



Full wwPDB X-ray Structure Validation Report ⓘ

May 30, 2020 – 02:18 pm BST

PDB ID : 1AWH
Title : NOVEL COVALENT THROMBIN INHIBITOR FROM PLANT EXTRACT
Authors : Jhoti, H.; Cleasby, A.; Wonacott, A.
Deposited on : 1997-10-02
Resolution : 3.00 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
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) : 2.11

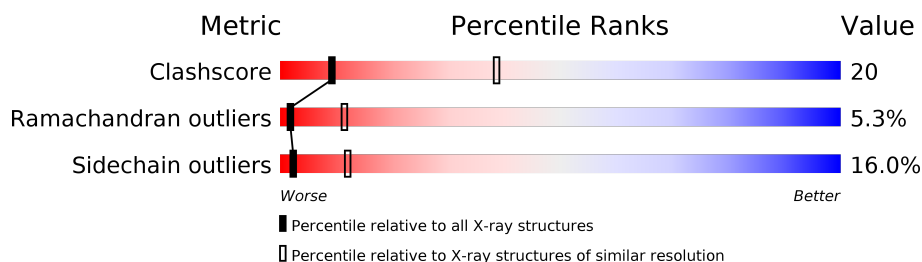
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)

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 on 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	36	<div> <div>28%</div> <div>42%</div> <div>25%</div> <div>6%</div> </div>
1	C	36	<div> <div>33%</div> <div>47%</div> <div>6%</div> <div>14%</div> </div>
2	B	259	<div> <div>49%</div> <div>36%</div> <div>12%</div> <div>•</div> </div>
2	D	259	<div> <div>43%</div> <div>43%</div> <div>12%</div> <div>•</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4840 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.

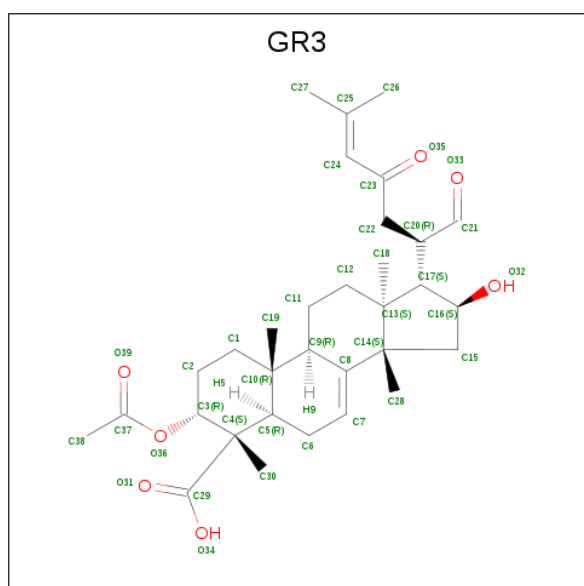
- Molecule 1 is a protein called ALPHA THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	36	Total	C	N	O	S	0	0	0
			287	177	48	61	1			
1	C	36	Total	C	N	O	S	0	0	0
			287	177	48	61	1			

- Molecule 2 is a protein called ALPHA THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	259	Total	C	N	O	S	0	0	0
			2093	1334	370	375	14			
2	D	259	Total	C	N	O	S	0	0	0
			2093	1334	370	375	14			

- Molecule 3 is 3-ACETOXY-17-(1-FORMYL-5-METHYL-3-OXO-HEX-4-ENYL)-16-HYDROXY-4,10,13,14-TETRAMETHYL-2,3,4,5,6,9,10,11,12,13,14,15,16,17-TETRADECAHYDRO-1H-CYCLOPENTA[A]PHENANTHRENE-4-CARBOXYLIC ACID (three-letter code: GR3) (formula: C₃₂H₄₆O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			39	32	7		
3	D	1	Total	C	O	0	0
			39	32	7		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	O	0	0
			1	1		
4	D	1	Total	O	0	0
			1	1		

R77A	E146	G211	R77A	E146	G211
R81	T147	E217	R81	T147	E217
I82	W148		I82	W148	
L85	T149	D221A	L85	T149	D221A
E86	A149A	R221	E86	A149A	R221
R87	M149B	Y225	R87	M149B	Y225
I88	K149E	G226	I88	K149E	G226
	G150	F227		G150	F227
R91	Q151		R91	Q151	
P92	P152	H230	P92	P152	H230
R93	S153	V231	R93	S153	V231
Y94	V157	F232	Y94	V157	F232
N95		R233	N95		R233
R97	P161	L234	R97	P161	L234
L99	I162	K235	L99	I162	K235
D100	V163	K236	D100	V163	K236
R101	E164	V237	R101	E164	V237
D102	R165	I238	D102	R165	I238
I103	R166	Q239	I103	R166	Q239
	V167	K240		V167	K240
	C168			C168	
	K169	D243		K169	D243
M106	R173	Q244	M106	R173	Q244
R107	I174	F245	R107	I174	F245
L108	R175	G246	L108	R175	G246
K109	I176	E247	K109	I176	E247
K110			K110		
P111	M180		P111	M180	
V112	F181		V112	F181	
A113	C182		A113	C182	
F114			F114		
S115	K185		S115	K185	
D116	P186		D116	P186	
Y117	D186A		Y117	D186A	
I118	E186B		I118	E186B	
H119			H119		
P120	R187		P120	R187	
V121			V121		
	A190			A190	
R126	C191		R126	C191	
E127			E127		
T128	S195		T128	S195	
A129			A129		
A129A	P198		A129A	P198	
S129B	F199		S129B	F199	
L129C	V200		L129C	V200	
L130	W201		L130	W201	
Q131	K202		Q131	K202	
Y134	F204A		Y134	F204A	
K135	N204B		K135	N204B	
G136	N205		G136	N205	
R137	R206		R137	R206	
V138	W207		V138	W207	
T139	Y208		T139	Y208	
G140	Q209		G140	Q209	
W141	M210		W141	M210	

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.60 Å 102.80 Å 119.70 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.202 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4840	wwPDB-VP
Average B, all atoms (Å ²)	25.0	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: GR3

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	1.05	0/290	1.89	3/384 (0.8%)
1	C	1.10	0/290	1.83	6/384 (1.6%)
2	B	1.01	5/2148 (0.2%)	1.77	26/2903 (0.9%)
2	D	1.01	5/2148 (0.2%)	1.79	32/2903 (1.1%)
All	All	1.02	10/4876 (0.2%)	1.79	67/6574 (1.0%)

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	7
1	C	0	4
2	B	0	9
2	D	0	9
All	All	0	29

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	195	SER	CB-OG	-9.10	1.30	1.42
2	D	195	SER	CB-OG	-8.60	1.31	1.42
2	B	148	TRP	CD2-CE2	6.58	1.49	1.41
2	D	96	TRP	CD2-CE2	6.03	1.48	1.41
2	B	237	TRP	CD2-CE2	5.95	1.48	1.41
2	D	29	TRP	CD2-CE2	5.90	1.48	1.41
2	B	29	TRP	CD2-CE2	5.67	1.48	1.41
2	D	148	TRP	CD2-CE2	5.25	1.47	1.41
2	B	51	TRP	CD2-CE2	5.07	1.47	1.41
2	D	60(D)	TRP	CD2-CE2	5.04	1.47	1.41

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	134	TYR	CB-CG-CD2	-13.41	112.96	121.00
2	B	134	TYR	CB-CG-CD1	12.84	128.70	121.00
1	A	15	ARG	NE-CZ-NH1	12.81	126.70	120.30
2	D	67	ARG	NE-CZ-NH1	8.03	124.31	120.30
2	B	221	ARG	NE-CZ-NH1	7.66	124.13	120.30
2	D	165	ARG	CD-NE-CZ	-7.64	112.91	123.60
2	D	126	ARG	NE-CZ-NH1	7.46	124.03	120.30
2	D	50	ARG	NE-CZ-NH2	-7.39	116.61	120.30
2	B	75	ARG	CA-CB-CG	7.24	129.32	113.40
2	B	220	CYS	CA-CB-SG	-7.23	100.99	114.00
2	D	175	ARG	NE-CZ-NH2	-7.05	116.77	120.30
2	D	151	GLN	CB-CA-C	-7.05	96.30	110.40
2	D	165	ARG	NE-CZ-NH2	-6.96	116.82	120.30
2	D	131	GLN	CA-CB-CG	-6.81	98.42	113.40
2	B	67	ARG	CD-NE-CZ	-6.65	114.28	123.60
2	B	97	ARG	NE-CZ-NH1	6.63	123.61	120.30
2	B	233	ARG	NE-CZ-NH1	6.58	123.59	120.30
2	D	134	TYR	CB-CG-CD2	-6.46	117.12	121.00
1	A	14(G)	LEU	CA-CB-CG	6.36	129.92	115.30
2	B	144	LEU	CA-CB-CG	-6.20	101.04	115.30
2	B	101	ARG	NE-CZ-NH2	-6.15	117.22	120.30
2	D	75	ARG	N-CA-CB	-6.13	99.56	110.60
1	C	14(G)	LEU	N-CA-CB	-6.07	98.26	110.40
1	C	14(K)	ILE	N-CA-CB	6.05	124.71	110.80
2	B	244	GLN	C-N-CA	-5.98	106.74	121.70
2	D	173	ARG	NE-CZ-NH1	5.95	123.27	120.30
2	D	33	LEU	CB-CA-C	-5.88	99.04	110.20
2	D	175	ARG	CB-CA-C	-5.81	98.78	110.40
1	C	14(G)	LEU	CB-CA-C	5.73	121.09	110.20
2	D	148	TRP	CB-CA-C	-5.73	98.94	110.40
2	D	35	ARG	NE-CZ-NH2	-5.70	117.45	120.30
2	D	165	ARG	NE-CZ-NH1	5.70	123.15	120.30
2	D	182	CYS	CA-CB-SG	5.69	124.25	114.00
2	B	65	LEU	CA-CB-CG	5.69	128.39	115.30
2	D	149(A)	ALA	N-CA-C	5.67	126.31	111.00
2	B	65	LEU	N-CA-CB	-5.54	99.32	110.40
2	D	77(A)	ARG	CA-CB-CG	5.51	125.53	113.40
1	A	14(M)	GLY	C-N-CA	5.48	135.39	121.70
2	B	205	ASN	CA-CB-CG	-5.46	101.40	113.40
2	D	221(A)	ASP	C-N-CA	-5.45	108.07	121.70
2	B	210	MET	CA-CB-CG	5.42	122.52	113.30
2	B	182	CYS	CA-CB-SG	5.42	123.75	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	14(K)	ILE	CB-CA-C	-5.32	100.96	111.60
2	D	76	TYR	C-N-CA	-5.32	108.40	121.70
2	D	180	MET	N-CA-CB	-5.31	101.04	110.60
2	D	115	SER	C-N-CA	-5.31	108.43	121.70
2	B	68	ILE	C-N-CA	-5.30	111.17	122.30
2	B	75	ARG	NE-CZ-NH1	5.28	122.94	120.30
2	B	149(A)	ALA	C-N-CA	-5.27	108.52	121.70
2	D	126	ARG	CB-CG-CD	5.25	125.26	111.60
2	B	94	TYR	CB-CG-CD2	-5.25	117.85	121.00
2	D	137	ARG	CD-NE-CZ	-5.25	116.26	123.60
2	D	116	ASP	CB-CA-C	5.23	120.86	110.40
1	C	14(H)	GLU	CA-CB-CG	5.22	124.88	113.40
2	D	175	ARG	NE-CZ-NH1	5.21	122.90	120.30
2	B	149(A)	ALA	CB-CA-C	5.16	117.84	110.10
2	D	112	VAL	CG1-CB-CG2	5.15	119.14	110.90
2	D	153	SER	CB-CA-C	-5.13	100.36	110.10
2	D	38	GLN	CA-C-N	-5.12	105.93	117.20
2	B	228	TYR	CB-CG-CD1	-5.10	117.94	121.00
2	B	52	VAL	CG1-CB-CG2	-5.10	102.74	110.90
2	D	77(A)	ARG	C-N-CA	-5.09	108.97	121.70
2	B	72	SER	CB-CA-C	-5.08	100.45	110.10
1	C	1(C)	GLU	CA-CB-CG	5.05	124.50	113.40
2	B	149(E)	LYS	C-N-CA	-5.04	111.71	122.30
2	B	233	ARG	NE-CZ-NH2	-5.03	117.78	120.30
2	D	71	HIS	CA-CB-CG	5.01	122.11	113.60

There are no chirality outliers.

All (29) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1(D)	GLY	Peptide
1	A	1(G)	PHE	Peptide
1	A	1(H)	THR	Peptide
1	A	14(J)	TYR	Sidechain,Peptide
1	A	14(L)	ASP	Peptide
1	A	15	ARG	Sidechain
2	B	137	ARG	Sidechain
2	B	142	GLY	Peptide
2	B	148	TRP	Peptide
2	B	149(A)	ALA	Peptide
2	B	187	ARG	Sidechain
2	B	221	ARG	Sidechain

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Mol	Chain	Res	Type	Group
2	B	246	GLY	Peptide
2	B	67	ARG	Sidechain
2	B	77	GLU	Peptide
1	C	1(F)	GLY	Peptide
1	C	1(G)	PHE	Peptide
1	C	14(J)	TYR	Sidechain
1	C	14(M)	GLY	Peptide
2	D	134	TYR	Sidechain
2	D	146	GLU	Peptide
2	D	147	THR	Peptide
2	D	165	ARG	Sidechain
2	D	173	ARG	Sidechain
2	D	186(A)	ASP	Peptide
2	D	246	GLY	Peptide
2	D	81	LYS	Peptide
2	D	93	ARG	Sidechain

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	287	0	278	22	0
1	C	287	0	278	19	0
2	B	2093	0	2063	80	0
2	D	2093	0	2063	86	0
3	B	39	0	44	5	0
3	D	39	0	44	4	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
All	All	4840	0	4770	191	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 20.

All (191) 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)
2:B:60(F):LYS:HG3	2:B:60(H):PHE:HE1	1.33	0.94
1:A:14(G):LEU:HB3	1:A:15:ARG:NE	1.90	0.87
2:D:230:HIS:CD2	2:D:233:ARG:HG2	2.16	0.80
2:D:200:VAL:HG12	2:D:209:GLN:HA	1.62	0.80
2:D:190:ALA:HB3	3:D:1:GR3:H262	1.64	0.80
2:D:67:ARG:HG2	2:D:82:ILE:HG12	1.64	0.79
2:B:185:LYS:HB2	2:B:186(B):GLU:HG3	1.65	0.77
2:B:60(F):LYS:HG3	2:B:60(H):PHE:CE1	2.20	0.74
1:C:14(G):LEU:HB3	1:C:15:ARG:CZ	2.18	0.73
2:B:144:LEU:HD11	2:B:152:PRO:HB3	1.72	0.71
2:D:163:VAL:HG21	2:D:225:TYR:CD2	2.28	0.68
2:B:130:LEU:HD23	2:B:162:ILE:HD13	1.75	0.68
2:D:49:ASP:HB3	2:D:114:PHE:CZ	2.28	0.68
1:C:14(G):LEU:O	1:C:15:ARG:HD3	1.95	0.67
2:D:163:VAL:HG21	2:D:225:TYR:CE2	2.31	0.65
2:B:35:ARG:HH21	2:B:37(A):PRO:HD2	1.62	0.65
1:C:14(G):LEU:HB3	1:C:15:ARG:NH1	2.10	0.65
1:A:1(G):PHE:CD1	1:A:1(E):SER:HB2	2.34	0.63
1:A:3:LEU:HD11	2:B:206:ARG:NH1	2.14	0.63
2:B:154:VAL:O	2:B:156:GLN:HG3	1.98	0.62
2:B:91:HIS:HB2	2:B:103:ILE:HG22	1.80	0.62
2:B:144:LEU:HD13	2:B:149(E):LYS:HB3	1.82	0.61
2:B:60(A):TYR:HD2	2:B:60(D):TRP:HB2	1.67	0.60
2:B:51:TRP:CE2	2:B:242:ILE:HG12	2.37	0.60
2:B:60(A):TYR:CD2	2:B:60(D):TRP:HB2	2.36	0.60
2:B:81:LYS:HD3	2:B:112:VAL:HG23	1.83	0.59
2:D:128:THR:HG23	2:D:129(C):LEU:HD12	1.85	0.58
2:D:33:LEU:O	2:D:40:LEU:HD23	2.03	0.58
1:A:14(G):LEU:HD21	2:B:202:LYS:HD3	1.84	0.58
1:C:14(F):LEU:HD21	2:D:207:TRP:HH2	1.68	0.57
2:D:68:ILE:HG22	2:D:118:ILE:HG12	1.87	0.57
2:B:57:HIS:O	2:B:60(F):LYS:HE2	2.04	0.57
1:C:4:ARG:HD3	2:D:26:MET:O	2.05	0.57
2:D:17:VAL:HG12	2:D:18:GLU:HG2	1.87	0.57
2:B:67:ARG:HG2	2:B:82:ILE:HG12	1.86	0.57
2:B:147:THR:HA	2:B:149:THR:O	2.04	0.57
1:C:1(H):THR:HA	2:D:243:ASP:HA	1.87	0.56
2:D:165:ARG:HH22	2:D:176:ILE:HG22	1.69	0.56
2:B:42:CYS:HB3	2:B:195:SER:O	2.06	0.56
2:D:91:HIS:ND1	2:D:101:ARG:HD3	2.20	0.56
1:C:14:ASP:HB2	2:D:23:GLU:OE1	2.05	0.55
3:D:1:GR3:H192	3:D:1:GR3:H303	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:211:GLY:HA2	2:B:231:VAL:HG23	1.89	0.55
2:D:180:MET:HG2	2:D:227:PHE:HD2	1.72	0.54
2:B:190:ALA:HB3	3:B:1:GR3:H262	1.90	0.54
2:B:61:GLU:CD	2:B:61:GLU:H	2.11	0.53
1:C:1(H):THR:N	2:D:243:ASP:HA	2.23	0.53
2:D:91:HIS:CE1	2:D:101:ARG:HD3	2.44	0.53
2:B:94:TYR:CZ	2:B:96:TRP:HB3	2.44	0.53
2:D:182:CYS:HB2	2:D:226:GLY:O	2.09	0.53
2:D:32:MET:HE2	2:D:34:PHE:HE1	1.74	0.53
2:B:73:ARG:CZ	2:B:151:GLN:HG2	2.40	0.52
1:C:1(H):THR:HA	2:D:243:ASP:OD1	2.10	0.52
2:D:22:ALA:O	2:D:71:HIS:HE1	1.94	0.51
2:B:35:ARG:HE	2:B:37(A):PRO:HD2	1.75	0.51
2:B:70:LYS:HB2	2:B:70:LYS:NZ	2.26	0.51
2:B:21:ASP:HA	2:B:156:GLN:HG2	1.92	0.51
2:B:144:LEU:CD1	2:B:149(E):LYS:HB3	2.41	0.51
3:B:1:GR3:H303	3:B:1:GR3:H192	1.92	0.51
2:D:204(B):ASN:ND2	2:D:206:ARG:HD2	2.26	0.51
2:D:165:ARG:HG2	2:D:169:LYS:HE2	1.93	0.51
1:A:14(F):LEU:HB3	1:A:14(G):LEU:HD22	1.94	0.50
1:A:14(G):LEU:HB3	1:A:15:ARG:CZ	2.41	0.50
2:B:176:ILE:HD11	2:B:227:PHE:CE2	2.46	0.50
2:D:60(B):PRO:HB2	2:D:60(C):PRO:HD3	1.93	0.50
2:B:213:VAL:HG22	2:B:228:TYR:CE2	2.47	0.50
2:D:135:LYS:HA	2:D:161:PRO:HA	1.93	0.50
1:C:5:PRO:HA	1:C:9:LYS:HB2	1.94	0.49
1:A:1(G):PHE:HD1	1:A:1(E):SER:HB2	1.76	0.49
2:D:185:LYS:HB2	2:D:186(B):GLU:HG3	1.93	0.49
2:B:77:GLU:HB2	2:B:80:GLU:HG2	1.95	0.49
2:D:73:ARG:HD3	2:D:152:PRO:O	2.12	0.48
2:D:126:ARG:HA	2:D:232:PHE:CZ	2.48	0.48
1:A:3:LEU:HD11	2:B:206:ARG:HH12	1.77	0.48
2:D:50:ARG:HD3	2:D:111:PRO:HD3	1.95	0.48
1:A:15:ARG:HB2	1:A:15:ARG:HH11	1.78	0.48
2:B:70:LYS:HZ3	2:B:70:LYS:HB2	1.79	0.47
2:D:96:TRP:CH2	2:D:97:ARG:HG3	2.49	0.47
2:B:201:MET:SD	2:B:210:MET:HG3	2.55	0.47
1:A:1(H):THR:H3	2:B:246:GLY:HA3	1.79	0.47
2:B:17:VAL:O	2:B:188:GLY:HA2	2.13	0.47
2:B:41:LEU:O	2:B:42:CYS:SG	2.72	0.47
2:D:20:SER:O	2:D:157:VAL:HG12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:73:ARG:HG3	2:D:141:TRP:CD1	2.50	0.47
2:D:32:MET:HE2	2:D:34:PHE:CE1	2.48	0.47
1:C:14(K):ILE:HG13	2:D:134:TYR:CE1	2.49	0.47
2:D:165:ARG:N	2:D:166:PRO:HD2	2.30	0.47
2:B:145:LYS:HE3	2:B:149(C):VAL:N	2.31	0.46
2:B:32:MET:HG3	2:B:40:LEU:HD23	1.97	0.46
2:D:202:LYS:HE2	2:D:205:ASN:HB2	1.98	0.46
2:D:168:CYS:HB3	2:D:227:PHE:CZ	2.50	0.46
1:A:3:LEU:O	1:A:5:PRO:HD3	2.15	0.46
2:D:211:GLY:HA2	2:D:231:VAL:HG23	1.98	0.46
1:A:14(G):LEU:HD12	1:A:15:ARG:HH21	1.81	0.46
1:A:7:PHE:CE2	2:B:26:MET:HA	2.51	0.46
2:D:199:PHE:O	2:D:210:MET:N	2.49	0.46
2:D:60(B):PRO:HG2	2:D:96:TRP:CE2	2.51	0.46
2:D:94:TYR:CZ	2:D:96:TRP:HB3	2.51	0.46
1:C:1(G):PHE:CE1	1:C:1(C):GLU:HB3	2.51	0.46
2:D:85:LEU:HD13	2:D:88:ILE:HD11	1.97	0.46
2:B:60(A):TYR:CZ	2:B:60(C):PRO:HG2	2.50	0.45
2:D:243:ASP:O	2:D:245:PHE:N	2.49	0.45
2:D:91:HIS:HD2	2:D:237:TRP:CD2	2.34	0.45
1:A:5:PRO:HG2	2:B:115:SER:O	2.15	0.45
2:B:47:ILE:HD12	2:B:123:LEU:HD11	1.98	0.45
2:D:191:CYS:HA	3:D:1:GR3:H273	1.98	0.45
1:C:14(G):LEU:C	1:C:15:ARG:HD3	2.37	0.45
2:D:95:ASN:HB3	2:D:100:ASP:HB3	1.98	0.45
2:B:151:GLN:CD	2:B:151:GLN:H	2.19	0.45
2:B:163:VAL:HG12	2:B:168:CYS:SG	2.56	0.45
2:D:136:GLY:HA3	2:D:199:PHE:CZ	2.51	0.45
2:B:185:LYS:HD3	2:B:186(B):GLU:OE2	2.17	0.45
1:A:1(H):THR:N	2:B:246:GLY:HA3	2.32	0.45
2:D:146:GLU:O	2:D:149(B):ASN:OD1	2.35	0.45
2:B:105:LEU:HD13	2:B:241:VAL:HG22	1.99	0.44
2:B:135:LYS:HA	2:B:161:PRO:HA	1.98	0.44
1:A:14(G):LEU:HB3	1:A:15:ARG:CD	2.46	0.44
3:B:1:GR3:C12	3:B:1:GR3:H221	2.48	0.44
2:D:42:CYS:SG	2:D:195:SER:HB2	2.57	0.44
2:B:213:VAL:HG22	2:B:228:TYR:HE2	1.83	0.44
2:D:60(C):PRO:HD3	2:D:96:TRP:CZ3	2.52	0.44
2:D:91:HIS:HA	2:D:92:PRO:HD2	1.85	0.44
2:D:18:GLU:HG3	2:D:187:ARG:HB2	2.00	0.43
2:B:177:THR:O	2:B:180:MET:HG2	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:29:TRP:CG	2:B:121:VAL:HB	2.52	0.43
1:C:6:LEU:HD21	2:D:116:ASP:O	2.18	0.43
2:D:50:ARG:NH1	2:D:111:PRO:HG3	2.32	0.43
1:A:1(G):PHE:HD1	1:A:1(E):SER:CB	2.32	0.43
2:B:70:LYS:HB2	2:B:77:GLU:OE1	2.17	0.43
2:D:149(E):LYS:HE2	2:D:151:GLN:O	2.17	0.43
2:B:160:LEU:HD22	2:B:184(A):GLY:HA2	2.01	0.43
2:D:101:ARG:O	2:D:103:ILE:N	2.51	0.43
2:D:110:LYS:HA	2:D:111:PRO:HD2	1.75	0.43
2:D:60(C):PRO:HD3	2:D:96:TRP:CE3	2.53	0.43
2:B:91:HIS:HB2	2:B:103:ILE:CG2	2.48	0.43
2:D:162:ILE:HD13	2:D:162:ILE:HG21	1.78	0.43
2:B:39:GLU:OE2	2:D:169:LYS:HE3	2.18	0.43
2:D:180:MET:HB2	2:D:180:MET:HE3	1.91	0.43
2:B:148:TRP:CZ3	2:B:219:GLY:HA2	2.54	0.43
2:D:138:VAL:CG2	2:D:199:PHE:HD1	2.32	0.43
2:B:25:GLY:O	2:B:28:PRO:HD3	2.19	0.42
2:B:64:LEU:HD23	2:B:85:LEU:CD1	2.49	0.42
1:C:14(K):ILE:HG13	2:D:134:TYR:HE1	1.84	0.42
2:B:72:SER:HA	2:B:154:VAL:HG23	2.02	0.42
2:D:60(F):LYS:HG2	2:D:60(H):PHE:CE2	2.54	0.42
1:A:14:ASP:OD1	1:A:14(C):GLU:HB2	2.20	0.42
2:B:31:VAL:HG13	2:B:66:VAL:HG13	2.02	0.42
2:B:50:ARG:HG2	2:B:51:TRP:CD1	2.54	0.42
1:C:14(J):TYR:CD2	1:C:15:ARG:HD2	2.53	0.42
2:D:129(B):SER:HA	2:D:131:GLN:NE2	2.34	0.42
1:A:1(G):PHE:HE1	1:A:1(C):GLU:HB3	1.84	0.42
2:B:101:ARG:O	2:B:103:ILE:N	2.52	0.42
2:B:191:CYS:HA	3:B:1:GR3:H273	2.02	0.42
2:B:64:LEU:HD23	2:B:85:LEU:HD12	2.02	0.42
2:D:237:TRP:O	2:D:238:ILE:C	2.57	0.42
2:B:144:LEU:HA	2:B:144:LEU:HD23	1.79	0.42
2:B:202:LYS:NZ	2:B:205:ASN:HD21	2.18	0.42
2:B:33:LEU:HD13	2:B:41:LEU:HD12	2.01	0.42
1:A:1(G):PHE:CE1	1:A:1(E):SER:HB2	2.55	0.41
2:B:131:GLN:O	2:B:132:ALA:C	2.59	0.41
2:B:37:SER:HA	2:B:37(A):PRO:HA	1.75	0.41
2:B:37(A):PRO:HG3	2:D:165:ARG:HD2	2.02	0.41
2:D:232:PHE:HA	2:D:235:LYS:HB2	2.03	0.41
2:B:125:ASP:O	2:B:126:ARG:C	2.58	0.41
2:D:88:ILE:HD13	2:D:106:MET:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:182:CYS:SG	2:D:225:TYR:HB3	2.60	0.41
3:B:1:GR3:H111	3:B:1:GR3:H283	2.03	0.41
2:D:198:PRO:HB2	2:D:200:VAL:HG13	2.03	0.41
1:C:1(H):THR:CA	2:D:243:ASP:HA	2.51	0.41
2:D:41:LEU:HA	2:D:41:LEU:HD23	1.88	0.41
2:D:88:ILE:CD1	2:D:106:MET:HE2	2.51	0.41
2:B:31:VAL:HG11	2:B:52:VAL:HG11	2.02	0.41
1:C:3:LEU:HD23	1:C:8:GLU:HG2	2.02	0.41
2:D:31:VAL:O	2:D:44:ALA:N	2.53	0.41
2:B:147:THR:C	2:B:148:TRP:HD1	2.23	0.41
2:D:70:LYS:HE3	2:D:72:SER:O	2.21	0.41
2:B:230:HIS:ND1	2:B:233:ARG:HB2	2.35	0.41
2:B:46:LEU:HD22	2:B:48:SER:O	2.21	0.41
1:A:7:PHE:HB3	1:A:12:LEU:O	2.21	0.40
2:B:68:ILE:HD13	2:B:112:VAL:HG21	2.03	0.40
1:A:14(D):ARG:O	1:A:14(H):GLU:HB2	2.22	0.40
2:B:145:LYS:HB2	2:B:145:LYS:HE2	1.66	0.40
2:B:191:CYS:SG	2:B:192:GLU:N	2.94	0.40
2:D:127:GLU:O	2:D:129(A):ALA:HB3	2.22	0.40
2:B:60(A):TYR:CE2	2:B:60(C):PRO:HG2	2.57	0.40
1:C:14(F):LEU:HD21	2:D:207:TRP:CH2	2.52	0.40
2:D:121:VAL:HG11	2:D:209:GLN:HB2	2.04	0.40
2:D:230:HIS:O	2:D:231:VAL:C	2.59	0.40
2:D:57:HIS:NE2	3:D:1:GR3:H281	2.37	0.40
2:D:119:HIS:HA	2:D:120:PRO:HD3	1.98	0.40
2:D:174:ILE:O	2:D:176:ILE:N	2.55	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	34/36 (94%)	16 (47%)	10 (29%)	8 (24%)	0	0
1	C	34/36 (94%)	20 (59%)	9 (26%)	5 (15%)	0	1
2	B	257/259 (99%)	220 (86%)	32 (12%)	5 (2%)	8	36
2	D	257/259 (99%)	210 (82%)	34 (13%)	13 (5%)	2	12
All	All	582/590 (99%)	466 (80%)	85 (15%)	31 (5%)	2	11

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1(B)	ALA
1	A	14(K)	ILE
1	A	14(L)	ASP
2	B	62	ASN
2	B	93	ARG
2	B	102	ASP
2	D	102	ASP
2	D	149(A)	ALA
2	D	149(B)	ASN
2	D	244	GLN
1	A	1(D)	GLY
1	A	1(C)	GLU
1	A	14(C)	GLU
1	C	1(E)	SER
1	C	1(B)	ALA
1	C	14(H)	GLU
1	C	14(K)	ILE
2	D	60(E)	ASP
2	D	61	GLU
2	D	148	TRP
2	D	175	ARG
2	B	61	GLU
1	C	1(G)	PHE
2	D	75	ARG
2	D	231	VAL
2	B	147	THR
2	D	117	TYR
1	A	14(A)	LYS
1	A	8	GLU
2	D	92	PRO
2	D	111	PRO

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	31/31 (100%)	23 (74%)	8 (26%)	0	2
1	C	31/31 (100%)	22 (71%)	9 (29%)	0	2
2	B	225/225 (100%)	192 (85%)	33 (15%)	3	15
2	D	225/225 (100%)	193 (86%)	32 (14%)	3	16
All	All	512/512 (100%)	430 (84%)	82 (16%)	2	12

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1(H)	THR
1	A	6	LEU
1	A	9	LYS
1	A	11	SER
1	A	14(F)	LEU
1	A	14(H)	GLU
1	A	14(K)	ILE
1	A	15	ARG
2	B	16	ILE
2	B	18	GLU
2	B	27	SER
2	B	33	LEU
2	B	40	LEU
2	B	45	SER
2	B	46	LEU
2	B	50	ARG
2	B	60(D)	TRP
2	B	64	LEU
2	B	70	LYS
2	B	72	SER
2	B	80	GLU
2	B	87	LYS
2	B	88	ILE
2	B	93	ARG

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Mol	Chain	Res	Type
2	B	101	ARG
2	B	109	LYS
2	B	115	SER
2	B	128	THR
2	B	135	LYS
2	B	143	ASN
2	B	149(B)	ASN
2	B	149(C)	VAL
2	B	163	VAL
2	B	165	ARG
2	B	175	ARG
2	B	185	LYS
2	B	205	ASN
2	B	241	VAL
2	B	243	ASP
2	B	245	PHE
2	B	247	GLU
1	C	1(H)	THR
1	C	1(G)	PHE
1	C	1(C)	GLU
1	C	1(A)	ASP
1	C	12	LEU
1	C	14(D)	ARG
1	C	14(H)	GLU
1	C	14(I)	SER
1	C	14(J)	TYR
2	D	26	MET
2	D	27	SER
2	D	29	TRP
2	D	33	LEU
2	D	40	LEU
2	D	45	SER
2	D	46	LEU
2	D	48	SER
2	D	53	LEU
2	D	66	VAL
2	D	70	LYS
2	D	75	ARG
2	D	86	GLU
2	D	99	LEU
2	D	108	LEU
2	D	109	LYS

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Mol	Chain	Res	Type
2	D	121	VAL
2	D	127	GLU
2	D	129(B)	SER
2	D	139	THR
2	D	147	THR
2	D	149(B)	ASN
2	D	165	ARG
2	D	186(A)	ASP
2	D	204(A)	PHE
2	D	210	MET
2	D	217	GLU
2	D	221	ARG
2	D	236	LYS
2	D	240	LYS
2	D	244	GLN
2	D	245	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	205	ASN
2	B	244	GLN
2	D	71	HIS
2	D	91	HIS
2	D	131	GLN
2	D	143	ASN
2	D	151	GLN
2	D	159	ASN
2	D	230	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	GR3	D	1	2	40,42,42	1.08	3 (7%)	56,68,68	1.59	8 (14%)
3	GR3	B	1	2	40,42,42	1.08	4 (10%)	56,68,68	1.40	9 (16%)

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	GR3	D	1	2	-	4/18/99/99	0/4/4/4
3	GR3	B	1	2	-	5/18/99/99	0/4/4/4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1	GR3	C4-C29	2.81	1.55	1.50
3	D	1	GR3	C14-C8	-2.79	1.48	1.53
3	B	1	GR3	C4-C3	-2.30	1.51	1.56
3	B	1	GR3	C10-C5	-2.15	1.53	1.56
3	D	1	GR3	C4-C29	2.11	1.54	1.50
3	D	1	GR3	C4-C3	-2.02	1.52	1.56
3	B	1	GR3	C11-C9	-2.00	1.50	1.53

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	GR3	C15-C14-C8	4.80	120.99	117.19
3	D	1	GR3	C28-C14-C8	-4.63	101.65	107.83
3	B	1	GR3	C22-C20-C21	-4.40	98.83	109.51
3	D	1	GR3	C3-O36-C37	4.28	124.42	117.86
3	D	1	GR3	C22-C20-C21	-3.77	100.37	109.51
3	B	1	GR3	C3-O36-C37	3.42	123.10	117.86
3	B	1	GR3	C28-C14-C8	-2.66	104.28	107.83
3	D	1	GR3	C28-C14-C15	-2.57	104.76	112.16
3	D	1	GR3	O36-C3-C4	2.26	111.33	107.26
3	B	1	GR3	C1-C10-C5	2.19	110.71	108.02
3	B	1	GR3	O36-C3-C2	-2.12	104.84	108.48
3	D	1	GR3	O36-C37-O39	2.12	127.17	122.96
3	B	1	GR3	C19-C10-C5	-2.08	109.06	112.92
3	B	1	GR3	C18-C13-C14	2.07	115.10	111.10
3	B	1	GR3	O36-C37-O39	2.06	127.05	122.96
3	B	1	GR3	C18-C13-C17	-2.05	106.56	110.97
3	D	1	GR3	O33-C21-C20	-2.02	120.43	125.16

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1	GR3	C17-C20-C22-C23
3	D	1	GR3	C21-C20-C22-C23
3	B	1	GR3	C17-C20-C22-C23
3	B	1	GR3	C21-C20-C22-C23
3	B	1	GR3	C38-C37-O36-C3
3	B	1	GR3	O39-C37-O36-C3
3	D	1	GR3	O39-C37-O36-C3
3	D	1	GR3	C38-C37-O36-C3
3	B	1	GR3	C4-C3-O36-C37

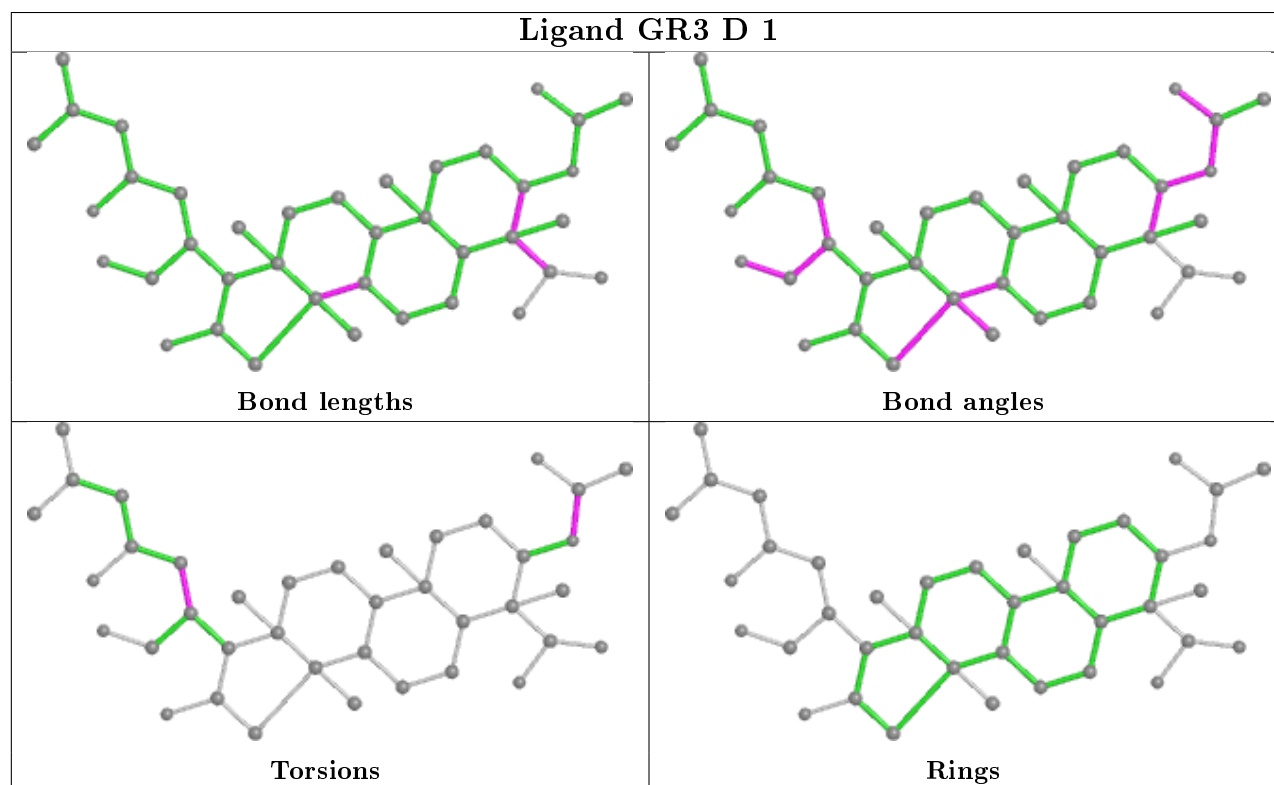
There are no ring outliers.

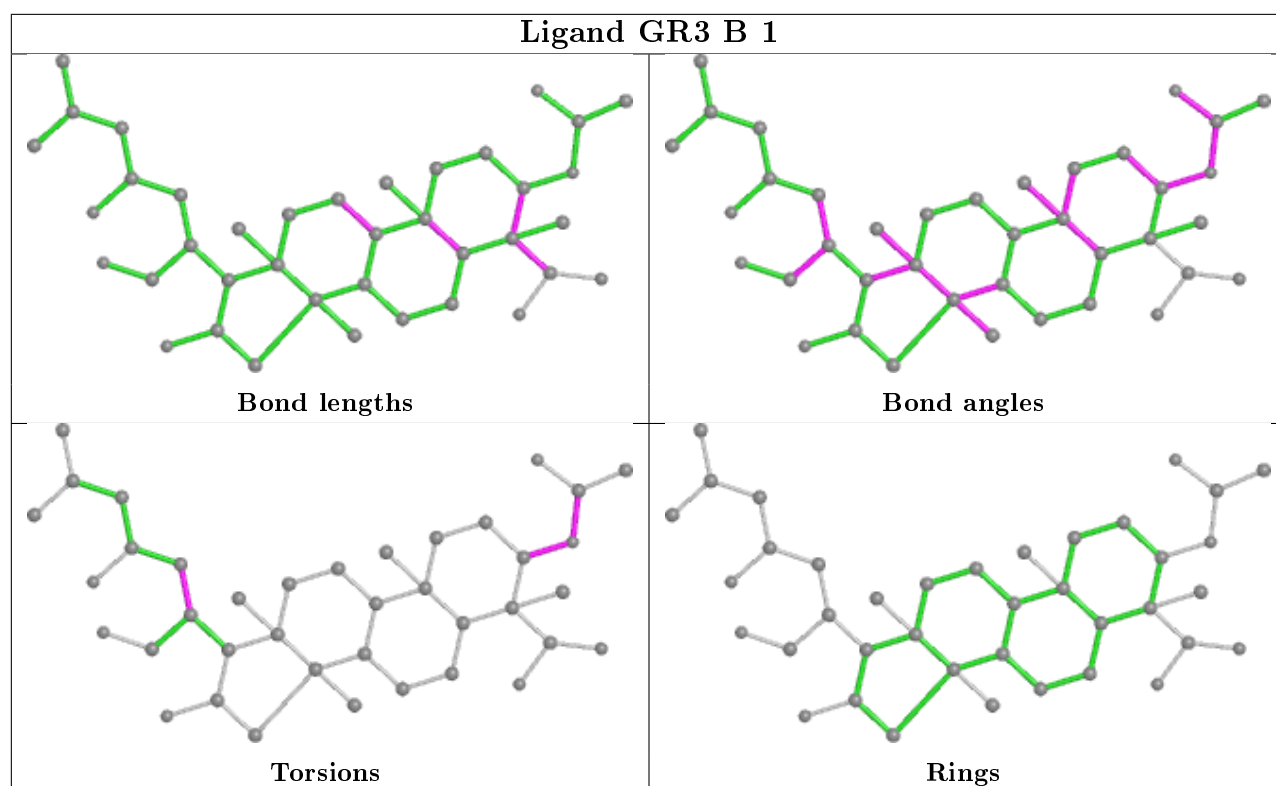
2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1	GR3	4	0
3	B	1	GR3	5	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 ⓘ

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.