



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 14, 2023 – 08:34 PM EDT

PDB ID : 1RPW
Title : Crystal Structure Of The Multidrug Binding Protein Qacr Bound To The Diamidine Hexamidine
Authors : Murray, D.S.; Schumacher, M.A.; Brennan, R.G.
Deposited on : 2003-12-03
Resolution : 2.90 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

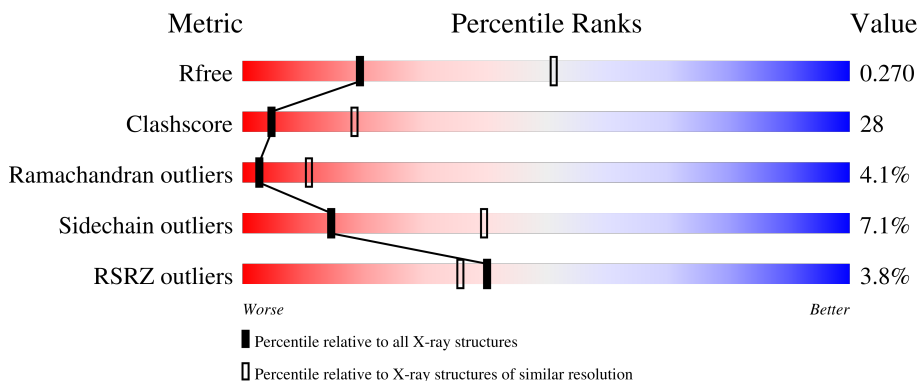
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	188	
1	B	188	
1	C	188	
1	D	188	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	190	-	-	-	X
2	SO4	C	194	-	-	-	X
2	SO4	D	191	-	-	X	-
3	DID	B	193	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6333 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 Transcriptional regulator qacR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	186	1547	998	252	295	2	0	0	0
1	B	186	1547	998	252	295	2	0	0	0
1	C	186	1547	998	252	295	2	0	0	0
1	D	186	1547	998	252	295	2	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	72	ALA	CYS	engineered mutation	UNP P0A0N4
A	141	SER	CYS	engineered mutation	UNP P0A0N4
B	72	ALA	CYS	engineered mutation	UNP P0A0N4
B	141	SER	CYS	engineered mutation	UNP P0A0N4
C	72	ALA	CYS	engineered mutation	UNP P0A0N4
C	141	SER	CYS	engineered mutation	UNP P0A0N4
D	72	ALA	CYS	engineered mutation	UNP P0A0N4
D	141	SER	CYS	engineered mutation	UNP P0A0N4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



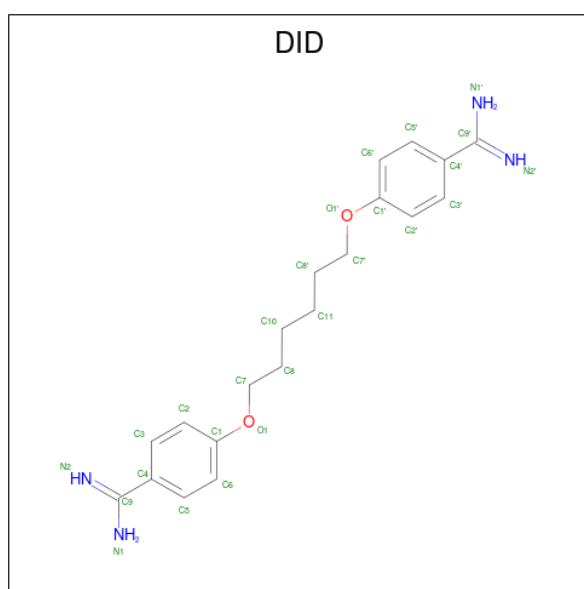
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 4,4'[1,6-HEXANEDIYLBIS(OXY)]BISBENZENECARBOXIMIDAMIDE (three-letter code: DID) (formula: C₂₀H₂₆N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			26	20	4	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	O	0	0
			3	3		
4	B	10	Total	O	0	0
			10	10		

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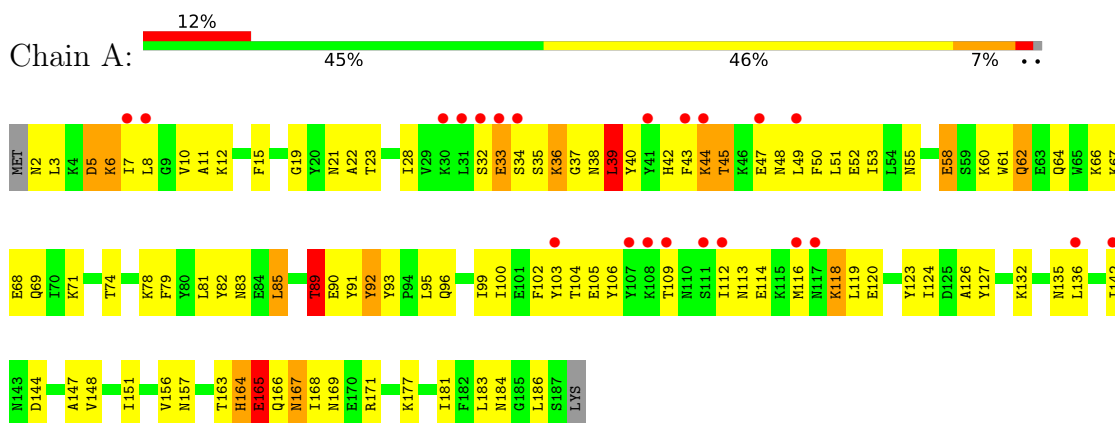
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	O	0	0
			1	1		
4	D	5	Total	O	0	0
			5	5		

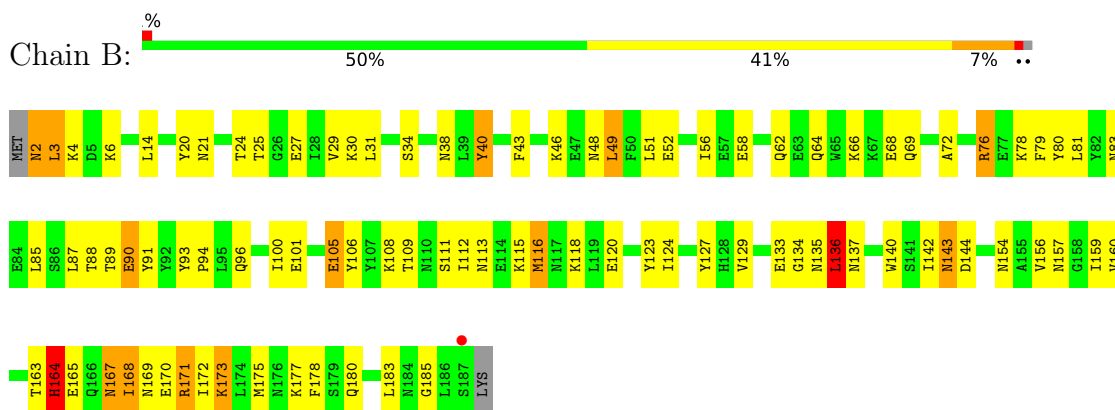
3 Residue-property plots [i](#)

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.

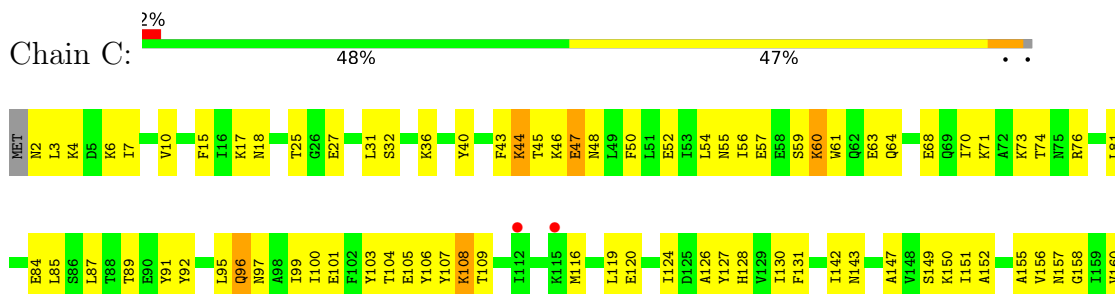
- Molecule 1: Transcriptional regulator qacR



- Molecule 1: Transcriptional regulator qacR

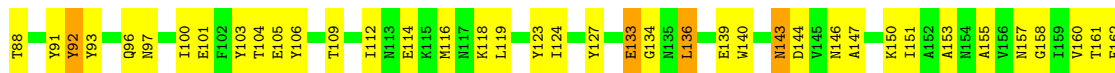
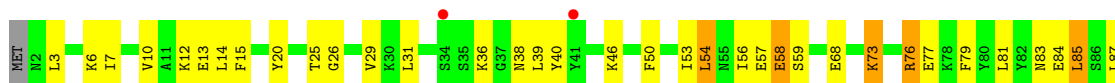


- Molecule 1: Transcriptional regulator qacR





● Molecule 1: Transcriptional regulator qacR



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	171.80Å 171.80Å 94.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.98 – 2.90 48.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.3 (48.98-2.90) 98.4 (48.98-2.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.81Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.229 , 0.280 0.225 , 0.270	Depositor DCC
R_{free} test set	1568 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	75.9	Xtrriage
Anisotropy	0.174	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 60.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6333	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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: DID, SO4

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	0.39	0/1578	0.60	0/2126
1	B	0.50	0/1578	0.66	1/2126 (0.0%)
1	C	0.41	0/1578	0.63	0/2126
1	D	0.45	0/1578	0.61	0/2126
All	All	0.44	0/6312	0.63	1/8504 (0.0%)

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	B	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	3	LEU	CA-CB-CG	6.78	130.90	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	40	TYR	Sidechain
1	B	91	TYR	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	1547	0	1536	113	1
1	B	1547	0	1536	78	0
1	C	1547	0	1536	84	0
1	D	1547	0	1536	97	0
2	A	10	0	0	0	0
2	B	20	0	0	1	0
2	C	30	0	0	1	0
2	D	40	0	0	4	0
3	B	26	0	25	7	0
4	A	3	0	0	0	0
4	B	10	0	0	0	0
4	C	1	0	0	0	0
4	D	5	0	0	1	0
All	All	6333	0	6169	348	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 28.

All (348) 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:C:165:GLU:HB3	1:D:103:TYR:HE1	1.07	1.16
1:C:142:ILE:HD11	1:C:186:LEU:HD13	1.37	1.04
1:C:165:GLU:HB3	1:D:103:TYR:CE1	1.94	1.02
1:A:96:GLN:HG3	1:A:157:ASN:HD21	1.23	1.01
1:B:135:ASN:HD21	1:B:142:ILE:H	1.19	0.90
1:B:116:MET:O	1:B:120:GLU:HG3	1.72	0.89
1:A:66:LYS:HA	1:A:69:GLN:HE21	1.36	0.89
1:A:44:LYS:H	1:A:48:ASN:HD22	1.19	0.88
1:A:6:LYS:HE3	1:A:6:LYS:HA	1.54	0.87
1:D:163:THR:O	1:D:171:ARG:HD3	1.76	0.86
1:B:164:HIS:HA	1:B:171:ARG:NH2	1.92	0.84
1:C:55:ASN:OD1	1:C:119:LEU:HD21	1.78	0.84
1:D:96:GLN:NE2	1:D:157:ASN:HD21	1.76	0.84
1:D:143:ASN:ND2	1:D:144:ASP:H	1.76	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:ILE:HD13	1:B:168:ILE:O	1.80	0.81
1:C:97:ASN:HD22	1:D:100:ILE:HD11	1.44	0.81
1:D:14:LEU:HD11	1:D:31:LEU:HD12	1.63	0.81
1:C:97:ASN:HB2	1:D:100:ILE:HD11	1.63	0.80
1:D:3:LEU:O	1:D:7:ILE:HG12	1.82	0.79
1:B:83:ASN:ND2	1:B:156:VAL:HG21	1.99	0.78
1:C:97:ASN:HD21	1:D:101:GLU:HG3	1.48	0.77
1:A:106:TYR:HB3	1:A:112:ILE:HG13	1.66	0.76
1:B:154:ASN:OD1	3:B:193:DID:H3	1.85	0.76
1:B:120:GLU:HB3	3:B:193:DID:H5	1.68	0.75
1:D:25:THR:HG21	1:D:40:TYR:OH	1.87	0.75
1:A:39:LEU:HD22	1:A:39:LEU:H	1.52	0.74
1:A:58:GLU:HG3	1:A:119:LEU:HD21	1.68	0.74
1:B:134:GLY:HA3	1:B:140:TRP:CE2	2.22	0.74
1:A:119:LEU:O	1:A:119:LEU:HD23	1.87	0.73
1:C:97:ASN:HD22	1:D:100:ILE:CD1	2.00	0.73
1:A:7:ILE:HD13	1:A:32:SER:OG	1.89	0.73
1:C:73:LYS:HG3	1:C:74:THR:H	1.54	0.72
1:C:76:ARG:HG2	1:C:183:LEU:HD13	1.70	0.72
1:C:147:ALA:O	1:C:151:ILE:HG12	1.88	0.72
1:A:19:GLY:H	1:A:22:ALA:HB3	1.55	0.72
1:A:44:LYS:H	1:A:48:ASN:ND2	1.88	0.71
1:D:81:LEU:O	1:D:85:LEU:HD22	1.90	0.71
1:D:106:TYR:HB3	1:D:112:ILE:HG21	1.73	0.71
1:A:34:SER:HB3	1:A:38:ASN:OD1	1.92	0.69
1:A:19:GLY:O	1:A:23:THR:HG22	1.93	0.69
1:C:45:THR:HG23	1:C:47:GLU:HG3	1.74	0.69
1:C:97:ASN:ND2	1:D:101:GLU:HG3	2.07	0.69
1:C:177:LYS:O	1:C:181:ILE:HG13	1.93	0.69
1:D:163:THR:C	1:D:165:GLU:H	1.96	0.69
1:A:36:LYS:H	1:A:36:LYS:HD3	1.59	0.68
1:B:68:GLU:HG2	1:B:85:LEU:HD21	1.76	0.68
1:B:76:ARG:HG3	1:B:183:LEU:HD23	1.76	0.68
1:A:120:GLU:O	1:A:124:ILE:HG23	1.94	0.68
1:A:100:ILE:O	1:A:104:THR:HG23	1.94	0.67
1:C:97:ASN:HD22	1:D:100:ILE:CG1	2.07	0.67
1:C:70:ILE:H	1:C:70:ILE:HD12	1.60	0.67
1:D:20:TYR:CZ	1:D:46:LYS:HE2	2.29	0.67
1:C:120:GLU:O	1:C:124:ILE:HG12	1.95	0.66
1:D:100:ILE:O	1:D:104:THR:HG23	1.95	0.66
1:A:60:LYS:HB3	1:A:91:TYR:CD1	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:HIS:HA	1:B:171:ARG:CZ	2.25	0.65
1:A:177:LYS:HE3	1:A:181:ILE:HD11	1.77	0.65
1:C:165:GLU:HG3	1:C:166:GLN:OE1	1.96	0.65
1:D:158:GLY:O	1:D:162:PHE:HB2	1.96	0.65
1:C:57:GLU:HG3	1:C:95:LEU:HD12	1.77	0.65
1:D:29:VAL:HG21	1:D:36:LYS:HA	1.77	0.65
1:C:96:GLN:O	1:C:100:ILE:HG12	1.97	0.64
1:C:3:LEU:HD13	1:C:3:LEU:O	1.97	0.64
1:D:109:THR:HG21	1:D:112:ILE:HD13	1.78	0.64
1:D:68:GLU:HG2	1:D:85:LEU:HD21	1.79	0.64
1:B:120:GLU:CB	3:B:193:DID:H5	2.27	0.64
1:C:151:ILE:CG2	1:D:178:PHE:HB2	2.27	0.64
1:A:8:LEU:HD21	1:A:52:GLU:HG2	1.79	0.63
1:A:6:LYS:HA	1:A:6:LYS:CE	2.28	0.63
1:A:89:THR:HG22	1:A:90:GLU:H	1.64	0.63
1:A:47:GLU:HB2	1:A:102:PHE:HE2	1.65	0.62
1:B:58:GLU:HG2	1:B:123:TYR:CZ	2.35	0.62
1:B:90:GLU:HG2	1:B:160:VAL:CG1	2.30	0.62
1:B:96:GLN:HE21	1:B:100:ILE:HD11	1.65	0.61
1:D:12:LYS:HE3	1:D:57:GLU:OE2	2.00	0.61
1:A:47:GLU:O	1:A:51:LEU:HB2	2.00	0.61
1:B:134:GLY:HA3	1:B:140:TRP:CZ2	2.34	0.61
1:D:10:VAL:HG11	1:D:31:LEU:HB2	1.82	0.61
1:C:70:ILE:HD12	1:C:70:ILE:N	2.15	0.60
1:A:44:LYS:N	1:A:48:ASN:HD22	1.94	0.60
1:B:157:ASN:HB2	3:B:193:DID:H2	1.82	0.60
1:A:38:ASN:H	1:A:38:ASN:ND2	1.99	0.60
1:A:38:ASN:HB2	1:A:39:LEU:HD22	1.84	0.60
1:D:76:ARG:CD	1:D:183:LEU:HD23	2.33	0.59
1:B:76:ARG:HG3	1:B:183:LEU:CD2	2.32	0.59
1:D:73:LYS:HB3	1:D:77:GLU:OE1	2.03	0.59
1:A:66:LYS:HG2	1:A:69:GLN:NE2	2.18	0.59
1:A:8:LEU:HG	1:A:43:PHE:HZ	1.68	0.59
1:A:112:ILE:O	1:A:116:MET:HB3	2.02	0.59
1:A:103:TYR:HD2	1:A:116:MET:HG2	1.68	0.58
1:C:96:GLN:HG2	1:C:161:THR:HG21	1.83	0.58
1:D:96:GLN:O	1:D:100:ILE:HG23	2.03	0.58
1:B:72:ALA:HB3	1:B:78:LYS:HG2	1.85	0.58
1:A:38:ASN:N	1:A:38:ASN:HD22	1.99	0.58
1:A:39:LEU:HD22	1:A:39:LEU:N	2.19	0.58
1:A:142:ILE:HD11	1:A:186:LEU:HB3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:45:THR:CG2	1:C:47:GLU:HG3	2.33	0.57
1:B:135:ASN:ND2	1:B:142:ILE:H	1.95	0.57
1:C:151:ILE:HG23	1:D:178:PHE:HB2	1.85	0.57
1:D:81:LEU:HG	1:D:85:LEU:CD2	2.34	0.57
1:A:38:ASN:H	1:A:38:ASN:HD22	1.50	0.57
1:A:60:LYS:HB3	1:A:91:TYR:CE1	2.39	0.57
1:B:83:ASN:ND2	1:B:127:TYR:OH	2.37	0.57
1:C:155:ALA:HB1	1:D:155:ALA:HB1	1.86	0.57
1:B:2:ASN:N	1:B:2:ASN:HD22	2.00	0.57
1:C:59:SER:C	1:C:61:TRP:H	2.08	0.57
1:C:103:TYR:CE1	1:D:165:GLU:HG2	2.40	0.57
1:C:43:PHE:O	1:C:44:LYS:HB2	2.05	0.56
1:D:103:TYR:HD2	1:D:116:MET:HG2	1.70	0.56
1:A:47:GLU:HB2	1:A:102:PHE:CE2	2.41	0.56
1:C:163:THR:O	1:C:171:ARG:HD2	2.05	0.56
1:C:97:ASN:ND2	1:D:100:ILE:HD11	2.17	0.56
1:A:81:LEU:HG	1:A:85:LEU:HD22	1.88	0.56
1:A:90:GLU:HG2	1:A:91:TYR:CD2	2.41	0.56
1:A:44:LYS:HB2	1:A:48:ASN:ND2	2.21	0.55
1:D:96:GLN:NE2	1:D:157:ASN:ND2	2.53	0.55
1:D:114:GLU:O	1:D:118:LYS:HG3	2.06	0.55
1:A:36:LYS:H	1:A:36:LYS:CD	2.18	0.55
1:B:143:ASN:HD22	1:B:144:ASP:N	2.04	0.55
1:D:92:TYR:CD2	1:D:123:TYR:CZ	2.94	0.55
1:C:36:LYS:HG2	1:C:40:TYR:CE1	2.41	0.55
1:B:143:ASN:HD22	1:B:143:ASN:C	2.09	0.55
1:C:10:VAL:HG11	1:C:31:LEU:HB3	1.88	0.55
1:A:15:PHE:CD2	1:A:23:THR:HG21	2.41	0.54
1:A:21:ASN:OD1	1:A:105:GLU:OE2	2.25	0.54
1:A:15:PHE:CE1	1:A:50:PHE:HD2	2.25	0.54
1:A:60:LYS:HD3	1:A:91:TYR:CZ	2.41	0.54
1:A:66:LYS:C	1:A:68:GLU:H	2.11	0.54
1:B:79:PHE:CD2	1:B:183:LEU:HD13	2.42	0.54
1:C:184:ASN:O	1:D:185:GLY:HA2	2.08	0.54
1:D:112:ILE:O	1:D:116:MET:HB2	2.07	0.54
1:A:11:ALA:HB3	1:A:53:ILE:HD11	1.89	0.54
1:D:134:GLY:HA3	1:D:140:TRP:CZ2	2.43	0.54
1:D:143:ASN:CG	1:D:144:ASP:H	2.11	0.54
1:A:45:THR:HG23	1:A:48:ASN:HB2	1.88	0.54
1:B:66:LYS:HA	1:B:69:GLN:HE21	1.73	0.54
1:B:133:GLU:HA	1:B:133:GLU:OE1	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ILE:C	1:A:114:GLU:H	2.10	0.54
1:C:25:THR:HG22	1:C:46:LYS:HB2	1.90	0.54
1:B:120:GLU:CD	3:B:193:DID:HN11	2.10	0.53
1:D:68:GLU:HG2	1:D:85:LEU:CD2	2.38	0.53
1:B:62:GLN:HA	1:B:62:GLN:NE2	2.22	0.53
1:D:83:ASN:O	1:D:175:MET:HG2	2.09	0.53
1:A:40:TYR:HA	1:A:43:PHE:O	2.09	0.53
1:C:64:GLN:O	1:C:68:GLU:HG3	2.09	0.53
1:A:132:LYS:O	1:A:136:LEU:HD13	2.09	0.53
1:D:109:THR:HB	1:D:112:ILE:HB	1.90	0.52
1:D:76:ARG:HD2	1:D:183:LEU:HD23	1.91	0.52
1:C:6:LYS:HG3	1:C:32:SER:HB2	1.91	0.52
1:B:58:GLU:HG2	1:B:123:TYR:CE2	2.44	0.52
1:A:81:LEU:HG	1:A:85:LEU:CD2	2.40	0.52
1:D:163:THR:C	1:D:165:GLU:N	2.64	0.51
1:D:168:ILE:O	1:D:172:ILE:HG13	2.10	0.51
1:A:62:GLN:OE1	1:A:126:ALA:HB2	2.11	0.51
1:B:25:THR:O	1:B:29:VAL:HG23	2.11	0.51
1:D:3:LEU:HD13	1:D:38:ASN:ND2	2.25	0.51
1:D:103:TYR:CD2	1:D:116:MET:HG2	2.43	0.51
1:D:96:GLN:CG	1:D:161:THR:HG21	2.40	0.51
1:B:168:ILE:O	1:B:172:ILE:HG12	2.10	0.51
1:C:97:ASN:CB	1:D:100:ILE:HD11	2.38	0.51
1:A:164:HIS:HA	1:A:171:ARG:CZ	2.41	0.51
1:C:70:ILE:H	1:C:70:ILE:CD1	2.23	0.51
1:D:84:GLU:HG3	1:D:172:ILE:HG23	1.92	0.51
1:D:116:MET:HE3	1:D:119:LEU:HD12	1.92	0.51
1:B:170:GLU:HA	1:B:170:GLU:OE2	2.10	0.51
1:A:90:GLU:HG2	1:A:91:TYR:CE2	2.46	0.51
1:A:39:LEU:H	1:A:39:LEU:CD2	2.19	0.51
1:A:91:TYR:O	1:A:93:TYR:N	2.44	0.51
1:C:25:THR:CG2	1:C:46:LYS:HB2	2.41	0.51
1:D:163:THR:HA	1:D:165:GLU:OE1	2.10	0.51
1:D:134:GLY:HA3	1:D:140:TRP:CE2	2.45	0.50
1:C:157:ASN:O	1:C:161:THR:HG23	2.11	0.50
1:C:183:LEU:C	1:C:185:GLY:H	2.15	0.50
1:D:147:ALA:O	1:D:151:ILE:HG12	2.12	0.50
1:B:96:GLN:O	1:B:100:ILE:HG12	2.12	0.50
1:D:124:ILE:HD13	1:D:150:LYS:HG2	1.94	0.50
1:D:153:ALA:O	1:D:157:ASN:HB2	2.11	0.50
1:B:120:GLU:CD	3:B:193:DID:N1	2.65	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:LEU:O	1:B:52:GLU:HB3	2.12	0.49
1:A:114:GLU:O	1:A:118:LYS:HB2	2.12	0.49
1:C:17:LYS:HE2	1:C:18:ASN:OD1	2.11	0.49
1:C:97:ASN:HD22	1:D:100:ILE:HG12	1.78	0.49
1:B:106:TYR:C	1:B:108:LYS:H	2.14	0.49
1:A:164:HIS:O	1:A:165:GLU:HB3	2.12	0.49
1:A:36:LYS:HD3	1:A:36:LYS:N	2.25	0.49
1:D:105:GLU:OE1	1:D:106:TYR:HE2	1.96	0.49
1:C:4:LYS:HG3	1:C:43:PHE:CZ	2.48	0.49
1:C:7:ILE:HD13	1:C:32:SER:OG	2.12	0.49
1:D:96:GLN:HA	1:D:96:GLN:OE1	2.12	0.49
1:B:20:TYR:CZ	1:B:46:LYS:HE3	2.48	0.49
1:D:105:GLU:OE1	1:D:106:TYR:CE2	2.66	0.49
1:A:147:ALA:O	1:A:151:ILE:HG13	2.12	0.49
1:D:106:TYR:CB	1:D:112:ILE:HG21	2.42	0.49
1:B:129:VAL:O	1:B:133:GLU:HG2	2.13	0.48
1:C:54:LEU:HD23	1:C:119:LEU:CD1	2.43	0.48
1:D:26:GLY:HA3	2:D:189:SO4:O2	2.13	0.48
1:C:131:PHE:CE1	1:C:186:LEU:HD11	2.48	0.48
1:A:177:LYS:HE2	1:B:144:ASP:OD2	2.14	0.48
1:A:6:LYS:HE3	1:A:6:LYS:CA	2.33	0.48
1:A:6:LYS:O	1:A:10:VAL:HG23	2.14	0.48
1:C:103:TYR:CE1	1:C:107:TYR:HB3	2.49	0.48
1:C:106:TYR:C	1:C:108:LYS:N	2.65	0.48
1:D:53:ILE:O	1:D:57:GLU:HG2	2.13	0.48
1:A:37:GLY:O	1:A:38:ASN:C	2.51	0.48
1:D:88:THR:HG22	1:D:168:ILE:HD11	1.96	0.48
1:D:83:ASN:HB3	1:D:175:MET:SD	2.54	0.48
1:A:19:GLY:N	1:A:22:ALA:HB3	2.27	0.47
1:C:105:GLU:O	1:C:105:GLU:HG2	2.14	0.47
1:C:126:ALA:O	1:C:130:ILE:HD13	2.14	0.47
1:B:43:PHE:O	1:B:48:ASN:HB3	2.14	0.47
1:B:173:LYS:HD3	1:B:173:LYS:C	2.35	0.47
1:D:54:LEU:HD23	1:D:116:MET:HE3	1.96	0.47
1:B:24:THR:OG1	1:B:27:GLU:HG3	2.13	0.47
1:D:39:LEU:C	1:D:39:LEU:HD23	2.34	0.47
1:D:83:ASN:HD21	1:D:127:TYR:HE1	1.62	0.47
1:C:27:GLU:O	1:C:31:LEU:HB2	2.15	0.47
1:A:96:GLN:NE2	1:A:157:ASN:OD1	2.45	0.47
1:A:114:GLU:O	1:A:118:LYS:N	2.48	0.47
1:A:64:GLN:OE1	1:A:89:THR:HG23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:ASP:OD1	1:B:177:LYS:HE2	2.15	0.46
1:B:144:ASP:N	1:B:144:ASP:OD1	2.48	0.46
1:A:64:GLN:O	1:A:68:GLU:HG3	2.15	0.46
1:B:88:THR:CG2	1:B:172:ILE:HD11	2.45	0.46
1:C:158:GLY:HA3	1:D:158:GLY:HA3	1.97	0.46
1:A:66:LYS:CA	1:A:69:GLN:HE21	2.18	0.46
1:A:103:TYR:CD2	1:A:116:MET:HG2	2.49	0.46
1:C:96:GLN:NE2	1:C:157:ASN:HD21	2.14	0.46
1:C:156:VAL:O	1:C:160:VAL:HG23	2.14	0.46
1:B:14:LEU:HD11	1:B:31:LEU:HD12	1.98	0.46
1:C:100:ILE:O	1:C:104:THR:HG23	2.16	0.46
1:D:56:ILE:HG21	2:D:191:SO4:O1	2.16	0.46
1:A:40:TYR:O	1:A:44:LYS:HD2	2.16	0.46
1:D:56:ILE:HD13	2:D:191:SO4:O1	2.15	0.46
1:B:163:THR:HG22	1:B:171:ARG:HD3	1.97	0.46
1:D:15:PHE:CE1	1:D:50:PHE:HD2	2.33	0.46
1:D:133:GLU:HA	1:D:136:LEU:HD11	1.97	0.46
1:A:2:ASN:N	1:A:6:LYS:HD2	2.30	0.46
1:B:87:LEU:C	1:B:87:LEU:HD23	2.37	0.46
1:C:95:LEU:O	1:C:99:ILE:HG13	2.16	0.46
1:D:58:GLU:HG3	1:D:59:SER:N	2.30	0.46
1:D:79:PHE:CD2	1:D:183:LEU:HD13	2.50	0.46
1:A:74:THR:O	1:A:78:LYS:HG3	2.16	0.45
1:C:128:HIS:HA	1:C:149:SER:OG	2.16	0.45
1:A:33:GLU:N	1:A:33:GLU:OE2	2.48	0.45
1:A:167:ASN:ND2	1:A:169:ASN:H	2.13	0.45
1:A:38:ASN:ND2	1:A:38:ASN:N	2.59	0.45
1:A:58:GLU:HG2	1:A:92:TYR:CD2	2.50	0.45
1:B:38:ASN:HD22	1:D:13:GLU:CD	2.20	0.45
1:A:35:SER:OG	1:A:38:ASN:ND2	2.50	0.45
1:D:164:HIS:HA	1:D:171:ARG:NH1	2.31	0.45
1:A:163:THR:O	1:A:166:GLN:HG2	2.16	0.45
1:C:163:THR:C	1:C:165:GLU:HG2	2.36	0.45
1:D:76:ARG:HB2	1:D:139:GLU:OE1	2.17	0.45
1:A:36:LYS:HG2	1:A:37:GLY:N	2.31	0.45
1:B:124:ILE:HD11	3:B:193:DID:N2	2.32	0.45
1:C:73:LYS:HG3	1:C:74:THR:N	2.26	0.45
1:A:167:ASN:HD22	1:A:168:ILE:N	2.15	0.45
1:B:3:LEU:HD12	1:B:34:SER:HB2	1.99	0.45
1:B:163:THR:O	1:B:165:GLU:CD	2.55	0.45
1:A:47:GLU:CB	1:A:102:PHE:HE2	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:TRP:C	1:C:63:GLU:H	2.21	0.44
1:A:151:ILE:CG2	1:B:178:PHE:HB2	2.47	0.44
1:B:4:LYS:N	2:B:191:SO4:O3	2.48	0.44
1:A:83:ASN:ND2	1:A:127:TYR:OH	2.50	0.44
1:B:106:TYR:CD1	1:B:112:ILE:HG13	2.53	0.44
1:B:52:GLU:O	1:B:56:ILE:HG12	2.18	0.44
1:B:159:ILE:CG2	1:B:175:MET:HE2	2.48	0.44
1:C:46:LYS:HE2	2:C:192:SO4:O3	2.18	0.44
1:A:167:ASN:HD22	1:A:167:ASN:C	2.20	0.44
1:B:2:ASN:N	1:B:2:ASN:ND2	2.66	0.44
1:C:150:LYS:HE3	1:C:150:LYS:HB2	1.85	0.43
1:D:96:GLN:HG3	1:D:161:THR:HG21	2.00	0.43
1:B:21:ASN:ND2	1:B:101:GLU:OE1	2.50	0.43
1:A:156:VAL:O	1:A:157:ASN:C	2.57	0.43
1:B:81:LEU:CD1	1:B:85:LEU:HD13	2.47	0.43
1:B:136:LEU:N	1:B:136:LEU:CD1	2.80	0.43
1:B:168:ILE:HD13	1:B:168:ILE:C	2.39	0.43
1:B:135:ASN:C	1:B:137:ASN:H	2.22	0.43
1:C:84:GLU:CG	1:C:176:ASN:HD21	2.31	0.43
1:D:39:LEU:HD23	1:D:39:LEU:O	2.19	0.43
1:A:142:ILE:CG2	1:A:148:VAL:HG21	2.49	0.43
1:B:85:LEU:O	1:B:89:THR:HG23	2.19	0.43
1:C:15:PHE:CD1	1:C:50:PHE:HD2	2.37	0.43
1:C:60:LYS:HE2	1:C:91:TYR:CE2	2.54	0.43
1:A:112:ILE:C	1:A:114:GLU:N	2.73	0.43
1:D:123:TYR:HB3	4:D:201:HOH:O	2.18	0.43
1:A:7:ILE:HD13	1:A:32:SER:CB	2.48	0.42
1:C:59:SER:C	1:C:61:TRP:N	2.72	0.42
1:C:84:GLU:HG3	1:C:176:ASN:HD21	1.84	0.42
1:A:79:PHE:CD2	1:A:183:LEU:HG	2.54	0.42
1:C:127:TYR:O	1:C:131:PHE:CD2	2.72	0.42
1:A:21:ASN:ND2	1:A:105:GLU:OE1	2.45	0.42
1:C:52:GLU:O	1:C:56:ILE:HG12	2.19	0.42
1:C:61:TRP:C	1:C:63:GLU:N	2.72	0.42
1:A:12:LYS:HB2	1:A:53:ILE:HD13	2.01	0.42
1:B:93:TYR:HA	1:B:94:PRO:HD3	1.88	0.42
1:A:85:LEU:HD12	1:A:85:LEU:HA	1.91	0.42
1:C:158:GLY:O	1:C:162:PHE:HD1	2.02	0.42
1:D:183:LEU:HA	1:D:183:LEU:HD12	1.74	0.42
1:A:144:ASP:OD2	1:A:144:ASP:N	2.53	0.42
1:B:51:LEU:HD13	1:B:115:LYS:HD2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:ILE:HB	1:B:175:MET:HE1	2.00	0.42
1:C:70:ILE:N	1:C:70:ILE:CD1	2.82	0.42
1:A:28:ILE:HD13	1:A:49:LEU:HD21	2.02	0.42
1:C:45:THR:O	1:C:48:ASN:N	2.44	0.42
1:C:87:LEU:HD11	1:C:175:MET:HB2	2.01	0.42
1:C:183:LEU:HD23	1:C:183:LEU:HA	1.79	0.42
1:A:51:LEU:HD12	1:A:51:LEU:HA	1.82	0.42
1:A:184:ASN:ND2	1:B:185:GLY:HA2	2.35	0.41
1:A:39:LEU:HD12	1:A:43:PHE:HD1	1.85	0.41
1:D:163:THR:O	1:D:165:GLU:N	2.53	0.41
1:C:101:GLU:CG	1:D:97:ASN:HD21	2.34	0.41
1:C:156:VAL:HG13	1:C:175:MET:CE	2.51	0.41
1:D:68:GLU:CG	1:D:85:LEU:HD21	2.47	0.41
1:B:105:GLU:OE1	1:B:106:TYR:CE2	2.74	0.41
1:A:58:GLU:HG3	1:A:119:LEU:CD2	2.44	0.41
1:A:61:TRP:CZ2	1:A:82:TYR:CE1	3.08	0.41
1:A:168:ILE:HA	1:A:171:ARG:HG3	2.03	0.41
1:B:3:LEU:HA	1:B:6:LYS:HB3	2.01	0.41
1:B:96:GLN:NE2	1:B:100:ILE:HD11	2.31	0.41
1:A:5:ASP:O	1:A:6:LYS:HE3	2.21	0.41
1:A:132:LYS:O	1:A:135:ASN:HB2	2.21	0.41
1:C:186:LEU:HD23	1:C:186:LEU:HA	1.91	0.41
1:D:56:ILE:HG21	2:D:191:SO4:S	2.61	0.41
1:A:2:ASN:OD1	1:A:3:LEU:N	2.54	0.41
1:B:133:GLU:OE1	1:B:133:GLU:CA	2.68	0.41
1:D:87:LEU:HD23	1:D:160:VAL:HG22	2.01	0.41
1:A:15:PHE:HD2	1:A:23:THR:HG21	1.82	0.41
1:A:66:LYS:HA	1:A:69:GLN:NE2	2.19	0.41
1:A:66:LYS:C	1:A:68:GLU:N	2.73	0.41
1:B:167:ASN:ND2	1:B:169:ASN:H	2.18	0.41
1:C:81:LEU:O	1:C:85:LEU:HG	2.20	0.41
1:D:91:TYR:O	1:D:93:TYR:N	2.52	0.41
1:A:58:GLU:HG2	1:A:92:TYR:CE2	2.56	0.41
1:D:14:LEU:HD23	1:D:14:LEU:HA	1.84	0.41
1:D:81:LEU:HG	1:D:85:LEU:HD21	2.02	0.41
1:C:167:ASN:HB3	1:C:170:GLU:HB2	2.01	0.40
1:A:52:GLU:O	1:A:55:ASN:HB3	2.21	0.40
1:B:40:TYR:CZ	1:D:6:LYS:HD3	2.56	0.40
1:B:163:THR:O	1:B:165:GLU:N	2.54	0.40
1:D:54:LEU:HA	1:D:54:LEU:HD12	1.89	0.40
1:B:109:THR:HG22	1:B:111:SER:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:LEU:HD11	1:C:99:ILE:HG12	2.04	0.40
1:A:95:LEU:O	1:A:99:ILE:HG13	2.22	0.40
1:A:123:TYR:OH	1:A:157:ASN:HB2	2.21	0.40
1:B:80:TYR:OH	1:B:180:GLN:HG3	2.22	0.40
1:D:29:VAL:HG21	1:D:36:LYS:CA	2.49	0.40
1:D:83:ASN:N	1:D:83:ASN:HD22	2.19	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:42:HIS:CE1	1:A:42:HIS:CE1[8_445]	2.17	0.03

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	184/188 (98%)	145 (79%)	29 (16%)	10 (5%)	2	6
1	B	184/188 (98%)	168 (91%)	14 (8%)	2 (1%)	14	42
1	C	184/188 (98%)	146 (79%)	29 (16%)	9 (5%)	2	8
1	D	184/188 (98%)	161 (88%)	14 (8%)	9 (5%)	2	8
All	All	736/752 (98%)	620 (84%)	86 (12%)	30 (4%)	3	11

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	THR
1	A	165	GLU
1	B	164	HIS
1	C	165	GLU

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Mol	Chain	Res	Type
1	D	143	ASN
1	D	166	GLN
1	A	33	GLU
1	A	44	LYS
1	A	92	TYR
1	A	164	HIS
1	C	108	LYS
1	C	109	THR
1	D	92	TYR
1	B	136	LEU
1	C	92	TYR
1	A	113	ASN
1	C	71	LYS
1	D	73	LYS
1	D	164	HIS
1	A	39	LEU
1	A	67	LYS
1	A	71	LYS
1	C	44	LYS
1	C	60	LYS
1	C	152	ALA
1	D	136	LEU
1	D	180	GLN
1	D	186	LEU
1	C	143	ASN
1	D	168	ILE

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	169/171 (99%)	156 (92%)	13 (8%)	13 35
1	B	169/171 (99%)	152 (90%)	17 (10%)	7 23
1	C	169/171 (99%)	161 (95%)	8 (5%)	26 59
1	D	169/171 (99%)	159 (94%)	10 (6%)	19 49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	676/684 (99%)	628 (93%)	48 (7%)	14	40

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASP
1	A	6	LYS
1	A	36	LYS
1	A	39	LEU
1	A	45	THR
1	A	58	GLU
1	A	62	GLN
1	A	85	LEU
1	A	89	THR
1	A	109	THR
1	A	118	LYS
1	A	165	GLU
1	A	167	ASN
1	B	2	ASN
1	B	30	LYS
1	B	49	LEU
1	B	64	GLN
1	B	76	ARG
1	B	90	GLU
1	B	105	GLU
1	B	113	ASN
1	B	116	MET
1	B	118	LYS
1	B	136	LEU
1	B	143	ASN
1	B	164	HIS
1	B	167	ASN
1	B	168	ILE
1	B	171	ARG
1	B	173	LYS
1	C	2	ASN
1	C	47	GLU
1	C	89	THR
1	C	96	GLN
1	C	116	MET
1	C	166	GLN
1	C	168	ILE

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Mol	Chain	Res	Type
1	C	170	GLU
1	D	54	LEU
1	D	58	GLU
1	D	76	ARG
1	D	85	LEU
1	D	133	GLU
1	D	146	ASN
1	D	164	HIS
1	D	165	GLU
1	D	167	ASN
1	D	170	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	38	ASN
1	A	48	ASN
1	A	62	GLN
1	A	69	GLN
1	A	75	ASN
1	A	83	ASN
1	A	96	GLN
1	A	113	ASN
1	A	128	HIS
1	A	154	ASN
1	A	157	ASN
1	A	166	GLN
1	A	167	ASN
1	A	184	ASN
1	B	21	ASN
1	B	38	ASN
1	B	55	ASN
1	B	62	GLN
1	B	69	GLN
1	B	83	ASN
1	B	96	GLN
1	B	113	ASN
1	B	121	ASN
1	B	135	ASN
1	B	143	ASN
1	B	166	GLN

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Mol	Chain	Res	Type
1	B	167	ASN
1	B	180	GLN
1	C	75	ASN
1	C	83	ASN
1	C	96	GLN
1	C	97	ASN
1	C	117	ASN
1	C	143	ASN
1	C	154	ASN
1	C	167	ASN
1	C	176	ASN
1	D	18	ASN
1	D	38	ASN
1	D	62	GLN
1	D	75	ASN
1	D	83	ASN
1	D	96	GLN
1	D	97	ASN
1	D	121	ASN
1	D	143	ASN

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](#)

21 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	SO4	D	192	-	4,4,4	0.24	0	6,6,6	0.10	0
2	SO4	D	191	-	4,4,4	0.22	0	6,6,6	0.06	0
2	SO4	C	193	-	4,4,4	0.25	0	6,6,6	0.09	0
2	SO4	C	192	-	4,4,4	0.25	0	6,6,6	0.08	0
2	SO4	D	190	-	4,4,4	0.23	0	6,6,6	0.14	0
2	SO4	C	194	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	B	191	-	4,4,4	0.28	0	6,6,6	0.09	0
2	SO4	D	194	-	4,4,4	0.25	0	6,6,6	0.06	0
2	SO4	B	190	-	4,4,4	0.28	0	6,6,6	0.11	0
2	SO4	D	195	-	4,4,4	0.25	0	6,6,6	0.05	0
2	SO4	D	189	-	4,4,4	0.25	0	6,6,6	0.14	0
3	DID	B	193	-	27,27,27	3.38	12 (44%)	26,34,34	1.10	2 (7%)
2	SO4	D	196	-	4,4,4	0.25	0	6,6,6	0.06	0
2	SO4	B	189	-	4,4,4	0.27	0	6,6,6	0.08	0
2	SO4	C	191	-	4,4,4	0.26	0	6,6,6	0.05	0
2	SO4	A	189	-	4,4,4	0.21	0	6,6,6	0.11	0
2	SO4	B	192	-	4,4,4	0.25	0	6,6,6	0.07	0
2	SO4	D	193	-	4,4,4	0.25	0	6,6,6	0.07	0
2	SO4	C	189	-	4,4,4	0.26	0	6,6,6	0.05	0
2	SO4	A	190	-	4,4,4	0.26	0	6,6,6	0.07	0
2	SO4	C	190	-	4,4,4	0.25	0	6,6,6	0.07	0

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	DID	B	193	-	-	6/19/19/19	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	193	DID	C3'-C4'	7.26	1.51	1.39
3	B	193	DID	C6-C5	6.76	1.51	1.38
3	B	193	DID	C2'-C1'	6.37	1.51	1.38
3	B	193	DID	C2-C1	5.73	1.50	1.38
3	B	193	DID	C3-C4	5.71	1.49	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	193	DID	C5'-C6'	5.46	1.48	1.38
3	B	193	DID	C9-N1	3.68	1.43	1.33
3	B	193	DID	O1-C1	3.51	1.45	1.37
3	B	193	DID	C9'-N2'	3.04	1.41	1.28
3	B	193	DID	O1'-C1'	2.60	1.43	1.37
3	B	193	DID	C9'-N1'	-2.40	1.27	1.33
3	B	193	DID	C5'-C4'	-2.33	1.35	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	193	DID	C8'-C11-C10	-2.20	103.25	114.42
3	B	193	DID	O1'-C7'-C8'	-2.13	100.40	108.33

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	193	DID	C5-C4-C9-N1
3	B	193	DID	C3-C4-C9-N1
3	B	193	DID	C5'-C4'-C9'-N1'
3	B	193	DID	C3'-C4'-C9'-N1'
3	B	193	DID	C5'-C4'-C9'-N2'
3	B	193	DID	C3'-C4'-C9'-N2'

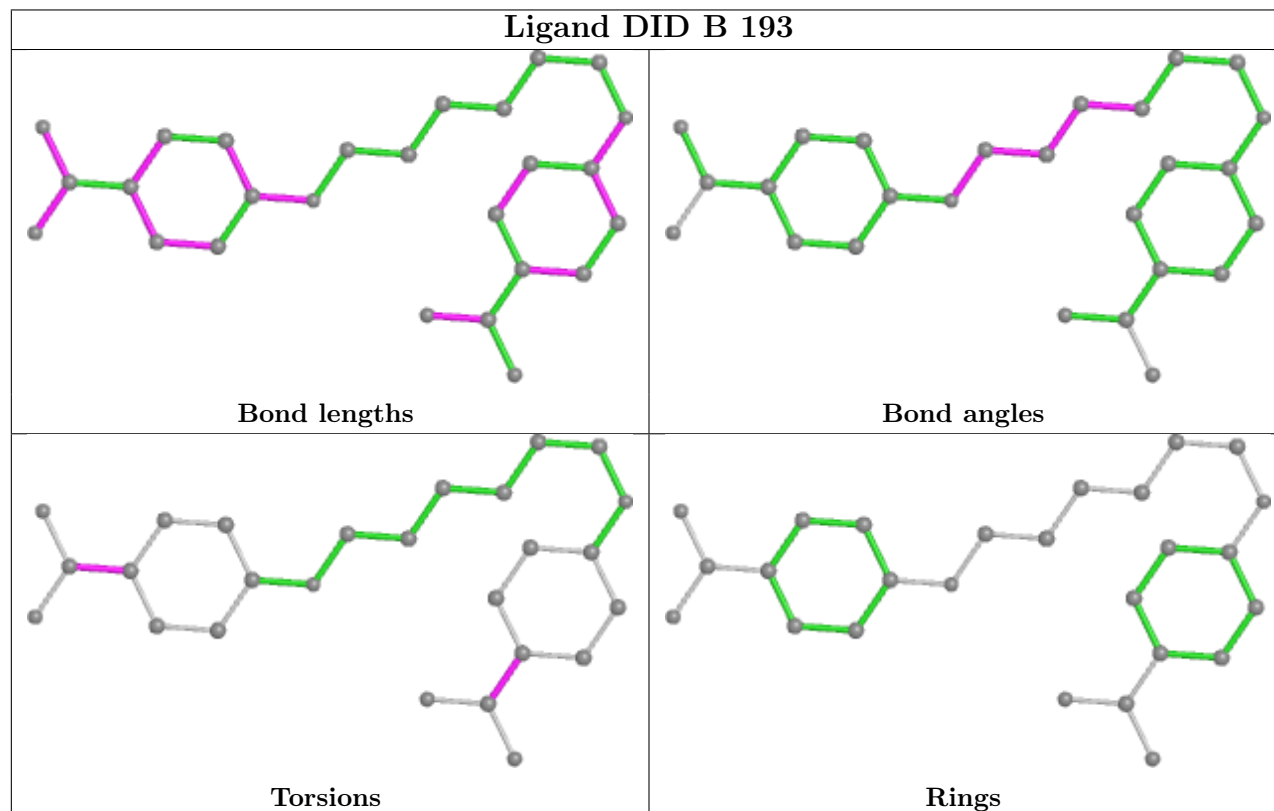
There are no ring outliers.

5 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	191	SO4	3	0
2	C	192	SO4	1	0
2	B	191	SO4	1	0
2	D	189	SO4	1	0
3	B	193	DID	7	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	186/188 (98%)	0.52	22 (11%) 4 3	41, 83, 159, 170	0
1	B	186/188 (98%)	-0.21	1 (0%) 91 91	33, 52, 89, 109	0
1	C	186/188 (98%)	0.23	3 (1%) 72 71	48, 82, 114, 130	0
1	D	186/188 (98%)	0.09	2 (1%) 80 80	45, 66, 98, 111	0
All	All	744/752 (98%)	0.16	28 (3%) 40 36	33, 70, 121, 170	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	112	ILE	6.8
1	A	116	MET	6.4
1	A	108	LYS	5.7
1	A	107	TYR	5.4
1	A	32	SER	5.2
1	A	41	TYR	5.1
1	A	111	SER	4.0
1	A	33	GLU	3.5
1	D	41	TYR	3.5
1	A	43	PHE	3.4
1	A	142	ILE	3.4
1	A	117	ASN	3.2
1	C	112	ILE	3.1
1	A	31	LEU	3.1
1	A	44	LYS	2.6
1	A	30	LYS	2.6
1	A	103	TYR	2.5
1	A	49	LEU	2.5
1	A	34	SER	2.5
1	A	136	LEU	2.4
1	B	187	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	34	SER	2.3
1	A	7	ILE	2.3
1	A	8	LEU	2.3
1	A	109	THR	2.3
1	A	47	GLU	2.3
1	C	115	LYS	2.2
1	C	161	THR	2.1

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 monosaccharides 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.

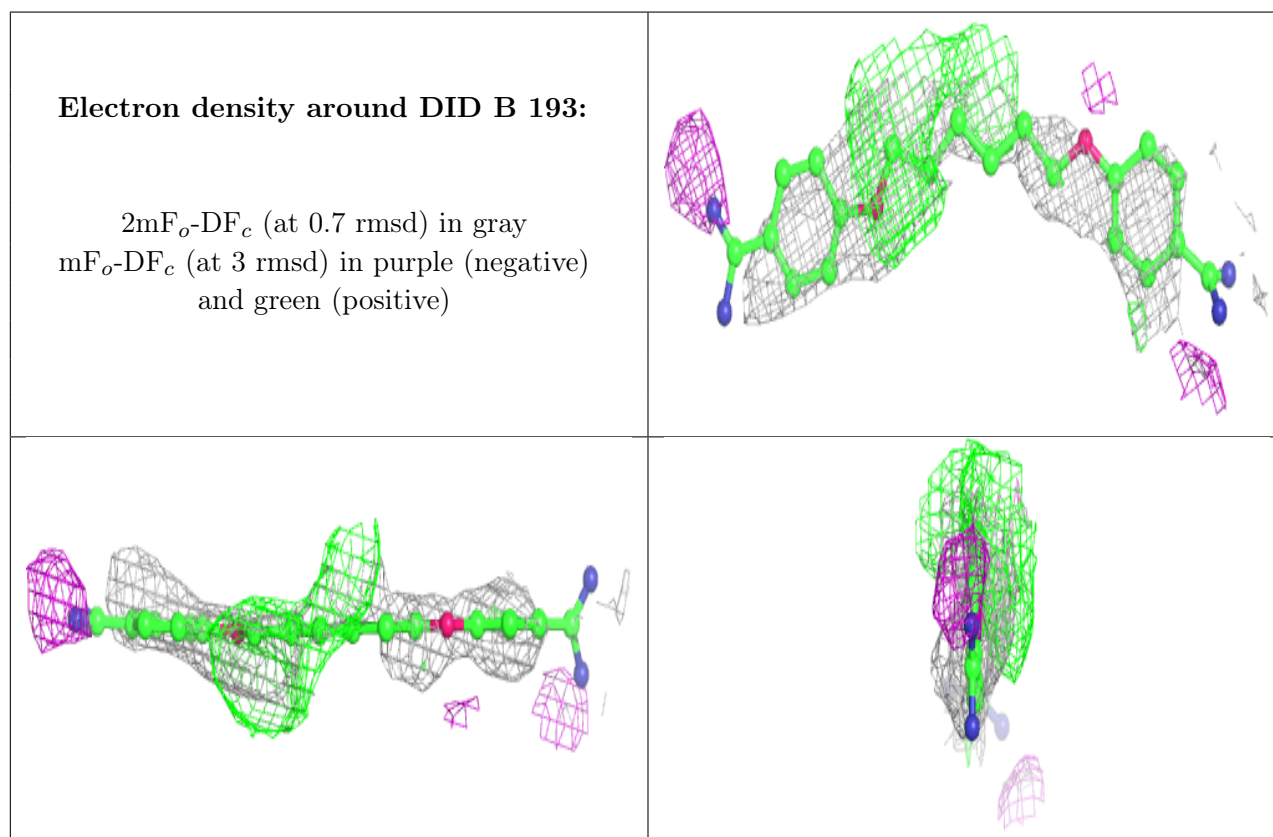
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	DID	B	193	26/26	0.50	0.57	138,143,150,152	0
2	SO4	C	191	5/5	0.59	0.27	182,182,182,183	0
2	SO4	D	190	5/5	0.62	0.34	137,137,137,138	0
2	SO4	B	192	5/5	0.63	0.31	196,196,197,197	0
2	SO4	C	190	5/5	0.69	0.29	179,179,180,180	0
2	SO4	A	190	5/5	0.70	0.55	185,186,186,186	0
2	SO4	C	194	5/5	0.70	0.44	185,185,185,186	0
2	SO4	D	194	5/5	0.72	0.39	194,194,194,195	0
2	SO4	D	193	5/5	0.74	0.24	158,159,159,159	0
2	SO4	D	195	5/5	0.77	0.39	190,190,191,191	0
2	SO4	C	193	5/5	0.78	0.32	176,177,177,177	0
2	SO4	C	192	5/5	0.81	0.18	152,152,153,153	0
2	SO4	C	189	5/5	0.82	0.38	179,179,179,179	0
2	SO4	D	196	5/5	0.83	0.19	186,187,188,188	0
2	SO4	B	191	5/5	0.87	0.20	133,134,135,135	0
2	SO4	B	190	5/5	0.89	0.19	121,121,122,123	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	D	189	5/5	0.89	0.17	129,130,130,130	0
2	SO4	A	189	5/5	0.90	0.24	139,139,139,140	0
2	SO4	B	189	5/5	0.91	0.14	128,128,129,129	0
2	SO4	D	192	5/5	0.94	0.50	133,133,134,135	0
2	SO4	D	191	5/5	0.96	0.33	156,157,157,158	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.