



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 14, 2024 – 04:52 PM EST

PDB ID : 2GYV
Title : Crystal structure of Mus musculus Acetylcholinesterase in complex with Ortho-7
Authors : Pang, Y.P.; Boman, M.; Artursson, E.; Akfur, C.; Lundberg, S.
Deposited on : 2006-05-10
Resolution : 2.50 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

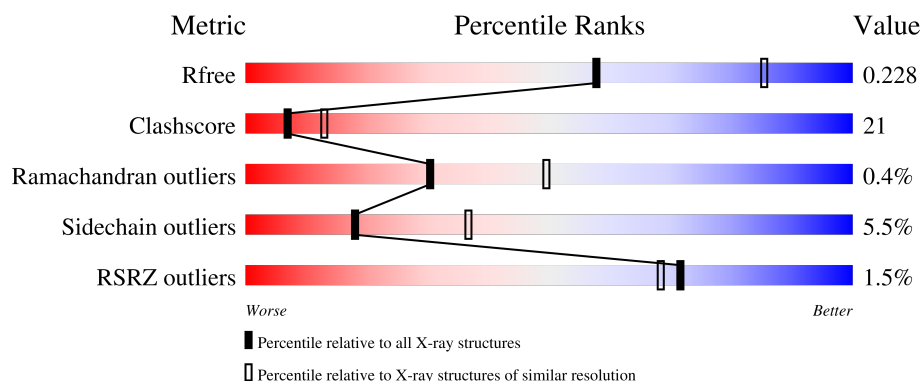
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	164625	5504 (2.50-2.50)
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (2.50-2.50)
RSRZ outliers	164620	5504 (2.50-2.50)

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	<div> <div>%</div> <div> <div></div> <div>56%</div> <div>41%</div> <div>..</div> </div> </div>
1	B	543	<div> <div>%</div> <div> <div></div> <div>62%</div> <div>34%</div> <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
3	CO3	A	1408	-	-	X	-
3	CO3	B	544	-	-	X	-
4	HBP	A	1409	-	-	X	-
4	HBP	B	545	-	-	X	-

2 Entry composition [i](#)

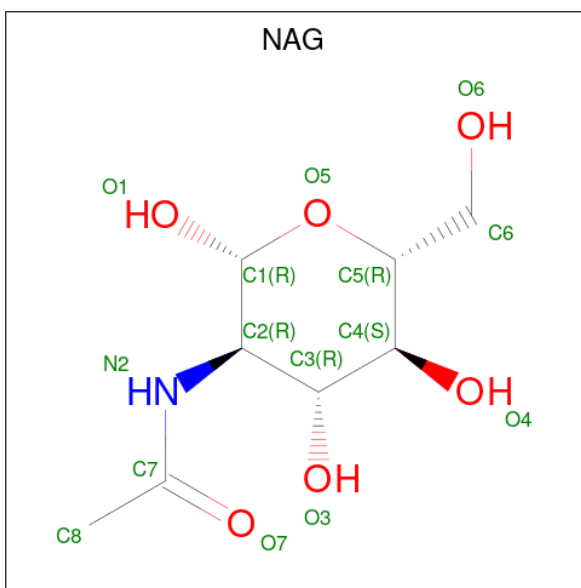
There are 6 unique types of molecules in this entry. The entry contains 8771 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

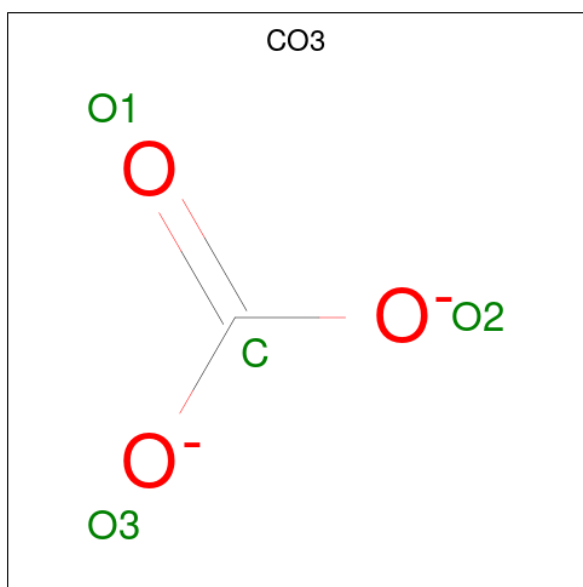
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	0	0	0
			4177	2679	725	759	14			
1	B	533	Total	C	N	O	S	0	0	0
			4159	2670	718	757	14			

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



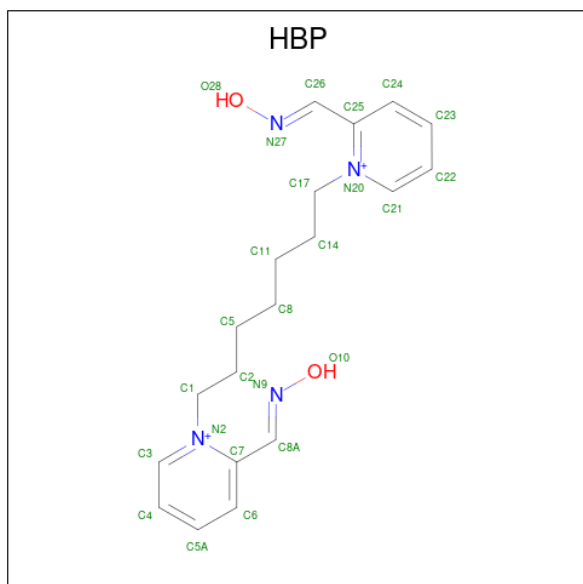
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			13	8	1	4		

- Molecule 3 is CARBONATE ION (three-letter code: CO3) (formula: CO_3).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	1	3		
3	B	1	Total	C	O	0	0
			4	1	3		

- Molecule 4 is 1,7-HEPTYLENE-BIS-N,N'-SYN-2-PYRIDINIUMALDOXIME (three-letter code: HBP) (formula: C₁₉H₂₆N₄O₂).



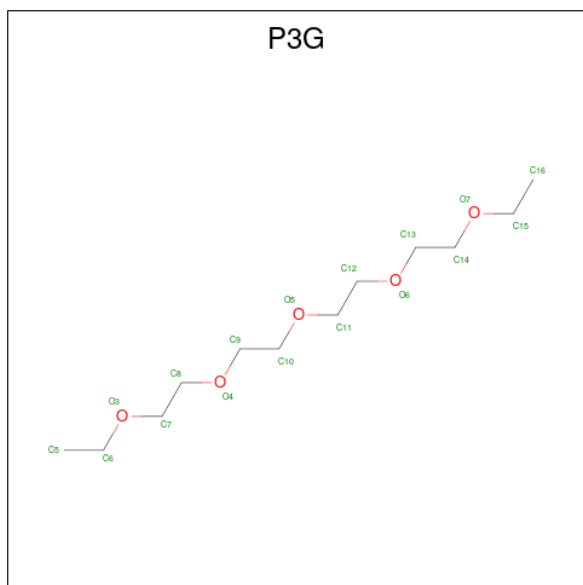
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			25	19	4	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			25	19	4	2		

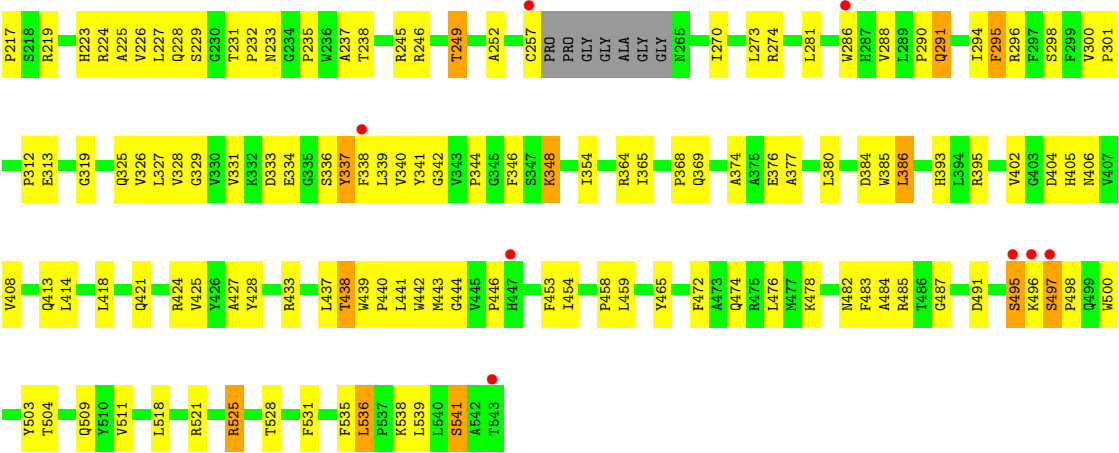
- Molecule 5 is 3,6,9,12,15-PENTAOXAHEPTADECANE (three-letter code: P3G) (formula: $C_{12}H_{26}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	5	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	194	Total	O	0	0
			194	194		
6	B	149	Total	O	0	0
			149	149		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.73Å 108.58Å 220.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.00 – 2.50 29.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (29.00-2.50) 99.2 (29.00-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.50 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.197 , 0.241 0.188 , 0.228	Depositor DCC
R_{free} test set	1276 reflections (1.99%)	wwPDB-VP
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8771	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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: P3G, CO3, HBP, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.75	0/4300	0.77	1/5875 (0.0%)
1	B	0.68	0/4282	0.71	0/5851
All	All	0.71	0/8582	0.74	1/11726 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	161	LEU	CA-CB-CG	-5.63	102.36	115.30

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	4177	0	4063	184	0
1	B	4159	0	4047	156	0
2	A	27	0	26	0	0
3	A	4	0	0	5	0
3	B	4	0	0	3	0
4	A	25	0	26	15	0
4	B	25	0	26	9	0
5	B	7	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	194	0	0	13	0
6	B	149	0	0	11	0
All	All	8771	0	8194	345	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 21.

All (345) 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:203:SER:OG	3:A:1408:CO3:C	1.74	1.31
4:A:1409:HBP:H26	4:A:1409:HBP:H16	1.18	1.10
1:B:203:SER:OG	3:B:544:CO3:C	2.04	1.06
1:B:245:ARG:O	1:B:249:THR:HG23	1.57	1.03
1:B:197:VAL:H	1:B:223:HIS:HD2	1.07	1.00
1:A:203:SER:HG	3:A:1408:CO3:C	1.82	0.93
4:A:1409:HBP:H16	4:A:1409:HBP:C26	2.00	0.92
1:B:197:VAL:H	1:B:223:HIS:CD2	1.89	0.90
4:A:1409:HBP:H26	4:A:1409:HBP:C14	2.03	0.88
1:A:48:MET:HE1	1:A:166:GLU:HA	1.52	0.88
4:B:545:HBP:H1B	4:B:545:HBP:N9	1.90	0.84
1:A:245:ARG:O	1:A:249:THR:HG23	1.78	0.83
1:A:197:VAL:H	1:A:223:HIS:CD2	1.96	0.83
1:A:353:LEU:HB3	1:A:391:PRO:HB2	1.62	0.80
1:B:136:ARG:HD2	6:B:686:HOH:O	1.81	0.79
4:B:545:HBP:H16	4:B:545:HBP:H26	1.64	0.79
1:B:296:ARG:NH2	1:B:406:ASN:OD1	2.17	0.78
1:A:266:ASP:O	1:A:270:ILE:HD12	1.84	0.78
1:A:197:VAL:H	1:A:223:HIS:HD2	1.30	0.75
1:B:369:GLN:HB2	6:B:562:HOH:O	1.86	0.75
1:A:203:SER:OG	3:A:1408:CO3:O3	2.04	0.74
1:A:535:PHE:HB2	1:B:380:LEU:HD12	1.70	0.73
1:A:203:SER:OG	3:A:1408:CO3:O2	2.06	0.73
1:A:224:ARG:HG2	1:A:325:GLN:HB2	1.69	0.73
1:B:112:THR:HG21	1:B:143:GLY:O	1.90	0.71
1:B:207:ALA:O	1:B:211:MET:HG2	1.90	0.71
1:B:525:ARG:HH11	1:B:525:ARG:CG	2.02	0.71
1:A:227:LEU:HD22	1:A:414:LEU:HD21	1.71	0.71
1:B:497:SER:H	1:B:498:PRO:HA	1.56	0.70
4:A:1409:HBP:H1B	4:A:1409:HBP:N9	2.07	0.70
1:B:328:VAL:O	1:B:427:ALA:HA	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:TRP:HB2	4:A:1409:HBP:C6	2.22	0.69
1:A:48:MET:HE1	1:A:166:GLU:CA	2.23	0.69
1:A:48:MET:HE1	1:A:165:ARG:O	1.92	0.68
1:B:525:ARG:HH11	1:B:525:ARG:HG2	1.58	0.67
1:A:424:ARG:HD2	6:A:1440:HOH:O	1.96	0.66
1:B:286:TRP:HB2	4:B:545:HBP:C6	2.26	0.65
1:A:103:THR:HG21	1:A:190:PHE:HB3	1.77	0.65
1:A:458:PRO:HA	1:A:465:TYR:CD2	2.32	0.65
4:B:545:HBP:H26	4:B:545:HBP:C14	2.24	0.65
1:B:458:PRO:HA	1:B:465:TYR:CD2	2.31	0.65
1:B:77:TYR:CZ	1:B:348:LYS:HG2	2.32	0.64
1:B:104:PRO:HG3	1:B:143:GLY:HA2	1.79	0.64
1:A:211:MET:HG2	1:A:308:LEU:HD21	1.81	0.63
1:A:176:GLN:OE1	1:A:208:SER:HB3	1.98	0.63
1:B:224:ARG:HD3	1:B:487:GLY:CA	2.29	0.62
4:A:1409:HBP:C26	4:A:1409:HBP:C14	2.68	0.61
1:A:112:THR:HG21	1:A:143:GLY:O	2.00	0.61
1:B:224:ARG:HD3	1:B:487:GLY:HA2	1.82	0.61
1:A:337:TYR:HA	1:A:443:MET:CE	2.30	0.61
1:B:88:PRO:HG2	1:B:92:LEU:HD21	1.83	0.61
1:B:224:ARG:HG2	1:B:325:GLN:HB2	1.81	0.61
1:A:277:PRO:HG2	1:A:280:ASP:HB2	1.83	0.61
1:A:522:ARG:HB3	1:B:386:LEU:HD21	1.81	0.61
1:B:80:PHE:CE1	1:B:438:THR:HB	2.35	0.61
1:B:231:THR:HB	1:B:233:ASN:OD1	2.01	0.61
1:A:141:VAL:HG21	1:A:459:LEU:HD23	1.83	0.61
1:A:511:VAL:HB	1:A:518:LEU:HD22	1.82	0.61
1:A:141:VAL:HG21	1:A:459:LEU:CD2	2.30	0.60
1:A:538:LYS:HE3	1:B:376:GLU:OE2	2.00	0.60
1:B:200:PHE:CB	1:B:226:VAL:HB	2.31	0.60
1:B:496:LYS:HB2	6:B:581:HOH:O	2.01	0.60
1:B:141:VAL:HG21	1:B:459:LEU:CD2	2.32	0.59
1:B:176:GLN:OE1	1:B:208:SER:HB3	2.03	0.59
1:B:203:SER:HG	3:B:544:CO3:C	2.13	0.59
1:B:374:ALA:HA	1:B:539:LEU:HD23	1.85	0.59
1:B:197:VAL:N	1:B:223:HIS:HD2	1.90	0.59
4:A:1409:HBP:N9	4:A:1409:HBP:C1	2.65	0.58
1:A:38:ALA:HB2	1:A:178:LEU:HD23	1.86	0.58
1:A:337:TYR:HA	1:A:443:MET:HE1	1.84	0.58
1:A:328:VAL:O	1:A:427:ALA:HA	2.03	0.58
1:B:184:GLN:OE1	1:B:194:PRO:HB2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:ARG:HD3	6:B:568:HOH:O	2.04	0.58
4:B:545:HBP:N9	4:B:545:HBP:C1	2.63	0.58
1:A:29:VAL:HG21	1:A:136:ARG:HB2	1.86	0.58
1:A:253:ARG:NH1	6:A:1547:HOH:O	2.33	0.58
1:B:96:CYS:O	1:B:150:ASN:HB2	2.04	0.58
1:B:459:LEU:HD11	1:B:474:GLN:HG3	1.86	0.58
1:A:374:ALA:HB1	1:A:536:LEU:HD21	1.86	0.58
1:A:235:PRO:HB2	1:A:296:ARG:NH2	2.18	0.57
1:B:68:VAL:HG23	1:B:90:ARG:HB2	1.85	0.57
1:B:203:SER:OG	3:B:544:CO3:O1	2.22	0.57
1:A:118:ILE:O	1:A:205:GLY:HA3	2.04	0.57
1:A:424:ARG:CD	6:A:1440:HOH:O	2.53	0.57
1:A:200:PHE:CB	1:A:226:VAL:HB	2.35	0.56
1:A:84:GLU:HG3	6:A:1593:HOH:O	2.03	0.56
1:A:166:GLU:OE2	1:A:267:THR:HG23	2.05	0.56
1:A:337:TYR:CE2	1:A:341:TYR:HE1	2.24	0.56
1:B:245:ARG:O	1:B:249:THR:CG2	2.44	0.56
1:B:341:TYR:CE2	4:B:545:HBP:H12	2.40	0.56
1:A:227:LEU:HD22	1:A:414:LEU:CD2	2.36	0.56
1:B:313:GLU:HG2	6:B:580:HOH:O	2.06	0.56
1:B:113:PRO:HG2	1:B:485:ARG:HG2	1.86	0.55
1:B:497:SER:N	1:B:498:PRO:HA	2.19	0.55
1:B:511:VAL:HB	1:B:518:LEU:HD22	1.88	0.55
1:B:10:VAL:HG22	1:B:107:ARG:HH11	1.72	0.55
1:B:326:VAL:HG21	1:B:418:LEU:HD13	1.88	0.55
1:B:197:VAL:N	1:B:223:HIS:CD2	2.70	0.54
1:A:200:PHE:HB2	1:A:226:VAL:HB	1.89	0.54
1:B:504:THR:HB	6:B:602:HOH:O	2.07	0.54
4:B:545:HBP:C14	4:B:545:HBP:C26	2.84	0.54
1:B:177:ARG:CZ	1:B:217:PRO:HB2	2.37	0.54
1:A:227:LEU:HB2	1:A:328:VAL:HG12	1.90	0.54
1:A:339:LEU:HD11	1:A:399:SER:HA	1.90	0.54
1:A:197:VAL:HB	1:A:222:PHE:HA	1.89	0.54
1:B:252:ALA:HA	1:B:273:LEU:HD21	1.89	0.54
1:B:453:PHE:HB3	1:B:476:LEU:HD12	1.88	0.54
1:A:203:SER:CB	3:A:1408:CO3:C	2.80	0.53
1:B:118:ILE:O	1:B:205:GLY:HA3	2.07	0.53
1:A:355:SER:OG	1:A:358:GLN:HG3	2.08	0.53
1:B:81:GLU:HB3	1:B:438:THR:HG21	1.90	0.53
1:B:444:GLY:O	1:B:446:PRO:HD3	2.09	0.53
1:A:414:LEU:HG	1:A:418:LEU:HD22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:GLU:HG3	6:A:1558:HOH:O	2.08	0.53
1:A:453:PHE:HB3	1:A:476:LEU:HD12	1.90	0.53
1:B:210:GLY:HA3	1:B:232:PRO:HG3	1.91	0.53
1:B:339:LEU:HD13	1:B:346:PHE:CE2	2.44	0.52
1:A:514:ASN:HB2	6:A:1461:HOH:O	2.09	0.52
1:B:319:GLY:O	1:B:421:GLN:HG3	2.09	0.52
1:B:341:TYR:CZ	4:B:545:HBP:H12	2.45	0.52
1:A:380:LEU:HD12	1:B:535:PHE:HB2	1.92	0.52
1:A:104:PRO:HG3	1:A:143:GLY:HA2	1.92	0.52
1:A:139:ALA:HA	1:A:144:ALA:O	2.10	0.52
1:A:430:PHE:HE2	1:A:476:LEU:HD11	1.75	0.52
1:A:380:LEU:HA	1:A:385:TRP:HZ2	1.76	0.52
1:A:337:TYR:HE2	1:A:341:TYR:HE1	1.56	0.51
1:B:404:ASP:HA	1:B:408:VAL:HB	1.92	0.51
1:A:228:GLN:HG2	1:A:428:TYR:OH	2.10	0.51
1:A:337:TYR:CE2	4:A:1409:HBP:C26	2.94	0.51
1:B:295:PHE:CE2	1:B:338:PHE:CE1	2.98	0.51
1:A:128:ALA:HB1	1:A:148:SER:OG	2.10	0.51
1:A:337:TYR:CD2	4:A:1409:HBP:N27	2.78	0.51
1:A:352:SER:O	1:A:395:ARG:HG3	2.11	0.51
1:A:376:GLU:HG2	1:B:538:LYS:HD2	1.93	0.51
1:A:424:ARG:HG2	1:A:426:TYR:CE1	2.46	0.51
1:B:346:PHE:CE1	1:B:354:ILE:HG21	2.46	0.51
1:B:472:PHE:CZ	1:B:476:LEU:HD11	2.46	0.51
1:B:495:SER:OG	1:B:496:LYS:N	2.43	0.51
1:B:374:ALA:HA	1:B:539:LEU:CD2	2.41	0.50
1:A:286:TRP:CE3	4:A:1409:HBP:C3	2.94	0.50
1:A:479:TYR:OH	1:A:518:LEU:HG	2.12	0.50
1:B:166:GLU:HG2	1:B:270:ILE:HD13	1.92	0.50
1:B:213:ILE:O	1:B:219:ARG:HD3	2.11	0.50
1:B:393:HIS:HB3	6:B:626:HOH:O	2.10	0.50
1:A:21:ARG:HG3	1:A:105:TYR:CE1	2.47	0.50
1:A:50:PRO:HD2	1:A:174:LEU:HB3	1.94	0.50
1:B:200:PHE:HB2	1:B:226:VAL:HB	1.92	0.50
1:B:227:LEU:HD22	1:B:414:LEU:HD21	1.94	0.50
1:A:76:LEU:HD22	1:A:341:TYR:CD2	2.47	0.50
1:A:166:GLU:HB2	1:A:270:ILE:HD13	1.92	0.50
1:A:471:ILE:HA	1:A:474:GLN:HE21	1.77	0.50
1:A:541:SER:O	1:A:542:ALA:HB2	2.10	0.50
1:A:382:TYR:O	1:A:393:HIS:HE1	1.94	0.50
1:B:29:VAL:HG21	1:B:136:ARG:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:431:GLU:CD	1:A:521:ARG:HH12	2.14	0.49
1:B:46:ARG:HB3	1:B:274:ARG:HG2	1.95	0.49
1:A:337:TYR:CG	4:A:1409:HBP:N27	2.80	0.49
1:B:340:VAL:HG11	1:B:443:MET:CE	2.42	0.49
1:A:86:TRP:CG	4:A:1409:HBP:H22	2.48	0.49
1:A:124:TYR:CD1	1:A:124:TYR:C	2.86	0.49
1:A:137:PHE:HB2	1:A:455:PHE:O	2.13	0.49
1:A:397:ALA:O	1:A:401:VAL:HG23	2.13	0.49
1:B:336:SER:HB2	1:B:443:MET:HA	1.95	0.49
1:A:122:GLY:O	1:A:123:PHE:HB2	2.13	0.49
1:B:337:TYR:HB2	6:B:604:HOH:O	2.13	0.49
1:A:177:ARG:CZ	1:A:217:PRO:HB2	2.43	0.48
1:B:525:ARG:CG	1:B:525:ARG:NH1	2.71	0.48
1:B:103:THR:HG21	1:B:190:PHE:HB3	1.94	0.48
1:A:252:ALA:HA	1:A:273:LEU:HD21	1.94	0.48
1:B:181:GLN:HG2	6:B:673:HOH:O	2.13	0.48
1:B:380:LEU:HD22	1:B:385:TRP:CZ2	2.49	0.48
1:A:86:TRP:CB	4:A:1409:HBP:H22	2.44	0.48
1:A:354:ILE:O	1:A:391:PRO:HB3	2.14	0.48
1:A:433:ARG:NH2	1:A:439:TRP:O	2.47	0.48
1:B:376:GLU:O	1:B:380:LEU:HG	2.13	0.48
1:B:478:LYS:HD2	1:B:491:ASP:OD2	2.14	0.48
1:B:119:TYR:CE1	1:B:151:TYR:CE1	3.02	0.47
1:B:141:VAL:HG21	1:B:459:LEU:HD23	1.96	0.47
1:A:50:PRO:HB2	1:A:178:LEU:HD22	1.95	0.47
1:A:376:GLU:O	1:A:380:LEU:HG	2.14	0.47
1:B:348:LYS:HA	1:B:440:PRO:HG3	1.96	0.47
1:A:167:ALA:HB2	1:A:270:ILE:HG12	1.96	0.47
1:A:213:ILE:HA	1:A:219:ARG:HD3	1.95	0.47
1:A:231:THR:HG21	1:A:411:VAL:HA	1.97	0.47
1:B:340:VAL:HG11	1:B:443:MET:HE1	1.96	0.47
1:A:5:ASP:C	1:A:5:ASP:OD1	2.53	0.47
1:A:293:SER:HB3	1:A:368:PRO:HB3	1.97	0.47
1:A:360:LEU:CD2	1:A:379:VAL:HG11	2.44	0.47
1:B:107:ARG:HB3	6:B:583:HOH:O	2.14	0.47
1:B:224:ARG:HD2	1:B:483:PHE:O	2.15	0.47
1:B:238:THR:HA	1:B:301:PRO:O	2.14	0.47
1:A:338:PHE:HZ	1:A:447:HIS:CD2	2.33	0.47
1:A:374:ALA:HA	1:A:539:LEU:HD23	1.97	0.47
1:A:183:VAL:HG13	1:A:187:ILE:HB	1.97	0.47
1:B:425:VAL:O	1:B:503:TYR:N	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:ARG:HD3	1:B:47:PHE:CZ	2.51	0.46
1:B:327:LEU:HD11	1:B:500:TRP:CH2	2.49	0.46
1:A:224:ARG:CD	1:A:487:GLY:HA2	2.46	0.46
1:A:369:GLN:HB2	6:A:1435:HOH:O	2.16	0.46
1:A:504:THR:HB	6:A:1585:HOH:O	2.16	0.46
1:B:296:ARG:HG2	6:B:641:HOH:O	2.15	0.46
1:A:468:GLU:CD	1:A:468:GLU:H	2.18	0.46
1:B:337:TYR:HD2	1:B:337:TYR:O	1.99	0.46
1:A:338:PHE:CZ	1:A:447:HIS:CD2	3.04	0.46
1:B:130:LEU:HD12	1:B:133:TYR:CE2	2.50	0.46
1:B:213:ILE:HD11	1:B:225:ALA:CB	2.46	0.46
1:A:224:ARG:HD3	1:A:487:GLY:HA2	1.98	0.46
1:A:330:VAL:HG11	1:A:408:VAL:HA	1.98	0.46
1:B:100:ASN:O	1:B:147:VAL:HA	2.16	0.46
1:A:428:TYR:HB3	1:A:500:TRP:CZ2	2.51	0.46
1:B:528:THR:O	1:B:531:PHE:HB3	2.16	0.46
1:B:116:ILE:HD11	1:B:183:VAL:HG11	1.98	0.45
1:B:210:GLY:CA	1:B:232:PRO:HG3	2.45	0.45
1:B:509:GLN:HA	1:B:521:ARG:O	2.16	0.45
1:A:337:TYR:HA	1:A:443:MET:HE3	1.98	0.45
1:B:536:LEU:HD22	1:B:536:LEU:HA	1.87	0.45
1:B:187:ILE:HD12	1:B:187:ILE:HA	1.83	0.45
1:B:227:LEU:HD22	1:B:414:LEU:CD2	2.47	0.45
1:A:22:LEU:HD13	1:A:136:ARG:NH1	2.32	0.45
1:A:24:ALA:HB3	1:A:140:GLN:HG3	1.99	0.45
1:A:101:VAL:HG22	1:A:147:VAL:HG22	1.98	0.45
1:A:119:TYR:CD1	1:A:126:GLY:HA3	2.52	0.45
1:A:317:ASN:OD1	1:A:417:ARG:HG3	2.17	0.45
1:B:202:GLU:HA	1:B:228:GLN:O	2.17	0.45
1:B:337:TYR:O	1:B:340:VAL:HG22	2.17	0.45
1:A:177:ARG:NH1	1:A:217:PRO:HB2	2.32	0.45
1:A:226:VAL:HG22	1:A:327:LEU:HB3	1.99	0.45
1:A:329:GLY:HA3	1:A:428:TYR:CE2	2.52	0.45
1:A:334:GLU:CD	1:A:407:VAL:HG11	2.37	0.45
1:A:337:TYR:HE2	1:A:341:TYR:CE1	2.34	0.45
1:A:339:LEU:HD11	1:A:399:SER:CA	2.47	0.45
1:A:291:GLN:HE22	1:A:369:GLN:HE21	1.64	0.44
1:A:326:VAL:O	1:A:425:VAL:HA	2.17	0.44
1:A:364:ARG:N	1:A:364:ARG:HD2	2.32	0.44
1:B:325:GLN:HG3	1:B:424:ARG:HB3	1.98	0.44
1:A:33:LEU:HD22	1:A:65:PHE:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:MET:HA	1:A:48:MET:CE	2.47	0.44
1:A:337:TYR:CD2	4:A:1409:HBP:C26	3.01	0.44
1:B:210:GLY:HA3	1:B:232:PRO:HD3	2.00	0.44
1:B:294:ILE:HG12	1:B:365:ILE:HG22	2.00	0.44
1:A:86:TRP:HB3	4:A:1409:HBP:H22	1.98	0.44
1:A:160:ALA:O	1:A:162:PRO:HD3	2.18	0.44
1:B:48:MET:SD	1:B:166:GLU:HA	2.58	0.44
1:A:17:LEU:HD23	1:A:60:LEU:HB3	1.99	0.44
1:B:300:VAL:HB	1:B:301:PRO:HD2	2.00	0.44
1:B:511:VAL:CB	1:B:518:LEU:HD22	2.47	0.44
1:A:170:ASN:OD1	1:A:301:PRO:HA	2.17	0.44
1:A:274:ARG:HD3	6:A:1411:HOH:O	2.16	0.44
1:B:291:GLN:HE22	1:B:369:GLN:HE21	1.66	0.44
1:B:340:VAL:CG1	1:B:443:MET:CE	2.96	0.44
1:A:187:ILE:HG13	1:A:192:GLY:HA3	2.00	0.44
1:A:440:PRO:HD2	1:A:443:MET:SD	2.58	0.44
1:B:364:ARG:O	1:B:368:PRO:HA	2.18	0.44
1:B:454:ILE:HD13	1:B:476:LEU:HB3	2.00	0.44
1:A:67:ASN:O	1:A:150:ASN:ND2	2.51	0.44
1:A:85:MET:CE	1:A:132:VAL:HG11	2.48	0.44
1:A:89:ASN:ND2	1:A:129:SER:O	2.41	0.44
1:B:117:TRP:HA	1:B:200:PHE:O	2.17	0.44
1:B:331:VAL:HG22	1:B:334:GLU:OE2	2.17	0.44
1:A:476:LEU:HD22	1:A:513:LEU:HD13	2.00	0.43
1:B:337:TYR:HB3	4:B:545:HBP:O28	2.17	0.43
1:A:61:ASP:OD1	1:A:61:ASP:C	2.57	0.43
1:A:130:LEU:HD12	1:A:133:TYR:CE2	2.53	0.43
1:A:512:SER:HB3	1:A:521:ARG:CZ	2.48	0.43
1:B:237:ALA:O	1:B:301:PRO:HD2	2.18	0.43
1:A:10:VAL:HG23	1:A:107:ARG:NH1	2.33	0.43
1:A:5:ASP:OD2	1:A:105:TYR:OH	2.27	0.43
1:B:346:PHE:HE2	1:B:395:ARG:HG2	1.84	0.43
1:B:329:GLY:HA3	1:B:428:TYR:CZ	2.53	0.43
1:A:48:MET:CE	1:A:165:ARG:O	2.62	0.43
1:A:202:GLU:HA	1:A:228:GLN:O	2.19	0.43
1:A:172:GLY:O	1:A:176:GLN:HG3	2.19	0.43
1:B:99:LEU:HA	1:B:149:MET:HA	2.00	0.43
1:B:395:ARG:CZ	1:B:442:TRP:HB2	2.48	0.43
1:A:48:MET:CE	1:A:166:GLU:HA	2.35	0.43
1:A:152:ARG:HG2	6:A:1416:HOH:O	2.18	0.43
1:A:294:ILE:HD11	1:A:402:VAL:HG21	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:GLY:HA3	1:A:428:TYR:CD2	2.54	0.43
1:A:373:LEU:CD2	1:A:539:LEU:HD11	2.48	0.43
1:A:441:LEU:HD23	1:A:441:LEU:HA	1.87	0.43
1:A:5:ASP:HB3	1:A:8:LEU:HD12	1.99	0.43
1:A:360:LEU:HD12	6:A:1514:HOH:O	2.18	0.43
1:B:4:GLU:OE2	1:B:18:ARG:NH1	2.51	0.42
1:B:139:ALA:HA	1:B:144:ALA:O	2.19	0.42
1:B:333:ASP:OD1	1:B:444:GLY:HA3	2.19	0.42
1:B:380:LEU:HD22	1:B:385:TRP:HZ2	1.84	0.42
1:A:535:PHE:CE1	1:B:377:ALA:HB2	2.54	0.42
1:A:210:GLY:HA3	1:A:232:PRO:HD3	2.01	0.42
1:A:224:ARG:HD3	1:A:487:GLY:CA	2.49	0.42
1:B:433:ARG:NH2	1:B:439:TRP:O	2.44	0.42
1:A:13:ARG:NH2	6:A:1490:HOH:O	2.48	0.42
1:A:216:LEU:HB3	1:A:217:PRO:HD3	2.00	0.42
1:A:227:LEU:N	1:A:227:LEU:HD12	2.34	0.42
1:B:211:MET:HE2	1:B:232:PRO:HA	2.02	0.42
1:A:528:THR:O	1:A:531:PHE:HB3	2.20	0.42
1:B:203:SER:HA	1:B:229:SER:O	2.20	0.42
1:B:235:PRO:HG3	1:B:405:HIS:CE1	2.55	0.42
1:B:290:PRO:HG2	1:B:291:GLN:NE2	2.35	0.42
1:B:294:ILE:HD11	1:B:402:VAL:HG21	2.02	0.42
1:B:440:PRO:HG2	1:B:443:MET:HG3	2.00	0.42
1:B:5:ASP:HA	1:B:6:PRO:HD2	1.94	0.42
1:B:81:GLU:O	1:B:85:MET:HG2	2.20	0.42
1:A:128:ALA:HB1	1:A:148:SER:CB	2.50	0.41
1:A:265:ASN:C	1:A:265:ASN:OD1	2.59	0.41
1:B:50:PRO:HG3	1:B:175:ASP:HA	2.02	0.41
1:B:187:ILE:HG13	1:B:192:GLY:HA3	2.02	0.41
1:B:198:THR:OG1	1:B:484:ALA:HA	2.20	0.41
1:B:346:PHE:CZ	1:B:354:ILE:HD13	2.56	0.41
1:B:478:LYS:O	1:B:482:ASN:HB2	2.20	0.41
1:A:472:PHE:CZ	1:A:476:LEU:HD11	2.54	0.41
1:B:252:ALA:HB1	1:B:257:CYS:HB2	2.03	0.41
1:A:80:PHE:CG	1:A:348:LYS:HE2	2.54	0.41
1:A:116:ILE:HA	1:A:147:VAL:O	2.20	0.41
1:A:119:TYR:CE1	1:A:151:TYR:CE1	3.07	0.41
1:A:226:VAL:HA	1:A:327:LEU:O	2.21	0.41
1:A:519:GLU:HB2	6:A:1551:HOH:O	2.20	0.41
1:B:329:GLY:HA3	1:B:428:TYR:CE1	2.56	0.41
1:A:4:GLU:HG2	1:A:9:LEU:HD21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:ILE:N	1:A:454:ILE:HD13	2.35	0.41
1:B:384:ASP:HA	5:B:546:P3G:H72	2.03	0.41
1:A:5:ASP:HA	1:A:6:PRO:HD2	1.84	0.41
1:B:342:GLY:O	1:B:344:PRO:HD3	2.20	0.41
1:A:88:PRO:CG	1:A:92:LEU:HD21	2.51	0.41
1:A:276:ARG:HD3	1:A:276:ARG:HA	1.84	0.41
1:A:337:TYR:O	1:A:340:VAL:HG22	2.21	0.41
1:A:406:ASN:C	1:A:410:PRO:HG2	2.41	0.41
1:A:512:SER:OG	1:A:519:GLU:HB3	2.20	0.41
1:B:9:LEU:HD23	1:B:9:LEU:HA	1.93	0.41
1:A:38:ALA:HB3	1:A:97:LEU:HB3	2.02	0.41
1:A:459:LEU:HD11	1:A:474:GLN:HG3	2.03	0.40
1:B:17:LEU:HD23	1:B:60:LEU:HB3	2.02	0.40
1:B:21:ARG:NH1	1:B:28:PRO:HB3	2.36	0.40
1:B:32:PHE:CG	1:B:190:PHE:HE2	2.39	0.40
1:A:12:VAL:HG11	1:A:182:TRP:CZ2	2.56	0.40
1:A:203:SER:HA	1:A:229:SER:O	2.21	0.40
1:B:211:MET:CE	1:B:301:PRO:HG2	2.52	0.40
1:B:227:LEU:N	1:B:227:LEU:HD12	2.36	0.40
1:B:233:ASN:HA	1:B:312:PRO:HB2	2.03	0.40
1:B:288:VAL:CG2	1:B:298:SER:HB3	2.51	0.40
1:A:491:ASP:HB3	1:A:494:ASP:HB3	2.03	0.40
1:A:10:VAL:CG2	1:A:107:ARG:NH1	2.84	0.40
1:A:99:LEU:HD12	1:A:99:LEU:C	2.41	0.40
1:A:354:ILE:HB	1:A:358:GLN:HB2	2.04	0.40
1:A:367:VAL:HG12	1:A:370:ALA:HB2	2.01	0.40
1:A:430:PHE:HA	1:A:513:LEU:HB2	2.02	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	531/543 (98%)	510 (96%)	21 (4%)	0	100	100
1	B	529/543 (97%)	501 (95%)	24 (4%)	4 (1%)	16	31
All	All	1060/1086 (98%)	1011 (95%)	45 (4%)	4 (0%)	30	49

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	495	SER
1	B	541	SER
1	B	111	PRO
1	B	497	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	439/443 (99%)	411 (94%)	28 (6%)	14	30
1	B	438/443 (99%)	418 (95%)	20 (5%)	23	45
All	All	877/886 (99%)	829 (94%)	48 (6%)	18	37

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ARG
1	A	4	GLU
1	A	13	ARG
1	A	21	ARG
1	A	23	LYS
1	A	45	ARG
1	A	54	ARG
1	A	70	TYR
1	A	184	GLN
1	A	246	ARG
1	A	249	THR
1	A	265	ASN

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Mol	Chain	Res	Type
1	A	267	THR
1	A	291	GLN
1	A	295	PHE
1	A	301	PRO
1	A	322	GLN
1	A	323	ASP
1	A	337	TYR
1	A	356	ARG
1	A	386	LEU
1	A	413	GLN
1	A	418	LEU
1	A	437	LEU
1	A	497	SER
1	A	524	LEU
1	A	536	LEU
1	A	540	LEU
1	B	45	ARG
1	B	115	LEU
1	B	181	GLN
1	B	184	GLN
1	B	200	PHE
1	B	246	ARG
1	B	249	THR
1	B	281	LEU
1	B	291	GLN
1	B	295	PHE
1	B	337	TYR
1	B	348	LYS
1	B	386	LEU
1	B	413	GLN
1	B	437	LEU
1	B	438	THR
1	B	441	LEU
1	B	525	ARG
1	B	536	LEU
1	B	541	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	184	GLN
1	A	223	HIS

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Mol	Chain	Res	Type
1	A	291	GLN
1	A	322	GLN
1	A	393	HIS
1	A	474	GLN
1	B	223	HIS
1	B	287	HIS
1	B	291	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

7 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	$\# Z > 2$
5	P3G	B	546	-	6,6,16	0.55	0	5,5,15	0.16	0
2	NAG	A	1402	1	14,14,15	1.00	1 (7%)	17,19,21	1.43	4 (23%)
2	NAG	A	1407	1	13,13,15	1.03	1 (7%)	16,17,21	2.67	5 (31%)
4	HBP	A	1409	-	26,26,26	0.71	0	27,31,31	1.91	5 (18%)
3	CO3	B	544	-	3,3,3	0.90	0	2,3,3	0.12	0
4	HBP	B	545	-	26,26,26	0.82	1 (3%)	27,31,31	1.98	6 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	CO3	A	1408	-	3,3,3	1.20	1 (33%)	2,3,3	0.80	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
5	P3G	B	546	-	-	2/4/4/14	-
2	NAG	A	1402	1	-	4/6/23/26	0/1/1/1
2	NAG	A	1407	1	-	4/5/22/26	0/1/1/1
4	HBP	A	1409	-	-	8/16/16/16	0/2/2/2
4	HBP	B	545	-	-	7/16/16/16	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	545	HBP	C25-N20	-2.72	1.34	1.37
2	A	1402	NAG	C1-C2	2.70	1.56	1.52
2	A	1407	NAG	C1-C2	2.45	1.55	1.52
3	A	1408	CO3	O1-C	2.04	1.32	1.25

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1407	NAG	C1-O5-C5	7.83	122.68	112.19
4	B	545	HBP	O10-N9-C8A	6.53	123.12	111.83
4	A	1409	HBP	O10-N9-C8A	5.04	120.54	111.83
2	A	1407	NAG	C7-N2-C2	4.64	121.50	114.43
4	A	1409	HBP	C21-N20-C25	-4.12	118.74	121.54
4	A	1409	HBP	C24-C25-N20	3.71	122.61	117.25
4	B	545	HBP	C21-N20-C25	-3.70	119.02	121.54
4	B	545	HBP	C24-C25-N20	3.67	122.55	117.25
2	A	1402	NAG	C4-C3-C2	3.09	115.55	111.02
4	A	1409	HBP	C3-N2-C7	-2.92	119.55	121.54
4	B	545	HBP	C25-C26-N27	2.72	123.01	117.75
4	A	1409	HBP	C22-C23-C24	-2.61	117.02	120.24
4	B	545	HBP	C22-C23-C24	-2.57	117.07	120.24
4	B	545	HBP	C3-N2-C7	-2.49	119.85	121.54
2	A	1407	NAG	O3-C3-C4	-2.47	104.56	110.38
2	A	1402	NAG	C2-N2-C7	2.24	125.90	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1402	NAG	O5-C5-C6	2.21	111.97	107.66
2	A	1407	NAG	C4-C3-C2	2.18	114.21	111.02
2	A	1407	NAG	O3-C3-C2	2.18	113.92	109.40
2	A	1402	NAG	O7-C7-C8	-2.04	118.43	122.05

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1402	NAG	C1-C2-N2-C7
4	A	1409	HBP	C2-C1-N2-C3
4	A	1409	HBP	N2-C7-C8A-N9
4	A	1409	HBP	C14-C17-N20-C21
4	A	1409	HBP	C14-C17-N20-C25
4	A	1409	HBP	N2-C1-C2-C5
2	A	1407	NAG	O5-C5-C6-O6
2	A	1407	NAG	C4-C5-C6-O6
5	B	546	P3G	O3-C7-C8-O4
4	B	545	HBP	C11-C14-C17-N20
2	A	1407	NAG	C3-C2-N2-C7
4	B	545	HBP	C1-C2-C5-C8
4	B	545	HBP	C2-C5-C8-C11
2	A	1402	NAG	C8-C7-N2-C2
2	A	1407	NAG	C1-C2-N2-C7
2	A	1402	NAG	C3-C2-N2-C7
4	B	545	HBP	C6-C7-C8A-N9
4	A	1409	HBP	C2-C1-N2-C7
4	B	545	HBP	C14-C17-N20-C21
4	B	545	HBP	C14-C17-N20-C25
2	A	1402	NAG	O7-C7-N2-C2
4	A	1409	HBP	C2-C5-C8-C11
4	B	545	HBP	N2-C7-C8A-N9
4	A	1409	HBP	C8-C11-C14-C17
5	B	546	P3G	C7-C8-O4-C9

There are no ring outliers.

5 monomers are involved in 33 short contacts:

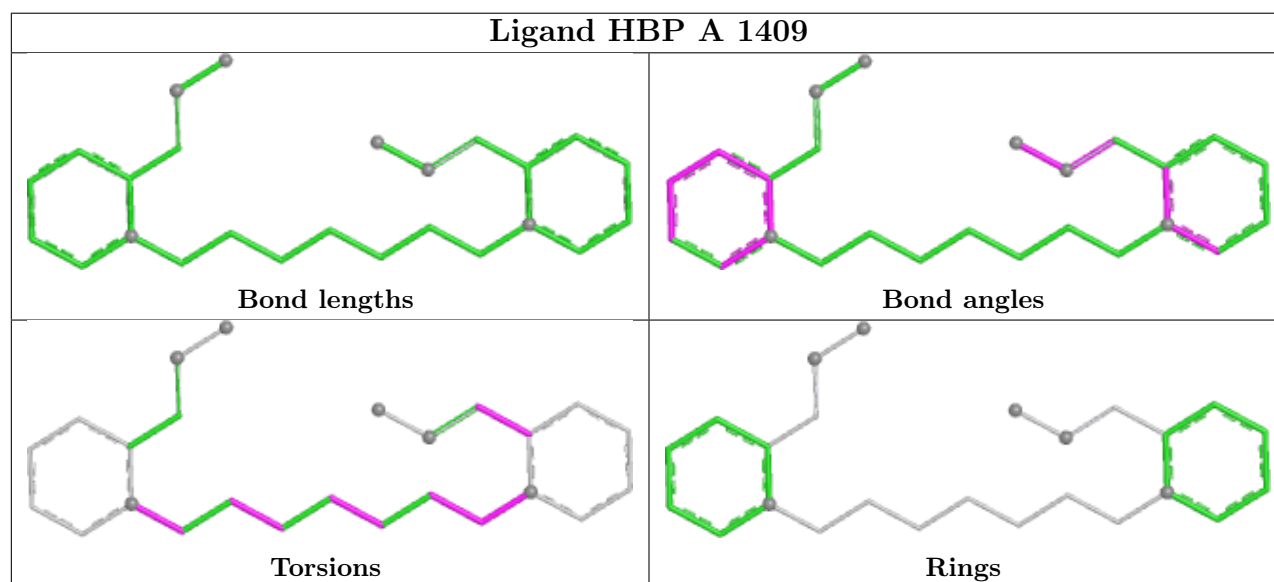
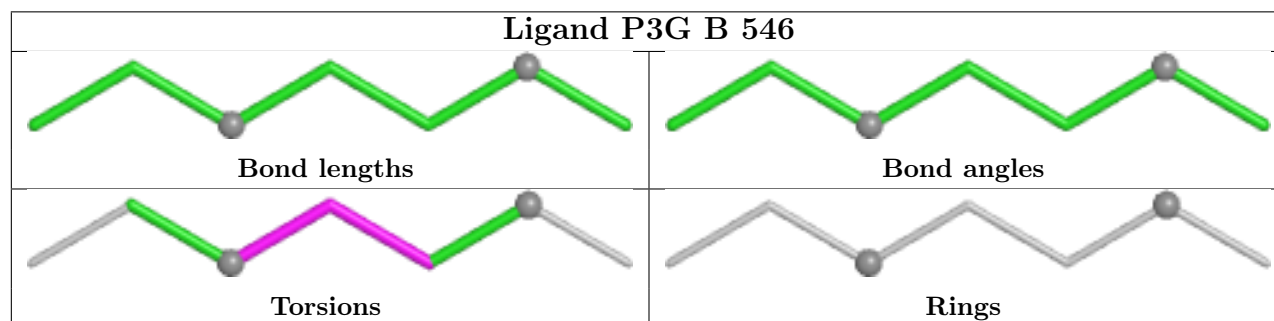
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	546	P3G	1	0
4	A	1409	HBP	15	0

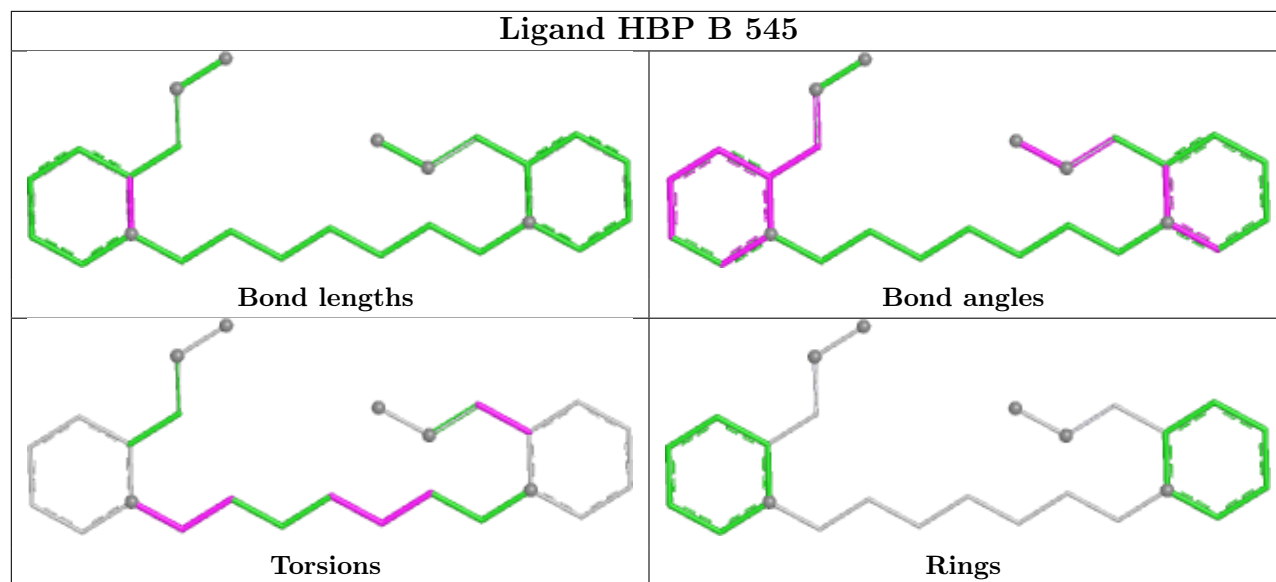
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	544	CO3	3	0
4	B	545	HBP	9	0
3	A	1408	CO3	5	0

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 [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	535/543 (98%)	-0.25	8 (1%) 71 68	29, 39, 55, 71	0
1	B	533/543 (98%)	-0.12	8 (1%) 71 68	32, 40, 53, 74	0
All	All	1068/1086 (98%)	-0.19	16 (1%) 71 68	29, 39, 54, 74	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	542	ALA	4.2
1	B	495	SER	3.2
1	B	497	SER	3.1
1	B	496	LYS	3.0
1	A	496	LYS	2.9
1	B	257	CYS	2.8
1	B	286	TRP	2.6
1	A	540	LEU	2.6
1	A	286	TRP	2.5
1	B	543	THR	2.5
1	A	491	ASP	2.5
1	A	492	PRO	2.4
1	B	338	PHE	2.2
1	A	447	HIS	2.0
1	A	338	PHE	2.0
1	B	447	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

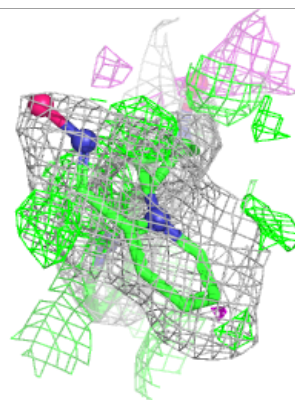
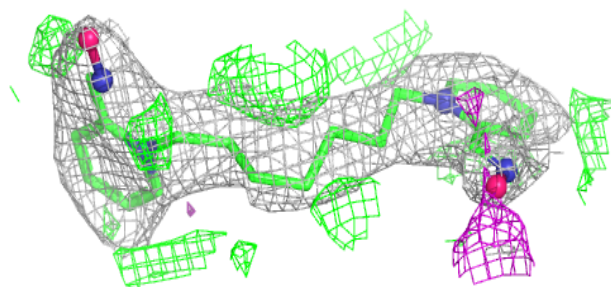
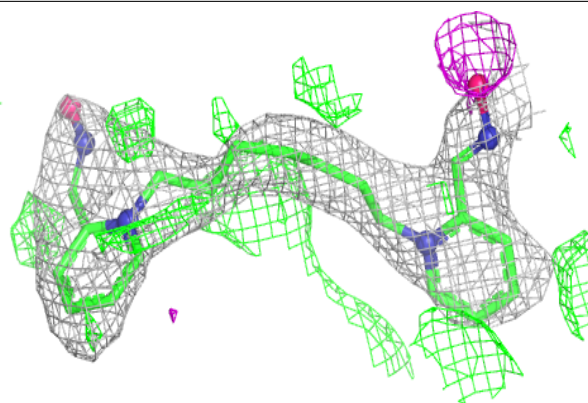
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	A	1402	14/15	0.65	0.19	74,82,82,84	0
2	NAG	A	1407	13/15	0.69	0.22	64,68,69,69	0
4	HBP	A	1409	25/25	0.83	0.21	53,55,56,59	25
4	HBP	B	545	25/25	0.83	0.28	61,67,78,79	25
5	P3G	B	546	7/17	0.84	0.26	78,80,85,86	0
3	CO3	B	544	4/4	0.95	0.12	42,42,42,42	0
3	CO3	A	1408	4/4	0.96	0.10	36,37,37,37	0

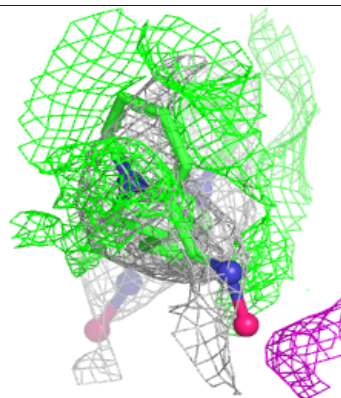
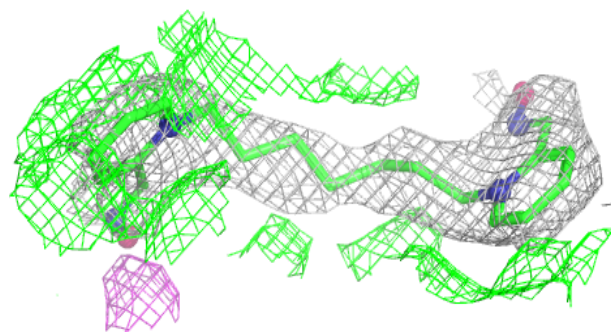
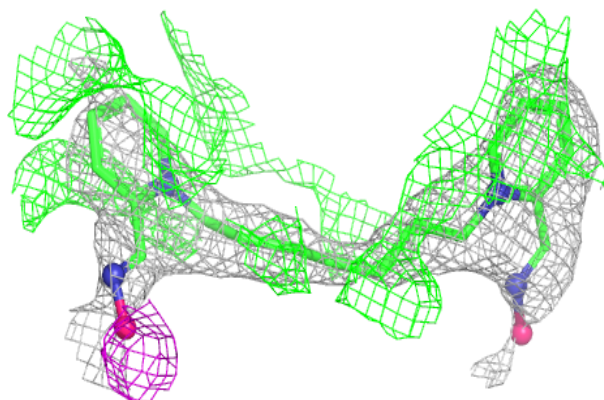
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

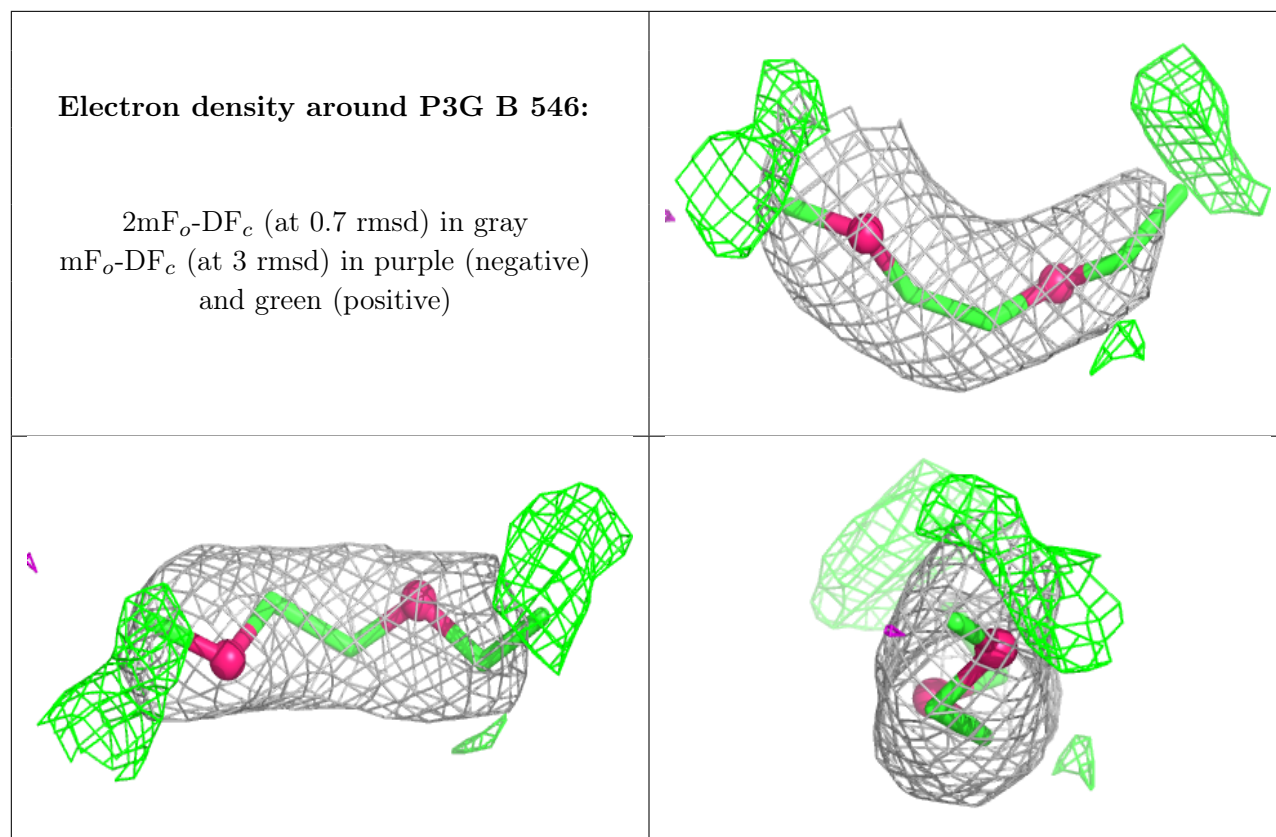
Electron density around HBP A 1409:

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

**Electron density around HBP B 545:**

$2mF_o-DF_c$ (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.