Full wwPDB X-ray Structure Validation Report (i)

Aug 6, 2023 – 06:34 AM EDT

PDB ID	:	1JUM
Title	:	Crystal structure of the multidrug binding transcriptional repressor QacR bound to the natural drug berberine
Authors	:	Schumacher, M.A.; Miller, M.C.; Grkovic, S.; Brown, M.H.; Skurray, R.A.; Brennan, R.G.
Deposited on	:	2001-08-24
Resolution	:	2.98 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 (i)) were used in the production of this report:

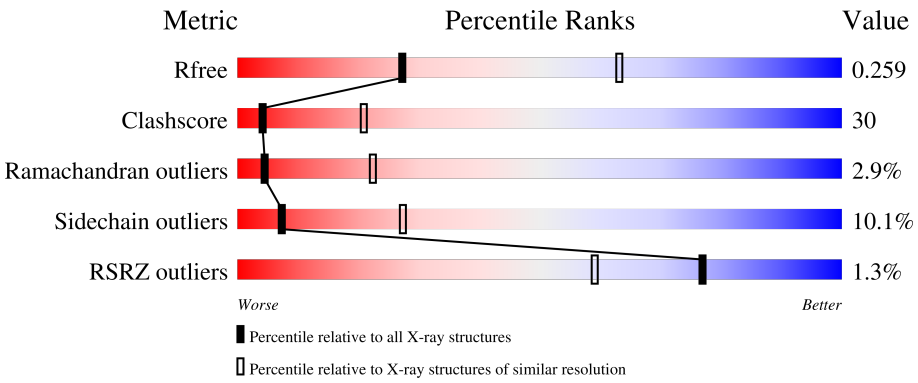
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	2754 (3.00-2.96)
Clashscore	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)
RSRZ outliers	127900	2644 (3.00-2.96)

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% 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	194	50% 38% 8% •
1	B	194	5% 37% 48% 10% • •
1	D	194	48% 41% 6% •
1	E	194	% 42% 48% 6% • •

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	B	700	-	-	-	X
2	SO4	D	1000	-	-	-	X
2	SO4	E	639	-	-	-	X
2	SO4	E	903	-	-	-	X
3	BER	A	803	-	-	-	X

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6330 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 HYPOTHETICAL TRANSCRIPTIONAL REGULATOR IN QACA 5'REGION.

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

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	72	ALA	CYS	engineered mutation	UNP P0A0N4
B	141	SER	CYS	engineered mutation	UNP P0A0N4
B	189	HIS	-	expression tag	UNP P0A0N4
B	190	HIS	-	expression tag	UNP P0A0N4
B	191	HIS	-	expression tag	UNP P0A0N4
B	192	HIS	-	expression tag	UNP P0A0N4
B	193	HIS	-	expression tag	UNP P0A0N4
B	194	HIS	-	expression tag	UNP P0A0N4
D	72	ALA	CYS	engineered mutation	UNP P0A0N4
D	141	SER	CYS	engineered mutation	UNP P0A0N4
D	189	HIS	-	expression tag	UNP P0A0N4
D	190	HIS	-	expression tag	UNP P0A0N4
D	191	HIS	-	expression tag	UNP P0A0N4
D	192	HIS	-	expression tag	UNP P0A0N4
D	193	HIS	-	expression tag	UNP P0A0N4
D	194	HIS	-	expression tag	UNP P0A0N4
A	72	ALA	CYS	engineered mutation	UNP P0A0N4
A	141	SER	CYS	engineered mutation	UNP P0A0N4
A	189	HIS	-	expression tag	UNP P0A0N4
A	190	HIS	-	expression tag	UNP P0A0N4

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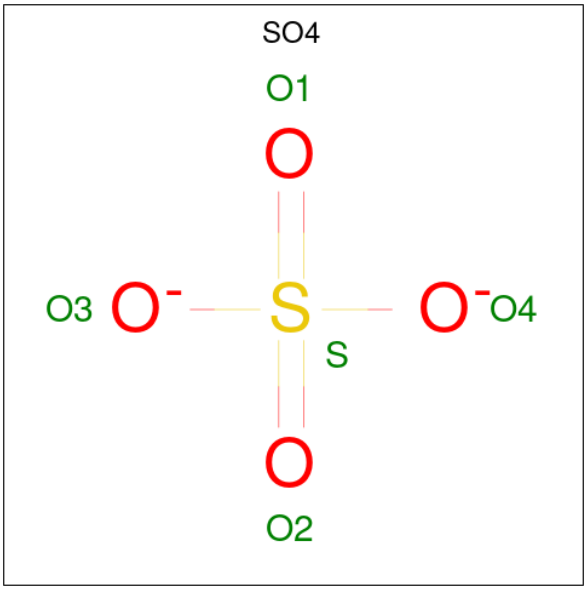
Chain	Residue	Modelled	Actual	Comment	Reference
A	191	HIS	-	expression tag	UNP P0A0N4
A	192	HIS	-	expression tag	UNP P0A0N4
A	193	HIS	-	expression tag	UNP P0A0N4
A	194	HIS	-	expression tag	UNP P0A0N4
E	72	ALA	CYS	engineered mutation	UNP P0A0N4
E	141	SER	CYS	engineered mutation	UNP P0A0N4
E	189	HIS	-	expression tag	UNP P0A0N4
E	190	HIS	-	expression tag	UNP P0A0N4
E	191	HIS	-	expression tag	UNP P0A0N4
E	192	HIS	-	expression tag	UNP P0A0N4
E	193	HIS	-	expression tag	UNP P0A0N4
E	194	HIS	-	expression tag	UNP P0A0N4

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

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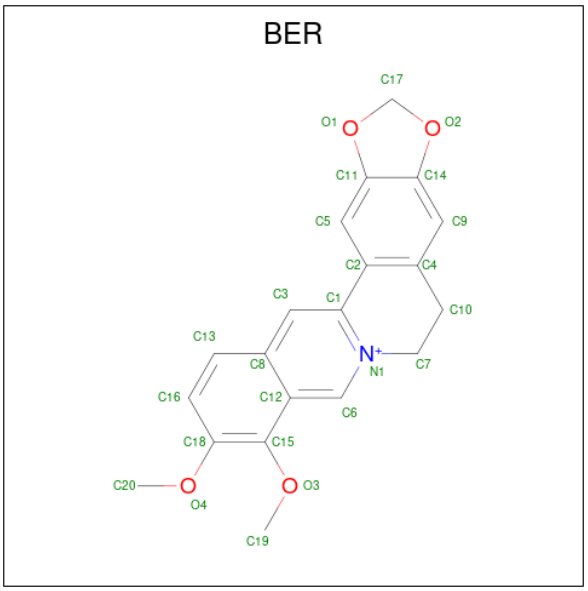
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	D	1	5	4	1	0	0
			Total	O	S		
2	D	1	5	4	1	0	0
			Total	O	S		
2	D	1	5	4	1	0	0
			Total	O	S		
2	D	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O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

- Molecule 3 is BERBERINE (three-letter code: BER) (formula: $C_{20}H_{18}NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	25	20	1	4	0	0

- Molecule 4 is water.

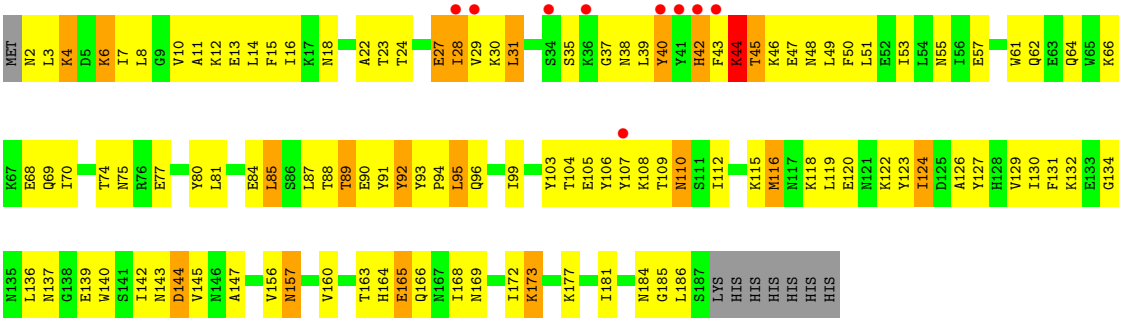
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	D	3	3	3	0	0
			Total	O		
4	A	3	3	3	0	0
			Total	O		
4	E	1	1	1	0	0

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 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.

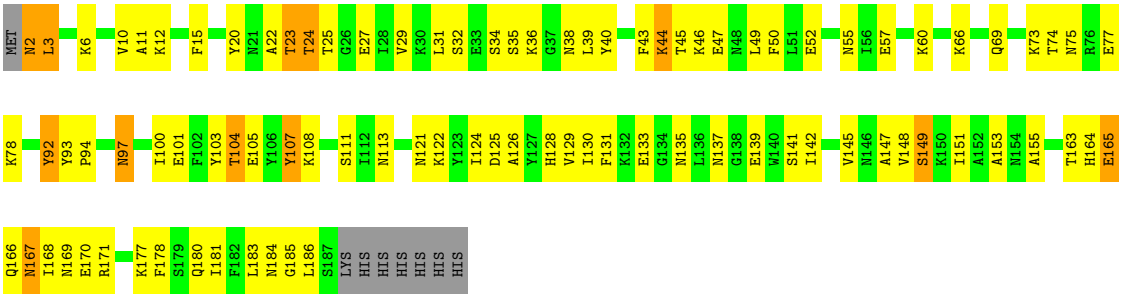
• Molecule 1: HYPOTHETICAL TRANSCRIPTIONAL REGULATOR IN QACA 5'REGION

Chain B:



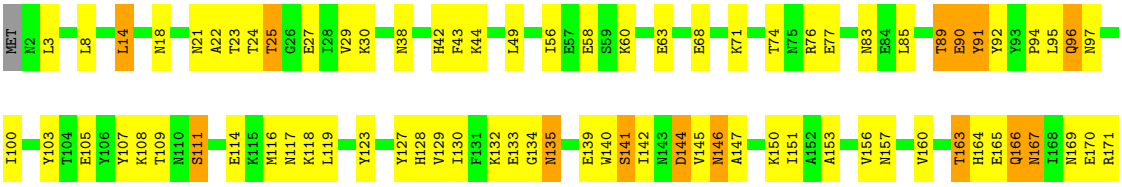
• Molecule 1: HYPOTHETICAL TRANSCRIPTIONAL REGULATOR IN QACA 5'REGION

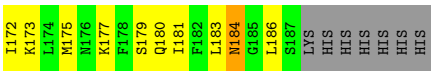
Chain D:



• Molecule 1: HYPOTHETICAL TRANSCRIPTIONAL REGULATOR IN QACA 5'REGION

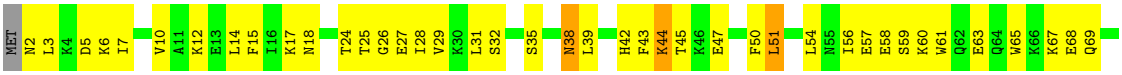
Chain A:





• Molecule 1: HYPOTHETICAL TRANSCRIPTIONAL REGULATOR IN QACA 5'REGION

Chain E:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α, β, γ	174.30Å 174.30Å 96.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.29 – 2.98 84.29 – 2.57	Depositor EDS
% Data completeness (in resolution range)	92.7 (84.29-2.98) 67.3 (84.29-2.57)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) >$ [1]	1.92 (at 2.58Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.236 , 0.284 0.208 , 0.259	Depositor DCC
R_{free} test set	2871 reflections (8.92%)	wwPDB-VP
Wilson B-factor (Å^2)	49.0	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 62.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6330	wwPDB-VP
Average B, all atoms (Å^2)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SO4, BER

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	0.66	0/1578	0.77	0/2126
1	B	0.60	0/1578	0.76	0/2126
1	D	0.63	0/1578	0.75	0/2126
1	E	0.70	0/1578	0.78	0/2126
All	All	0.65	0/6312	0.76	0/8504

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	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	40	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	74	0
1	B	1547	0	1536	125	0
1	D	1547	0	1536	89	0
1	E	1547	0	1536	97	0
2	A	30	0	0	1	0
2	B	25	0	0	1	0
2	D	20	0	0	0	0
2	E	35	0	0	2	0
3	A	25	0	18	3	0
4	A	3	0	0	0	0
4	D	3	0	0	1	0
4	E	1	0	0	0	0
All	All	6330	0	6162	369	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 30.

All (369) 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:D:2:ASN:ND2	1:D:3:LEU:H	1.49	1.10
1:B:45:THR:HG23	1:B:48:ASN:HB2	1.37	1.01
1:B:115:LYS:O	1:B:118:LYS:HG2	1.64	0.96
1:A:109:THR:HG22	1:A:111:SER:HB3	1.48	0.95
1:D:2:ASN:HD22	1:D:3:LEU:N	1.68	0.92
1:D:2:ASN:HD22	1:D:3:LEU:H	1.02	0.92
1:B:18:ASN:HD22	1:B:23:THR:HG22	1.33	0.92
1:E:24:THR:HG23	1:E:27:GLU:HB2	1.54	0.86
1:B:89:THR:HG22	1:B:90:GLU:H	1.38	0.86
1:D:24:THR:HG23	1:D:27:GLU:HB2	1.56	0.86
1:B:164:HIS:HB3	1:A:107:TYR:OH	1.75	0.85
1:B:39:LEU:H	1:B:39:LEU:HD22	1.43	0.84
1:B:62:GLN:HE22	1:B:126:ALA:HB2	1.43	0.84
1:A:68:GLU:HG2	1:A:85:LEU:HD11	1.58	0.84
1:B:29:VAL:HG22	1:B:39:LEU:HD21	1.59	0.83
1:D:165:GLU:HG2	1:D:166:GLN:OE1	1.78	0.83
1:A:103:TYR:O	1:A:107:TYR:HB2	1.78	0.83
1:B:66:LYS:HA	1:B:69:GLN:HE21	1.44	0.82
1:D:163:THR:HA	1:D:165:GLU:OE2	1.80	0.81
1:B:177:LYS:O	1:B:181:ILE:HG12	1.79	0.81
1:D:147:ALA:O	1:D:151:ILE:HG13	1.81	0.80
1:E:95:LEU:O	1:E:99:ILE:HG13	1.81	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:163:THR:HA	1:E:165:GLU:OE2	1.82	0.80
1:B:39:LEU:H	1:B:39:LEU:CD2	1.95	0.80
1:A:180:GLN:O	1:A:184:ASN:HB2	1.82	0.79
1:E:24:THR:CG2	1:E:27:GLU:HB2	2.14	0.78
1:A:30:LYS:O	1:A:30:LYS:HD3	1.84	0.78
1:A:147:ALA:O	1:A:151:ILE:HG12	1.84	0.77
1:E:143:ASN:CG	1:E:144:ASP:H	1.90	0.76
1:A:24:THR:OG1	1:A:27:GLU:HG3	1.86	0.75
1:D:2:ASN:ND2	1:D:3:LEU:N	2.28	0.75
1:E:3:LEU:O	1:E:7:ILE:HD13	1.86	0.74
1:A:167:ASN:ND2	1:A:169:ASN:H	1.86	0.74
1:D:10:VAL:HG21	1:D:32:SER:HB3	1.70	0.73
1:D:131:PHE:HB3	1:D:142:ILE:HD11	1.71	0.73
1:B:39:LEU:HD22	1:B:39:LEU:N	2.04	0.72
1:D:177:LYS:O	1:D:181:ILE:HG12	1.90	0.71
1:A:166:GLN:N	1:A:166:GLN:HE21	1.86	0.71
1:B:120:GLU:O	1:B:124:ILE:HG23	1.91	0.71
1:B:18:ASN:HB3	1:B:22:ALA:HB1	1.73	0.70
1:D:125:ASP:O	1:D:128:HIS:HB3	1.90	0.70
1:B:2:ASN:HB2	2:B:700:SO4:S	2.32	0.70
1:A:132:LYS:O	1:A:135:ASN:HB2	1.90	0.70
1:B:10:VAL:HG12	1:B:14:LEU:HD13	1.73	0.69
1:B:28:ILE:HD13	1:B:49:LEU:HD21	1.74	0.69
1:D:145:VAL:O	1:D:149:SER:HB2	1.92	0.69
1:B:62:GLN:NE2	1:B:126:ALA:HB2	2.07	0.69
1:A:114:GLU:HG2	1:A:118:LYS:HE3	1.75	0.68
1:B:106:TYR:O	1:B:112:ILE:HG13	1.93	0.68
1:E:26:GLY:O	1:E:29:VAL:HG22	1.93	0.68
1:E:58:GLU:HA	1:E:123:TYR:OH	1.92	0.68
1:B:96:GLN:HG2	1:B:157:ASN:HD21	1.59	0.68
1:A:179:SER:O	1:A:183:LEU:HB2	1.93	0.67
1:E:142:ILE:HD12	1:E:186:LEU:HD22	1.75	0.67
1:B:124:ILE:HG12	1:B:124:ILE:O	1.95	0.67
1:B:64:GLN:O	1:B:68:GLU:HG3	1.94	0.67
1:B:89:THR:HG22	1:B:90:GLU:N	2.09	0.66
1:B:11:ALA:HB3	1:B:53:ILE:HD11	1.77	0.66
1:A:76:ARG:HB3	1:A:76:ARG:NH1	2.11	0.66
1:B:4:LYS:HE3	1:B:42:HIS:CE1	2.31	0.66
1:D:23:THR:HG22	1:D:23:THR:O	1.94	0.65
1:E:24:THR:HG23	1:E:27:GLU:H	1.62	0.65
1:B:120:GLU:OE1	1:A:165:GLU:HG3	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:THR:H	1:B:27:GLU:HG3	1.61	0.65
1:B:89:THR:HG22	1:B:91:TYR:H	1.60	0.65
1:E:142:ILE:CD1	1:E:186:LEU:HD22	2.27	0.65
1:B:106:TYR:HB3	1:B:112:ILE:HD11	1.78	0.64
1:B:18:ASN:HD22	1:B:23:THR:CG2	2.07	0.64
1:B:44:LYS:HE3	1:B:44:LYS:HA	1.80	0.64
1:D:155:ALA:HB1	1:E:155:ALA:HB1	1.80	0.64
1:B:3:LEU:HD11	1:B:38:ASN:HD22	1.62	0.64
1:E:164:HIS:HA	1:E:171:ARG:NH1	2.13	0.63
1:B:18:ASN:HB2	1:B:23:THR:HG23	1.79	0.63
1:D:184:ASN:O	1:E:185:GLY:HA2	1.97	0.63
1:A:146:ASN:O	1:A:150:LYS:HG3	1.99	0.63
1:D:73:LYS:HB2	1:D:73:LYS:NZ	2.14	0.63
1:D:6:LYS:O	1:D:10:VAL:HG23	1.99	0.63
1:B:105:GLU:HG2	1:B:105:GLU:O	1.99	0.62
1:D:97:ASN:ND2	1:E:100:ILE:HD12	2.14	0.62
1:A:25:THR:O	1:A:29:VAL:HG23	1.99	0.62
1:B:4:LYS:O	1:B:8:LEU:HD12	1.99	0.62
1:A:83:ASN:HB3	1:A:175:MET:CE	2.29	0.62
1:B:110:ASN:C	1:B:112:ILE:H	2.03	0.62
1:E:14:LEU:HD11	1:E:31:LEU:HD12	1.80	0.62
1:E:112:ILE:O	1:E:116:MET:HB2	1.99	0.62
1:B:24:THR:O	1:B:27:GLU:HB2	2.00	0.61
1:E:167:ASN:HD21	1:E:169:ASN:CB	2.13	0.61
1:D:107:TYR:HA	1:D:113:ASN:OD1	2.01	0.61
1:B:74:THR:OG1	1:B:77:GLU:HG3	2.01	0.61
1:D:100:ILE:O	1:D:104:THR:HG23	2.01	0.60
1:D:121:ASN:O	1:D:125:ASP:HB2	2.02	0.60
1:A:165:GLU:HG2	1:A:166:GLN:HE22	1.66	0.60
1:B:14:LEU:C	1:B:23:THR:HG21	2.22	0.60
1:E:143:ASN:CG	1:E:144:ASP:N	2.56	0.59
1:D:167:ASN:ND2	1:D:169:ASN:H	2.00	0.59
1:A:83:ASN:HB3	1:A:175:MET:HE3	1.84	0.59
1:D:164:HIS:HD2	1:E:104:THR:CG2	2.15	0.59
1:B:93:TYR:O	1:B:96:GLN:HB2	2.03	0.59
1:B:66:LYS:HA	1:B:69:GLN:NE2	2.16	0.58
1:B:7:ILE:HG21	1:B:43:PHE:HE1	1.68	0.58
1:B:70:ILE:N	1:B:70:ILE:HD13	2.18	0.58
1:B:156:VAL:O	1:B:160:VAL:HG23	2.04	0.58
1:E:165:GLU:O	1:E:166:GLN:O	2.21	0.58
1:D:73:LYS:HB2	1:D:73:LYS:HZ2	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:163:THR:C	1:E:171:ARG:HH11	2.06	0.58
1:B:6:LYS:O	1:B:10:VAL:HG23	2.05	0.57
1:A:132:LYS:HD3	2:A:801:SO4:O1	2.05	0.57
1:E:133:GLU:O	1:E:137:ASN:HB2	2.05	0.57
1:B:24:THR:HG22	1:B:27:GLU:HG3	1.87	0.56
1:B:75:ASN:HB2	1:B:139:GLU:OE2	2.04	0.56
1:D:164:HIS:HD2	1:E:104:THR:HG22	1.70	0.56
1:E:103:TYR:CD2	1:E:116:MET:HG2	2.40	0.56
1:B:15:PHE:CE1	1:B:50:PHE:HD2	2.24	0.56
1:B:3:LEU:HD11	1:B:38:ASN:ND2	2.21	0.56
1:B:14:LEU:HB3	1:B:23:THR:HG21	1.88	0.56
1:D:171:ARG:HG2	1:D:171:ARG:HH11	1.71	0.56
1:A:90:GLU:HG3	1:A:91:TYR:N	2.19	0.55
1:D:66:LYS:HD2	1:D:69:GLN:HE21	1.69	0.55
1:D:125:ASP:O	1:D:128:HIS:CB	2.55	0.55
1:D:167:ASN:C	1:D:167:ASN:HD22	2.09	0.55
1:B:103:TYR:O	1:B:107:TYR:HB3	2.06	0.55
1:B:81:LEU:HG	1:B:85:LEU:CD2	2.36	0.55
1:D:180:GLN:HG2	1:D:184:ASN:HD21	1.72	0.55
1:A:114:GLU:O	1:A:118:LYS:HG3	2.07	0.55
1:A:153:ALA:O	1:A:157:ASN:HB2	2.07	0.55
1:D:103:TYR:HE1	1:E:165:GLU:HB3	1.72	0.55
1:B:144:ASP:OD2	1:B:144:ASP:N	2.33	0.54
1:B:15:PHE:CD1	1:B:50:PHE:HD2	2.25	0.54
1:B:116:MET:HA	1:B:116:MET:CE	2.37	0.53
1:D:12:LYS:HD2	1:D:57:GLU:OE1	2.07	0.53
1:A:170:GLU:OE2	1:A:170:GLU:HA	2.07	0.53
1:B:66:LYS:HA	1:B:69:GLN:HG2	1.89	0.53
1:B:91:TYR:O	1:B:93:TYR:N	2.42	0.53
1:B:44:LYS:HA	1:B:44:LYS:CE	2.38	0.53
1:D:43:PHE:O	1:D:44:LYS:C	2.47	0.53
1:A:76:ARG:HG3	1:A:139:GLU:OE2	2.09	0.53
1:D:180:GLN:CG	1:D:184:ASN:HD21	2.22	0.53
1:E:126:ALA:O	1:E:130:ILE:HG12	2.09	0.53
1:B:28:ILE:HD13	1:B:49:LEU:CD2	2.39	0.52
1:A:144:ASP:N	1:A:144:ASP:OD1	2.43	0.52
1:D:66:LYS:HD2	1:D:69:GLN:NE2	2.24	0.52
1:A:165:GLU:HG2	1:A:166:GLN:NE2	2.24	0.52
1:E:43:PHE:O	1:E:45:THR:N	2.44	0.51
1:A:18:ASN:O	1:A:22:ALA:HB3	2.10	0.51
1:B:18:ASN:ND2	1:B:23:THR:HG22	2.15	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:15:PHE:CE1	1:E:50:PHE:HD2	2.27	0.51
1:D:183:LEU:C	1:D:185:GLY:H	2.14	0.51
1:A:89:THR:HB	3:A:803:BER:H161	1.93	0.51
1:B:80:TYR:O	1:B:84:GLU:HG3	2.11	0.51
1:A:134:GLY:HA3	1:A:140:TRP:CZ2	2.47	0.50
1:E:25:THR:O	1:E:29:VAL:HG13	2.10	0.50
1:E:103:TYR:HD2	1:E:116:MET:HG2	1.76	0.50
1:B:51:LEU:HD21	1:B:115:LYS:HB2	1.93	0.50
1:D:6:LYS:HE2	1:D:32:SER:HA	1.93	0.50
1:B:2:ASN:OD1	1:B:3:LEU:N	2.38	0.50
1:B:11:ALA:O	1:B:12:LYS:C	2.50	0.50
1:B:29:VAL:CG2	1:B:39:LEU:HD21	2.38	0.50
1:B:123:TYR:OH	1:B:157:ASN:HB2	2.11	0.50
1:D:135:ASN:OD1	1:D:141:SER:HA	2.11	0.50
1:B:107:TYR:CE1	1:B:108:LYS:HG3	2.47	0.50
1:D:75:ASN:HB2	1:D:139:GLU:OE2	2.11	0.50
1:D:177:LYS:HE3	1:D:181:ILE:HD11	1.94	0.50
1:E:106:TYR:N	1:E:106:TYR:CD2	2.79	0.50
1:E:136:LEU:HD12	1:E:136:LEU:O	2.12	0.50
1:B:14:LEU:O	1:B:23:THR:HG21	2.11	0.49
1:D:145:VAL:O	1:D:145:VAL:HG12	2.11	0.49
1:B:30:LYS:CD	1:B:31:LEU:HD13	2.42	0.49
1:B:87:LEU:HD13	1:B:172:ILE:HD13	1.94	0.49
1:E:144:ASP:OD2	1:E:146:ASN:HB3	2.12	0.49
1:B:115:LYS:HE2	1:B:118:LYS:NZ	2.27	0.49
1:B:45:THR:CG2	1:B:48:ASN:HB2	2.26	0.49
1:A:166:GLN:N	1:A:166:GLN:NE2	2.59	0.49
1:B:106:TYR:O	1:B:112:ILE:CG1	2.60	0.49
1:D:20:TYR:O	1:D:22:ALA:N	2.45	0.49
1:B:30:LYS:HD2	1:B:31:LEU:HD13	1.93	0.49
1:B:185:GLY:HA3	1:A:181:ILE:O	2.13	0.49
1:B:35:SER:HB3	1:B:38:ASN:OD1	2.13	0.48
1:B:134:GLY:HA3	1:B:140:TRP:CE2	2.48	0.48
1:D:163:THR:O	1:D:171:ARG:NH1	2.45	0.48
1:A:127:TYR:O	1:A:128:HIS:C	2.49	0.48
1:E:177:LYS:O	1:E:181:ILE:HG12	2.13	0.48
1:E:17:LYS:HG2	1:E:18:ASN:ND2	2.28	0.48
1:E:76:ARG:HD2	1:E:183:LEU:HD23	1.95	0.48
1:B:89:THR:CG2	1:B:91:TYR:H	2.25	0.48
1:B:109:THR:OG1	1:B:112:ILE:HG12	2.13	0.48
1:D:177:LYS:O	1:D:181:ILE:CG1	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:VAL:O	1:A:133:GLU:HG2	2.12	0.48
1:E:116:MET:HA	1:E:116:MET:CE	2.44	0.48
1:E:167:ASN:HD21	1:E:169:ASN:HB3	1.78	0.48
1:E:10:VAL:HG11	1:E:31:LEU:HB2	1.95	0.48
1:A:173:LYS:O	1:A:173:LYS:HD3	2.14	0.48
1:D:46:LYS:O	1:D:49:LEU:HB3	2.14	0.48
1:E:12:LYS:HD2	2:E:399:SO4:O4	2.14	0.48
1:D:167:ASN:O	1:D:169:ASN:N	2.47	0.47
1:E:159:ILE:HB	1:E:175:MET:HE1	1.96	0.47
1:B:89:THR:CG2	1:B:90:GLU:H	2.20	0.47
1:B:116:MET:HA	1:B:116:MET:HE2	1.96	0.47
1:B:132:LYS:O	1:B:136:LEU:HD13	2.14	0.47
1:A:135:ASN:HD21	1:A:142:ILE:H	1.63	0.47
1:A:58:GLU:OE1	3:A:803:BER:H102	2.15	0.47
1:E:160:VAL:HG12	1:E:161:THR:N	2.29	0.47
1:B:18:ASN:HB3	1:B:22:ALA:CB	2.42	0.47
1:B:61:TRP:HB2	1:B:91:TYR:HB2	1.96	0.47
1:D:126:ALA:O	1:D:130:ILE:HG12	2.15	0.47
1:D:142:ILE:HG13	1:D:145:VAL:HG22	1.97	0.47
1:A:43:PHE:O	1:A:44:LYS:HB2	2.13	0.47
1:E:58:GLU:HG3	1:E:59:SER:N	2.30	0.47
1:D:97:ASN:HD21	1:E:100:ILE:HD12	1.80	0.47
1:B:137:ASN:HD22	1:B:139:GLU:HG3	1.79	0.47
1:D:74:THR:HG23	1:D:77:GLU:OE2	2.15	0.47
1:A:83:ASN:HB3	1:A:175:MET:HE2	1.97	0.47
1:A:156:VAL:O	1:A:160:VAL:HG23	2.14	0.47
1:E:119:LEU:O	1:E:121:ASN:N	2.48	0.47
1:B:110:ASN:C	1:B:112:ILE:N	2.65	0.47
1:D:103:TYR:CE1	1:E:165:GLU:HG2	2.50	0.47
1:E:83:ASN:ND2	1:E:127:TYR:OH	2.48	0.47
1:B:18:ASN:C	1:B:22:ALA:HB3	2.35	0.46
1:D:100:ILE:HD11	1:E:162:PHE:CE2	2.50	0.46
1:A:172:ILE:HG22	1:A:172:ILE:O	2.13	0.46
1:D:39:LEU:HD21	1:D:49:LEU:HD22	1.98	0.46
1:A:96:GLN:O	1:A:100:ILE:HG13	2.15	0.46
1:E:47:GLU:CG	1:E:106:TYR:CE1	2.99	0.46
1:E:60:LYS:HD3	1:E:91:TYR:CZ	2.50	0.46
1:D:10:VAL:CG2	1:D:32:SER:HB3	2.40	0.46
1:A:68:GLU:O	1:A:71:LYS:HG2	2.15	0.46
1:B:173:LYS:HA	1:B:173:LYS:HE2	1.98	0.46
1:B:43:PHE:O	1:B:44:LYS:O	2.33	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3:LEU:HD12	1:D:38:ASN:HD21	1.80	0.46
1:E:2:ASN:HB2	1:E:5:ASP:HB2	1.98	0.46
1:D:20:TYR:C	1:D:22:ALA:H	2.20	0.46
1:B:81:LEU:O	1:B:85:LEU:HD22	2.16	0.46
1:A:117:ASN:C	1:A:119:LEU:N	2.69	0.46
1:A:123:TYR:CE2	3:A:803:BER:H172	2.50	0.46
1:B:24:THR:HG22	1:B:27:GLU:CG	2.46	0.46
1:E:65:TRP:O	1:E:69:GLN:HG2	2.15	0.46
1:B:130:ILE:O	1:B:140:TRP:HZ2	1.99	0.45
1:A:38:ASN:OD1	1:A:42:HIS:HD2	2.00	0.45
1:B:53:ILE:O	1:B:57:GLU:HG2	2.15	0.45
1:B:144:ASP:HB2	1:B:147:ALA:HB3	1.98	0.45
1:B:39:LEU:O	1:B:43:PHE:O	2.34	0.45
1:B:45:THR:OG1	1:B:47:GLU:HG2	2.16	0.45
1:E:54:LEU:HD12	1:E:54:LEU:HA	1.76	0.45
1:A:135:ASN:HD22	1:A:135:ASN:HA	1.50	0.45
1:D:34:SER:OG	1:D:35:SER:N	2.47	0.45
1:D:131:PHE:CE2	1:D:148:VAL:HG12	2.52	0.45
1:D:20:TYR:C	1:D:22:ALA:N	2.69	0.45
1:A:83:ASN:ND2	1:A:156:VAL:HG21	2.31	0.45
1:D:103:TYR:C	1:D:105:GLU:H	2.19	0.45
1:A:165:GLU:N	1:A:165:GLU:OE2	2.49	0.45
1:D:25:THR:HG23	1:D:39:LEU:HD13	1.99	0.45
1:A:76:ARG:HB3	1:A:76:ARG:HH11	1.81	0.44
1:E:50:PHE:CE1	1:E:54:LEU:HD22	2.52	0.44
1:B:3:LEU:HD13	1:B:4:LYS:N	2.32	0.44
1:A:95:LEU:O	1:A:97:ASN:N	2.50	0.44
1:E:164:HIS:HA	1:E:171:ARG:HH12	1.82	0.44
1:D:125:ASP:O	1:D:128:HIS:N	2.50	0.44
1:D:3:LEU:HD11	1:D:34:SER:HB3	1.99	0.44
1:D:164:HIS:CG	1:D:164:HIS:O	2.71	0.44
1:E:57:GLU:HG3	1:E:95:LEU:HD12	1.99	0.44
1:B:45:THR:OG1	1:B:46:LYS:N	2.51	0.44
1:B:104:THR:HG21	1:A:100:ILE:HD13	1.99	0.44
1:B:144:ASP:O	1:B:147:ALA:N	2.45	0.44
1:E:92:TYR:CE2	1:E:123:TYR:CZ	3.06	0.44
1:E:132:LYS:O	1:E:135:ASN:HB2	2.18	0.44
1:D:180:GLN:O	1:D:181:ILE:C	2.55	0.44
1:E:140:TRP:CD1	1:E:142:ILE:HG12	2.53	0.44
1:A:109:THR:C	1:A:111:SER:N	2.71	0.44
1:A:167:ASN:ND2	1:A:167:ASN:C	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:LYS:HE2	1:B:118:LYS:HD3	2.00	0.44
1:B:137:ASN:HD22	1:B:139:GLU:CG	2.31	0.44
1:D:101:GLU:HG3	1:E:97:ASN:HD21	1.83	0.44
1:E:47:GLU:HG2	1:E:106:TYR:CE1	2.52	0.43
1:E:107:TYR:CD2	1:E:107:TYR:O	2.71	0.43
1:B:3:LEU:O	1:B:7:ILE:HG13	2.18	0.43
1:B:92:TYR:C	1:B:94:PRO:HD2	2.39	0.43
1:B:142:ILE:HD11	1:B:186:LEU:HD13	2.00	0.43
1:D:167:ASN:O	1:D:170:GLU:N	2.51	0.43
1:D:180:GLN:HG2	1:D:184:ASN:ND2	2.32	0.43
1:A:56:ILE:O	1:A:60:LYS:HG2	2.18	0.43
1:E:183:LEU:HD12	1:E:183:LEU:HA	1.77	0.43
1:D:15:PHE:CD1	1:D:50:PHE:HD2	2.36	0.43
1:A:105:GLU:C	1:A:107:TYR:H	2.21	0.43
1:E:162:PHE:O	1:E:163:THR:HG23	2.18	0.43
1:D:36:LYS:HG2	1:D:40:TYR:HE2	1.84	0.43
1:D:167:ASN:ND2	1:D:167:ASN:C	2.71	0.43
1:A:117:ASN:O	1:A:119:LEU:N	2.52	0.43
1:A:183:LEU:HD12	1:A:183:LEU:HA	1.68	0.43
1:B:126:ALA:O	1:B:130:ILE:HG13	2.19	0.43
1:A:8:LEU:HD13	1:A:56:ILE:CD1	2.49	0.43
1:A:95:LEU:HD23	1:A:95:LEU:HA	1.84	0.43
1:D:29:VAL:CG2	1:D:39:LEU:HD12	2.49	0.43
1:A:30:LYS:HD3	1:A:30:LYS:C	2.36	0.43
1:E:43:PHE:O	1:E:44:LYS:C	2.56	0.43
1:B:55:ASN:CG	1:B:119:LEU:HD21	2.39	0.43
1:D:133:GLU:O	1:D:137:ASN:HB2	2.18	0.42
1:E:27:GLU:O	1:E:28:ILE:C	2.57	0.42
1:E:65:TRP:HZ3	1:E:81:LEU:HD23	1.84	0.42
1:B:7:ILE:HG21	1:B:43:PHE:CE1	2.52	0.42
1:B:166:GLN:H	1:B:166:GLN:HG2	1.72	0.42
1:A:163:THR:HB	1:A:171:ARG:HD2	2.01	0.42
1:B:55:ASN:HD22	1:B:122:LYS:HE3	1.84	0.42
1:B:95:LEU:O	1:B:99:ILE:HG12	2.19	0.42
1:E:65:TRP:CZ3	1:E:81:LEU:HD23	2.54	0.42
1:E:106:TYR:H	1:E:106:TYR:HD2	1.67	0.42
1:B:168:ILE:O	1:B:172:ILE:HG12	2.19	0.42
1:E:68:GLU:HG2	1:E:85:LEU:HD21	2.00	0.42
1:E:168:ILE:H	1:E:168:ILE:HG13	1.68	0.42
1:B:18:ASN:HB2	1:B:23:THR:CG2	2.45	0.42
1:D:73:LYS:HE3	1:D:77:GLU:OE1	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LYS:O	1:B:69:GLN:HG2	2.19	0.42
1:B:165:GLU:OE2	1:B:166:GLN:NE2	2.53	0.42
1:E:43:PHE:N	1:E:43:PHE:CD1	2.86	0.42
1:D:135:ASN:HD21	1:D:142:ILE:HG12	1.85	0.42
1:A:156:VAL:HG22	1:A:175:MET:CE	2.50	0.41
1:B:31:LEU:HD12	1:B:31:LEU:HA	1.80	0.41
1:D:92:TYR:O	1:D:94:PRO:N	2.53	0.41
1:D:163:THR:O	1:D:171:ARG:HD2	2.20	0.41
1:E:121:ASN:HD22	1:E:121:ASN:HA	1.68	0.41
1:E:167:ASN:HD21	1:E:169:ASN:HB2	1.84	0.41
1:A:74:THR:O	1:A:77:GLU:HB2	2.20	0.41
1:D:11:ALA:O	1:D:12:LYS:C	2.59	0.41
1:A:68:GLU:O	1:A:68:GLU:HG3	2.19	0.41
1:A:94:PRO:O	1:A:95:LEU:C	2.58	0.41
1:A:8:LEU:HD23	1:A:8:LEU:HA	1.69	0.41
1:E:6:LYS:O	1:E:10:VAL:HG23	2.21	0.41
1:D:57:GLU:HB3	1:D:92:TYR:H	1.85	0.41
1:D:78:LYS:HB3	1:D:130:ILE:HD12	2.02	0.41
1:E:132:LYS:HD2	2:E:903:SO4:O3	2.21	0.41
1:B:16:ILE:HG13	1:B:95:LEU:HD23	2.01	0.41
1:B:186:LEU:HD23	1:B:186:LEU:HA	1.86	0.41
1:E:51:LEU:HD22	1:E:102:PHE:CZ	2.55	0.41
1:E:74:THR:H	1:E:77:GLU:HG3	1.86	0.41
1:B:127:TYR:O	1:B:131:PHE:CD2	2.74	0.41
1:D:52:GLU:O	1:D:55:ASN:N	2.50	0.41
1:D:124:ILE:HD12	1:D:153:ALA:CB	2.51	0.41
1:E:35:SER:OG	1:E:38:ASN:HB2	2.21	0.41
1:E:99:ILE:HD13	1:E:99:ILE:HG21	1.87	0.41
1:E:155:ALA:O	1:E:159:ILE:HG12	2.21	0.41
1:B:7:ILE:O	1:B:10:VAL:N	2.52	0.41
1:D:122:LYS:O	1:D:125:ASP:HB3	2.21	0.41
1:D:171:ARG:HG2	1:D:171:ARG:NH1	2.34	0.41
1:A:14:LEU:HG	1:A:23:THR:HG23	2.02	0.41
1:A:92:TYR:CD1	1:A:92:TYR:O	2.74	0.41
1:A:135:ASN:ND2	1:A:141:SER:HA	2.36	0.41
1:A:142:ILE:HD11	1:A:186:LEU:HD13	2.03	0.41
1:A:156:VAL:HG22	1:A:175:MET:HE1	2.03	0.41
1:E:56:ILE:O	1:E:57:GLU:C	2.59	0.41
1:E:61:TRP:HH2	1:E:85:LEU:HB3	1.86	0.41
1:E:100:ILE:O	1:E:101:GLU:C	2.59	0.41
1:B:4:LYS:HA	1:B:7:ILE:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:LYS:HE3	4:D:1002:HOH:O	2.20	0.40
1:E:38:ASN:O	1:E:42:HIS:CD2	2.74	0.40
1:E:74:THR:OG1	1:E:77:GLU:CG	2.70	0.40
1:E:111:SER:C	1:E:113:ASN:N	2.75	0.40
1:E:119:LEU:C	1:E:121:ASN:N	2.73	0.40
1:E:174:LEU:HA	1:E:174:LEU:HD23	1.75	0.40
1:E:74:THR:O	1:E:78:LYS:HG3	2.21	0.40
1:B:37:GLY:O	1:B:40:TYR:N	2.54	0.40
1:A:177:LYS:O	1:A:180:GLN:HB3	2.21	0.40
1:E:109:THR:O	1:E:110:ASN:HB2	2.22	0.40
1:B:44:LYS:N	1:B:48:ASN:HD22	2.20	0.40
1:D:178:PHE:HB2	1:E:151:ILE:CG2	2.51	0.40
1:E:110:ASN:HB3	1:E:111:SER:H	1.52	0.40
1:B:61:TRP:HB2	1:B:91:TYR:CB	2.51	0.40
1:B:163:THR:O	1:B:166:GLN:HG2	2.22	0.40
1:D:121:ASN:N	1:D:121:ASN:HD22	2.19	0.40
1:D:186:LEU:HD23	1:E:181:ILE:HG23	2.03	0.40
1:E:47:GLU:HG2	1:E:106:TYR:CD1	2.57	0.40
1:E:119:LEU:O	1:E:120:GLU:C	2.60	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	184/194 (95%)	155 (84%)	24 (13%)	5 (3%)	5 24
1	B	184/194 (95%)	148 (80%)	30 (16%)	6 (3%)	4 19
1	D	184/194 (95%)	154 (84%)	24 (13%)	6 (3%)	4 19
1	E	184/194 (95%)	153 (83%)	27 (15%)	4 (2%)	6 29
All	All	736/776 (95%)	610 (83%)	105 (14%)	21 (3%)	4 22

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	44	LYS
1	B	165	GLU
1	A	89	THR
1	A	144	ASP
1	E	44	LYS
1	E	166	GLN
1	B	28	ILE
1	B	92	TYR
1	B	145	VAL
1	A	96	GLN
1	A	164	HIS
1	B	27	GLU
1	D	44	LYS
1	D	93	TYR
1	D	104	THR
1	D	92	TYR
1	D	111	SER
1	E	143	ASN
1	E	120	GLU
1	D	168	ILE
1	A	145	VAL

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	169/177 (96%)	150 (89%)	19 (11%)	6 23
1	B	169/177 (96%)	148 (88%)	21 (12%)	4 19
1	D	169/177 (96%)	155 (92%)	14 (8%)	11 37
1	E	169/177 (96%)	155 (92%)	14 (8%)	11 37
All	All	676/708 (96%)	608 (90%)	68 (10%)	7 27

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	4	LYS
1	B	6	LYS
1	B	13	GLU
1	B	31	LEU
1	B	42	HIS
1	B	44	LYS
1	B	45	THR
1	B	85	LEU
1	B	88	THR
1	B	89	THR
1	B	95	LEU
1	B	110	ASN
1	B	116	MET
1	B	124	ILE
1	B	129	VAL
1	B	143	ASN
1	B	144	ASP
1	B	157	ASN
1	B	169	ASN
1	B	173	LYS
1	B	184	ASN
1	D	2	ASN
1	D	3	LEU
1	D	23	THR
1	D	24	THR
1	D	31	LEU
1	D	45	THR
1	D	47	GLU
1	D	97	ASN
1	D	107	TYR
1	D	108	LYS
1	D	129	VAL
1	D	149	SER
1	D	165	GLU
1	D	167	ASN
1	A	3	LEU
1	A	14	LEU
1	A	21	ASN
1	A	25	THR
1	A	49	LEU
1	A	63	GLU
1	A	90	GLU
1	A	91	TYR

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Mol	Chain	Res	Type
1	A	108	LYS
1	A	111	SER
1	A	116	MET
1	A	130	ILE
1	A	135	ASN
1	A	141	SER
1	A	146	ASN
1	A	163	THR
1	A	166	GLN
1	A	167	ASN
1	A	184	ASN
1	E	32	SER
1	E	38	ASN
1	E	39	LEU
1	E	51	LEU
1	E	63	GLU
1	E	67	LYS
1	E	76	ARG
1	E	77	GLU
1	E	106	TYR
1	E	111	SER
1	E	141	SER
1	E	165	GLU
1	E	166	GLN
1	E	183	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	18	ASN
1	B	42	HIS
1	B	48	ASN
1	B	55	ASN
1	B	62	GLN
1	B	69	GLN
1	B	83	ASN
1	B	96	GLN
1	B	110	ASN
1	B	137	ASN
1	B	146	ASN
1	B	154	ASN
1	B	157	ASN

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Mol	Chain	Res	Type
1	B	166	GLN
1	D	2	ASN
1	D	38	ASN
1	D	48	ASN
1	D	55	ASN
1	D	69	GLN
1	D	83	ASN
1	D	121	ASN
1	D	154	ASN
1	D	164	HIS
1	D	167	ASN
1	D	176	ASN
1	D	180	GLN
1	D	184	ASN
1	A	42	HIS
1	A	48	ASN
1	A	55	ASN
1	A	83	ASN
1	A	96	GLN
1	A	121	ASN
1	A	135	ASN
1	A	166	GLN
1	A	167	ASN
1	A	180	GLN
1	A	184	ASN
1	E	18	ASN
1	E	48	ASN
1	E	64	GLN
1	E	83	ASN
1	E	97	ASN
1	E	110	ASN
1	E	117	ASN
1	E	121	ASN
1	E	166	GLN
1	E	167	ASN
1	E	184	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)

23 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	A	801	-	4,4,4	0.22	0	6,6,6	0.10	0
2	SO4	A	299	-	4,4,4	0.23	0	6,6,6	0.19	0
2	SO4	E	901	-	4,4,4	0.32	0	6,6,6	0.19	0
2	SO4	B	1000	-	4,4,4	0.24	0	6,6,6	0.13	0
2	SO4	E	903	-	4,4,4	0.30	0	6,6,6	0.07	0
2	SO4	E	399	-	4,4,4	0.22	0	6,6,6	0.18	0
2	SO4	B	195	-	4,4,4	0.24	0	6,6,6	0.12	0
2	SO4	D	999	-	4,4,4	0.45	0	6,6,6	0.20	0
2	SO4	B	199	-	4,4,4	0.22	0	6,6,6	0.12	0
3	BER	A	803	-	29,29,29	3.94	21 (72%)	40,43,43	1.97	12 (30%)
2	SO4	A	699	-	4,4,4	0.28	0	6,6,6	0.10	0
2	SO4	A	800	-	4,4,4	0.29	0	6,6,6	0.11	0
2	SO4	D	1000	-	4,4,4	0.27	0	6,6,6	0.12	0
2	SO4	D	899	-	4,4,4	0.23	0	6,6,6	0.10	0
2	SO4	E	579	-	4,4,4	0.30	0	6,6,6	0.07	0
2	SO4	E	900	-	4,4,4	0.27	0	6,6,6	0.17	0
2	SO4	A	802	-	4,4,4	0.16	0	6,6,6	0.15	0
2	SO4	E	902	-	4,4,4	0.26	0	6,6,6	0.06	0
2	SO4	A	799	-	4,4,4	0.29	0	6,6,6	0.05	0
2	SO4	E	639	-	4,4,4	0.26	0	6,6,6	0.13	0
2	SO4	B	700	-	4,4,4	0.30	0	6,6,6	0.09	0
2	SO4	B	499	-	4,4,4	0.30	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SO4	D	900	-	4,4,4	0.26	0	6,6,6	0.12	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	BER	A	803	-	-	2/4/19/19	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	803	BER	C6-N1	9.21	1.45	1.33
3	A	803	BER	C10-C4	6.18	1.55	1.50
3	A	803	BER	C18-C15	5.93	1.48	1.38
3	A	803	BER	C2-C1	5.80	1.51	1.42
3	A	803	BER	C7-C10	5.70	1.58	1.51
3	A	803	BER	C1-N1	5.21	1.46	1.39
3	A	803	BER	O4-C18	4.83	1.44	1.37
3	A	803	BER	C9-C4	4.61	1.45	1.37
3	A	803	BER	C9-C14	4.55	1.47	1.38
3	A	803	BER	C5-C11	3.92	1.43	1.36
3	A	803	BER	C4-C2	3.90	1.51	1.42
3	A	803	BER	C13-C16	3.53	1.44	1.36
3	A	803	BER	C7-N1	3.35	1.55	1.48
3	A	803	BER	O3-C15	3.32	1.43	1.38
3	A	803	BER	C14-C11	3.31	1.47	1.39
3	A	803	BER	C3-C1	3.18	1.45	1.39
3	A	803	BER	C16-C18	3.02	1.45	1.39
3	A	803	BER	C12-C8	2.43	1.47	1.43
3	A	803	BER	C6-C12	2.34	1.45	1.41
3	A	803	BER	O1-C17	2.34	1.47	1.43
3	A	803	BER	C15-C12	2.25	1.48	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	803	BER	C10-C7-N1	5.63	116.60	109.47
3	A	803	BER	C19-O3-C15	3.67	124.34	114.88
3	A	803	BER	O4-C18-C15	3.64	121.59	116.49
3	A	803	BER	C20-O4-C18	3.60	122.97	117.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	803	BER	O1-C11-C5	3.49	130.90	128.11
3	A	803	BER	C7-N1-C1	3.09	121.32	117.91
3	A	803	BER	C7-C10-C4	2.79	112.98	109.79
3	A	803	BER	O2-C14-C9	2.72	131.49	127.85
3	A	803	BER	C16-C18-C15	-2.40	118.07	119.76
3	A	803	BER	O4-C18-C16	-2.35	120.34	124.37
3	A	803	BER	C13-C8-C3	-2.05	118.64	122.00
3	A	803	BER	C10-C4-C2	2.01	121.47	119.08

There are no chirality outliers.

All (2) torsion outliers are listed below:

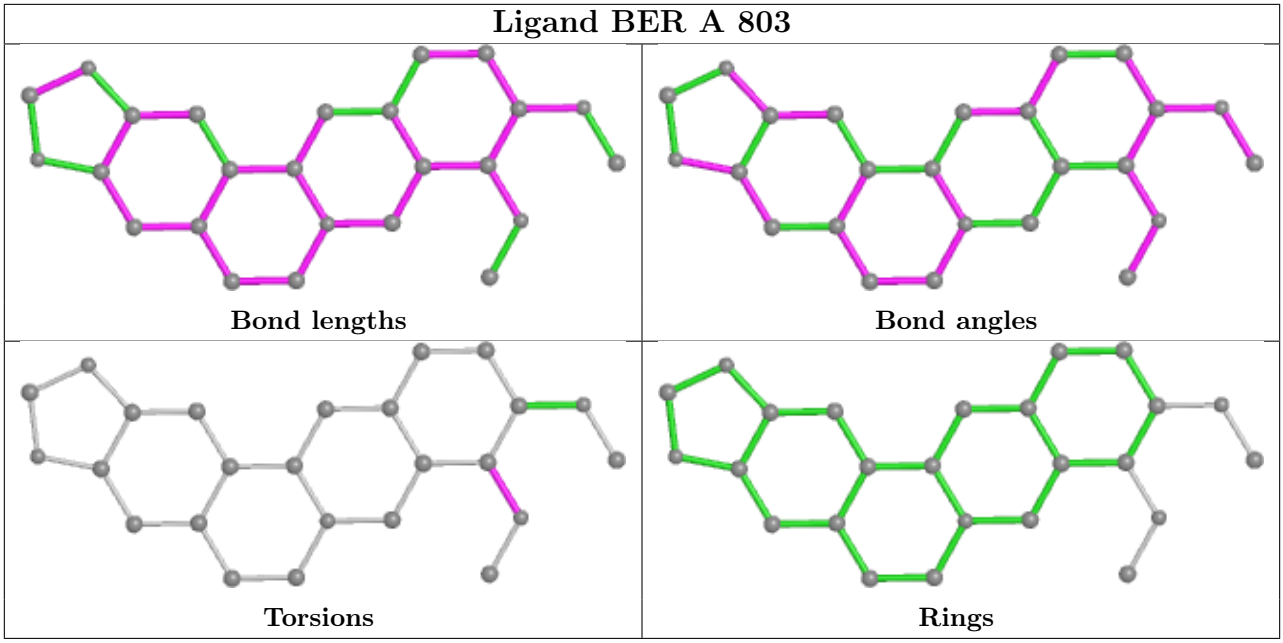
Mol	Chain	Res	Type	Atoms
3	A	803	BER	C18-C15-O3-C19
3	A	803	BER	C12-C15-O3-C19

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	SO4	1	0
2	E	903	SO4	1	0
2	E	399	SO4	1	0
3	A	803	BER	3	0
2	B	700	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.



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)

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	186/194 (95%)	-0.27	0 100 100	25, 46, 82, 103	0
1	B	186/194 (95%)	0.17	9 (4%) 30 18	29, 60, 115, 130	0
1	D	186/194 (95%)	-0.17	0 100 100	34, 57, 91, 105	0
1	E	186/194 (95%)	-0.17	1 (0%) 91 80	28, 48, 85, 97	0
All	All	744/776 (95%)	-0.11	10 (1%) 77 59	25, 52, 98, 130	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	41	TYR	6.1
1	B	42	HIS	5.3
1	B	40	TYR	4.8
1	E	123	TYR	4.4
1	B	43	PHE	3.0
1	B	34	SER	2.9
1	B	28	ILE	2.9
1	B	36	LYS	2.6
1	B	107	TYR	2.5
1	B	29	VAL	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)

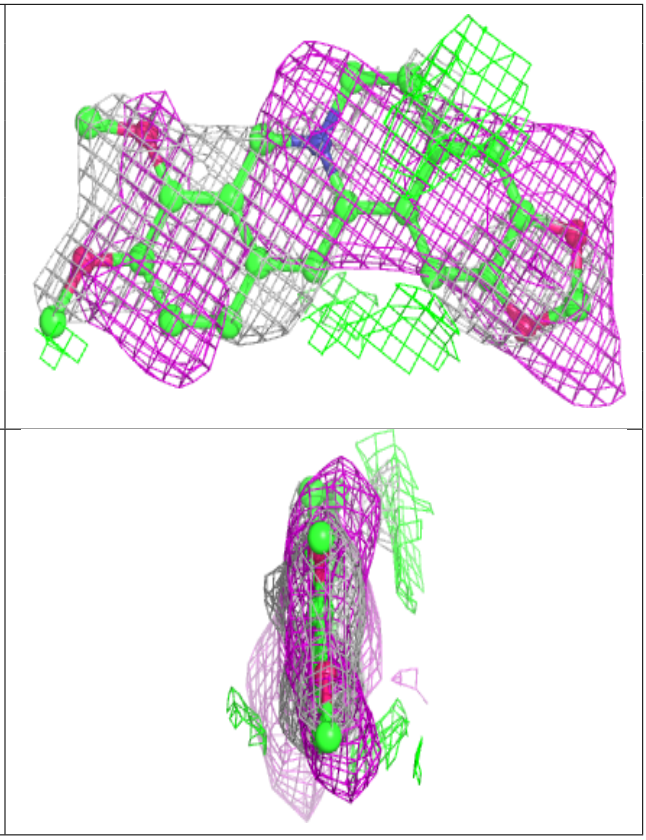
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SO4	D	1000	5/5	0.20	1.17	173,173,173,174	0
2	SO4	B	700	5/5	0.45	0.60	199,200,200,200	0
2	SO4	E	579	5/5	0.48	0.32	176,176,176,177	0
2	SO4	E	903	5/5	0.60	0.47	174,174,174,175	0
2	SO4	E	639	5/5	0.64	0.68	188,188,189,189	0
2	SO4	E	902	5/5	0.71	0.24	181,182,182,182	0
2	SO4	D	999	5/5	0.76	0.31	123,123,124,126	0
3	BER	A	803	25/25	0.77	0.62	47,53,61,62	0
2	SO4	E	901	5/5	0.79	0.31	145,145,145,146	0
2	SO4	D	899	5/5	0.79	0.23	147,148,148,149	0
2	SO4	A	299	5/5	0.79	0.23	148,148,148,149	0
2	SO4	B	1000	5/5	0.81	0.22	176,176,176,176	0
2	SO4	A	800	5/5	0.83	0.20	134,135,135,135	0
2	SO4	A	699	5/5	0.84	0.18	143,143,143,143	0
2	SO4	B	499	5/5	0.84	0.47	157,157,157,157	0
2	SO4	E	399	5/5	0.86	0.77	194,194,195,195	0
2	SO4	A	801	5/5	0.86	0.39	161,161,161,161	0
2	SO4	B	195	5/5	0.86	0.30	177,177,177,178	0
2	SO4	D	900	5/5	0.88	0.16	135,135,136,136	0
2	SO4	B	199	5/5	0.89	0.57	146,147,147,147	0
2	SO4	E	900	5/5	0.90	0.21	138,138,138,138	0
2	SO4	A	799	5/5	0.90	0.51	134,134,134,134	0
2	SO4	A	802	5/5	0.95	0.30	152,153,154,154	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around BER A 803:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.