



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 04:34 AM UTC

PDB ID : 1A0Q / pdb_00001a0q
Title : 29G11 COMPLEXED WITH PHENYL [1-(1-N-SUCCINYLAMINO)PENTYL] PHOSPHONATE
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Deposited on : 1997-12-05
Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

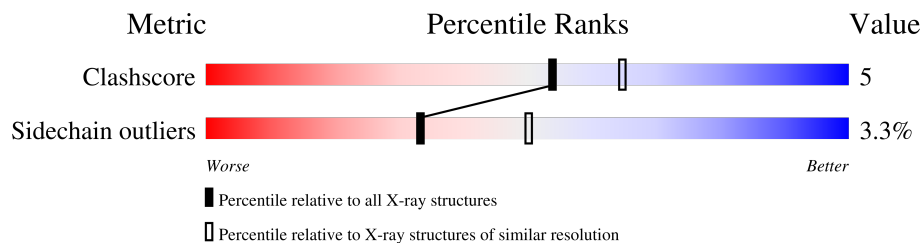
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	190562	6919 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	212	
2	H	97	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3301 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 29G11 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	211	Total	C	N	O	S	0	0	0
			1643	1032	278	328	5			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	4	LEU	MET	conflict	GB 12002896
L	22	THR	ALA	conflict	GB 12002896
L	26	SER	ASN	conflict	GB 12002896
L	30	LYS	ASN	conflict	GB 12002896
L	34	GLY	ALA	conflict	GB 12002896
L	43	GLN	GLY	conflict	GB 12002896
L	55	LEU	GLN	conflict	GB 12002896
L	63	ARG	SER	conflict	GB 12002896
L	92	TYR	ASP	conflict	GB 12002896
L	141	SER	PRO	conflict	GB 12002896
L	187	GLU	GLY	conflict	GB 12002896

- Molecule 2 is a protein called 29G11 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	205	Total	C	N	O	S	0	0	0
			1540	977	249	308	6			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	8	ASP	GLY	conflict	GB 3399661
H	9	ALA	PRO	conflict	GB 3399661
H	20	ILE	MET	conflict	GB 3399661
H	31	ASP	SER	conflict	GB 3399661
H	32	HIS	TYR	conflict	GB 3399661

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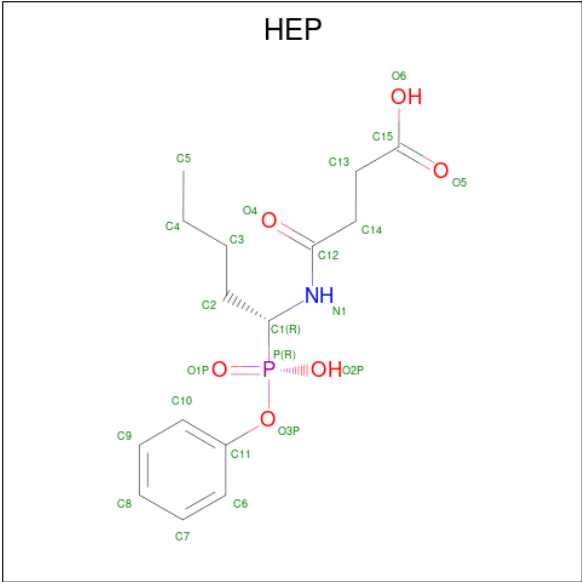
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Chain	Residue	Modelled	Actual	Comment	Reference
H	34	ILE	MET	conflict	GB 3399661
H	42	GLU	GLY	conflict	GB 3399661
H	52	SER	ASN	conflict	GB 3399661
H	53	GLY	TYR	conflict	GB 3399661
H	55	GLY	ASP	conflict	GB 3399661
H	56	ASP	GLY	conflict	GB 3399661
H	57	ILE	THR	conflict	GB 3399661
H	71	ALA	SER	conflict	GB 3399661
H	81	GLN	GLU	conflict	GB 3399661
H	82A	ASN	SER	conflict	GB 3399661
H	91	LEU	TYR	conflict	GB 3399661
H	93	LYS	VAL	conflict	GB 3399661
H	?	-	GLY	deletion	GB 3399661
H	97	TYR	ARG	conflict	GB 3399661
H	98	GLY	PRO	conflict	GB 3399661
H	99	ARG	TYR	conflict	GB 3399661
H	100	SER	TYR	conflict	GB 3399661
H	100A	ASN	ALA	conflict	GB 3399661
H	100B	VAL	MET	conflict	GB 3399661
H	108	THR	SER	conflict	GB 3399661
H	109	LEU	VAL	conflict	GB 3399661

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	L	1	Total Zn 1 1	0	0
3	H	2	Total Zn 2 2	0	0

- Molecule 4 is PHENYL[1-(N-SUCCINYLAMINO)PENTYL]PHOSPHONATE (CCD ID: HEP) (formula: C₁₅H₂₂NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	H	1	Total	C	N	O	P	0	0
			23	15	1	6	1		

- Molecule 5 is water.

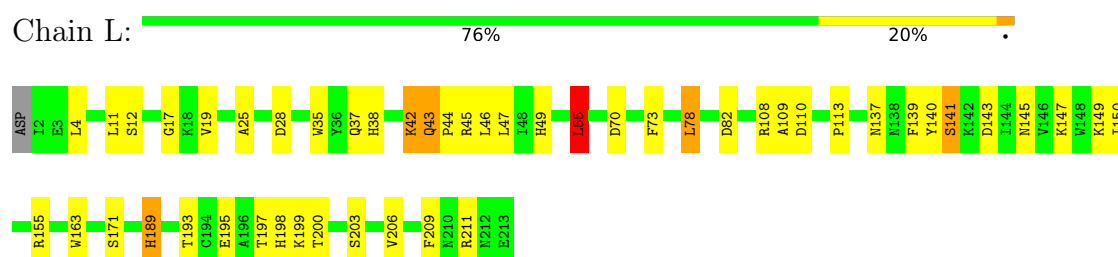
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	45	Total	O	0	0
			45	45		
5	H	47	Total	O	0	0
			47	47		

3 Residue-property plots [i](#)

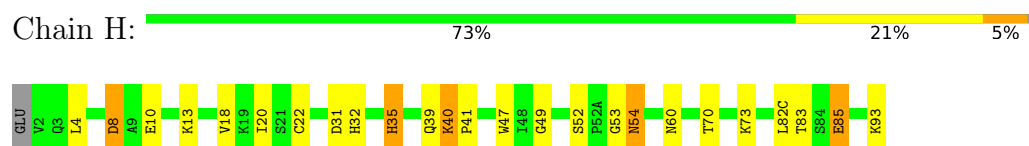
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: 29G11 FAB (LIGHT CHAIN)



- Molecule 2: 29G11 FAB (HEAVY CHAIN)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	36.78Å 82.61Å 132.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30	Depositor
% Data completeness (in resolution range)	93.0 (50.00-2.30)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	0.09	Depositor
Refinement program	X-PLOR 3.843	Depositor
R, R_{free}	0.203 , 0.275	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3301	wwPDB-VP
Average B, all atoms (Å ²)	25.0	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: HEP, ZN

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	L	0.82	4/1681 (0.2%)	1.54	27/2277 (1.2%)
2	H	0.83	4/1577 (0.3%)	1.53	15/2150 (0.7%)
All	All	0.83	8/3258 (0.2%)	1.54	42/4427 (0.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	L	0	1

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	49	HIS	CD2-NE2	-7.95	1.29	1.37
2	H	32	HIS	CD2-NE2	-7.26	1.29	1.37
1	L	189	HIS	CD2-NE2	-7.06	1.30	1.37
1	L	198	HIS	CD2-NE2	-6.85	1.30	1.37
2	H	35	HIS	CD2-NE2	-6.74	1.30	1.37

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	31	ASP	CA-CB-CG	9.80	122.40	112.60
1	L	141	SER	N-CA-C	-8.75	92.17	110.80
1	L	49	HIS	CB-CG-CD2	-8.33	120.37	131.20
1	L	140	TYR	O-C-N	7.88	132.88	123.27
1	L	137	ASN	N-CA-C	6.96	121.66	109.96

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	43	GLN	Peptide

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	L	1643	0	1590	18	0
2	H	1540	0	1503	13	0
3	H	2	0	0	0	0
3	L	1	0	0	0	0
4	H	23	0	20	1	0
5	H	47	0	0	0	0
5	L	45	0	0	0	0
All	All	3301	0	3113	31	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.

The worst 5 of 31 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:L:149:LYS:HB2	1:L:193:THR:HB	1.77	0.66
2:H:35:HIS:HD2	2:H:47:TRP:HE1	1.49	0.59
1:L:43:GLN:HG3	1:L:45:ARG:HB2	1.85	0.58
2:H:4:LEU:HG	2:H:22:CYS:SG	2.48	0.53
1:L:145:ASN:HB3	1:L:197:THR:HB	1.92	0.51

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	L	185/189 (98%)	180 (97%)	5 (3%)	39	58
2	H	175/83 (211%)	168 (96%)	7 (4%)	28	42
All	All	360/272 (132%)	348 (97%)	12 (3%)	33	50

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	H	70	THR
2	H	85	GLU
2	H	150	VAL
2	H	147	PRO
1	L	141	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	5	GLN
2	H	35	HIS
2	H	82(A)	ASN
1	L	138	ASN
1	L	137	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 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	$\# Z > 2$
4	HEP	H	214	-	22,23,23	2.06	4 (18%)	23,30,30	1.31	3 (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
4	HEP	H	214	-	-	3/24/24/24	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	214	HEP	P-C1	-7.43	1.74	1.84
4	H	214	HEP	P-O2P	-3.53	1.48	1.56
4	H	214	HEP	P-O3P	3.49	1.62	1.58
4	H	214	HEP	O6-C15	-2.17	1.23	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	214	HEP	O2P-P-O1P	3.34	119.73	111.58
4	H	214	HEP	O1P-P-C1	-3.22	109.80	114.19
4	H	214	HEP	O3P-P-O1P	-2.15	110.20	114.81

There are no chirality outliers.

All (3) torsion outliers are listed below:

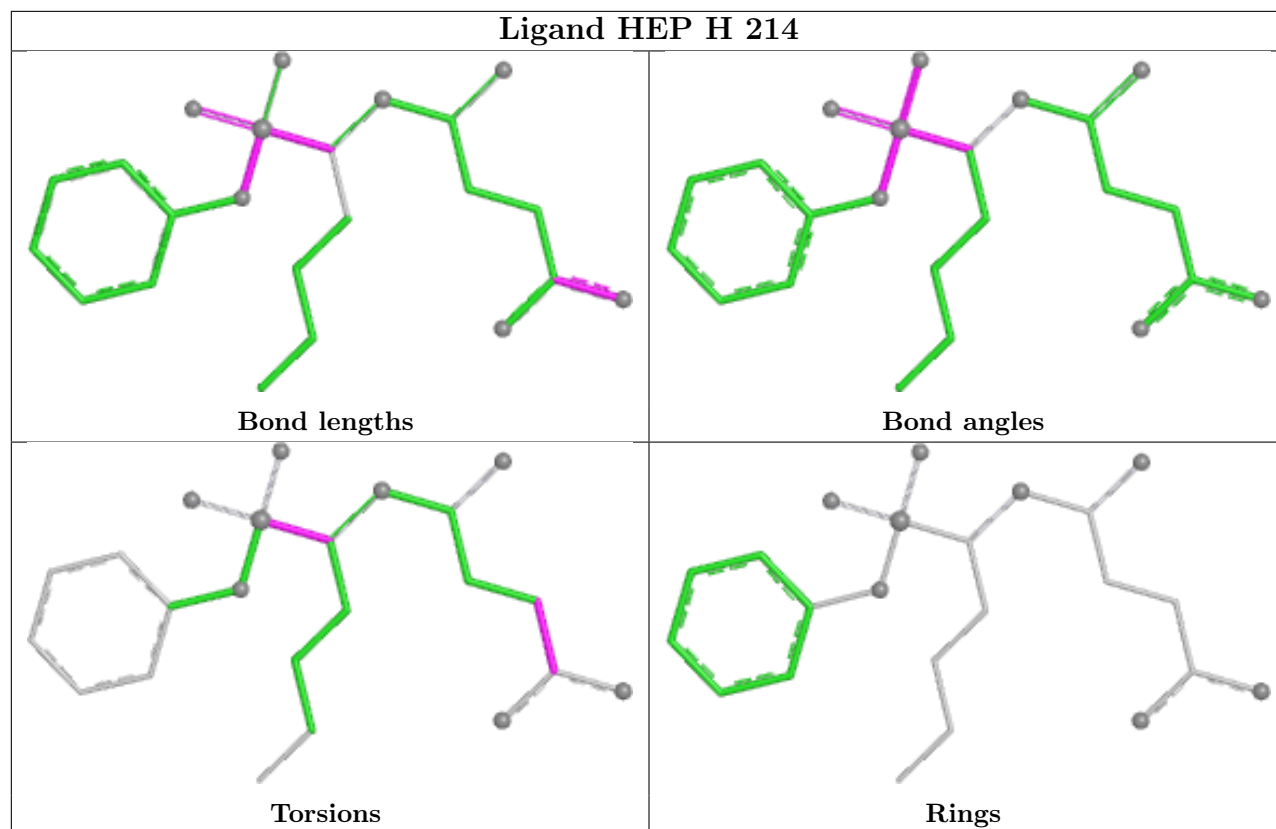
Mol	Chain	Res	Type	Atoms
4	H	214	HEP	C14-C13-C15-O6
4	H	214	HEP	C14-C13-C15-O5
4	H	214	HEP	N1-C1-P-O1P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	214	HEP	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.