



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 11:29 pm BST

PDB ID : 3W5C
Title : Crystal structure of the calcium pump in the E2 state free from exogenous inhibitors
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Deposited on : 2013-01-27
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 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 : 2.11
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) : 2.11

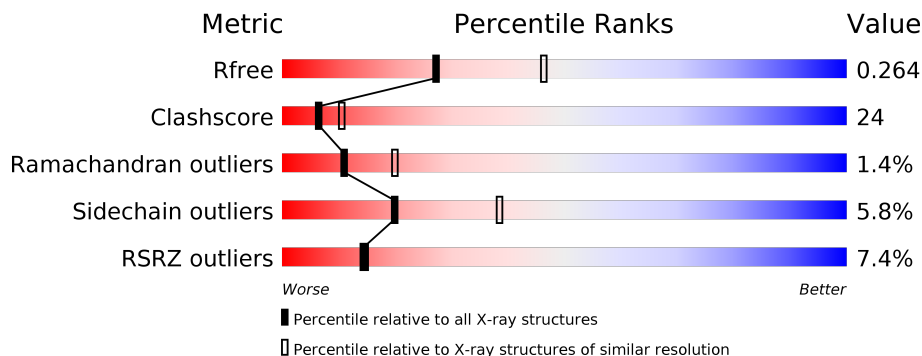
1 Overall quality at a glance i

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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (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 on 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	995	

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	PTY	A	1002	-	-	-	X
3	PTY	A	1004	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7782 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 SERCA1a.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	995	7674	4878	1287	1452	57	0	0	0

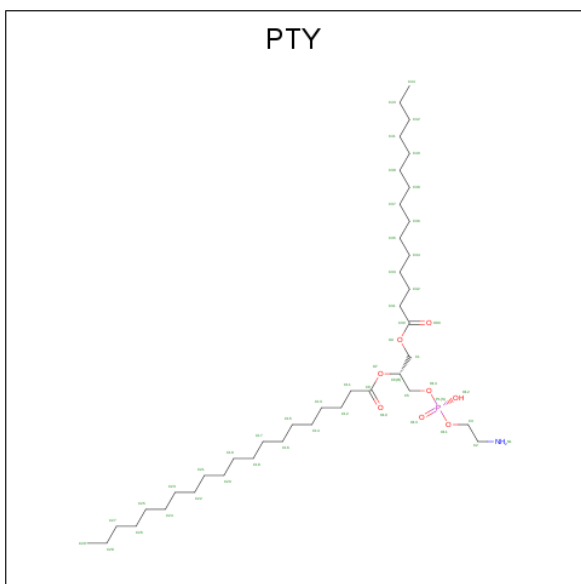
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ACE	-	ACETYLATION	UNP B6CAM1

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is PHOSPHATIDYLETHANOLAMINE (three-letter code: PTY) (formula: C₄₀H₈₀NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			19	9	1	8	1		
3	A	1	Total	C	N	O	P	0	0
			19	9	1	8	1		
3	A	1	Total	C	N	O	P	0	0
			19	9	1	8	1		
3	A	1	Total	C	N	O	P	0	0
			19	9	1	8	1		
3	A	1	Total	C	N	O	P	0	0
			19	9	1	8	1		

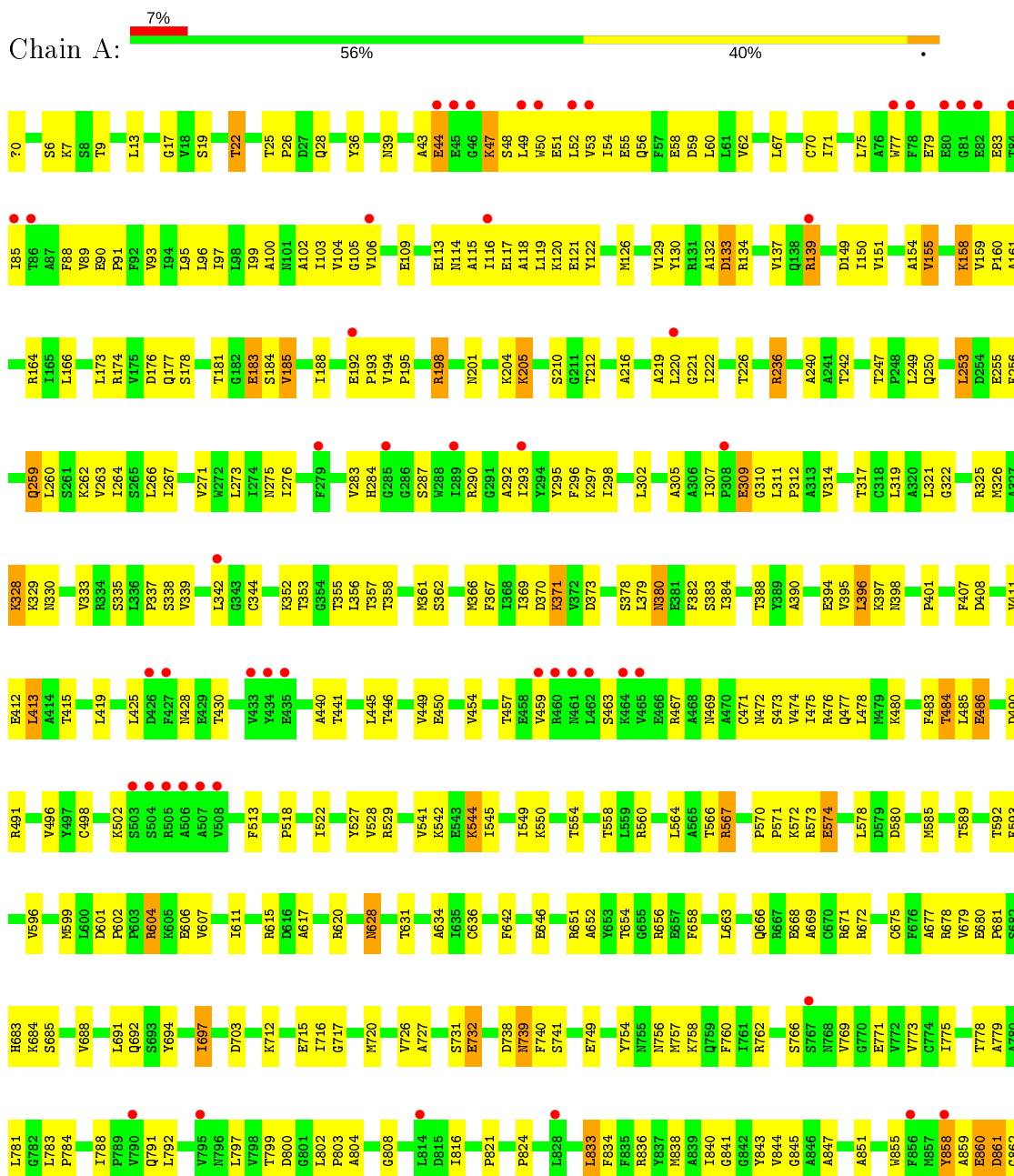
- Molecule 4 is water.

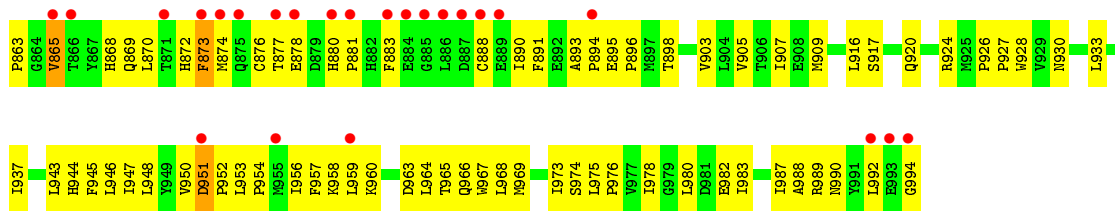
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: SERCA1a





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	71.23Å 71.23Å 586.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 48.07 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.7 (15.00-2.50) 98.7 (48.07-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.95 (at 2.51Å)	Xtrriage
Refinement program	CNS, REFMAC 5.6.0117	Depositor
R, R_{free}	0.270 , 0.276 0.263 , 0.264	Depositor DCC
R_{free} test set	2715 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	58.2	Xtrriage
Anisotropy	0.082	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7782	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.76% 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: NA, PTY, ACE

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	0.42	1/7813 (0.0%)	0.63	1/10594 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	994	GLY	C-OXT	-7.44	1.09	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	413	LEU	CA-CB-CG	5.89	128.85	115.30

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	7674	0	7765	372	0
2	A	1	0	0	1	0
3	A	95	0	55	8	0
4	A	12	0	0	0	0
All	All	7782	0	7820	372	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 24.

All (372) 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:990:ASN:ND2	3:A:1003:PTY:HC31	1.66	1.09
1:A:990:ASN:HD21	3:A:1003:PTY:HC31	0.86	1.01
1:A:990:ASN:HD21	3:A:1003:PTY:C3	1.73	1.00
1:A:484:THR:HB	1:A:496:VAL:HG12	1.53	0.91
1:A:328:LYS:HA	1:A:328:LYS:HE2	1.54	0.88
1:A:739:ASN:HD22	1:A:740:PHE:N	1.73	0.85
1:A:679:VAL:HG13	1:A:683:HIS:HB2	1.59	0.84
1:A:205:LYS:HB3	1:A:205:LYS:NZ	1.93	0.83
1:A:628:ASN:C	1:A:628:ASN:HD22	1.83	0.82
1:A:604:ARG:HH11	1:A:604:ARG:HG3	1.44	0.81
1:A:865:VAL:HB	1:A:868:HIS:HB2	1.62	0.80
1:A:791:GLN:HE21	1:A:959:LEU:HD21	1.45	0.80
1:A:663:LEU:HD12	1:A:663:LEU:H	1.45	0.80
1:A:628:ASN:HD21	1:A:631:THR:H	1.30	0.79
1:A:305:ALA:HB1	1:A:771:GLU:HB3	1.64	0.79
1:A:441:THR:HG23	1:A:599:MET:SD	2.22	0.79
1:A:989:ARG:NH2	3:A:1003:PTY:HC51	2.00	0.77
1:A:441:THR:HG21	1:A:560:ARG:NH1	2.00	0.76
1:A:873:PHE:CE2	1:A:876:CYS:HA	2.19	0.76
1:A:963:ASP:HB3	1:A:966:GLN:NE2	2.00	0.75
1:A:873:PHE:HE2	1:A:876:CYS:HA	1.52	0.74
1:A:775:ILE:O	1:A:778:THR:HG22	1.88	0.74
1:A:326:MET:HG2	1:A:749:GLU:HG2	1.69	0.73
1:A:369:ILE:HG13	1:A:528:VAL:CG1	2.18	0.73
1:A:379:LEU:HG	1:A:544:LYS:NZ	2.04	0.73
1:A:116:ILE:HD11	1:A:240:ALA:HB2	1.71	0.72
1:A:52:LEU:HD11	1:A:109:GLU:HG3	1.69	0.72
1:A:720:MET:HB3	1:A:738:ASP:OD2	1.90	0.72
1:A:395:VAL:O	1:A:396:LEU:HD23	1.88	0.72
1:A:413:LEU:HD22	1:A:564:LEU:HD12	1.71	0.72
1:A:865:VAL:HB	1:A:868:HIS:CB	2.19	0.72
1:A:472:ASN:HB3	1:A:476:ARG:HH12	1.54	0.72
1:A:326:MET:HE1	1:A:339:VAL:HG22	1.71	0.72
1:A:963:ASP:HB3	1:A:966:GLN:CD	2.09	0.72
1:A:370:ASP:HB3	1:A:378:SER:OG	1.91	0.71
1:A:44:GLU:HB3	1:A:117:GLU:OE2	1.91	0.71
1:A:305:ALA:HB2	1:A:792:LEU:HD13	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:678:ARG:HH11	1:A:678:ARG:HG3	1.55	0.70
1:A:271:VAL:HG11	1:A:298:ILE:CD1	2.22	0.70
1:A:472:ASN:HB3	1:A:476:ARG:NH1	2.05	0.70
1:A:176:ASP:OD1	1:A:178:SER:HB2	1.92	0.70
1:A:680:GLU:HB3	1:A:681:PRO:HD2	1.73	0.69
1:A:192:GLU:CD	1:A:193:PRO:HD2	2.12	0.69
1:A:963:ASP:OD2	1:A:965:THR:HB	1.91	0.69
1:A:126:MET:HG3	1:A:139:ARG:NH1	2.08	0.69
1:A:567:ARG:HD2	1:A:570:PRO:HA	1.75	0.69
1:A:791:GLN:NE2	1:A:959:LEU:HD21	2.08	0.69
1:A:295:TYR:O	1:A:298:ILE:HG12	1.92	0.69
1:A:784:PRO:HG2	1:A:870:LEU:HD22	1.75	0.68
1:A:283:VAL:HG13	1:A:284:HIS:ND1	2.07	0.68
1:A:59:ASP:HB3	1:A:62:VAL:HG22	1.73	0.68
1:A:271:VAL:HG11	1:A:298:ILE:HD11	1.75	0.67
1:A:604:ARG:HH11	1:A:604:ARG:CG	2.08	0.67
1:A:335:SER:HB3	1:A:338:SER:OG	1.95	0.66
1:A:762:ARG:NH1	1:A:833:LEU:HD21	2.09	0.66
1:A:19:SER:HB3	1:A:22:THR:CG2	2.24	0.66
1:A:843:TYR:OH	1:A:976:PRO:HG2	1.96	0.66
1:A:585:MET:O	1:A:589:THR:HG23	1.95	0.66
1:A:473:SER:O	1:A:477:GLN:HG2	1.95	0.66
1:A:654:THR:HA	1:A:677:ALA:O	1.96	0.66
1:A:628:ASN:ND2	1:A:631:THR:H	1.92	0.66
1:A:784:PRO:CG	1:A:870:LEU:HD22	2.26	0.65
1:A:739:ASN:ND2	1:A:741:SER:H	1.95	0.65
1:A:860:GLU:H	1:A:860:GLU:CD	2.00	0.65
1:A:739:ASN:C	1:A:739:ASN:HD22	1.98	0.64
1:A:52:LEU:HG	1:A:106:VAL:HG22	1.79	0.64
1:A:668:GLU:O	1:A:671:ARG:HG2	1.97	0.64
1:A:361:MET:HG2	1:A:440:ALA:HB1	1.78	0.64
1:A:55:GLU:HA	1:A:58:GLU:HG2	1.79	0.63
1:A:679:VAL:CG1	1:A:683:HIS:HB2	2.28	0.63
1:A:161:ALA:HA	1:A:210:SER:HB2	1.79	0.63
1:A:311:LEU:N	1:A:312:PRO:HD2	2.14	0.62
1:A:369:ILE:HG13	1:A:528:VAL:HG13	1.79	0.62
1:A:518:PRO:HB3	1:A:549:ILE:HD13	1.80	0.62
1:A:567:ARG:CZ	1:A:571:PRO:HD3	2.30	0.62
1:A:974:SER:C	1:A:976:PRO:HD2	2.20	0.62
1:A:669:ALA:HA	1:A:672:ARG:NH1	2.14	0.61
1:A:950:VAL:O	1:A:954:PRO:HD2	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:684:LYS:O	1:A:688:VAL:HG23	2.01	0.61
1:A:205:LYS:HB3	1:A:205:LYS:HZ3	1.64	0.61
1:A:860:GLU:N	1:A:860:GLU:CD	2.55	0.60
1:A:527:TYR:O	1:A:592:THR:HA	2.01	0.60
1:A:952:PRO:O	1:A:956:ILE:HG13	2.01	0.60
1:A:205:LYS:HB3	1:A:205:LYS:HZ2	1.66	0.60
1:A:50:TRP:O	1:A:54:ILE:HG12	2.02	0.60
1:A:379:LEU:HG	1:A:544:LYS:HZ2	1.66	0.59
1:A:394:GLU:HG3	1:A:396:LEU:HD21	1.84	0.59
1:A:851:ALA:HB2	1:A:903:VAL:HG21	1.83	0.59
1:A:198:ARG:HD3	1:A:198:ARG:O	2.02	0.59
1:A:290:ARG:HH11	1:A:290:ARG:HB3	1.65	0.59
1:A:310:GLY:O	1:A:314:VAL:HG23	2.01	0.59
1:A:67:LEU:O	1:A:70:CYS:HB2	2.02	0.59
1:A:870:LEU:HD12	1:A:870:LEU:H	1.67	0.59
1:A:732:GLU:OE1	2:A:1001:NA:NA	1.75	0.59
1:A:19:SER:HB3	1:A:22:THR:HG23	1.83	0.59
1:A:964:LEU:O	1:A:968:LEU:HB2	2.03	0.59
1:A:836:ARG:O	1:A:840:ILE:HG12	2.02	0.59
1:A:273:LEU:HD23	1:A:276:ILE:HD11	1.84	0.58
1:A:567:ARG:CD	1:A:570:PRO:HA	2.32	0.58
1:A:572:LYS:HB3	1:A:574:GLU:OE2	2.02	0.58
1:A:573:ARG:HG2	1:A:573:ARG:HH11	1.68	0.58
1:A:855:TRP:HA	1:A:859:ALA:CB	2.33	0.58
1:A:412:GLU:OE1	1:A:529:ARG:HD2	2.03	0.58
1:A:419:LEU:HD12	1:A:513:PHE:CE2	2.39	0.58
1:A:679:VAL:HG13	1:A:683:HIS:CB	2.32	0.58
1:A:844:VAL:HG22	1:A:907:ILE:HG21	1.85	0.58
1:A:703:ASP:HA	1:A:727:ALA:CB	2.34	0.58
1:A:628:ASN:ND2	1:A:628:ASN:C	2.57	0.57
1:A:678:ARG:NH1	1:A:678:ARG:HG3	2.19	0.57
1:A:6:SER:HA	1:A:194:VAL:O	2.04	0.57
1:A:860:GLU:OE2	1:A:860:GLU:N	2.37	0.57
1:A:989:ARG:HH22	3:A:1003:PTY:HC51	1.69	0.57
1:A:457:THR:O	1:A:459:VAL:HG13	2.04	0.57
1:A:174:ARG:CZ	1:A:188:ILE:HD11	2.34	0.57
1:A:800:ASP:C	1:A:803:PRO:HD2	2.25	0.57
1:A:697:ILE:HD11	1:A:824:PRO:HG2	1.86	0.56
1:A:259:GLN:O	1:A:263:VAL:HG23	2.05	0.56
1:A:449:VAL:HG21	1:A:472:ASN:OD1	2.04	0.56
1:A:287:SER:HB2	1:A:290:ARG:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:ILE:HB	1:A:309:GLU:OE2	2.06	0.56
1:A:119:LEU:HD13	1:A:236:ARG:HB2	1.87	0.56
1:A:352:LYS:HA	1:A:356:LEU:HD12	1.88	0.56
1:A:236:ARG:HD3	1:A:236:ARG:C	2.25	0.56
1:A:408:ASP:HA	1:A:411:VAL:HG23	1.87	0.56
1:A:943:LEU:O	1:A:946:LEU:HB3	2.05	0.56
1:A:48:SER:OG	1:A:51:GLU:HG3	2.06	0.55
1:A:604:ARG:NH1	1:A:604:ARG:CG	2.67	0.55
1:A:322:GLY:O	1:A:325:ARG:HB3	2.06	0.55
1:A:758:LYS:O	1:A:762:ARG:HG3	2.06	0.55
1:A:71:ILE:HG21	1:A:296:PHE:HB3	1.89	0.55
1:A:60:LEU:O	1:A:60:LEU:HD23	2.06	0.55
1:A:905:VAL:O	1:A:909:MET:HG2	2.07	0.55
1:A:541:VAL:O	1:A:545:ILE:HG13	2.06	0.55
1:A:858:TYR:N	1:A:858:TYR:CD1	2.75	0.54
1:A:259:GLN:HA	1:A:259:GLN:HE21	1.72	0.54
1:A:975:LEU:N	1:A:976:PRO:HD2	2.22	0.54
1:A:56:GLN:OE1	1:A:105:GLY:HA3	2.07	0.54
1:A:950:VAL:HG12	1:A:952:PRO:HD2	1.89	0.54
1:A:847:ALA:O	1:A:973:ILE:HD11	2.07	0.54
1:A:697:ILE:HD11	1:A:824:PRO:CG	2.38	0.54
1:A:118:ALA:O	1:A:121:GLU:HB3	2.07	0.54
1:A:43:ALA:HA	1:A:120:LYS:NZ	2.23	0.54
1:A:174:ARG:HB2	1:A:216:ALA:HB3	1.90	0.54
1:A:895:GLU:N	1:A:896:PRO:HD2	2.23	0.53
1:A:256:PHE:CE2	1:A:260:LEU:HD22	2.44	0.53
1:A:415:THR:HA	1:A:475:ILE:HG21	1.90	0.53
1:A:668:GLU:OE2	1:A:671:ARG:HD3	2.09	0.53
1:A:428:ASN:OD1	1:A:430:THR:HB	2.09	0.53
1:A:329:LYS:O	1:A:330:ASN:HB2	2.08	0.53
1:A:425:LEU:HB2	1:A:469:ASN:OD1	2.08	0.53
1:A:314:VAL:HG11	1:A:804:ALA:HB1	1.91	0.53
1:A:155:VAL:O	1:A:155:VAL:HG13	2.08	0.53
1:A:99:ILE:O	1:A:103:ILE:HG13	2.08	0.53
1:A:358:THR:OG1	1:A:602:PRO:HG2	2.09	0.53
1:A:604:ARG:HB2	1:A:607:VAL:HG23	1.91	0.52
1:A:894:PRO:HG2	1:A:895:GLU:OE1	2.09	0.52
1:A:262:LYS:O	1:A:266:LEU:HD23	2.10	0.52
1:A:550:LYS:O	1:A:554:THR:HB	2.09	0.52
1:A:166:LEU:H	1:A:221:GLY:HA2	1.75	0.52
1:A:174:ARG:NH2	1:A:188:ILE:HD11	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:ILE:HG21	1:A:302:LEU:HD11	1.92	0.52
1:A:716:ILE:N	1:A:716:ILE:HD12	2.25	0.52
1:A:129:VAL:HG12	1:A:151:VAL:HG12	1.91	0.52
1:A:441:THR:HG21	1:A:560:ARG:HH12	1.73	0.52
1:A:293:ILE:O	1:A:297:LYS:HB2	2.10	0.52
1:A:611:ILE:O	1:A:615:ARG:HG3	2.09	0.51
1:A:71:ILE:HG22	1:A:75:LEU:HD13	1.90	0.51
1:A:754:TYR:HD1	1:A:757:MET:HE3	1.74	0.51
1:A:96:LEU:HD23	1:A:96:LEU:C	2.30	0.51
1:A:260:LEU:O	1:A:264:ILE:HG13	2.10	0.51
1:A:483:PHE:HE2	1:A:485:LEU:HD21	1.76	0.51
1:A:89:VAL:O	1:A:93:VAL:HG23	2.10	0.51
1:A:873:PHE:CE1	1:A:881:PRO:HG3	2.46	0.51
1:A:983:ILE:O	1:A:987:ILE:HG13	2.10	0.51
1:A:446:THR:O	1:A:449:VAL:HG22	2.10	0.51
1:A:917:SER:OG	1:A:920:GLN:HB2	2.11	0.50
1:A:396:LEU:HD22	1:A:401:PRO:HG3	1.93	0.50
1:A:880:HIS:N	1:A:881:PRO:HD2	2.26	0.50
1:A:944:HIS:O	1:A:947:ILE:HG12	2.12	0.50
1:A:958:LYS:O	1:A:959:LEU:HD23	2.11	0.50
1:A:255:GLU:O	1:A:259:GLN:HG2	2.10	0.50
1:A:471:CYS:O	1:A:475:ILE:HG13	2.12	0.50
1:A:55:GLU:HA	1:A:58:GLU:CG	2.42	0.50
1:A:309:GLU:HB3	1:A:797:LEU:HD11	1.92	0.50
1:A:893:ALA:O	1:A:896:PRO:HD2	2.12	0.50
1:A:788:ILE:HG12	1:A:791:GLN:OE1	2.12	0.50
1:A:878:GLU:HB3	1:A:880:HIS:CD2	2.48	0.49
1:A:384:ILE:HA	1:A:394:GLU:O	2.12	0.49
1:A:860:GLU:O	1:A:862:GLY:N	2.46	0.49
1:A:989:ARG:NH2	3:A:1003:PTY:O13	2.43	0.49
1:A:328:LYS:CE	1:A:328:LYS:HA	2.35	0.49
1:A:413:LEU:CD2	1:A:564:LEU:HD12	2.41	0.49
1:A:870:LEU:HB3	1:A:891:PHE:CE2	2.48	0.49
1:A:25:THR:HB	1:A:26:PRO:HD2	1.95	0.49
1:A:863:PRO:HB2	1:A:865:VAL:HG22	1.94	0.49
1:A:161:ALA:CA	1:A:210:SER:HB2	2.44	0.48
1:A:247:THR:H	1:A:250:GLN:NE2	2.11	0.48
1:A:450:GLU:OE2	1:A:450:GLU:HA	2.13	0.48
1:A:863:PRO:HD3	1:A:890:ILE:CD1	2.44	0.48
1:A:96:LEU:HD23	1:A:96:LEU:O	2.14	0.48
1:A:606:GLU:OE1	1:A:606:GLU:N	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:ALA:O	1:A:104:VAL:HG23	2.13	0.48
1:A:855:TRP:O	1:A:859:ALA:HB3	2.13	0.48
1:A:67:LEU:O	1:A:67:LEU:HD12	2.14	0.48
1:A:863:PRO:CG	1:A:890:ILE:HD13	2.43	0.48
1:A:916:LEU:HD21	1:A:930:ASN:CB	2.44	0.48
1:A:71:ILE:CG2	1:A:296:PHE:HB3	2.42	0.48
1:A:371:LYS:HE2	1:A:373:ASP:HB2	1.94	0.48
1:A:408:ASP:O	1:A:411:VAL:N	2.45	0.48
1:A:924:ARG:NE	1:A:924:ARG:HA	2.28	0.48
1:A:326:MET:CE	1:A:339:VAL:HG22	2.42	0.47
1:A:800:ASP:O	1:A:803:PRO:HD2	2.14	0.47
1:A:388:THR:CG2	1:A:390:ALA:HB3	2.44	0.47
1:A:122:TYR:O	1:A:158:LYS:HD3	2.14	0.47
1:A:878:GLU:HB3	1:A:880:HIS:NE2	2.29	0.47
1:A:361:MET:SD	1:A:601:ASP:HB2	2.54	0.47
1:A:757:MET:HA	1:A:760:PHE:CE2	2.50	0.47
1:A:951:ASP:O	1:A:952:PRO:C	2.50	0.47
1:A:988:ALA:HA	1:A:992:LEU:HD12	1.96	0.47
1:A:769:VAL:O	1:A:773:VAL:HG23	2.13	0.47
1:A:52:LEU:CD1	1:A:109:GLU:HG3	2.40	0.47
1:A:384:ILE:HD12	1:A:384:ILE:N	2.28	0.47
1:A:122:TYR:HE1	1:A:726:VAL:HG21	1.79	0.47
1:A:671:ARG:HD2	1:A:694:TYR:CZ	2.49	0.47
1:A:739:ASN:C	1:A:739:ASN:ND2	2.68	0.47
1:A:834:PHE:O	1:A:838:MET:HB2	2.15	0.47
1:A:56:GLN:HB3	1:A:102:ALA:HA	1.97	0.47
1:A:652:ALA:HA	1:A:675:CYS:O	2.15	0.47
1:A:259:GLN:CA	1:A:259:GLN:HE21	2.27	0.47
1:A:90:GLU:HB3	1:A:91:PRO:CD	2.45	0.47
1:A:0:ACE:H2	1:A:36:TYR:CE1	2.49	0.46
1:A:474:VAL:O	1:A:478:LEU:HD13	2.15	0.46
1:A:703:ASP:HA	1:A:727:ALA:HB1	1.96	0.46
1:A:395:VAL:C	1:A:396:LEU:HD23	2.35	0.46
1:A:342:LEU:HD13	1:A:716:ILE:HG21	1.97	0.46
1:A:802:LEU:HB2	1:A:803:PRO:HD3	1.97	0.46
1:A:397:LYS:O	1:A:398:ASN:HB2	2.15	0.46
1:A:39:ASN:OD1	1:A:226:THR:HB	2.15	0.46
1:A:183:GLU:O	1:A:185:VAL:N	2.49	0.46
1:A:290:ARG:HB3	1:A:290:ARG:NH1	2.30	0.46
1:A:383:SER:O	1:A:395:VAL:HA	2.15	0.46
1:A:773:VAL:HB	1:A:845:GLY:HA3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:870:LEU:N	1:A:870:LEU:HD12	2.29	0.46
1:A:898:THR:HG21	1:A:960:LYS:O	2.16	0.46
1:A:47:LYS:HD3	1:A:52:LEU:HD13	1.97	0.46
1:A:838:MET:HA	1:A:838:MET:HE3	1.97	0.46
1:A:412:GLU:OE2	1:A:566:THR:HG21	2.16	0.46
1:A:573:ARG:NH1	1:A:573:ARG:HG2	2.30	0.45
1:A:680:GLU:HB2	1:A:683:HIS:NE2	2.30	0.45
1:A:55:GLU:O	1:A:58:GLU:HG2	2.15	0.45
1:A:717:GLY:O	1:A:731:SER:HB2	2.17	0.45
1:A:17:GLY:O	1:A:28:GLN:NE2	2.44	0.45
1:A:13:LEU:HD23	1:A:222:ILE:HD12	1.97	0.45
1:A:369:ILE:HG13	1:A:528:VAL:HG11	1.98	0.45
1:A:567:ARG:NH2	1:A:571:PRO:HD3	2.31	0.45
1:A:201:ASN:HA	1:A:204:LYS:HD2	1.98	0.45
1:A:636:CYS:HB3	1:A:642:PHE:CD2	2.51	0.45
1:A:771:GLU:O	1:A:775:ILE:HG12	2.16	0.45
1:A:43:ALA:HA	1:A:120:LYS:HZ1	1.82	0.45
1:A:527:TYR:O	1:A:593:PHE:N	2.46	0.45
1:A:779:ALA:C	1:A:781:LEU:H	2.20	0.45
1:A:916:LEU:HD21	1:A:930:ASN:HB2	1.99	0.45
1:A:292:ALA:O	1:A:296:PHE:HD1	2.00	0.45
1:A:483:PHE:HE1	1:A:573:ARG:HD3	1.82	0.45
1:A:77:TRP:C	1:A:79:GLU:H	2.20	0.45
1:A:116:ILE:CD1	1:A:240:ALA:HB2	2.42	0.45
1:A:326:MET:CE	1:A:333:VAL:HG21	2.46	0.45
1:A:480:LYS:O	1:A:498:CYS:HA	2.17	0.45
1:A:847:ALA:CA	1:A:973:ILE:HD11	2.47	0.45
1:A:50:TRP:CZ2	1:A:54:ILE:HD11	2.52	0.44
1:A:766:SER:O	1:A:769:VAL:HB	2.17	0.44
1:A:150:ILE:CG2	1:A:220:LEU:HD11	2.47	0.44
1:A:522:ILE:HG22	1:A:542:LYS:HE3	1.99	0.44
1:A:877:THR:O	1:A:877:THR:HG22	2.18	0.44
1:A:957:PHE:O	1:A:958:LYS:HB2	2.18	0.44
1:A:119:LEU:CD1	1:A:236:ARG:HB2	2.47	0.44
1:A:192:GLU:CG	1:A:193:PRO:HD2	2.48	0.44
1:A:947:ILE:HG22	1:A:953:LEU:HD13	1.98	0.44
1:A:256:PHE:CZ	1:A:260:LEU:HD22	2.52	0.44
1:A:88:PHE:O	1:A:91:PRO:HG2	2.18	0.44
1:A:926:PRO:HA	1:A:927:PRO:HD3	1.88	0.44
1:A:93:VAL:O	1:A:97:ILE:HG13	2.18	0.44
1:A:486:GLU:O	1:A:491:ARG:NH2	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:567:ARG:HD2	1:A:570:PRO:CA	2.44	0.44
1:A:756:ASN:HB3	1:A:808:GLY:HA2	2.00	0.44
1:A:963:ASP:CB	1:A:966:GLN:NE2	2.77	0.44
1:A:273:LEU:HA	1:A:276:ILE:HD11	1.99	0.44
1:A:367:PHE:CD2	1:A:379:LEU:HD13	2.53	0.44
1:A:840:ILE:HD13	1:A:980:LEU:CD2	2.47	0.44
1:A:47:LYS:NZ	1:A:55:GLU:OE2	2.48	0.43
1:A:275:ASN:HB2	1:A:295:TYR:OH	2.17	0.43
1:A:916:LEU:CD1	1:A:927:PRO:HA	2.49	0.43
1:A:951:ASP:HA	1:A:954:PRO:HG2	2.00	0.43
1:A:9:THR:HB	1:A:166:LEU:HD23	1.99	0.43
1:A:328:LYS:CA	1:A:328:LYS:HE2	2.38	0.43
1:A:412:GLU:O	1:A:415:THR:N	2.51	0.43
1:A:411:VAL:HA	1:A:454:VAL:HG11	2.00	0.43
1:A:267:ILE:CG2	1:A:302:LEU:HD21	2.48	0.43
1:A:59:ASP:HB3	1:A:62:VAL:CG2	2.46	0.43
1:A:855:TRP:HA	1:A:859:ALA:HB2	1.98	0.43
1:A:412:GLU:OE2	1:A:529:ARG:NH1	2.51	0.43
1:A:688:VAL:O	1:A:692:GLN:HG3	2.18	0.43
1:A:865:VAL:HB	1:A:868:HIS:CG	2.54	0.43
1:A:275:ASN:N	1:A:275:ASN:HD22	2.17	0.43
1:A:49:LEU:O	1:A:53:VAL:HG23	2.18	0.43
1:A:502:LYS:O	1:A:502:LYS:HG3	2.18	0.43
1:A:863:PRO:HD3	1:A:890:ILE:HD11	1.99	0.43
1:A:978:ILE:O	1:A:982:GLU:HB2	2.19	0.43
1:A:858:TYR:HD1	1:A:858:TYR:H	1.62	0.43
1:A:926:PRO:CB	1:A:928:TRP:CE2	3.02	0.43
1:A:945:PHE:O	1:A:948:LEU:N	2.48	0.43
1:A:173:LEU:HD22	1:A:219:ALA:HB2	2.00	0.42
1:A:249:LEU:O	1:A:253:LEU:HB2	2.19	0.42
1:A:259:GLN:NE2	1:A:259:GLN:CA	2.82	0.42
1:A:317:THR:O	1:A:321:LEU:HG	2.18	0.42
1:A:380:ASN:HD22	1:A:382:PHE:HZ	1.67	0.42
1:A:769:VAL:HG12	1:A:841:GLY:HA3	2.01	0.42
1:A:335:SER:OG	1:A:337:PRO:HD2	2.18	0.42
1:A:926:PRO:HB3	1:A:928:TRP:CZ2	2.54	0.42
1:A:783:LEU:CD2	1:A:870:LEU:HD13	2.49	0.42
1:A:19:SER:HB3	1:A:22:THR:HG22	2.00	0.42
1:A:663:LEU:CD1	1:A:663:LEU:H	2.23	0.42
1:A:366:MET:HA	1:A:596:VAL:O	2.19	0.42
1:A:130:TYR:CZ	1:A:137:VAL:HB	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:ALA:O	1:A:155:VAL:C	2.58	0.42
1:A:355:THR:HA	1:A:738:ASP:O	2.19	0.42
1:A:558:THR:HG22	1:A:634:ALA:HB1	2.01	0.42
1:A:969:MET:O	1:A:973:ILE:HG22	2.20	0.42
1:A:181:THR:OG1	1:A:183:GLU:HG3	2.20	0.42
1:A:159:VAL:HG13	1:A:160:PRO:HD2	2.02	0.42
1:A:188:ILE:HD12	1:A:188:ILE:N	2.35	0.42
1:A:754:TYR:CD1	1:A:757:MET:HE3	2.55	0.41
1:A:762:ARG:O	1:A:766:SER:HB3	2.20	0.41
1:A:132:ALA:C	1:A:134:ARG:H	2.24	0.41
1:A:617:ALA:O	1:A:821:PRO:HD3	2.20	0.41
1:A:861:ASP:OD2	1:A:966:GLN:NE2	2.52	0.41
1:A:104:VAL:HG21	3:A:1002:PTY:HC11	2.01	0.41
1:A:353:THR:HA	1:A:357:THR:OG1	2.20	0.41
1:A:783:LEU:HD23	1:A:870:LEU:HD13	2.02	0.41
1:A:408:ASP:HA	1:A:411:VAL:CG2	2.50	0.41
1:A:933:LEU:O	1:A:937:ILE:HG13	2.20	0.41
1:A:122:TYR:HE1	1:A:726:VAL:CG2	2.32	0.41
1:A:267:ILE:O	1:A:271:VAL:HG23	2.21	0.41
1:A:388:THR:HG21	1:A:390:ALA:HB3	2.02	0.41
1:A:85:ILE:O	1:A:85:ILE:HG13	2.20	0.41
1:A:973:ILE:O	1:A:973:ILE:HG12	2.20	0.41
1:A:115:ALA:O	1:A:118:ALA:HB3	2.20	0.41
1:A:177:GLN:HG2	1:A:212:THR:HG21	2.02	0.41
1:A:369:ILE:HD13	1:A:379:LEU:HD22	2.02	0.41
1:A:874:MET:HG2	1:A:891:PHE:CE2	2.56	0.41
1:A:247:THR:OG1	1:A:250:GLN:HG3	2.21	0.41
1:A:314:VAL:HG11	1:A:804:ALA:CB	2.50	0.41
1:A:326:MET:HE1	1:A:333:VAL:HG21	2.02	0.41
1:A:715:GLU:O	1:A:732:GLU:HG2	2.21	0.41
1:A:0:ACE:CH3	1:A:36:TYR:CE1	3.03	0.41
1:A:298:ILE:O	1:A:302:LEU:HB2	2.21	0.41
1:A:380:ASN:HB3	1:A:382:PHE:CZ	2.56	0.41
1:A:658:PHE:CZ	1:A:666:GLN:HB3	2.55	0.41
1:A:773:VAL:CG1	1:A:845:GLY:HA3	2.50	0.41
1:A:989:ARG:HH21	3:A:1003:PTY:HC51	1.78	0.41
1:A:273:LEU:HA	1:A:276:ILE:HG13	2.02	0.41
1:A:873:PHE:CD2	1:A:876:CYS:HA	2.53	0.41
1:A:242:THR:HB	1:A:712:LYS:HD3	2.02	0.41
1:A:769:VAL:CG1	1:A:841:GLY:HA3	2.51	0.40
1:A:799:THR:HG21	1:A:905:VAL:HG22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:874:MET:HG2	1:A:891:PHE:HE2	1.86	0.40
1:A:114:ASN:OD1	1:A:117:GLU:HG2	2.22	0.40
1:A:194:VAL:HA	1:A:195:PRO:HD2	1.67	0.40
1:A:283:VAL:HG13	1:A:284:HIS:N	2.36	0.40
1:A:860:GLU:O	1:A:861:ASP:C	2.59	0.40
1:A:407:PHE:O	1:A:411:VAL:HG23	2.22	0.40
1:A:646:GLU:OE1	1:A:651:ARG:NH1	2.54	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	993/995 (100%)	891 (90%)	88 (9%)	14 (1%)	11 20

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	463	SER
1	A	883	PHE
1	A	184	SER
1	A	309	GLU
1	A	858	TYR
1	A	861	ASP
1	A	869	GLN
1	A	133	ASP
1	A	155	VAL
1	A	865	VAL
1	A	872	HIS
1	A	951	ASP
1	A	47	LYS
1	A	185	VAL

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	840/840 (100%)	791 (94%)	49 (6%)	20 38

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	22	THR
1	A	44	GLU
1	A	83	GLU
1	A	95	LEU
1	A	113	GLU
1	A	133	ASP
1	A	139	ARG
1	A	149	ASP
1	A	158	LYS
1	A	164	ARG
1	A	183	GLU
1	A	198	ARG
1	A	205	LYS
1	A	236	ARG
1	A	253	LEU
1	A	259	GLN
1	A	319	LEU
1	A	328	LYS
1	A	344	CYS
1	A	362	SER
1	A	371	LYS
1	A	380	ASN
1	A	396	LEU
1	A	445	LEU
1	A	467	ARG
1	A	484	THR
1	A	486	GLU
1	A	490	ASP
1	A	544	LYS

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Mol	Chain	Res	Type
1	A	567	ARG
1	A	574	GLU
1	A	578	LEU
1	A	580	ASP
1	A	604	ARG
1	A	620	ARG
1	A	628	ASN
1	A	656	ARG
1	A	685	SER
1	A	691	LEU
1	A	697	ILE
1	A	732	GLU
1	A	739	ASN
1	A	816	ILE
1	A	833	LEU
1	A	860	GLU
1	A	873	PHE
1	A	888	CYS
1	A	967	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	ASN
1	A	138	GLN
1	A	202	GLN
1	A	250	GLN
1	A	275	ASN
1	A	359	ASN
1	A	380	ASN
1	A	461	ASN
1	A	510	ASN
1	A	628	ASN
1	A	739	ASN
1	A	919	ASN
1	A	990	ASN

5.3.3 RNA

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PTY	A	1005	-	18,18,49	1.47	4 (22%)	21,23,54	1.41	2 (9%)
3	PTY	A	1006	-	18,18,49	1.47	4 (22%)	21,23,54	1.41	2 (9%)
3	PTY	A	1003	-	18,18,49	1.31	3 (16%)	21,23,54	1.19	2 (9%)
3	PTY	A	1004	-	18,18,49	1.21	2 (11%)	21,23,54	1.41	2 (9%)
3	PTY	A	1002	-	18,18,49	1.48	4 (22%)	21,23,54	1.41	2 (9%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PTY	A	1005	-	-	9/20/20/53	-
3	PTY	A	1006	-	-	9/20/20/53	-
3	PTY	A	1003	-	-	9/20/20/53	-
3	PTY	A	1004	-	-	11/20/20/53	-
3	PTY	A	1002	-	-	9/20/20/53	-

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	PTY	C1-C6	3.41	1.61	1.50
3	A	1005	PTY	C1-C6	3.39	1.61	1.50
3	A	1006	PTY	C1-C6	3.38	1.61	1.50
3	A	1003	PTY	C1-C6	2.93	1.59	1.50
3	A	1002	PTY	P1-O13	2.82	1.60	1.50
3	A	1006	PTY	P1-O13	2.80	1.60	1.50
3	A	1005	PTY	P1-O13	2.79	1.60	1.50
3	A	1003	PTY	P1-O13	2.65	1.60	1.50
3	A	1003	PTY	C5-C6	2.35	1.57	1.50
3	A	1004	PTY	P1-O13	2.34	1.59	1.50
3	A	1002	PTY	C5-C6	2.24	1.57	1.50
3	A	1005	PTY	C5-C6	2.24	1.57	1.50
3	A	1006	PTY	C5-C6	2.24	1.57	1.50
3	A	1004	PTY	C5-C6	2.15	1.57	1.50
3	A	1006	PTY	C2-C3	2.13	1.58	1.50
3	A	1002	PTY	C2-C3	2.13	1.58	1.50
3	A	1005	PTY	C2-C3	2.12	1.58	1.50

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1004	PTY	O7-C8-C11	3.75	117.99	111.09
3	A	1006	PTY	O7-C8-C11	3.52	117.56	111.09
3	A	1005	PTY	O7-C8-C11	3.50	117.53	111.09
3	A	1002	PTY	O7-C8-C11	3.49	117.51	111.09
3	A	1005	PTY	O4-C1-C6	2.77	116.48	108.43
3	A	1006	PTY	O4-C1-C6	2.76	116.47	108.43
3	A	1002	PTY	O4-C1-C6	2.75	116.44	108.43
3	A	1003	PTY	O4-C1-C6	2.52	115.77	108.43
3	A	1003	PTY	O7-C8-C11	2.27	115.26	111.09
3	A	1004	PTY	O12-P1-O11	2.03	117.19	107.75

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1005	PTY	C11-C8-O7-C6
3	A	1004	PTY	N1-C2-C3-O11
3	A	1004	PTY	C5-O14-P1-O11
3	A	1004	PTY	C5-O14-P1-O12
3	A	1004	PTY	C5-O14-P1-O13
3	A	1002	PTY	C11-C8-O7-C6

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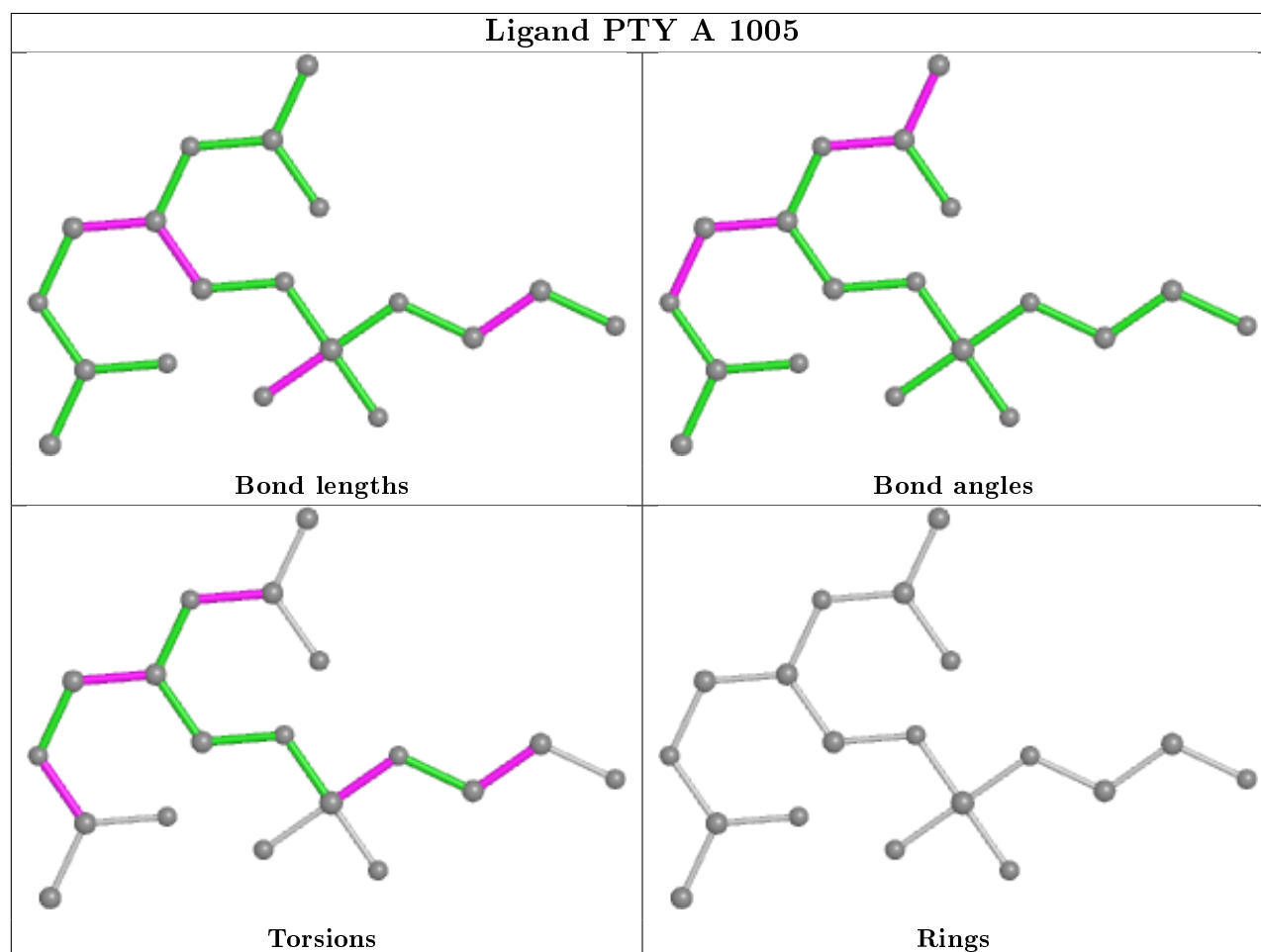
Mol	Chain	Res	Type	Atoms
3	A	1006	PTY	C11-C8-O7-C6
3	A	1003	PTY	C11-C8-O7-C6
3	A	1005	PTY	O10-C8-O7-C6
3	A	1004	PTY	C11-C8-O7-C6
3	A	1002	PTY	O10-C8-O7-C6
3	A	1006	PTY	O10-C8-O7-C6
3	A	1004	PTY	C31-C30-O4-C1
3	A	1003	PTY	C31-C30-O4-C1
3	A	1005	PTY	C31-C30-O4-C1
3	A	1002	PTY	C31-C30-O4-C1
3	A	1006	PTY	C31-C30-O4-C1
3	A	1004	PTY	O10-C8-O7-C6
3	A	1004	PTY	O30-C30-O4-C1
3	A	1003	PTY	O10-C8-O7-C6
3	A	1003	PTY	O30-C30-O4-C1
3	A	1005	PTY	O30-C30-O4-C1
3	A	1002	PTY	O30-C30-O4-C1
3	A	1006	PTY	O30-C30-O4-C1
3	A	1004	PTY	O4-C1-C6-C5
3	A	1003	PTY	O4-C1-C6-O7
3	A	1003	PTY	C3-O11-P1-O14
3	A	1005	PTY	N1-C2-C3-O11
3	A	1002	PTY	N1-C2-C3-O11
3	A	1006	PTY	N1-C2-C3-O11
3	A	1005	PTY	O4-C1-C6-O7
3	A	1002	PTY	O4-C1-C6-O7
3	A	1006	PTY	O4-C1-C6-O7
3	A	1004	PTY	O4-C1-C6-O7
3	A	1005	PTY	C3-O11-P1-O14
3	A	1004	PTY	C3-O11-P1-O14
3	A	1002	PTY	C3-O11-P1-O14
3	A	1006	PTY	C3-O11-P1-O14
3	A	1003	PTY	O4-C1-C6-C5
3	A	1005	PTY	O4-C1-C6-C5
3	A	1002	PTY	O4-C1-C6-C5
3	A	1006	PTY	O4-C1-C6-C5
3	A	1005	PTY	C3-O11-P1-O13
3	A	1002	PTY	C3-O11-P1-O13
3	A	1006	PTY	C3-O11-P1-O13
3	A	1003	PTY	C3-O11-P1-O13
3	A	1003	PTY	C2-C3-O11-P1

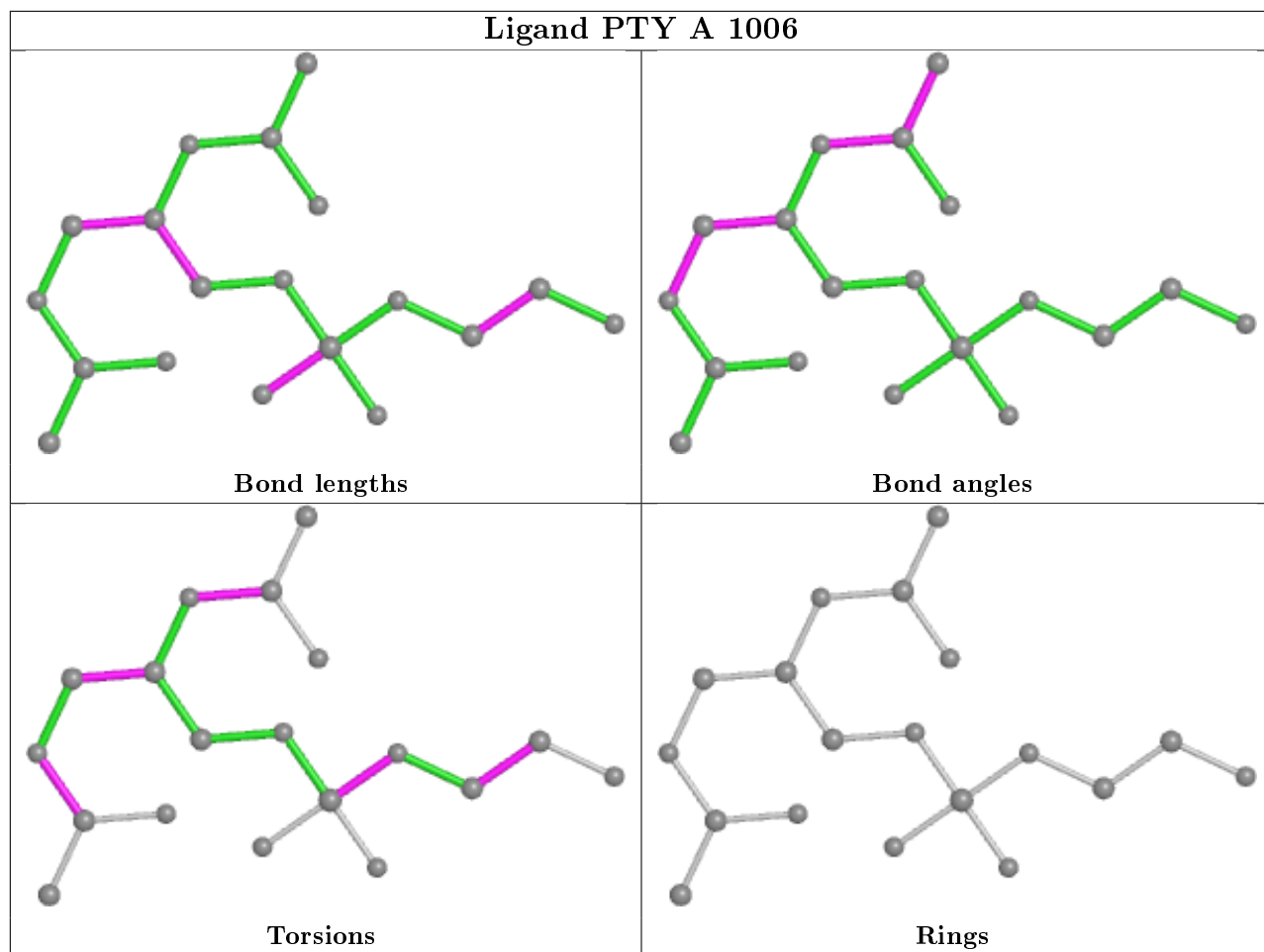
There are no ring outliers.

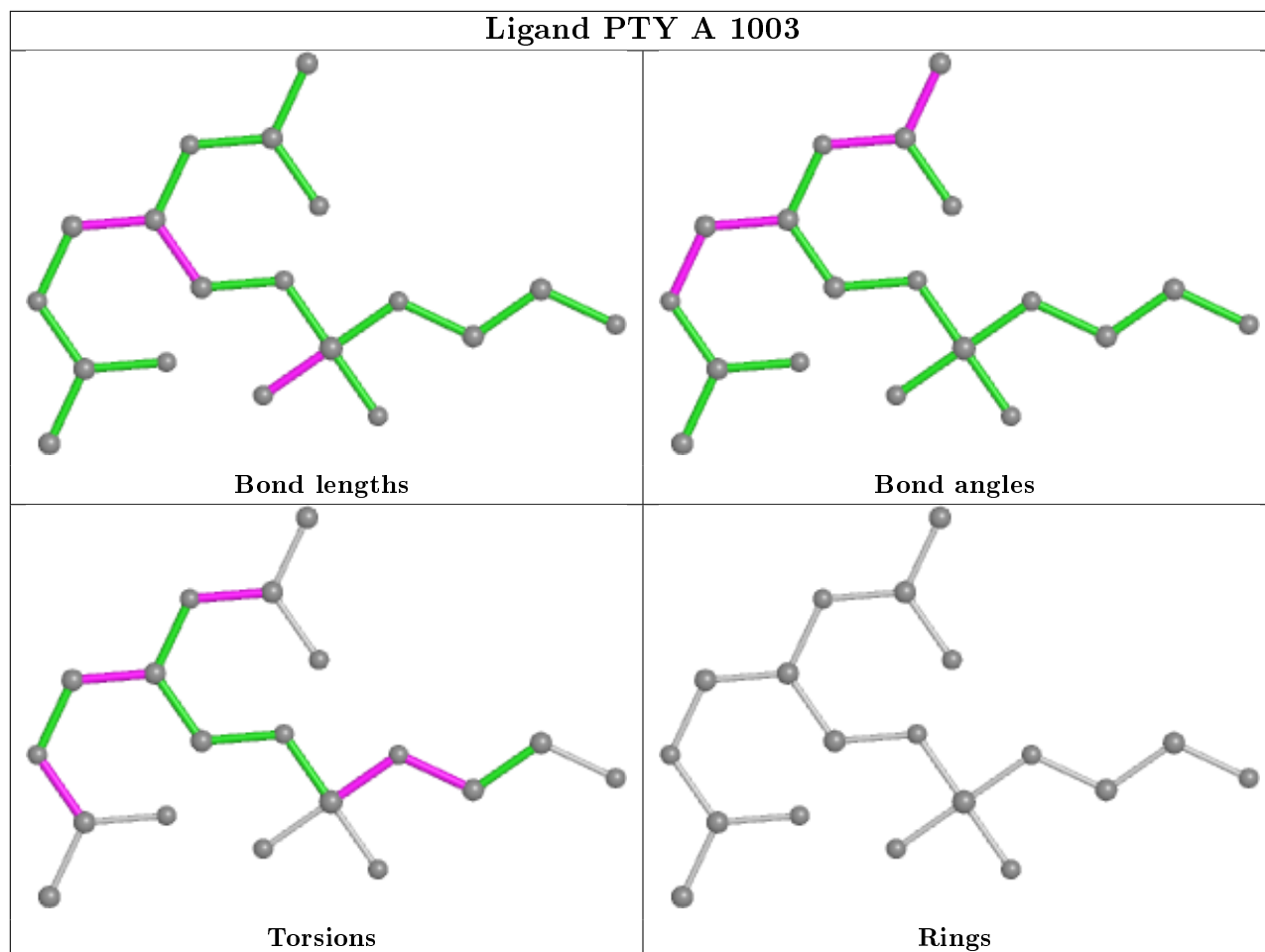
2 monomers are involved in 8 short contacts:

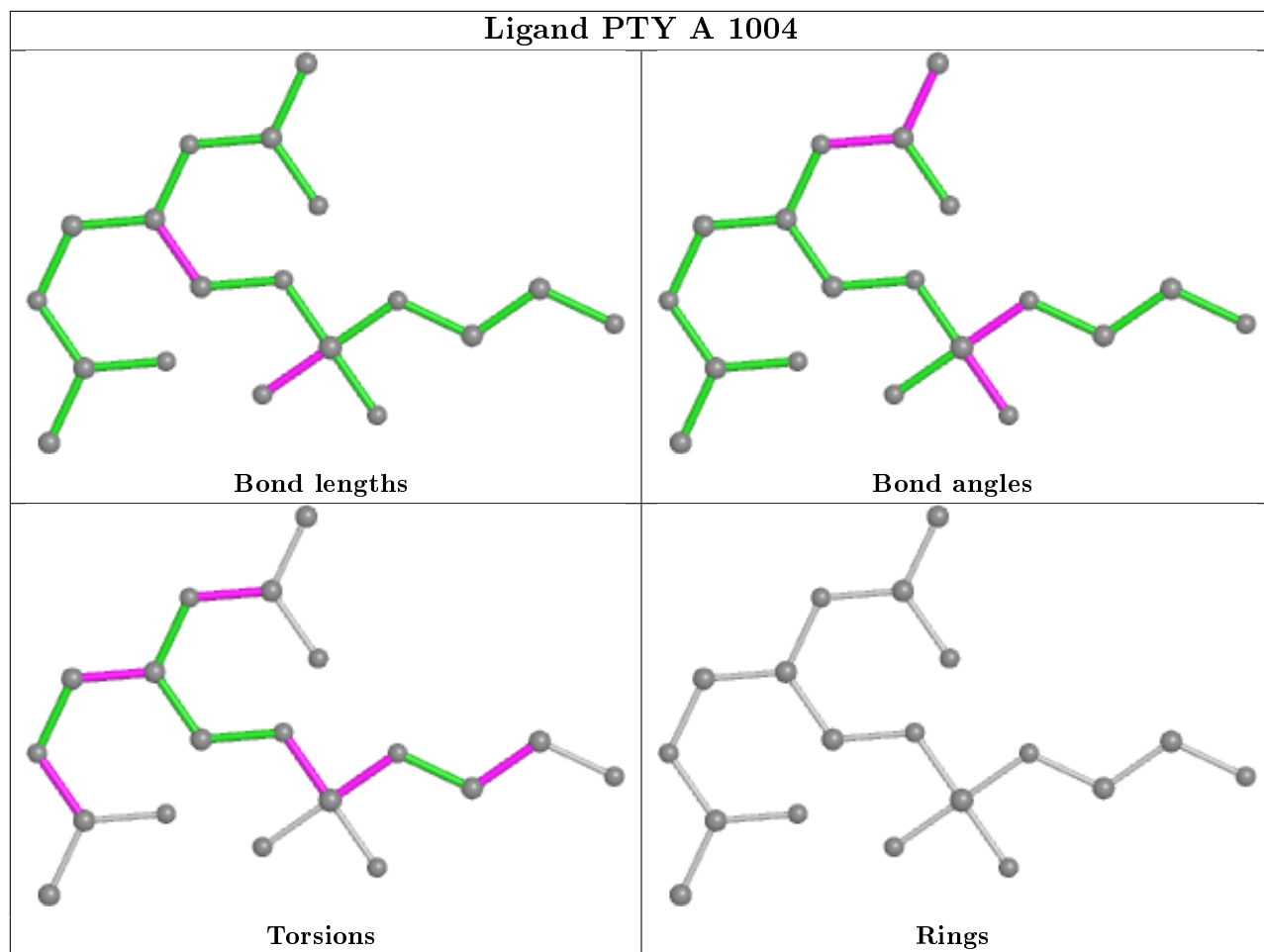
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1003	PTY	7	0
3	A	1002	PTY	1	0

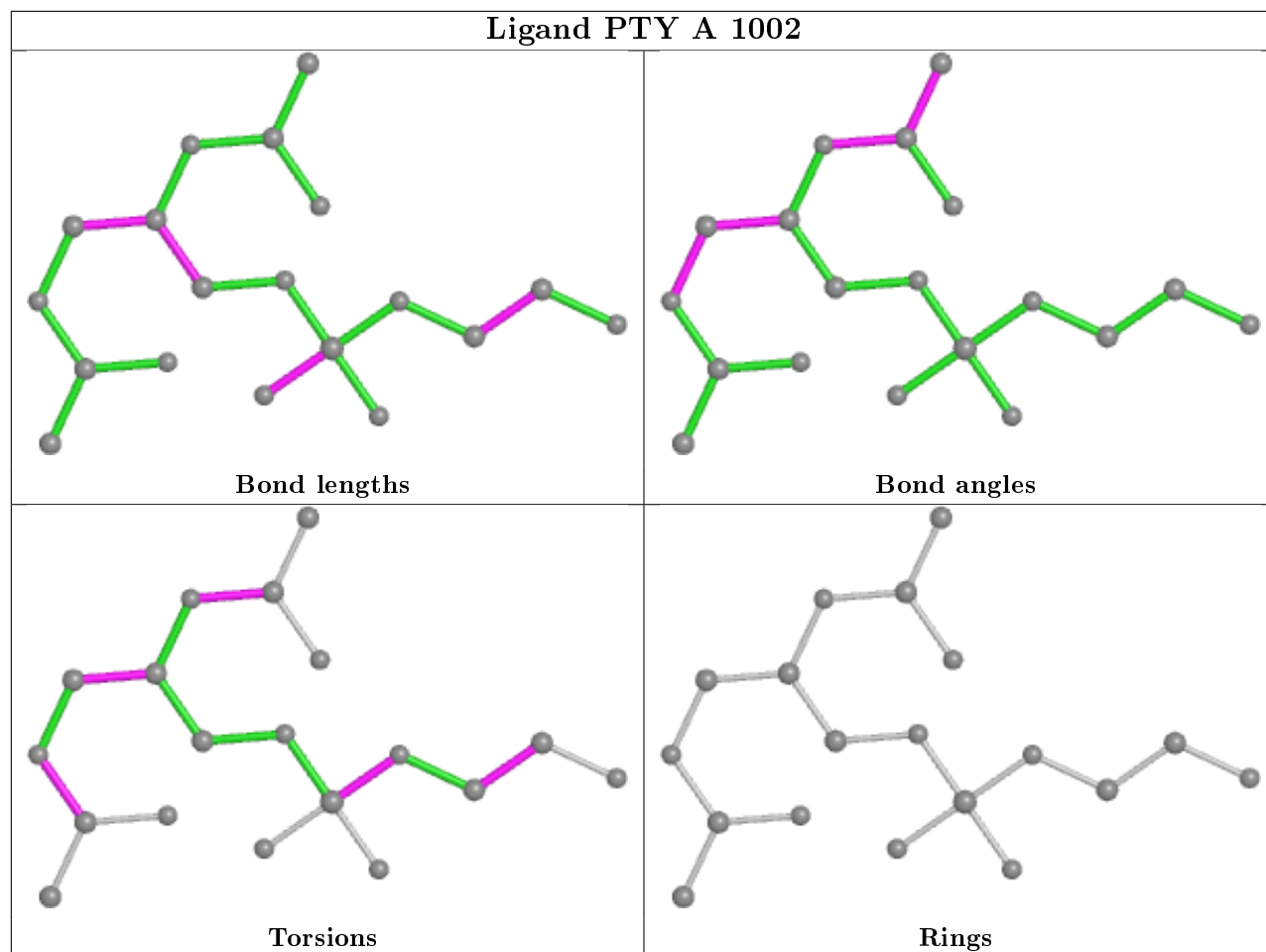
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	994/995 (99%)	0.53	74 (7%) 14 15	33, 70, 125, 150	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	504	SER	13.3
1	A	883	PHE	11.7
1	A	506	ALA	10.8
1	A	505	ARG	10.8
1	A	886	LEU	9.8
1	A	81	GLY	9.0
1	A	994	GLY	8.4
1	A	507	ALA	8.3
1	A	82	GLU	8.3
1	A	80	GLU	8.3
1	A	508	VAL	7.7
1	A	885	GLY	7.7
1	A	78	PHE	7.1
1	A	84	THR	7.1
1	A	877	THR	7.1
1	A	85	ILE	7.1
1	A	46	GLY	6.2
1	A	993	GLU	5.8
1	A	427	PHE	5.1
1	A	462	LEU	4.8
1	A	49	LEU	4.7
1	A	279	PHE	4.6
1	A	878	GLU	4.5
1	A	461	ASN	4.4
1	A	426	ASP	4.4
1	A	459	VAL	4.1
1	A	503	SER	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	884	GLU	4.1
1	A	53	VAL	4.1
1	A	856	PHE	4.0
1	A	50	TRP	3.7
1	A	86	THR	3.6
1	A	881	PRO	3.6
1	A	889	GLU	3.5
1	A	874	MET	3.5
1	A	45	GLU	3.3
1	A	106	VAL	3.2
1	A	289	ILE	3.1
1	A	955	MET	3.1
1	A	434	TYR	3.0
1	A	77	TRP	3.0
1	A	865	VAL	2.9
1	A	139	ARG	2.9
1	A	871	THR	2.9
1	A	52	LEU	2.8
1	A	433	VAL	2.7
1	A	828	LEU	2.7
1	A	959	LEU	2.7
1	A	44	GLU	2.7
1	A	873	PHE	2.6
1	A	875	GLN	2.5
1	A	888	CYS	2.4
1	A	464	LYS	2.4
1	A	894	PRO	2.4
1	A	790	VAL	2.4
1	A	866	THR	2.4
1	A	951	ASP	2.4
1	A	465	VAL	2.4
1	A	814	LEU	2.4
1	A	767	SER	2.3
1	A	293	ILE	2.3
1	A	220	LEU	2.3
1	A	308	PRO	2.2
1	A	858	TYR	2.2
1	A	460	ARG	2.2
1	A	192	GLU	2.1
1	A	285	GLY	2.1
1	A	116	ILE	2.1
1	A	887	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	795	VAL	2.1
1	A	992	LEU	2.1
1	A	342	LEU	2.0
1	A	435	GLU	2.0
1	A	880	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 carbohydrates in this entry.

6.4 Ligands [i](#)

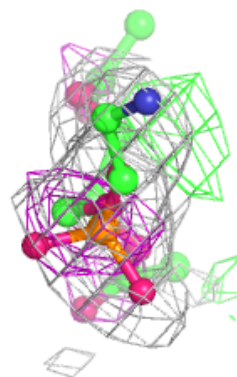
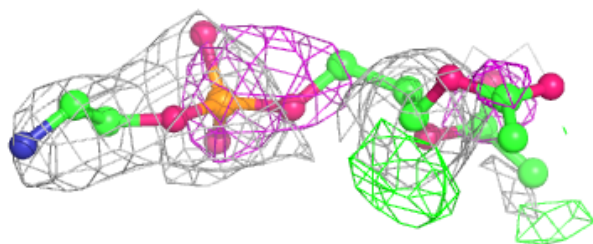
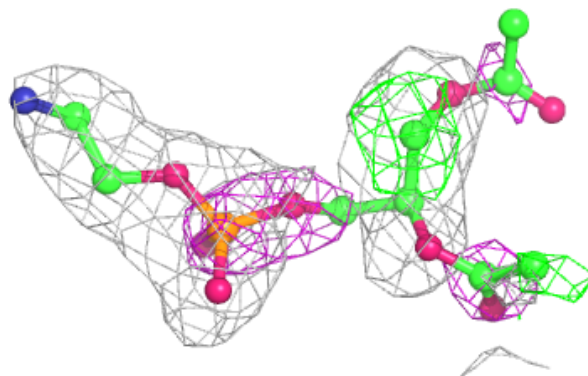
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PTY	A	1004	19/50	0.53	0.43	85,92,97,97	0
3	PTY	A	1002	19/50	0.60	0.50	108,110,113,114	0
3	PTY	A	1003	19/50	0.75	0.37	113,120,123,125	0
3	PTY	A	1005	19/50	0.76	0.35	108,110,113,114	0
3	PTY	A	1006	19/50	0.78	0.39	108,110,113,114	0
2	NA	A	1001	1/1	0.86	0.11	30,30,30,30	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.

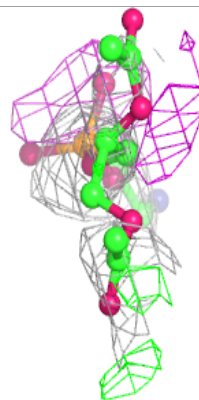
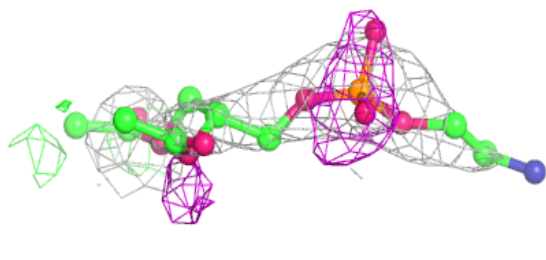
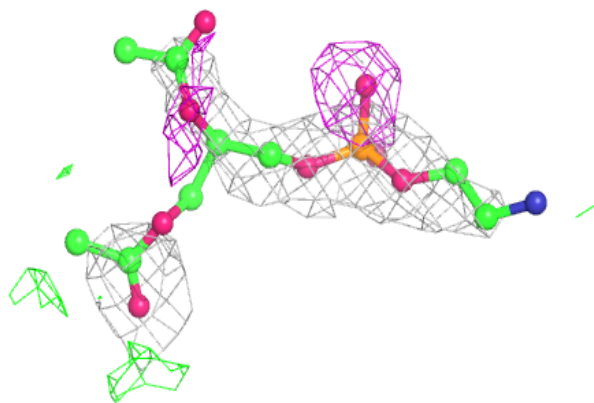
Electron density around PTY A 1004:

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

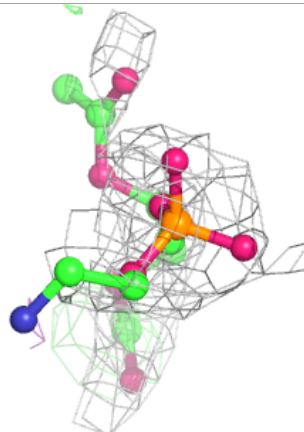
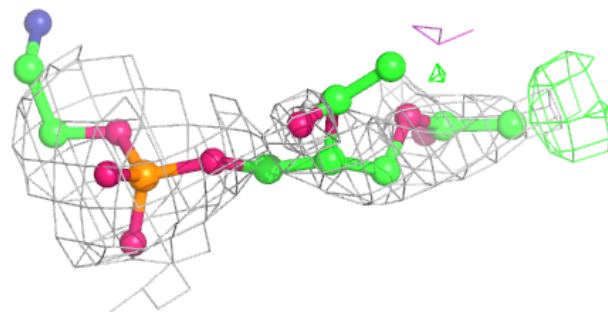
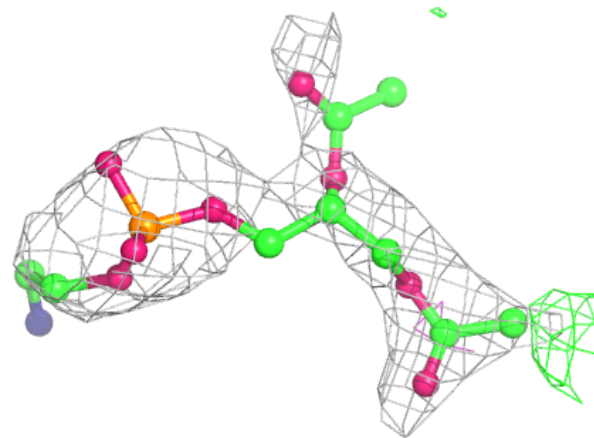


Electron density around PTY A 1002:

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

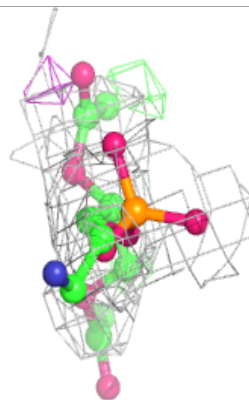
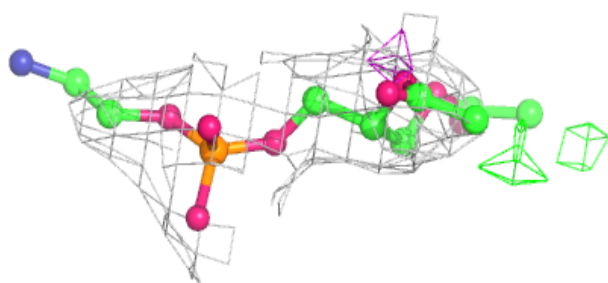
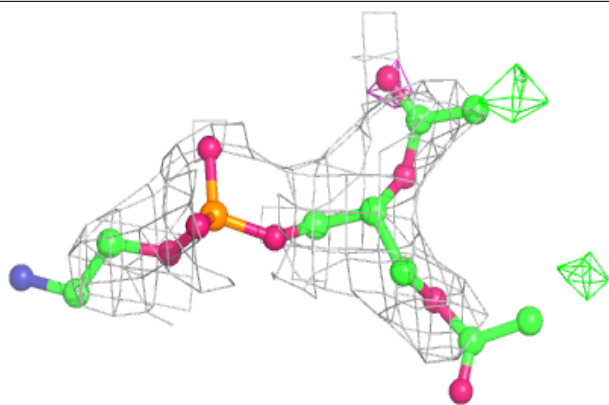
**Electron density around PTY A 1003:**

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

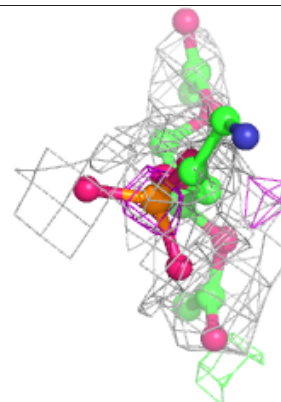
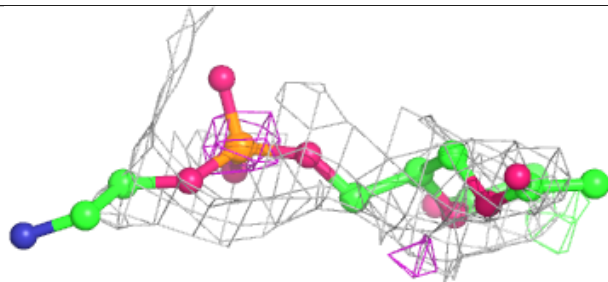
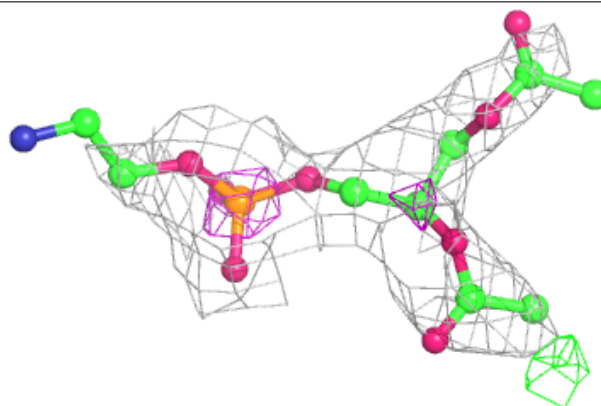


Electron density around PTY A 1005:

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

**Electron density around PTY A 1006:**

$2mF_o-DF_c$ (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.