



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:37 AM UTC

PDB ID : 2BU5 / pdb_00002bu5
Title : crystal structures of human pyruvate dehydrogenase kinase 2 containing physiological and synthetic ligands
Authors : Knoechel, T.R.; Tucker, A.D.; Robinson, C.M.; Phillips, C.; Taylor, W.; Bungay, P.J.; Kasten, S.A.; Roche, T.E.; Brown, D.G.
Deposited on : 2005-06-08
Resolution : 2.35 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)	:	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

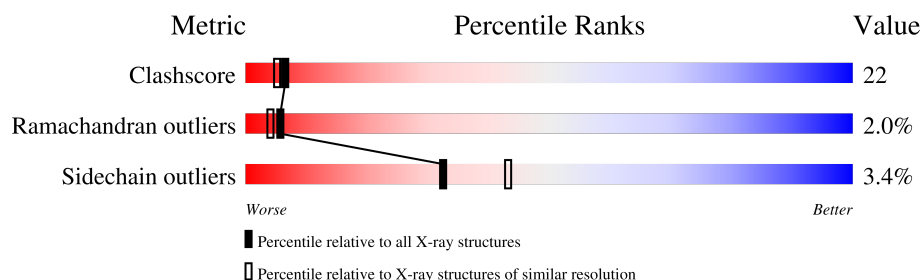
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)

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	A	394	 52% 34% • 10%

2 Entry composition [i](#)

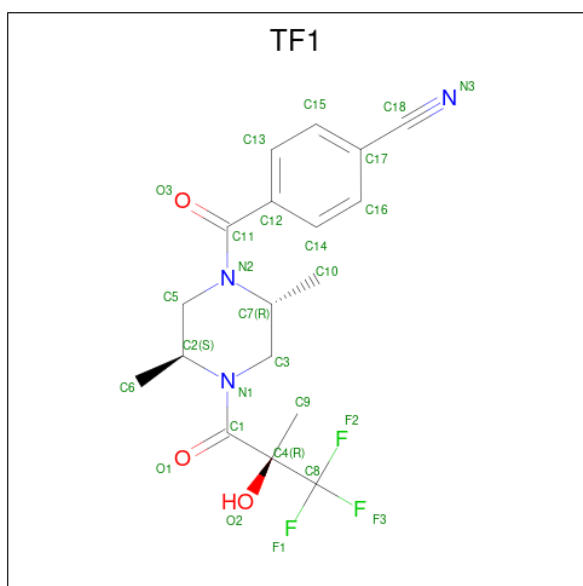
There are 3 unique types of molecules in this entry. The entry contains 2995 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PYRUVATE DEHYDROGENASE KINASE ISOENZYME 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	356	2860	1834	475	533	18	0	0	1

- Molecule 2 is 4-({(2R,5S)-2,5-DIMETHYL-4-[(2R)-3,3,3-TRIFLUORO-2-HYDROXY-2-METHYLPROPANOYL]PIPERAZIN-1-YL}CARBONYL)BENZONITRILE (CCD ID: TF1) (formula: C₁₈H₂₀F₃N₃O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
2	A	1	27	18	3	3	3	0	0

- Molecule 3 is water.

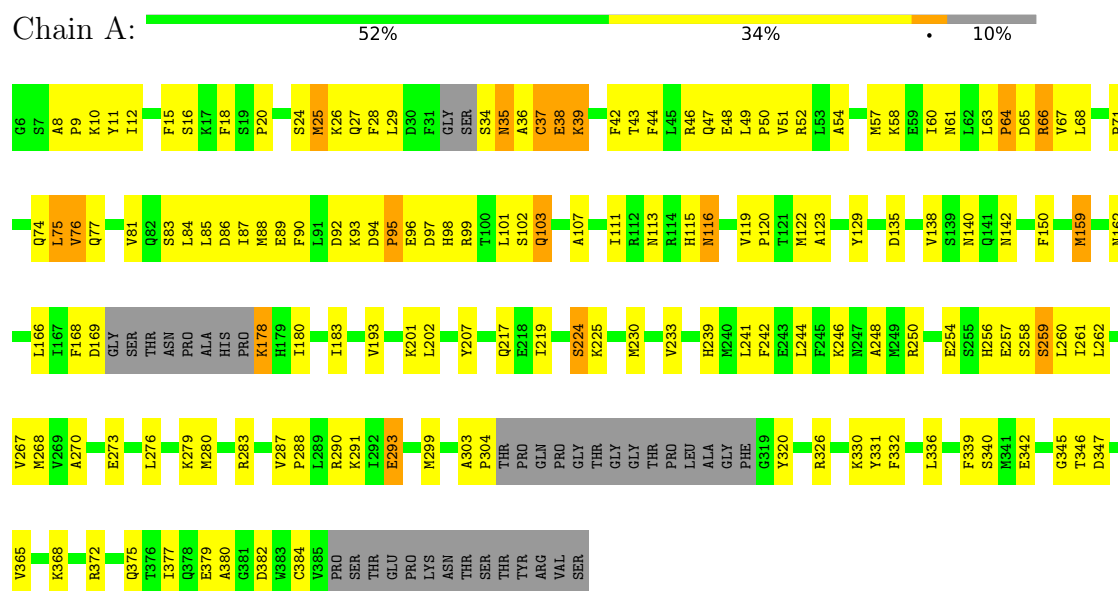
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	108	Total	O	0	0
			108	108		

3 Residue-property plots

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: PYRUVATE DEHYDROGENASE KINASE ISOENZYME 2



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	108.07 Å 108.07 Å 84.39 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.35	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-2.35)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNX	Depositor
R, R_{free}	0.237 , 0.255	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2995	wwPDB-VP
Average B, all atoms (Å ²)	65.0	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: TF1

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	0.65	5/2927 (0.2%)	1.00	12/3963 (0.3%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	260	LEU	CG-CD1	12.89	1.95	1.52
1	A	257	GLU	C-O	8.93	1.37	1.24
1	A	260	LEU	CB-CG	8.12	1.69	1.53
1	A	258	SER	CA-C	-7.45	1.43	1.53
1	A	384	CYS	C-N	-5.72	1.25	1.33

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	36	ALA	N-CA-C	-13.79	93.96	112.72
1	A	39	LYS	N-CA-C	-8.82	101.61	111.14
1	A	259	SER	N-CA-C	-8.54	96.36	109.85
1	A	258	SER	N-CA-C	-7.16	101.73	112.54
1	A	103	GLN	N-CA-C	-6.92	104.10	112.54
1	A	38	GLU	N-CA-C	-6.61	105.18	113.18
1	A	102	SER	N-CA-C	-6.43	105.26	113.23
1	A	260	LEU	N-CA-C	-5.74	105.95	113.12
1	A	368	LYS	N-CA-C	-5.54	105.78	112.54
1	A	116	ASN	N-CA-C	5.47	118.75	111.75
1	A	257	GLU	CA-C-N	-5.00	111.89	121.15
1	A	257	GLU	C-N-CA	-5.00	111.89	121.15

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	2860	0	2819	125	1
2	A	27	0	19	0	0
3	A	108	0	0	8	0
All	All	2995	0	2838	125	1

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 22.

All (125) 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:76:VAL:HG21	1:A:122:MET:HE2	1.39	1.00
1:A:273:GLU:HG2	3:A:2073:HOH:O	1.68	0.91
1:A:64:PRO:HB3	1:A:66:ARG:HD3	1.55	0.89
1:A:35:ASN:C	1:A:37:CYS:H	1.81	0.89
1:A:250:ARG:HD3	3:A:2064:HOH:O	1.77	0.84
1:A:244:LEU:C	1:A:280:MET:HE1	2.03	0.84
1:A:60:ILE:O	1:A:63:LEU:HG	1.80	0.82
1:A:244:LEU:O	1:A:280:MET:HE1	1.80	0.81
1:A:140:ASN:HB2	1:A:299:MET:HE2	1.64	0.77
1:A:303:ALA:HB1	1:A:304:PRO:HA	1.67	0.75
1:A:11:TYR:CG	1:A:88:MET:HE3	2.26	0.71
1:A:83:SER:OG	1:A:111:ILE:HG23	1.93	0.69
1:A:35:ASN:C	1:A:37:CYS:N	2.51	0.68
1:A:71:PRO:HB3	3:A:2010:HOH:O	1.93	0.68
1:A:372:ARG:HA	1:A:375:GLN:HG2	1.76	0.68
1:A:248:ALA:HB2	1:A:280:MET:HE3	1.76	0.68
1:A:24:SER:OG	1:A:27:GLN:HG3	1.94	0.67
1:A:47:GLN:O	1:A:51:VAL:HG23	1.95	0.66
1:A:268:MET:HE3	1:A:270:ALA:HB2	1.79	0.65
1:A:201:LYS:HE2	3:A:2039:HOH:O	1.96	0.64
1:A:63:LEU:HD13	1:A:67:VAL:HB	1.79	0.63
1:A:119:VAL:HB	1:A:120:PRO:HD3	1.81	0.63
1:A:248:ALA:CB	1:A:280:MET:HE3	2.30	0.61
1:A:326:ARG:HD3	1:A:330:LYS:HE2	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:PHE:CE1	1:A:88:MET:HE1	2.36	0.61
1:A:71:PRO:O	1:A:75:LEU:HB2	2.01	0.61
1:A:29:LEU:HD23	1:A:180:ILE:HD12	1.83	0.60
1:A:63:LEU:HD12	1:A:68:LEU:HG	1.84	0.59
1:A:49:LEU:HB2	1:A:50:PRO:HD3	1.83	0.58
1:A:135:ASP:HB2	1:A:299:MET:HE3	1.86	0.58
1:A:268:MET:CE	1:A:270:ALA:HB2	2.34	0.58
1:A:95:PRO:C	1:A:97:ASP:H	2.12	0.57
1:A:140:ASN:HB2	1:A:299:MET:CE	2.34	0.57
1:A:230:MET:SD	1:A:276:LEU:HD13	2.44	0.57
1:A:37:CYS:SG	1:A:39:LYS:HE2	2.45	0.56
1:A:89:GLU:O	1:A:93:LYS:HD3	2.05	0.56
1:A:168:PHE:O	1:A:169:ASP:C	2.48	0.56
1:A:379:GLU:O	1:A:380:ALA:HB3	2.06	0.55
1:A:8:ALA:HB3	1:A:9:PRO:HD3	1.87	0.55
1:A:42:PHE:CZ	1:A:46:ARG:HD2	2.42	0.55
1:A:99:ARG:O	1:A:103:GLN:HG2	2.07	0.55
1:A:95:PRO:O	1:A:97:ASP:N	2.33	0.54
1:A:87:ILE:HD13	3:A:2008:HOH:O	2.07	0.54
1:A:340:SER:HB2	1:A:345:GLY:O	2.07	0.54
1:A:12:ILE:HD11	1:A:57:MET:HE2	1.88	0.54
1:A:42:PHE:CZ	1:A:90:PHE:HB3	2.43	0.54
1:A:242:PHE:CZ	1:A:246:LYS:HD2	2.43	0.53
1:A:77:GLN:O	1:A:81:VAL:HG23	2.08	0.53
1:A:16:SER:OG	1:A:58:LYS:HG3	2.09	0.53
1:A:46:ARG:O	1:A:50:PRO:HG2	2.09	0.52
1:A:12:ILE:O	1:A:16:SER:HB3	2.08	0.52
1:A:25:MET:HE1	1:A:233:VAL:HG23	1.91	0.52
1:A:340:SER:HB3	1:A:346:THR:HG23	1.92	0.52
1:A:330:LYS:HE3	3:A:2091:HOH:O	2.09	0.52
1:A:34:SER:O	1:A:35:ASN:CB	2.58	0.52
1:A:38:GLU:HB3	1:A:168:PHE:CZ	2.45	0.51
1:A:28:PHE:HB3	1:A:183:ILE:HD12	1.91	0.51
1:A:339:PHE:CE1	1:A:347:ASP:HB2	2.46	0.51
1:A:207:TYR:HB2	3:A:2036:HOH:O	2.10	0.51
1:A:42:PHE:HZ	1:A:90:PHE:HB3	1.76	0.50
1:A:375:GLN:O	1:A:377:ILE:HG13	2.10	0.50
1:A:38:GLU:HB3	1:A:168:PHE:CE1	2.46	0.50
1:A:20:PRO:HD2	1:A:365:VAL:HG22	1.92	0.50
1:A:256:HIS:HE1	1:A:283:ARG:O	1.94	0.50
1:A:107:ALA:O	1:A:111:ILE:HG13	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:PHE:CD1	1:A:347:ASP:HB2	2.47	0.50
1:A:42:PHE:CE2	1:A:46:ARG:HD2	2.46	0.50
1:A:239:HIS:HB2	3:A:2059:HOH:O	2.12	0.49
1:A:95:PRO:C	1:A:97:ASP:N	2.71	0.49
1:A:12:ILE:O	1:A:16:SER:CB	2.61	0.48
1:A:162:ASN:O	1:A:166:LEU:HD13	2.13	0.48
1:A:283:ARG:HH11	1:A:283:ARG:HG3	1.79	0.48
1:A:11:TYR:HB3	1:A:88:MET:CE	2.44	0.48
1:A:202:LEU:C	1:A:202:LEU:HD23	2.39	0.48
1:A:43:THR:HA	1:A:46:ARG:HD3	1.95	0.47
1:A:287:VAL:HG13	1:A:288:PRO:HD2	1.97	0.47
1:A:224:SER:O	1:A:225:LYS:HD3	2.14	0.47
1:A:250:ARG:O	1:A:254:GLU:HG3	2.15	0.47
1:A:25:MET:HE2	1:A:183:ILE:HG23	1.95	0.47
1:A:84:LEU:HG	1:A:88:MET:HE2	1.96	0.47
1:A:94:ASP:O	1:A:95:PRO:O	2.33	0.47
1:A:111:ILE:O	1:A:115:HIS:HD2	1.97	0.47
1:A:290:ARG:HH12	1:A:291:LYS:HE3	1.78	0.47
1:A:54:ALA:O	1:A:58:LYS:HB2	2.15	0.47
1:A:95:PRO:O	1:A:96:GLU:HB2	2.15	0.46
1:A:65:ASP:HA	1:A:68:LEU:HB2	1.96	0.46
1:A:28:PHE:HB3	1:A:183:ILE:CD1	2.46	0.46
1:A:97:ASP:C	1:A:99:ARG:H	2.24	0.46
1:A:250:ARG:NH1	1:A:254:GLU:OE1	2.49	0.46
1:A:377:ILE:HG22	1:A:377:ILE:O	2.15	0.46
1:A:283:ARG:HG3	1:A:283:ARG:NH1	2.31	0.46
1:A:11:TYR:CD2	1:A:88:MET:HE3	2.50	0.46
1:A:84:LEU:O	1:A:88:MET:HG2	2.16	0.46
1:A:261:ILE:HG22	1:A:262:LEU:N	2.31	0.45
1:A:12:ILE:CD1	1:A:57:MET:HE2	2.47	0.45
1:A:81:VAL:O	1:A:85:LEU:HG	2.17	0.45
1:A:15:PHE:HE1	1:A:88:MET:HE1	1.80	0.44
1:A:20:PRO:HB3	1:A:52:ARG:HG2	1.98	0.44
1:A:193:VAL:HA	1:A:241:LEU:HD13	1.97	0.44
1:A:169:ASP:O	1:A:169:ASP:OD1	2.36	0.44
1:A:44:PHE:CZ	1:A:48:GLU:HG2	2.52	0.44
1:A:64:PRO:HB2	1:A:67:VAL:HG23	2.00	0.44
1:A:293:GLU:H	1:A:293:GLU:HG2	1.60	0.44
1:A:159:MET:HE2	1:A:332:PHE:HZ	1.83	0.43
1:A:113:ASN:HA	1:A:116:ASN:ND2	2.34	0.43
1:A:159:MET:HG3	1:A:331:TYR:HE2	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:PHE:CE1	1:A:347:ASP:CB	3.02	0.43
1:A:122:MET:HE3	1:A:150:PHE:CE2	2.54	0.42
1:A:178:LYS:N	1:A:178:LYS:HD2	2.34	0.42
1:A:92:ASP:OD1	1:A:93:LYS:HD2	2.19	0.42
1:A:18:PHE:CE2	1:A:47:GLN:NE2	2.88	0.42
1:A:129:TYR:C	1:A:129:TYR:CD1	2.97	0.42
1:A:267:VAL:HA	1:A:279:LYS:O	2.20	0.42
1:A:63:LEU:O	1:A:64:PRO:C	2.62	0.42
1:A:66:ARG:NH2	1:A:129:TYR:OH	2.52	0.42
1:A:159:MET:HG3	1:A:331:TYR:CE2	2.55	0.42
1:A:202:LEU:HD23	1:A:202:LEU:O	2.20	0.42
1:A:123:ALA:HB2	1:A:320:TYR:CE2	2.55	0.41
1:A:288:PRO:HG3	1:A:342:GLU:OE1	2.20	0.41
1:A:12:ILE:HD11	1:A:57:MET:CE	2.49	0.41
1:A:68:LEU:O	1:A:74:GLN:HG3	2.20	0.41
1:A:138:VAL:HG12	1:A:142:ASN:HD21	1.85	0.41
1:A:217:GLN:HE21	1:A:219:ILE:HD11	1.86	0.40
1:A:379:GLU:O	1:A:380:ALA:CB	2.70	0.40
1:A:336:LEU:C	1:A:336:LEU:HD23	2.46	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:LYS:NZ	1:A:259:SER:O[3_665]	2.03	0.17

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	348/394 (88%)	321 (92%)	20 (6%)	7 (2%)	6 4

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	95	PRO
1	A	224	SER
1	A	37	CYS
1	A	61	ASN
1	A	98	HIS
1	A	64	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	320/353 (91%)	309 (97%)	11 (3%)	32 43

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	MET
1	A	26	LYS
1	A	66	ARG
1	A	75	LEU
1	A	76	VAL
1	A	86	ASP
1	A	101	LEU
1	A	159	MET
1	A	178	LYS
1	A	293	GLU
1	A	382	ASP

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

Mol	Chain	Res	Type
1	A	55	ASN
1	A	77	GLN

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Mol	Chain	Res	Type
1	A	103	GLN
1	A	113	ASN
1	A	116	ASN
1	A	142	ASN
1	A	179	HIS
1	A	217	GLN
1	A	229	HIS
1	A	236	HIS
1	A	375	GLN

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](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	TF1	A	1386	-	27,28,28	2.48	13 (48%)	38,43,43	1.67	6 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.
'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TF1	A	1386	-	-	3/29/45/45	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1386	TF1	C8-C4	5.42	1.60	1.52
2	A	1386	TF1	O2-C4	-4.78	1.34	1.42
2	A	1386	TF1	C1-N1	4.12	1.42	1.34
2	A	1386	TF1	C13-C12	3.99	1.45	1.39
2	A	1386	TF1	C3-N1	3.71	1.51	1.46
2	A	1386	TF1	C15-C13	2.96	1.43	1.38
2	A	1386	TF1	C16-C14	2.87	1.43	1.38
2	A	1386	TF1	C9-C4	2.55	1.56	1.53
2	A	1386	TF1	C11-N2	2.39	1.40	1.34
2	A	1386	TF1	C15-C17	2.36	1.44	1.39
2	A	1386	TF1	C5-N2	2.34	1.49	1.46
2	A	1386	TF1	C16-C17	2.20	1.43	1.39
2	A	1386	TF1	C14-C12	2.18	1.42	1.39

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1386	TF1	C8-C4-C1	4.94	114.60	108.54
2	A	1386	TF1	O2-C4-C8	4.30	112.52	105.94
2	A	1386	TF1	C7-C3-N1	2.76	113.81	110.80
2	A	1386	TF1	C14-C12-C13	2.34	121.54	118.57
2	A	1386	TF1	C9-C4-C8	-2.06	106.31	109.04
2	A	1386	TF1	C3-N1-C1	2.00	129.74	123.20

There are no chirality outliers.

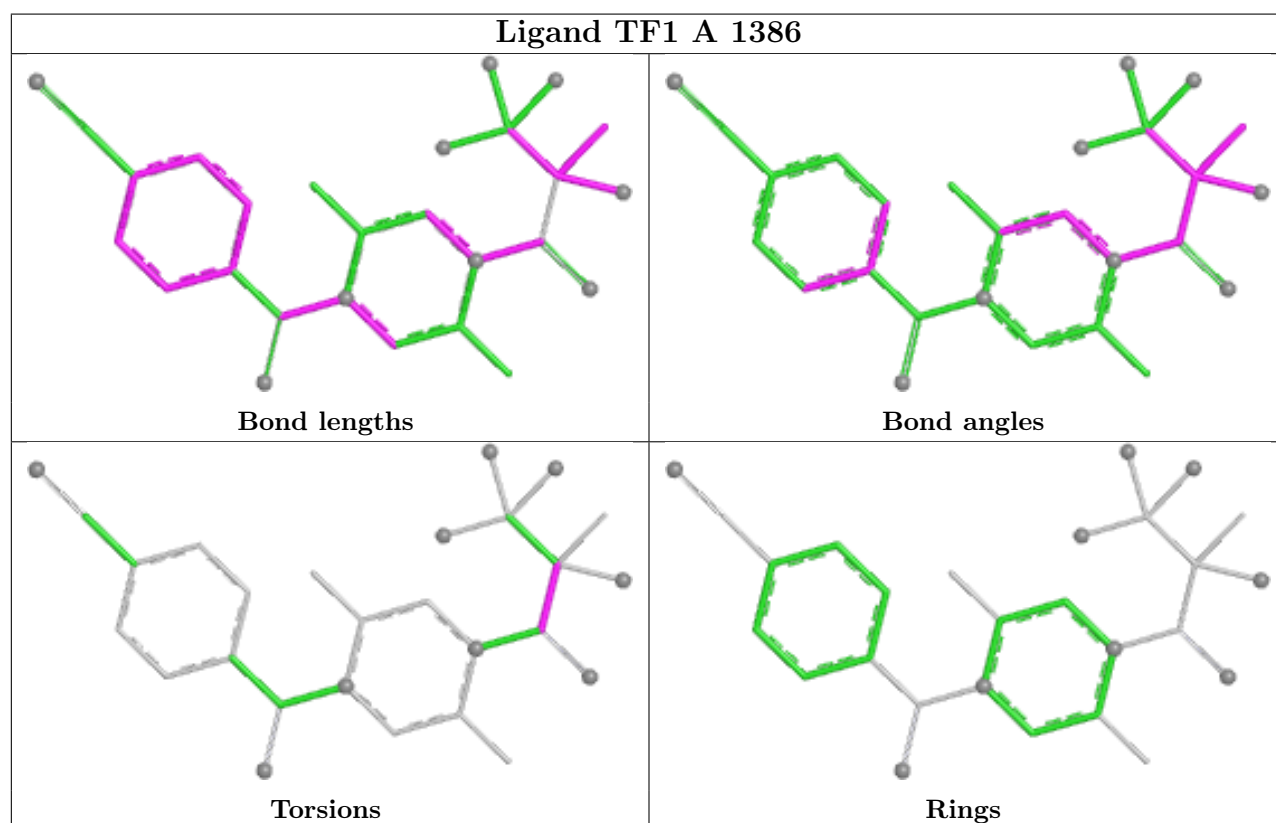
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1386	TF1	N1-C1-C4-C8
2	A	1386	TF1	O1-C1-C4-C8
2	A	1386	TF1	O1-C1-C4-O2

There are no ring outliers.

No monomer is involved in short contacts.

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.