



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 07:22 am BST

PDB ID : 2ZBD  
Title : Crystal Structure of the SR Calcium Pump with Bound Aluminium Fluoride, ADP and Calcium  
Authors : Toyoshima, C.; Nomura, H.; Tsuda, T.; Ogawa, H.; Norimatsu, Y.  
Deposited on : 2007-10-20  
Resolution : 2.40 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](mailto: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.

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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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

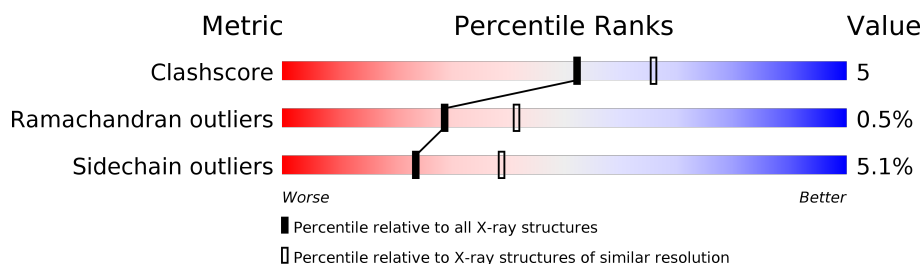
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)

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  $\geq 3$ , 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	995	

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7916 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 Sarcoplasmic/endoplasmic reticulum calcium ATPase 1.

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

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	994	GLY	-	SEE REMARK 999	UNP P04191

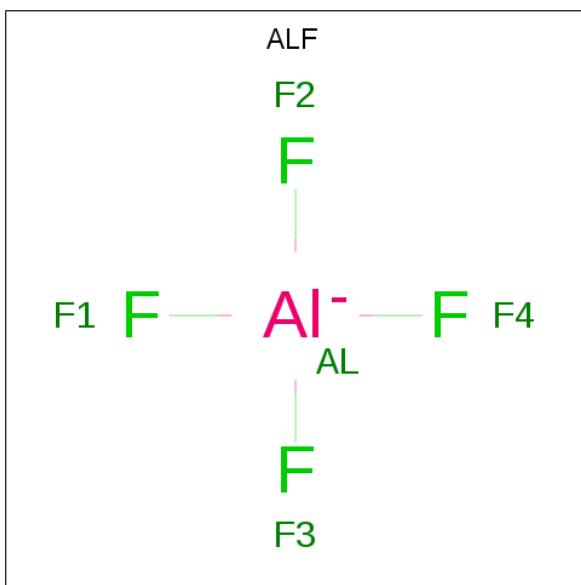
- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

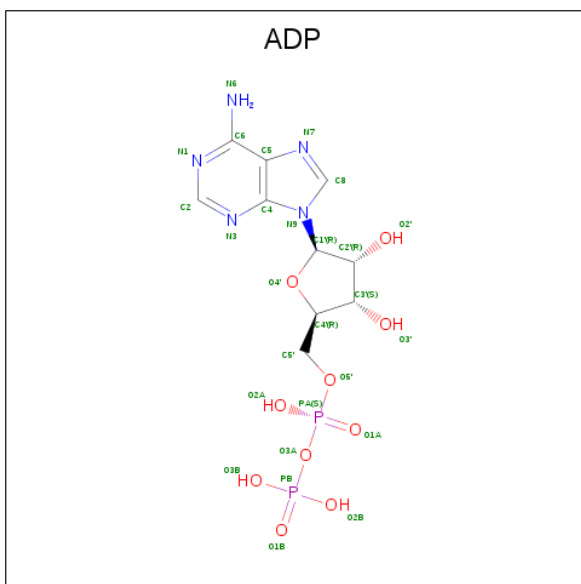
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is TETRAFLUOROALUMINATE ION (three-letter code: ALF) (formula:  $\text{AlF}_4$ ).



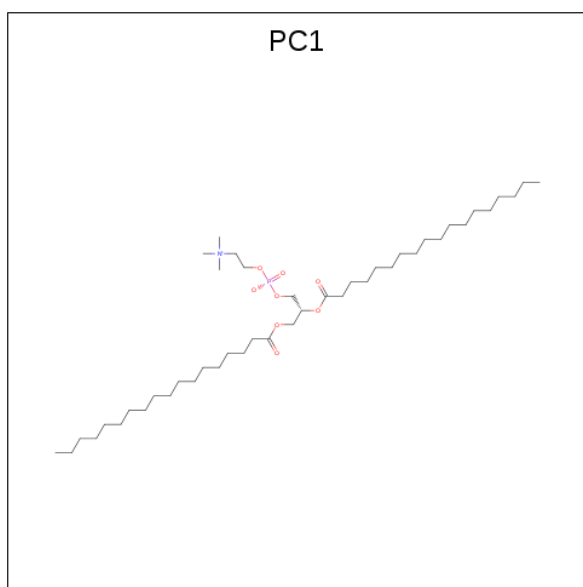
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	Al	F	0	0
			5	1	4		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{10}\text{P}_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 6 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PC1) (formula:  $\text{C}_{44}\text{H}_{88}\text{NO}_8\text{P}$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			54	44	1	8	1		
6	A	1	Total	C	N	O	P	0	0
			54	44	1	8	1		

- Molecule 7 is water.

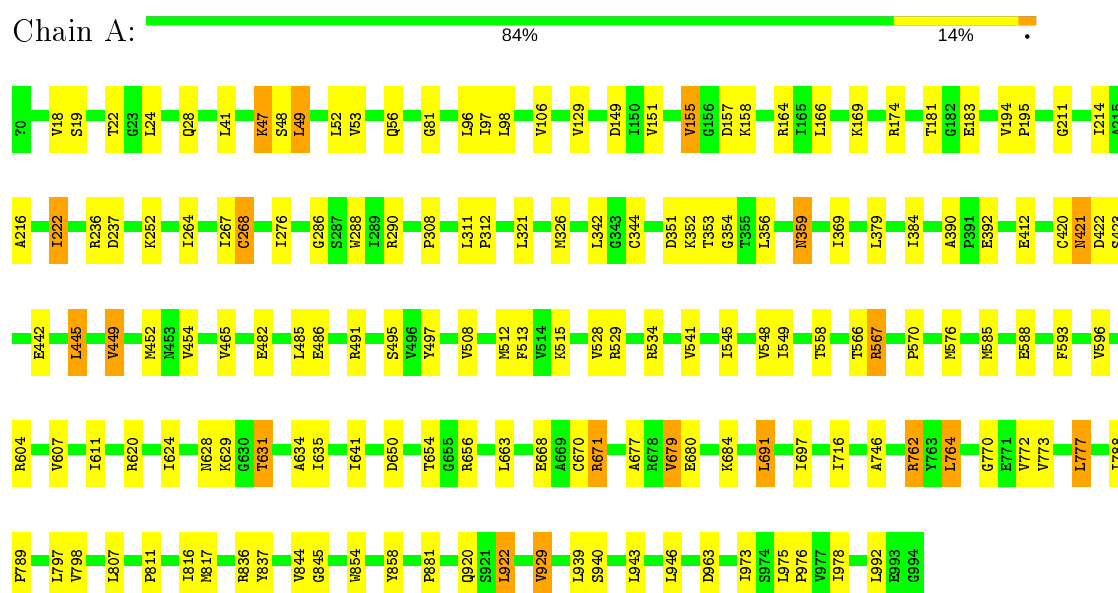
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	98	Total	O	0	0
			98	98		

### 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. 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. 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.

Note EDS was not executed.

- Molecule 1: Sarcoplasmic/endoplasmic reticulum calcium ATPase 1



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	162.90 Å   75.10 Å   152.10 Å 90.00°   109.30°   90.00°	Depositor
Resolution (Å)	12.00 – 2.40	Depositor
% Data completeness (in resolution range)	100.0 (12.00-2.40)	Depositor
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS, REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.246 , 0.271	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7916	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	92.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ADP, ACE, ALF, CA, PC1

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	0.35	0/7813	0.51	0/10594

There are no bond length outliers.

There are no bond angle outliers.

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	7674	0	7764	85	0
2	A	2	0	0	0	0
3	A	2	0	0	0	0
4	A	5	0	0	0	0
5	A	27	0	12	1	0
6	A	108	0	176	1	0
7	A	98	0	0	1	0
All	All	7916	0	7952	86	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 5.

All (86) 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:369:ILE:HD11	1:A:545:ILE:HD11	1.66	0.78
1:A:308:PRO:HB3	1:A:764:LEU:HD23	1.67	0.77
1:A:412:GLU:OE1	1:A:529:ARG:HD2	1.86	0.76
1:A:495:SER:HB2	1:A:588:GLU:OE1	1.90	0.72
1:A:770:GLY:HA3	1:A:844:VAL:HG23	1.75	0.67
1:A:354:GLY:HA2	1:A:359:ASN:HB2	1.80	0.64
1:A:286:GLY:HA3	1:A:290:ARG:HD3	1.81	0.63
1:A:421:ASN:HD22	1:A:423:SER:H	1.46	0.63
1:A:421:ASN:ND2	1:A:423:SER:H	1.99	0.60
1:A:129:VAL:HG12	1:A:151:VAL:HG22	1.84	0.59
1:A:19:SER:HB2	1:A:22:THR:HB	1.85	0.58
1:A:798:VAL:HG13	1:A:940:SER:HB3	1.86	0.57
1:A:515:LYS:HE3	5:A:1002:ADP:N1	2.20	0.57
1:A:975:LEU:N	1:A:976:PRO:HD2	2.21	0.56
1:A:549:ILE:HD11	1:A:596:VAL:HG21	1.87	0.56
1:A:52:LEU:HD23	1:A:106:VAL:HG13	1.87	0.55
1:A:491:ARG:NH1	1:A:588:GLU:OE1	2.40	0.54
1:A:513:PHE:HD1	1:A:566:THR:HG22	1.72	0.54
1:A:449:VAL:O	1:A:452:MET:O	2.26	0.53
1:A:157:ASP:HB2	1:A:214:ILE:HD13	1.91	0.53
1:A:670:CYS:HB3	1:A:691:LEU:HD13	1.89	0.53
1:A:56:GLN:HG3	1:A:106:VAL:HG23	1.91	0.53
1:A:773:VAL:HG23	1:A:845:GLY:HA3	1.90	0.53
1:A:788:ILE:HB	1:A:789:PRO:HD2	1.92	0.52
1:A:96:LEU:HD23	1:A:797:LEU:HD11	1.90	0.52
1:A:773:VAL:O	1:A:777:LEU:HB2	2.09	0.52
1:A:267:ILE:HD12	1:A:772:VAL:HG11	1.92	0.52
1:A:679:VAL:HG13	1:A:680:GLU:N	2.25	0.51
1:A:311:LEU:HB3	1:A:312:PRO:HD3	1.94	0.50
1:A:158:LYS:HE3	1:A:211:GLY:HA2	1.94	0.50
1:A:811:PRO:HG2	1:A:929:VAL:HG22	1.93	0.49
1:A:816:ILE:HG23	1:A:817:MET:HG2	1.94	0.49
1:A:47:LYS:HG3	1:A:48:SER:H	1.78	0.49
1:A:166:LEU:HD11	1:A:222:ILE:HB	1.94	0.48
1:A:24:LEU:HG	1:A:28:GLN:HB3	1.95	0.48
1:A:770:GLY:HA2	1:A:773:VAL:HG22	1.95	0.48
1:A:445:LEU:O	1:A:449:VAL:HG12	2.13	0.48
1:A:264:ILE:O	1:A:268:CYS:HB2	2.14	0.47
1:A:620:ARG:NH2	1:A:671:ARG:HA	2.29	0.47
1:A:762:ARG:HG2	1:A:837:TYR:HE1	1.80	0.47
1:A:326:MET:CE	1:A:746:ALA:HB2	2.45	0.47
1:A:558:THR:HG22	1:A:634:ALA:HB1	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:770:GLY:HA3	1:A:844:VAL:CG2	2.44	0.46
1:A:352:LYS:HA	1:A:356:LEU:HD12	1.97	0.46
1:A:326:MET:HE1	1:A:746:ALA:HB2	1.98	0.46
1:A:369:ILE:CD1	1:A:545:ILE:HD11	2.39	0.46
1:A:420:CYS:SG	1:A:515:LYS:HE2	2.56	0.45
1:A:922:LEU:HD13	1:A:978:ILE:HG23	1.97	0.45
1:A:41:LEU:HD13	1:A:236:ARG:HG2	1.98	0.45
1:A:352:LYS:HD2	1:A:635:ILE:HD12	1.99	0.45
1:A:379:LEU:HD12	1:A:548:VAL:HG21	1.99	0.45
1:A:351:ASP:HA	1:A:624:ILE:O	2.17	0.45
1:A:486:GLU:O	1:A:491:ARG:NH2	2.51	0.44
1:A:528:VAL:HG11	1:A:541:VAL:HG11	2.00	0.44
1:A:567:ARG:HD2	1:A:570:PRO:HA	1.98	0.44
1:A:24:LEU:HD22	1:A:149:ASP:HB3	1.99	0.44
1:A:342:LEU:HG	1:A:716:ILE:HD13	2.00	0.44
1:A:654:THR:HA	1:A:677:ALA:O	2.18	0.44
1:A:194:VAL:HA	1:A:195:PRO:HD2	1.76	0.44
1:A:482:GLU:HG3	1:A:497:TYR:HD1	1.83	0.44
1:A:545:ILE:HD12	1:A:593:PHE:CE1	2.52	0.44
1:A:491:ARG:HG3	1:A:585:MET:HB2	1.99	0.43
1:A:174:ARG:HB2	1:A:216:ALA:HB3	1.99	0.43
1:A:344:CYS:HB3	1:A:697:ILE:HG13	1.99	0.43
1:A:541:VAL:O	1:A:545:ILE:HG12	2.18	0.43
1:A:611:ILE:HD13	1:A:641:ILE:HG13	2.00	0.43
1:A:512:MET:HB2	1:A:567:ARG:HB2	1.99	0.43
1:A:624:ILE:CG2	1:A:684:LYS:HG2	2.49	0.43
1:A:854:TRP:HE1	1:A:858:TYR:HD2	1.67	0.43
1:A:631:THR:HG21	7:A:1074:HOH:O	2.19	0.42
6:A:1011:PC1:H2F2	6:A:1011:PC1:H391	2.01	0.42
1:A:628:ASN:HB3	1:A:631:THR:HG23	2.00	0.42
1:A:49:LEU:O	1:A:53:VAL:HG23	2.19	0.42
1:A:629:LYS:HA	1:A:677:ALA:CB	2.50	0.42
1:A:390:ALA:C	1:A:392:GLU:H	2.24	0.41
1:A:352:LYS:HD2	1:A:635:ILE:CD1	2.49	0.41
1:A:939:LEU:O	1:A:943:LEU:HB2	2.20	0.41
1:A:442:GLU:O	1:A:445:LEU:HB2	2.20	0.41
1:A:485:LEU:HB2	1:A:495:SER:HB3	2.02	0.41
1:A:576:MET:HB2	1:A:576:MET:HE3	1.92	0.41
1:A:421:ASN:HD22	1:A:422:ASP:N	2.19	0.41
1:A:679:VAL:CG1	1:A:680:GLU:N	2.83	0.41
1:A:181:THR:OG1	1:A:183:GLU:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:604:ARG:HB2	1:A:607:VAL:HG23	2.03	0.40
1:A:214:ILE:H	1:A:214:ILE:HD12	1.85	0.40
1:A:155:VAL:HG13	1:A:216:ALA:HA	2.04	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%)	946 (95%)	42 (4%)	5 (0%)	29	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	GLY
1	A	992	LEU
1	A	47	LYS
1	A	508	VAL
1	A	881	PRO

### 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%)	797 (95%)	43 (5%)	24	39

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	VAL
1	A	49	LEU
1	A	97	ILE
1	A	98	LEU
1	A	155	VAL
1	A	164	ARG
1	A	169	LYS
1	A	222	ILE
1	A	237	ASP
1	A	252	LYS
1	A	268	CYS
1	A	276	ILE
1	A	288	TRP
1	A	321	LEU
1	A	353	THR
1	A	359	ASN
1	A	384	ILE
1	A	421	ASN
1	A	445	LEU
1	A	449	VAL
1	A	454	VAL
1	A	465	VAL
1	A	534	ARG
1	A	567	ARG
1	A	631	THR
1	A	650	ASP
1	A	656	ARG
1	A	663	LEU
1	A	668	GLU
1	A	671	ARG
1	A	679	VAL
1	A	691	LEU
1	A	762	ARG
1	A	764	LEU
1	A	777	LEU
1	A	807	LEU
1	A	836	ARG
1	A	920	GLN
1	A	922	LEU
1	A	929	VAL
1	A	946	LEU
1	A	963	ASP

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Mol	Chain	Res	Type
1	A	973	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	ASN
1	A	114	ASN
1	A	250	GLN
1	A	359	ASN
1	A	380	ASN
1	A	421	ASN
1	A	526	ASN
1	A	628	ASN
1	A	875	GLN
1	A	919	ASN
1	A	920	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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
6	PC1	A	1011	-	53,53,53	1.46	5 (9%)	59,61,61	1.28	5 (8%)
6	PC1	A	1012	-	53,53,53	1.46	5 (9%)	59,61,61	1.27	3 (5%)
4	ALF	A	998	-	0,4,4	0.00	-	-	-	-
5	ADP	A	1002	3	24,29,29	0.93	1 (4%)	29,45,45	1.31	4 (13%)

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
6	PC1	A	1011	-	-	30/57/57/57	-
6	PC1	A	1012	-	-	31/57/57/57	-
5	ADP	A	1002	3	-	2/12/32/32	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1011	PC1	O22-C21	7.19	1.43	1.22
6	A	1012	PC1	O22-C21	7.17	1.43	1.22
6	A	1011	PC1	O31-C31	4.60	1.46	1.33
6	A	1012	PC1	O31-C31	4.52	1.46	1.33
6	A	1012	PC1	O21-C21	3.29	1.43	1.34
6	A	1011	PC1	O21-C21	3.20	1.43	1.34
6	A	1012	PC1	C3A-C39	-3.09	1.34	1.51
6	A	1012	PC1	C2A-C29	-3.07	1.34	1.51
6	A	1011	PC1	C2A-C29	-3.05	1.34	1.51
6	A	1011	PC1	C3A-C39	-3.04	1.34	1.51
5	A	1002	ADP	C5-C4	2.42	1.47	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1012	PC1	O21-C21-O22	-6.11	108.94	123.70
6	A	1011	PC1	O21-C21-O22	-5.93	109.36	123.70
5	A	1002	ADP	N3-C2-N1	-3.80	122.74	128.68
6	A	1012	PC1	O22-C21-C22	-3.64	109.52	123.73
6	A	1011	PC1	O22-C21-C22	-3.63	109.58	123.73
6	A	1011	PC1	O31-C31-C32	2.74	120.52	111.91
5	A	1002	ADP	O3B-PB-O2B	2.57	117.47	107.64
6	A	1012	PC1	O31-C31-C32	2.52	119.82	111.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1002	ADP	PA-O3A-PB	-2.14	125.47	132.83
6	A	1011	PC1	C2-O21-C21	-2.10	112.62	117.79
6	A	1011	PC1	C3A-C39-C38	2.09	125.03	114.42
5	A	1002	ADP	C2-N1-C6	2.01	122.19	118.75

There are no chirality outliers.

All (63) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1012	PC1	O13-C11-C12-N
6	A	1012	PC1	O22-C21-O21-C2
6	A	1011	PC1	O22-C21-O21-C2
6	A	1012	PC1	C28-C29-C2A-C2B
6	A	1012	PC1	C38-C39-C3A-C3B
6	A	1011	PC1	C28-C29-C2A-C2B
6	A	1011	PC1	C38-C39-C3A-C3B
6	A	1012	PC1	C31-C32-C33-C34
6	A	1012	PC1	C21-C22-C23-C24
6	A	1012	PC1	C22-C21-O21-C2
6	A	1011	PC1	C11-O13-P-O11
6	A	1012	PC1	C34-C35-C36-C37
6	A	1011	PC1	C29-C2A-C2B-C2C
6	A	1012	PC1	C24-C25-C26-C27
6	A	1012	PC1	C2A-C2B-C2C-C2D
6	A	1012	PC1	C32-C33-C34-C35
6	A	1012	PC1	C37-C38-C39-C3A
6	A	1011	PC1	C2B-C2C-C2D-C2E
6	A	1011	PC1	C2C-C2D-C2E-C2F
6	A	1011	PC1	C22-C23-C24-C25
6	A	1011	PC1	C3D-C3E-C3F-C3G
6	A	1011	PC1	C23-C24-C25-C26
6	A	1011	PC1	C3A-C3B-C3C-C3D
6	A	1012	PC1	C25-C26-C27-C28
6	A	1012	PC1	C3C-C3D-C3E-C3F
6	A	1012	PC1	C22-C23-C24-C25
6	A	1012	PC1	C33-C34-C35-C36
6	A	1012	PC1	C35-C36-C37-C38
6	A	1011	PC1	C22-C21-O21-C2
6	A	1011	PC1	C25-C26-C27-C28
6	A	1011	PC1	C36-C37-C38-C39
6	A	1011	PC1	C27-C28-C29-C2A
6	A	1012	PC1	C3A-C3B-C3C-C3D

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms
6	A	1012	PC1	C2D-C2E-C2F-C2G
6	A	1012	PC1	O11-C1-C2-O21
6	A	1011	PC1	C26-C27-C28-C29
6	A	1012	PC1	C3F-C3G-C3H-C3I
6	A	1012	PC1	O11-C1-C2-C3
6	A	1011	PC1	C24-C25-C26-C27
6	A	1011	PC1	C2A-C2B-C2C-C2D
6	A	1011	PC1	C37-C38-C39-C3A
6	A	1012	PC1	C2B-C2C-C2D-C2E
6	A	1011	PC1	C3F-C3G-C3H-C3I
6	A	1011	PC1	C11-O13-P-O14
6	A	1011	PC1	O11-C1-C2-O21
6	A	1011	PC1	C21-C22-C23-C24
6	A	1012	PC1	C3D-C3E-C3F-C3G
6	A	1012	PC1	C27-C28-C29-C2A
6	A	1011	PC1	C34-C35-C36-C37
6	A	1011	PC1	C3E-C3F-C3G-C3H
5	A	1002	ADP	PA-O3A-PB-O1B
6	A	1011	PC1	C35-C36-C37-C38
6	A	1012	PC1	C36-C37-C38-C39
6	A	1011	PC1	C3C-C3D-C3E-C3F
6	A	1011	PC1	O11-C1-C2-C3
5	A	1002	ADP	PA-O3A-PB-O3B
6	A	1012	PC1	O31-C31-C32-C33
6	A	1012	PC1	C29-C2A-C2B-C2C
6	A	1012	PC1	C3E-C3F-C3G-C3H
6	A	1012	PC1	O22-C21-C22-C23
6	A	1012	PC1	O32-C31-C32-C33
6	A	1011	PC1	O31-C31-C32-C33
6	A	1011	PC1	C3B-C3C-C3D-C3E

There are no ring outliers.

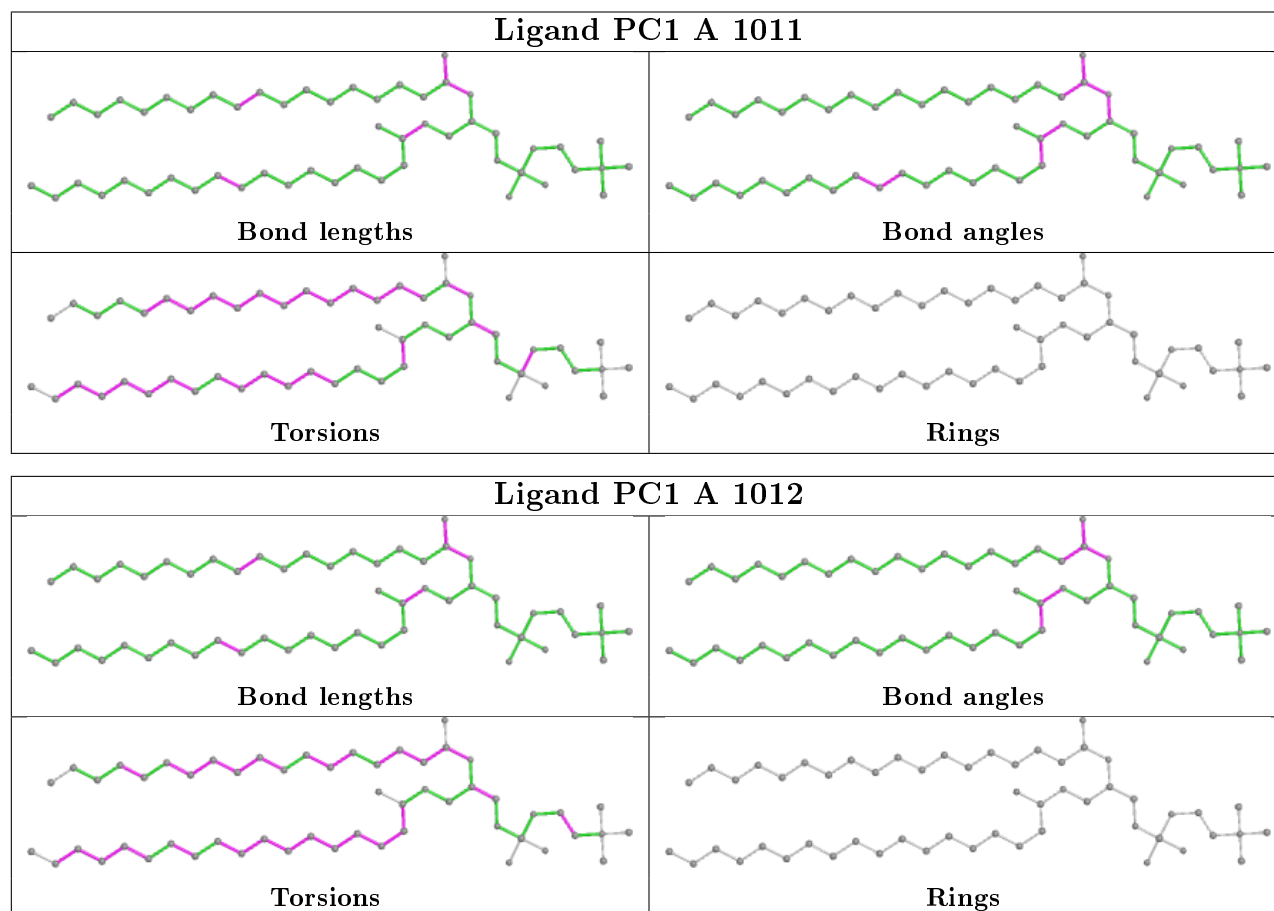
2 monomers are involved in 2 short contacts:

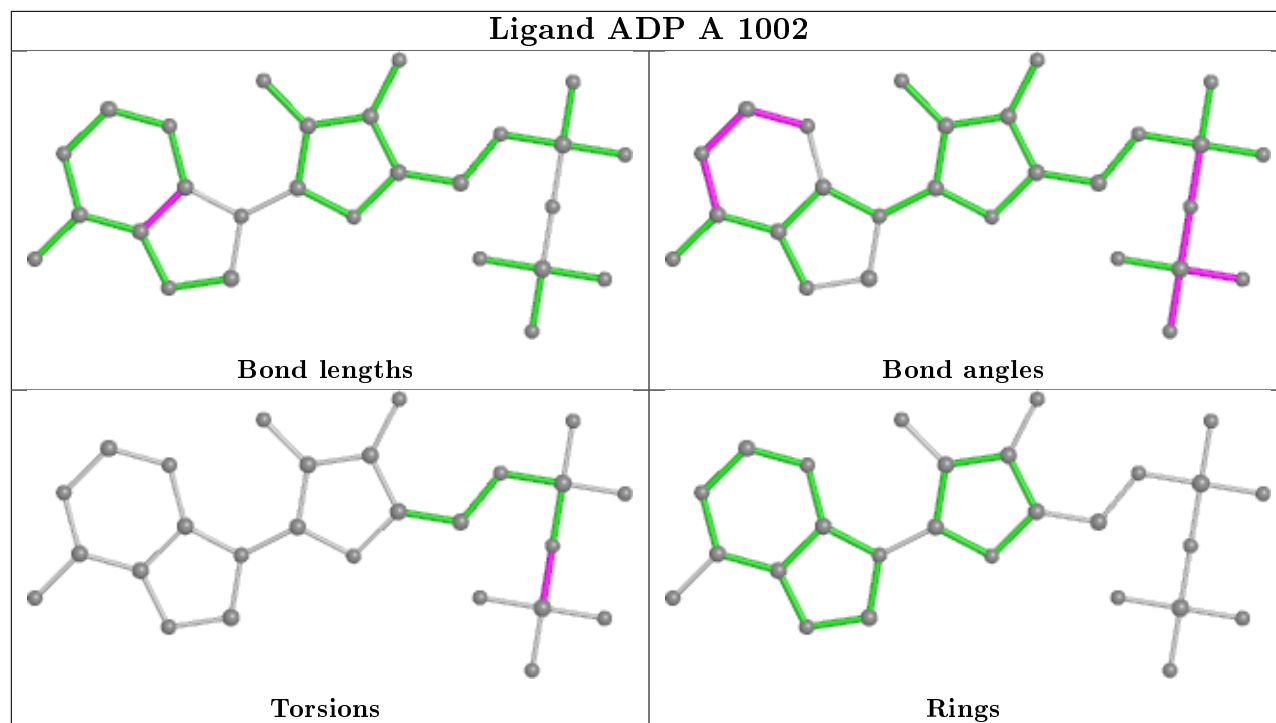
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1011	PC1	1	0
5	A	1002	ADP	1	0

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 ⓘ

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.