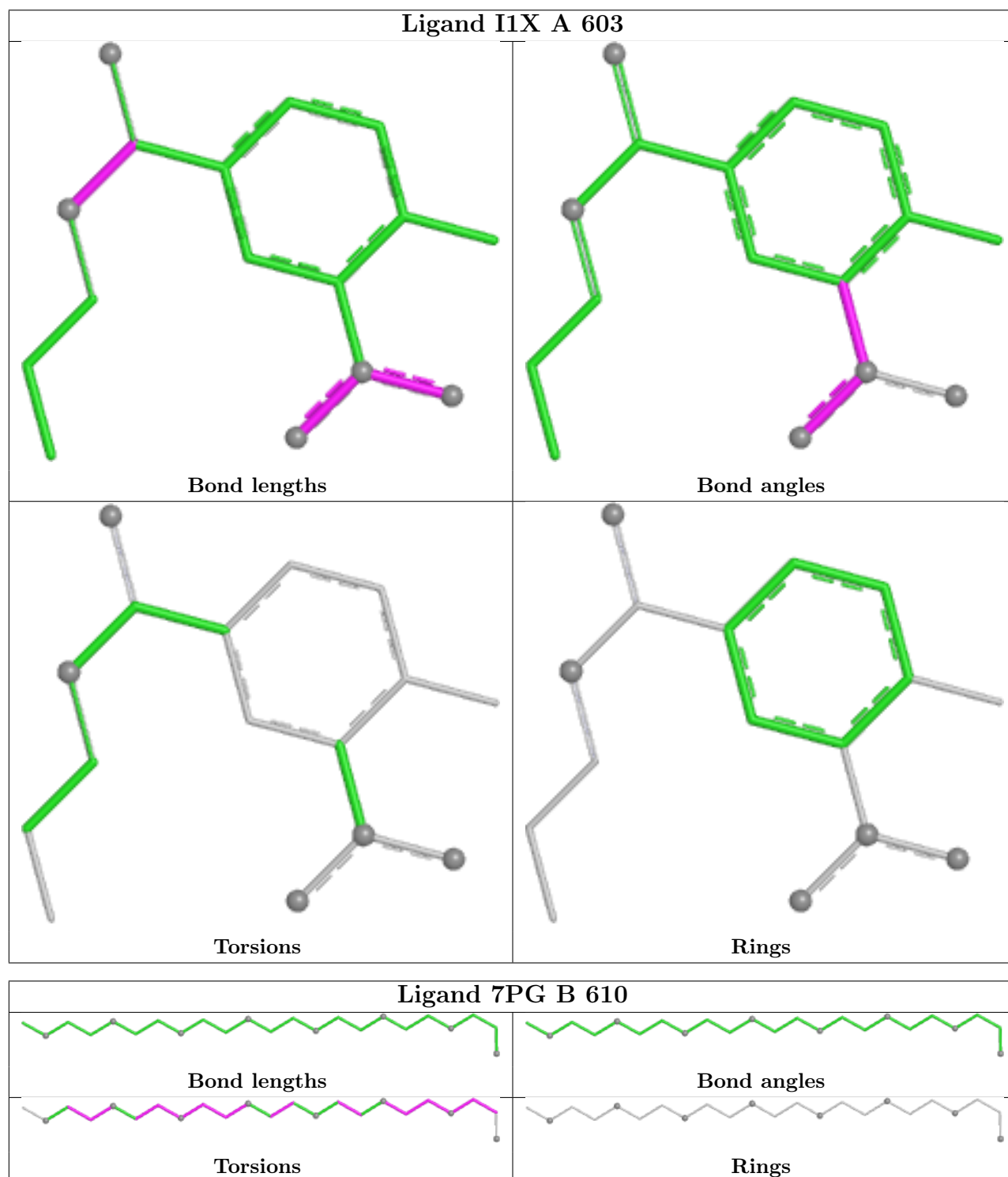
There are no ring outliers.

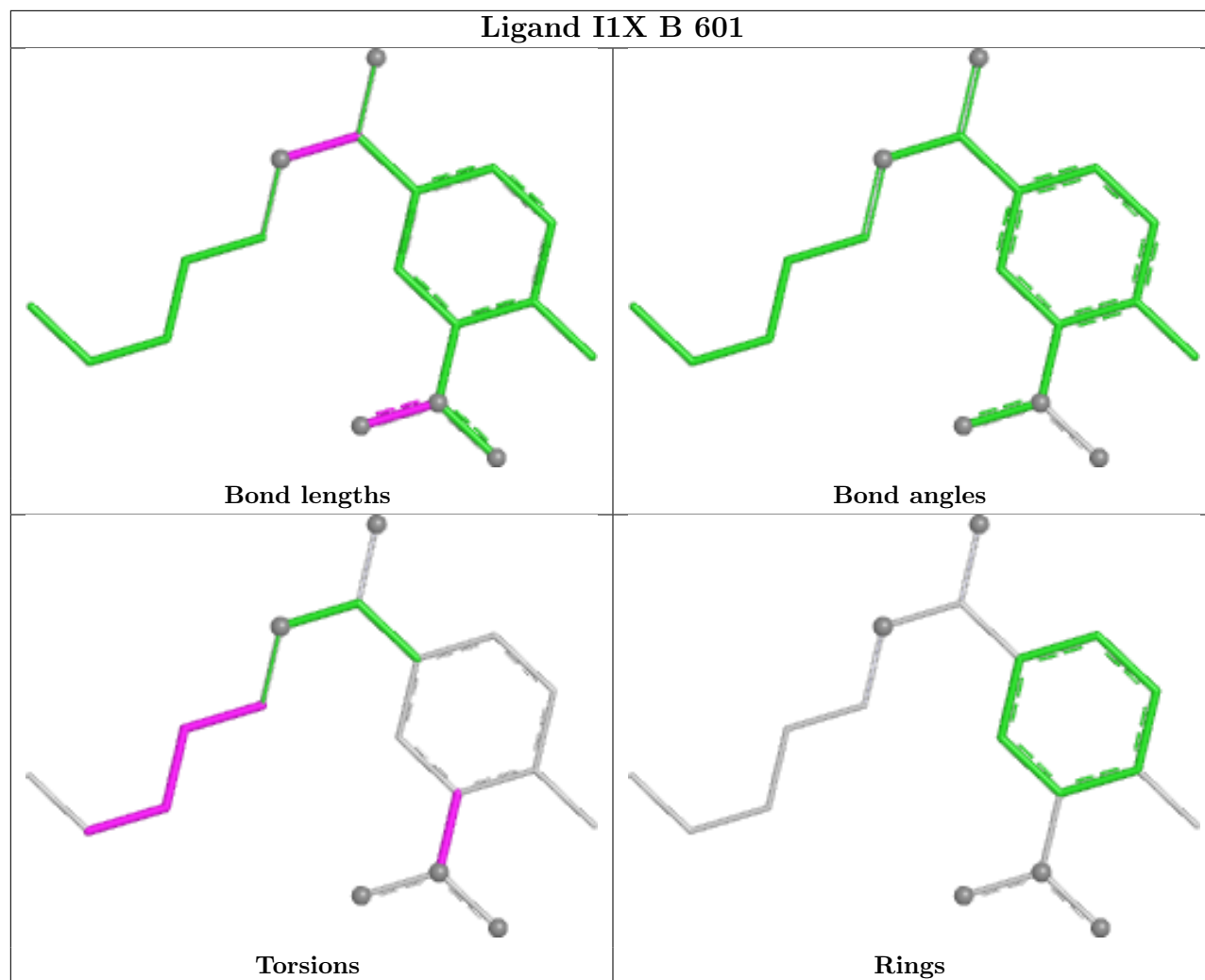
5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	605	PG0	3	0
4	A	607	PG0	2	0
4	A	605	PG0	2	0
5	B	610	7PG	4	0
4	B	603	PG0	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	535/543 (98%)	-0.09	11 (2%) 63 65	18, 49, 84, 135	42 (7%)
1	B	532/543 (97%)	0.04	17 (3%) 50 52	21, 55, 88, 143	42 (7%)
All	All	1067/1086 (98%)	-0.02	28 (2%) 57 59	18, 52, 85, 143	84 (7%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	497	SER	7.0
1	A	258	PRO	6.8
1	B	543	THR	5.3
1	B	257	CYS	4.8
1	B	498	PRO	4.7
1	B	386	LEU	4.3
1	A	493	ARG	3.6
1	B	323	ASP	3.6
1	A	542	ALA	3.6
1	A	541	SER	3.5
1	B	542	ALA	3.4
1	B	496	LYS	3.3
1	A	540	LEU	3.2
1	B	286	TRP	3.1
1	A	338[A]	PHE	3.0
1	A	109	ALA	2.8
1	B	338	PHE	2.7
1	B	495	SER	2.6
1	A	257	CYS	2.5
1	A	337	TYR	2.4
1	A	75	THR	2.3
1	B	360[A]	LEU	2.3
1	B	313	GLU	2.3
1	A	3	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	493	ARG	2.2
1	B	265	ASN	2.2
1	B	4	GLU	2.2
1	B	540	LEU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SUN	B	203	14/15	0.96	0.10	37,46,74,84	0
1	SUN	A	203	14/15	0.98	0.08	34,44,68,76	0

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
2	NAG	A	602	14/15	0.62	0.18	100,111,120,120	0
4	PG0	B	607	8/8	0.69	0.26	83,88,95,96	0
4	PG0	B	606	8/8	0.76	0.23	86,88,91,93	0
2	NAG	A	601	14/15	0.76	0.20	83,96,111,112	0
4	PG0	B	609	8/8	0.77	0.30	94,100,104,104	0
4	PG0	A	607	6/8	0.78	0.33	80,89,92,94	0
4	PG0	B	608	8/8	0.78	0.24	82,89,101,102	0
3	I1X	B	601	18/27	0.78	0.21	68,103,114,122	0
4	PG0	A	608	8/8	0.79	0.22	77,86,90,91	0
4	PG0	B	603	8/8	0.79	0.26	60,78,99,101	0
4	PG0	B	605	8/8	0.80	0.26	86,94,97,100	0
4	PG0	B	602	8/8	0.83	0.21	74,80,86,92	0

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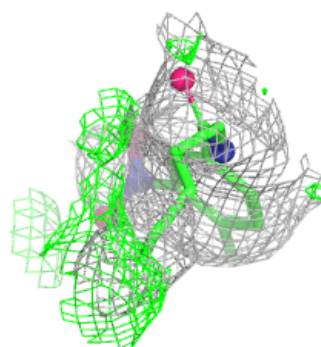
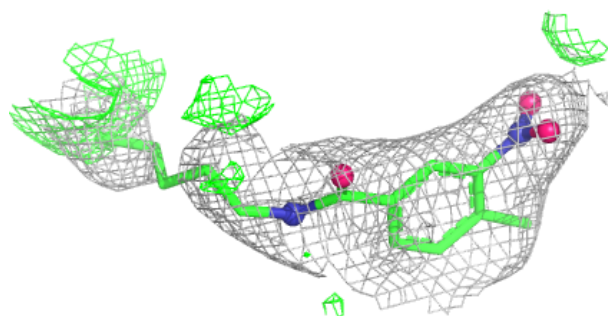
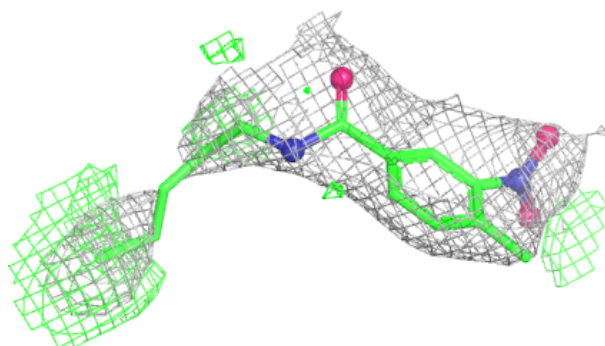
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	PG0	A	609	8/8	0.84	0.26	85,91,109,111	0
4	PG0	A	606	8/8	0.84	0.27	85,88,95,97	0
4	PG0	A	605	7/8	0.85	0.28	61,69,86,86	0
3	I1X	A	603	16/27	0.86	0.20	69,76,85,95	0
4	PG0	B	604	8/8	0.87	0.21	79,84,91,93	0
4	PG0	A	604	8/8	0.87	0.19	69,76,88,90	0
5	7PG	B	610	23/26	0.89	0.20	46,75,90,91	0

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

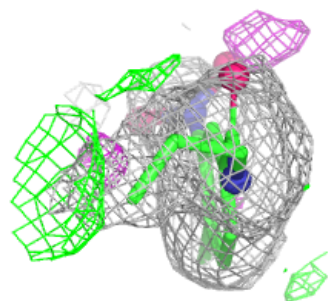
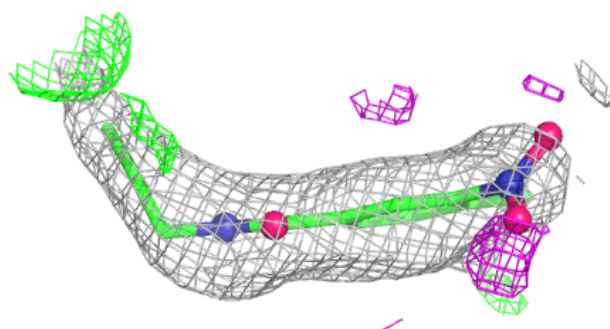
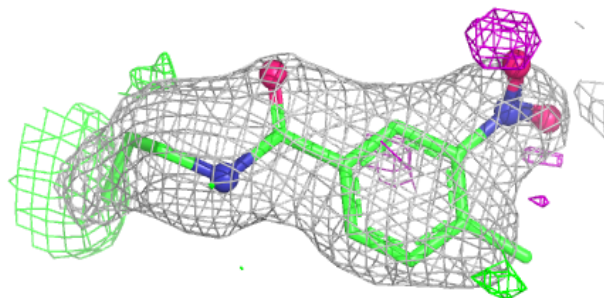
Electron density around I1X B 601:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



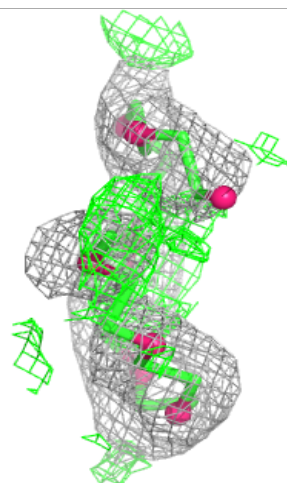
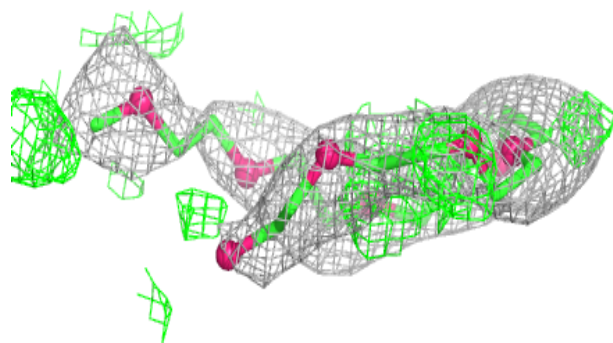
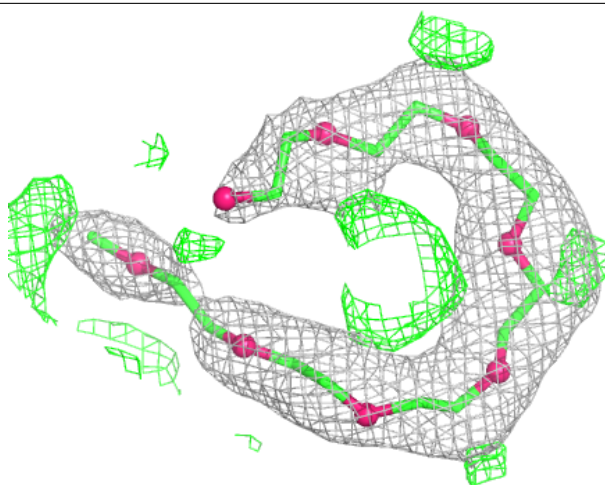
Electron density around I1X A 603:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 7PG B 610:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.