



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:43 PM UTC

PDB ID : 3FO9 / pdb_00003fo9
Title : Crystal structure of aldolase antibody 33F12 Fab' in complex with hapten 1,3-diketone
Authors : Zhu, X.; Wilson, I.A.
Deposited on : 2008-12-29
Resolution : 1.90 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

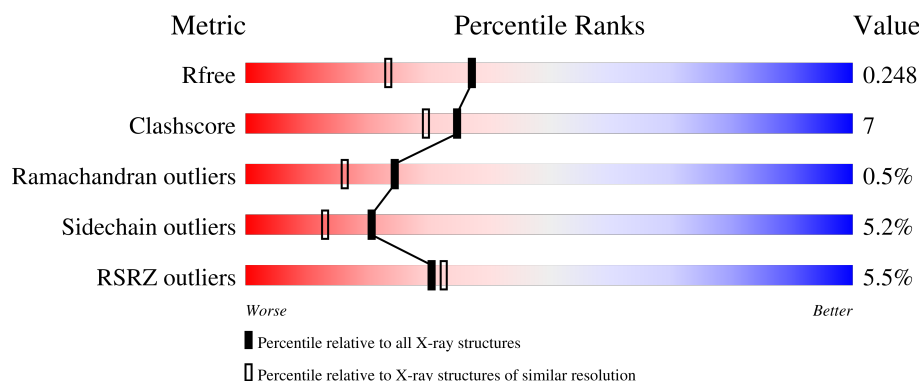
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	219	5% 84% 14% .
1	L	219	8% 81% 16% .
2	B	218	4% 84% 12% .
2	H	218	6% 84% 14% .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7028 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.

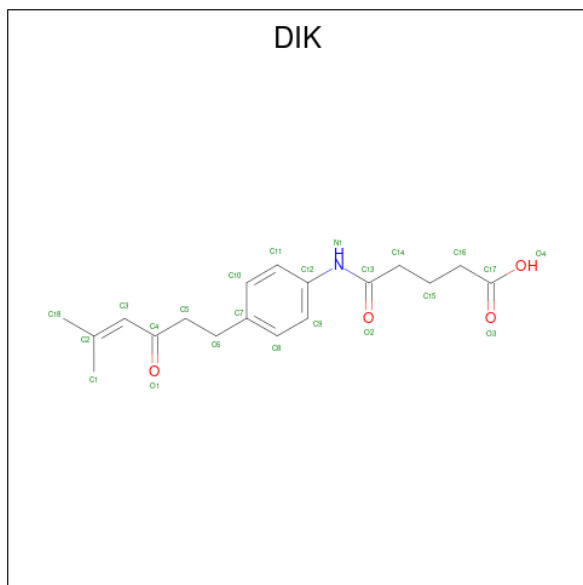
- Molecule 1 is a protein called Immunoglobulin IGG2A - light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	219	Total	C	N	O	S	0	0	0
			1697	1060	287	343	7			
1	A	219	Total	C	N	O	S	0	0	0
			1697	1060	287	343	7			

- Molecule 2 is a protein called Immunoglobulin IGG2A - heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	218	Total	C	N	O	S	0	0	0
			1676	1066	274	328	8			
2	B	218	Total	C	N	O	S	0	0	0
			1676	1066	274	328	8			

- Molecule 3 is 5-{{4-(5-methyl-3-oxohex-4-en-1-yl)phenyl}amino}-5-oxopentanoic acid (CCD ID: DIK) (formula: C₁₈H₂₃NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	H	1	Total	C	N	O	0	0
			22	17	1	4		
3	B	1	Total	C	N	O	0	0
			22	17	1	4		

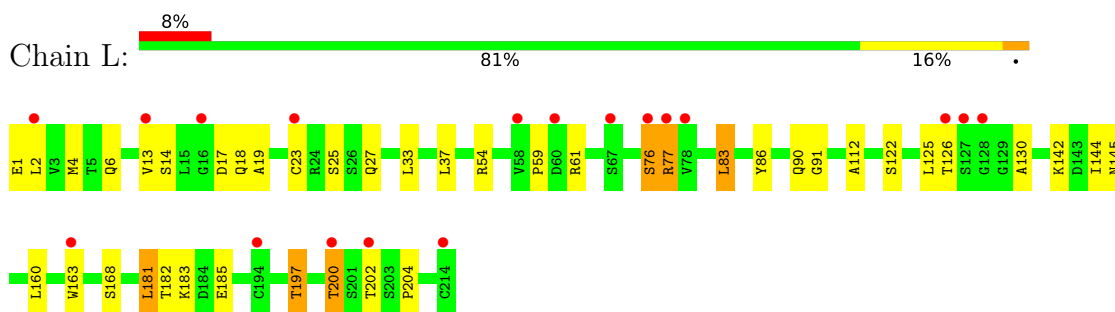
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	231	Total	O	0	0
			231	231		
4	H	4	Total	O	0	0
			4	4		
4	A	3	Total	O	0	0
			3	3		

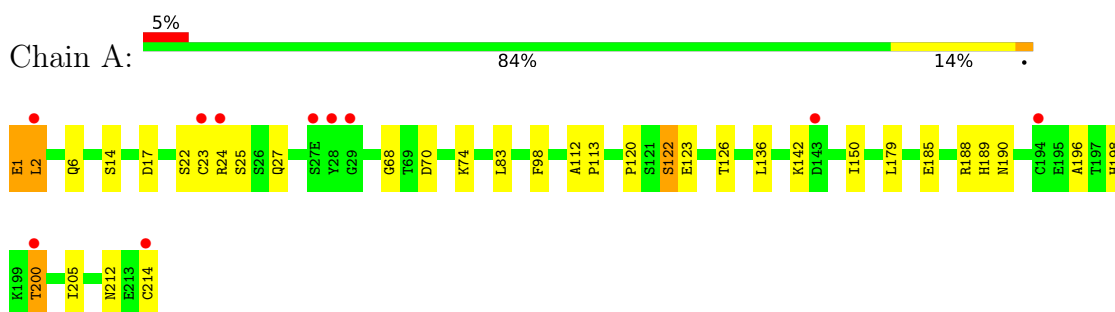
3 Residue-property plots [i](#)

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.

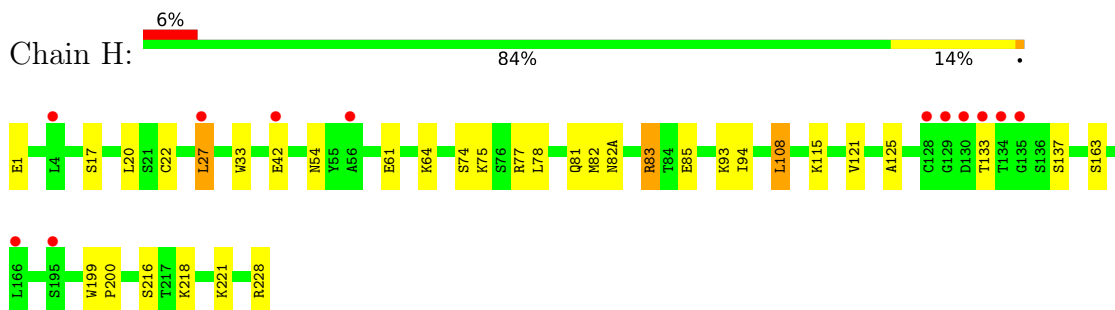
- Molecule 1: Immunoglobulin IGG2A - light chain



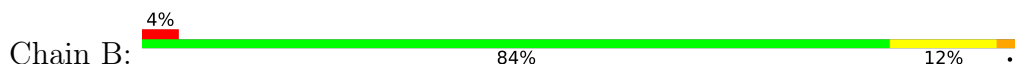
- Molecule 1: Immunoglobulin IGG2A - light chain

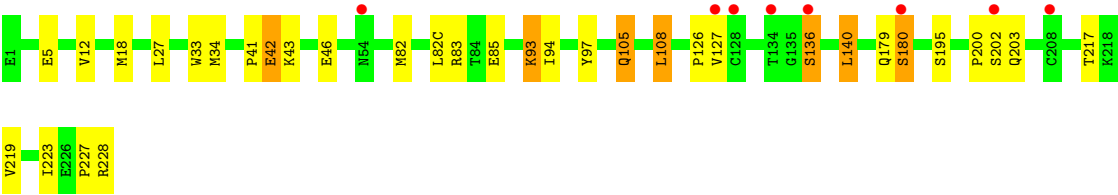


- Molecule 2: Immunoglobulin IGG2A - heavy chain



- Molecule 2: Immunoglobulin IGG2A - heavy chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.83Å 81.04Å 174.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.56 – 1.90 43.56 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.1 (43.56-1.90) 96.1 (43.56-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.215 , 0.249 0.213 , 0.248	Depositor DCC
R_{free} test set	3601 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.383	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7028	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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: DIK

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	0.57	0/1735	0.74	0/2351
1	L	0.65	0/1735	0.78	0/2351
2	B	0.65	0/1719	0.83	0/2339
2	H	0.63	0/1719	0.80	0/2339
All	All	0.63	0/6908	0.78	0/9380

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	1697	0	1637	19	0
1	L	1697	0	1637	37	0
2	B	1676	0	1645	27	0
2	H	1676	0	1645	17	0
3	B	22	0	19	6	0
3	H	22	0	19	4	0
4	A	3	0	0	0	0
4	H	4	0	0	0	0
4	L	231	0	0	4	0
All	All	7028	0	6602	100	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 7.

All (100) 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:2:LEU:CD2	1:A:27:GLN:HG2	2.01	0.91
1:L:77:ARG:HH11	1:L:77:ARG:HG3	1.38	0.89
1:L:77:ARG:HH11	1:L:77:ARG:CG	1.89	0.84
1:L:112:ALA:HB2	1:L:200:THR:HG21	1.58	0.84
1:A:2:LEU:HD11	1:A:25:SER:HB2	1.58	0.84
2:H:33:TRP:CH2	3:H:500:DIK:H16A	2.13	0.82
2:B:5:GLU:HA	2:B:105:GLN:HE22	1.43	0.81
2:B:136:SER:HA	2:B:195:SER:HB2	1.65	0.79
4:L:276:HOH:O	2:H:108:LEU:HD13	1.85	0.76
1:L:2:LEU:HD12	1:L:27:GLN:HG2	1.68	0.75
1:A:24:ARG:NH1	1:A:70:ASP:OD1	2.20	0.74
2:B:105:GLN:HE21	2:B:105:GLN:H	1.35	0.74
1:L:142:LYS:O	1:L:163:TRP:CZ3	2.41	0.73
1:A:2:LEU:HD23	1:A:27:GLN:HG2	1.71	0.73
1:L:2:LEU:CD1	1:L:27:GLN:HG2	2.19	0.72
1:L:144:ILE:HG22	1:L:163:TRP:CH2	2.23	0.72
2:H:83:ARG:HD3	2:H:85:GLU:OE2	1.89	0.72
1:L:2:LEU:HD22	1:L:90:GLN:HG2	1.78	0.66
1:L:77:ARG:HG3	1:L:77:ARG:NH1	2.13	0.63
1:A:122:SER:O	1:A:126:THR:HG23	1.99	0.63
2:B:41:PRO:HD2	2:B:42:GLU:OE2	1.98	0.63
2:H:27:LEU:HD11	2:H:94:ILE:HD13	1.80	0.62
2:B:97:TYR:CZ	3:B:500:DIK:H15A	2.36	0.61
2:B:136:SER:CA	2:B:195:SER:HB2	2.32	0.60
1:L:91:GLY:HA3	3:H:500:DIK:H5A	1.83	0.60
1:L:17:ASP:O	1:L:76:SER:O	2.20	0.59
1:A:190:ASN:HD21	1:A:212:ASN:H	1.50	0.59
2:B:5:GLU:HA	2:B:105:GLN:NE2	2.17	0.59
1:L:197:THR:HG22	1:L:204:PRO:HB3	1.85	0.58
2:B:136:SER:C	2:B:195:SER:HB2	2.29	0.58
1:A:214:CYS:SG	2:B:127:VAL:HG13	2.45	0.57
1:L:2:LEU:HD21	1:L:25:SER:HB2	1.87	0.57
1:A:112:ALA:HA	1:A:200:THR:HG21	1.87	0.56
1:L:142:LYS:O	1:L:163:TRP:HZ3	1.86	0.56
2:H:17:SER:HB2	2:H:82(A):ASN:HD22	1.70	0.56
1:A:2:LEU:HD22	1:A:27:GLN:HG2	1.83	0.56
1:L:142:LYS:O	1:L:163:TRP:CH2	2.59	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:284:HOH:O	2:B:108:LEU:HD13	2.06	0.56
2:B:126:PRO:HD3	2:B:140:LEU:HD12	1.90	0.54
2:B:136:SER:HA	2:B:195:SER:CB	2.36	0.54
1:L:59:PRO:HB2	1:L:61:ARG:HG3	1.89	0.54
1:L:130:ALA:O	1:L:181:LEU:HD23	2.06	0.54
1:L:182:THR:OG1	1:L:185:GLU:HG3	2.07	0.53
2:B:43:LYS:HE2	2:B:46:GLU:OE1	2.08	0.53
2:B:200:PRO:HB3	2:B:227:PRO:HG3	1.91	0.53
1:A:120:PRO:O	2:B:228:ARG:NH2	2.42	0.52
2:B:140:LEU:HG	2:B:223:ILE:HG21	1.93	0.51
2:B:18:MET:HB3	2:B:82:MET:HE3	1.93	0.51
1:L:18:GLN:HG3	1:L:76:SER:HA	1.92	0.51
2:B:179:GLN:O	2:B:180:SER:HB2	2.10	0.51
2:B:105:GLN:H	2:B:105:GLN:NE2	2.07	0.50
1:L:91:GLY:HA3	3:H:500:DIK:C5	2.40	0.50
1:L:1:GLU:HB2	4:L:235:HOH:O	2.11	0.49
2:B:93:LYS:NZ	3:B:500:DIK:O1	2.45	0.49
1:A:136:LEU:CD2	1:A:196:ALA:HB2	2.43	0.49
1:A:185:GLU:O	1:A:189:HIS:HD2	1.96	0.49
1:L:2:LEU:HD12	1:L:27:GLN:CG	2.42	0.49
2:H:61:GLU:O	2:H:64:LYS:HG2	2.13	0.48
2:B:34:MET:HE3	2:B:94:ILE:HB	1.95	0.48
1:L:14:SER:O	1:L:17:ASP:HB2	2.15	0.47
2:H:75:LYS:HE2	2:H:77:ARG:HE	1.80	0.47
2:B:33:TRP:CH2	3:B:500:DIK:H16A	2.50	0.47
1:L:181:LEU:HD23	1:L:181:LEU:H	1.79	0.47
1:L:144:ILE:HG22	1:L:163:TRP:HH2	1.77	0.47
1:L:2:LEU:CD2	1:L:4:MET:SD	3.02	0.46
2:B:12:VAL:HG11	2:B:82(C):LEU:HD13	1.98	0.46
3:B:500:DIK:H10	3:B:500:DIK:H5	1.59	0.46
2:H:81:GLN:HE21	2:H:82(A):ASN:HD21	1.64	0.46
2:H:83:ARG:NH1	2:H:85:GLU:OE2	2.50	0.45
1:A:98:PHE:CE2	3:B:500:DIK:H1B	2.52	0.45
1:A:198:HIS:ND1	1:A:200:THR:HB	2.32	0.45
3:H:500:DIK:H10	3:H:500:DIK:H5	1.80	0.45
4:L:437:HOH:O	3:B:500:DIK:H14	2.17	0.45
1:L:122:SER:O	1:L:126:THR:HG23	2.17	0.44
2:H:199:TRP:CG	2:H:200:PRO:HA	2.52	0.44
2:H:20:LEU:HG	2:H:82:MET:HE2	2.00	0.43
1:L:54:ARG:HD2	1:L:59:PRO:O	2.18	0.43
1:A:113:PRO:HG2	1:A:205:ILE:HD12	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:2:LEU:HD11	1:L:27:GLN:HG2	1.97	0.43
1:L:37:LEU:HD13	1:L:86:TYR:CZ	2.54	0.43
2:H:81:GLN:HE21	2:H:82(A):ASN:ND2	2.16	0.43
2:H:216:SER:O	1:A:1:GLU:N	2.43	0.43
1:L:83:LEU:HD21	1:L:168:SER:HB3	2.01	0.42
2:B:27:LEU:HD11	2:B:94:ILE:HD13	2.02	0.42
1:L:13:VAL:HG21	1:L:19:ALA:HB2	2.02	0.42
2:H:125:ALA:O	2:H:228:ARG:NH1	2.52	0.42
1:A:150:ILE:HD11	1:A:179:LEU:HD21	2.02	0.42
1:L:181:LEU:N	1:L:181:LEU:CD2	2.83	0.42
2:B:83:ARG:HD3	2:B:85:GLU:OE2	2.20	0.42
2:B:200:PRO:CB	2:B:227:PRO:HG3	2.49	0.41
1:L:181:LEU:HD23	1:L:181:LEU:N	2.36	0.41
2:H:27:LEU:C	2:H:27:LEU:HD12	2.46	0.41
2:H:121:VAL:HG12	2:H:221:LYS:HG3	2.01	0.41
2:B:202:SER:OG	2:B:203:GLN:HG3	2.21	0.41
1:L:6:GLN:HB2	1:L:23:CYS:SG	2.61	0.41
1:A:74:LYS:HE3	1:A:74:LYS:HB2	1.89	0.41
1:L:77:ARG:CG	1:L:77:ARG:NH1	2.60	0.40
1:L:125:LEU:O	1:L:183:LYS:HD2	2.20	0.40
1:A:6:GLN:HG3	1:A:23:CYS:SG	2.61	0.40
2:H:22:CYS:HB3	2:H:78:LEU:HB3	2.03	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	217/219 (99%)	212 (98%)	4 (2%)	1 (0%)	24	16
1	L	217/219 (99%)	214 (99%)	2 (1%)	1 (0%)	24	16
2	B	216/218 (99%)	214 (99%)	1 (0%)	1 (0%)	24	16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	216/218 (99%)	212 (98%)	3 (1%)	1 (0%)	24	16
All	All	866/874 (99%)	852 (98%)	10 (1%)	4 (0%)	24	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	180	SER
1	L	76	SER
2	H	133	THR
1	A	68	GLY

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	196/196 (100%)	185 (94%)	11 (6%)	19	11
1	L	196/196 (100%)	187 (95%)	9 (5%)	24	16
2	B	191/191 (100%)	183 (96%)	8 (4%)	26	19
2	H	191/191 (100%)	179 (94%)	12 (6%)	16	8
All	All	774/774 (100%)	734 (95%)	40 (5%)	21	13

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	33	LEU
1	L	77	ARG
1	L	83	LEU
1	L	145	ASN
1	L	160	LEU
1	L	181	LEU
1	L	197	THR
1	L	200	THR
1	L	202	THR
2	H	1	GLU

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Mol	Chain	Res	Type
2	H	27	LEU
2	H	42	GLU
2	H	54	ASN
2	H	74	SER
2	H	83	ARG
2	H	93	LYS
2	H	108	LEU
2	H	115	LYS
2	H	137	SER
2	H	163	SER
2	H	218	LYS
1	A	1	GLU
1	A	2	LEU
1	A	14	SER
1	A	17	ASP
1	A	22	SER
1	A	83	LEU
1	A	122	SER
1	A	123	GLU
1	A	142	LYS
1	A	188	ARG
1	A	200	THR
2	B	42	GLU
2	B	93	LYS
2	B	105	GLN
2	B	108	LEU
2	B	136	SER
2	B	140	LEU
2	B	217	THR
2	B	219	VAL

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

Mol	Chain	Res	Type
1	L	27(D)	HIS
1	L	38	GLN
1	L	53	ASN
1	L	137	ASN
1	L	157	ASN
1	L	189	HIS
2	H	13	GLN
2	H	39	GLN

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Mol	Chain	Res	Type
2	H	82(A)	ASN
1	A	137	ASN
1	A	189	HIS
1	A	190	ASN
2	B	81	GLN
2	B	105	GLN
2	B	179	GLN
2	B	203	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](#)

2 ligands are 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	DIK	H	500	2	22,22,23	0.88	1 (4%)	26,27,29	1.31	4 (15%)
3	DIK	B	500	2	22,22,23	0.66	0	26,27,29	1.16	1 (3%)

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	DIK	H	500	2	-	11/18/18/19	0/1/1/1
3	DIK	B	500	2	-	9/18/18/19	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	500	DIK	C14-C13	2.78	1.56	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	500	DIK	C1-C2-C3	-4.19	116.85	125.28
3	B	500	DIK	C1-C2-C3	-3.36	118.53	125.28
3	H	500	DIK	C12-N1-C13	-2.87	122.45	127.52
3	H	500	DIK	C15-C16-C17	-2.19	108.79	114.51
3	H	500	DIK	C14-C13-N1	2.11	118.26	114.59

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	500	DIK	C13-C14-C15-C16
3	H	500	DIK	C4-C5-C6-C7
3	B	500	DIK	C4-C5-C6-C7
3	B	500	DIK	C11-C12-N1-C13
3	H	500	DIK	C14-C15-C16-C17
3	B	500	DIK	C9-C12-N1-C13
3	B	500	DIK	C14-C15-C16-C17
3	H	500	DIK	N1-C13-C14-C15
3	H	500	DIK	O2-C13-C14-C15
3	B	500	DIK	N1-C13-C14-C15
3	B	500	DIK	O2-C13-C14-C15
3	H	500	DIK	C11-C12-N1-C13
3	H	500	DIK	C9-C12-N1-C13
3	H	500	DIK	C5-C6-C7-C10
3	H	500	DIK	C15-C16-C17-O3
3	H	500	DIK	C15-C16-C17-O4
3	H	500	DIK	C5-C6-C7-C8
3	B	500	DIK	C15-C16-C17-O4

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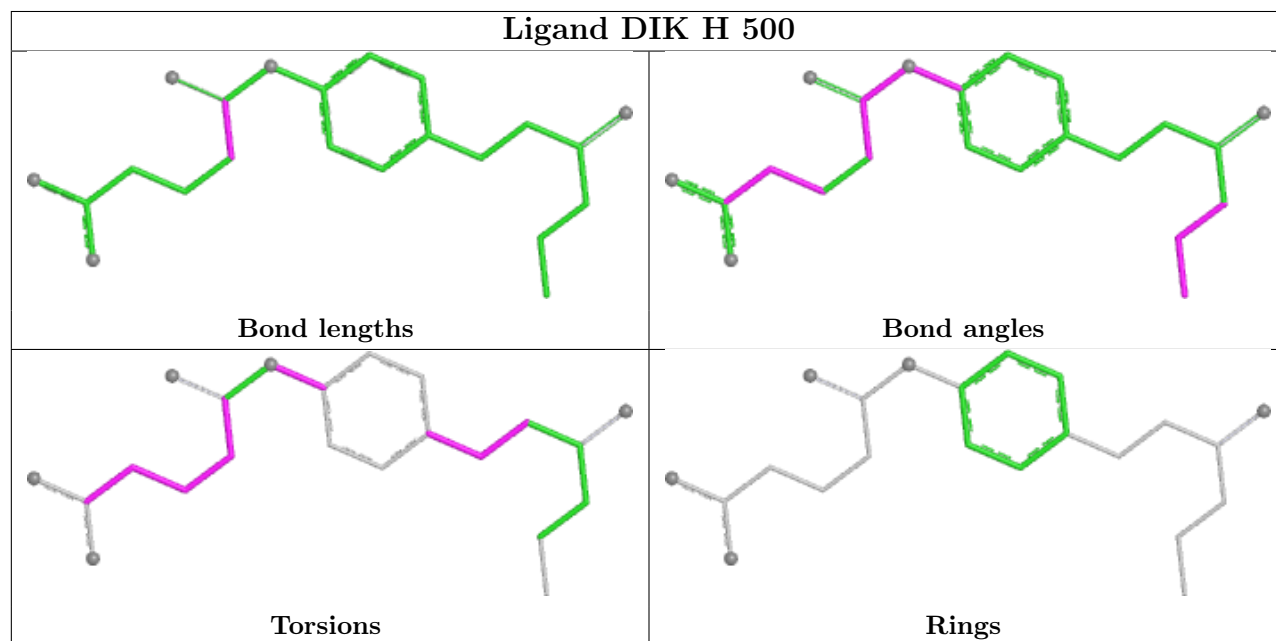
Mol	Chain	Res	Type	Atoms
3	H	500	DIK	C13-C14-C15-C16
3	B	500	DIK	C15-C16-C17-O3

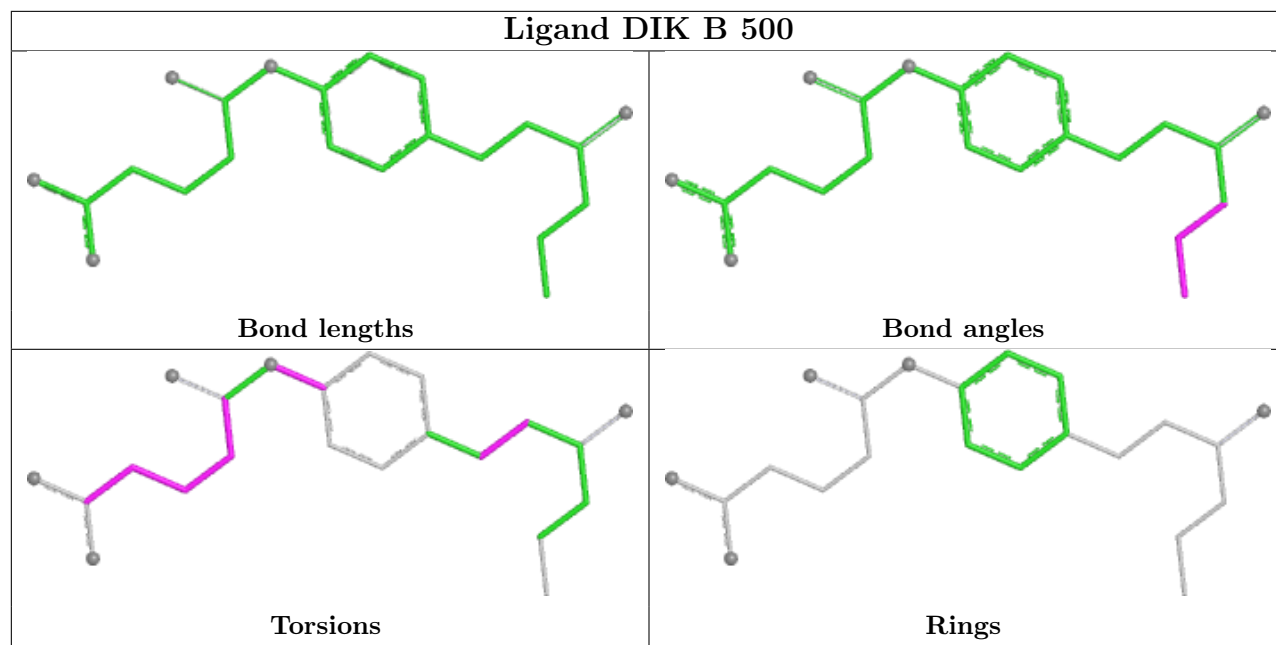
There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	500	DIK	4	0
3	B	500	DIK	6	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	219/219 (100%)	0.68	10 (4%) 37 40	34, 43, 51, 67	0
1	L	219/219 (100%)	0.67	18 (8%) 17 18	29, 39, 50, 71	0
2	B	218/218 (100%)	0.61	8 (3%) 45 48	32, 39, 52, 58	0
2	H	218/218 (100%)	0.64	12 (5%) 30 32	27, 42, 52, 67	0
All	All	874/874 (100%)	0.65	48 (5%) 30 32	27, 41, 51, 71	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	133	THR	5.8
1	A	28	TYR	5.3
1	L	76	SER	4.5
1	A	27(E)	SER	4.1
2	H	130	ASP	3.9
1	L	23	CYS	3.9
1	A	23	CYS	3.9
1	A	214	CYS	3.9
1	L	2	LEU	3.6
1	L	163	TRP	3.5
2	H	134	THR	3.3
2	B	134	THR	3.2
2	H	128	CYS	3.2
2	H	56	ALA	3.1
1	A	194	CYS	3.0
2	H	129	GLY	3.0
1	L	16	GLY	2.9
1	A	2	LEU	2.9
2	B	136	SER	2.9
1	A	29	GLY	2.9
1	L	127	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	200	THR	2.8
1	L	78	VAL	2.8
1	L	214	CYS	2.7
1	L	194	CYS	2.7
1	A	24	ARG	2.6
1	L	126	THR	2.6
1	L	202	THR	2.6
1	L	77	ARG	2.4
2	H	42	GLU	2.4
2	B	208	CYS	2.4
2	H	135	GLY	2.4
1	L	67	SER	2.3
2	B	128	CYS	2.3
2	B	127	VAL	2.3
2	H	27	LEU	2.3
1	L	58	VAL	2.2
1	L	128	GLY	2.2
2	H	195	SER	2.2
2	H	4	LEU	2.2
2	H	166	LEU	2.2
2	B	202	SER	2.2
1	A	143	ASP	2.2
1	L	13	VAL	2.1
2	B	54	ASN	2.1
1	L	200	THR	2.0
1	L	60	ASP	2.0
2	B	180	SER	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,

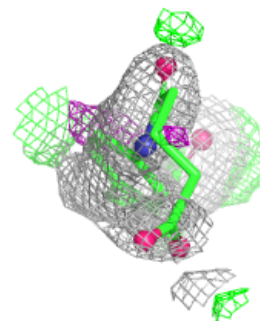
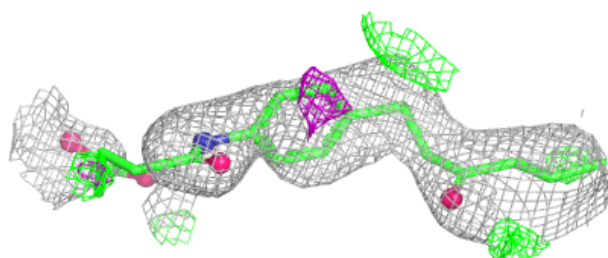
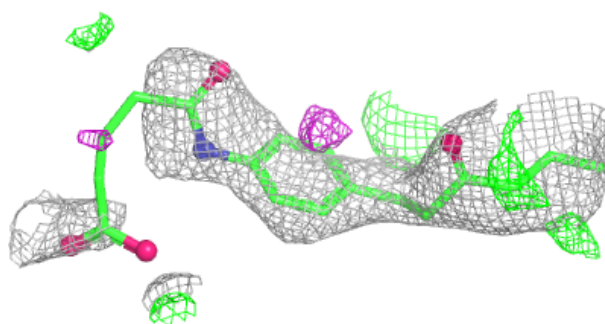
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	DIK	H	500	22/23	0.70	0.21	68,74,84,85	0
3	DIK	B	500	22/23	0.73	0.20	67,72,80,80	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.

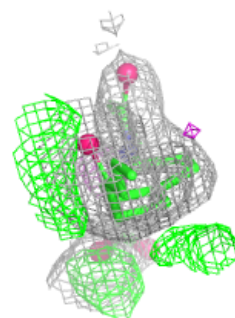
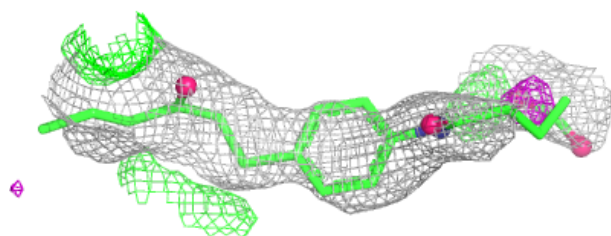
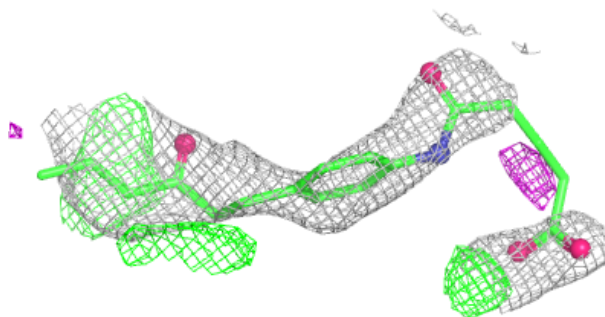
Electron density around DIK H 500:

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 DIK B 500:

$2mF_o-DF_c$ (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.