



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2023 – 07:45 AM EDT

PDB ID : 2ZVT
Title : Cys285Ser mutant PPARgamma ligand-binding domain complexed with 15-d
eoxy-delta12,14-prostaglandin J2
Authors : Waku, T.; Oyama, T.; Shiraki, T.; Morikawa, K.
Deposited on : 2008-11-19
Resolution : 1.90 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

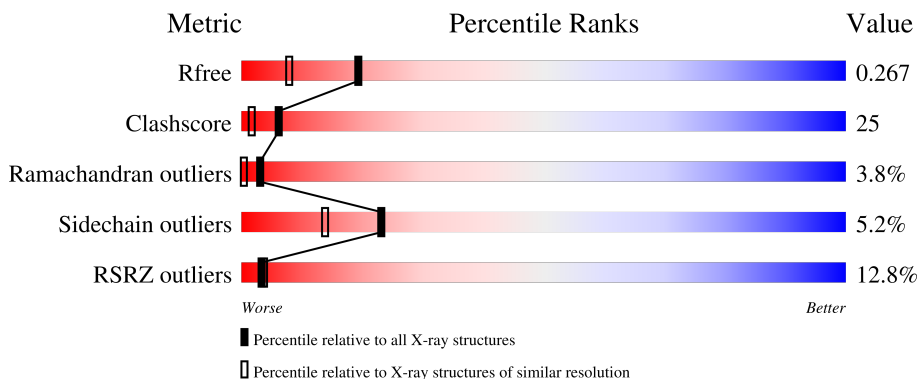
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	286	
1	B	286	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PTG	B	2	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4538 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.

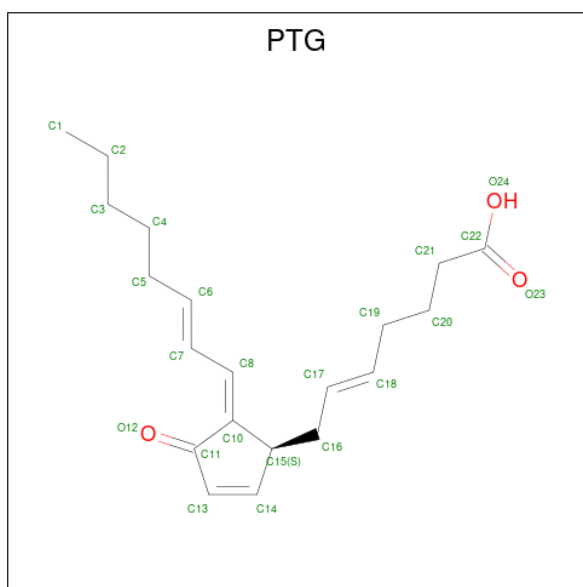
- Molecule 1 is a protein called Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	274	Total	C	N	O	S	0	0	0
			2198	1417	360	412	9			
1	B	263	Total	C	N	O	S	0	0	0
			2110	1363	346	393	8			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	191	GLY	-	expression tag	UNP P37231
A	192	SER	-	expression tag	UNP P37231
A	193	HIS	-	expression tag	UNP P37231
A	194	MET	-	expression tag	UNP P37231
A	285	SER	CYS	engineered mutation	UNP P37231
B	191	GLY	-	expression tag	UNP P37231
B	192	SER	-	expression tag	UNP P37231
B	193	HIS	-	expression tag	UNP P37231
B	194	MET	-	expression tag	UNP P37231
B	285	SER	CYS	engineered mutation	UNP P37231

- Molecule 2 is (5E,14E)-11-oxoprostanoic acid (three-letter code: PTG) (formula: C₂₀H₂₈O₃).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	A	1	Total	C	O	0	0
			23	20	3		
2	B	1	Total	C	O	0	0
			23	20	3		

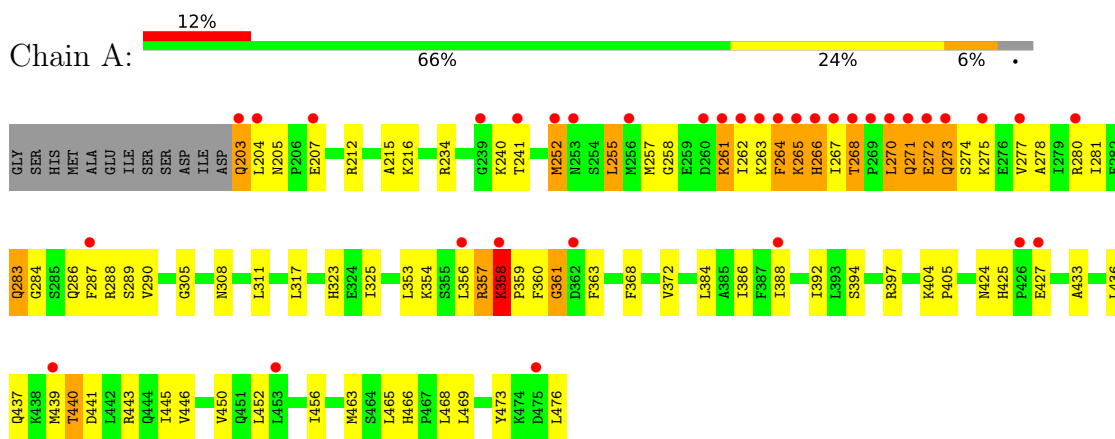
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	92	Total	O	0	0
			92	92		
3	B	92	Total	O	0	0
			92	92		

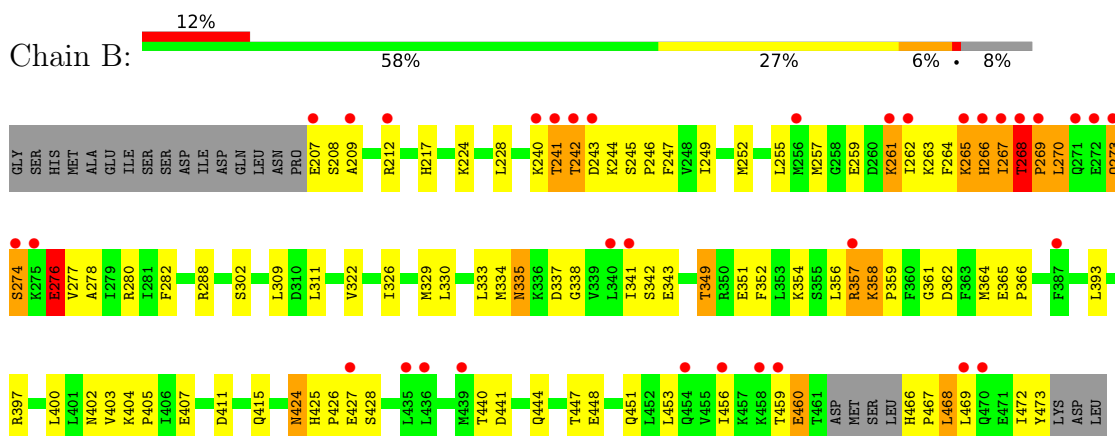
3 Residue-property plots i

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: Peroxisome proliferator-activated receptor gamma



- Molecule 1: Peroxisome proliferator-activated receptor gamma



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	92.98Å 61.39Å 118.54Å 90.00° 102.89° 90.00°	Depositor
Resolution (Å)	29.67 – 1.90 29.67 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.9 (29.67-1.90) 94.0 (29.67-1.90)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 1.91Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.251 , 0.275 0.244 , 0.267	Depositor DCC
R_{free} test set	2364 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtrriage
Anisotropy	0.535	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4538	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% 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: PTG

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	0/2236	0.63	0/3013
1	B	0.36	0/2146	0.65	2/2891 (0.1%)
All	All	0.37	0/4382	0.64	2/5904 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	267	ILE	N-CA-C	-9.20	86.16	111.00
1	B	266	HIS	N-CA-C	6.25	127.86	111.00

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	2198	0	2264	86	0
1	B	2110	0	2174	133	0
2	A	23	0	27	2	0
2	B	23	0	27	10	0
3	A	92	0	0	1	0
3	B	92	0	0	6	0
All	All	4538	0	4492	217	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 25.

All (217) 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:349:THR:HG22	1:B:352:PHE:H	1.13	1.13
1:B:266:HIS:ND1	1:B:268:THR:HB	1.65	1.12
1:A:240:LYS:HG3	1:A:241:THR:HG23	1.35	1.05
1:A:358:LYS:HD2	1:A:358:LYS:H	1.28	0.98
1:B:459:THR:HG22	1:B:460:GLU:H	1.30	0.96
1:A:205:ASN:HD22	1:A:207:GLU:HB2	1.29	0.95
1:B:358:LYS:HB3	1:B:359:PRO:CD	1.95	0.94
1:B:278:ALA:HB3	1:B:357:ARG:NH1	1.85	0.91
1:B:427:GLU:H	1:B:427:GLU:CD	1.73	0.91
1:A:205:ASN:ND2	1:A:207:GLU:HB2	1.86	0.90
1:A:273:GLN:C	1:A:275:LYS:H	1.76	0.86
1:A:272:GLU:HG2	1:A:273:GLN:H	1.42	0.85
1:A:272:GLU:HG2	1:A:273:GLN:HG3	1.59	0.85
1:B:266:HIS:C	1:B:268:THR:H	1.77	0.84
1:B:326:ILE:HG23	2:B:2:PTG:H3	1.60	0.83
1:B:335:ASN:C	1:B:335:ASN:HD22	1.85	0.80
1:B:268:THR:OG1	1:B:270:LEU:HG	1.84	0.78
1:B:335:ASN:ND2	1:B:337:ASP:H	1.82	0.78
1:A:358:LYS:HB2	1:A:359:PRO:HD3	1.66	0.77
1:B:358:LYS:HB3	1:B:359:PRO:HD3	1.66	0.77
1:A:240:LYS:HG3	1:A:241:THR:H	1.49	0.76
1:B:266:HIS:HB3	1:B:268:THR:HG22	1.65	0.76
1:A:267:ILE:O	1:A:287:PHE:HB2	1.86	0.76
1:A:358:LYS:HD2	1:A:358:LYS:N	2.02	0.74
1:B:311:LEU:H	1:B:311:LEU:HD22	1.51	0.74
1:B:242:THR:O	1:B:244:LYS:N	2.21	0.74
1:B:349:THR:HG22	1:B:352:PHE:N	1.98	0.73
1:B:266:HIS:C	1:B:268:THR:N	2.43	0.71
1:B:349:THR:CG2	1:B:352:PHE:H	1.98	0.71
1:A:252:MET:HE1	1:A:277:VAL:HG21	1.72	0.71
1:A:240:LYS:HG3	1:A:241:THR:CG2	2.19	0.69
1:B:264:PHE:O	1:B:265:LYS:C	2.31	0.69
1:A:273:GLN:C	1:A:275:LYS:N	2.47	0.68
1:B:459:THR:HG22	1:B:460:GLU:N	2.06	0.67
1:A:270:LEU:O	1:A:271:GLN:HG3	1.94	0.67
1:B:278:ALA:HB3	1:B:357:ARG:HH12	1.56	0.67
1:A:358:LYS:CB	1:A:359:PRO:HD3	2.25	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:ARG:CZ	1:B:358:LYS:HD2	2.24	0.66
1:B:244:LYS:O	1:B:244:LYS:HG3	1.96	0.66
1:B:425:HIS:HA	1:B:427:GLU:OE1	1.96	0.66
1:B:357:ARG:HH21	1:B:358:LYS:CE	2.09	0.65
1:B:357:ARG:NH2	1:B:358:LYS:HD2	2.10	0.65
1:B:311:LEU:HD22	1:B:311:LEU:N	2.11	0.65
1:B:208:SER:O	1:B:212:ARG:HG3	1.97	0.65
1:B:459:THR:CG2	1:B:460:GLU:H	2.03	0.64
1:A:273:GLN:O	1:A:275:LYS:N	2.31	0.63
1:B:358:LYS:HD3	1:B:460:GLU:OE1	1.98	0.63
1:B:252:MET:SD	1:B:277:VAL:HG21	2.38	0.63
1:B:261:LYS:HE3	1:B:261:LYS:HA	1.80	0.63
1:B:466:HIS:HD2	1:B:468:LEU:H	1.46	0.63
1:B:354:LYS:HA	1:B:361:GLY:O	1.99	0.63
1:B:447:THR:O	1:B:451:GLN:HG3	1.98	0.62
1:B:267:ILE:O	1:B:269:PRO:HD3	1.99	0.62
1:B:266:HIS:HB3	1:B:268:THR:CG2	2.29	0.62
1:A:441:ASP:O	1:A:445:ILE:HG12	2.00	0.61
1:A:452:LEU:O	1:A:456:ILE:HG12	2.01	0.61
1:B:335:ASN:HD22	1:B:337:ASP:H	1.50	0.60
1:B:427:GLU:CD	1:B:427:GLU:N	2.53	0.60
1:B:357:ARG:NH2	1:B:358:LYS:CE	2.65	0.59
1:A:288:ARG:HD2	1:A:288:ARG:O	2.01	0.59
1:B:278:ALA:CB	1:B:357:ARG:NH1	2.64	0.59
1:B:311:LEU:H	1:B:311:LEU:CD2	2.15	0.59
1:B:357:ARG:NH2	1:B:358:LYS:NZ	2.51	0.59
1:B:228:LEU:HD23	1:B:333:LEU:HD21	1.84	0.59
1:A:240:LYS:HG3	1:A:241:THR:N	2.18	0.58
1:A:270:LEU:HD23	1:A:270:LEU:C	2.22	0.58
1:A:325:ILE:HG12	1:A:388:ILE:HG23	1.84	0.58
1:A:433:ALA:O	1:A:437:GLN:HG3	2.03	0.58
1:B:357:ARG:HH21	1:B:358:LYS:HE3	1.68	0.58
1:B:259:GLU:O	1:B:263:LYS:HA	2.04	0.57
1:B:330:LEU:HD13	2:B:2:PTG:H4	1.86	0.57
1:A:288:ARG:HG3	2:A:1:PTG:O12	2.04	0.57
1:B:357:ARG:HD3	1:B:357:ARG:C	2.24	0.57
1:A:212:ARG:O	1:A:216:LYS:HG2	2.04	0.56
1:A:368:PHE:O	1:A:372:VAL:HG23	2.05	0.56
1:B:252:MET:SD	1:B:277:VAL:HG11	2.46	0.56
1:B:335:ASN:ND2	1:B:338:GLY:H	2.03	0.56
1:A:272:GLU:CG	1:A:273:GLN:H	2.09	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:LYS:HG3	3:B:1090:HOH:O	2.06	0.55
1:A:323:HIS:CE1	1:A:473:TYR:OH	2.60	0.55
1:A:268:THR:CG2	1:A:280:ARG:HH11	2.18	0.55
1:B:266:HIS:O	1:B:268:THR:HG22	2.07	0.55
1:B:278:ALA:HB3	1:B:357:ARG:HH11	1.68	0.55
1:B:288:ARG:HG3	2:B:2:PTG:H2A	1.89	0.54
1:A:275:LYS:O	1:A:275:LYS:HD3	2.07	0.54
1:B:278:ALA:CB	1:B:357:ARG:HH11	2.20	0.54
1:A:325:ILE:HD11	1:A:392:ILE:CG1	2.38	0.54
1:B:269:PRO:O	1:B:270:LEU:O	2.26	0.54
1:A:252:MET:CE	1:A:252:MET:HA	2.39	0.53
1:B:262:ILE:O	1:B:262:ILE:HG22	2.07	0.53
1:A:436:LEU:O	1:A:439:MET:HB2	2.09	0.53
1:A:271:GLN:O	1:A:280:ARG:NH2	2.42	0.53
1:A:437:GLN:O	1:A:440:THR:HG23	2.09	0.52
1:B:349:THR:HG23	3:B:1068:HOH:O	2.08	0.52
1:B:270:LEU:HD22	1:B:469:LEU:HD22	1.91	0.52
1:A:263:LYS:NZ	1:A:265:LYS:HE2	2.25	0.52
1:B:402:ASN:O	1:B:405:PRO:HD2	2.09	0.52
1:B:427:GLU:N	1:B:427:GLU:OE2	2.38	0.52
1:A:252:MET:CE	1:A:255:LEU:HD12	2.40	0.52
1:B:349:THR:HG21	3:B:1081:HOH:O	2.09	0.52
1:B:335:ASN:C	1:B:335:ASN:ND2	2.59	0.51
1:B:269:PRO:O	1:B:270:LEU:C	2.48	0.51
1:A:354:LYS:HA	1:A:361:GLY:O	2.10	0.51
1:B:257:MET:CG	1:B:261:LYS:HG3	2.41	0.51
1:B:342:SER:H	2:B:2:PTG:H2O	1.75	0.51
1:A:272:GLU:CG	1:A:273:GLN:N	2.71	0.50
1:B:264:PHE:HB2	1:B:267:ILE:HG12	1.93	0.50
1:A:263:LYS:HD2	1:A:265:LYS:HG2	1.92	0.50
1:A:443:ARG:HG3	1:A:443:ARG:HH11	1.75	0.50
1:B:255:LEU:O	1:B:259:GLU:HG3	2.11	0.50
1:B:329:MET:HB2	3:B:1109:HOH:O	2.12	0.50
1:B:357:ARG:NH2	1:B:358:LYS:CD	2.75	0.50
1:B:466:HIS:CD2	1:B:467:PRO:HD2	2.47	0.50
1:A:443:ARG:NH1	1:B:440:THR:CG2	2.75	0.50
1:A:277:VAL:O	1:A:281:ILE:HG12	2.11	0.50
1:A:443:ARG:NH1	1:B:440:THR:HG21	2.28	0.49
1:A:446:VAL:O	1:A:450:VAL:HG23	2.12	0.49
1:B:404:LYS:HB3	1:B:405:PRO:HD3	1.95	0.49
1:A:277:VAL:HG13	1:A:278:ALA:N	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:473:TYR:HA	1:A:476:LEU:HD12	1.95	0.49
1:A:353:LEU:O	1:A:356:LEU:HG	2.12	0.48
1:A:258:GLY:O	1:A:262:ILE:N	2.42	0.48
1:A:264:PHE:O	1:A:266:HIS:N	2.46	0.48
1:B:276:GLU:OE1	1:B:357:ARG:NH2	2.47	0.48
1:B:425:HIS:N	1:B:426:PRO:HD3	2.30	0.47
1:B:466:HIS:CD2	1:B:468:LEU:H	2.29	0.47
1:B:247:PHE:HB2	1:B:262:ILE:HD12	1.95	0.47
1:B:268:THR:OG1	1:B:270:LEU:CG	2.60	0.47
1:B:357:ARG:NH2	1:B:358:LYS:HZ2	2.12	0.47
1:A:272:GLU:HG2	1:A:273:GLN:N	2.18	0.47
1:A:394:SER:O	1:A:397:ARG:HG2	2.15	0.47
1:B:335:ASN:ND2	1:B:337:ASP:N	2.58	0.47
1:A:263:LYS:HE3	1:A:263:LYS:HB3	1.62	0.46
1:B:358:LYS:CB	1:B:359:PRO:HD3	2.39	0.46
1:B:342:SER:H	2:B:2:PTG:C21	2.28	0.46
1:A:289:SER:OG	2:A:1:PTG:H19	2.16	0.46
1:A:443:ARG:HG3	1:A:443:ARG:NH1	2.30	0.46
1:B:357:ARG:HG2	1:B:358:LYS:N	2.30	0.46
1:B:249:ILE:HD11	1:B:264:PHE:CE1	2.51	0.46
1:B:357:ARG:HD2	3:B:1078:HOH:O	2.15	0.46
1:A:252:MET:HA	1:A:252:MET:HE3	1.97	0.46
1:A:463:MET:HE3	1:A:465:LEU:HD21	1.98	0.46
1:A:203:GLN:OE1	1:A:203:GLN:O	2.33	0.46
1:B:274:SER:CB	1:B:280:ARG:HD3	2.46	0.45
1:B:276:GLU:OE2	1:B:276:GLU:C	2.54	0.45
1:B:362:ASP:CG	1:B:362:ASP:O	2.55	0.45
1:B:270:LEU:HB3	1:B:273:GLN:OE1	2.16	0.45
1:B:330:LEU:HD21	1:B:364:MET:HE1	1.98	0.45
1:B:269:PRO:C	1:B:270:LEU:O	2.55	0.45
1:B:265:LYS:O	1:B:266:HIS:CG	2.70	0.45
1:B:456:ILE:HG22	1:B:456:ILE:O	2.17	0.45
1:A:280:ARG:HD3	1:A:280:ARG:HA	1.89	0.45
1:B:330:LEU:HD22	2:B:2:PTG:H4A	1.98	0.45
1:B:342:SER:H	2:B:2:PTG:H21A	1.83	0.44
1:A:317:LEU:HD22	1:A:392:ILE:O	2.17	0.44
1:B:277:VAL:HG12	1:B:278:ALA:N	2.33	0.44
1:B:402:ASN:OD1	1:B:405:PRO:HD3	2.18	0.44
1:B:207:GLU:C	1:B:209:ALA:H	2.20	0.44
1:B:365:GLU:HB3	1:B:366:PRO:HD3	1.99	0.44
1:B:217:HIS:HE1	1:B:302:SER:O	2.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:403:VAL:HG12	1:B:407:GLU:HG3	2.00	0.44
1:A:358:LYS:CB	1:A:359:PRO:CD	2.96	0.44
1:B:343:GLU:H	2:B:2:PTG:C22	2.31	0.44
1:B:270:LEU:CB	1:B:273:GLN:OE1	2.66	0.43
1:B:357:ARG:CG	1:B:358:LYS:N	2.81	0.43
1:A:255:LEU:HD11	1:A:277:VAL:HG23	2.00	0.43
1:B:411:ASP:O	1:B:415:GLN:HG3	2.19	0.43
1:A:271:GLN:HB2	1:A:280:ARG:HH22	1.84	0.43
1:A:261:LYS:N	1:A:261:LYS:HD2	2.33	0.43
1:A:404:LYS:N	1:A:405:PRO:HD2	2.33	0.43
1:A:234:ARG:NH1	3:A:1009:HOH:O	2.51	0.43
1:B:444:GLN:O	1:B:448:GLU:HB2	2.18	0.43
1:B:425:HIS:HB3	1:B:428:SER:OG	2.18	0.43
1:A:360:PHE:O	1:A:363:PHE:HD2	2.02	0.43
1:A:215:ALA:CB	1:A:386:ILE:HD11	2.48	0.42
1:A:384:LEU:O	1:A:388:ILE:HG13	2.18	0.42
1:B:224:LYS:O	1:B:224:LYS:HG2	2.18	0.42
1:B:393:LEU:HD22	1:B:393:LEU:N	2.34	0.42
1:B:342:SER:H	2:B:2:PTG:C20	2.32	0.42
1:B:257:MET:HG3	1:B:261:LYS:HG3	2.01	0.42
1:A:290:VAL:HG13	1:A:468:LEU:HD23	2.01	0.42
1:A:325:ILE:HD11	1:A:392:ILE:HG13	2.01	0.42
1:A:356:LEU:HD23	1:A:356:LEU:HA	1.81	0.42
1:B:397:ARG:HB2	1:B:400:LEU:HD11	2.00	0.42
1:A:305:GLY:HA2	1:A:308:ASN:HD22	1.85	0.42
1:B:341:ILE:HA	2:B:2:PTG:H21A	2.01	0.42
1:A:283:GLN:O	1:A:286:GLN:HB2	2.20	0.42
1:B:264:PHE:HD1	1:B:267:ILE:HG13	1.85	0.42
1:B:424:ASN:ND2	3:B:1086:HOH:O	2.53	0.42
1:A:265:LYS:H	1:A:265:LYS:HG3	1.49	0.42
1:B:472:ILE:HB	1:B:473:TYR:CD2	2.55	0.42
1:B:311:LEU:N	1:B:311:LEU:CD2	2.79	0.41
1:A:204:LEU:HB3	1:A:205:ASN:H	1.38	0.41
1:A:271:GLN:CB	1:A:280:ARG:NH2	2.83	0.41
1:B:309:LEU:CD2	1:B:405:PRO:HB2	2.50	0.41
1:B:257:MET:HG2	1:B:261:LYS:HG3	2.01	0.41
1:B:282:PHE:HZ	1:B:453:LEU:HD23	1.84	0.41
1:B:245:SER:HA	1:B:246:PRO:HD3	1.92	0.41
1:B:288:ARG:HD2	1:B:288:ARG:O	2.20	0.41
1:A:356:LEU:O	1:A:357:ARG:HB2	2.20	0.41
1:B:207:GLU:C	1:B:209:ALA:N	2.72	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:LYS:O	1:A:265:LYS:HG3	2.20	0.41
1:A:268:THR:HG22	1:A:280:ARG:HH11	1.83	0.41
1:B:330:LEU:O	1:B:334:MET:HG3	2.21	0.41
1:A:325:ILE:HD11	1:A:392:ILE:HG12	2.03	0.41
1:B:263:LYS:O	1:B:264:PHE:C	2.59	0.41
1:B:241:THR:O	1:B:242:THR:C	2.59	0.41
1:B:349:THR:HG23	1:B:351:GLU:H	1.86	0.41
1:B:356:LEU:O	1:B:361:GLY:HA3	2.21	0.41
1:B:273:GLN:CD	1:B:273:GLN:N	2.74	0.40
1:B:358:LYS:HB3	1:B:359:PRO:HD2	1.93	0.40
1:B:466:HIS:HA	1:B:467:PRO:HD3	1.85	0.40
1:A:270:LEU:O	1:A:270:LEU:HD23	2.21	0.40
1:B:322:VAL:O	1:B:326:ILE:HG13	2.21	0.40
1:A:267:ILE:HG22	1:A:284:GLY:O	2.22	0.40
1:A:290:VAL:HG21	1:A:466:HIS:CD2	2.57	0.40
1:A:424:ASN:HB3	1:A:425:HIS:ND1	2.36	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	272/286 (95%)	251 (92%)	11 (4%)	10 (4%)	3 0
1	B	259/286 (91%)	232 (90%)	17 (7%)	10 (4%)	3 0
All	All	531/572 (93%)	483 (91%)	28 (5%)	20 (4%)	3 0

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	265	LYS
1	A	270	LEU

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Mol	Chain	Res	Type
1	A	272	GLU
1	A	358	LYS
1	B	243	ASP
1	B	270	LEU
1	B	274	SER
1	B	358	LYS
1	A	271	GLN
1	A	274	SER
1	B	240	LYS
1	B	265	LYS
1	B	269	PRO
1	B	276	GLU
1	B	460	GLU
1	A	264	PHE
1	A	266	HIS
1	A	357	ARG
1	A	361	GLY
1	B	268	THR

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	247/257 (96%)	234 (95%)	13 (5%)	22	13
1	B	236/257 (92%)	224 (95%)	12 (5%)	24	14
All	All	483/514 (94%)	458 (95%)	25 (5%)	23	14

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	203	GLN
1	A	252	MET
1	A	255	LEU
1	A	257	MET
1	A	261	LYS

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Mol	Chain	Res	Type
1	A	268	THR
1	A	273	GLN
1	A	283	GLN
1	A	311	LEU
1	A	358	LYS
1	A	427	GLU
1	A	440	THR
1	A	469	LEU
1	B	241	THR
1	B	242	THR
1	B	261	LYS
1	B	268	THR
1	B	273	GLN
1	B	276	GLU
1	B	335	ASN
1	B	349	THR
1	B	357	ARG
1	B	424	ASN
1	B	441	ASP
1	B	468	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	203	GLN
1	A	283	GLN
1	A	308	ASN
1	A	314	GLN
1	A	323	HIS
1	A	425	HIS
1	A	430	GLN
1	A	449	HIS
1	A	470	GLN
1	B	217	HIS
1	B	253	ASN
1	B	283	GLN
1	B	308	ASN
1	B	335	ASN
1	B	424	ASN
1	B	430	GLN
1	B	454	GLN
1	B	466	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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
2	PTG	A	1	-	19,23,23	1.55	4 (21%)	19,27,27	1.35	4 (21%)
2	PTG	B	2	-	19,23,23	1.46	4 (21%)	19,27,27	1.32	4 (21%)

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	PTG	A	1	-	-	4/17/30/30	0/1/1/1
2	PTG	B	2	-	-	7/17/30/30	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	PTG	C15-C10	-3.96	1.46	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	PTG	C15-C10	-3.77	1.47	1.54
2	A	1	PTG	C7-C8	3.47	1.54	1.43
2	B	2	PTG	C7-C8	2.89	1.52	1.43
2	A	1	PTG	O12-C11	2.69	1.31	1.24
2	B	2	PTG	C16-C17	2.46	1.57	1.50
2	B	2	PTG	O12-C11	2.42	1.31	1.24
2	A	1	PTG	C16-C17	2.33	1.57	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	PTG	O24-C22-C21	2.55	122.24	114.03
2	A	1	PTG	C8-C7-C6	2.41	129.27	123.63
2	B	2	PTG	C21-C20-C19	-2.40	108.95	113.23
2	B	2	PTG	O24-C22-C21	2.38	121.67	114.03
2	A	1	PTG	O23-C22-C21	-2.35	115.55	123.08
2	A	1	PTG	C15-C14-C13	2.32	113.89	111.56
2	B	2	PTG	O23-C22-C21	-2.17	116.11	123.08
2	B	2	PTG	C15-C10-C11	2.12	110.12	107.05

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	PTG	C2-C3-C4-C5
2	B	2	PTG	C2-C3-C4-C5
2	A	1	PTG	C1-C2-C3-C4
2	B	2	PTG	C1-C2-C3-C4
2	B	2	PTG	C18-C19-C20-C21
2	B	2	PTG	C14-C15-C16-C17
2	A	1	PTG	C20-C21-C22-O23
2	A	1	PTG	C20-C21-C22-O24
2	B	2	PTG	C15-C16-C17-C18
2	B	2	PTG	C17-C18-C19-C20
2	B	2	PTG	C3-C4-C5-C6

There are no ring outliers.

2 monomers are involved in 12 short contacts:

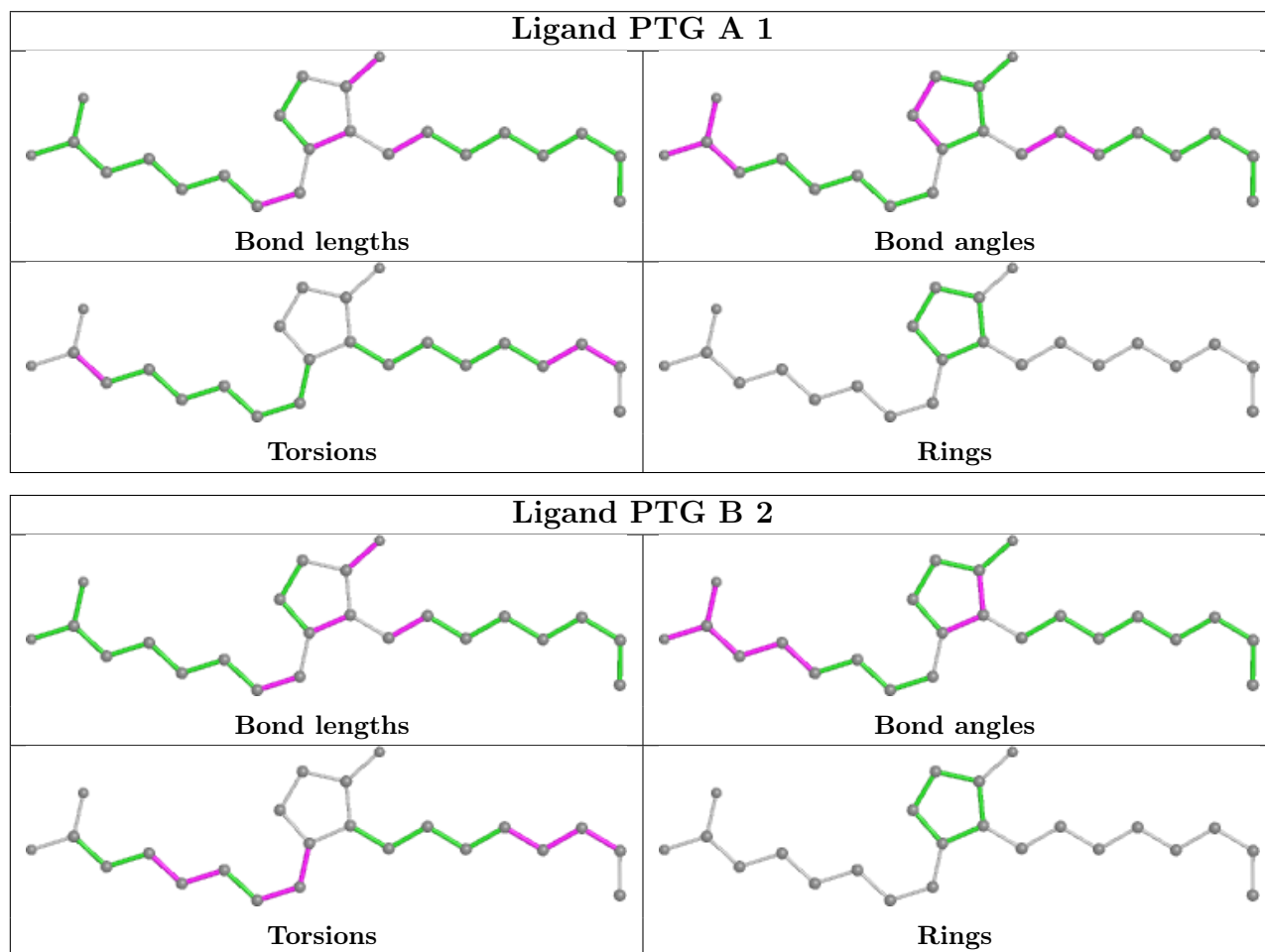
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	PTG	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2	PTG	10	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

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	274/286 (95%)	0.81	35 (12%) 3 4	26, 38, 64, 68	0
1	B	263/286 (91%)	0.76	34 (12%) 3 3	24, 38, 61, 68	0
All	All	537/572 (93%)	0.79	69 (12%) 3 4	24, 38, 63, 68	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	270	LEU	10.2
1	A	267	ILE	10.2
1	A	269	PRO	9.8
1	A	266	HIS	8.4
1	A	268	THR	7.8
1	B	268	THR	7.4
1	A	265	LYS	7.3
1	B	241	THR	7.2
1	B	242	THR	6.8
1	B	243	ASP	5.8
1	A	203	GLN	5.3
1	A	261	LYS	5.3
1	B	267	ILE	4.9
1	B	458	LYS	4.6
1	A	204	LEU	4.6
1	A	239	GLY	4.3
1	B	274	SER	4.2
1	B	261	LYS	4.2
1	A	262	ILE	4.1
1	A	264	PHE	4.1
1	B	265	LYS	4.0
1	B	273	GLN	4.0
1	A	271	GLN	3.9
1	B	272	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	260	ASP	3.8
1	B	240	LYS	3.7
1	B	341	ILE	3.7
1	A	275	LYS	3.7
1	A	272	GLU	3.6
1	B	275	LYS	3.6
1	A	356	LEU	3.6
1	B	459	THR	3.6
1	A	475	ASP	3.4
1	B	209	ALA	3.3
1	B	271	GLN	3.3
1	B	454	GLN	3.3
1	B	262	ILE	3.2
1	A	263	LYS	3.0
1	B	207	GLU	2.9
1	A	256	MET	2.8
1	A	273	GLN	2.8
1	A	453	LEU	2.7
1	B	436	LEU	2.7
1	A	252	MET	2.7
1	B	357	ARG	2.7
1	A	426	PRO	2.6
1	A	358	LYS	2.5
1	B	439	MET	2.4
1	A	362	ASP	2.4
1	A	287	PHE	2.4
1	B	456	ILE	2.3
1	A	253	ASN	2.3
1	B	387	PHE	2.3
1	B	470	GLN	2.3
1	B	469	LEU	2.3
1	B	256	MET	2.2
1	B	266	HIS	2.2
1	B	435	LEU	2.2
1	B	269	PRO	2.1
1	A	280	ARG	2.1
1	A	207	GLU	2.1
1	A	439	MET	2.1
1	B	212	ARG	2.0
1	A	388	ILE	2.0
1	A	277	VAL	2.0
1	A	427	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	427	GLU	2.0
1	B	340	LEU	2.0
1	A	241	THR	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](#)

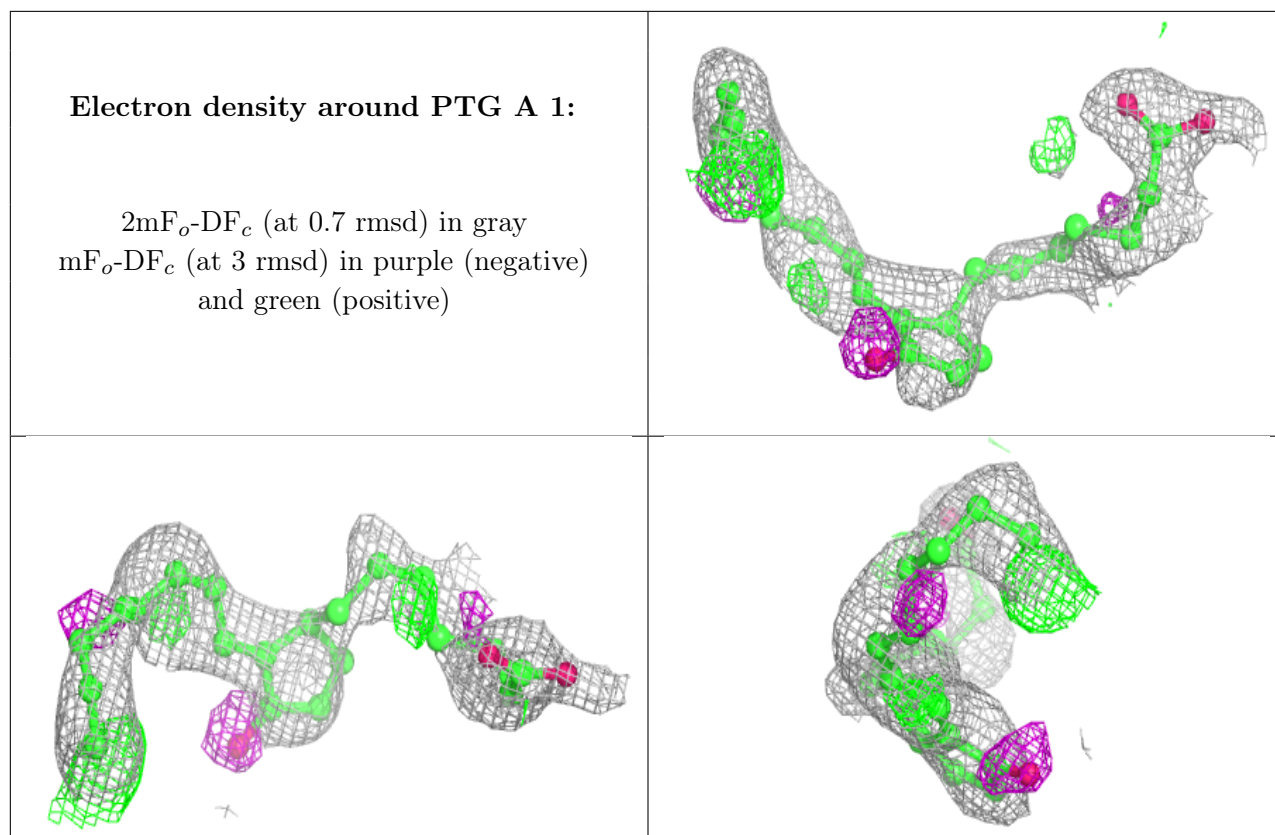
There are no monosaccharides in this entry.

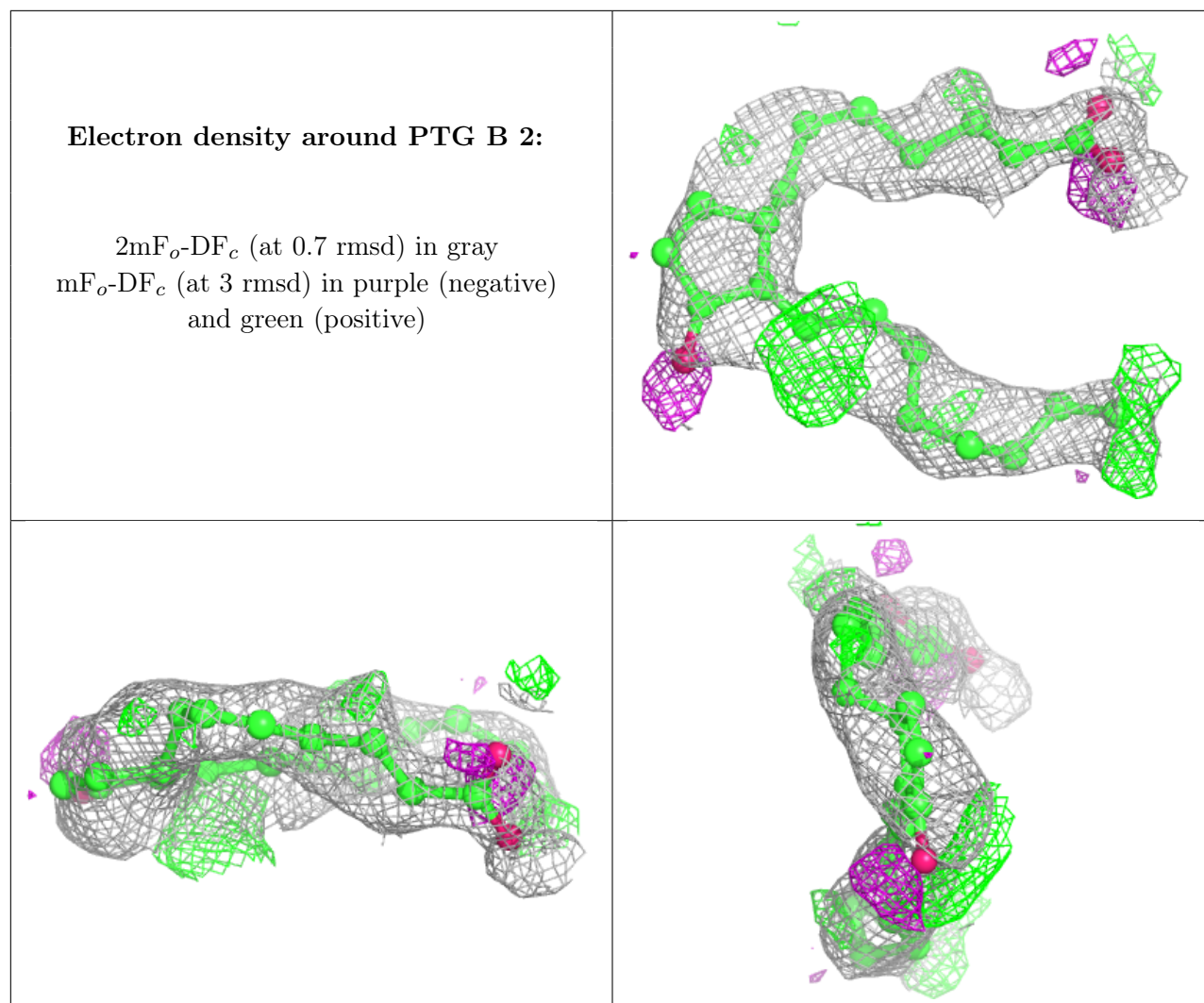
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
2	PTG	A	1	23/23	0.52	0.35	53,60,61,63	0
2	PTG	B	2	23/23	0.59	0.37	49,58,62,64	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.