Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 01:57 AM EDT

PDB ID	:	2DHF
Title	:	CRYSTAL STRUCTURES OF RECOMBINANT HUMAN DIHYDROFOLATE REDUCTASE COMPLEXED WITH FOLATE AND 5-DEAZOFOLATE
Authors	:	Davies /II, J.F.; Kraut, J.
Deposited on	:	1989-10-25
Resolution	:	2.30 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 (i) 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 (1)) were used in the production of this report:

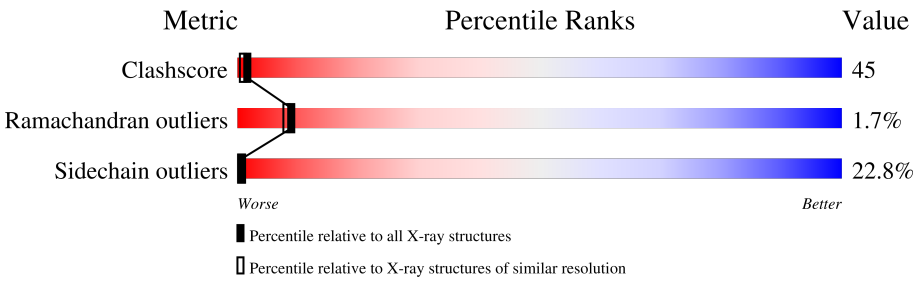
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	186	30% 46% 19% • •
1	B	186	30% 47% 18% • •

2 Entry composition (i)

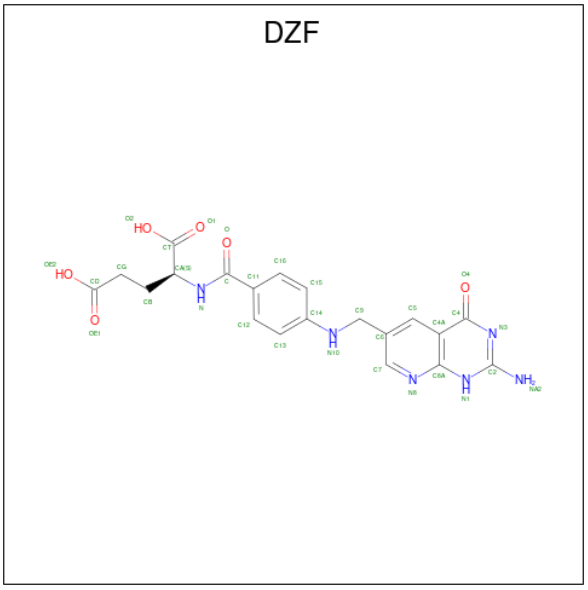
There are 3 unique types of molecules in this entry. The entry contains 3123 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 DIHYDROFOLATE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	182	1476	949	249	271	7	0	0	0
1	B	182	1472	947	249	269	7	0	0	0

- Molecule 2 is 5-DEAZAFOLIC ACID (three-letter code: DZF) (formula: $C_{20}H_{20}N_6O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	32	20	6	6	0	0
2	B	1	32	20	6	6	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	52	52	52	0	0
			Total	O		
3	B	59	59	59	0	0

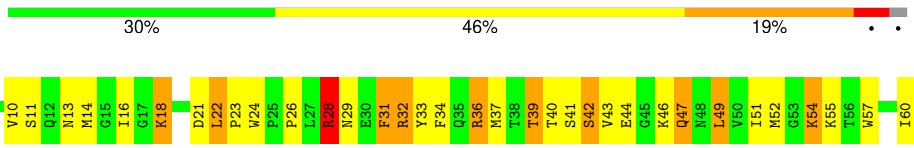
3 Residue-property plots (i)

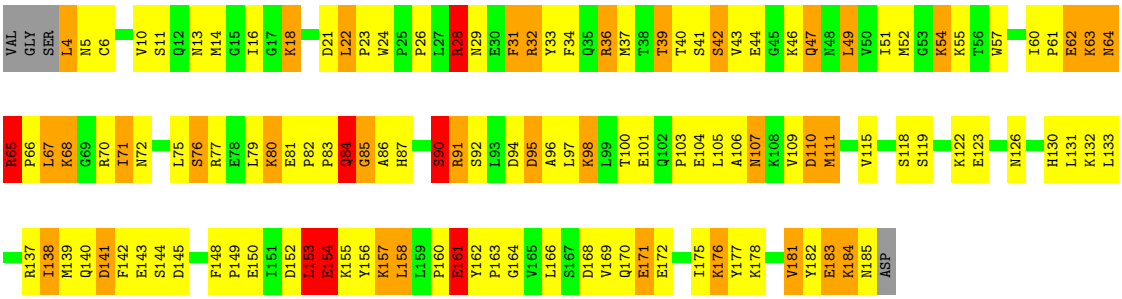
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

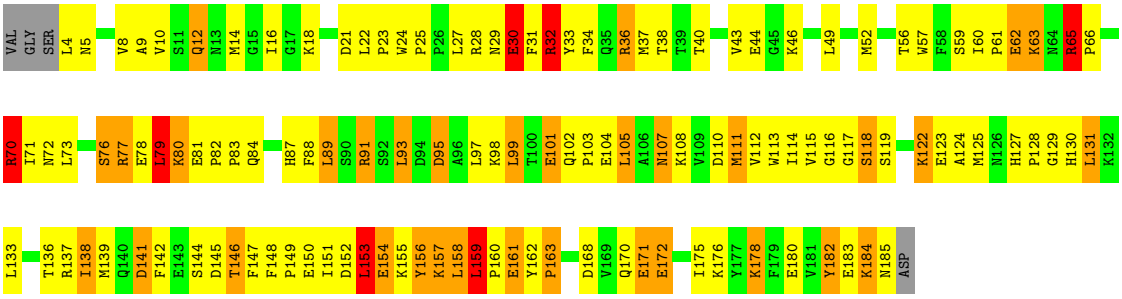
Chain A:





• Molecule 1: DIHYDROFOLATE REDUCTASE

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	36.10Å 38.50Å 76.80Å 93.90° 91.40° 111.50°	Depositor
Resolution (Å)	(Not available) – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.194 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3123	wwPDB-VP
Average B, all atoms ($Å^2$)	44.0	wwPDB-VP

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: DZF

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	1.12	2/1511 (0.1%)	1.84	28/2039 (1.4%)
1	B	1.20	1/1507 (0.1%)	1.96	41/2034 (2.0%)
All	All	1.16	3/3018 (0.1%)	1.90	69/4073 (1.7%)

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	1	4
1	B	0	5
All	All	1	9

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	70	ARG	CD-NE	-6.88	1.34	1.46
1	A	92	SER	CB-OG	-6.82	1.33	1.42
1	A	144	SER	CB-OG	5.53	1.49	1.42

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	70	ARG	CD-NE-CZ	17.64	148.30	123.60
1	A	65	ARG	NE-CZ-NH2	-16.75	111.92	120.30
1	A	28	ARG	NE-CZ-NH1	9.67	125.14	120.30
1	B	95	ASP	CB-CG-OD1	9.64	126.98	118.30
1	B	70	ARG	NE-CZ-NH2	9.39	124.99	120.30
1	B	36	ARG	NE-CZ-NH1	9.30	124.95	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	141	ASP	CB-CG-OD1	9.21	126.59	118.30
1	B	30	GLU	CA-CB-CG	8.55	132.22	113.40
1	B	36	ARG	NE-CZ-NH2	-8.29	116.15	120.30
1	B	65	ARG	NE-CZ-NH1	-8.28	116.16	120.30
1	A	110	ASP	CB-CG-OD1	8.15	125.64	118.30
1	B	144	SER	N-CA-CB	-7.96	98.56	110.50
1	A	4	LEU	CA-CB-CG	7.53	132.62	115.30
1	B	95	ASP	CB-CG-OD2	-7.13	111.88	118.30
1	B	9	ALA	CB-CA-C	-7.12	99.42	110.10
1	B	65	ARG	NE-CZ-NH2	7.07	123.84	120.30
1	B	93	LEU	CB-CA-C	6.95	123.40	110.20
1	B	93	LEU	CA-CB-CG	6.67	130.65	115.30
1	A	67	LEU	CB-CA-C	6.66	122.85	110.20
1	B	182	TYR	N-CA-CB	6.61	122.50	110.60
1	B	44	GLU	CA-CB-CG	6.54	127.78	113.40
1	A	153	LEU	CB-CA-C	6.50	122.56	110.20
1	A	65	ARG	NE-CZ-NH1	6.44	123.52	120.30
1	A	161	GLU	CA-CB-CG	6.39	127.46	113.40
1	B	111	MET	CA-CB-CG	6.22	123.88	113.30
1	A	28	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	B	97	LEU	CB-CA-C	6.18	121.94	110.20
1	A	84	GLN	CB-CA-C	-6.14	98.12	110.40
1	B	145	ASP	CB-CG-OD1	6.09	123.78	118.30
1	A	62	GLU	CA-CB-CG	6.07	126.75	113.40
1	B	77	ARG	NE-CZ-NH2	-6.07	117.27	120.30
1	A	181	VAL	CA-CB-CG2	6.06	119.99	110.90
1	A	31	PHE	CB-CG-CD2	-6.04	116.57	120.80
1	B	168	ASP	CB-CG-OD1	6.03	123.73	118.30
1	A	95	ASP	CB-CG-OD1	6.03	123.73	118.30
1	A	90	SER	N-CA-CB	6.02	119.53	110.50
1	A	111	MET	CA-CB-CG	5.99	123.48	113.30
1	B	159	LEU	CA-CB-CG	5.88	128.81	115.30
1	A	21	ASP	CB-CG-OD1	5.87	123.58	118.30
1	B	172	GLU	CA-CB-CG	5.84	126.25	113.40
1	B	76	SER	N-CA-CB	5.76	119.14	110.50
1	B	29	ASN	CA-CB-CG	5.74	126.02	113.40
1	B	171	GLU	CB-CA-C	-5.72	98.96	110.40
1	B	154	GLU	CA-CB-CG	5.70	125.94	113.40
1	B	65	ARG	N-CA-CB	5.69	120.84	110.60
1	B	118	SER	N-CA-CB	5.65	118.97	110.50
1	B	59	SER	N-CA-CB	5.62	118.93	110.50
1	A	39	THR	CA-CB-CG2	-5.54	104.64	112.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	156	TYR	CB-CG-CD2	-5.54	117.67	121.00
1	A	22	LEU	CB-CA-C	5.54	120.72	110.20
1	B	30	GLU	CB-CA-C	5.53	121.46	110.40
1	A	171	GLU	OE1-CD-OE2	5.44	129.82	123.30
1	B	99	LEU	CB-CA-C	5.41	120.49	110.20
1	A	110	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	A	183	GLU	N-CA-CB	5.36	120.24	110.60
1	A	171	GLU	CG-CD-OE2	-5.30	107.69	118.30
1	B	110	ASP	CB-CG-OD1	-5.25	113.58	118.30
1	B	129	GLY	C-N-CA	5.22	134.75	121.70
1	B	185	ASN	CA-C-O	5.16	130.93	120.10
1	B	71	ILE	N-CA-C	-5.15	97.08	111.00
1	B	141	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	B	79	LEU	O-C-N	5.14	130.92	122.70
1	A	182	TYR	CB-CG-CD2	-5.11	117.93	121.00
1	B	142	PHE	CB-CA-C	5.11	120.61	110.40
1	A	21	ASP	CB-CG-OD2	-5.04	113.77	118.30
1	B	153	LEU	CB-CA-C	5.04	119.77	110.20
1	A	154	GLU	CA-CB-CG	5.01	124.42	113.40
1	A	60	ILE	CB-CA-C	-5.00	101.59	111.60
1	A	92	SER	N-CA-CB	-5.00	102.99	110.50

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	4	LEU	CA

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	28	ARG	Sidechain
1	A	32	ARG	Sidechain
1	A	36	ARG	Sidechain
1	A	91	ARG	Sidechain
1	B	137	ARG	Sidechain
1	B	28	ARG	Sidechain
1	B	32	ARG	Sidechain
1	B	65	ARG	Sidechain
1	B	91	ARG	Sidechain

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	1476	0	1487	140	1
1	B	1472	0	1483	135	1
2	A	32	0	18	0	0
2	B	32	0	17	4	0
3	A	52	0	0	6	0
3	B	59	0	0	16	0
All	All	3123	0	3005	270	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 45.

All (270) 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:171:GLU:OE2	1:A:176:LYS:HE2	1.38	1.22
1:B:80:LYS:NZ	1:B:80:LYS:HA	1.56	1.19
1:B:80:LYS:NZ	1:B:80:LYS:CA	2.10	1.15
1:B:80:LYS:HA	1:B:80:LYS:HZ2	1.08	1.13
1:B:80:LYS:CA	1:B:80:LYS:HZ2	1.63	1.11
1:A:66:PRO:O	1:A:68:LYS:HD2	1.51	1.09
1:B:80:LYS:HE3	1:B:80:LYS:C	1.74	1.08
1:B:130:HIS:HE1	1:B:183:GLU:OE1	1.35	1.07
1:A:156:TYR:C	1:A:157:LYS:HD3	1.76	1.05
1:A:14:MET:HE2	1:A:14:MET:HA	1.40	1.02
1:B:76:SER:O	1:B:91:ARG:HD3	1.59	1.01
1:A:64:ASN:N	1:A:64:ASN:HD22	1.58	1.01
1:B:107:ASN:HB3	3:B:493:HOH:O	1.58	1.01
1:A:77:ARG:HA	1:A:91:ARG:HD2	1.44	0.99
1:A:156:TYR:O	1:A:157:LYS:HD3	1.65	0.96
1:B:130:HIS:CE1	1:B:183:GLU:OE1	2.19	0.96
1:B:80:LYS:C	1:B:80:LYS:CE	2.35	0.95
1:B:104:GLU:HA	3:B:468:HOH:O	1.66	0.94
1:A:153:LEU:O	1:A:153:LEU:CD1	2.15	0.94
1:A:153:LEU:O	1:A:153:LEU:HD13	1.67	0.94
1:B:146:THR:HG21	3:B:408:HOH:O	1.65	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:ASN:N	1:A:64:ASN:ND2	2.11	0.92
1:B:23:PRO:HG2	1:B:24:TRP:CZ3	2.05	0.92
1:B:146:THR:HG23	3:B:421:HOH:O	1.70	0.91
1:B:81:GLU:HB3	1:B:82:PRO:HD2	1.52	0.90
1:B:22:LEU:HD12	3:B:428:HOH:O	1.70	0.90
1:B:80:LYS:HE3	1:B:80:LYS:O	1.72	0.89
1:A:160:PRO:HG2	1:A:161:GLU:OE2	1.71	0.89
1:A:54:LYS:HG3	1:A:79:LEU:HD11	1.60	0.84
1:A:171:GLU:OE2	1:A:176:LYS:CE	2.25	0.84
1:B:139:MET:HG2	3:B:501:HOH:O	1.78	0.84
1:B:24:TRP:HB2	1:B:25:PRO:HD2	1.60	0.83
1:B:99:LEU:HD11	1:B:105:LEU:CD2	2.07	0.83
1:B:99:LEU:HD11	1:B:105:LEU:HD21	1.60	0.82
1:B:104:GLU:HG2	3:B:468:HOH:O	1.80	0.80
1:A:14:MET:HA	1:A:14:MET:CE	2.10	0.80
1:B:99:LEU:CD1	1:B:105:LEU:CD2	2.60	0.79
1:A:98:LYS:O	1:A:101:GLU:CB	2.31	0.78
1:A:94:ASP:O	1:A:98:LYS:HD3	1.83	0.78
1:B:80:LYS:HA	1:B:80:LYS:HZ1	1.46	0.78
1:B:89:LEU:HD21	1:B:91:ARG:NH2	1.99	0.77
1:B:118:SER:HB2	3:B:440:HOH:O	1.84	0.77
1:B:80:LYS:NZ	1:B:80:LYS:C	2.36	0.77
1:B:80:LYS:CE	1:B:80:LYS:O	2.32	0.77
1:A:64:ASN:HD22	1:A:64:ASN:H	1.31	0.76
1:B:99:LEU:O	1:B:105:LEU:HD23	1.85	0.75
1:B:76:SER:O	1:B:91:ARG:CD	2.34	0.75
1:A:171:GLU:HG3	1:A:176:LYS:HD3	1.69	0.75
1:A:98:LYS:O	1:A:101:GLU:HB2	1.86	0.75
1:B:40:THR:O	1:B:111:MET:HE1	1.86	0.75
1:A:80:LYS:HA	1:A:80:LYS:HE2	1.69	0.75
1:B:81:GLU:CB	1:B:82:PRO:HD2	2.16	0.75
1:A:68:LYS:N	1:A:70:ARG:HH11	1.86	0.74
1:A:54:LYS:HG3	1:A:79:LEU:CD1	2.18	0.73
1:A:66:PRO:O	1:A:68:LYS:CD	2.33	0.73
1:A:6:CYS:HB2	1:A:133:LEU:HD23	1.70	0.72
1:A:84:GLN:CG	1:A:85:GLY:N	2.49	0.72
1:A:153:LEU:O	1:A:153:LEU:HD12	1.88	0.72
1:A:68:LYS:H	1:A:70:ARG:HH11	1.34	0.71
1:B:89:LEU:HD21	1:B:91:ARG:HH22	1.53	0.71
1:B:114:ILE:HD13	1:B:124:ALA:CB	2.18	0.71
1:B:154:GLU:CD	3:B:487:HOH:O	2.28	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:GLU:O	1:B:79:LEU:C	2.27	0.71
1:A:154:GLU:OE1	1:A:155:LYS:HD3	1.92	0.70
1:A:80:LYS:HE2	1:A:80:LYS:O	1.91	0.70
1:A:152:ASP:OD1	1:A:154:GLU:N	2.25	0.70
1:B:136:THR:HG22	1:B:138:ILE:CD1	2.22	0.69
1:B:23:PRO:HG2	1:B:24:TRP:CE3	2.27	0.69
1:B:81:GLU:HB3	1:B:82:PRO:CD	2.22	0.69
1:A:76:SER:O	1:A:91:ARG:HD3	1.93	0.69
1:B:62:GLU:CG	1:B:62:GLU:O	2.39	0.69
1:A:14:MET:HE1	3:A:426:HOH:O	1.92	0.69
1:A:111:MET:HE1	3:A:404:HOH:O	1.93	0.69
1:B:79:LEU:HD13	1:B:83:PRO:HD3	1.74	0.69
1:B:159:LEU:HD21	1:B:183:GLU:HB3	1.73	0.68
1:B:40:THR:O	1:B:111:MET:CE	2.42	0.68
1:A:84:GLN:O	1:A:84:GLN:HG2	1.76	0.68
1:A:95:ASP:HA	1:A:98:LYS:HZ3	1.59	0.67
1:B:79:LEU:CD1	1:B:83:PRO:HD3	2.24	0.67
1:A:18:LYS:HD2	1:A:143:GLU:O	1.95	0.67
1:B:122:LYS:NZ	3:B:440:HOH:O	2.26	0.67
1:A:61:PRO:HB2	1:A:64:ASN:HD21	1.58	0.67
1:B:80:LYS:CA	1:B:80:LYS:HZ1	1.99	0.67
1:B:114:ILE:HD13	1:B:124:ALA:HB2	1.76	0.67
1:A:137:ARG:O	1:A:177:TYR:HA	1.97	0.65
1:B:76:SER:HB3	1:B:79:LEU:HB2	1.79	0.65
1:A:171:GLU:HG3	1:A:176:LYS:CD	2.27	0.64
1:A:80:LYS:O	1:A:80:LYS:CE	2.45	0.64
1:B:80:LYS:CA	1:B:80:LYS:CE	2.76	0.64
1:B:62:GLU:O	1:B:62:GLU:HG2	1.98	0.63
1:A:71:ILE:CD1	1:A:71:ILE:N	2.62	0.63
1:A:80:LYS:HE2	1:A:80:LYS:CA	2.28	0.63
1:A:183:GLU:HG2	1:A:184:LYS:N	2.13	0.63
1:A:139:MET:HE2	1:A:177:TYR:HA	1.80	0.62
1:A:18:LYS:HA	1:A:145:ASP:OD1	1.98	0.62
1:A:95:ASP:HA	1:A:98:LYS:NZ	2.13	0.62
1:A:171:GLU:HA	1:A:175:ILE:O	1.99	0.62
1:B:61:PRO:O	1:B:63:LYS:N	2.30	0.61
1:A:14:MET:CE	1:A:14:MET:CA	2.78	0.61
1:A:68:LYS:H	1:A:70:ARG:NH1	1.99	0.61
1:A:170:GLN:O	1:A:176:LYS:HA	2.01	0.61
1:A:139:MET:CE	1:A:177:TYR:HA	2.30	0.61
1:B:151:ILE:O	1:B:153:LEU:HD22	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:PRO:O	1:A:64:ASN:ND2	2.34	0.61
1:A:83:PRO:O	1:A:84:GLN:O	2.19	0.61
1:B:99:LEU:CD1	1:B:105:LEU:HD23	2.30	0.61
1:B:24:TRP:HB2	1:B:25:PRO:CD	2.29	0.61
1:B:81:GLU:CB	1:B:82:PRO:CD	2.78	0.61
1:A:71:ILE:N	1:A:71:ILE:HD13	2.15	0.60
1:A:104:GLU:OE1	1:A:104:GLU:HA	2.01	0.60
1:B:30:GLU:OE1	2:B:187:DZF:NA2	2.30	0.60
1:A:139:MET:CE	1:A:177:TYR:CA	2.79	0.60
1:B:139:MET:HB2	1:B:176:LYS:O	2.01	0.60
1:A:14:MET:HE2	1:A:14:MET:CA	2.23	0.60
1:A:63:LYS:CD	1:A:63:LYS:O	2.49	0.59
1:B:5:ASN:O	1:B:113:TRP:HA	2.02	0.59
1:B:98:LYS:O	1:B:101:GLU:HG2	2.02	0.59
1:B:171:GLU:HA	1:B:175:ILE:O	2.03	0.58
1:A:81:GLU:HB3	1:A:82:PRO:HD2	1.85	0.58
1:B:80:LYS:NZ	1:B:80:LYS:O	2.33	0.58
1:B:180:GLU:OE1	1:B:182:TYR:OH	2.15	0.58
1:A:130:HIS:NE2	1:A:183:GLU:OE1	2.24	0.58
1:B:22:LEU:HD12	3:B:454:HOH:O	2.03	0.58
1:A:98:LYS:O	1:A:101:GLU:HB3	2.04	0.58
1:A:63:LYS:O	1:A:63:LYS:HD3	2.04	0.57
1:A:23:PRO:HD2	1:A:24:TRP:CE3	2.40	0.57
1:A:157:LYS:HD3	1:A:157:LYS:N	2.12	0.57
1:A:161:GLU:OE2	1:A:161:GLU:N	2.38	0.56
1:B:176:LYS:NZ	3:B:445:HOH:O	2.38	0.56
1:A:84:GLN:C	1:A:86:ALA:H	2.08	0.56
1:B:93:LEU:HD23	1:B:123:GLU:HB3	1.88	0.56
1:A:36:ARG:HH12	1:A:164:GLY:C	2.10	0.55
1:B:73:LEU:HG	1:B:88:PHE:HB2	1.87	0.55
1:A:77:ARG:HG3	1:A:91:ARG:HB3	1.86	0.55
1:A:80:LYS:HE2	1:A:80:LYS:C	2.26	0.55
1:A:77:ARG:HA	1:A:91:ARG:CD	2.29	0.55
1:A:84:GLN:O	1:A:86:ALA:N	2.38	0.55
1:B:148:PHE:CD2	1:B:149:PRO:O	2.59	0.55
1:B:130:HIS:HE1	1:B:183:GLU:CD	2.07	0.55
1:A:75:LEU:HA	1:A:90:SER:O	2.06	0.54
1:A:72:ASN:O	1:A:87:HIS:HB2	2.07	0.54
1:A:57:TRP:CE2	1:A:65:ARG:HD3	2.42	0.54
1:A:10:VAL:HG12	1:A:16:ILE:CG2	2.38	0.54
1:B:22:LEU:CD1	3:B:454:HOH:O	2.53	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:MET:HB3	1:B:115:VAL:CG2	2.38	0.54
1:B:34:PHE:O	1:B:38:THR:HG23	2.08	0.53
1:B:80:LYS:C	1:B:80:LYS:HZ1	2.03	0.53
1:A:26:PRO:HA	3:A:488:HOH:O	2.08	0.53
1:B:72:ASN:O	1:B:87:HIS:HB2	2.09	0.53
1:A:84:GLN:HG2	1:A:85:GLY:N	1.93	0.52
1:A:29:ASN:HB2	1:A:172:GLU:OE1	2.10	0.52
1:A:158:LEU:HD23	1:A:181:VAL:O	2.09	0.51
1:B:61:PRO:HB2	1:B:63:LYS:HE3	1.92	0.51
1:A:43:VAL:HG11	1:A:46:LYS:HG3	1.92	0.51
1:A:6:CYS:HB2	1:A:133:LEU:CD2	2.40	0.51
1:B:4:LEU:HB3	1:B:131:LEU:HG	1.91	0.51
1:A:40:THR:O	1:A:111:MET:CE	2.58	0.51
1:A:11:SER:O	1:A:14:MET:HE2	2.10	0.51
1:A:154:GLU:OE1	1:A:154:GLU:O	2.28	0.51
1:A:140:GLN:HG2	1:A:141:ASP:N	2.25	0.50
1:A:13:ASN:ND2	3:A:411:HOH:O	2.44	0.50
1:A:183:GLU:HG2	1:A:184:LYS:H	1.77	0.50
1:B:104:GLU:CA	3:B:468:HOH:O	2.42	0.50
1:B:60:ILE:O	1:B:65:ARG:NH2	2.38	0.50
1:B:99:LEU:HD12	1:B:105:LEU:HD23	1.94	0.50
1:B:31:PHE:HB3	2:B:187:DZF:CD	2.42	0.50
1:B:56:THR:O	1:B:57:TRP:C	2.50	0.50
1:A:47:GLN:O	1:A:109:VAL:HA	2.12	0.50
1:A:51:ILE:CG2	1:A:75:LEU:HD11	2.42	0.50
1:B:104:GLU:CG	3:B:468:HOH:O	2.48	0.49
1:B:152:ASP:OD1	1:B:154:GLU:CB	2.60	0.49
1:B:161:GLU:OE1	1:B:163:PRO:HD3	2.12	0.49
1:A:139:MET:HE3	1:A:176:LYS:C	2.32	0.49
1:B:114:ILE:HD13	1:B:124:ALA:HB1	1.94	0.49
1:B:8:VAL:HG11	1:B:148:PHE:CE1	2.47	0.49
1:A:10:VAL:HG12	1:A:16:ILE:HG22	1.94	0.49
1:B:66:PRO:HB2	1:B:72:ASN:ND2	2.27	0.49
1:B:153:LEU:HD13	1:B:153:LEU:HA	1.48	0.49
1:A:11:SER:O	1:A:14:MET:CE	2.61	0.49
1:A:80:LYS:HE3	1:A:91:ARG:NH1	2.27	0.49
1:A:160:PRO:HG2	1:A:161:GLU:CD	2.32	0.48
1:B:66:PRO:HB2	1:B:72:ASN:CG	2.33	0.48
1:B:114:ILE:CD1	1:B:124:ALA:HB1	2.43	0.48
1:B:49:LEU:HD23	1:B:112:VAL:HG13	1.95	0.48
1:B:63:LYS:C	1:B:63:LYS:CD	2.82	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:LEU:HD23	1:B:4:LEU:HA	1.49	0.48
1:B:27:LEU:HD22	1:B:172:GLU:HG2	1.95	0.48
1:B:70:ARG:O	3:B:465:HOH:O	2.20	0.48
1:B:102:GLN:HB2	1:B:105:LEU:HB2	1.95	0.48
1:A:52:MET:HB3	1:A:115:VAL:CG2	2.45	0.47
1:A:148:PHE:CD1	1:A:149:PRO:HD2	2.50	0.47
1:B:99:LEU:HD12	1:B:105:LEU:CD2	2.42	0.47
1:B:158:LEU:HD23	1:B:158:LEU:HA	1.62	0.47
1:B:80:LYS:HE3	1:B:81:GLU:N	2.26	0.47
1:A:119:SER:N	3:A:456:HOH:O	2.49	0.46
1:B:117:GLY:O	1:B:118:SER:C	2.52	0.46
1:A:158:LEU:C	1:A:158:LEU:CD2	2.83	0.46
1:B:60:ILE:HG23	1:B:61:PRO:HD2	1.98	0.46
1:A:169:VAL:CG2	1:A:176:LYS:HG3	2.45	0.46
1:A:24:TRP:O	3:A:477:HOH:O	2.20	0.46
1:A:162:TYR:HA	1:A:163:PRO:HD2	1.69	0.46
1:A:96:ALA:O	1:A:100:THR:HG23	2.15	0.46
1:B:12:GLN:HB3	1:B:141:ASP:OD1	2.16	0.45
1:B:22:LEU:HD11	1:B:31:PHE:CZ	2.51	0.45
1:A:33:TYR:O	1:A:37:MET:HG2	2.16	0.45
1:B:127:HIS:HA	1:B:128:PRO:HD3	1.81	0.45
1:B:152:ASP:OD1	1:B:154:GLU:HB2	2.16	0.45
1:A:46:LYS:NZ	1:A:106:ALA:O	2.49	0.45
1:A:68:LYS:N	1:A:70:ARG:NH1	2.57	0.45
1:B:24:TRP:CB	1:B:25:PRO:CD	2.94	0.45
1:A:5:ASN:OD1	1:A:132:LYS:HB2	2.17	0.45
1:A:107:ASN:OD1	1:A:107:ASN:O	2.34	0.45
1:B:105:LEU:HD12	1:B:105:LEU:HA	1.65	0.45
1:A:40:THR:O	1:A:111:MET:HE3	2.17	0.45
1:A:183:GLU:CG	1:A:184:LYS:N	2.77	0.45
1:B:78:GLU:O	1:B:79:LEU:O	2.34	0.45
1:B:14:MET:O	1:B:147:PHE:HA	2.17	0.45
1:A:42:SER:N	1:A:110:ASP:OD2	2.49	0.44
1:A:172:GLU:N	1:A:175:ILE:O	2.44	0.44
1:A:139:MET:HE2	1:A:177:TYR:CA	2.44	0.44
1:B:33:TYR:O	1:B:37:MET:HG2	2.17	0.44
1:A:14:MET:CE	1:A:14:MET:N	2.80	0.44
1:A:83:PRO:O	1:A:84:GLN:C	2.56	0.44
1:A:138:ILE:O	1:A:138:ILE:HG22	2.18	0.44
1:A:166:LEU:N	1:A:166:LEU:HD23	2.33	0.44
1:A:152:ASP:OD1	1:A:152:ASP:C	2.55	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:LYS:C	1:A:157:LYS:HE2	2.39	0.43
1:B:125:MET:HE1	1:B:151:ILE:HA	2.00	0.43
1:B:157:LYS:HB2	1:B:157:LYS:HE3	1.61	0.43
1:A:87:HIS:ND1	1:B:104:GLU:CG	2.82	0.43
1:A:171:GLU:CG	1:A:176:LYS:CD	2.96	0.43
1:A:171:GLU:HB2	1:A:176:LYS:HD2	2.01	0.43
1:B:62:GLU:HA	1:B:65:ARG:HG3	2.00	0.43
1:A:67:LEU:HB3	1:A:70:ARG:CZ	2.49	0.43
1:A:80:LYS:NZ	1:B:98:LYS:CE	2.82	0.43
1:A:49:LEU:HD12	1:A:71:ILE:HB	2.01	0.43
1:A:153:LEU:CD1	1:A:153:LEU:C	2.76	0.43
1:B:52:MET:HB2	1:B:56:THR:HB	2.01	0.43
1:B:89:LEU:HD21	1:B:91:ARG:CZ	2.48	0.43
1:A:66:PRO:O	1:A:67:LEU:C	2.57	0.42
1:A:107:ASN:OD1	1:A:107:ASN:C	2.56	0.42
1:B:52:MET:HA	1:B:116:GLY:O	2.19	0.42
1:B:57:TRP:O	1:B:65:ARG:HD3	2.19	0.42
1:B:162:TYR:HA	1:B:163:PRO:HD2	1.81	0.42
1:B:43:VAL:HB	1:B:46:LYS:HB2	2.00	0.42
1:A:23:PRO:HB2	1:A:142:PHE:HB3	2.02	0.42
1:B:36:ARG:HH21	1:B:36:ARG:HD3	1.54	0.42
1:A:80:LYS:NZ	1:B:98:LYS:HE2	2.34	0.42
1:B:16:ILE:CG2	1:B:148:PHE:HB2	2.50	0.42
1:A:158:LEU:C	1:A:158:LEU:HD22	2.40	0.41
1:A:183:GLU:CG	1:A:184:LYS:H	2.33	0.41
1:B:21:ASP:OD2	1:B:22:LEU:N	2.51	0.41
1:B:115:VAL:O	2:B:187:DZF:H7	2.19	0.41
1:B:170:GLN:O	1:B:176:LYS:HA	2.19	0.41
1:A:22:LEU:HA	1:A:23:PRO:HD3	1.86	0.41
1:A:97:LEU:HD23	1:A:97:LEU:HA	1.78	0.41
1:A:103:PRO:HG2	1:B:87:HIS:O	2.19	0.41
1:A:80:LYS:HZ2	1:B:98:LYS:CE	2.33	0.41
1:B:178:LYS:HE3	1:B:178:LYS:HB3	1.79	0.41
1:A:43:VAL:CG1	1:A:46:LYS:HG3	2.50	0.41
1:A:83:PRO:C	1:A:84:GLN:O	2.59	0.41
1:B:34:PHE:CD2	1:B:34:PHE:C	2.94	0.41
1:A:169:VAL:HG23	1:A:176:LYS:HG3	2.02	0.41
1:B:31:PHE:HB3	2:B:187:DZF:CG	2.50	0.41
1:B:156:TYR:CZ	1:B:184:LYS:HD3	2.56	0.41
1:B:124:ALA:O	1:B:131:LEU:HD11	2.21	0.40
1:B:159:LEU:HA	1:B:160:PRO:HD2	1.82	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLU:OE1	1:A:154:GLU:C	2.59	0.40
1:A:94:ASP:C	1:A:98:LYS:HD3	2.40	0.40
1:B:10:VAL:HG12	1:B:16:ILE:HG22	2.03	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:141:ASP:OD1	1:B:32:ARG:NH1[1_554]	2.08	0.12

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	180/186 (97%)	165 (92%)	12 (7%)	3 (2%)	9 8
1	B	180/186 (97%)	167 (93%)	10 (6%)	3 (2%)	9 8
All	All	360/372 (97%)	332 (92%)	22 (6%)	6 (2%)	9 8

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	85	GLY
1	B	62	GLU
1	A	62	GLU
1	B	79	LEU
1	B	163	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	165/168 (98%)	121 (73%)	44 (27%)	0 0
1	B	164/168 (98%)	133 (81%)	31 (19%)	1 1
All	All	329/336 (98%)	254 (77%)	75 (23%)	1 0

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	18	LYS
1	A	28	ARG
1	A	31	PHE
1	A	32	ARG
1	A	34	PHE
1	A	39	THR
1	A	41	SER
1	A	42	SER
1	A	44	GLU
1	A	47	GLN
1	A	49	LEU
1	A	54	LYS
1	A	55	LYS
1	A	63	LYS
1	A	64	ASN
1	A	65	ARG
1	A	68	LYS
1	A	71	ILE
1	A	76	SER
1	A	80	LYS
1	A	84	GLN
1	A	90	SER
1	A	98	LYS
1	A	105	LEU
1	A	107	ASN
1	A	118	SER

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Mol	Chain	Res	Type
1	A	122	LYS
1	A	123	GLU
1	A	126	ASN
1	A	131	LEU
1	A	138	ILE
1	A	141	ASP
1	A	150	GLU
1	A	153	LEU
1	A	154	GLU
1	A	157	LYS
1	A	158	LEU
1	A	161	GLU
1	A	168	ASP
1	A	176	LYS
1	A	178	LYS
1	A	184	LYS
1	A	185	ASN
1	B	12	GLN
1	B	18	LYS
1	B	30	GLU
1	B	32	ARG
1	B	63	LYS
1	B	70	ARG
1	B	77	ARG
1	B	80	LYS
1	B	84	GLN
1	B	89	LEU
1	B	95	ASP
1	B	101	GLU
1	B	103	PRO
1	B	105	LEU
1	B	107	ASN
1	B	108	LYS
1	B	119	SER
1	B	122	LYS
1	B	131	LEU
1	B	133	LEU
1	B	138	ILE
1	B	146	THR
1	B	150	GLU
1	B	153	LEU
1	B	155	LYS

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Mol	Chain	Res	Type
1	B	157	LYS
1	B	158	LEU
1	B	159	LEU
1	B	161	GLU
1	B	178	LYS
1	B	184	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	126	ASN
1	B	12	GLN
1	B	107	ASN
1	B	130	HIS

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 monosaccharides 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	$\#\|Z\| > 2$
2	DZF	A	187	-	33,34,34	2.11	11 (33%)	45,47,47	2.85	19 (42%)
2	DZF	B	187	-	33,34,34	2.17	10 (30%)	45,47,47	3.37	23 (51%)

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	DZF	A	187	-	-	4/22/22/22	0/3/3/3
2	DZF	B	187	-	-	1/22/22/22	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	187	DZF	C4A-C8A	-6.80	1.35	1.42
2	A	187	DZF	C4A-C8A	-5.72	1.36	1.42
2	B	187	DZF	C2-N3	4.38	1.43	1.35
2	A	187	DZF	CA-N	4.03	1.54	1.45
2	A	187	DZF	C8A-N8	-3.94	1.31	1.37
2	B	187	DZF	CB-CG	3.51	1.63	1.52
2	B	187	DZF	CG-CD	3.45	1.58	1.50
2	A	187	DZF	CG-CD	3.37	1.58	1.50
2	A	187	DZF	C4-N3	3.06	1.38	1.33
2	A	187	DZF	CB-CG	2.98	1.61	1.52
2	A	187	DZF	C2-NA2	-2.83	1.28	1.33
2	B	187	DZF	C2-NA2	-2.76	1.28	1.33
2	A	187	DZF	C2-N3	2.58	1.39	1.35
2	B	187	DZF	C15-C14	2.48	1.43	1.39
2	A	187	DZF	C13-C14	-2.46	1.35	1.39
2	A	187	DZF	OE2-CD	-2.43	1.22	1.30
2	B	187	DZF	C13-C14	-2.32	1.35	1.39
2	B	187	DZF	C-N	-2.32	1.28	1.34
2	B	187	DZF	C5-C4A	-2.24	1.36	1.41
2	B	187	DZF	C7-C6	2.22	1.43	1.38
2	A	187	DZF	C5-C4A	-2.03	1.37	1.41

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	187	DZF	C4A-C4-N3	-8.71	115.77	124.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	187	DZF	C4A-C4-N3	-8.43	116.06	124.68
2	B	187	DZF	N1-C2-N3	-7.54	117.63	127.21
2	B	187	DZF	C5-C4A-C8A	7.03	123.09	117.02
2	A	187	DZF	C5-C4A-C8A	6.72	122.82	117.02
2	B	187	DZF	C2-N1-C8A	5.69	121.62	115.48
2	A	187	DZF	OE1-CD-CG	-5.67	105.09	123.09
2	B	187	DZF	C2-N3-C4	5.60	123.75	115.96
2	A	187	DZF	N1-C2-N3	-5.55	120.16	127.21
2	B	187	DZF	NA2-C2-N1	5.39	126.20	117.79
2	B	187	DZF	O-C-N	-5.22	112.55	122.47
2	A	187	DZF	C2-N3-C4	4.99	122.89	115.96
2	B	187	DZF	C5-C4A-C4	-4.40	117.67	123.66
2	B	187	DZF	C7-N8-C8A	-4.31	111.19	117.20
2	A	187	DZF	C6-C5-C4A	-4.31	115.43	122.59
2	B	187	DZF	CB-CA-N	4.29	119.40	110.91
2	B	187	DZF	OE1-CD-CG	-4.16	109.89	123.09
2	B	187	DZF	C11-C-N	4.09	124.62	117.04
2	B	187	DZF	C6-C5-C4A	-3.88	116.13	122.59
2	A	187	DZF	CB-CA-N	-3.81	103.36	110.91
2	A	187	DZF	NA2-C2-N1	3.72	123.59	117.79
2	A	187	DZF	C2-N1-C8A	3.70	119.48	115.48
2	B	187	DZF	CA-N-C	3.43	129.78	121.56
2	A	187	DZF	OE2-CD-CG	3.35	124.58	114.00
2	B	187	DZF	C13-C12-C11	3.29	124.32	120.80
2	A	187	DZF	C5-C4A-C4	-3.20	119.30	123.66
2	A	187	DZF	CB-CG-CD	-3.19	103.98	112.49
2	B	187	DZF	C6-C7-N8	3.09	128.60	124.01
2	A	187	DZF	C7-N8-C8A	-2.93	113.11	117.20
2	B	187	DZF	C16-C11-C12	-2.85	114.95	118.57
2	A	187	DZF	C12-C13-C14	-2.71	117.18	120.30
2	A	187	DZF	OE2-CD-OE1	2.68	130.23	123.33
2	A	187	DZF	N8-C8A-N1	-2.64	112.88	115.77
2	A	187	DZF	C13-C12-C11	2.48	123.45	120.80
2	B	187	DZF	CB-CG-CD	-2.48	105.87	112.49
2	B	187	DZF	OE2-CD-CG	2.39	121.54	114.00
2	A	187	DZF	C8A-C4A-C4	-2.14	117.82	119.99
2	B	187	DZF	C16-C11-C	2.08	127.34	120.60
2	B	187	DZF	C15-C14-N10	-2.08	116.27	120.96
2	B	187	DZF	C4A-C8A-N1	-2.07	121.60	123.45
2	A	187	DZF	C15-C14-C13	2.04	121.76	119.04
2	B	187	DZF	OE2-CD-OE1	2.02	128.53	123.33

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	187	DZF	N-CA-CB-CG
2	A	187	DZF	CT-CA-CB-CG
2	A	187	DZF	CA-CB-CG-CD
2	A	187	DZF	C6-C9-N10-C14
2	B	187	DZF	OE2-CD-CG-CB

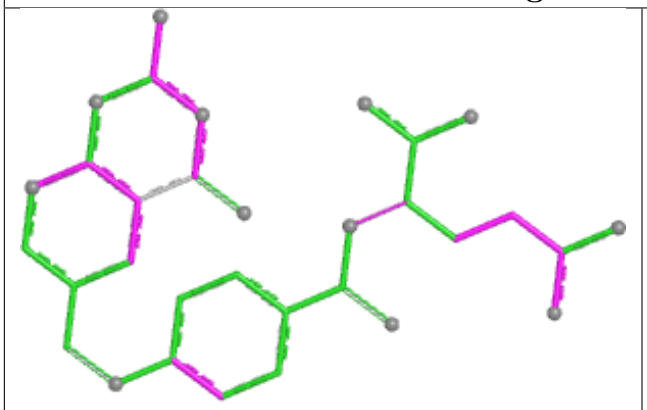
There are no ring outliers.

1 monomer is involved in 4 short contacts:

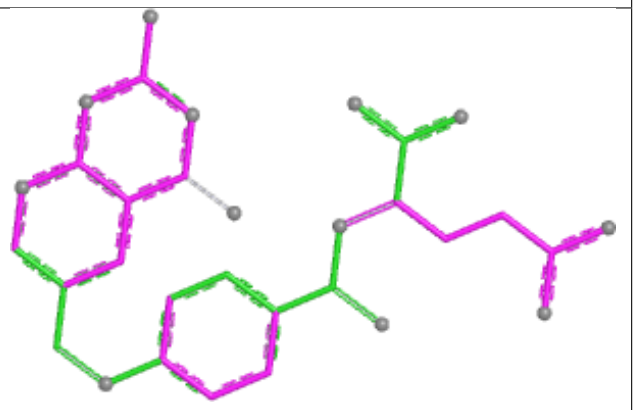
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	187	DZF	4	0

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 then 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.

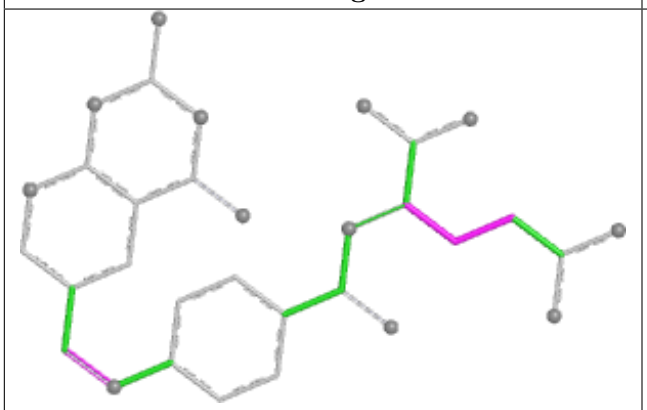
Ligand DZF A 187



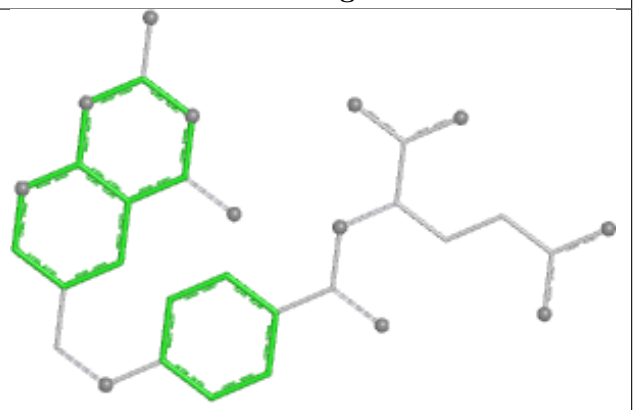
Bond lengths



Bond angles

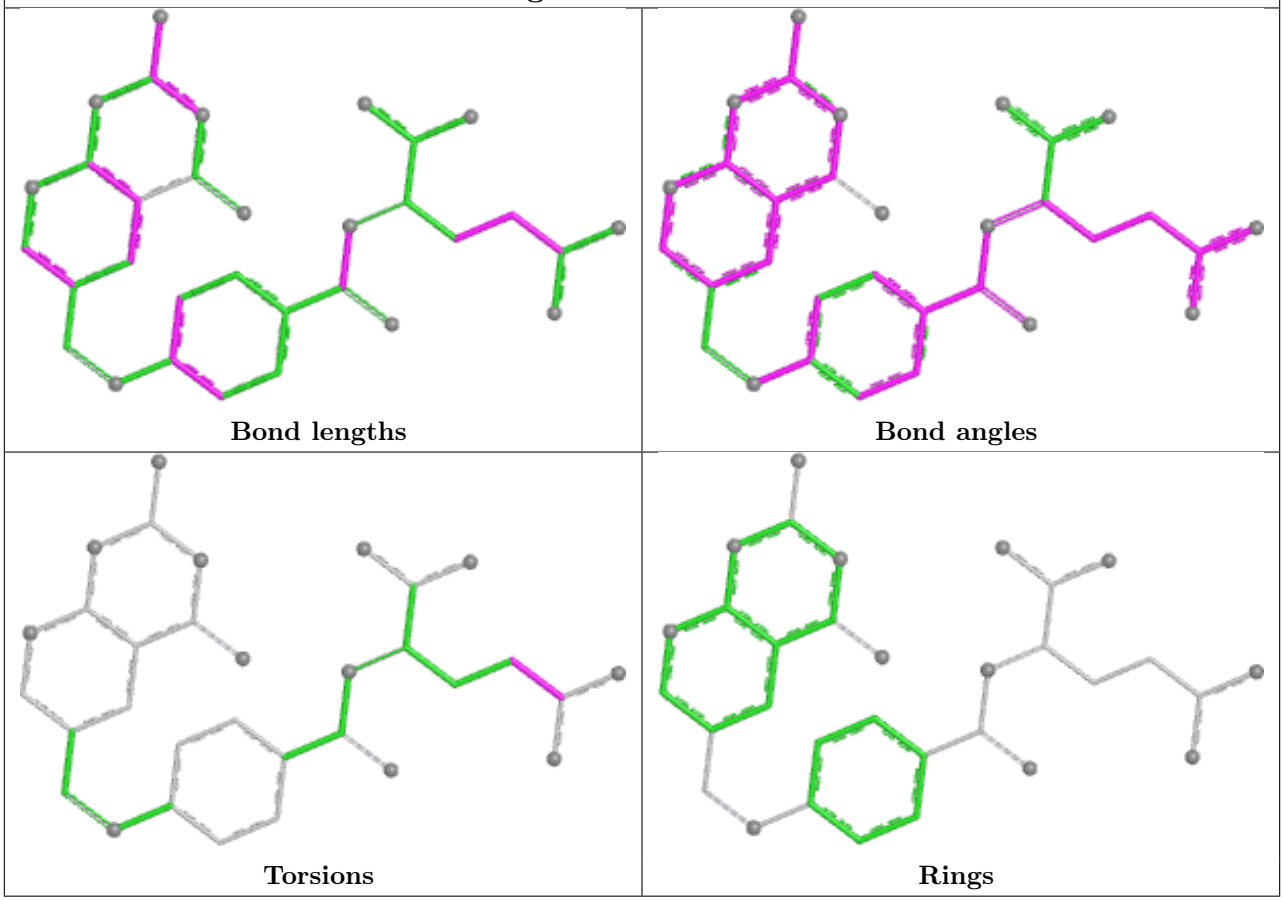


Torsions



Rings

Ligand DZF B 187



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.