



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:50 PM UTC

PDB ID : 1UNL / pdb_00001unl
Title : Structural mechanism for the inhibition of CD5-p25 from the roscovitine, aloisine and indirubin.
Authors : Mapelli, M.; Crovace, C.; Massimiliano, L.; Musacchio, A.
Deposited on : 2003-09-10
Resolution : 2.20 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

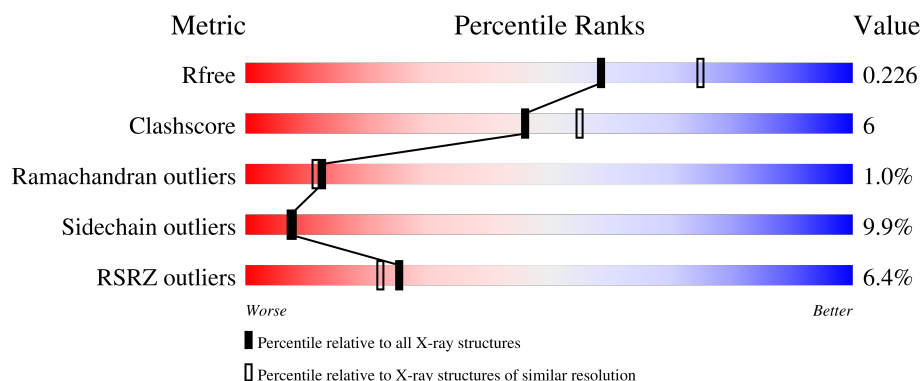
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	292	6% 80% 16% ..
1	B	292	9% 78% 15% ..
2	D	208	64% 7% 28% .
2	E	208	5% 56% 14% 28% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7424 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 CYCLIN-DEPENDENT KINASE 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2346	1505	403	426	12			
1	B	292	Total	C	N	O	S	0	0	0
			2346	1505	403	426	12			

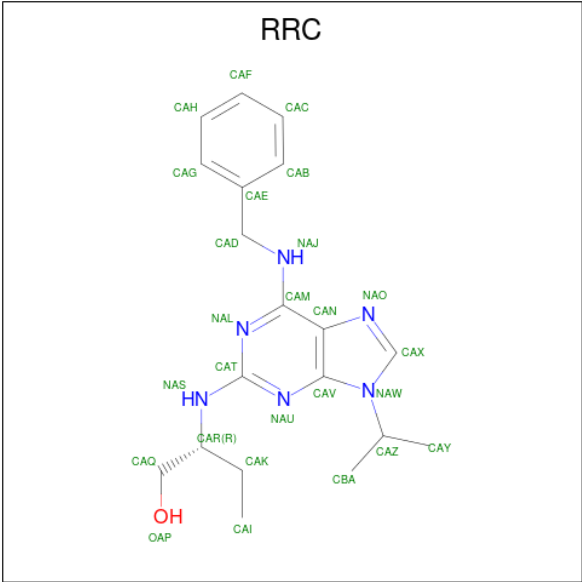
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	144	ASN	ASP	engineered mutation	UNP Q00535
B	144	ASN	ASP	engineered mutation	UNP Q00535

- Molecule 2 is a protein called CYCLIN-DEPENDENT KINASE 5 ACTIVATOR 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	150	Total	C	N	O	S	0	0	1
			1202	771	199	221	11			
2	E	150	Total	C	N	O	S	0	0	1
			1202	771	199	221	11			

- Molecule 3 is R-ROSCOVITINE (CCD ID: RRC) (formula: C₁₉H₂₆N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			26	19	6	1		

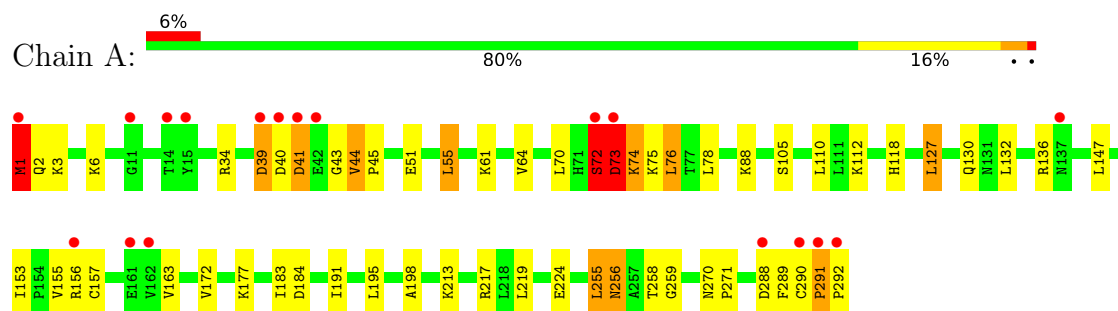
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	142	Total	O	0	0
			142	142		
4	B	80	Total	O	0	0
			80	80		
4	D	58	Total	O	0	0
			58	58		
4	E	22	Total	O	0	0
			22	22		

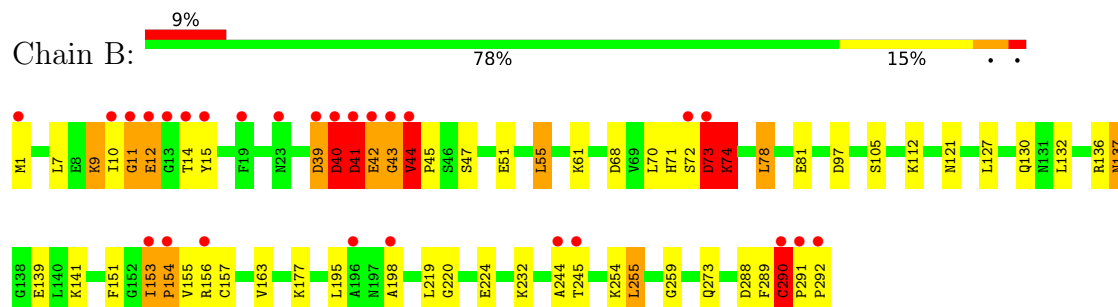
3 Residue-property plots [i](#)

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.

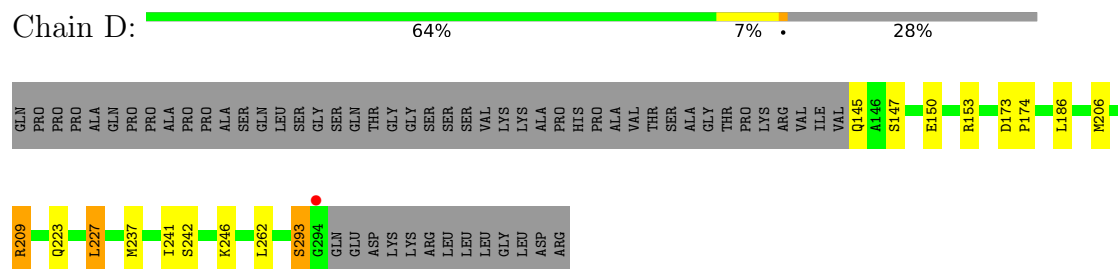
• Molecule 1: CYCLIN-DEPENDENT KINASE 5



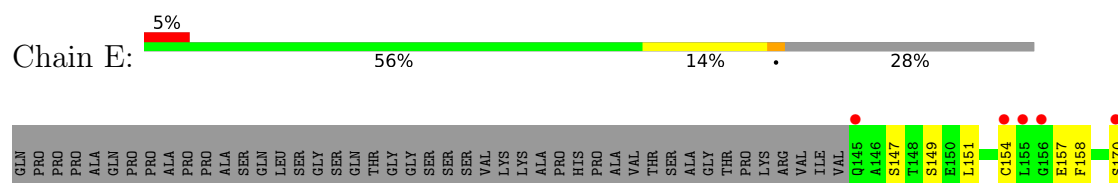
• Molecule 1: CYCLIN-DEPENDENT KINASE 5

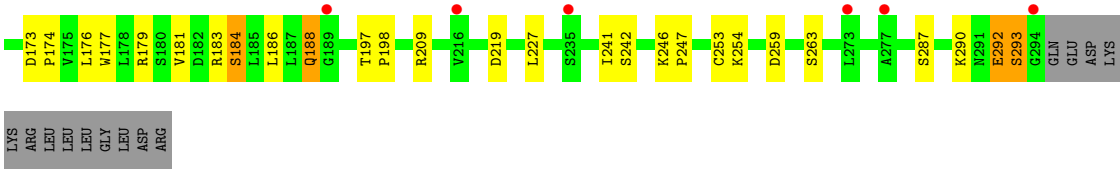


• Molecule 2: CYCLIN-DEPENDENT KINASE 5 ACTIVATOR 1



• Molecule 2: CYCLIN-DEPENDENT KINASE 5 ACTIVATOR 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.99Å 117.99Å 156.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.76 – 2.20 19.76 – 2.20	Depositor EDS
% Data completeness (in resolution range)	94.4 (19.76-2.20) 94.3 (19.76-2.20)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.21Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.216 , 0.219 0.218 , 0.226	Depositor DCC
R_{free} test set	3057 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	36.4	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7424	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: RRC

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	1.09	6/2402 (0.2%)	1.11	10/3251 (0.3%)
1	B	0.91	5/2402 (0.2%)	1.12	15/3251 (0.5%)
2	D	1.03	2/1230 (0.2%)	0.95	0/1669
2	E	0.58	1/1230 (0.1%)	0.83	0/1669
All	All	0.95	14/7264 (0.2%)	1.04	25/9840 (0.3%)

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	A	0	2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	74	LYS	N-CA	-11.69	1.35	1.46
1	B	44	VAL	CA-CB	10.11	1.59	1.54
1	A	73	ASP	C-N	-9.50	1.18	1.34
1	A	198	ALA	N-CA	7.79	1.56	1.46
1	B	43	GLY	N-CA	7.32	1.55	1.45
2	E	293	SER	C-N	-6.62	1.24	1.33
2	D	293	SER	C-N	-6.42	1.24	1.33
1	A	74	LYS	C-N	6.19	1.42	1.33
1	A	256	ASN	CG-OD1	6.09	1.35	1.23
2	D	237	MET	SD-CE	-6.03	1.64	1.79
1	B	39	ASP	C-N	-5.96	1.25	1.33
1	A	1	MET	SD-CE	5.75	1.94	1.79
1	B	1	MET	SD-CE	5.30	1.92	1.79
1	B	291	PRO	CA-C	5.12	1.56	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	40	ASP	CA-CB-CG	-13.49	99.11	112.60
1	A	74	LYS	N-CA-C	-12.88	92.25	112.99
1	A	72	SER	CA-C-N	-11.72	99.15	121.54
1	A	72	SER	C-N-CA	-11.72	99.15	121.54
1	B	41	ASP	CA-CB-CG	-10.91	101.69	112.60
1	A	198	ALA	N-CA-CB	9.37	126.33	110.49
1	B	72	SER	N-CA-C	8.77	124.48	113.43
1	B	42	GLU	N-CA-C	-8.22	93.29	110.80
1	B	198	ALA	N-CA-CB	7.10	122.48	110.49
1	B	291	PRO	N-CA-C	6.78	118.97	110.70
1	B	71	HIS	CA-C-N	-6.30	109.53	122.58
1	B	71	HIS	C-N-CA	-6.30	109.53	122.58
1	A	74	LYS	CA-C-N	-6.08	113.13	122.62
1	A	74	LYS	C-N-CA	-6.08	113.13	122.62
1	A	72	SER	N-CA-C	5.79	119.47	111.71
1	B	73	ASP	CA-CB-CG	5.75	118.35	112.60
1	B	41	ASP	CA-C-O	5.75	128.73	120.51
1	A	291	PRO	N-CA-C	5.57	117.49	110.70
1	B	41	ASP	O-C-N	5.49	129.89	122.59
1	A	40	ASP	N-CA-C	-5.38	105.54	113.61
1	B	74	LYS	N-CA-C	-5.35	103.92	110.61
1	B	41	ASP	CB-CA-C	5.35	121.06	110.42
1	B	68	ASP	N-CA-C	5.31	116.35	108.60
1	B	40	ASP	CB-CA-C	-5.26	99.95	110.42
1	A	39	ASP	N-CA-C	5.17	117.79	110.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	41	ASP	Peptide
1	A	72	SER	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	A	2346	0	2365	37	0
1	B	2346	0	2365	33	0
2	D	1202	0	1187	8	0
2	E	1202	0	1187	10	0
3	A	26	0	26	1	0
4	A	142	0	0	3	0
4	B	80	0	0	2	0
4	D	58	0	0	1	0
4	E	22	0	0	1	0
All	All	7424	0	7130	83	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 6.

All (83) 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:73:ASP:OD1	1:A:74:LYS:N	1.85	1.08
1:A:73:ASP:OD1	1:A:75:LYS:N	1.89	1.04
1:A:73:ASP:CG	1:A:74:LYS:H	1.69	0.97
1:A:72:SER:O	1:A:73:ASP:HB3	1.62	0.96
1:B:40:ASP:N	1:B:40:ASP:OD1	1.81	0.96
1:B:40:ASP:CG	1:B:41:ASP:H	1.71	0.88
1:B:40:ASP:CG	1:B:41:ASP:N	2.30	0.84
1:A:1:MET:HG3	1:A:70:LEU:HD13	1.61	0.82
1:B:39:ASP:HB3	1:B:40:ASP:OD1	1.83	0.77
1:A:73:ASP:CG	1:A:74:LYS:N	2.30	0.76
1:B:220:GLY:HA3	1:B:244:ALA:HB2	1.66	0.76
2:E:181:VAL:HA	4:E:2006:HOH:O	1.86	0.75
1:B:81:GLU:OE1	1:B:141:LYS:HE2	1.88	0.74
1:A:256:ASN:HD21	1:A:258:THR:HB	1.57	0.70
1:A:157:CYS:HB2	1:A:177:LYS:HB3	1.75	0.68
1:A:73:ASP:OD1	1:A:74:LYS:C	2.37	0.68
1:B:9:LYS:HB3	1:B:12:GLU:OE2	1.94	0.67
1:B:43:GLY:HA3	2:E:242:SER:HB3	1.76	0.66
2:D:209:ARG:HD2	4:D:2021:HOH:O	1.95	0.66
1:B:177:LYS:HE2	4:B:2040:HOH:O	1.96	0.66
1:A:72:SER:O	1:A:73:ASP:CB	2.29	0.64
1:A:290:CYS:HB2	1:A:291:PRO:HD3	1.82	0.61
1:A:291:PRO:HB2	1:A:292:PRO:HD3	1.83	0.61
1:B:137:ASN:HD22	1:B:139:GLU:HG3	1.66	0.60
2:E:173:ASP:HB2	2:E:174:PRO:HD3	1.84	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:ASP:O	1:B:44:VAL:CG2	2.50	0.59
1:A:43:GLY:HA3	2:D:242:SER:HB2	1.84	0.58
1:A:256:ASN:ND2	1:A:258:THR:HB	2.17	0.58
2:E:174:PRO:HA	2:E:177:TRP:CE3	2.40	0.57
2:D:145:GLN:HA	2:D:150:GLU:HG2	1.87	0.57
1:B:39:ASP:C	1:B:40:ASP:OD1	2.47	0.55
1:B:157:CYS:HB2	1:B:177:LYS:HB3	1.90	0.54
1:A:270:ASN:HD22	1:A:270:ASN:C	2.16	0.52
2:E:184:SER:O	2:E:188:GLN:HB2	2.08	0.52
2:E:246:LYS:HB3	2:E:247:PRO:HD3	1.91	0.52
1:B:121:ASN:HA	1:B:151:PHE:CE1	2.45	0.52
1:B:40:ASP:O	1:B:42:GLU:N	2.41	0.52
1:B:41:ASP:O	1:B:44:VAL:HG22	2.10	0.51
1:A:73:ASP:H	1:A:75:LYS:H	1.58	0.51
1:B:41:ASP:O	1:B:44:VAL:HG23	2.11	0.51
1:A:172:VAL:HG21	1:A:183:ILE:HD12	1.93	0.51
2:D:241:ILE:HG13	2:D:242:SER:N	2.26	0.50
1:B:156:ARG:HH12	2:E:197:THR:HG21	1.76	0.50
1:B:290:CYS:C	1:B:292:PRO:HD2	2.37	0.50
1:A:64:VAL:HG21	3:A:1293:RRC:HAY2	1.94	0.49
1:A:6:LYS:HE3	1:A:34:ARG:HD2	1.96	0.48
2:D:173:ASP:HB2	2:D:174:PRO:HD3	1.96	0.48
1:B:15:TYR:OH	1:B:51:GLU:OE2	2.23	0.48
1:A:288:ASP:HB3	4:A:2141:HOH:O	2.13	0.48
1:B:44:VAL:N	1:B:45:PRO:CD	2.77	0.47
2:E:151:LEU:HD12	2:E:198:PRO:HA	1.96	0.47
1:B:43:GLY:CA	2:E:242:SER:HB3	2.45	0.47
1:B:105:SER:HA	1:B:289:PHE:CZ	2.50	0.47
1:A:255:LEU:HG	1:A:259:GLY:HA3	1.97	0.47
2:E:158:PHE:HB2	2:E:292:GLU:HG2	1.97	0.47
1:A:118:HIS:HE1	1:A:184:ASP:OD1	1.99	0.46
1:B:10:ILE:O	1:B:12:GLU:N	2.47	0.46
1:A:51:GLU:HG3	1:A:55:LEU:HD22	1.98	0.46
1:B:273:GLN:NE2	4:B:2071:HOH:O	2.48	0.46
1:B:43:GLY:H	1:B:45:PRO:HD2	1.80	0.46
2:D:223:GLN:HG2	2:D:227:LEU:HD22	1.97	0.46
1:A:289:PHE:HB3	4:A:2142:HOH:O	2.16	0.45
1:B:97:ASP:OD2	1:B:254:LYS:NZ	2.44	0.45
1:A:44:VAL:N	1:A:45:PRO:CD	2.80	0.45
1:B:51:GLU:O	1:B:55:LEU:HB2	2.17	0.45
1:A:76:LEU:HD11	2:D:262:LEU:HD11	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:ASP:HB3	1:B:74:LYS:H	1.16	0.44
2:D:153:ARG:HH12	2:D:293:SER:HB2	1.83	0.44
1:A:127:LEU:HD22	1:A:191:ILE:CD1	2.48	0.43
1:B:10:ILE:O	1:B:11:GLY:C	2.61	0.43
1:B:255:LEU:HG	1:B:259:GLY:HA3	2.01	0.43
1:A:110:LEU:HD11	1:A:127:LEU:HD21	2.01	0.42
1:A:73:ASP:N	1:A:75:LYS:H	2.17	0.42
1:A:73:ASP:OD1	1:A:75:LYS:HG3	2.19	0.42
1:B:78:LEU:N	1:B:78:LEU:HD23	2.33	0.42
1:A:88:LYS:NZ	4:A:2042:HOH:O	2.47	0.42
1:A:213:LYS:O	1:A:217:ARG:HB2	2.20	0.42
1:A:270:ASN:HD22	1:A:271:PRO:N	2.17	0.42
1:A:270:ASN:HD22	1:A:271:PRO:HD2	1.85	0.42
1:A:1:MET:HE3	1:A:3:LYS:H	1.85	0.41
1:A:105:SER:HA	1:A:289:PHE:CZ	2.56	0.41
1:A:105:SER:HA	1:A:289:PHE:HZ	1.86	0.41
1:B:153:ILE:O	1:B:154:PRO:C	2.63	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	290/292 (99%)	273 (94%)	15 (5%)	2 (1%)	18	19
1	B	290/292 (99%)	265 (91%)	19 (7%)	6 (2%)	5	3
2	D	148/208 (71%)	146 (99%)	2 (1%)	0	100	100
2	E	148/208 (71%)	142 (96%)	5 (3%)	1 (1%)	18	19
All	All	876/1000 (88%)	826 (94%)	41 (5%)	9 (1%)	12	11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	40	ASP
1	B	11	GLY
1	B	163	VAL
1	A	73	ASP
1	A	163	VAL
2	E	293	SER
1	B	41	ASP
1	B	154	PRO
1	B	290	CYS

5.3.2 Protein sidechains ⓘ

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	260/260 (100%)	238 (92%)	22 (8%)	10	11
1	B	260/260 (100%)	231 (89%)	29 (11%)	6	5
2	D	138/186 (74%)	132 (96%)	6 (4%)	26	35
2	E	138/186 (74%)	116 (84%)	22 (16%)	2	2
All	All	796/892 (89%)	717 (90%)	79 (10%)	7	8

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	GLN
1	A	39	ASP
1	A	41	ASP
1	A	44	VAL
1	A	55	LEU
1	A	61	LYS
1	A	76	LEU
1	A	78	LEU
1	A	112	LYS
1	A	127	LEU
1	A	130	GLN
1	A	132	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	136	ARG
1	A	147	LEU
1	A	153	ILE
1	A	155	VAL
1	A	156	ARG
1	A	195	LEU
1	A	219	LEU
1	A	224	GLU
1	A	255	LEU
1	B	7	LEU
1	B	9	LYS
1	B	12	GLU
1	B	14	THR
1	B	41	ASP
1	B	44	VAL
1	B	47	SER
1	B	55	LEU
1	B	61	LYS
1	B	70	LEU
1	B	73	ASP
1	B	74	LYS
1	B	78	LEU
1	B	112	LYS
1	B	127	LEU
1	B	130	GLN
1	B	132	LEU
1	B	136	ARG
1	B	137	ASN
1	B	153	ILE
1	B	155	VAL
1	B	195	LEU
1	B	219	LEU
1	B	224	GLU
1	B	232	LYS
1	B	245	THR
1	B	255	LEU
1	B	288	ASP
1	B	290	CYS
2	D	147	SER
2	D	186	LEU
2	D	206	MET
2	D	209	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	D	227	LEU
2	D	246	LYS
2	E	147	SER
2	E	149	SER
2	E	154	CYS
2	E	157	GLU
2	E	170	SER
2	E	176	LEU
2	E	179	ARG
2	E	183	ARG
2	E	184	SER
2	E	186	LEU
2	E	188	GLN
2	E	209	ARG
2	E	219	ASP
2	E	227	LEU
2	E	241	ILE
2	E	253	CYS
2	E	254	LYS
2	E	259	ASP
2	E	263	SER
2	E	287	SER
2	E	290	LYS
2	E	292	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	85	GLN
1	A	118	HIS
1	A	130	GLN
1	A	135	ASN
1	A	206	ASN
1	A	270	ASN
1	B	62	ASN
1	B	130	GLN
1	B	135	ASN
1	B	137	ASN
1	B	206	ASN
1	B	270	ASN
1	B	273	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	282	GLN
2	E	193	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
3	RRC	A	1293	-	28,28,28	1.14	3 (10%)	38,38,38	2.26	13 (34%)

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	RRC	A	1293	-	-	3/17/17/17	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1293	RRC	CAX-NAO	3.65	1.38	1.31
3	A	1293	RRC	CAV-NAW	-2.86	1.31	1.37
3	A	1293	RRC	CAN-CAV	-2.10	1.35	1.39

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1293	RRC	CAZ-NAW-CAX	6.98	136.35	128.05
3	A	1293	RRC	NAL-CAT-NAU	-4.97	118.12	126.26
3	A	1293	RRC	NAS-CAT-NAL	4.02	123.41	117.09
3	A	1293	RRC	CAN-CAV-NAU	-3.86	123.11	127.18
3	A	1293	RRC	CAV-NAW-CAZ	-3.61	116.33	125.74
3	A	1293	RRC	CAN-CAV-NAW	3.23	109.33	105.81
3	A	1293	RRC	NAJ-CAM-NAL	3.02	122.38	118.33
3	A	1293	RRC	CAQ-CAR-NAS	-3.00	103.14	108.68
3	A	1293	RRC	CBA-CAZ-CAY	2.83	120.15	112.46
3	A	1293	RRC	CAT-NAL-CAM	2.52	122.55	116.36
3	A	1293	RRC	CAN-CAM-NAJ	-2.21	117.76	122.03
3	A	1293	RRC	CAK-CAR-CAQ	-2.19	107.39	112.12
3	A	1293	RRC	NAW-CAX-NAO	-2.10	110.96	113.94

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1293	RRC	CAI-CAK-CAR-CAQ
3	A	1293	RRC	CAI-CAK-CAR-NAS
3	A	1293	RRC	CAQ-CAR-NAS-CAT

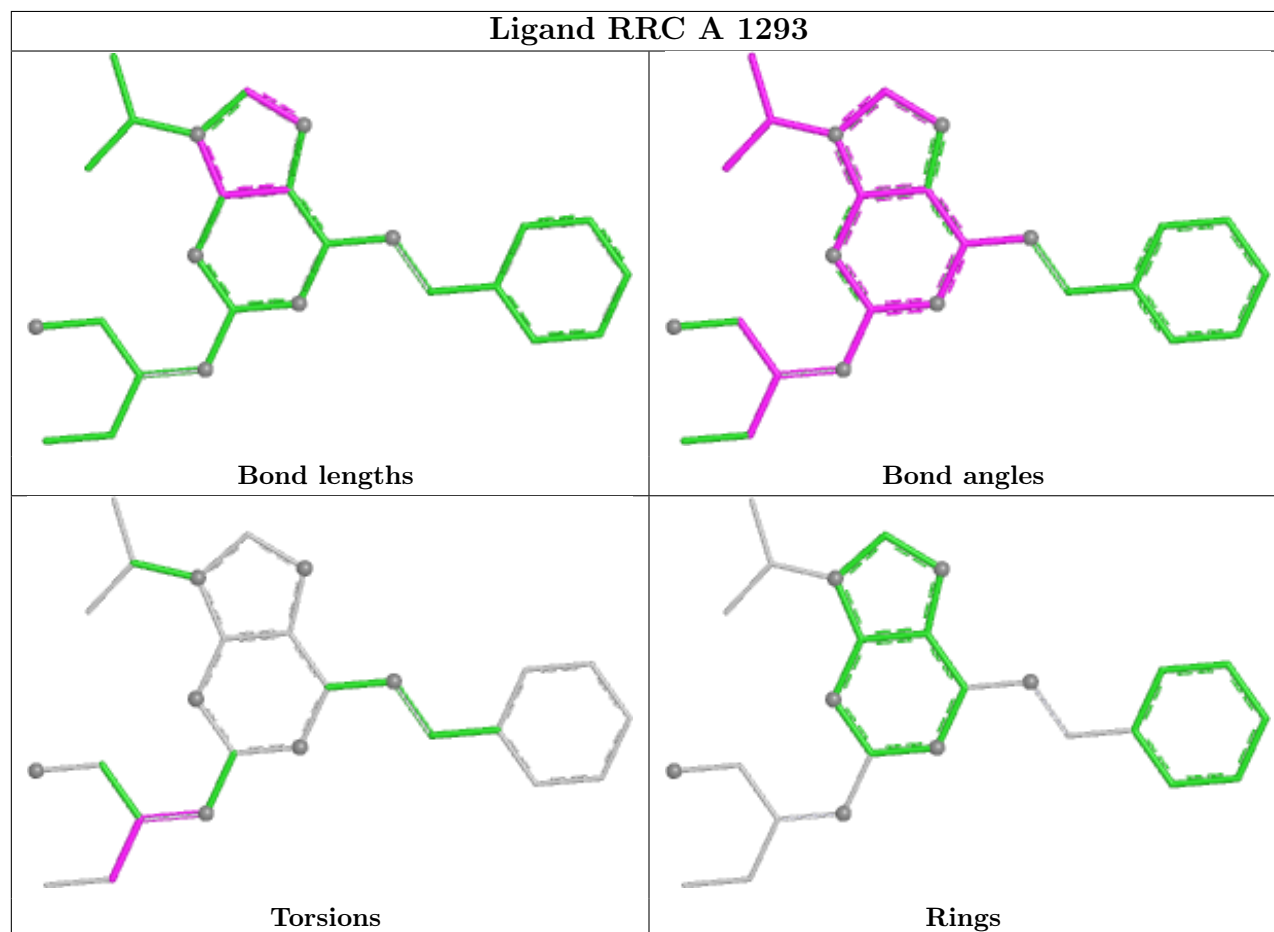
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1293	RRC	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	73:ASP	C	74:LYS	N	1.18

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	292/292 (100%)	0.19	18 (6%) 26 23	18, 23, 33, 53	0
1	B	292/292 (100%)	0.81	27 (9%) 14 12	18, 24, 32, 53	0
2	D	150/208 (72%)	-0.00	1 (0%) 84 82	19, 23, 29, 32	0
2	E	150/208 (72%)	0.68	11 (7%) 21 18	20, 24, 25, 26	0
All	All	884/1000 (88%)	0.44	57 (6%) 25 22	18, 24, 31, 53	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	PRO	8.8
1	B	292	PRO	7.5
1	B	43	GLY	6.8
2	E	294	GLY	6.4
1	A	290	CYS	6.4
1	B	291	PRO	6.3
1	A	291	PRO	6.1
1	B	13	GLY	5.9
1	B	1	MET	5.9
1	B	290	CYS	5.0
1	B	10	ILE	4.6
1	B	244	ALA	4.6
1	B	153	ILE	4.6
1	B	41	ASP	4.3
1	B	15	TYR	4.2
1	B	39	ASP	4.1
1	A	39	ASP	4.0
1	A	40	ASP	4.0
1	B	73	ASP	3.8
1	B	40	ASP	3.6
2	D	294	GLY	3.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	11	GLY	3.4
1	B	14	THR	3.4
1	B	245	THR	3.4
1	A	42	GLU	3.4
1	B	12	GLU	3.4
1	A	11	GLY	3.3
1	B	72	SER	3.3
1	A	72	SER	3.1
1	A	288	ASP	3.1
1	A	41	ASP	3.1
1	B	156	ARG	3.1
1	B	44	VAL	3.0
1	B	154	PRO	2.9
1	B	198	ALA	2.9
2	E	273	LEU	2.9
2	E	216	VAL	2.8
2	E	189	GLY	2.8
1	A	137	ASN	2.7
1	A	14	THR	2.6
2	E	156	GLY	2.6
1	B	23	ASN	2.5
1	A	161	GLU	2.4
2	E	155	LEU	2.4
1	A	15	TYR	2.4
1	A	73	ASP	2.4
1	A	162	VAL	2.3
1	B	42	GLU	2.3
1	B	19	PHE	2.3
1	A	1	MET	2.3
2	E	170	SER	2.3
2	E	277	ALA	2.2
2	E	235	SER	2.2
2	E	154	CYS	2.2
1	A	156	ARG	2.1
2	E	145	GLN	2.1
1	B	196	ALA	2.1

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

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](#)

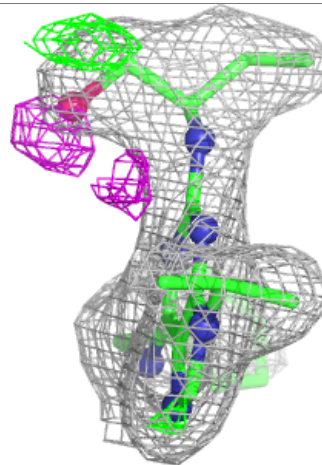
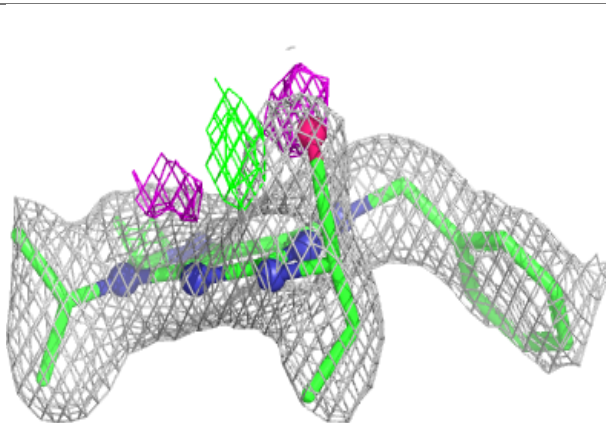
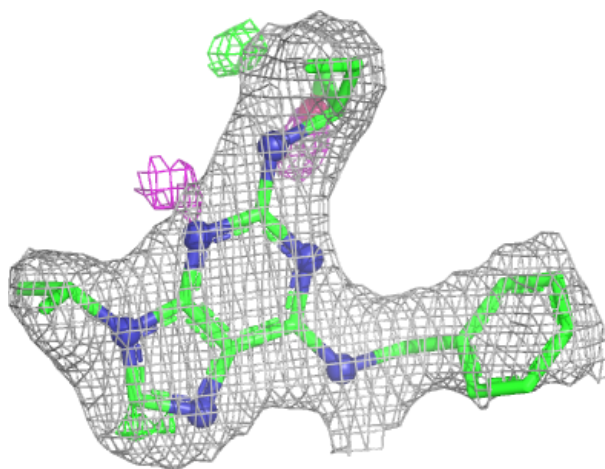
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	RRC	A	1293	26/26	0.87	0.10	19,23,24,25	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 RRC A 1293:

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.