



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:31 PM UTC

PDB ID : 3AR2 / pdb_00003ar2
Title : Calcium pump crystal structure with bound AMPPCP and Ca²⁺
Authors : Toyoshima, C.; Yonekura, S.; Tsueda, J.; Iwasawa, S.
Deposited on : 2010-11-24
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

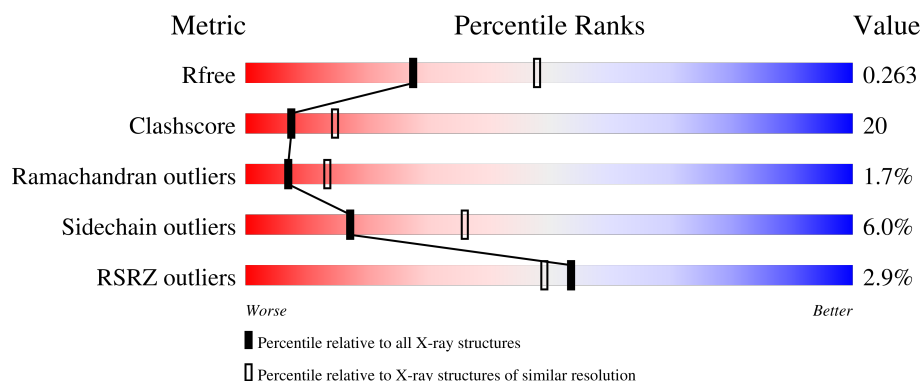
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

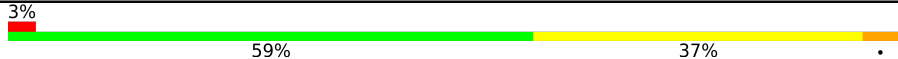
The reported resolution of this entry is 2.50 Å.

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	995	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7808 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			

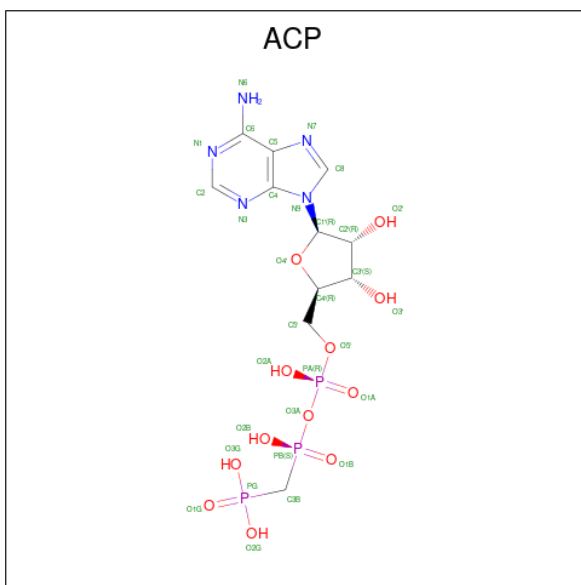
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ACE	-	acetylation	UNP P04191
A	994	GLY	ASP	SEE REMARK 999	UNP P04191

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 3 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (CCD ID: ACP) (formula: C₁₁H₁₈N₅O₁₂P₃).

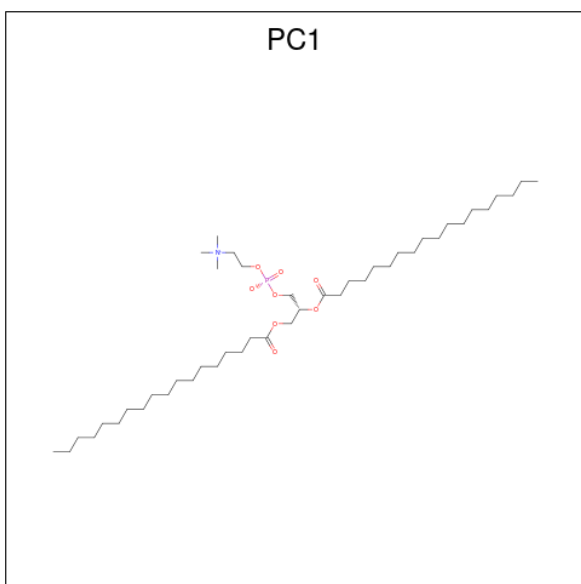


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

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

- Molecule 5 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PC1) (formula: $C_{44}H_{88}NO_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			20	10	1	8	1		

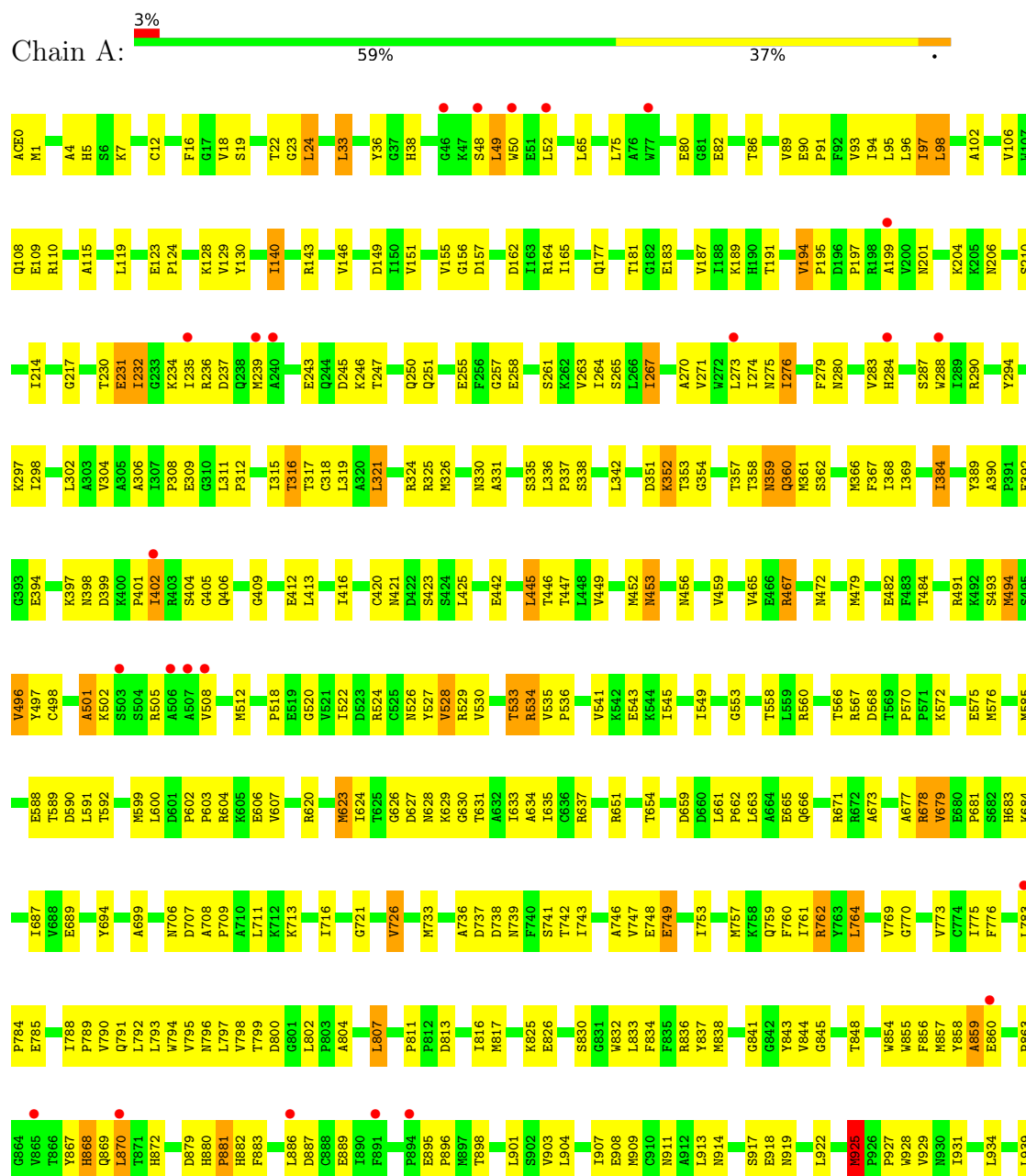
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	79	Total	O	0	0
			79	79		

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: Sarcoplasmic/endoplasmic reticulum calcium ATPase 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.00Å 75.31Å 151.54Å 90.00° 108.97° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 15.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (15.00-2.50) 98.8 (15.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 2.51Å)	Xtriage
Refinement program	CNS, REFMAC 5.5.0109	Depositor
R, R_{free}	0.247 , 0.272 0.239 , 0.263	Depositor DCC
R_{free} test set	3001 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	63.3	Xtriage
Anisotropy	0.568	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 87.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7808	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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, CA, ACP, PC1, 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	0.55	0/7813	0.89	8/10594 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	661	LEU	CA-C-N	-7.19	112.91	120.03
1	A	661	LEU	C-N-CA	-7.19	112.91	120.03
1	A	232	ILE	N-CA-C	-6.81	106.40	112.12
1	A	194	VAL	CA-C-N	5.93	125.41	119.24
1	A	194	VAL	C-N-CA	5.93	125.41	119.24
1	A	870	LEU	N-CA-C	-5.43	107.02	113.97
1	A	925	MET	CA-C-N	5.13	125.67	120.38
1	A	925	MET	C-N-CA	5.13	125.67	120.38

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	7764	315	0
2	A	3	0	0	0	0
3	A	31	0	14	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	1	0	0	0	0
5	A	20	0	18	0	0
6	A	79	0	0	2	0
All	All	7808	0	7796	315	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 20.

All (315) 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:361:MET:HE1	1:A:560:ARG:HD3	1.41	1.00
1:A:369:ILE:HD11	1:A:545:ILE:HD11	1.45	0.98
1:A:567:ARG:HD3	1:A:570:PRO:HA	1.46	0.96
1:A:624:ILE:HG21	1:A:684:LYS:HG2	1.47	0.96
1:A:342:LEU:HG	1:A:716:ILE:HD13	1.53	0.91
1:A:352:LYS:HB2	1:A:623:MET:HE2	1.54	0.90
1:A:624:ILE:CG2	1:A:684:LYS:HG2	2.03	0.89
1:A:65:LEU:HD22	1:A:309:GLU:HG2	1.56	0.88
1:A:528:VAL:HG11	1:A:541:VAL:HG11	1.54	0.87
1:A:308:PRO:HB3	1:A:764:LEU:HD23	1.57	0.85
1:A:773:VAL:HG23	1:A:845:GLY:HA3	1.58	0.84
1:A:917:SER:HB2	1:A:925:MET:HE3	1.63	0.80
1:A:129:VAL:HG12	1:A:151:VAL:HG22	1.62	0.79
1:A:94:ILE:HD11	1:A:793:LEU:HD22	1.65	0.79
1:A:631:THR:O	1:A:635:ILE:HG12	1.84	0.78
1:A:679:VAL:HG22	1:A:683:HIS:CB	2.14	0.77
1:A:75:LEU:HD22	1:A:297:LYS:HD3	1.64	0.77
1:A:361:MET:HE3	1:A:599:MET:SD	2.25	0.76
1:A:65:LEU:CD2	1:A:309:GLU:HG2	2.17	0.74
1:A:330:ASN:HB2	1:A:737:ASP:HB2	1.70	0.74
1:A:951:ASP:HB3	1:A:952:PRO:HD3	1.70	0.74
1:A:90:GLU:HB2	1:A:91:PRO:HD3	1.68	0.73
1:A:311:LEU:HB3	1:A:312:PRO:HD3	1.71	0.71
1:A:762:ARG:HG2	1:A:837:TYR:HE1	1.55	0.71
1:A:90:GLU:HG2	1:A:790:VAL:CG1	2.21	0.71
1:A:369:ILE:CD1	1:A:545:ILE:HD11	2.21	0.70
1:A:412:GLU:OE1	1:A:529:ARG:HD2	1.92	0.70
1:A:770:GLY:HA3	1:A:844:VAL:CG2	2.22	0.70
1:A:679:VAL:HG22	1:A:683:HIS:CG	2.27	0.70
1:A:572:LYS:HB2	1:A:575:GLU:HG2	1.72	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:770:GLY:HA3	1:A:844:VAL:HG23	1.75	0.69
1:A:96:LEU:HD23	1:A:797:LEU:HD11	1.74	0.69
1:A:928:TRP:HA	1:A:934:LEU:HD11	1.76	0.68
1:A:210:SER:O	1:A:232:ILE:HG12	1.94	0.68
1:A:671:ARG:HD3	1:A:694:TYR:CZ	2.29	0.67
1:A:518:PRO:O	1:A:522:ILE:HD13	1.96	0.66
1:A:115:ALA:HB2	1:A:324:ARG:HH12	1.59	0.66
1:A:904:LEU:O	1:A:908:GLU:HG3	1.96	0.66
1:A:791:GLN:HG2	1:A:901:LEU:HD22	1.78	0.66
1:A:308:PRO:HB3	1:A:764:LEU:CD2	2.25	0.66
1:A:361:MET:HE1	1:A:560:ARG:CD	2.24	0.66
1:A:898:THR:HG22	1:A:962:LEU:HD11	1.78	0.66
1:A:869:GLN:NE2	1:A:872:HIS:H	1.94	0.65
1:A:679:VAL:HG22	1:A:683:HIS:HB3	1.78	0.65
1:A:922:LEU:HD22	1:A:927:PRO:HG3	1.78	0.65
1:A:748:GLU:HG3	1:A:817:MET:HG2	1.78	0.64
1:A:129:VAL:HG22	1:A:140:ILE:HD12	1.78	0.64
1:A:230:THR:C	1:A:232:ILE:H	2.04	0.64
1:A:743:ILE:O	1:A:747:VAL:HG23	1.98	0.64
1:A:880:HIS:N	1:A:881:PRO:HD2	2.12	0.64
1:A:770:GLY:HA2	1:A:773:VAL:HG22	1.79	0.64
1:A:421:ASN:HD21	1:A:423:SER:HB2	1.61	0.64
1:A:94:ILE:HD12	1:A:304:VAL:HG11	1.79	0.63
1:A:108:GLN:HG2	1:A:317:THR:HG23	1.80	0.63
1:A:883:PHE:HB3	1:A:886:LEU:HD12	1.80	0.63
1:A:90:GLU:HG2	1:A:790:VAL:HG13	1.80	0.62
1:A:604:ARG:HB2	1:A:607:VAL:HG23	1.81	0.62
1:A:352:LYS:HE2	1:A:353:THR:OG1	1.98	0.62
1:A:181:THR:OG1	1:A:183:GLU:HG2	2.00	0.62
1:A:629:LYS:HA	1:A:677:ALA:CB	2.30	0.62
1:A:798:VAL:HG13	1:A:940:SER:HB3	1.82	0.61
1:A:115:ALA:HB2	1:A:324:ARG:NH1	2.15	0.61
1:A:939:LEU:O	1:A:943:LEU:HB2	2.01	0.61
1:A:983:ILE:O	1:A:987:ILE:HG12	2.01	0.61
1:A:366:MET:HE2	1:A:384:ILE:HD11	1.84	0.60
1:A:261:SER:O	1:A:265:SER:HB2	2.01	0.60
1:A:367:PHE:C	1:A:367:PHE:CD2	2.80	0.60
1:A:389:TYR:HA	1:A:447:THR:HG21	1.84	0.60
1:A:793:LEU:O	1:A:796:ASN:HB3	2.02	0.59
1:A:94:ILE:CD1	1:A:793:LEU:HD22	2.32	0.59
1:A:947:ILE:HD13	1:A:953:LEU:HD23	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:762:ARG:HH21	1:A:836:ARG:HH12	1.48	0.59
1:A:94:ILE:HD11	1:A:793:LEU:HD13	1.85	0.59
1:A:762:ARG:HG2	1:A:837:TYR:CE1	2.38	0.59
1:A:762:ARG:NH2	1:A:836:ARG:HH12	2.00	0.58
1:A:311:LEU:HD13	1:A:761:ILE:HD11	1.86	0.58
1:A:302:LEU:HD13	1:A:775:ILE:HD12	1.85	0.58
1:A:287:SER:HB2	1:A:290:ARG:HB2	1.86	0.58
1:A:280:ASN:HD21	1:A:288:TRP:HA	1.69	0.58
1:A:446:THR:HG23	1:A:472:ASN:ND2	2.19	0.57
1:A:368:ILE:HD12	1:A:409:GLY:HA3	1.86	0.57
1:A:162:ASP:OD2	1:A:230:THR:HA	2.04	0.57
1:A:258:GLU:HA	1:A:261:SER:HB3	1.87	0.57
1:A:491:ARG:NH1	1:A:588:GLU:OE1	2.38	0.57
1:A:155:VAL:HG22	1:A:217:GLY:N	2.20	0.56
1:A:708:ALA:HB3	1:A:709:PRO:HD3	1.88	0.56
1:A:354:GLY:HA2	1:A:359:ASN:HB2	1.87	0.56
1:A:326:MET:HE2	1:A:746:ALA:HB2	1.88	0.55
1:A:792:LEU:O	1:A:795:VAL:HG12	2.05	0.55
1:A:963:ASP:HB2	1:A:966:GLN:HG3	1.88	0.55
1:A:683:HIS:O	1:A:687:ILE:HG12	2.07	0.55
1:A:352:LYS:NZ	1:A:627:ASP:OD2	2.40	0.55
1:A:366:MET:HE2	1:A:384:ILE:CD1	2.37	0.55
1:A:38:HIS:CD2	1:A:143:ARG:HH21	2.24	0.55
1:A:412:GLU:OE2	1:A:529:ARG:NH1	2.39	0.55
1:A:491:ARG:HD2	1:A:493:SER:OG	2.08	0.54
1:A:416:ILE:HD11	1:A:566:THR:CG2	2.37	0.54
1:A:856:PHE:HD1	1:A:870:LEU:HD21	1.73	0.54
1:A:357:THR:HA	1:A:603:PRO:HA	1.89	0.54
1:A:530:VAL:HB	1:A:533:THR:HG22	1.90	0.54
1:A:681:PRO:HG3	1:A:706:ASN:HB2	1.90	0.54
1:A:965:THR:HG23	1:A:968:LEU:HD12	1.88	0.54
1:A:788:ILE:HB	1:A:789:PRO:HD2	1.90	0.54
1:A:1:MET:HG2	1:A:4:ALA:HB2	1.89	0.53
1:A:102:ALA:O	1:A:106:VAL:HG12	2.09	0.53
1:A:156:GLY:HA2	1:A:726:VAL:CG1	2.38	0.53
1:A:165:ILE:HG22	1:A:191:THR:HG22	1.90	0.53
1:A:267:ILE:HG22	1:A:776:PHE:CE1	2.43	0.53
1:A:760:PHE:HA	1:A:807:LEU:HD23	1.89	0.53
1:A:5:HIS:HD2	1:A:194:VAL:HG13	1.74	0.53
1:A:155:VAL:HG22	1:A:217:GLY:H	1.74	0.53
1:A:880:HIS:C	1:A:882:HIS:H	2.17	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:ASN:ND2	1:A:442:GLU:HB3	2.23	0.53
1:A:201:ASN:ND2	1:A:231:GLU:OE2	2.42	0.53
1:A:813:ASP:HA	1:A:925:MET:HE1	1.91	0.53
1:A:529:ARG:HH22	1:A:568:ASP:CG	2.17	0.53
1:A:312:PRO:O	1:A:315:ILE:HG22	2.09	0.52
1:A:518:PRO:HB3	1:A:549:ILE:HD13	1.92	0.52
1:A:308:PRO:CB	1:A:764:LEU:HD23	2.32	0.52
1:A:834:PHE:O	1:A:838:MET:HG2	2.10	0.52
1:A:952:PRO:O	1:A:956:ILE:HG13	2.10	0.52
1:A:271:VAL:HA	1:A:274:ILE:HG12	1.91	0.52
1:A:397:LYS:CB	1:A:402:ILE:HD12	2.40	0.51
1:A:123:GLU:OE1	1:A:124:PRO:HD2	2.11	0.51
1:A:210:SER:HB3	1:A:232:ILE:HD13	1.93	0.51
1:A:352:LYS:CB	1:A:623:MET:HE2	2.36	0.51
1:A:604:ARG:HB2	1:A:607:VAL:CG2	2.40	0.51
1:A:843:TYR:OH	1:A:976:PRO:HG2	2.10	0.51
1:A:264:ILE:HD12	1:A:306:ALA:HB1	1.93	0.51
1:A:811:PRO:HG3	1:A:929:VAL:CG1	2.40	0.51
1:A:629:LYS:HD2	1:A:654:THR:CG2	2.41	0.51
1:A:802:LEU:HD12	1:A:939:LEU:HD23	1.93	0.51
1:A:491:ARG:HD2	1:A:588:GLU:OE2	2.11	0.51
1:A:520:GLY:O	1:A:524:ARG:HG3	2.11	0.51
1:A:1:MET:HE2	1:A:16:PHE:CE1	2.46	0.51
1:A:230:THR:C	1:A:232:ILE:N	2.70	0.50
1:A:589:THR:HG22	1:A:590:ASP:N	2.27	0.50
1:A:739:ASN:HD21	1:A:741:SER:HB2	1.77	0.49
1:A:491:ARG:HH11	1:A:588:GLU:CD	2.20	0.49
1:A:895:GLU:HG2	1:A:960:LYS:HD3	1.94	0.49
1:A:895:GLU:N	1:A:896:PRO:HD2	2.27	0.49
1:A:979:GLY:O	1:A:983:ILE:HG12	2.12	0.49
1:A:276:ILE:HA	1:A:279:PHE:CE2	2.47	0.49
1:A:748:GLU:HG3	1:A:817:MET:CG	2.43	0.49
1:A:342:LEU:HD22	1:A:747:VAL:HG22	1.93	0.49
1:A:335:SER:OG	1:A:337:PRO:HD2	2.12	0.49
1:A:589:THR:HG22	1:A:590:ASP:H	1.77	0.49
1:A:318:CYS:SG	1:A:757:MET:HE3	2.53	0.49
1:A:533:THR:HG23	1:A:534:ARG:N	2.28	0.49
1:A:624:ILE:HG22	1:A:684:LYS:HG2	1.91	0.49
1:A:651:ARG:HA	1:A:673:ALA:HA	1.94	0.49
1:A:773:VAL:CG2	1:A:845:GLY:HA3	2.37	0.49
1:A:975:LEU:N	1:A:976:PRO:HD2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ALA:HB2	1:A:742:THR:HG21	1.95	0.48
1:A:351:ASP:OD2	3:A:1002:ACP:O3G	2.30	0.48
1:A:716:ILE:HD11	1:A:733:MET:HE2	1.96	0.48
1:A:1:MET:HE1	1:A:12:CYS:HA	1.95	0.48
1:A:553:GLY:CA	1:A:631:THR:HG22	2.43	0.48
1:A:848:THR:HG21	1:A:904:LEU:HD13	1.95	0.48
1:A:497:TYR:CE2	1:A:576:MET:HE1	2.48	0.48
1:A:659:ASP:HA	1:A:666:GLN:NE2	2.28	0.48
1:A:75:LEU:CD2	1:A:297:LYS:HD3	2.41	0.48
1:A:336:LEU:N	1:A:337:PRO:CD	2.77	0.48
1:A:572:LYS:HB2	1:A:575:GLU:CG	2.43	0.48
1:A:830:SER:HB3	1:A:833:LEU:HB2	1.95	0.48
1:A:89:VAL:HG21	1:A:956:ILE:HG21	1.95	0.47
1:A:97:ILE:HD11	1:A:800:ASP:O	2.13	0.47
1:A:524:ARG:HB2	1:A:591:LEU:HD12	1.96	0.47
1:A:854:TRP:O	1:A:859:ALA:HA	2.14	0.47
1:A:287:SER:HB2	1:A:290:ARG:CB	2.44	0.47
1:A:246:LYS:HE3	1:A:251:GLN:HE21	1.79	0.47
1:A:404:SER:C	1:A:406:GLN:H	2.21	0.47
1:A:412:GLU:CD	1:A:529:ARG:HH11	2.22	0.47
1:A:770:GLY:HA2	1:A:841:GLY:O	2.14	0.47
1:A:558:THR:HG22	1:A:634:ALA:HB1	1.96	0.47
1:A:757:MET:HE2	1:A:757:MET:HA	1.97	0.47
1:A:887:ASP:OD2	1:A:889:GLU:HB2	2.15	0.47
1:A:689:GLU:HG3	1:A:713:LYS:HE3	1.95	0.47
1:A:445:LEU:HD13	1:A:599:MET:HE1	1.96	0.47
1:A:903:VAL:HG22	1:A:970:VAL:HA	1.96	0.47
1:A:358:THR:HG23	1:A:602:PRO:O	2.15	0.47
1:A:23:GLY:HA3	1:A:130:TYR:O	2.15	0.46
1:A:247:THR:HG23	1:A:250:GLN:HE21	1.80	0.46
1:A:352:LYS:CE	1:A:353:THR:OG1	2.62	0.46
1:A:24:LEU:HD22	1:A:149:ASP:HB3	1.96	0.46
1:A:360:GLN:HE21	1:A:360:GLN:HA	1.79	0.46
1:A:7:LYS:HB2	1:A:7:LYS:NZ	2.31	0.46
1:A:389:TYR:O	1:A:425:LEU:HD21	2.15	0.46
1:A:880:HIS:N	1:A:881:PRO:CD	2.79	0.46
1:A:974:SER:C	1:A:976:PRO:HD2	2.41	0.46
1:A:981:ASP:O	1:A:985:LYS:N	2.45	0.46
1:A:140:ILE:H	1:A:140:ILE:HD13	1.79	0.46
1:A:425:LEU:HD22	1:A:447:THR:HG22	1.98	0.46
1:A:155:VAL:HA	1:A:214:ILE:HG22	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:527:TYR:O	1:A:592:THR:HA	2.16	0.46
1:A:1:MET:CG	1:A:4:ALA:HB2	2.45	0.45
1:A:232:ILE:N	1:A:232:ILE:HD12	2.32	0.45
1:A:624:ILE:O	1:A:684:LYS:HE2	2.15	0.45
1:A:530:VAL:O	1:A:533:THR:HB	2.15	0.45
1:A:390:ALA:C	1:A:392:GLU:H	2.25	0.45
1:A:629:LYS:HD2	1:A:654:THR:HG23	1.98	0.45
1:A:681:PRO:HG3	1:A:706:ASN:CB	2.47	0.45
1:A:913:LEU:O	1:A:922:LEU:HD21	2.16	0.45
1:A:366:MET:HE1	1:A:452:MET:SD	2.56	0.45
1:A:389:TYR:CA	1:A:447:THR:HG21	2.47	0.45
1:A:526:ASN:HD22	1:A:590:ASP:HA	1.82	0.45
1:A:816:ILE:HG23	1:A:817:MET:N	2.32	0.45
1:A:911:ASN:O	1:A:914:ASN:HB2	2.16	0.45
1:A:804:ALA:O	1:A:807:LEU:HB2	2.16	0.45
1:A:119:LEU:HD11	1:A:330:ASN:HA	1.99	0.45
1:A:491:ARG:HG3	1:A:585:MET:HB2	1.98	0.45
1:A:620:ARG:NH2	1:A:671:ARG:HA	2.32	0.45
1:A:237:ASP:C	1:A:239:MET:H	2.25	0.44
1:A:662:PRO:HG2	1:A:665:GLU:HG2	1.99	0.44
1:A:957:PHE:O	1:A:959:LEU:HG	2.17	0.44
1:A:19:SER:HB2	1:A:22:THR:HB	1.98	0.44
1:A:737:ASP:O	1:A:738:ASP:HB2	2.17	0.44
1:A:33:LEU:HD12	1:A:146:VAL:CG1	2.47	0.44
1:A:48:SER:C	1:A:50:TRP:H	2.26	0.44
1:A:283:VAL:HG13	1:A:284:HIS:H	1.82	0.44
1:A:482:GLU:HG3	1:A:497:TYR:HD1	1.81	0.44
1:A:315:ILE:HG23	1:A:316:THR:N	2.33	0.44
1:A:397:LYS:HB3	1:A:402:ILE:HD12	2.00	0.44
1:A:501:ALA:HB3	1:A:505:ARG:HG2	2.00	0.44
1:A:606:GLU:H	1:A:606:GLU:CD	2.26	0.44
1:A:128:LYS:NZ	1:A:157:ASP:OD2	2.33	0.44
1:A:978:ILE:O	1:A:982:GLU:HB2	2.18	0.44
1:A:49:LEU:HD22	1:A:110:ARG:NE	2.33	0.44
1:A:231:GLU:HA	1:A:234:LYS:HG2	1.98	0.44
1:A:267:ILE:O	1:A:271:VAL:HG23	2.18	0.44
1:A:353:THR:HG1	3:A:1002:ACP:PG	2.40	0.44
1:A:624:ILE:HG22	1:A:679:VAL:HG11	2.00	0.44
1:A:0:ACE:H3	1:A:36:TYR:CE1	2.52	0.43
1:A:246:LYS:HE3	1:A:251:GLN:NE2	2.33	0.43
1:A:273:LEU:C	1:A:275:ASN:H	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:THR:HA	1:A:357:THR:OG1	2.18	0.43
1:A:479:MET:HE3	1:A:498:CYS:HB3	2.00	0.43
1:A:397:LYS:HB2	1:A:402:ILE:HD12	2.00	0.43
1:A:762:ARG:HH11	1:A:762:ARG:HB3	1.84	0.43
1:A:797:LEU:HD23	1:A:798:VAL:HG23	2.00	0.43
1:A:94:ILE:HD11	1:A:793:LEU:CD2	2.42	0.43
1:A:96:LEU:HB3	1:A:797:LEU:CD1	2.49	0.43
1:A:342:LEU:HG	1:A:716:ILE:CD1	2.35	0.43
1:A:553:GLY:HA2	1:A:631:THR:HG22	2.01	0.43
1:A:530:VAL:HB	1:A:533:THR:CG2	2.48	0.43
1:A:880:HIS:O	1:A:882:HIS:N	2.52	0.43
1:A:1:MET:CE	1:A:12:CYS:HA	2.49	0.43
1:A:879:ASP:HB3	1:A:882:HIS:CE1	2.54	0.43
1:A:255:GLU:C	1:A:257:GLY:H	2.26	0.43
1:A:508:VAL:HG12	1:A:508:VAL:O	2.19	0.43
1:A:567:ARG:HD3	1:A:570:PRO:CA	2.33	0.43
1:A:633:ILE:O	1:A:637:ARG:HG3	2.19	0.43
1:A:783:LEU:HB3	1:A:784:PRO:HD2	1.99	0.42
1:A:799:THR:HG22	1:A:909:MET:HE3	2.01	0.42
1:A:529:ARG:NH2	1:A:568:ASP:OD1	2.52	0.42
1:A:707:ASP:O	1:A:711:LEU:HD12	2.19	0.42
1:A:795:VAL:O	1:A:799:THR:OG1	2.37	0.42
1:A:836:ARG:HD3	1:A:984:LEU:HD13	2.00	0.42
1:A:361:MET:CE	1:A:599:MET:SD	3.01	0.42
1:A:52:LEU:HB3	1:A:106:VAL:HG23	2.01	0.42
1:A:195:PRO:O	1:A:197:PRO:HD3	2.20	0.42
1:A:267:ILE:HA	1:A:270:ALA:HB3	2.02	0.42
1:A:459:VAL:HB	1:A:467:ARG:CD	2.49	0.42
1:A:628:ASN:OD1	1:A:630:GLY:N	2.51	0.42
1:A:367:PHE:CD2	1:A:367:PHE:O	2.73	0.42
1:A:394:GLU:OE1	1:A:401:PRO:HB3	2.19	0.42
1:A:677:ALA:O	1:A:678:ARG:C	2.63	0.42
1:A:784:PRO:O	1:A:785:GLU:C	2.63	0.42
1:A:964:LEU:H	1:A:964:LEU:HD12	1.84	0.42
1:A:1:MET:HE2	1:A:16:PHE:HE1	1.83	0.42
1:A:535:VAL:HB	1:A:536:PRO:HD2	2.00	0.42
1:A:879:ASP:C	1:A:881:PRO:HD2	2.44	0.41
1:A:553:GLY:O	1:A:558:THR:HA	2.20	0.41
1:A:749:GLU:O	1:A:753:ILE:HG12	2.21	0.41
1:A:255:GLU:C	1:A:257:GLY:N	2.76	0.41
1:A:529:ARG:NH2	1:A:592:THR:HG21	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:699:ALA:HA	1:A:716:ILE:O	2.20	0.41
1:A:119:LEU:CD1	1:A:330:ASN:HA	2.51	0.41
1:A:452:MET:O	1:A:453:ASN:C	2.63	0.41
1:A:533:THR:CG2	1:A:534:ARG:N	2.83	0.41
1:A:420:CYS:SG	1:A:494:MET:HG2	2.61	0.41
1:A:811:PRO:HG3	1:A:929:VAL:HG13	2.03	0.41
1:A:904:LEU:O	1:A:907:ILE:HG22	2.21	0.41
1:A:38:HIS:CG	1:A:143:ARG:HH21	2.38	0.41
1:A:903:VAL:HA	1:A:970:VAL:HG13	2.03	0.41
1:A:97:ILE:CG2	1:A:98:LEU:N	2.83	0.41
1:A:321:LEU:HD22	1:A:325:ARG:CZ	2.51	0.41
1:A:626:GLY:O	1:A:678:ARG:HA	2.21	0.41
1:A:86:THR:O	1:A:86:THR:HG22	2.19	0.41
1:A:263:VAL:O	1:A:267:ILE:HG13	2.20	0.41
1:A:279:PHE:C	1:A:280:ASN:HD22	2.29	0.41
1:A:326:MET:CE	1:A:746:ALA:HB2	2.50	0.41
1:A:788:ILE:HB	1:A:789:PRO:CD	2.51	0.41
1:A:825:LYS:O	1:A:826:GLU:C	2.63	0.41
1:A:855:TRP:HH2	1:A:863:PRO:HD2	1.85	0.41
1:A:624:ILE:CG2	1:A:679:VAL:HG11	2.51	0.41
1:A:721:GLY:HA3	1:A:736:ALA:HA	2.03	0.41
1:A:89:VAL:O	1:A:93:VAL:HG23	2.21	0.40
1:A:832:TRP:CE2	1:A:988:ALA:HB2	2.56	0.40
1:A:858:TYR:O	1:A:860:GLU:N	2.54	0.40
1:A:867:TYR:CE2	1:A:868:HIS:ND1	2.89	0.40
1:A:918:GLU:HG2	1:A:919:ASN:ND2	2.36	0.40
1:A:189:LYS:HD3	1:A:189:LYS:HA	1.87	0.40
1:A:235:ILE:O	1:A:239:MET:HB2	2.22	0.40
1:A:794:TRP:CZ2	1:A:947:ILE:HG13	2.57	0.40
1:A:898:THR:HG21	1:A:960:LYS:O	2.21	0.40
1:A:52:LEU:HD11	1:A:109:GLU:HG2	2.04	0.40
1:A:177:GLN:NE2	1:A:187:VAL:HG21	2.35	0.40
1:A:362:SER:HA	6:A:4506:HOH:O	2.20	0.40
1:A:496:VAL:O	1:A:512:MET:HA	2.21	0.40
1:A:800:ASP:OD2	6:A:3401:HOH:O	2.22	0.40
1:A:398:ASN:O	1:A:399:ASP:HB2	2.20	0.40
1:A:204:LYS:C	1:A:206:ASN:H	2.28	0.40
1:A:294:TYR:CZ	1:A:298:ILE:HD11	2.56	0.40
1:A:528:VAL:HG11	1:A:541:VAL:CG1	2.39	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%)	878 (88%)	98 (10%)	17 (2%)	7 13

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	857	MET
1	A	951	ASP
1	A	245	ASP
1	A	859	ALA
1	A	243	GLU
1	A	338	SER
1	A	501	ALA
1	A	502	LYS
1	A	868	HIS
1	A	231	GLU
1	A	236	ARG
1	A	82	GLU
1	A	453	ASN
1	A	456	ASN
1	A	199	ALA
1	A	405	GLY
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%)	790 (94%)	50 (6%)	17	36

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	VAL
1	A	24	LEU
1	A	33	LEU
1	A	49	LEU
1	A	80	GLU
1	A	95	LEU
1	A	97	ILE
1	A	98	LEU
1	A	140	ILE
1	A	164	ARG
1	A	267	ILE
1	A	276	ILE
1	A	316	THR
1	A	319	LEU
1	A	321	LEU
1	A	352	LYS
1	A	359	ASN
1	A	360	GLN
1	A	384	ILE
1	A	402	ILE
1	A	413	LEU
1	A	445	LEU
1	A	449	VAL
1	A	465	VAL
1	A	467	ARG
1	A	484	THR
1	A	494	MET
1	A	496	VAL
1	A	528	VAL
1	A	533	THR
1	A	534	ARG
1	A	543	GLU
1	A	600	LEU
1	A	623	MET
1	A	663	LEU
1	A	678	ARG
1	A	679	VAL
1	A	726	VAL

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Mol	Chain	Res	Type
1	A	749	GLU
1	A	759	GLN
1	A	762	ARG
1	A	764	LEU
1	A	769	VAL
1	A	807	LEU
1	A	925	MET
1	A	931	ILE
1	A	953	LEU
1	A	955	MET
1	A	973	ILE
1	A	975	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	GLN
1	A	108	GLN
1	A	111	ASN
1	A	114	ASN
1	A	201	ASN
1	A	202	GLN
1	A	250	GLN
1	A	251	GLN
1	A	280	ASN
1	A	359	ASN
1	A	360	GLN
1	A	380	ASN
1	A	421	ASN
1	A	526	ASN
1	A	755	ASN
1	A	869	GLN
1	A	875	GLN
1	A	914	ASN
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 [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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$
5	PC1	A	1011	-	19,19,53	1.96	3 (15%)	21,25,61	1.34	3 (14%)
3	ACP	A	1002	2	31,33,33	2.75	10 (32%)	47,52,52	2.09	14 (29%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PC1	A	1011	-	-	4/21/21/57	-
3	ACP	A	1002	2	-	0/19/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	ACP	PA-O3A	9.57	1.69	1.59
5	A	1011	PC1	O21-C21	6.93	1.47	1.33
3	A	1002	ACP	PG-O1G	6.81	1.64	1.50
3	A	1002	ACP	PB-O3A	5.83	1.64	1.58
3	A	1002	ACP	C5-C4	4.26	1.46	1.39
5	A	1011	PC1	O31-C31	3.73	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	ACP	C5-C6	2.74	1.48	1.41
3	A	1002	ACP	C4-N9	-2.48	1.32	1.37
3	A	1002	ACP	PB-O2B	-2.26	1.50	1.56
3	A	1002	ACP	PB-O1B	2.11	1.56	1.51
3	A	1002	ACP	C5-N7	-2.10	1.35	1.39
5	A	1011	PC1	O21-C2	-2.09	1.43	1.46
3	A	1002	ACP	C8-N7	2.08	1.35	1.31

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1002	ACP	C5-C4-N3	-5.72	118.83	126.72
3	A	1002	ACP	O1B-PB-C3B	4.85	121.82	109.05
3	A	1002	ACP	N3-C4-N9	4.58	134.96	127.17
3	A	1002	ACP	N3-C2-N1	-4.27	122.11	128.58
3	A	1002	ACP	C2-N3-C4	4.14	121.93	111.83
5	A	1011	PC1	O21-C21-O22	-4.06	120.49	125.70
5	A	1011	PC1	C3-C2-C1	-3.27	104.16	111.78
3	A	1002	ACP	C4-C5-N7	-3.18	106.94	110.58
3	A	1002	ACP	C4-N9-C8	2.98	108.87	105.74
3	A	1002	ACP	PB-O3A-PA	-2.62	123.82	132.37
3	A	1002	ACP	O2A-PA-O1A	2.51	124.14	112.44
3	A	1002	ACP	O2G-PG-O1G	-2.42	106.12	112.39
5	A	1011	PC1	O31-C3-C2	2.29	114.99	108.40
3	A	1002	ACP	C6-C5-N7	2.24	136.40	132.09
3	A	1002	ACP	C5-N7-C8	2.12	106.78	103.45
3	A	1002	ACP	O3G-PG-O2G	2.05	113.82	107.96
3	A	1002	ACP	O2'-C2'-C1'	-2.01	103.19	110.10

There are no chirality outliers.

All (4) torsion outliers are listed below:

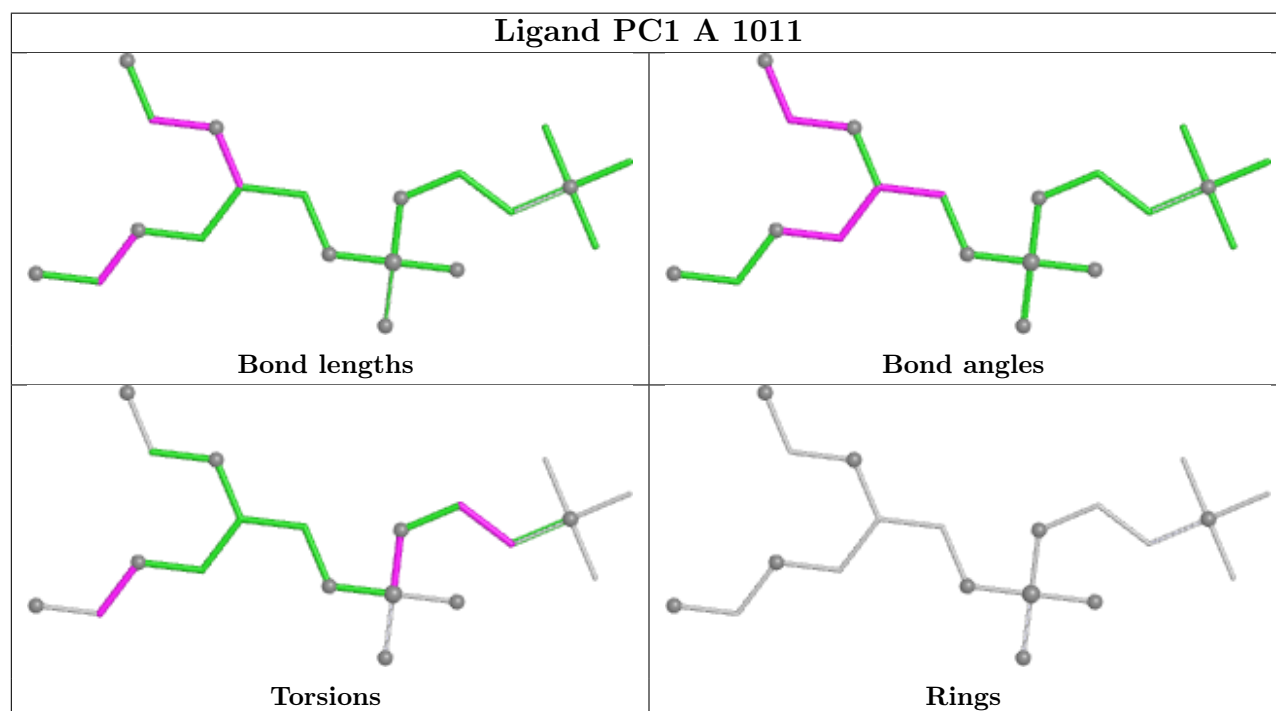
Mol	Chain	Res	Type	Atoms
5	A	1011	PC1	C11-O13-P-O12
5	A	1011	PC1	O13-C11-C12-N
5	A	1011	PC1	O32-C31-O31-C3
5	A	1011	PC1	C11-O13-P-O11

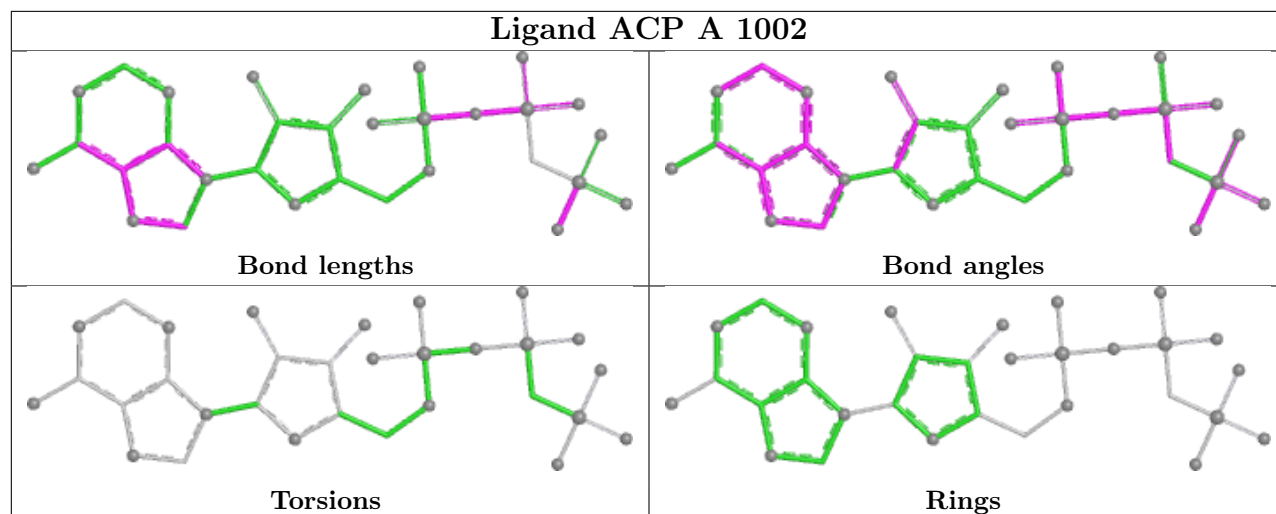
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	ACP	2	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%)	0.21	29 (2%) 53 49	48, 111, 207, 257	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	992	LEU	3.8
1	A	77	TRP	3.6
1	A	506	ALA	3.4
1	A	48	SER	3.0
1	A	288	TRP	2.9
1	A	951	ASP	2.9
1	A	508	VAL	2.8
1	A	240	ALA	2.8
1	A	865	VAL	2.7
1	A	46	GLY	2.6
1	A	964	LEU	2.6
1	A	870	LEU	2.5
1	A	52	LEU	2.5
1	A	886	LEU	2.5
1	A	199	ALA	2.5
1	A	507	ALA	2.4
1	A	284	HIS	2.4
1	A	860	GLU	2.3
1	A	235	ILE	2.3
1	A	273	LEU	2.2
1	A	503	SER	2.2
1	A	50	TRP	2.1
1	A	891	PHE	2.1
1	A	952	PRO	2.1
1	A	783	LEU	2.1
1	A	402	ILE	2.1
1	A	894	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	965	THR	2.0
1	A	239	MET	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

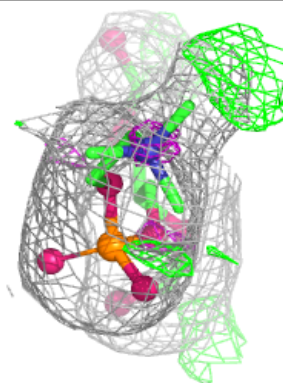
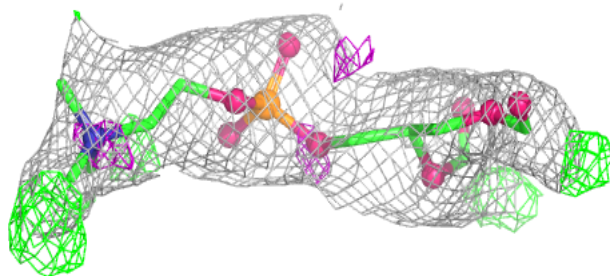
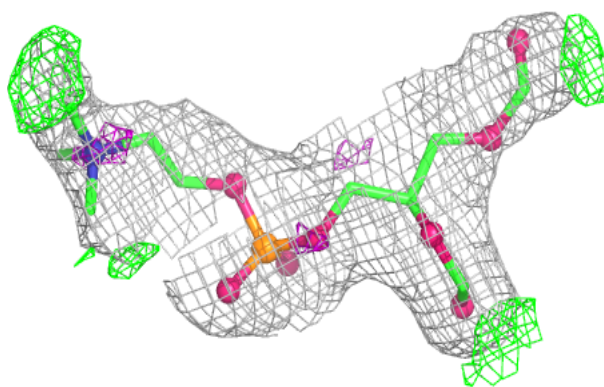
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	PC1	A	1011	20/54	0.77	0.15	116,118,124,124	0
4	NA	A	1000	1/1	0.87	0.09	74,74,74,74	0
3	ACP	A	1002	31/31	0.94	0.09	44,55,60,65	0
2	CA	A	1001	1/1	0.96	0.04	67,67,67,67	0
2	CA	A	996	1/1	0.97	0.03	85,85,85,85	0
2	CA	A	995	1/1	0.99	0.02	87,87,87,87	0

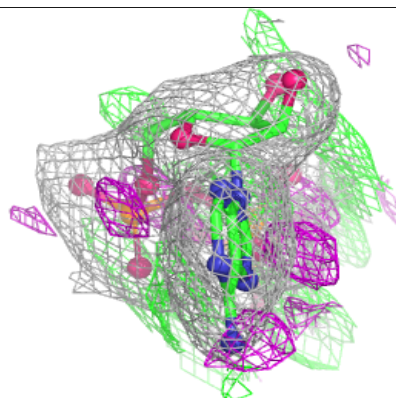
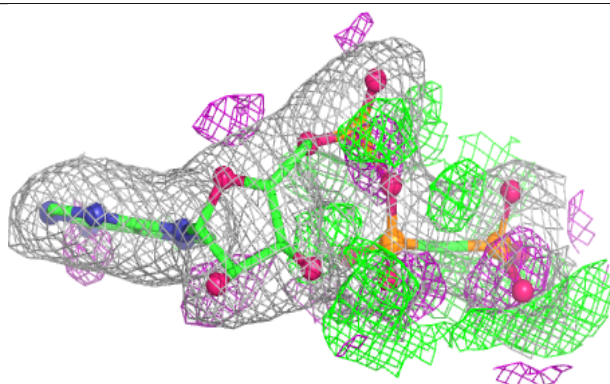
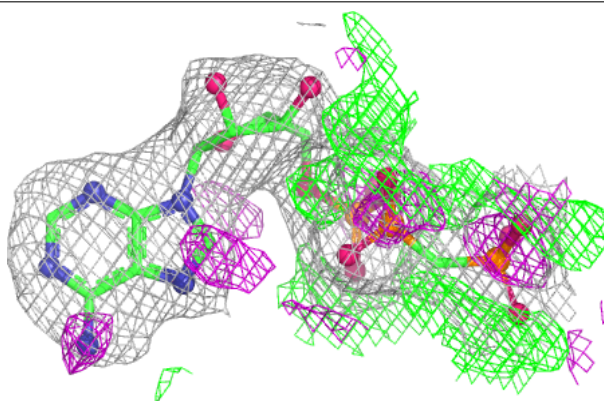
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PC1 A 1011:

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

**Electron density around ACP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.